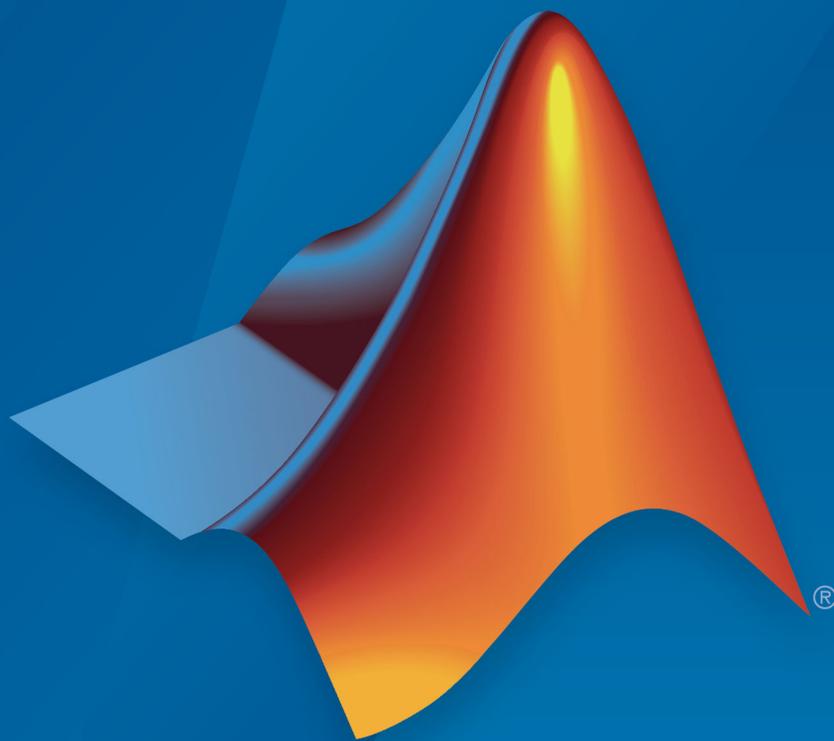


# System Identification Toolbox™

Reference

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# MATLAB® & SIMULINK®

R2018b



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**1** | Functions - Alphabetical List

**2** | Blocks – Alphabetical List



# Functions - Alphabetical List

---

## absorbDelay

Replace time delays by poles at  $z = 0$  or phase shift

### Syntax

```
sysnd = absorbDelay(sysd)  
[sysnd,G] = absorbDelay(sysd)
```

### Description

`sysnd = absorbDelay(sysd)` absorbs all time delays of the dynamic system model `sysd` into the system dynamics or the frequency response data.

For discrete-time models (other than frequency response data models), a delay of  $k$  sampling periods is replaced by  $k$  poles at  $z = 0$ . For continuous-time models (other than frequency response data models), time delays have no exact representation with a finite number of poles and zeros. Therefore, use `pade` to compute a rational approximation of the time delay.

For frequency response data models in both continuous and discrete time, `absorbDelay` absorbs all time delays into the frequency response data as a phase shift.

`[sysnd,G] = absorbDelay(sysd)` returns the matrix  $G$  that maps the initial states of the ss model `sysd` to the initial states of the `sysnd`.

### Examples

#### Absorb Time Delay into System Dynamics

Create a discrete-time transfer function that has a time delay.

```
z = tf('z',-1);  
sysd = (-0.4*z -0.1)/(z^2 + 1.05*z + 0.08);  
sysd.InputDelay = 3
```

```
sysd =
```

$$z^{-3} * \frac{-0.4 z - 0.1}{z^2 + 1.05 z + 0.08}$$

```
Sample time: unspecified
Discrete-time transfer function.
```

The display of `sysd` represents the `InputDelay` as a factor of  $z^{-3}$ , separate from the system poles that appear in the transfer function denominator.

Absorb the time delay into the system dynamics as poles at  $z = 0$ .

```
sysnd = absorbDelay(sysd)
```

```
sysnd =
```

$$\frac{-0.4 z - 0.1}{z^5 + 1.05 z^4 + 0.08 z^3}$$

```
Sample time: unspecified
Discrete-time transfer function.
```

The display of `sysnd` shows that the factor of  $z^{-3}$  has been absorbed as additional poles in the denominator.

Verify that `sysnd` has no input delay.

```
sysnd.InputDelay
```

```
ans = 0
```

### Convert Leading Structural Zeros of Polynomial Model to Regular Coefficients

Create a discrete-time polynomial model.

```
m = idpoly(1,[0 0 0 2 3]);
```

Convert `m` to a transfer function model.

```
sys = tf(m)
```

```
sys =
```

```
z^(-2) * (2 z^-1 + 3 z^-2)
```

```
Sample time: unspecified  
Discrete-time transfer function.
```

The numerator of the transfer function, `sys`, is `[0 2 3]` and the transport delay, `sys.IODelay`, is 2. This is because the value of the B polynomial, `m.B`, has 3 leading zeros. The first fixed zero shows lack of feedthrough in the model. The two zeros after that are treated as input-output delays.

Use `absorbDelay` to treat the leading zeros as regular B coefficients.

```
m2 = absorbDelay(m);  
sys2 = tf(m2)
```

```
sys2 =
```

```
2 z^-3 + 3 z^-4
```

```
Sample time: unspecified  
Discrete-time transfer function.
```

The numerator of `sys2` is `[0 0 0 2 3]` and transport delay is 0. The model `m2` treats the leading zeros as regular coefficients by freeing their values. `m2.Structure.B.Free(2:3)` is TRUE while `m.Structure.B.Free(2:3)` is FALSE.

## See Also

`hasdelay` | `pade` | `totaldelay`

**Introduced in R2012a**

# advice

Analysis and recommendations for data or estimated linear models

## Syntax

```
advice(data)  
advice(model,data)
```

## Description

`advice(data)` displays the following information about the data in the MATLAB® Command Window:

- What are the excitation levels of the signals and how does this affect the model orders? See also `pexcit`.
- Does it make sense to remove constant offsets and linear trends from the data? See also `detrend`.
- Is there an indication of output feedback in the data? See also `feedback`.
- Would a nonlinear ARX model perform better than a linear ARX model?

`advice(model,data)` displays the following information about the estimated linear model in the MATLAB Command Window:

- Does the model capture essential dynamics of the system and the disturbance characteristics?
- Is the model order higher than necessary?
- Is there potential output feedback in the validation data?

## Input Arguments

### **data**

Specify data as an `iddata` object.

### **model**

Specify `model` as an `idtf`, `idgrey`, `idpoly`, `idproc`, or `idss` model object.

### **See Also**

`detrend` | `feedback` | `iddata` | `pexcit`

**Introduced before R2006a**

# addreg

Add custom regressors to nonlinear ARX model

## Syntax

```
m = addreg(model,regressors)  
m = addreg(model,regressors,output)
```

## Description

*m* = addreg(*model*,*regressors*) adds custom regressors to a nonlinear ARX model by appending the CustomRegressors *model* property. *model* and *m* are idnlarx objects. For single-output models, *regressors* is an object array of regressors you create using customreg or polyreg, or a cell array of character vectors. For multiple-output models, *regressors* is 1-by-ny cell array of customreg objects or 1-by-ny cell array of cell arrays of character vectors. addreg adds each element of the ny cells to the corresponding *model* output channel. If *regressors* is a single regressor, addreg adds this regressor to all output channels.

*m* = addreg(*model*,*regressors*,*output*) adds regressors *regressors* to specific output channels *output* of a multiple-output model. *output* is a scalar integer or vector of integers, where each integer is the index of a model output channel. Specify several pairs of *regressors* and *output* values to add different regressor variables to the corresponding output channels.

## Examples

### Add Regressors to a Nonlinear ARX Model

Create nonlinear ARX model with standard regressors.

```
m1 = idnlarx([4 2 1], 'wavenet', 'nlr', [1:3]);
```

Create model with additional custom regressors, specified as a cell array of character vectors.

```
m2 = addreg(m1,{'y1(t-2)^2';'u1(t)*y1(t-7)'});
```

List all standard and custom regressors of m2.

```
getreg(m2)
```

Regressors:

```
y1(t-1)
y1(t-2)
y1(t-3)
y1(t-4)
u1(t-1)
u1(t-2)
y1(t-2)^2
u1(t)*y1(t-7)
```

## Add Regressors to a Nonlinear ARX Model as customreg Objects

Create nonlinear ARX model with standard regressors.

```
m1 = idnlarx([4 2 1], 'wavenet', 'nlr', [1:3]);
```

Create customreg objects.

```
r1 = customreg(@(x)x^2,{'y1'},2)
```

Custom Regressor:

```
Expression: y1(t-2)^2
Function: @(x)x^2
Arguments: {'y1'}
Delays: 2
Vectorized: 0
TimeVariable: 't'
```

```
r2 = customreg(@(x,y)x*y,{'u1','y1'},[0 7])
```

Custom Regressor:

```
Expression: u1(t)*y1(t-7)
Function: @(x,y)x*y
Arguments: {'u1' 'y1'}
```

```
Delays: [0 7]
Vectorized: 0
TimeVariable: 't'
```

Create a model based on m1 with custom regressors.

```
m2 = addreg(m1, [r1 r2]);
```

List all standard and custom regressors of m2.

```
getreg(m2)
```

Regressors:

```
y1(t-1)
y1(t-2)
y1(t-3)
y1(t-4)
u1(t-1)
u1(t-2)
y1(t-2)^2
u1(t)*y1(t-7)
```

## See Also

[customreg](#) | [getreg](#) | [nlarx](#) | [polyreg](#)

## Topics

“Identifying Nonlinear ARX Models”

**Introduced in R2007a**

## aic

Akaike's Information Criterion for estimated model

### Syntax

```
value = aic(model)
value = aic(model1,...,modeln)
value = aic( ____,measure)
```

### Description

`value = aic(model)` returns the normalized "Akaike's Information Criterion (AIC)" on page 1-14 value for the estimated model.

`value = aic(model1,...,modeln)` returns the normalized AIC values for multiple estimated models.

`value = aic( ____,measure)` specifies the type of AIC.

### Examples

#### Compute Normalized Akaike's Information Criterion of Estimated Model

Estimate a transfer function model.

```
load iddata1 z1;
np = 2;
sys = tfest(z1,np);
```

Compute the normalized Akaike's Information Criterion value.

```
value = aic(sys)
value = 0.5453
```

The value is also computed during model estimation. Alternatively, use the `Report` property of the model to access this value.

```
sys.Report.Fit.nAIC  
ans = 0.5453
```

### Compute Akaike's Information Criterion Metrics of Estimated Model

Estimate a transfer function model.

```
load iddata1 z1;  
np = 2;  
sys = tfest(z1,np);
```

Compute the normalized Akaike's Information Criterion (AIC) value. This syntax is equivalent to `aic_raw = aic(sys)`.

```
aic_raw = aic(sys, 'nAIC')  
aic_raw = 0.5453
```

Compute the raw AIC value.

```
aic_raw = aic(sys, 'aic')  
aic_raw = 1.0150e+03
```

Compute the sample-size corrected AIC value.

```
aic_c = aic(sys, 'AICc')  
aic_c = 1.0153e+03
```

Compute the Bayesian Information Criteria (BIC) value.

```
bic = aic(sys, 'BIC')  
bic = 1.0372e+03
```

These values are also computed during model estimation. Alternatively, use the `Report.Fit` property of the model to access these values.

```
sys.Report.Fit
```

```
ans = struct with fields:
  FitPercent: 70.7720
  LossFcn: 1.6575
  MSE: 1.6575
  FPE: 1.7252
  AIC: 1.0150e+03
  AICc: 1.0153e+03
  nAIC: 0.5453
  BIC: 1.0372e+03
```

## Pick Model with Optimal Tradeoff Between Accuracy and Complexity Using AICc Criterion

Estimate multiple Output-Error (OE) models and use the small sample-size corrected Akaike's Information Criterion (AICc) value to pick the one with optimal tradeoff between accuracy and complexity.

Load the estimation data.

```
load iddata2
```

Specify model orders varying in 1:4 range.

```
nf = 1:4;
nb = 1:4;
nk = 0:4;
```

Estimate OE models with all possible combinations of chosen order ranges.

```
NN = struc(nf,nb,nk);
models = cell(size(NN,1),1);
for ct = 1:size(NN,1)
    models{ct} = oe(z2, NN(ct,:));
end
```

Compute the small sample-size corrected AIC values for the models, and return the smallest value.

```
V = aic(models{:}, 'AICc');
[Vmin,I] = min(V);
```

Return the optimal model that has the smallest AICc value.

```
models{I}
```

```
ans =
```

```
Discrete-time OE model:  $y(t) = [B(z)/F(z)]u(t) + e(t)$ 
```

```
  B(z) = 1.067 z-2
```

```
  F(z) = 1 - 1.824 z-1 + 1.195 z-2 - 0.2307 z-3
```

```
Sample time: 0.1 seconds
```

```
Parameterization:
```

```
  Polynomial orders:  nb=1  nf=3  nk=2
```

```
  Number of free coefficients: 4
```

```
  Use "polydata", "getpvec", "getcov" for parameters and their uncertainties.
```

```
Status:
```

```
Estimated using OE on time domain data "z2".
```

```
Fit to estimation data: 86.53%
```

```
FPE: 0.9809, MSE: 0.9615
```

## Input Arguments

### **model** — Identified model

```
idtf | idgrey | idpoly | idproc | idss | idnlarx | idnlhw | idnlgrey
```

Identified model, specified as one of the following model objects:

- `idtf`
- `idgrey`
- `idpoly`
- `idproc`
- `idss`
- `idnlarx`, except nonlinear ARX model that includes a binary-tree or neural network nonlinearity estimator
- `idnlhw`
- `idnlgrey`

**measure — Type of AIC**`'nAIC' (default) | 'aic' | 'AICc' | 'BIC'`

Type of AIC, specified as one of the following values:

- `'nAIC'` — Normalized AIC
- `'aic'` — Raw AIC
- `'AICc'` — Small sample-size corrected AIC
- `'BIC'` — Bayesian Information Criteria

See “Akaike's Information Criterion (AIC)” on page 1-14 for more information.

## Output Arguments

**value — Value of quality metric**`scalar | vector`

Value of the quality measure, returned as a scalar or vector. For multiple models, `value` is a row vector where `value(k)` corresponds to the *k*th estimated model `modelk`.

## Definitions

### Akaike's Information Criterion (AIC)

Akaike's Information Criterion (AIC) provides a measure of model quality obtained by simulating the situation where the model is tested on a different data set. After computing several different models, you can compare them using this criterion. According to Akaike's theory, the most accurate model has the smallest AIC. If you use the same data set for both model estimation and validation, the fit always improves as you increase the model order and, therefore, the flexibility of the model structure.

Akaike's Information Criterion (AIC) includes the following quality metrics:

- Raw AIC, defined as:

$$AIC = N * \log \left( \det \left( \frac{1}{N} \sum_1^N \varepsilon(t, \theta_N) (\varepsilon(t, \theta_N))^T \right) \right) + 2n_p + N * (n_y * (\log(2\pi) + 1))$$

where:

- $N$  is the number of values in the estimation data set
- $\varepsilon(t)$  is a  $n_y$ -by-1 vector of prediction errors
- $\theta_N$  represents the estimated parameters
- $n_p$  is the number of estimated parameters
- $n_y$  is the number of model outputs
- Small sample-size corrected AIC, defined as:

$$AICc = AIC + 2n_p * \frac{n_p + 1}{N - n_p - 1}$$

- Normalized AIC, defined as:

$$nAIC = \log \left( \det \left( \frac{1}{N} \sum_1^N \varepsilon(t, \theta_N) (\varepsilon(t, \theta_N))^T \right) \right) + \frac{2n_p}{N}$$

- Bayesian Information Criteria, defined as:

$$BIC = N * \log \left( \det \left( \frac{1}{N} \sum_1^N \varepsilon(t, \theta_N) (\varepsilon(t, \theta_N))^T \right) \right) + N * (n_y * \log(2\pi) + 1) + n_p * \log(N)$$

## Tips

- The software computes and stores all types of Akaike's Information Criterion metrics during model estimation. If you want to access these values, see the `Report.Fit` property of the model.

## References

- [1] Ljung, L. *System Identification: Theory for the User*, Upper Saddle River, NJ, Prentice-Hall PTR, 1999. See sections about the statistical framework for parameter estimation and maximum likelihood method and comparing model structures.

## **See Also**

fpe | goodnessOfFit

## **Topics**

“Loss Function and Model Quality Metrics”

**Introduced before R2006a**

# append

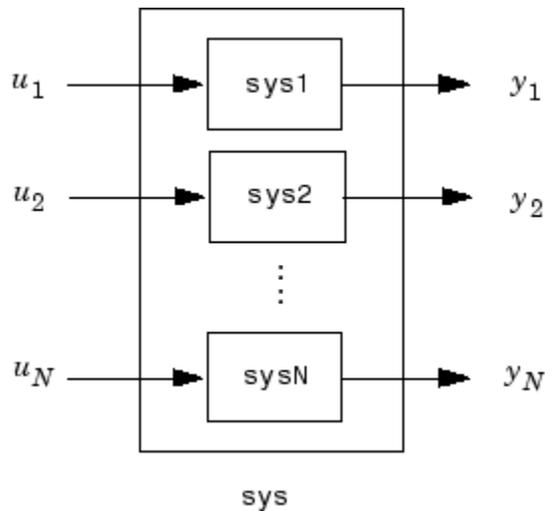
Group models by appending their inputs and outputs

## Syntax

```
sys = append(sys1,sys2,...,sysN)
```

## Description

`sys = append(sys1,sys2,...,sysN)` appends the inputs and outputs of the models `sys1,...,sysN` to form the augmented model `sys` depicted below.



For systems with transfer functions  $H_1(s), \dots, H_N(s)$ , the resulting system `sys` has the block-diagonal transfer function

$$\begin{bmatrix} H_1(s) & 0 & \dots & 0 \\ 0 & H_2(s) & \dots & \vdots \\ \vdots & \vdots & \ddots & 0 \\ 0 & \dots & 0 & H_N(s) \end{bmatrix}$$

For state-space models `sys1` and `sys2` with data  $(A_1, B_1, C_1, D_1)$  and  $(A_2, B_2, C_2, D_2)$ , `append(sys1, sys2)` produces the following state-space model:

$$\begin{aligned} \begin{bmatrix} \dot{x}_1 \\ \dot{x}_2 \end{bmatrix} &= \begin{bmatrix} A_1 & 0 \\ 0 & A_2 \end{bmatrix} \begin{bmatrix} x_1 \\ x_2 \end{bmatrix} + \begin{bmatrix} B_1 & 0 \\ 0 & B_2 \end{bmatrix} \begin{bmatrix} u_1 \\ u_2 \end{bmatrix} \\ \begin{bmatrix} y_1 \\ y_2 \end{bmatrix} &= \begin{bmatrix} C_1 & 0 \\ 0 & C_2 \end{bmatrix} \begin{bmatrix} x_1 \\ x_2 \end{bmatrix} + \begin{bmatrix} D_1 & 0 \\ 0 & D_2 \end{bmatrix} \begin{bmatrix} u_1 \\ u_2 \end{bmatrix} \end{aligned}$$

## Arguments

The input arguments `sys1, ..., sysN` can be model objects of any type. Regular matrices are also accepted as a representation of static gains, but there should be at least one model in the input list. The models should be either all continuous, or all discrete with the same sample time. When appending models of different types, the resulting type is determined by the precedence rules (see “Rules That Determine Model Type” (Control System Toolbox) for details).

There is no limitation on the number of inputs.

## Examples

### Append Inputs and Outputs of Models

Create a SISO transfer function.

```
sys1 = tf(1,[1 0]);  
size(sys1)
```

Transfer function with 1 outputs and 1 inputs.

Create a SISO continuous-time state-space model.

```
sys2 = ss(1,2,3,4);  
size(sys2)
```

State-space model with 1 outputs, 1 inputs, and 1 states.

Append the inputs and outputs of `sys1`, a SISO static gain system, and `sys2`. The resulting model should be a 3-input, 3-output state-space model.

```
sys = append(sys1,10,sys2)
```

```
sys =
```

```
A =
```

	x1	x2
x1	0	0
x2	0	1

```
B =
```

	u1	u2	u3
x1	1	0	0
x2	0	0	2

```
C =
```

	x1	x2
y1	1	0
y2	0	0
y3	0	3

```
D =
```

	u1	u2	u3
y1	0	0	0
y2	0	10	0
y3	0	0	4

Continuous-time state-space model.

```
size(sys)
```

State-space model with 3 outputs, 3 inputs, and 2 states.

## See Also

[connect](#) | [feedback](#) | [parallel](#) | [series](#)

**Introduced in R2012a**

## ar

Estimate parameters of AR model for scalar time series

### Syntax

```
m = ar(y,n)
[m,refl] = ar(y,n,approach>window)
m= ar(y,n,Name,Value)
m= ar(y,n, ____,opt)
```

### Description

---

**Note** Use for scalar time series only. For multivariate data, use `arx`.

---

`m = ar(y,n)` returns an `idpoly` model `m`.

`[m,refl] = ar(y,n,approach>window)` returns an `idpoly` model `m` and the variable `refl`. For the two lattice-based approaches, 'burg' and 'gl', `refl` stores the reflection coefficients in the first row, and the corresponding loss function values in the second row. The first column of `refl` is the zeroth-order model, and the (2,1) element of `refl` is the norm of the time series itself.

`m= ar(y,n,Name,Value)` specifies model structure attributes using one or more `Name,Value` pair arguments.

`m= ar(y,n, ____,opt)` specifies the estimations options using `opt`.

### Input Arguments

**y**

`iddata` object that contains the time-series data (one output channel).

## **Default:**

### **n**

Scalar that specifies the order of the model you want to estimate (the number of  $A$  parameters in the AR model).

### **approach**

Algorithm for computing the least squares AR model, specified as one of the following values:

- 'burg': Burg's lattice-based method. Solves the lattice filter equations using the harmonic mean of forward and backward squared prediction errors.
- 'fb': (Default) Forward-backward approach. Minimizes the sum of a least-squares criterion for a forward model, and the analogous criterion for a time-reversed model.
- 'gl': Geometric lattice approach. Similar to Burg's method, but uses the geometric mean instead of the harmonic mean during minimization.
- 'ls': Least-squares approach. Minimizes the standard sum of squared forward-prediction errors.
- 'yw': Yule-Walker approach. Solves the Yule-Walker equations, formed from sample covariances.

### **window**

Use of information about the data outside the measured time interval (past and future values), specified as one of the following values:

- 'now': (Default) No windowing. This value is the default except when the **approach** argument is 'yw'. Only measured data is used to form regression vectors. The summation in the criteria starts at the sample index equal to  $n+1$ .
- 'pow': Postwindowing. Missing end values are replaced with zeros and the summation is extended to time  $N+n$  ( $N$  is the number of observations).
- 'ppw': Pre- and postwindowing. Used in the Yule-Walker approach.
- 'prw': Prewindowing. Missing past values are replaced with zeros so that the summation in the criteria can start at time equal to zero.

### **opt**

Estimation options.

opt is an options set that specifies the following:

- data offsets
- covariance handling
- estimation approach
- estimation window

Use `arOptions` to create the options set.

## Name-Value Pair Arguments

Specify optional comma-separated pairs of `Name`, `Value` arguments. `Name` is the argument name and `Value` is the corresponding value. `Name` must appear inside quotes. You can specify several name and value pair arguments in any order as `Name1, Value1, ..., NameN, ValueN`.

### **Ts**

Positive scalar that specifies the sample time. Use when you specify `Y` as double vector rather than an `IDDATA` object.

### **Default:**

### **IntegrateNoise**

Boolean value that specifies whether the noise source contains an integrator or not. Use it

to create "ARI" structure models:  $Ay = \frac{e}{(1-z^{-1})}$

**Default:** false

## Output Arguments

**m**

An `idpoly` model.

## refl

An 2-by-2 array. The first row stores the reflection coefficients, and the second row stores the corresponding loss function values. The first column of `refl` is the zeroth-order model, and the (2,1) element of `refl` is the norm of the time series itself.

## Examples

Given a sinusoidal signal with noise, compare the spectral estimates of Burg's method with those found from the forward-backward approach and no-windowing method on a Bode plot.

```
y = sin([1:300]') + 0.5*randn(300,1);
y = iddata(y);
mb = ar(y,4,'burg');
mfb = ar(y,4);
bode(mb,mfb)
```

Estimate an ARI model.

```
load iddata9 z9
Ts = z9.Ts;
y = cumsum(z9.y);
model = ar(y, 4, 'ls', 'Ts', Ts, 'IntegrateNoise', true)
compare(y,model,5) % 5 step ahead prediction
```

Use option set to choose 'ls' estimation approach and to specify that covariance matrix should not be estimated.

```
y = rand(100,1);
opt = arOptions('Approach', 'ls', 'EstimateCovariance', false);
model = ar(y, N, opt);
```

## Algorithms

The AR model structure is given by the following equation:

$$A(q)y(t) = e(t)$$

AR model parameters are estimated using variants of the least-squares method. The following table summarizes the common names for methods with a specific combination of approach and window argument values.

Method	Approach and Windowing
Modified Covariance Method	(Default) Forward-backward approach and no windowing.
Correlation Method	Yule-Walker approach, which corresponds to least squares plus pre- and postwindowing.
Covariance Method	Least squares approach with no windowing. <code>arx</code> uses this routine.

## References

Marple, Jr., S.L., *Digital Spectral Analysis with Applications*, Prentice Hall, Englewood Cliffs, 1987, Chapter 8.

## See Also

`arOptions` | `arx` | `etfe` | `forecast` | `idpoly` | `ivar` | `pem` | `spa`

**Introduced before R2006a**

## armax

Estimate parameters of ARMAX model using time-domain data

### Syntax

```
sys = armax(data,[na nb nc nk])  
sys = armax(data,[na nb nc nk],Name,Value)  
  
sys = armax(data,init_sys)  
  
sys = armax(data, ____,opt)
```

### Description

`sys = armax(data,[na nb nc nk])` returns an `idpoly` model, `sys`, with estimated parameters and covariance (parameter uncertainties). Estimates the parameters using the prediction-error method and specified polynomial orders.

`sys = armax(data,[na nb nc nk],Name,Value)` returns an `idpoly` model, `sys`, with additional options specified by one or more `Name,Value` pair arguments.

`sys = armax(data,init_sys)` estimates a polynomial model using a discrete-time linear model `init_sys` to configure the initial parameterization.

`sys = armax(data, ____,opt)` specifies estimation options using the option set `opt`.

### Input Arguments

#### **data** — Time-domain estimation data

`iddata` object

Time-domain estimation data, specified as an `iddata` object. You cannot use frequency-domain data for estimating ARMAX models.

**[na nb nc nk] — Polynomial orders**

1-by-4 vector of positive integers | 1-by-4 vector of matrices

Polynomial orders of an “ARMAX Model” on page 1-39, specified as a 1-by-4 vector, [na nb nc nk].

For a model with  $N_y$  outputs and  $N_u$  inputs:

- na is the order of the polynomial  $A(q)$ , specified as an  $N_y$ -by- $N_y$  matrix of nonnegative integers.
- nb is the order of the polynomial  $B(q) + 1$ , specified as an  $N_y$ -by- $N_u$  matrix of nonnegative integers.
- nc is the order of the polynomial  $C(q)$ , specified as a column vector of nonnegative integers of length  $N_y$ .
- nk is the input-output delay expressed as fixed leading zeros of the  $B$  polynomial.

Specify nk as an  $N_y$ -by- $N_u$  matrix of nonnegative integers.

**init\_sys — System for configuring initial parametrization**

discrete-time linear model

System for configuring initial parametrization of `sys`, specified as a discrete-time linear model. You obtain `init_sys` by either performing an estimation using measured data or by direct construction using commands such as `idpoly` and `idss`.

If `init_sys` is an ARMAX model, `armax` uses the parameter values of `init_sys` as the initial guess for estimation. To configure initial guesses and constraints for  $A(q)$ ,  $B(q)$ , and  $C(q)$ , use the `Structure` property of `init_sys`. For example:

- To specify an initial guess for the  $A(q)$  term of `init_sys`, set `init_sys.Structure.A.Value` as the initial guess.
- To specify constraints for the  $B(q)$  term of `init_sys`:
  - set `init_sys.Structure.B.Minimum` to the minimum  $B(q)$  coefficient values.
  - set `init_sys.Structure.B.Maximum` to the maximum  $B(q)$  coefficient values.
  - set `init_sys.Structure.B.Free` to indicate which  $B(q)$  coefficients are free for estimation.

If `init_sys` is not a polynomial model of ARMAX structure, the software first converts `init_sys` to an ARMAX model. `armax` uses the parameters of the resulting model as the initial guess for estimating `sys`.

If `opt` is not specified, and `init_sys` was obtained by estimation, then the estimation options from `init_sys.Report.OptionsUsed` are used.

## **opt — Estimation options**

`armaxOptions` option set

Estimation options for ARMAX model identification, specified as an `armaxOptions` option set.

## **Name-Value Pair Arguments**

Specify optional comma-separated pairs of `Name`, `Value` arguments. `Name` is the argument name and `Value` is the corresponding value. `Name` must appear inside quotes. You can specify several name and value pair arguments in any order as `Name1, Value1, ..., NameN, ValueN`.

## **InputDelay — Input delays**

0 (default) | scalar | vector of positive integers

Input delays, specified as the comma-separated pair consisting of 'InputDelay' and one of the following:

- `Nu`-by-1 vector, where `Nu` is the number of inputs — Each entry is a numerical value representing the input delay for the corresponding input channel. Specify input delays as integer multiples of the sample time `Ts`. For example, `InputDelay = 3` means a delay of three sampling periods.
- Scalar value — The same delay is applied to all input channels.

## **IODelay — Transport delays**

0 (default) | scalar | matrix

Transport delays for each input/output pair, specified as the comma-separated pair consisting of 'IODelay' and one of the following:

- `Ny`-by-`Nu` matrix, where `Ny` is the number of outputs and `Nu` is the number of inputs — Each entry is an integer value representing the transport delay for the corresponding input/output pair. Specify transport delays as integers denoting delay of a multiple of the sample time `Ts`.
- Scalar value — The same delay is applied to all input/output pairs.

'IODElay' is useful as a replacement for the  $nk$  order. You can factor out  $\max(nk-1, 0)$  lags as the IODElay value. For  $nk > 1$ ,  $\text{armax}(na, nb, nk)$  is equivalent to  $\text{armax}(na, nb, 1, \text{'IODElay'}, nk-1)$ .

### **IntegrateNoise — Addition of integrators in noise channel**

`false(Ny, 1)`, where  $Ny$  is the number of outputs (default) | logical vector

Addition of integrators in noise channel, specified as the comma-separated pair consisting of 'IntegrateNoise' and a logical vector of length  $Ny$ , where  $Ny$  is the number of outputs.

Setting IntegrateNoise to true for a particular output results in the model:

$$A(q)y(t) = B(q)u(t - nk) + \frac{C(q)}{1 - q^{-1}}e(t)$$

Where,  $\frac{1}{1 - q^{-1}}$  is the integrator in the noise channel,  $e(t)$ .

Use IntegrateNoise to create ARIMA or ARIMAX models.

For example,

```
load iddata9 z9;
z9.y = cumsum(z9.y); % integrated data
sys = armax(z9, [4 1], 'IntegrateNoise', true);
compare(z9, sys, 10) % 10-step ahead prediction
```

## **Output Arguments**

### **sys — ARMAX model**

idpoly object

ARMAX model that fits the given estimation data, returned as a discrete-time idpoly object. This model is created using the specified model orders, delays, and estimation options.

Information about the estimation results and options used is stored in the Report property of the model. Report has the following fields:

<b>Report Field</b>	<b>Description</b>																		
Status	Summary of the model status, which indicates whether the model was created by construction or obtained by estimation.																		
Method	Estimation command used.																		
InitialCondition	<p>Handling of initial conditions during model estimation, returned as one of the following values:</p> <ul style="list-style-type: none"> <li>• 'zero' — The initial conditions were set to zero.</li> <li>• 'estimate' — The initial conditions were treated as independent estimation parameters.</li> <li>• 'backcast' — The initial conditions were estimated using the best least squares fit.</li> </ul> <p>This field is especially useful to view how the initial conditions were handled when the InitialCondition option in the estimation option set is 'auto'.</p>																		
Fit	<p>Quantitative assessment of the estimation, returned as a structure. See “Loss Function and Model Quality Metrics” for more information on these quality metrics. The structure has the following fields:</p> <table border="1"> <thead> <tr> <th><b>Field</b></th> <th><b>Description</b></th> </tr> </thead> <tbody> <tr> <td>FitPercent</td> <td>Normalized root mean squared error (NRMSE) measure of how well the response of the model fits the estimation data, expressed as a percentage.</td> </tr> <tr> <td>LossFcn</td> <td>Value of the loss function when the estimation completes.</td> </tr> <tr> <td>MSE</td> <td>Mean squared error (MSE) measure of how well the response of the model fits the estimation data.</td> </tr> <tr> <td>FPE</td> <td>Final prediction error for the model.</td> </tr> <tr> <td>AIC</td> <td>Raw Akaike Information Criteria (AIC) measure of model quality.</td> </tr> <tr> <td>AICc</td> <td>Small sample-size corrected AIC.</td> </tr> <tr> <td>nAIC</td> <td>Normalized AIC.</td> </tr> <tr> <td>BIC</td> <td>Bayesian Information Criteria (BIC).</td> </tr> </tbody> </table>	<b>Field</b>	<b>Description</b>	FitPercent	Normalized root mean squared error (NRMSE) measure of how well the response of the model fits the estimation data, expressed as a percentage.	LossFcn	Value of the loss function when the estimation completes.	MSE	Mean squared error (MSE) measure of how well the response of the model fits the estimation data.	FPE	Final prediction error for the model.	AIC	Raw Akaike Information Criteria (AIC) measure of model quality.	AICc	Small sample-size corrected AIC.	nAIC	Normalized AIC.	BIC	Bayesian Information Criteria (BIC).
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nAIC	Normalized AIC.																		
BIC	Bayesian Information Criteria (BIC).																		
Parameters	Estimated values of model parameters.																		

Report Field	Description																
OptionsUsed	Option set used for estimation. If no custom options were configured, this is a set of default options. See <code>armaxOptions</code> for more information.																
RandState	State of the random number stream at the start of estimation. Empty, <code>[]</code> , if randomization was not used during estimation. For more information, see <code>rng</code> in the MATLAB documentation.																
DataUsed	Attributes of the data used for estimation, returned as a structure with the following fields: <table border="1" data-bbox="387 604 1332 1295"> <thead> <tr> <th>Field</th> <th>Description</th> </tr> </thead> <tbody> <tr> <td>Name</td> <td>Name of the data set.</td> </tr> <tr> <td>Type</td> <td>Data type.</td> </tr> <tr> <td>Length</td> <td>Number of data samples.</td> </tr> <tr> <td>Ts</td> <td>Sample time.</td> </tr> <tr> <td>InterSample</td> <td>Input intersample behavior, returned as one of the following values: <ul style="list-style-type: none"> <li>'zoh' — Zero-order hold maintains a piecewise-constant input signal between samples.</li> <li>'foh' — First-order hold maintains a piecewise-linear input signal between samples.</li> <li>'bl' — Band-limited behavior specifies that the continuous-time input signal has zero power above the Nyquist frequency.</li> </ul> </td> </tr> <tr> <td>InputOffset</td> <td>Offset removed from time-domain input data during estimation. For nonlinear models, it is <code>[]</code>.</td> </tr> <tr> <td>OutputOffset</td> <td>Offset removed from time-domain output data during estimation. For nonlinear models, it is <code>[]</code>.</td> </tr> </tbody> </table>	Field	Description	Name	Name of the data set.	Type	Data type.	Length	Number of data samples.	Ts	Sample time.	InterSample	Input intersample behavior, returned as one of the following values: <ul style="list-style-type: none"> <li>'zoh' — Zero-order hold maintains a piecewise-constant input signal between samples.</li> <li>'foh' — First-order hold maintains a piecewise-linear input signal between samples.</li> <li>'bl' — Band-limited behavior specifies that the continuous-time input signal has zero power above the Nyquist frequency.</li> </ul>	InputOffset	Offset removed from time-domain input data during estimation. For nonlinear models, it is <code>[]</code> .	OutputOffset	Offset removed from time-domain output data during estimation. For nonlinear models, it is <code>[]</code> .
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OutputOffset	Offset removed from time-domain output data during estimation. For nonlinear models, it is <code>[]</code> .																

Report Field	Description	
Termination	Termination conditions for the iterative search used for prediction error minimization. Structure with the following fields:	
	Field	Description
	WhyStop	Reason for terminating the numerical search.
	Iterations	Number of search iterations performed by the estimation algorithm.
	FirstOrderOptimality	$\infty$ -norm of the gradient search vector when the search algorithm terminates.
	FcnCount	Number of times the objective function was called.
	UpdateNorm	Norm of the gradient search vector in the last iteration. Omitted when the search method is 'lsqnonlin' or 'fmincon'.
	LastImprovement	Criterion improvement in the last iteration, expressed as a percentage. Omitted when the search method is 'lsqnonlin' or 'fmincon'.
	Algorithm	Algorithm used by 'lsqnonlin' or 'fmincon' search method. Omitted when other search methods are used.
For estimation methods that do not require numerical search optimization, the Termination field is omitted.		

For more information on using Report, see “Estimation Report”.

## Examples

### Estimate ARMAX Model Using Regularization

Estimate a regularized ARMAX model by converting a regularized ARX model.

Load data.

```
load regularizationExampleData.mat m0simdata;
```

Estimate an unregularized ARMAX model of order 15.

```
m1 = armax(m0simdata(1:150),[30 30 30 1]);
```

Estimate a regularized ARMAX model by determining Lambda value by trial and error.

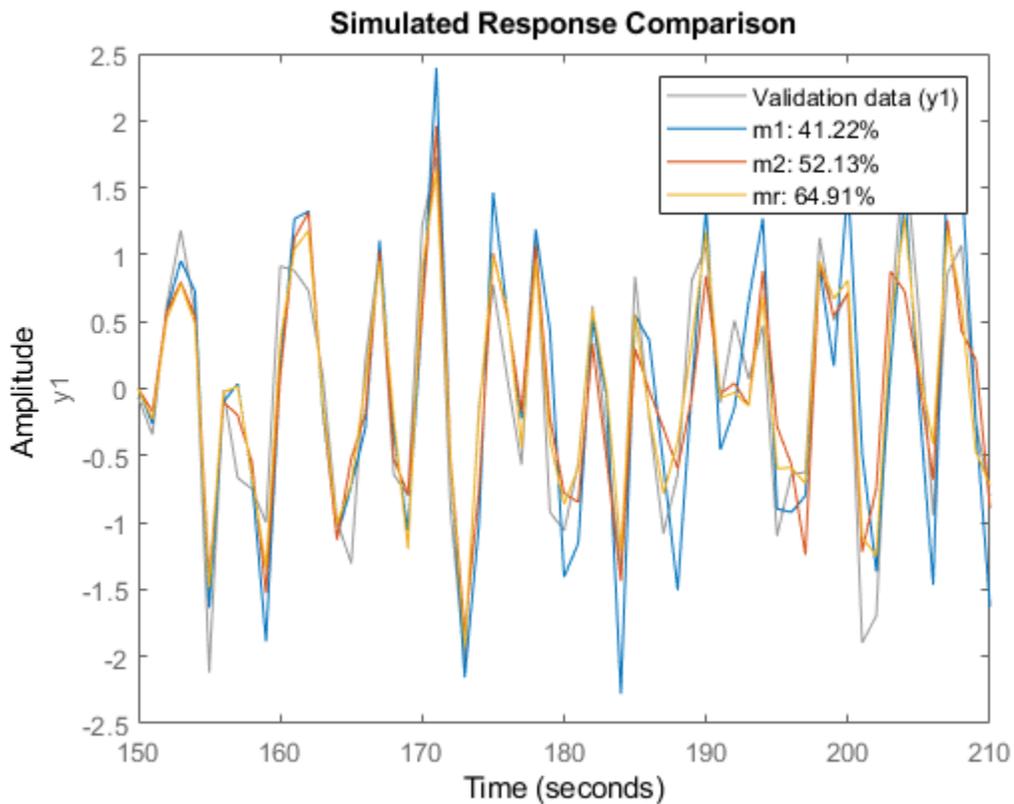
```
opt = armaxOptions;  
opt.Regularization.Lambda = 1;  
m2 = armax(m0simdata(1:150),[30 30 30 1],opt);
```

Obtain a lower-order ARMAX model by converting a regularized ARX model followed by order reduction.

```
opt1 = arxOptions;  
[L,R] = arxRegul(m0simdata(1:150),[30 30 1]);  
opt1.Regularization.Lambda = L;  
opt1.Regularization.R = R;  
m0 = arx(m0simdata(1:150),[30 30 1],opt1);  
mr = idpoly(balred(idss(m0),7));
```

Compare the model outputs against data.

```
opt2 = compareOptions('InitialCondition','z');  
compare(m0simdata(150:end),m1,m2,mr,opt2);
```



### Specify Estimation Options

Estimate an ARMAX model from measured data and specify the estimation options.

Estimate an ARMAX model with simulation focus, using 'lm' as the search method and maximum number of search iterations set to 10.

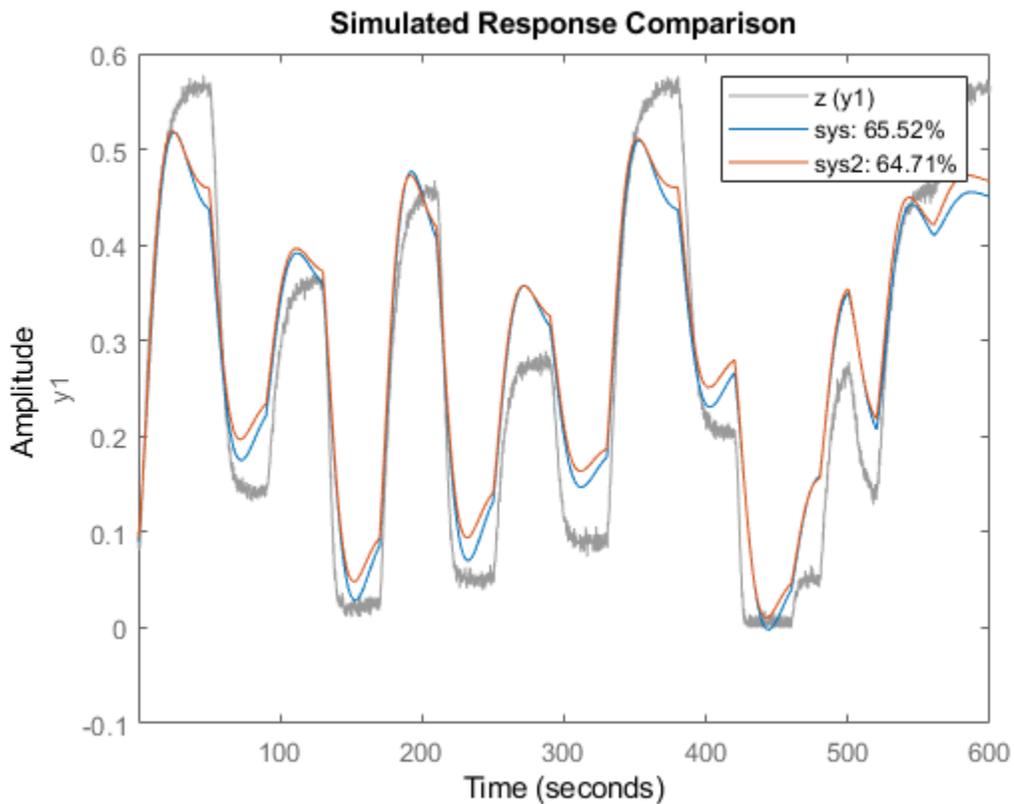
```
load twotankdata;  
z = iddata(y,u,0.2);  
opt = armaxOptions;  
opt.Focus = 'simulation';
```

```
opt.SearchMethod = 'lm';  
opt.SearchOptions.MaxIterations = 10;  
opt.Display = 'on';  
sys = armax(z,[2 2 2 1],opt);
```

The termination conditions for measured component of the model shown in the progress viewer is that the maximum number of iterations were reached.

To improve results, re-estimate the model using a greater value for MaxIterations, or continue iterations on the previously estimated model as follows:

```
sys2 = armax(z,sys);  
compare(z,sys,sys2)
```

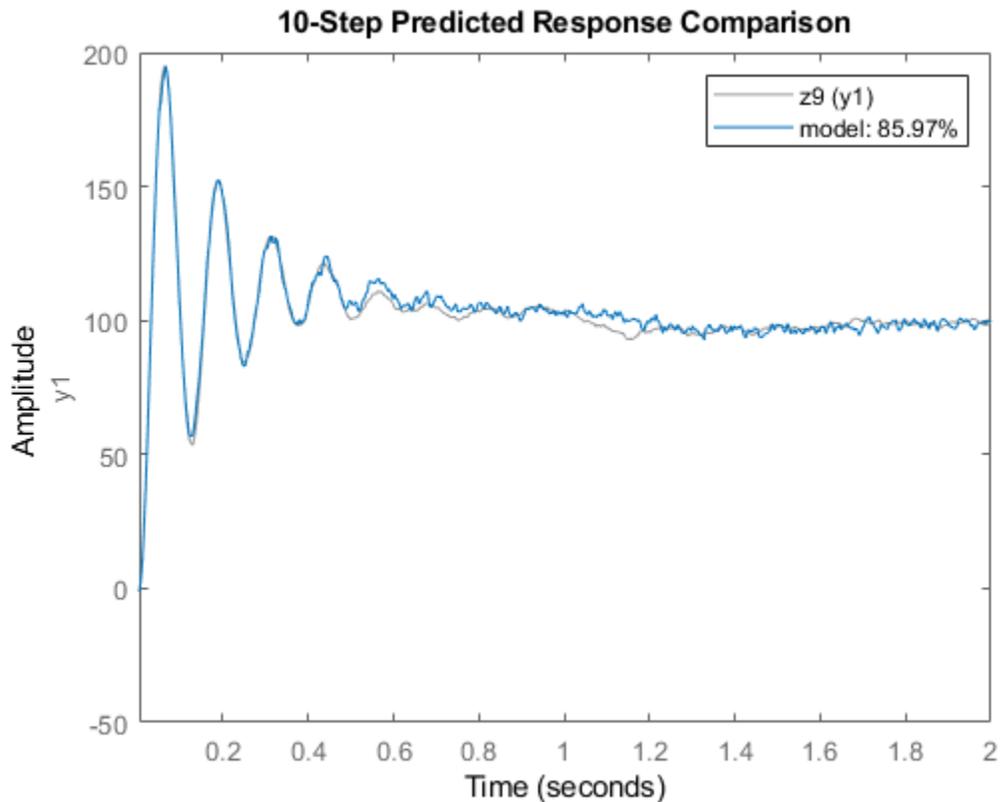


where `sys2` refines the parameters of `sys` to improve the fit to data.

### Estimate ARIMA Model

Estimate a 4th order ARIMA model for univariate time-series data.

```
load iddata9;  
z9.y = cumsum(z9.y); % integrated data  
model = armax(z9,[4 1], 'IntegrateNoise',true);  
compare(z9,model,10) % 10-step ahead prediction
```



## Estimate ARMAX Models Iteratively

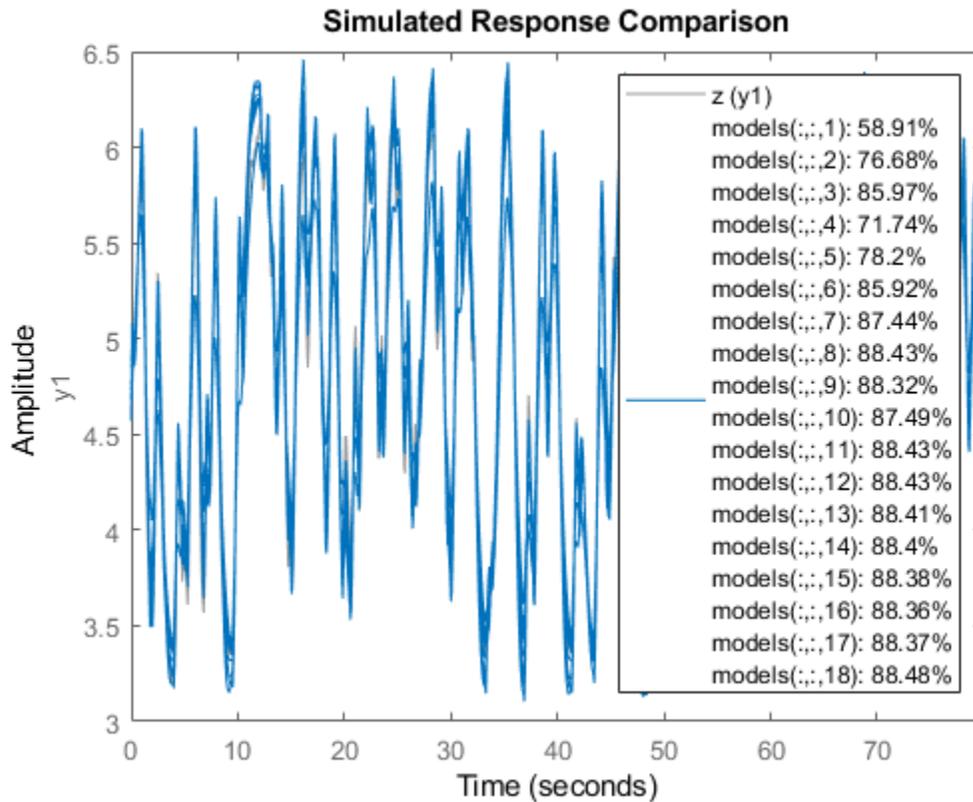
Estimate ARMAX models of varying orders iteratively from measured data.

Estimate ARMAX models of orders varying between 1 and 4 for dryer data.

```
load dryer2;
z = iddata(y2,u2,0.08, 'Tstart',0);
na = 2:4;
nc = 1:2;
nk = 0:2;
models = cell(1,18);
ct = 1;
for i = 1:3
    na_ = na(i);
    nb_ = na_;
    for j = 1:2
        nc_ = nc(j);
        for k = 1:3
            nk_ = nk(k);
            models{ct} = armax(z,[na_ nb_ nc_ nk_]);
            ct = ct+1;
        end
    end
end
end
```

Stack the estimated models and compare their simulated responses to estimation data z.

```
models = stack(1,models{:});
compare(z,models)
```



### Estimate ARMAX Model Using State-Space Model as Initial Guess

Load the estimation data.

```
load iddata2 z2
```

Estimate a state-space model of order 3 from the estimation data.

```
sys0 = n4sid(z2,3);
```

Estimate an ARMAX model using the previously estimated state-space model as an initial guess.

```
sys = armax(z2,sys0);
```

## Definitions

### ARMAX Model

The ARMAX model structure is:

$$y(t) + a_1 y(t-1) + \dots + a_{n_a} y(t-n_a) = b_1 u(t-n_k) + \dots + b_{n_b} u(t-n_k-n_b+1) + c_1 e(t-1) + \dots + c_{n_c} e(t-n_c) + e(t)$$

A more compact way to write the difference equation is:

$$A(q)y(t) = B(q)u(t-n_k) + C(q)e(t)$$

where,

- $y(t)$  — Output at time  $t$ .
- $n_a$  — Number of poles.
- $n_b$  — Number of zeroes plus 1.
- $n_c$  — Number of  $C$  coefficients.
- $n_k$  — Number of input samples that occur before the input affects the output, also called the *dead time* in the system.
- $y(t-1)\dots y(t-n_a)$  — Previous outputs on which the current output depends.
- $u(t-n_k)\dots u(t-n_k-n_b+1)$  — Previous and delayed inputs on which the current output depends.
- $e(t-1)\dots e(t-n_c)$  — White-noise disturbance value.

The parameters  $n_a$ ,  $n_b$ , and  $n_c$  are the orders of the ARMAX model, and  $n_k$  is the delay.  $q$  is the delay operator. Specifically,

$$A(q) = 1 + a_1q^{-1} + \dots + a_{n_a}q^{-n_a}$$

$$B(q) = b_1 + b_2q^{-1} + \dots + b_{n_b}q^{-n_b+1}$$

$$C(q) = 1 + c_1q^{-1} + \dots + c_{n_c}q^{-n_c}$$

If `data` is a time series that has no input channels and one output channel, then `armax` calculates an ARMA model for the time series

$$A(q)y(t) = C(q)e(t)$$

In this case,

`orders = [na nc]`

## ARIMAX Model

An ARIMAX model structure is similar to ARMAX, except that it contains an integrator in the noise source  $e(t)$ :

$$A(q)y(t) = B(q)u(t - nk) + \frac{C(q)}{(1 - q^{-1})}e(t)$$

If there are no inputs, the structure reduces to an ARIMA model:

$$A(q)y(t) = \frac{C(q)}{(1 - q^{-1})}e(t)$$

## Tips

- Use the `IntegrateNoise` property to add integrators to the noise source.

## Algorithms

An iterative search algorithm minimizes a robustified quadratic prediction error criterion. The iterations are terminated when any of the following is true:

- Maximum number of iterations is reached.
- Expected improvement is less than the specified tolerance.
- Lower value of the criterion cannot be found.

You can get information about the stopping criteria using `sys.Report.Termination`.

Use the `armaxOptions` option set to create and configure options affecting the estimation results. In particular, set the search algorithm attributes, such as `MaxIterations` and `Tolerance`, using the 'SearchOptions' property.

When you do not specify initial parameter values for the iterative search as an initial model, they are constructed in a special four-stage LS-IV algorithm.

The cutoff value for the robustification is based on the `Advanced.ErrorThreshold` estimation option and on the estimated standard deviation of the residuals from the initial parameter estimate. It is not recalculated during the minimization. By default, no robustification is performed; the default value of `ErrorThreshold` option is 0.

To ensure that only models corresponding to stable predictors are tested, the algorithm performs a stability test of the predictor. Generally, both  $C(q)$  and  $F(q)$  (if applicable) must have all zeros inside the unit circle.

Minimization information is displayed on the screen when the estimation option 'Display' is 'On' or 'Full'. With 'Display' = 'Full', both the current and the previous parameter estimates are displayed in column-vector form, listing parameters in alphabetical order. Also, the values of the criterion function (cost) are given and the Gauss-Newton vector and its norm are also displayed. With 'Display' = 'On', only the criterion values are displayed.

## Alternatives

`armax` does not support continuous-time model estimation. Use `tfest` to estimate a continuous-time transfer function model, or `ssest` to estimate a continuous-time state-space model.

armax supports only time-domain data. For frequency-domain data, use `oe` to estimate an Output-Error (OE) model.

## References

Ljung, L. *System Identification: Theory for the User*, Upper Saddle River, NJ: Prentice-Hall PTR, 1999. See chapter about computing the estimate.

## Extended Capabilities

### Automatic Parallel Support

Accelerate code by automatically running computation in parallel using Parallel Computing Toolbox™.

Parallel computing support is available for estimation using the `lsqnonlin` search method (requires Optimization Toolbox™). To enable parallel computing, use `armaxOptions`, set `SearchMethod` to `'lsqnonlin'`, and set `SearchOptions.Advanced.UseParallel` to `true`.

For example:

```
opt = armaxOptions;  
opt.SearchMethod = 'lsqnonlin';  
opt.SearchOptions.Advanced.UseParallel = true;
```

### See Also

[aic](#) | [armaxOptions](#) | [arx](#) | [bj](#) | [forecast](#) | [fpe](#) | [goodnessofFit](#) | [iddata](#) | [idfrd](#) | [idpoly](#) | [oe](#) | [polyest](#) | [srest](#) | [tfest](#)

### Topics

“Loss Function and Model Quality Metrics”

“Regularized Estimates of Model Parameters”

“Estimation Report”

**Introduced before R2006a**

## armaxOptions

Option set for armax

### Syntax

```
opt = armaxOptions  
opt = armaxOptions(Name,Value)
```

### Description

`opt = armaxOptions` creates the default options set for armax.

`opt = armaxOptions(Name,Value)` creates an option set with the options specified by one or more Name,Value pair arguments.

### Input Arguments

#### Name-Value Pair Arguments

Specify optional comma-separated pairs of Name,Value arguments. Name is the argument name and Value is the corresponding value. Name must appear inside quotes. You can specify several name and value pair arguments in any order as Name1,Value1,...,NameN,ValueN.

#### InitialCondition — Handling of initial conditions

'auto' (default) | 'zero' | 'estimate' | 'backcast'

Handling of initial conditions during estimation, specified as one of the following values:

- 'zero' — The initial conditions are set to zero.
- 'estimate' — The initial conditions are treated as independent estimation parameters.
- 'backcast' — The initial conditions are estimated using the best least squares fit.

- 'auto' — The software chooses the method to handle initial conditions based on the estimation data.

### Focus — Error to be minimized

'prediction' (default) | 'simulation'

Error to be minimized in the loss function during estimation, specified as the comma-separated pair consisting of 'Focus' and one of the following values:

- 'prediction' — The one-step ahead prediction error between measured and predicted outputs is minimized during estimation. As a result, the estimation focuses on producing a good predictor model.
- 'simulation' — The simulation error between measured and simulated outputs is minimized during estimation. As a result, the estimation focuses on making a good fit for simulation of model response with the current inputs.

The Focus option can be interpreted as a weighting filter in the loss function. For more information, see “Loss Function and Model Quality Metrics”.

### WeightingFilter — Weighting prefilter

[] (default) | vector | matrix | cell array | linear system

Weighting prefilter applied to the loss function to be minimized during estimation. To understand the effect of WeightingFilter on the loss function, see “Loss Function and Model Quality Metrics”.

Specify WeightingFilter as one of the following values:

- [] — No weighting prefilter is used.
- Passbands — Specify a row vector or matrix containing frequency values that define desired passbands. You select a frequency band where the fit between estimated model and estimation data is optimized. For example, [w<sub>l</sub>, w<sub>h</sub>], where w<sub>l</sub> and w<sub>h</sub> represent lower and upper limits of a passband. For a matrix with several rows defining frequency passbands, [w<sub>1l</sub>, w<sub>1h</sub>; w<sub>2l</sub>, w<sub>2h</sub>; w<sub>3l</sub>, w<sub>3h</sub>; . . . ], the estimation algorithm uses the union of the frequency ranges to define the estimation passband.

Passbands are expressed in rad/TimeUnit for time-domain data and in FrequencyUnit for frequency-domain data, where TimeUnit and FrequencyUnit are the time and frequency units of the estimation data.

- SISO filter — Specify a single-input-single-output (SISO) linear filter in one of the following ways:

- A SISO LTI model
- $\{A, B, C, D\}$  format, which specifies the state-space matrices of a filter with the same sample time as estimation data.
- $\{\text{numerator}, \text{denominator}\}$  format, which specifies the numerator and denominator of the filter as a transfer function with same sample time as estimation data.

This option calculates the weighting function as a product of the filter and the input spectrum to estimate the transfer function.

### **EnforceStability** — Control whether to enforce stability of model

`false` (default) | `true`

Control whether to enforce stability of estimated model, specified as the comma-separated pair consisting of 'EnforceStability' and either `true` or `false`.

This option is not available for multi-output models with a non-diagonal  $A$  polynomial array.

Data Types: `logical`

### **EstimateCovariance** — Control whether to generate parameter covariance data

`true` (default) | `false`

Controls whether parameter covariance data is generated, specified as `true` or `false`.

If `EstimateCovariance` is `true`, then use `getcov` to fetch the covariance matrix from the estimated model.

### **Display** — Specify whether to display the estimation progress

`'off'` (default) | `'on'`

Specify whether to display the estimation progress, specified as one of the following values:

- `'on'` — Information on model structure and estimation results are displayed in a progress-viewer window.
- `'off'` — No progress or results information is displayed.

### **InputOffset** — Removal of offset from time-domain input data during estimation

`[]` (default) | vector of positive integers | matrix

Removal of offset from time-domain input data during estimation, specified as the comma-separated pair consisting of 'InputOffset' and one of the following:

- A column vector of positive integers of length  $Nu$ , where  $Nu$  is the number of inputs.
- `[]` — Indicates no offset.
- $Nu$ -by- $Ne$  matrix — For multi-experiment data, specify InputOffset as an  $Nu$ -by- $Ne$  matrix.  $Nu$  is the number of inputs, and  $Ne$  is the number of experiments.

Each entry specified by InputOffset is subtracted from the corresponding input data.

### **OutputOffset — Removal of offset from time-domain output data during estimation**

`[]` (default) | vector | matrix

Removal of offset from time-domain output data during estimation, specified as the comma-separated pair consisting of 'OutputOffset' and one of the following:

- A column vector of length  $Ny$ , where  $Ny$  is the number of outputs.
- `[]` — Indicates no offset.
- $Ny$ -by- $Ne$  matrix — For multi-experiment data, specify OutputOffset as a  $Ny$ -by- $Ne$  matrix.  $Ny$  is the number of outputs, and  $Ne$  is the number of experiments.

Each entry specified by OutputOffset is subtracted from the corresponding output data.

### **Regularization — Options for regularized estimation of model parameters** structure

Options for regularized estimation of model parameters. For more information on regularization, see “Regularized Estimates of Model Parameters”.

Regularization is a structure with the following fields:

- **Lambda** — Constant that determines the bias versus variance tradeoff.

Specify a positive scalar to add the regularization term to the estimation cost.

The default value of zero implies no regularization.

**Default:** 0

- **R** — Weighting matrix.

Specify a vector of nonnegative numbers or a square positive semi-definite matrix. The length must be equal to the number of free parameters of the model.

For black-box models, using the default value is recommended. For structured and grey-box models, you can also specify a vector of `np` positive numbers such that each entry denotes the confidence in the value of the associated parameter.

The default value of 1 implies a value of `eye(npfree)`, where `npfree` is the number of free parameters.

**Default:** 1

- **Nominal** — The nominal value towards which the free parameters are pulled during estimation.

The default value of zero implies that the parameter values are pulled towards zero. If you are refining a model, you can set the value to `'model'` to pull the parameters towards the parameter values of the initial model. The initial parameter values must be finite for this setting to work.

**Default:** 0

## **SearchMethod** — Numerical search method used for iterative parameter estimation

`'auto'` (default) | `'gn'` | `'gna'` | `'lm'` | `'grad'` | `'lsqnonlin'` | `'fmincon'`

Numerical search method used for iterative parameter estimation, specified as the comma-separated pair consisting of `'SearchMethod'` and one of the following:

- `'auto'` — A combination of the line search algorithms, `'gn'`, `'lm'`, `'gna'`, and `'grad'` methods is tried in sequence at each iteration. The first descent direction leading to a reduction in estimation cost is used.
- `'gn'` — Subspace Gauss-Newton least squares search. Singular values of the Jacobian matrix less than `GnPinvConstant*eps*max(size(J))*norm(J)` are discarded when computing the search direction.  $J$  is the Jacobian matrix. The Hessian matrix is approximated as  $J^T J$ . If there is no improvement in this direction, the function tries the gradient direction.
- `'gna'` — Adaptive subspace Gauss-Newton search. Eigenvalues less than `gamma*max(sv)` of the Hessian are ignored, where `sv` contains the singular values of the Hessian. The Gauss-Newton direction is computed in the remaining subspace.

*gamma* has the initial value `InitialGnaTolerance` (see `Advanced in 'SearchOptions'` for more information). This value is increased by the factor `LMStep` each time the search fails to find a lower value of the criterion in fewer than five bisections. This value is decreased by the factor `2*LMStep` each time a search is successful without any bisections.

- `'lm'` — Levenberg-Marquardt least squares search, where the next parameter value is  $-\text{pinv}(H+d*I)*\text{grad}$  from the previous one.  $H$  is the Hessian,  $I$  is the identity matrix, and  $\text{grad}$  is the gradient.  $d$  is a number that is increased until a lower value of the criterion is found.
- `'grad'` — Steepest descent least squares search.
- `'lsqnonlin'` — Trust-region-reflective algorithm of `lsqnonlin`. Requires Optimization Toolbox software.
- `'fmincon'` — Constrained nonlinear solvers. You can use the sequential quadratic programming (SQP) and trust-region-reflective algorithms of the `fmincon` solver. If you have Optimization Toolbox software, you can also use the interior-point and active-set algorithms of the `fmincon` solver. Specify the algorithm in the `SearchOptions.Algorithm` option. The `fmincon` algorithms may result in improved estimation results in the following scenarios:
  - Constrained minimization problems when there are bounds imposed on the model parameters.
  - Model structures where the loss function is a nonlinear or non smooth function of the parameters.
  - Multi-output model estimation. A determinant loss function is minimized by default for multi-output model estimation. `fmincon` algorithms are able to minimize such loss functions directly. The other search methods such as `'lm'` and `'gn'` minimize the determinant loss function by alternately estimating the noise variance and reducing the loss value for a given noise variance value. Hence, the `fmincon` algorithms can offer better efficiency and accuracy for multi-output model estimations.

### **SearchOptions — Option set for the search algorithm**

`search option set`

Option set for the search algorithm, specified as the comma-separated pair consisting of `'SearchOptions'` and a search option set with fields that depend on the value of `SearchMethod`.

**SearchOptions Structure When SearchMethod is Specified as 'gn', 'gna', 'lm', 'grad', or 'auto'**

Field Name	Description	Default
Tolerance	Minimum percentage difference between the current value of the loss function and its expected improvement after the next iteration, specified as a positive scalar. When the percentage of expected improvement is less than <code>Tolerance</code> , the iterations stop. The estimate of the expected loss-function improvement at the next iteration is based on the Gauss-Newton vector computed for the current parameter value.	0.01
MaxIterations	<p>Maximum number of iterations during loss-function minimization, specified as a positive integer. The iterations stop when <code>MaxIterations</code> is reached or another stopping criterion is satisfied, such as <code>Tolerance</code>.</p> <p>Setting <code>MaxIterations = 0</code> returns the result of the start-up procedure.</p> <p>Use <code>sys.Report.Termination.Iterations</code> to get the actual number of iterations during an estimation, where <code>sys</code> is an <code>idtf</code> model.</p>	20

Field Name	Description	Default
Advanced	Advanced search settings, specified as a structure with the following fields:	
Field Name	Description	Default
GnPinvConstant	Jacobian matrix singular value threshold, specified as a positive scalar. Singular values of the Jacobian matrix that are smaller than $\text{GnPinvConstant} * \max(\text{size}(J) * \text{norm}(J) * \text{eps})$ are discarded when computing the search direction. Applicable when SearchMethod is 'gn'.	10000
InitialGnaTolerance	Initial value of $\gamma$ , specified as a positive scalar. Applicable when SearchMethod is 'gna'.	0.0001
LMStartValue	Starting value of search-direction length $d$ in the Levenberg-Marquardt method, specified as a positive scalar. Applicable when SearchMethod is 'lm'.	0.001
LMStep	Size of the Levenberg-Marquardt step, specified as a positive integer. The next value of the search-direction length $d$ in the Levenberg-Marquardt method is LMStep times the previous one. Applicable when SearchMethod is 'lm'.	2
MaxBisections	Maximum number of bisections used for line search along the search direction, specified as a positive integer.	25
MaxFunctionEvaluations	Maximum number of calls to the model file, specified as a positive integer. Iterations stop if the number of calls to the model file exceeds this value.	Inf
MinParameterChange	Smallest parameter update allowed per iteration, specified as a nonnegative scalar.	0
RelativeImprovement	Relative improvement threshold, specified as a nonnegative scalar. Iterations stop if the relative	0

Field Name	Description		Default
	Field Name	Description	Default
		improvement of the criterion function is less than this value.	
	StepReduction	<p>Step reduction factor, specified as a positive scalar that is greater than 1. The suggested parameter update is reduced by the factor <code>StepReduction</code> after each try. This reduction continues until <code>MaxBisections</code> tries are completed or a lower value of the criterion function is obtained.</p> <p><code>StepReduction</code> is not applicable for SearchMethod 'lm' (Levenberg-Marquardt method).</p>	2

**SearchOptions Structure When SearchMethod is Specified as 'lsqnonlin'**

<b>Field Name</b>	<b>Description</b>	<b>Default</b>
FunctionTolerance	Termination tolerance on the loss function that the software minimizes to determine the estimated parameter values, specified as a positive scalar.  The value of FunctionTolerance is the same as that of <code>opt.SearchOptions.Advanced.TolFun</code> .	1e-5
StepTolerance	Termination tolerance on the estimated parameter values, specified as a positive scalar.  The value of StepTolerance is the same as that of <code>opt.SearchOptions.Advanced.TolX</code> .	1e-6
MaxIterations	Maximum number of iterations during loss-function minimization, specified as a positive integer. The iterations stop when MaxIterations is reached or another stopping criterion is satisfied, such as FunctionTolerance.  The value of MaxIterations is the same as that of <code>opt.SearchOptions.Advanced.MaxIter</code> .	20
Advanced	Advanced search settings, specified as an option set for <code>lsqnonlin</code> .  For more information, see the Optimization Options table in "Optimization Options" (Optimization Toolbox).	Use <code>optimset('lsqnonlin')</code> to create a default option set.

**SearchOptions Structure When SearchMethod is Specified as 'fmincon'**

Field Name	Description	Default
Algorithm	<p>fmincon optimization algorithm, specified as one of the following:</p> <ul style="list-style-type: none"> <li>• 'sqp' — Sequential quadratic programming algorithm. The algorithm satisfies bounds at all iterations, and it can recover from NaN or Inf results. It is not a large-scale algorithm. For more information, see “Large-Scale vs. Medium-Scale Algorithms” (Optimization Toolbox).</li> <li>• 'trust-region-reflective' — Subspace trust-region method based on the interior-reflective Newton method. It is a large-scale algorithm.</li> <li>• 'interior-point' — Large-scale algorithm that requires Optimization Toolbox software. The algorithm satisfies bounds at all iterations, and it can recover from NaN or Inf results.</li> <li>• 'active-set' — Requires Optimization Toolbox software. The</li> </ul>	'sqp'

Field Name	Description	Default
	<p>algorithm can take large steps, which adds speed. It is not a large-scale algorithm.</p> <p>For more information about the algorithms, see “Constrained Nonlinear Optimization Algorithms” (Optimization Toolbox) and “Choosing the Algorithm” (Optimization Toolbox).</p>	
FunctionTolerance	Termination tolerance on the loss function that the software minimizes to determine the estimated parameter values, specified as a positive scalar.	1e-6
StepTolerance	Termination tolerance on the estimated parameter values, specified as a positive scalar.	1e-6
MaxIterations	Maximum number of iterations during loss function minimization, specified as a positive integer. The iterations stop when MaxIterations is reached or another stopping criterion is satisfied, such as FunctionTolerance.	100

### Advanced – Additional advanced options structure

Additional advanced options, specified as a structure with the following fields:

- `ErrorThreshold` — Specifies when to adjust the weight of large errors from quadratic to linear.

Errors larger than `ErrorThreshold` times the estimated standard deviation have a linear weight in the loss function. The standard deviation is estimated robustly as the median of the absolute deviations from the median of the prediction errors and divided by 0.7. For more information on robust norm choices, see section 15.2 of [2].

`ErrorThreshold = 0` disables robustification and leads to a purely quadratic loss function. When estimating with frequency-domain data, the software sets `ErrorThreshold` to zero. For time-domain data that contains outliers, try setting `ErrorThreshold` to 1.6.

**Default:** 0

- `MaxSize` — Specifies the maximum number of elements in a segment when input-output data is split into segments.

`MaxSize` must be a positive integer.

**Default:** 250000

- `StabilityThreshold` — Specifies thresholds for stability tests.

`StabilityThreshold` is a structure with the following fields:

- `s` — Specifies the location of the right-most pole to test the stability of continuous-time models. A model is considered stable when its right-most pole is to the left of `s`.

**Default:** 0

- `z` — Specifies the maximum distance of all poles from the origin to test stability of discrete-time models. A model is considered stable if all poles are within the distance `z` from the origin.

**Default:** `1+sqrt(eps)`

- `AutoInitThreshold` — Specifies when to automatically estimate the initial condition.

The initial condition is estimated when

$$\frac{\|y_{p,z} - y_{meas}\|}{\|y_{p,e} - y_{meas}\|} > \text{AutoInitThreshold}$$

- $y_{meas}$  is the measured output.
- $y_{p,z}$  is the predicted output of a model estimated using zero initial conditions.
- $y_{p,e}$  is the predicted output of a model estimated using estimated initial conditions.

Applicable when `InitialCondition` is 'auto'.

**Default:** 1.05

## Output Arguments

**opt** — Options set for armax

armaxOptions option set

Option set for armax, returned as an armaxOptions option set.

## Examples

### Create Default Options Set for ARMAX Estimation

```
opt = armaxOptions;
```

### Specify Options for ARMAX Estimation

Create an option set for armax to use the 'simulation' Focus and to set the Display to 'on'.

```
opt = armaxOptions('Focus','simulation','Display','on');
```

Alternatively, use dot notation to set the values of opt.

```
opt = armaxOptions;  
opt.Focus = 'simulation';  
opt.Display = 'on';
```

## References

- [1] Wills, Adrian, B. Ninness, and S. Gibson. "On Gradient-Based Search for Multivariable System Estimates". *Proceedings of the 16th IFAC World Congress, Prague, Czech Republic, July 3-8, 2005*. Oxford, UK: Elsevier Ltd., 2005.
- [2] Ljung, L. *System Identification: Theory for the User*. Upper Saddle River, NJ: Prentice-Hall PTR, 1999.

## See Also

armax | idfilt

## Topics

"Loss Function and Model Quality Metrics"

**Introduced in R2012a**

# arOptions

Option set for ar

## Syntax

```
opt = arOptions  
opt = arOptions(Name,Value)
```

## Description

`opt = arOptions` creates the default options set for ar.

`opt = arOptions(Name,Value)` creates an option set with the options specified by one or more Name,Value pair arguments.

## Input Arguments

### Name-Value Pair Arguments

Specify optional comma-separated pairs of Name,Value arguments. Name is the argument name and Value is the corresponding value. Name must appear inside quotes. You can specify several name and value pair arguments in any order as Name1,Value1,...,NameN,ValueN.

### Approach

Technique used for AR model estimation.

Approach is specified as one of the following values:

- 'fb' — Forward-backward approach.
- 'ls' — Least-squares method.
- 'yw' — Yule-Walker approach.

- 'burg' — Burg's method.
- 'gl' — Geometric lattice method.

**Default:** 'fb'

### **Window**

Data windowing technique.

Window determines how the data outside the measured time interval (past and future values) is handled.

Window is specified as one of the following values:

- 'now' — No windowing.
- 'prw' — Pre-windowing.
- 'pow' — Post-windowing.
- 'ppw' — Pre- and post-windowing.

This option is ignored when you use the Yule-Walker approach.

**Default:** 'now'

### **DataOffset**

Data offset level that is removed before estimation of parameters.

Specify `DataOffset` as a double scalar. For multiexperiment data, specify `DataOffset` as a vector of length  $N_e$ , where  $N_e$  is the number of experiments. Each entry of the vector is subtracted from the corresponding data.

**Default:** [] (no offsets)

### **MaxSize**

Specifies the maximum number of elements in a segment when input/output data is split into segments.

If larger matrices are needed, the software uses loops for calculations. Use this option to manage the trade-off between memory management and program execution speed. The original data matrix must be smaller than the matrix specified by `MaxSize`.

MaxSize must be a positive integer.

**Default:** 250000

## Output Arguments

**opt**

Option set containing the specified options for ar.

## Examples

### Create Default Options Set for AR Estimation

```
opt = arOptions;
```

### Specify Options for AR Estimation

Create an options set for ar using the least squares algorithm for estimation. Set Window to 'ppw'.

```
opt = arOptions('Approach','ls','Window','ppw');
```

Alternatively, use dot notation to set the values of opt.

```
opt = arOptions;  
opt.Approach = 'ls';  
opt.Window = 'ppw';
```

## See Also

ar

**Introduced in R2012a**

## arx

Estimate parameters of ARX or AR model using least squares

### Syntax

```
sys = arx(data,[na nb nk])  
sys = arx(data,[na nb nk],Name,Value)  
sys = arx(data,[na nb nk], ___,opt)
```

### Description

---

**Note** arx does not support continuous-time estimations. Use `tfest` instead.

---

`sys = arx(data,[na nb nk])` returns an ARX structure polynomial model, `sys`, with estimated parameters and covariances (parameter uncertainties) using the least-squares method and specified orders.

`sys = arx(data,[na nb nk],Name,Value)` estimates a polynomial model with additional options specified by one or more `Name,Value` pair arguments.

`sys = arx(data,[na nb nk], ___,opt)` specifies estimation options that configure the estimation objective, initial conditions and handle input/output data offsets.

### Input Arguments

#### **data**

Estimation data.

Specify `data` as an `iddata` object, an `frd` object, or an `idfrd` frequency-response-data object.

**[na nb nk]**

Polynomial orders.

[na nb nk] define the polynomial orders of an ARX model.

- na — Order of the polynomial  $A(q)$ .

Specify na as an  $N_y$ -by- $N_y$  matrix of nonnegative integers.  $N_y$  is the number of outputs.

- nb — Order of the polynomial  $B(q) + 1$ .

nb is an  $N_y$ -by- $N_u$  matrix of nonnegative integers.  $N_y$  is the number of outputs and  $N_u$  is the number of inputs.

- nk — Input-output delay expressed as fixed leading zeros of the  $B$  polynomial.

Specify nk as an  $N_y$ -by- $N_u$  matrix of nonnegative integers.  $N_y$  is the number of outputs and  $N_u$  is the number of inputs.

**opt**

Estimation options.

opt is an options set that specifies estimation options, including:

- input/output data offsets
- output weight

Use `arxOptions` to create the options set.

**Name-Value Pair Arguments**

Specify optional comma-separated pairs of `Name`, `Value` arguments. `Name` is the argument name and `Value` is the corresponding value. `Name` must appear inside quotes. You can specify several name and value pair arguments in any order as `Name1, Value1, . . . , NameN, ValueN`.

**InputDelay**

Input delays. `InputDelay` is a numeric vector specifying a time delay for each input channel. Specify input delays in integer multiples of the sample time  $T_s$ . For example, `InputDelay = 3` means a delay of three sampling periods.

For a system with  $N_u$  inputs, set `InputDelay` to an  $N_u$ -by-1 vector, where each entry is a numerical value representing the input delay for the corresponding input channel. You can also set `InputDelay` to a scalar value to apply the same delay to all channels.

**Default:** 0 for all input channels

## **IODelay**

Transport delays. `IODelay` is a numeric array specifying a separate transport delay for each input/output pair.

Specify transport delays as integers denoting delay of a multiple of the sample time,  $T_s$ .

For a MIMO system with  $N_y$  outputs and  $N_u$  inputs, set `IODelay` to a  $N_y$ -by- $N_u$  array, where each entry is a numerical value representing the transport delay for the corresponding input/output pair. You can also set `IODelay` to a scalar value to apply the same delay to all input/output pairs. Useful as a replacement for the `nk` order, you can factor out  $\max(nk - 1, 0)$  lags as the `IODelay` value.

**Default:** 0 for all input/output pairs

## **IntegrateNoise**

Specify integrators in the noise channels.

Adding an integrator creates an ARIX model represented by:

$$A(q)y(t) = B(q)u(t - nk) + \frac{1}{1 - q^{-1}} e(t)$$

where,  $\frac{1}{1 - q^{-1}}$  is the integrator in the noise channel,  $e(t)$ .

`IntegrateNoise` is a logical vector of length  $N_y$ , where  $N_y$  is the number of outputs.

**Default:** `false(Ny, 1)`, where  $N_y$  is the number of outputs

## Output Arguments

### sys

ARX model that fits the estimation data, returned as a discrete-time `idpoly` object. This model is created using the specified model orders, delays, and estimation options.

Information about the estimation results and options used is stored in the `Report` property of the model. `Report` has the following fields:

Report Field	Description
Status	Summary of the model status, which indicates whether the model was created by construction or obtained by estimation.
Method	Estimation command used.
InitialCondition	<p>Handling of initial conditions during model estimation, returned as one of the following values:</p> <ul style="list-style-type: none"> <li>'zero' — The initial conditions were set to zero.</li> <li>'estimate' — The initial conditions were treated as independent estimation parameters.</li> </ul> <p>This field is especially useful to view how the initial conditions were handled when the <code>InitialCondition</code> option in the estimation option set is 'auto'.</p>

Report Field	Description																		
Fit	<p>Quantitative assessment of the estimation, returned as a structure. See “Loss Function and Model Quality Metrics” for more information on these quality metrics. The structure has the following fields:</p> <table border="1" data-bbox="387 456 1338 951"> <thead> <tr> <th data-bbox="387 456 520 499">Field</th> <th data-bbox="520 456 1338 499">Description</th> </tr> </thead> <tbody> <tr> <td data-bbox="387 499 520 609">FitPercent</td> <td data-bbox="520 499 1338 609">Normalized root mean squared error (NRMSE) measure of how well the response of the model fits the estimation data, expressed as a percentage.</td> </tr> <tr> <td data-bbox="387 609 520 652">LossFcn</td> <td data-bbox="520 609 1338 652">Value of the loss function when the estimation completes.</td> </tr> <tr> <td data-bbox="387 652 520 730">MSE</td> <td data-bbox="520 652 1338 730">Mean squared error (MSE) measure of how well the response of the model fits the estimation data.</td> </tr> <tr> <td data-bbox="387 730 520 774">FPE</td> <td data-bbox="520 730 1338 774">Final prediction error for the model.</td> </tr> <tr> <td data-bbox="387 774 520 817">AIC</td> <td data-bbox="520 774 1338 817">Raw Akaike Information Criteria (AIC) measure of model quality.</td> </tr> <tr> <td data-bbox="387 817 520 861">AICc</td> <td data-bbox="520 817 1338 861">Small sample-size corrected AIC.</td> </tr> <tr> <td data-bbox="387 861 520 904">nAIC</td> <td data-bbox="520 861 1338 904">Normalized AIC.</td> </tr> <tr> <td data-bbox="387 904 520 951">BIC</td> <td data-bbox="520 904 1338 951">Bayesian Information Criteria (BIC).</td> </tr> </tbody> </table>	Field	Description	FitPercent	Normalized root mean squared error (NRMSE) measure of how well the response of the model fits the estimation data, expressed as a percentage.	LossFcn	Value of the loss function when the estimation completes.	MSE	Mean squared error (MSE) measure of how well the response of the model fits the estimation data.	FPE	Final prediction error for the model.	AIC	Raw Akaike Information Criteria (AIC) measure of model quality.	AICc	Small sample-size corrected AIC.	nAIC	Normalized AIC.	BIC	Bayesian Information Criteria (BIC).
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nAIC	Normalized AIC.																		
BIC	Bayesian Information Criteria (BIC).																		
Parameters	Estimated values of model parameters.																		
OptionsUsed	Option set used for estimation. If no custom options were configured, this is a set of default options. See <code>arxOptions</code> for more information.																		
RandState	State of the random number stream at the start of estimation. Empty, [], if randomization was not used during estimation. For more information, see <code>rng</code> in the MATLAB documentation.																		

Report Field	Description																
DataUsed	Attributes of the data used for estimation, returned as a structure with the following fields:																
	<table border="1"> <thead> <tr> <th>Field</th> <th>Description</th> </tr> </thead> <tbody> <tr> <td>Name</td> <td>Name of the data set.</td> </tr> <tr> <td>Type</td> <td>Data type.</td> </tr> <tr> <td>Length</td> <td>Number of data samples.</td> </tr> <tr> <td>Ts</td> <td>Sample time.</td> </tr> <tr> <td>InterSample</td> <td>Input intersample behavior, returned as one of the following values: <ul style="list-style-type: none"> <li>'zoh' — Zero-order hold maintains a piecewise-constant input signal between samples.</li> <li>'foh' — First-order hold maintains a piecewise-linear input signal between samples.</li> <li>'bl' — Band-limited behavior specifies that the continuous-time input signal has zero power above the Nyquist frequency.</li> </ul> </td> </tr> <tr> <td>InputOffset</td> <td>Offset removed from time-domain input data during estimation. For nonlinear models, it is [ ].</td> </tr> <tr> <td>OutputOffset</td> <td>Offset removed from time-domain output data during estimation. For nonlinear models, it is [ ].</td> </tr> </tbody> </table>	Field	Description	Name	Name of the data set.	Type	Data type.	Length	Number of data samples.	Ts	Sample time.	InterSample	Input intersample behavior, returned as one of the following values: <ul style="list-style-type: none"> <li>'zoh' — Zero-order hold maintains a piecewise-constant input signal between samples.</li> <li>'foh' — First-order hold maintains a piecewise-linear input signal between samples.</li> <li>'bl' — Band-limited behavior specifies that the continuous-time input signal has zero power above the Nyquist frequency.</li> </ul>	InputOffset	Offset removed from time-domain input data during estimation. For nonlinear models, it is [ ].	OutputOffset	Offset removed from time-domain output data during estimation. For nonlinear models, it is [ ].
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For more information on using `Report`, see “Estimation Report”.

## Examples

### Estimate ARX model

Generate input data based on a specified ARX model, and then use this data to estimate an ARX model.

```
A = [1 -1.5 0.7];
B = [0 1 0.5];
```

```
m0 = idpoly(A,B);  
u = iddata([],idinput(300,'rbs'));  
e = iddata([],randn(300,1));  
y = sim(m0,[u e]);  
z = [y,u];  
m = arx(z,[2 2 1]);
```

## Estimate ARX Model Using Regularization

Use `arxRegul` to automatically determine regularization constants and use the values for estimating an FIR model of order 50.

Obtain L and R values.

```
load regularizationExampleData eData;  
orders = [0 50 0];  
[L,R] = arxRegul(eData,orders);
```

By default, the TC kernel is used.

Use the returned Lambda and R values for regularized ARX model estimation.

```
opt = arxOptions;  
opt.Regularization.Lambda = L;  
opt.Regularization.R = R;  
model = arx(eData,orders,opt);
```

## Definitions

### ARX structure

The ARX model structure is :

$$y(t) + a_1 y(t-1) + \dots + a_{na} y(t-na) = b_1 u(t-nk) + \dots + b_{nb} u(t-nb-nk+1) + e(t)$$

The parameters  $na$  and  $nb$  are the orders of the ARX model, and  $nk$  is the delay.

- $y(t)$  — Output at time  $t$ .
- $n_a$  — Number of poles.
- $n_b$  — Number of zeroes plus 1.
- $n_k$  — Number of input samples that occur before the input affects the output, also called the *dead time* in the system.
- $y(t-1)\dots y(t-n_a)$  — Previous outputs on which the current output depends.
- $u(t-n_k)\dots u(t-n_k-n_b+1)$  — Previous and delayed inputs on which the current output depends.
- $e(t)$  — White-noise disturbance value.

A more compact way to write the difference equation is

$$A(q)y(t) = B(q)u(t - n_k) + e(t)$$

$q$  is the delay operator. Specifically,

$$A(q) = 1 + a_1q^{-1} + \dots + a_{n_a}q^{-n_a}$$

$$B(q) = b_1 + b_2q^{-1} + \dots + b_{n_b}q^{-n_b+1}$$

## Time Series Models

For time-series data that contains no inputs, one output and orders =  $n_a$ , the model has AR structure of order  $n_a$ .

The AR model structure is

$$A(q)y(t) = e(t)$$

## Multiple Inputs and Single-Output Models

For multiple-input systems,  $\mathbf{nb}$  and  $\mathbf{nk}$  are row vectors where the  $i$ th element corresponds to the order and delay associated with the  $i$ th input.

$$y(t) + A_1 y(t-1) + A_2 y(t-2) + \dots + A_{na} y(t-na) = B_0 u(t) + B_1 u(t-1) + \dots + B_{nb} u(t-nb) + e(t)$$

## Multi-Output Models

For models with multiple inputs and multiple outputs,  $\mathbf{na}$ ,  $\mathbf{nb}$ , and  $\mathbf{nk}$  contain one row for each output signal.

In the multiple-output case, `arx` minimizes the trace of the prediction error covariance matrix, or the norm

$$\sum_{t=1}^N e^T(t)e(t)$$

To transform this to an arbitrary quadratic norm using a weighting matrix  $\Lambda$

$$\sum_{t=1}^N e^T(t)\Lambda^{-1}e(t)$$

use the syntax

```
opt = arxOptions('OutputWeight',inv(lambda))  
m = arx(data,orders,opt)
```

## Estimating Initial Conditions

For time-domain data, the signals are shifted such that unmeasured signals are never required in the predictors. Therefore, there is no need to estimate initial conditions.

For frequency-domain data, it might be necessary to adjust the data by initial conditions that support circular convolution.

Set the `InitialCondition` estimation option (see `arxOptions`) to one the following values:

- 'zero' — No adjustment.
- 'estimate' — Perform adjustment to the data by initial conditions that support circular convolution.
- 'auto' — Automatically choose between 'zero' and 'estimate' based on the data.

## Algorithms

QR factorization solves the overdetermined set of linear equations that constitute the least-squares estimation problem.

Without regularization, the ARX model parameters vector  $\theta$  is estimated by solving the normal equation:

$$(J^T J)\theta = J^T y$$

where  $J$  is the regressor matrix and  $y$  is the measured output. Therefore,

$$\theta = (J^T J)^{-1} J^T y.$$

Using regularization adds a regularization term:

$$\theta = (J^T J + \lambda R)^{-1} J^T y$$

where,  $\lambda$  and  $R$  are the regularization constants. See `arxOptions` for more information on the regularization constants.

When the regression matrix is larger than the `MaxSize` specified in `arxOptions`, data is segmented and QR factorization is performed iteratively on these data segments.

## See Also

`ar` | `armax` | `arxOptions` | `arxRegul` | `arxstruc` | `bj` | `impulseest` | `iv4` | `n4sid` | `nlarx` | `oe`

**Topics**

“Initialize Nonlinear ARX Estimation Using Linear Model”

“Regularized Estimates of Model Parameters”

**Introduced before R2006a**

# arxOptions

Option set for arx

## Syntax

```
opt = arxOptions
opt = arxOptions(Name,Value)
```

## Description

`opt = arxOptions` creates the default options set for arx.

`opt = arxOptions(Name,Value)` creates an option set with the options specified by one or more `Name,Value` pair arguments.

## Input Arguments

### Name-Value Pair Arguments

Specify optional comma-separated pairs of `Name,Value` arguments. `Name` is the argument name and `Value` is the corresponding value. `Name` must appear inside quotes. You can specify several name and value pair arguments in any order as `Name1,Value1,...,NameN,ValueN`.

### InitialCondition — Handling of initial conditions

'auto' (default) | 'zero' | 'estimate'

Handling of initial conditions during estimation using frequency-domain data, specified as the comma-separated pair consisting of 'InitialCondition' and one of the following values:

- 'zero' — The initial conditions are set to zero.
- 'estimate' — The initial conditions are treated as independent estimation parameters.

- 'auto' — The software chooses the method to handle initial conditions based on the estimation data.

### **Focus — Error to be minimized**

'prediction' (default) | 'simulation'

Error to be minimized in the loss function during estimation, specified as the comma-separated pair consisting of 'Focus' and one of the following values:

- 'prediction' — The one-step ahead prediction error between measured and predicted outputs is minimized during estimation. As a result, the estimation focuses on producing a good predictor model.
- 'simulation' — The simulation error between measured and simulated outputs is minimized during estimation. As a result, the estimation focuses on making a good fit for simulation of model response with the current inputs.

The Focus option can be interpreted as a weighting filter in the loss function. For more information, see “Loss Function and Model Quality Metrics”.

### **WeightingFilter — Weighting prefilter**

[] (default) | vector | matrix | cell array | linear system

Weighting prefilter applied to the loss function to be minimized during estimation. To understand the effect of WeightingFilter on the loss function, see “Loss Function and Model Quality Metrics”.

Specify WeightingFilter as one of the following values:

- [] — No weighting prefilter is used.
- Passbands — Specify a row vector or matrix containing frequency values that define desired passbands. You select a frequency band where the fit between estimated model and estimation data is optimized. For example, [w<sub>l</sub>, w<sub>h</sub>] where w<sub>l</sub> and w<sub>h</sub> represent lower and upper limits of a passband. For a matrix with several rows defining frequency passbands, [w<sub>1l</sub>, w<sub>1h</sub>; w<sub>2l</sub>, w<sub>2h</sub>; w<sub>3l</sub>, w<sub>3h</sub>; . . . ], the estimation algorithm uses the union of the frequency ranges to define the estimation passband.

Passbands are expressed in rad/TimeUnit for time-domain data and in FrequencyUnit for frequency-domain data, where TimeUnit and FrequencyUnit are the time and frequency units of the estimation data.

- SISO filter — Specify a single-input-single-output (SISO) linear filter in one of the following ways:

- A SISO LTI model
- $\{A, B, C, D\}$  format, which specifies the state-space matrices of a filter with the same sample time as estimation data.
- $\{\text{numerator}, \text{denominator}\}$  format, which specifies the numerator and denominator of the filter as a transfer function with same sample time as estimation data.

This option calculates the weighting function as a product of the filter and the input spectrum to estimate the transfer function.

- **Weighting vector** — Applicable for frequency-domain data only. Specify a column vector of weights. This vector must have the same length as the frequency vector of the data set, `Data.Frequency`. Each input and output response in the data is multiplied by the corresponding weight at that frequency.

#### **EnforceStability** — Control whether to enforce stability of model

`false` (default) | `true`

Control whether to enforce stability of estimated model, specified as the comma-separated pair consisting of 'EnforceStability' and either `true` or `false`.

This option is not available for multi-output models with a non-diagonal  $A$  polynomial array.

Data Types: `logical`

#### **EstimateCovariance** — Control whether to generate parameter covariance data

`true` (default) | `false`

Controls whether parameter covariance data is generated, specified as `true` or `false`.

If `EstimateCovariance` is `true`, then use `getcov` to fetch the covariance matrix from the estimated model.

#### **Display** — Specify whether to display the estimation progress

'off' (default) | 'on'

Specify whether to display the estimation progress, specified as one of the following values:

- 'on' — Information on model structure and estimation results are displayed in a progress-viewer window.

- 'off' — No progress or results information is displayed.

## **InputOffset — Removal of offset from time-domain input data during estimation**

[ ] (default) | vector of positive integers | matrix

Removal of offset from time-domain input data during estimation, specified as the comma-separated pair consisting of 'InputOffset' and one of the following:

- A column vector of positive integers of length  $Nu$ , where  $Nu$  is the number of inputs.
- [ ] — Indicates no offset.
- $Nu$ -by- $Ne$  matrix — For multi-experiment data, specify InputOffset as an  $Nu$ -by- $Ne$  matrix.  $Nu$  is the number of inputs, and  $Ne$  is the number of experiments.

Each entry specified by InputOffset is subtracted from the corresponding input data.

## **OutputOffset — Removal of offset from time-domain output data during estimation**

[ ] (default) | vector | matrix

Removal of offset from time-domain output data during estimation, specified as the comma-separated pair consisting of 'OutputOffset' and one of the following:

- A column vector of length  $Ny$ , where  $Ny$  is the number of outputs.
- [ ] — Indicates no offset.
- $Ny$ -by- $Ne$  matrix — For multi-experiment data, specify OutputOffset as a  $Ny$ -by- $Ne$  matrix.  $Ny$  is the number of outputs, and  $Ne$  is the number of experiments.

Each entry specified by OutputOffset is subtracted from the corresponding output data.

## **OutputWeight — Weight of prediction errors in multi-output estimation**

[ ] (default) | positive semidefinite, symmetric matrix

Weight of prediction errors in multi-output estimation, specified as one of the following values:

- Positive semidefinite, symmetric matrix ( $W$ ). The software minimizes the trace of the weighted prediction error matrix  $\text{trace}(E' * E * W / N)$  where:
  - $E$  is the matrix of prediction errors, with one column for each output, and  $W$  is the positive semidefinite, symmetric matrix of size equal to the number of outputs. Use

W to specify the relative importance of outputs in multiple-output models, or the reliability of corresponding data.

- N is the number of data samples.
- [] — No weighting is used. Specifying as [] is the same as eye(Ny), where Ny is the number of outputs.

This option is relevant only for multi-output models.

### **Regularization — Options for regularized estimation of model parameters**

[] (default) | positive semidefinite, symmetric matrix

Options for regularized estimation of model parameters, specified as a structure with the following fields:

- Lambda — Constant that determines the bias versus variance tradeoff.

Specify a positive scalar to add the regularization term to the estimation cost.

The default value of zero implies no regularization.

**Default:** 0

- R — Weighting matrix.

Specify a positive scalar or a positive definite matrix. The length of the matrix must be equal to the number of free parameters (np) of the model. For ARX model, np = sum(sum([na nb])).

**Default:** 1

- Nominal — This option is not used for ARX models.

**Default:** 0

Use arxRegul to automatically determine Lambda and R values.

For more information on regularization, see “Regularized Estimates of Model Parameters”.

### **Advanced — Additional advanced options**

structure

Additional advanced options, specified as a structure with the following fields:

- **MaxSize** — Specifies the maximum number of elements in a segment when input-output data is split into segments.

MaxSize must be a positive integer.

**Default:** 250000

- **StabilityThreshold** — Specifies thresholds for stability tests.

StabilityThreshold is a structure with the following fields:

- **s** — Specifies the location of the right-most pole to test the stability of continuous-time models. A model is considered stable when its right-most pole is to the left of s.

**Default:** 0

- **z** — Specifies the maximum distance of all poles from the origin to test stability of discrete-time models. A model is considered stable if all poles are within the distance z from the origin.

**Default:** 1+sqrt(eps)

## Output Arguments

**opt** — Options set for arx

arxOptions option set

Option set for arx, returned as an arxOptions option set.

## Examples

**Create Default Options Set for ARX Estimation**

```
opt = arxOptions;
```

## Specify Options for ARX Estimation

Create an options set for arx using zero initial conditions for estimation. Set `Display` to 'on'.

```
opt = arxOptions('InitialCondition','zero','Display','on');
```

Alternatively, use dot notation to set the values of `opt`.

```
opt = arxOptions;  
opt.InitialCondition = 'zero';  
opt.Display = 'on';
```

## See Also

`arx` | `arxRegul` | `idfilt`

## Topics

“Loss Function and Model Quality Metrics”

## Introduced in R2012a

## arxRegul

Determine regularization constants for ARX model estimation

### Syntax

```
[lambda,R] = arxRegul(data,orders)
[lambda,R] = arxRegul(data,orders,options)
[lambda,R] = arxRegul(data,orders,Name,Value)
[lambda,R] = arxRegul(data,orders,options,Name,Value)
```

### Description

`[lambda,R] = arxRegul(data,orders)` returns the regularization constants used for ARX model estimation. Use the regularization constants in `arxOptions` to configure the regularization options for ARX model estimation.

`[lambda,R] = arxRegul(data,orders,options)` specifies regularization options such as regularization kernel and I/O offsets.

`[lambda,R] = arxRegul(data,orders,Name,Value)` specifies model structure attributes, such as noise integrator and input delay, using one or more `Name,Value` pair arguments.

`[lambda,R] = arxRegul(data,orders,options,Name,Value)` specifies both regularization options and model structure attributes.

### Examples

#### Determine Regularization Constants for ARX Model Estimation Using Default Kernel

```
load iddata1 z1;
orders = [10 10 1];
[Lambda,R] = arxRegul(z1,orders);
```

The ARX model is estimated using the default regularization kernel TC.

Use the Lambda and R values for ARX model estimation.

```
opt = arxOptions;  
opt.Regularization.Lambda = Lambda;  
opt.Regularization.R = R;  
model = arx(z1,orders,opt);
```

### Specify a Regularization Kernel

Specify 'DC' as the regularization kernel and obtain a regularized ARX model of order [10 10 1].

```
load iddata1 z1;  
orders = [10 10 1];  
option = arxRegulOptions('RegularizationKernel','DC');  
[Lambda,R] = arxRegul(z1,orders,option);
```

Use the Lambda and R values for ARX model estimation.

```
arxOpt = arxOptions;  
arxOpt.Regularization.Lambda = Lambda;  
arxOpt.Regularization.R = R;  
model = arx(z1,orders,arxOpt);
```

### Specify Noise Source Integrator

Specify to include a noise source integrator in the noise component of the model.

```
load iddata1 z1;  
orders = [10 10 1];  
[Lambda,R] = arxRegul(z1,orders,'IntegrateNoise',true);
```

## Specify Regularization Kernel And Noise Integrator

Specify the regularization kernel and include a noise source integrator in the noise component of the model.

```
load iddata1 z1;
orders = [10 10 1];
opt = arxRegulOptions('RegularizationKernel','DC');
[Lambda,R] = arxRegul(z1,orders,opt,'IntegrateNoise',true);
```

## Input Arguments

### data — Estimation data

iddata object

Estimation data, specified as an iddata object.

### orders — ARX model orders

matrix of nonnegative integers

ARX model orders [na nb nc], specified as a matrix of nonnegative integers. See the arx reference page for more information on model orders.

### options — Regularization options

arxRegulOptions options set

Regularization options, specified as an options set you create using arxRegulOptions.

## Name-Value Pair Arguments

Specify optional comma-separated pairs of Name, Value arguments. Name is the argument name and Value is the corresponding value. Name must appear inside quotes. You can specify several name and value pair arguments in any order as Name1, Value1, ..., NameN, ValueN.

Example: [Lambda, R] = arxRegul(z1,orders,option,'InputDelay',10);

### InputDelay — Input delay

0 (default) | positive integer

Input delay, specified as a positive, nonzero numeric value representing the number of samples.

Example: `[Lambda, R] = arxRegul(z1,orders,'InputDelay',10);`

Data Types: `double`

### **IntegrateNoise — Noise source integrator**

`false` (default) | `true`

Noise source integrator, specified as a logical. Specifies whether the noise source  $e(t)$  should contain an integrator. The default is `false`, indicating the noise integrator is off. To turn it on, change the value to `true`.

Example: `[Lambda, R] = arxRegul(z1,orders,'IntegrateNoise',true);`

Data Types: `logical`

## **Output Arguments**

### **lambda — Constant that determines bias versus variance trade-off**

positive scalar

Constant that determines the bias versus variance trade-off, returned as a positive scalar.

### **R — Weighting matrix**

vector of nonnegative numbers | square positive semi-definite matrix

Weighting matrix, returned as a vector of nonnegative numbers or a positive definite matrix.

## **Algorithms**

Without regularization, the ARX model parameters vector  $\theta$  is estimated by solving the normal equation:

$$(J^T J)\theta = J^T y$$

where  $J$  is the regressor matrix and  $y$  is the measured output. Therefore,

$$\theta = (J^T J)^{-1} J^T y.$$

Using regularization adds a regularization term:

$$\theta = (J^T J + \lambda R)^{-1} J^T y$$

where,  $\lambda$  and  $R$  are the regularization constants. See `arxOptions` for more information on the regularization constants.

### References

- [1] T. Chen, H. Ohlsson, and L. Ljung. "On the Estimation of Transfer Functions, Regularizations and Gaussian Processes - Revisited", *Automatica*, Volume 48, August 2012.

### See Also

`arx` | `arxOptions` | `arxRegulOptions`

### Topics

"Estimate Regularized ARX Model Using System Identification App"  
"Regularized Estimates of Model Parameters"

**Introduced in R2013b**

# arxRegulOptions

Option set for arxRegul

## Syntax

```
opt = arxRegulOptions  
opt = arxRegulOptions(Name,Value)
```

## Description

`opt = arxRegulOptions` creates a default option set for `arxRegul`.

`opt = arxRegulOptions(Name,Value)` creates an options set with the options specified by one or more name-value pair arguments.

## Examples

### Create Default Options Set for Determining Regularization Constants

```
opt = arxRegulOptions;
```

### Specify Regularizing Kernel for ARX Model Estimation

```
opt = arxRegulOptions('RegularizationKernel','DC');
```

## Input Arguments

### Name-Value Pair Arguments

Specify optional comma-separated pairs of `Name`, `Value` arguments. `Name` is the argument name and `Value` is the corresponding value. `Name` must appear inside quotes. You can specify several name and value pair arguments in any order as `Name1, Value1, ..., NameN, ValueN`.

Example: `option = arxRegulOptions('RegularizationKernel', 'DC')` specifies 'DC' as the regularization kernel.

#### **RegularizationKernel** — Regularization kernel

'TC' (default) | 'SE' | 'SS' | 'HF' | 'DI' | 'DC'

Regularization kernel, specified as one of the following values:

- 'TC' — Tuned and correlated kernel
- 'SE' — Squared exponential kernel
- 'SS' — Stable spline kernel
- 'HF' — High frequency stable spline kernel
- 'DI' — Diagonal kernel
- 'DC' — Diagonal and correlated kernel

The specified kernel is used for regularized estimation of impulse response for all input-output channels. Regularization reduces variance of estimated model coefficients and produces a smoother response by trading variance for bias.

For more information about these choices, see [1].

Data Types: `char`

#### **InputOffset** — Offset levels present in the input signals of estimation data

[] (default) | vector | matrix

Offset levels present in the input signals of time-domain estimation data, specified as one of the following:

- An `Nu`-element column vector, where `Nu` is the number of inputs. For multi-experiment data, specify a `Nu`-by-`Ne` matrix, where `Ne` is the number of experiments. The offset

value `InputOffset(i, j)` is subtracted from the  $i^{\text{th}}$  input signal of the  $j^{\text{th}}$  experiment.

- `[]` — No offsets.

Data Types: `double`

### **OutputOffset — Output signal offset levels**

`[]` (default) | `vector` | `matrix`

Output signal offset level of time-domain estimation data, specified as one of the following:

- An  $N_y$ -element column vector, where  $N_y$  is the number of outputs. For multi-experiment data, specify a  $N_y$ -by- $N_e$  matrix, where  $N_e$  is the number of experiments. The offset value `OputOffset(i, j)` is subtracted from the  $i^{\text{th}}$  output signal of the  $j^{\text{th}}$  experiment.
- `[]` — No offsets.

The specified values are subtracted from the output signals before using them for estimation.

Data Types: `double`

### **Advanced — Advanced estimation options**

structure

Advanced options for regularized estimation, specified as a structure with the following fields:

- `MaxSize` — Maximum allowable size of Jacobian matrices formed during estimation, specified as a large positive number.

**Default:** `250e3`

- `SearchMethod` — Search method for estimating regularization parameters, specified as one of the following values:
  - `'gn'`: Quasi-Newton line search.
  - `'fmincon'`: Trust-region-reflective constrained minimizer. In general, `'fmincon'` is better than `'gn'` for handling bounds on regularization parameters that are imposed automatically during estimation.

**Default:** `'fmincon'`

## Output Arguments

### **opt** — Regularization options

`arxRegulOptions` options set

Regularization options, returned as an `arxRegulOptions` options set.

## References

- [1] T. Chen, H. Ohlsson, and L. Ljung. “On the Estimation of Transfer Functions, Regularizations and Gaussian Processes - Revisited”, *Automatica*, Volume 48, August 2012.

## See Also

`arxRegul`

## Topics

“Regularized Estimates of Model Parameters”

**Introduced in R2014a**

# arxstruc

Compute loss functions for single-output ARX models

## Syntax

```
V = arxstruc(ze,zv,NN)
```

## Arguments

ze

Estimation data set can be `iddata` or `idfrd` object.

zv

Validation data set can be `iddata` or `idfrd` object.

NN

Matrix defines the number of different ARX-model structures. Each row of NN is of the form:

```
nn = [na nb nk]
```

## Description

---

**Note** Use `arxstruc` for single-output systems only. `arxstruc` supports both single-input and multiple-input systems.

---

`V = arxstruc(ze,zv,NN)` returns `V`, which contains the loss functions in its first row. The remaining rows of `V` contain the transpose of `NN`, so that the orders and delays are given just below the corresponding loss functions. The last column of `V` contains the number of data points in `ze`.

The output argument `V` is best analyzed using `selstruc`. The selection of a suitable model structure based on the information in `v` is normally done using `selstruc`.

## Examples

### Generate Model-Order Combinations and Estimate Single-Input ARX Model

Create an ARX model for generating data.

```
A = [1 -1.5 0.7];  
B = [0 1 0.5];  
m0 = idpoly(A,B);
```

Generate random input and additive noise signals.

```
u = iddata([],idinput(400,'rbs'));  
e = iddata([],0.1*randn(400,1));
```

Simulate the model output using the defined input and error signals.

```
y = sim(m0,[u e]);  
z = [y,u];
```

Generate model-order combinations for estimation. Specify a delay of 1 for all models, and a model order range between 1 and 5 for na and nb.

```
NN = struc(1:5,1:5,1);
```

Estimate ARX models and compute the loss function for each model order combination. The input data is split into estimation and validation data sets.

```
V = arxstruc(z(1:200),z(201:400),NN);
```

Select the model order with the best fit to the validation data.

```
order = selstruc(V,0);
```

Estimate an ARX model of selected order.

```
M = arx(z,order);
```

### Generate Model-Order Combinations and Estimate Multi-Input ARX Model

Create estimation and validation data sets.

```
load co2data;
Ts = 0.5; % Sample time is 0.5 min
ze = iddata(Output_exp1,Input_exp1,Ts);
zv = iddata(Output_exp2,Input_exp2,Ts);
```

Generate model-order combinations for:

- na = 2:4
- nb = 2:5 for the first input, and 1 or 4 for the second input.
- nk = 1:4 for the first input, and 0 for the second input.

```
NN = struc(2:4,2:5,[1 4],1:4,0);
```

Estimate an ARX model for each model order combination.

```
V = arxstruc(ze,zv,NN);
```

Select the model order with the best fit to the validation data.

```
order = selstruc(V,0);
```

Estimate an ARX model of selected order.

```
M = arx(ze,order);
```

## Tips

Each of `ze` and `zv` is an `iddata` object containing output-input data. Frequency-domain data and `idfrd` objects are also supported. Models for each of the model structures defined by `NN` are estimated using the data set `ze`. The loss functions (normalized sum of squared prediction errors) are then computed for these models when applied to the validation data set `zv`. The data sets `ze` and `zv` need not be of equal size. They could, however, be the same sets, in which case the computation is faster.

## See Also

`arx` | `idpoly` | `ivstruc` | `selstruc` | `struc`

**Introduced before R2006a**

## balred

Model order reduction

### Syntax

```
rsys = balred(sys,ORDERS)  
rsys = balred(sys,ORDERS,BALDATA)  
rsys = balred( ___,opts)
```

### Description

*rsys* = balred(*sys*,*ORDERS*) computes a reduced-order approximation *rsys* of the LTI model *sys*. The desired order (number of states) for *rsys* is specified by *ORDERS*. You can try multiple orders at once by setting *ORDERS* to a vector of integers, in which case *rsys* is a vector of reduced-order models. balred uses implicit balancing techniques to compute the reduced-order approximation *rsys*. Use *hsvd* to plot the Hankel singular values and pick an adequate approximation order. States with relatively small Hankel singular values can be safely discarded.

When *sys* has unstable poles, it is first decomposed into its stable and unstable parts using *stabsep*, and only the stable part is approximated. Use *balredOptions* to specify additional options for the stable/unstable decomposition.

This command requires Control System Toolbox™ license. *sys* can only be an identified state-space model (*idss*). The reduced-order model is also an *idss* model.

*rsys* = balred(*sys*,*ORDERS*,*BALDATA*) uses balancing data returned by *hsvd*. Because *hsvd* does most of the work needed to compute *rsys*, this syntax is more efficient when using *hsvd* and balred jointly.

*rsys* = balred( \_\_\_,*opts*) computes the model reduction using options that you specify using *balredOptions*. Options include offset and tolerance options for computing the stable-unstable decompositions. There also options for emphasizing particular time or frequency intervals. See *balredOptions* for details.

---

**Note** The order of the approximate model is always at least the number of unstable poles and at most the minimal order of the original model (number NNZ of nonzero Hankel singular values using an eps-level relative threshold)

---

## Examples

### Reduced-Order Approximation with Offset Option

Compute a reduced-order approximation of the system given by:

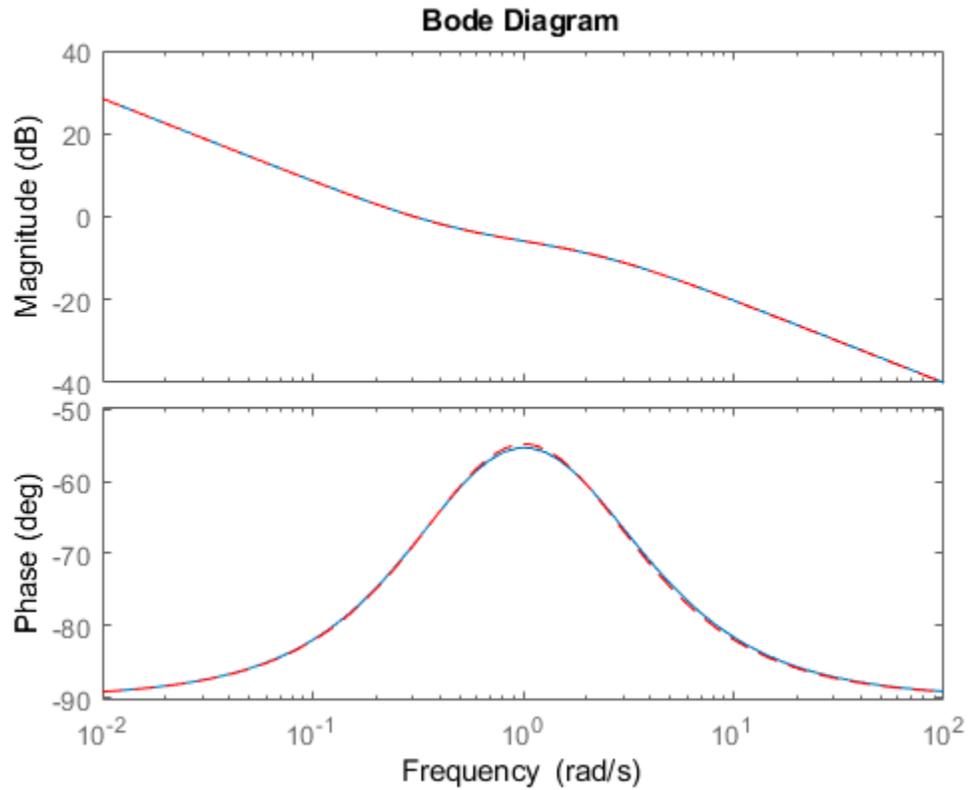
$$G(s) = \frac{(s + 0.5)(s + 1.1)(s + 2.9)}{(s + 10^{-6})(s + 1)(s + 2)(s + 3)}$$

Use the `Offset` option to exclude the pole at  $s = 10^{-6}$  from the stable term of the stable/unstable decomposition.

```
sys = zpk([- .5 -1.1 -2.9],[-1e-6 -2 -1 -3],1);
% Create balredOptions
opt = balredOptions('Offset',.001,'StateElimMethod','Truncate');
% Compute second-order approximation
rsys = balred(sys,2,opt);
```

Compare the responses of the original and reduced-order models.

```
bodeplot(sys,rsys,'r--')
```

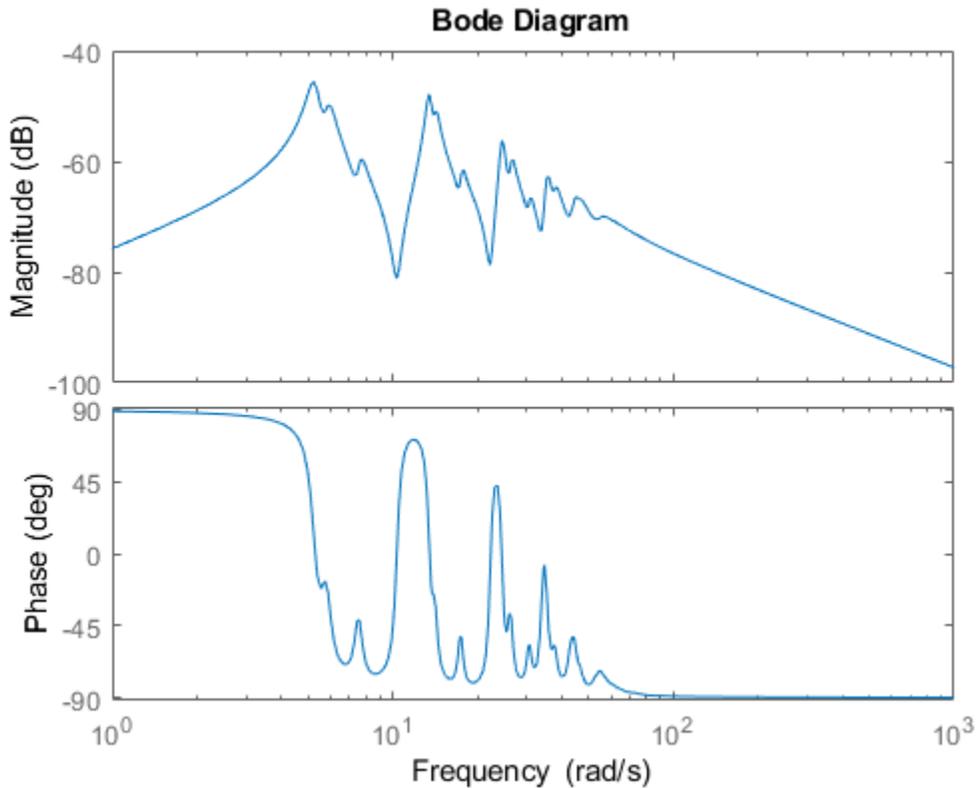


### Model Reduction in a Particular Frequency Band

Reduce a high-order model with a focus on the dynamics in a particular frequency range.

Load a model and examine its frequency response.

```
load(fullfile(matlabroot, 'examples', 'control', 'build.mat'), 'G')  
bodeplot(G)
```

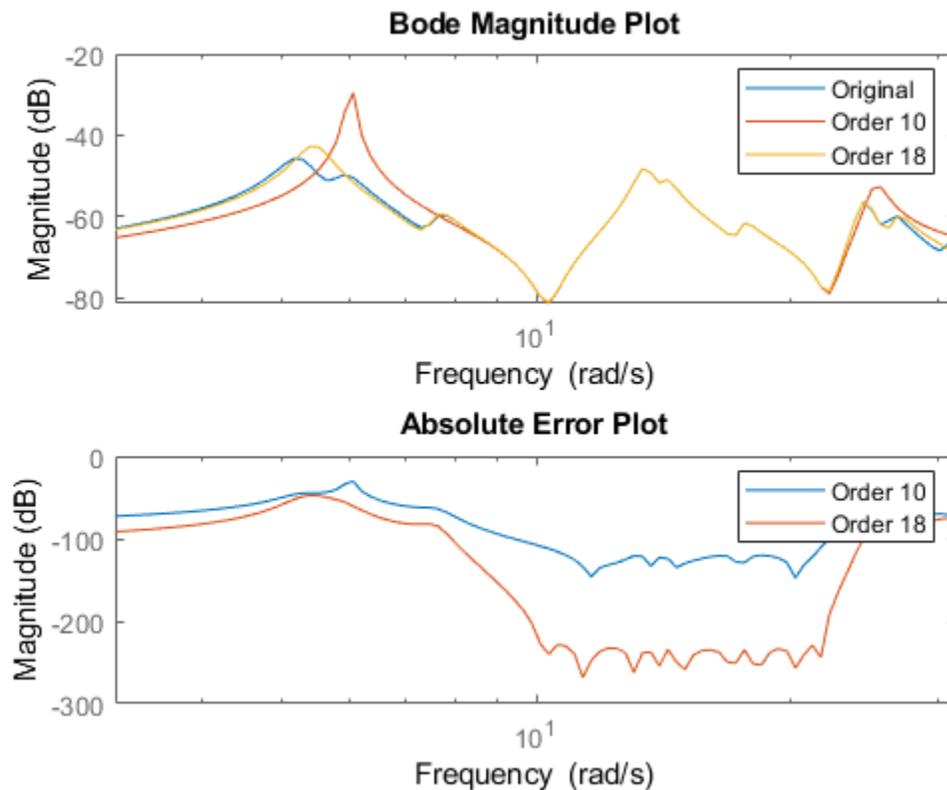


$G$  is a 48th-order model with several large peak regions around 5.2 rad/s, 13.5 rad/s, and 24.5 rad/s, and smaller peaks scattered across many frequencies. Suppose that for your application you are only interested in the dynamics near the second large peak, between 10 rad/s and 22 rad/s. Focus the model reduction on the region of interest to obtain a good match with a low-order approximation. Use `balredOptions` to specify the frequency interval for `balred`.

```
bopt = balredOptions('StateElimMethod','Truncate','FreqIntervals',[10,22]);
GLim10 = balred(G,10,bopt);
GLim18 = balred(G,18,bopt);
```

Examine the frequency responses of the reduced-order models. Also, examine the difference between those responses and the original response (the absolute error).

```
subplot(2,1,1);  
bodemag(G,GLim10,GLim18,logspace(0.5,1.5,100));  
title('Bode Magnitude Plot')  
legend('Original','Order 10','Order 18');  
subplot(2,1,2);  
bodemag(G-GLim10,G-GLim18,logspace(0.5,1.5,100));  
title('Absolute Error Plot')  
legend('Order 10','Order 18');
```



With the frequency-limited energy computation, even the 10th-order approximation is quite good in the region of interest.

## References

- [1] Varga, A., "Balancing-Free Square-Root Algorithm for Computing Singular Perturbation Approximations," Proc. of 30th IEEE CDC, Brighton, UK (1991), pp. 1062-1065.

## See Also

**Model Reducer** | `balredOptions` | `hsvd`

## Topics

"Balanced Truncation Model Reduction" (Control System Toolbox)

"Model Reduction Basics" (Control System Toolbox)

**Introduced before R2006a**

## bandwidth

Frequency response bandwidth

### Syntax

```
fb = bandwidth(sys)
fb = bandwidth(sys,dbdrop)
```

### Description

`fb = bandwidth(sys)` returns the bandwidth of the SISO dynamic system model `sys`. The bandwidth is the first frequency where the gain drops below 70.79% (-3 dB) of its DC value. The bandwidth is expressed in rad/TimeUnit, where TimeUnit is the TimeUnit property of `sys`.

This command requires a Control System Toolbox license.

`fb = bandwidth(sys,dbdrop)` returns the bandwidth for a specified gain drop.

### Examples

#### Compute System Bandwidth

Compute the bandwidth of the transfer function  $sys = 1/(s+1)$ .

```
sys = tf(1,[1 1]);
fb = bandwidth(sys)
```

```
fb = 0.9976
```

This result shows that the gain of `sys` drops to 3 dB below its DC value at around 1 rad/s.

### Find Bandwidth of System with Custom Gain Drop

Compute the frequency at which the gain of a system drops to 3.5 dB below its DC value. Create a state-space model.

```
A = [-2, -1; 1, 0];
B = [1; 0];
C = [1, 2];
D = 1;
sys = ss(A,B,C,D);
```

Find the 3.5 dB bandwidth of `sys`.

```
dbdrop = -3.5;
fb = bandwidth(sys,dbdrop)
```

```
fb = 0.8348
```

### Find Bandwidth of Model Array

Find the bandwidth of each entry in a 5-by-1 array of transfer function models. Use a `for` loop to create the array, and confirm its dimensions.

```
sys = tf(zeros(1,1,5));
s = tf('s');
for m = 1:5
    sys(:, :, m) = m/(s^2+s+m);
end
size(sys)
```

5x1 array of transfer functions.  
Each model has 1 outputs and 1 inputs.

Find the bandwidths.

```
fb = bandwidth(sys)
```

```
fb = 5x1

    1.2712
    1.9991
    2.5298
    2.9678
```

3.3493

`bandwidth` returns an array in which each entry is the bandwidth of the corresponding entry in `sys`. For instance, the bandwidth of `sys(:, :, 2)` is `fb(2)`.

## Input Arguments

### **sys** — Dynamic system

dynamic system model | model array

Dynamic system, specified as a SISO dynamic system model or an array of SISO dynamic system models. Dynamic systems that you can use include:

- Continuous-time or discrete-time numeric LTI models such as `tf`, `zpk`, or `ss` models.
- Frequency-response data models such as `frd` models. For such models, `bandwidth` uses the first frequency point to approximate the DC gain.

If `sys` is an array of models, `bandwidth` returns an array of the same size, where each entry is the bandwidth of the corresponding model in `sys`. For more information on model arrays, see “Model Arrays” (Control System Toolbox).

### **dbdrop** — Gain drop

-3 (default) | negative scalar

Gain drop in dB, specified as a real negative scalar.

## Output Arguments

### **fb** — Frequency response bandwidth

scalar | array

Frequency response bandwidth, returned as a scalar or an array. If `sys` is:

- A single model, then `fb` is the bandwidth of `sys`.
- A model array, then `fb` is an array of the same size as the model array `sys`. Each entry is the bandwidth of the corresponding entry in `sys`.

`fb` is expressed in `rad/TimeUnit`, where `TimeUnit` is the `TimeUnit` property of `sys`.

## See Also

bodeplot | dcgain | issiso

**Introduced before R2006a**

## bj

Estimate Box-Jenkins polynomial model using time domain data

### Syntax

```
sys = bj(data, [nb nc nd nf nk])  
sys = bj(data, [nb nc nd nf nk], Name, Value)  
sys = bj(data, init_sys)  
sys = bj(data, ____, opt)
```

### Description

`sys = bj(data, [nb nc nd nf nk])` estimates a Box-Jenkins polynomial model, `sys`, using the time-domain data, `data`. `[nb nc nd nf nk]` define the orders of the polynomials used for estimation.

`sys = bj(data, [nb nc nd nf nk], Name, Value)` estimates a polynomial model with additional options specified by one or more `Name, Value` pair arguments.

`sys = bj(data, init_sys)` estimates a Box-Jenkins polynomial using the polynomial model `init_sys` to configure the initial parameterization of `sys`.

`sys = bj(data, ____, opt)` estimates a Box-Jenkins polynomial using the option set, `opt`, to specify estimation behavior.

### Input Arguments

#### **data**

Estimation data.

`data` is an `iddata` object that contains time-domain input and output signal values.

You cannot use frequency-domain data for estimating Box-Jenkins models.

**Default:****[nb nc nd nf nk]**

A vector of matrices containing the orders and delays of the Box-Jenkins model. Matrices must contain nonnegative integers.

- `nb` is the order of the B polynomial plus 1 (Ny-by-Nu matrix)
- `nc` is the order of the C polynomial plus 1 (Ny-by-1 matrix)
- `nd` is the order of the D polynomial plus 1 (Ny-by-1 matrix)
- `nf` is the order of the F polynomial plus 1 (Ny-by-Nu matrix)
- `nk` is the input delay (in number of samples, Ny-by-Nu matrix) where Nu is the number of inputs and Ny is the number of outputs.

**opt**

Estimation options.

`opt` is an options set that configures, among others, the following:

- estimation objective
- initial conditions
- numerical search method to be used in estimation

Use `bjOptions` to create the options set.

**init\_sys**

Polynomial model that configures the initial parameterization of `sys`.

`init_sys` must be an `idpoly` model with the Box-Jenkins structure that has only *B*, *C*, *D* and *F* polynomials active. `bj` uses the parameters and constraints defined in `init_sys` as the initial guess for estimating `sys`.

Use the `Structure` property of `init_sys` to configure initial guesses and constraints for  $B(q)$ ,  $F(q)$ ,  $C(q)$  and  $D(q)$ .

To specify an initial guess for, say, the  $C(q)$  term of `init_sys`, set `init_sys.Structure.C.Value` as the initial guess.

To specify constraints for, say, the  $B(q)$  term of `init_sys`:

- set `init_sys.Structure.B.Minimum` to the minimum  $B(q)$  coefficient values
- set `init_sys.Structure.B.Maximum` to the maximum  $B(q)$  coefficient values
- set `init_sys.Structure.B.Free` to indicate which  $B(q)$  coefficients are free for estimation

You can similarly specify the initial guess and constraints for the other polynomials.

## Name-Value Pair Arguments

Specify optional comma-separated pairs of `Name`, `Value` arguments. `Name` is the argument name and `Value` is the corresponding value. `Name` must appear inside quotes. You can specify several name and value pair arguments in any order as `Name1, Value1, . . . , NameN, ValueN`.

### InputDelay

Input delays. `InputDelay` is a numeric vector specifying a time delay for each input channel. Specify input delays in integer multiples of the sample time  $T_s$ . For example, `InputDelay = 3` means a delay of three sampling periods.

For a system with  $N_u$  inputs, set `InputDelay` to an  $N_u$ -by-1 vector, where each entry is a numerical value representing the input delay for the corresponding input channel. You can also set `InputDelay` to a scalar value to apply the same delay to all channels.

**Default:** 0 for all input channels

### IODElay

Transport delays. `IODElay` is a numeric array specifying a separate transport delay for each input/output pair.

Specify transport delays as integers denoting delay of a multiple of the sample time  $T_s$ .

For a MIMO system with  $N_y$  outputs and  $N_u$  inputs, set `IODElay` to a  $N_y$ -by- $N_u$  array, where each entry is a numerical value representing the transport delay for the corresponding input/output pair. You can also set `IODElay` to a scalar value to apply the same delay to all input/output pairs.

**Default:** 0 for all input/output pairs

## IntegrateNoise

Logical specifying integrators in the noise channel.

IntegrateNoise is a logical vector of length  $N_y$ , where  $N_y$  is the number of outputs.

Setting IntegrateNoise to true for a particular output results in the model:

$$y(t) = \frac{B(q)}{F(q)} u(t - nk) + \frac{C(q)}{D(q)} \frac{e(t)}{1 - q^{-1}}$$

Where,  $\frac{1}{1 - q^{-1}}$  is the integrator in the noise channel,  $e(t)$ .

**Default:** false ( $N_y, 1$ ) ( $N_y$  is the number of outputs)

## Output Arguments

### sys

BJ model that fits the estimation data, returned as a discrete-time `idpoly` object. This model is created using the specified model orders, delays, and estimation options.

Information about the estimation results and options used is stored in the `Report` property of the model. `Report` has the following fields:

Report Field	Description
Status	Summary of the model status, which indicates whether the model was created by construction or obtained by estimation.
Method	Estimation command used.

Report Field	Description																		
InitialCondition	<p>Handling of initial conditions during model estimation, returned as one of the following values:</p> <ul style="list-style-type: none"> <li>• 'zero' — The initial conditions were set to zero.</li> <li>• 'estimate' — The initial conditions were treated as independent estimation parameters.</li> <li>• 'backcast' — The initial conditions were estimated using the best least squares fit.</li> </ul> <p>This field is especially useful to view how the initial conditions were handled when the InitialCondition option in the estimation option set is 'auto'.</p>																		
Fit	<p>Quantitative assessment of the estimation, returned as a structure. See “Loss Function and Model Quality Metrics” for more information on these quality metrics. The structure has the following fields:</p> <table border="1" data-bbox="387 829 1332 1322"> <thead> <tr> <th data-bbox="387 829 520 873">Field</th> <th data-bbox="520 829 1332 873">Description</th> </tr> </thead> <tbody> <tr> <td data-bbox="387 873 520 982">FitPercent</td> <td data-bbox="520 873 1332 982">Normalized root mean squared error (NRMSE) measure of how well the response of the model fits the estimation data, expressed as a percentage.</td> </tr> <tr> <td data-bbox="387 982 520 1025">LossFcn</td> <td data-bbox="520 982 1332 1025">Value of the loss function when the estimation completes.</td> </tr> <tr> <td data-bbox="387 1025 520 1104">MSE</td> <td data-bbox="520 1025 1332 1104">Mean squared error (MSE) measure of how well the response of the model fits the estimation data.</td> </tr> <tr> <td data-bbox="387 1104 520 1147">FPE</td> <td data-bbox="520 1104 1332 1147">Final prediction error for the model.</td> </tr> <tr> <td data-bbox="387 1147 520 1190">AIC</td> <td data-bbox="520 1147 1332 1190">Raw Akaike Information Criteria (AIC) measure of model quality.</td> </tr> <tr> <td data-bbox="387 1190 520 1234">AICc</td> <td data-bbox="520 1190 1332 1234">Small sample-size corrected AIC.</td> </tr> <tr> <td data-bbox="387 1234 520 1277">nAIC</td> <td data-bbox="520 1234 1332 1277">Normalized AIC.</td> </tr> <tr> <td data-bbox="387 1277 520 1322">BIC</td> <td data-bbox="520 1277 1332 1322">Bayesian Information Criteria (BIC).</td> </tr> </tbody> </table>	Field	Description	FitPercent	Normalized root mean squared error (NRMSE) measure of how well the response of the model fits the estimation data, expressed as a percentage.	LossFcn	Value of the loss function when the estimation completes.	MSE	Mean squared error (MSE) measure of how well the response of the model fits the estimation data.	FPE	Final prediction error for the model.	AIC	Raw Akaike Information Criteria (AIC) measure of model quality.	AICc	Small sample-size corrected AIC.	nAIC	Normalized AIC.	BIC	Bayesian Information Criteria (BIC).
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nAIC	Normalized AIC.																		
BIC	Bayesian Information Criteria (BIC).																		
Parameters	Estimated values of model parameters.																		
OptionsUsed	Option set used for estimation. If no custom options were configured, this is a set of default options. See <code>bjOptions</code> for more information.																		

Report Field	Description																
RandState	State of the random number stream at the start of estimation. Empty, [], if randomization was not used during estimation. For more information, see <code>rng</code> in the MATLAB documentation.																
DataUsed	Attributes of the data used for estimation, returned as a structure with the following fields:																
	<table border="1"> <thead> <tr> <th data-bbox="385 534 516 569">Field</th> <th data-bbox="516 534 1337 569">Description</th> </tr> </thead> <tbody> <tr> <td data-bbox="385 569 516 618">Name</td> <td data-bbox="516 569 1337 618">Name of the data set.</td> </tr> <tr> <td data-bbox="385 618 516 666">Type</td> <td data-bbox="516 618 1337 666">Data type.</td> </tr> <tr> <td data-bbox="385 666 516 715">Length</td> <td data-bbox="516 666 1337 715">Number of data samples.</td> </tr> <tr> <td data-bbox="385 715 516 756">Ts</td> <td data-bbox="516 715 1337 756">Sample time.</td> </tr> <tr> <td data-bbox="385 756 516 1069">InterSample</td> <td data-bbox="516 756 1337 1069">           Input intersample behavior, returned as one of the following values:           <ul style="list-style-type: none"> <li>• 'zoh' — Zero-order hold maintains a piecewise-constant input signal between samples.</li> <li>• 'foh' — First-order hold maintains a piecewise-linear input signal between samples.</li> <li>• 'bl' — Band-limited behavior specifies that the continuous-time input signal has zero power above the Nyquist frequency.</li> </ul> </td> </tr> <tr> <td data-bbox="385 1069 516 1138">InputOffset</td> <td data-bbox="516 1069 1337 1138">Offset removed from time-domain input data during estimation. For nonlinear models, it is [].</td> </tr> <tr> <td data-bbox="385 1138 516 1218">OutputOffset</td> <td data-bbox="516 1138 1337 1218">Offset removed from time-domain output data during estimation. For nonlinear models, it is [].</td> </tr> </tbody> </table>	Field	Description	Name	Name of the data set.	Type	Data type.	Length	Number of data samples.	Ts	Sample time.	InterSample	Input intersample behavior, returned as one of the following values: <ul style="list-style-type: none"> <li>• 'zoh' — Zero-order hold maintains a piecewise-constant input signal between samples.</li> <li>• 'foh' — First-order hold maintains a piecewise-linear input signal between samples.</li> <li>• 'bl' — Band-limited behavior specifies that the continuous-time input signal has zero power above the Nyquist frequency.</li> </ul>	InputOffset	Offset removed from time-domain input data during estimation. For nonlinear models, it is [].	OutputOffset	Offset removed from time-domain output data during estimation. For nonlinear models, it is [].
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OutputOffset	Offset removed from time-domain output data during estimation. For nonlinear models, it is [].																

Report Field	Description	
Termination	Termination conditions for the iterative search used for prediction error minimization. Structure with the following fields:	
	Field	Description
	WhyStop	Reason for terminating the numerical search.
	Iterations	Number of search iterations performed by the estimation algorithm.
	FirstOrderOptimality	$\infty$ -norm of the gradient search vector when the search algorithm terminates.
	FcnCount	Number of times the objective function was called.
	UpdateNorm	Norm of the gradient search vector in the last iteration. Omitted when the search method is 'lsqnonlin' or 'fmincon'.
	LastImprovement	Criterion improvement in the last iteration, expressed as a percentage. Omitted when the search method is 'lsqnonlin' or 'fmincon'.
	Algorithm	Algorithm used by 'lsqnonlin' or 'fmincon' search method. Omitted when other search methods are used.
For estimation methods that do not require numerical search optimization, the Termination field is omitted.		

For more information on using Report, see “Estimation Report”.

## Examples

### Identify SISO Box-Jenkins Model

Estimate the parameters of a single-input, single-output Box-Jenkins model from measured data.

```
load iddata1 z1;
nb = 2;
```

```
nc = 2;  
nd = 2;  
nf = 2;  
nk = 1;  
sys = bj(z1,[nb nc nd nf nk]);
```

sys is a discrete-time `idpoly` model with estimated coefficients. The order of sys is as described by nb, nc, nd, nf, and nk.

Use `getpvec` to obtain the estimated parameters and `getcov` to obtain the covariance associated with the estimated parameters.

### Estimate a Multi-Input, Single-Output Box-Jenkins Model

Estimate the parameters of a multi-input, single-output Box-Jenkins model from measured data.

```
load iddata8  
nb = [2 1 1];  
nc = 1;  
nd = 1;  
nf = [2 1 2];  
nk = [5 10 15];  
sys = bj(z8,[nb nc nd nf nk]);
```

sys estimates the parameters of a model with three inputs and one output. Each of the inputs has a delay associated with it.

### Estimate Box-Jenkins Model Using Regularization

Estimate a regularized BJ model by converting a regularized ARX model.

Load data.

```
load regularizationExampleData.mat m0simdata;
```

Estimate an unregularized BJ model of order 30.

```
m1 = bj(m0simdata(1:150),[15 15 15 15 1]);
```

Estimate a regularized BJ model by determining Lambda value by trial and error.

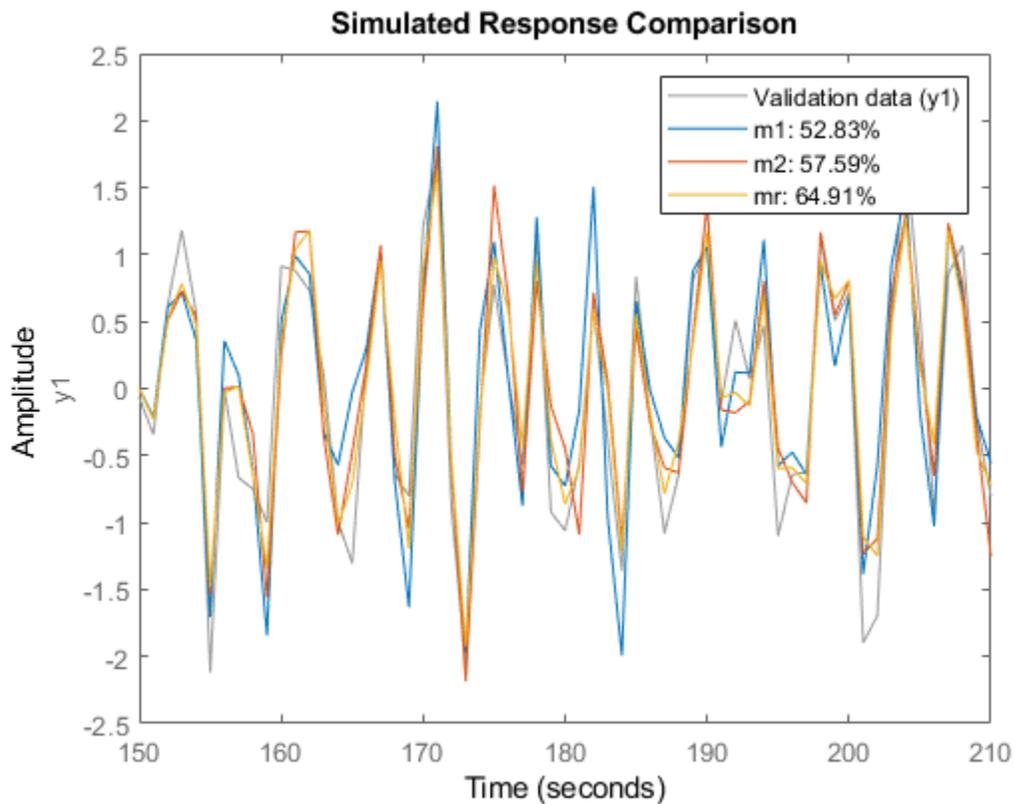
```
opt = bjOptions;  
opt.Regularization.Lambda = 1;  
m2 = bj(m0simdata(1:150),[15 15 15 15 1],opt);
```

Obtain a lower-order BJ model by converting a regularized ARX model followed by order reduction.

```
opt1 = arxOptions;  
[L,R] = arxRegul(m0simdata(1:150),[30 30 1]);  
opt1.Regularization.Lambda = L;  
opt1.Regularization.R = R;  
m0 = arx(m0simdata(1:150),[30 30 1],opt1);  
mr = idpoly(balred(idss(m0),7));
```

Compare the model outputs against data.

```
opt2 = compareOptions('InitialCondition','z');  
compare(m0simdata(150:end),m1,m2,mr,opt2);
```



### Configure Estimation Options

Estimate the parameters of a single-input, single-output Box-Jenkins model while configuring some estimation options.

Generate estimation data.

```
B = [0 1 0.5];  
C = [1 -1 0.2];  
D = [1 1.5 0.7];  
F = [1 -1.5 0.7];
```

```
sys0 = idpoly(1,B,C,D,F,0.1);  
e = iddata([],randn(200,1));  
u = iddata([],idinput(200));  
y = sim(sys0,[u e]);  
data = [y u];
```

`data` is a single-input, single-output data set created by simulating a known model.

Estimate initial Box-Jenkins model.

```
nb = 2;  
nc = 2;  
nd = 2;  
nf = 2;  
nk = 1;  
init_sys = bj(data,[2 2 2 2 1]);
```

Create an estimation option set to refine the parameters of the estimated model.

```
opt = bjOptions;  
opt.Display = 'on';  
opt.SearchOptions.MaxIterations = 50;
```

`opt` is an estimation option set that configures the estimation to iterate 50 times at most and display the estimation progress.

Reestimate the model parameters using the estimation option set.

```
sys = bj(data,init_sys,opt);
```

`sys` is estimated using `init_sys` for the initial parameterization for the polynomial coefficients.

To view the estimation result, enter `sys.Report`.

## **Estimate MIMO Box-Jenkins Model**

Estimate a multi-input, multi-output Box-Jenkins model from estimated data.

Load measured data.

```
load iddata1 z1
load iddata2 z2
data = [z1 z2(1:300)];
```

data contains the measured data for two inputs and two outputs.

Estimate the model.

```
nb = [2 2; 3 4];
nc = [2;2];
nd = [2;2];
nf = [1 0; 2 2];
nk = [1 1; 0 0];
sys = bj(data,[nb nc nd nf nk]);
```

The polynomial order coefficients contain one row for each output.

sys is a discrete-time idpoly model with two inputs and two outputs.

## Definitions

### Box-Jenkins Model Structure

The general Box-Jenkins model structure is:

$$y(t) = \sum_{i=1}^{nu} \frac{B_i(q)}{F_i(q)} u_i(t - nk_i) + \frac{C(q)}{D(q)} e(t)$$

where  $nu$  is the number of input channels.

The orders of Box-Jenkins model are defined as follows:

$$nb: B(q) = b_1 + b_2q^{-1} + \dots + b_{nb}q^{-nb+1}$$

$$nc: C(q) = 1 + c_1q^{-1} + \dots + c_{nc}q^{-nc}$$

$$nd: D(q) = 1 + d_1q^{-1} + \dots + d_{nd}q^{-nd}$$

$$nf: F(q) = 1 + f_1q^{-1} + \dots + f_{nf}q^{-nf}$$

## Alternatives

To estimate a continuous-time model, use:

- `tfest` — returns a transfer function model
- `ssest` — returns a state-space model
- `bj` to first estimate a discrete-time model and convert it a continuous-time model using `d2c`.

## References

- [1] Ljung, L. *System Identification: Theory for the User*, Upper Saddle River, NJ, Prentice-Hall PTR, 1999.

## Extended Capabilities

### Automatic Parallel Support

Accelerate code by automatically running computation in parallel using Parallel Computing Toolbox™.

Parallel computing support is available for estimation using the `lsqnonlin` search method (requires Optimization Toolbox). To enable parallel computing, use `bjOptions`, set `SearchMethod` to `'lsqnonlin'`, and set `SearchOptions.Advanced.UseParallel` to `true`.

For example:

```
opt = bjOptions;  
opt.SearchMethod = 'lsqnonlin';  
opt.SearchOptions.Advanced.UseParallel = true;
```

### See Also

[armax](#) | [arx](#) | [bjoptions](#) | [compare](#) | [d2c](#) | [forecast](#) | [iddata](#) | [idpoly](#) | [iv4](#) | [oe](#) | [polyest](#) | [sim](#) | [ssest](#) | [tfest](#)

## **Topics**

“Regularized Estimates of Model Parameters”

**Introduced before R2006a**

## bjOptions

Option set for bj

### Syntax

```
opt = bjOptions  
opt = bjOptions(Name,Value)
```

### Description

`opt = bjOptions` creates the default options set for `bj`.

`opt = bjOptions(Name,Value)` creates an option set with the options specified by one or more `Name,Value` pair arguments.

### Input Arguments

#### Name-Value Pair Arguments

Specify optional comma-separated pairs of `Name,Value` arguments. `Name` is the argument name and `Value` is the corresponding value. `Name` must appear inside quotes. You can specify several name and value pair arguments in any order as `Name1,Value1,...,NameN,ValueN`.

#### InitialCondition — Handling of initial conditions

'auto' (default) | 'zero' | 'estimate' | 'backcast'

Handling of initial conditions during estimation, specified as one of the following values:

- 'zero' — The initial conditions are set to zero.
- 'estimate' — The initial conditions are treated as independent estimation parameters.
- 'backcast' — The initial conditions are estimated using the best least squares fit.

- 'auto' — The software chooses the method to handle initial conditions based on the estimation data.

### Focus — Error to be minimized

'prediction' (default) | 'simulation'

Error to be minimized in the loss function during estimation, specified as the comma-separated pair consisting of 'Focus' and one of the following values:

- 'prediction' — The one-step ahead prediction error between measured and predicted outputs is minimized during estimation. As a result, the estimation focuses on producing a good predictor model.
- 'simulation' — The simulation error between measured and simulated outputs is minimized during estimation. As a result, the estimation focuses on making a good fit for simulation of model response with the current inputs.

The Focus option can be interpreted as a weighting filter in the loss function. For more information, see “Loss Function and Model Quality Metrics”.

### WeightingFilter — Weighting prefilter

[] (default) | vector | matrix | cell array | linear system

Weighting prefilter applied to the loss function to be minimized during estimation. To understand the effect of WeightingFilter on the loss function, see “Loss Function and Model Quality Metrics”.

Specify WeightingFilter as one of the following values:

- [] — No weighting prefilter is used.
- Passbands — Specify a row vector or matrix containing frequency values that define desired passbands. You select a frequency band where the fit between estimated model and estimation data is optimized. For example, [w<sub>l</sub>, w<sub>h</sub>], where w<sub>l</sub> and w<sub>h</sub> represent lower and upper limits of a passband. For a matrix with several rows defining frequency passbands, [w<sub>1l</sub>, w<sub>1h</sub>; w<sub>2l</sub>, w<sub>2h</sub>; w<sub>3l</sub>, w<sub>3h</sub>; . . . ], the estimation algorithm uses the union of the frequency ranges to define the estimation passband.

Passbands are expressed in rad/TimeUnit for time-domain data and in FrequencyUnit for frequency-domain data, where TimeUnit and FrequencyUnit are the time and frequency units of the estimation data.

- SISO filter — Specify a single-input-single-output (SISO) linear filter in one of the following ways:

- A SISO LTI model
- $\{A, B, C, D\}$  format, which specifies the state-space matrices of a filter with the same sample time as estimation data.
- $\{\text{numerator}, \text{denominator}\}$  format, which specifies the numerator and denominator of the filter as a transfer function with same sample time as estimation data.

This option calculates the weighting function as a product of the filter and the input spectrum to estimate the transfer function.

### **EnforceStability — Control whether to enforce stability of model**

`false` (default) | `true`

Control whether to enforce stability of estimated model, specified as the comma-separated pair consisting of 'EnforceStability' and either `true` or `false`.

Data Types: `logical`

### **EstimateCovariance — Control whether to generate parameter covariance data**

`true` (default) | `false`

Controls whether parameter covariance data is generated, specified as `true` or `false`.

If `EstimateCovariance` is `true`, then use `getcov` to fetch the covariance matrix from the estimated model.

### **Display — Specify whether to display the estimation progress**

`'off'` (default) | `'on'`

Specify whether to display the estimation progress, specified as one of the following values:

- `'on'` — Information on model structure and estimation results are displayed in a progress-viewer window.
- `'off'` — No progress or results information is displayed.

### **InputOffset — Removal of offset from time-domain input data during estimation**

`[]` (default) | vector of positive integers | matrix

Removal of offset from time-domain input data during estimation, specified as the comma-separated pair consisting of 'InputOffset' and one of the following:

- A column vector of positive integers of length  $Nu$ , where  $Nu$  is the number of inputs.
- `[]` — Indicates no offset.
- $Nu$ -by- $Ne$  matrix — For multi-experiment data, specify `InputOffset` as an  $Nu$ -by- $Ne$  matrix.  $Nu$  is the number of inputs, and  $Ne$  is the number of experiments.

Each entry specified by `InputOffset` is subtracted from the corresponding input data.

### **OutputOffset — Removal of offset from time-domain output data during estimation**

`[]` (default) | vector | matrix

Removal of offset from time-domain output data during estimation, specified as the comma-separated pair consisting of `'OutputOffset'` and one of the following:

- A column vector of length  $Ny$ , where  $Ny$  is the number of outputs.
- `[]` — Indicates no offset.
- $Ny$ -by- $Ne$  matrix — For multi-experiment data, specify `OutputOffset` as a  $Ny$ -by- $Ne$  matrix.  $Ny$  is the number of outputs, and  $Ne$  is the number of experiments.

Each entry specified by `OutputOffset` is subtracted from the corresponding output data.

### **Regularization — Options for regularized estimation of model parameters**

structure

Options for regularized estimation of model parameters. For more information on regularization, see “Regularized Estimates of Model Parameters”.

`Regularization` is a structure with the following fields:

- `Lambda` — Constant that determines the bias versus variance tradeoff.

Specify a positive scalar to add the regularization term to the estimation cost.

The default value of zero implies no regularization.

#### **Default:** 0

- `R` — Weighting matrix.

Specify a vector of nonnegative numbers or a square positive semi-definite matrix. The length must be equal to the number of free parameters of the model.

For black-box models, using the default value is recommended. For structured and grey-box models, you can also specify a vector of `np` positive numbers such that each entry denotes the confidence in the value of the associated parameter.

The default value of 1 implies a value of `eye(npfree)`, where `npfree` is the number of free parameters.

**Default:** 1

- **Nominal** — The nominal value towards which the free parameters are pulled during estimation.

The default value of zero implies that the parameter values are pulled towards zero. If you are refining a model, you can set the value to `'model'` to pull the parameters towards the parameter values of the initial model. The initial parameter values must be finite for this setting to work.

**Default:** 0

**SearchMethod** — Numerical search method used for iterative parameter estimation

`'auto'` (default) | `'gn'` | `'gna'` | `'lm'` | `'grad'` | `'lsqnonlin'` | `'fmincon'`

Numerical search method used for iterative parameter estimation, specified as the comma-separated pair consisting of `'SearchMethod'` and one of the following:

- `'auto'` — A combination of the line search algorithms, `'gn'`, `'lm'`, `'gna'`, and `'grad'` methods is tried in sequence at each iteration. The first descent direction leading to a reduction in estimation cost is used.
- `'gn'` — Subspace Gauss-Newton least squares search. Singular values of the Jacobian matrix less than `GnPinvConstant*eps*max(size(J))*norm(J)` are discarded when computing the search direction.  $J$  is the Jacobian matrix. The Hessian matrix is approximated as  $J^T J$ . If there is no improvement in this direction, the function tries the gradient direction.
- `'gna'` — Adaptive subspace Gauss-Newton search. Eigenvalues less than `gamma*max(sv)` of the Hessian are ignored, where `sv` contains the singular values of the Hessian. The Gauss-Newton direction is computed in the remaining subspace. `gamma` has the initial value `InitialGnaTolerance` (see `Advanced in 'SearchOptions'` for more information). This value is increased by the factor `LMStep` each time the search fails to find a lower value of the criterion in fewer than five bisections. This value is decreased by the factor `2*LMStep` each time a search is successful without any bisections.

- 'lm' — Levenberg-Marquardt least squares search, where the next parameter value is  $-\text{pinv}(H+d*I)*\text{grad}$  from the previous one.  $H$  is the Hessian,  $I$  is the identity matrix, and  $\text{grad}$  is the gradient.  $d$  is a number that is increased until a lower value of the criterion is found.
- 'grad' — Steepest descent least squares search.
- 'lsqnonlin' — Trust-region-reflective algorithm of `lsqnonlin`. Requires Optimization Toolbox software.
- 'fmincon' — Constrained nonlinear solvers. You can use the sequential quadratic programming (SQP) and trust-region-reflective algorithms of the `fmincon` solver. If you have Optimization Toolbox software, you can also use the interior-point and active-set algorithms of the `fmincon` solver. Specify the algorithm in the `SearchOptions.Algorithm` option. The `fmincon` algorithms may result in improved estimation results in the following scenarios:
  - Constrained minimization problems when there are bounds imposed on the model parameters.
  - Model structures where the loss function is a nonlinear or non smooth function of the parameters.
  - Multi-output model estimation. A determinant loss function is minimized by default for multi-output model estimation. `fmincon` algorithms are able to minimize such loss functions directly. The other search methods such as 'lm' and 'gn' minimize the determinant loss function by alternately estimating the noise variance and reducing the loss value for a given noise variance value. Hence, the `fmincon` algorithms can offer better efficiency and accuracy for multi-output model estimations.

### **SearchOptions — Option set for the search algorithm**

search option set

Option set for the search algorithm, specified as the comma-separated pair consisting of 'SearchOptions' and a search option set with fields that depend on the value of SearchMethod.

**SearchOptions Structure When SearchMethod is Specified as 'gn', 'gna', 'lm', 'grad', or 'auto'**

Field Name	Description	Default
Tolerance	Minimum percentage difference between the current value of the loss function and its expected improvement after the next iteration, specified as a positive scalar. When the percentage of expected improvement is less than <code>Tolerance</code> , the iterations stop. The estimate of the expected loss-function improvement at the next iteration is based on the Gauss-Newton vector computed for the current parameter value.	0.01
MaxIterations	<p>Maximum number of iterations during loss-function minimization, specified as a positive integer. The iterations stop when <code>MaxIterations</code> is reached or another stopping criterion is satisfied, such as <code>Tolerance</code>.</p> <p>Setting <code>MaxIterations = 0</code> returns the result of the start-up procedure.</p> <p>Use <code>sys.Report.Termination.Iterations</code> to get the actual number of iterations during an estimation, where <code>sys</code> is an <code>idtf</code> model.</p>	20

Field Name	Description	Default
Advanced	Advanced search settings, specified as a structure with the following fields:	
Field Name	Description	Default
GnPinvConstant	Jacobian matrix singular value threshold, specified as a positive scalar. Singular values of the Jacobian matrix that are smaller than $\text{GnPinvConstant} * \max(\text{size}(J) * \text{norm}(J) * \text{eps})$ are discarded when computing the search direction. Applicable when SearchMethod is 'gn'.	10000
InitialGnaTolerance	Initial value of $\gamma$ , specified as a positive scalar. Applicable when SearchMethod is 'gna'.	0.0001
LMStartValue	Starting value of search-direction length $d$ in the Levenberg-Marquardt method, specified as a positive scalar. Applicable when SearchMethod is 'lm'.	0.001
LMStep	Size of the Levenberg-Marquardt step, specified as a positive integer. The next value of the search-direction length $d$ in the Levenberg-Marquardt method is LMStep times the previous one. Applicable when SearchMethod is 'lm'.	2
MaxBisections	Maximum number of bisections used for line search along the search direction, specified as a positive integer.	25
MaxFunctionEvaluations	Maximum number of calls to the model file, specified as a positive integer. Iterations stop if the number of calls to the model file exceeds this value.	Inf
MinParameterChange	Smallest parameter update allowed per iteration, specified as a nonnegative scalar.	0
RelativeImprovement	Relative improvement threshold, specified as a nonnegative scalar. Iterations stop if the relative	0

<b>Field Name</b>	<b>Description</b>		<b>Default</b>
	<b>Field Name</b>	<b>Description</b>	<b>Default</b>
		improvement of the criterion function is less than this value.	
	StepReduction	<p>Step reduction factor, specified as a positive scalar that is greater than 1. The suggested parameter update is reduced by the factor <code>StepReduction</code> after each try. This reduction continues until <code>MaxBisections</code> tries are completed or a lower value of the criterion function is obtained.</p> <p><code>StepReduction</code> is not applicable for SearchMethod 'lm' (Levenberg-Marquardt method).</p>	2

**SearchOptions Structure When SearchMethod is Specified as 'lsqnonlin'**

<b>Field Name</b>	<b>Description</b>	<b>Default</b>
FunctionTolerance	Termination tolerance on the loss function that the software minimizes to determine the estimated parameter values, specified as a positive scalar.  The value of FunctionTolerance is the same as that of <code>opt.SearchOptions.Advanced.TolFun</code> .	1e-5
StepTolerance	Termination tolerance on the estimated parameter values, specified as a positive scalar.  The value of StepTolerance is the same as that of <code>opt.SearchOptions.Advanced.TolX</code> .	1e-6
MaxIterations	Maximum number of iterations during loss-function minimization, specified as a positive integer. The iterations stop when MaxIterations is reached or another stopping criterion is satisfied, such as FunctionTolerance.  The value of MaxIterations is the same as that of <code>opt.SearchOptions.Advanced.MaxIter</code> .	20
Advanced	Advanced search settings, specified as an option set for <code>lsqnonlin</code> .  For more information, see the Optimization Options table in "Optimization Options" (Optimization Toolbox).	Use <code>optimset('lsqnonlin')</code> to create a default option set.

**SearchOptions Structure When SearchMethod is Specified as 'fmincon'**

Field Name	Description	Default
Algorithm	<p>fmincon optimization algorithm, specified as one of the following:</p> <ul style="list-style-type: none"> <li>• 'sqp' — Sequential quadratic programming algorithm. The algorithm satisfies bounds at all iterations, and it can recover from NaN or Inf results. It is not a large-scale algorithm. For more information, see “Large-Scale vs. Medium-Scale Algorithms” (Optimization Toolbox).</li> <li>• 'trust-region-reflective' — Subspace trust-region method based on the interior-reflective Newton method. It is a large-scale algorithm.</li> <li>• 'interior-point' — Large-scale algorithm that requires Optimization Toolbox software. The algorithm satisfies bounds at all iterations, and it can recover from NaN or Inf results.</li> <li>• 'active-set' — Requires Optimization Toolbox software. The</li> </ul>	'sqp'

Field Name	Description	Default
	<p>algorithm can take large steps, which adds speed. It is not a large-scale algorithm.</p> <p>For more information about the algorithms, see “Constrained Nonlinear Optimization Algorithms” (Optimization Toolbox) and “Choosing the Algorithm” (Optimization Toolbox).</p>	
FunctionTolerance	Termination tolerance on the loss function that the software minimizes to determine the estimated parameter values, specified as a positive scalar.	1e-6
StepTolerance	Termination tolerance on the estimated parameter values, specified as a positive scalar.	1e-6
MaxIterations	Maximum number of iterations during loss function minimization, specified as a positive integer. The iterations stop when MaxIterations is reached or another stopping criterion is satisfied, such as FunctionTolerance.	100

### Advanced – Additional advanced options structure

Additional advanced options, specified as a structure with the following fields:

- `ErrorThreshold` — Specifies when to adjust the weight of large errors from quadratic to linear.

Errors larger than `ErrorThreshold` times the estimated standard deviation have a linear weight in the loss function. The standard deviation is estimated robustly as the median of the absolute deviations from the median of the prediction errors, divided by 0.7. For more information on robust norm choices, see section 15.2 of [2].

`ErrorThreshold = 0` disables robustification and leads to a purely quadratic loss function. When estimating with frequency-domain data, the software sets `ErrorThreshold` to zero. For time-domain data that contains outliers, try setting `ErrorThreshold` to 1.6.

**Default:** 0

- `MaxSize` — Specifies the maximum number of elements in a segment when input-output data is split into segments.

`MaxSize` must be a positive integer.

**Default:** 250000

- `StabilityThreshold` — Specifies thresholds for stability tests.

`StabilityThreshold` is a structure with the following fields:

- `s` — Specifies the location of the right-most pole to test the stability of continuous-time models. A model is considered stable when its right-most pole is to the left of `s`.

**Default:** 0

- `z` — Specifies the maximum distance of all poles from the origin to test stability of discrete-time models. A model is considered stable if all poles are within the distance `z` from the origin.

**Default:** `1+sqrt(eps)`

- `AutoInitThreshold` — Specifies when to automatically estimate the initial condition.

The initial condition is estimated when

$$\frac{\|y_{p,z} - y_{meas}\|}{\|y_{p,e} - y_{meas}\|} > \text{AutoInitThreshold}$$

- $y_{meas}$  is the measured output.
- $y_{p,z}$  is the predicted output of a model estimated using zero initial states.
- $y_{p,e}$  is the predicted output of a model estimated using estimated initial states.

Applicable when `InitialCondition` is 'auto'.

**Default:** 1.05

## Output Arguments

### **opt** — Options set for bj

bjOptions option set

Option set for bj, returned as an bjOptions option set.

## Examples

### **Create Default Options Set for Box-Jenkins Estimation**

```
opt = bjOptions;
```

### **Specify Options for Box-Jenkins Estimation**

Create an options set for bj using zero initial conditions for estimation. Set `Display` to 'on'.

```
opt = bjOptions('InitialCondition','zero','Display','on');
```

Alternatively, use dot notation to set the values of opt.

```
opt = bjOptions;  
opt.InitialCondition = 'zero';  
opt.Display = 'on';
```

## References

- [1] Wills, Adrian, B. Ninness, and S. Gibson. "On Gradient-Based Search for Multivariable System Estimates". *Proceedings of the 16th IFAC World Congress, Prague, Czech Republic, July 3-8, 2005*. Oxford, UK: Elsevier Ltd., 2005.
- [2] Ljung, L. *System Identification: Theory for the User*. Upper Saddle River, NJ: Prentice-Hall PTR, 1999.

## See Also

bj | idfilt

## Topics

"Loss Function and Model Quality Metrics"

**Introduced in R2012a**

# blkdiag

Block-diagonal concatenation of models

## Syntax

```
sys = blkdiag(sys1,sys2,...,sysN)
```

## Description

`sys = blkdiag(sys1,sys2,...,sysN)` produces the aggregate system

$$\begin{bmatrix} \text{sys1} & 0 & \dots & 0 \\ 0 & \text{sys2} & \dots & \vdots \\ \vdots & \dots & \dots & 0 \\ 0 & \dots & 0 & \text{sysN} \end{bmatrix}$$

`blkdiag` is equivalent to `append`.

## Examples

### Perform Block-Diagonal Concatenation

Perform block-diagonal concatenation of a transfer function model and a state-space model.

Create the SISO continuous-time transfer function model,  $1/s$ .

```
sys1 = tf(1,[1 0]);
```

Create a SISO continuous-time state-space model with state-space matrices 1,2,3, and 4.

```
sys2 = ss(1,2,3,4);
```

Concatenate `sys1`, a SISO static gain system, and `sys2`. The resulting model is a 3-input, 3-output state-space model.

```
sys = blkdiag(sys1,10,sys2)
```

```
sys =
```

```
A =
```

```
      x1  x2
x1    0   0
x2    0   1
```

```
B =
```

```
      u1  u2  u3
x1    1   0   0
x2    0   0   2
```

```
C =
```

```
      x1  x2
y1    1   0
y2    0   0
y3    0   3
```

```
D =
```

```
      u1  u2  u3
y1    0   0   0
y2    0  10   0
y3    0   0   4
```

Continuous-time state-space model.

Alternatively, use the `append` command.

```
sys = append(sys1,10,sys2);
```

## See Also

`append` | `feedback` | `parallel` | `series`

**Introduced in R2009a**

# bode

Bode plot of frequency response, or magnitude and phase data

## Syntax

```
bode(sys)
bode(sys1,sys2,...,sysN)
bode(sys1,LineStyle1,...,sysN,LineStyleN)
bode( ____,w)
```

```
[mag,phase,wout] = bode(sys)
[mag,phase,wout] = bode(sys,w)
[mag,phase,wout,sdmag,sdphase] = bode(sys,w)
```

## Description

`bode(sys)` creates a Bode plot of the frequency response of a dynamic system model `sys`. The plot displays the magnitude (in dB) and phase (in degrees) of the system response as a function of frequency. `bode` automatically determines frequencies to plot based on system dynamics.

If `sys` is a multi-input, multi-output (MIMO) model, then `bode` produces an array of Bode plots, each plot showing the frequency response of one I/O pair.

`bode(sys1,sys2,...,sysN)` plots the frequency response of multiple dynamic systems on the same plot. All systems must have the same number of inputs and outputs.

`bode(sys1,LineStyle1,...,sysN,LineStyleN)` specifies a color, linestyle, and marker for each system in the plot.

`bode( ____,w)` plots system responses for frequencies specified by `w`.

- If `w` is a cell array of the form `{wmin,wmax}`, then `bode` plots the response at frequencies ranging between `wmin` and `wmax`.
- If `w` is a vector of frequencies, then `bode` plots the response at each specified frequency.

You can use `w` with any of the input-argument combinations in previous syntaxes.

`[mag, phase, wout] = bode(sys)` returns the magnitude and phase of the response at each frequency in the vector `wout`. The function automatically determines frequencies in `wout` based on system dynamics. This syntax does not draw a plot.

`[mag, phase, wout] = bode(sys, w)` returns the response data at the frequencies specified by `w`.

- If `w` is a cell array of the form `{wmin, wmax}`, then `wout` contains frequencies ranging between `wmin` and `wmax`.
- If `w` is a vector of frequencies, then `wout = w`.

`[mag, phase, wout, sdmag, sdphase] = bode(sys, w)` also returns the estimated standard deviation of the magnitude and phase values for the identified model `sys`. If you omit `w`, then the function automatically determines frequencies in `wout` based on system dynamics.

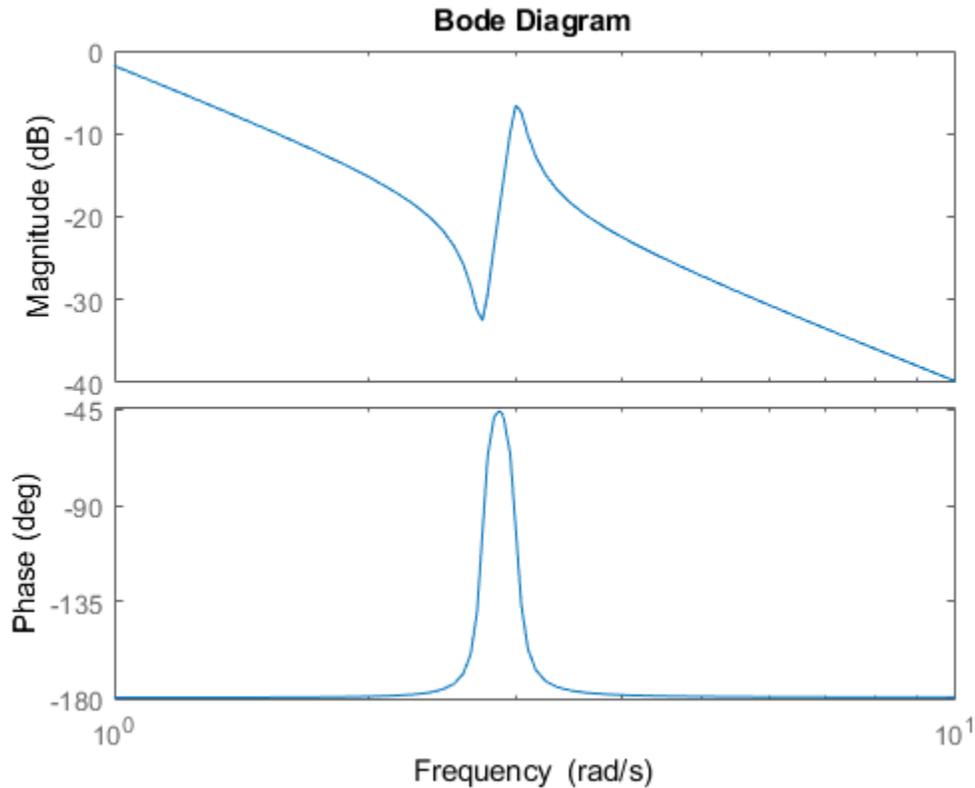
## Examples

### Bode Plot of Dynamic System

Create a Bode plot of the following continuous-time SISO dynamic system.

$$H(s) = \frac{s^2 + 0.1s + 7.5}{s^4 + 0.12s^3 + 9s^2}$$

```
H = tf([1 0.1 7.5],[1 0.12 9 0 0]);  
bode(H)
```

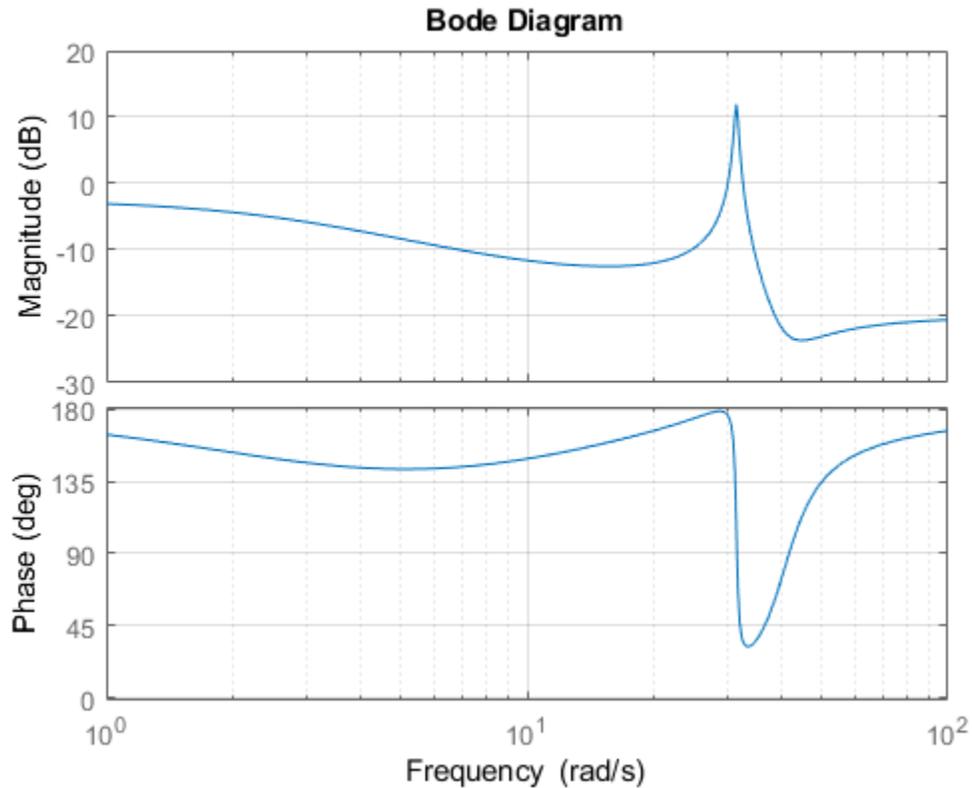


bode automatically selects the plot range based on the system dynamics.

### Bode Plot at Specified Frequencies

Create a Bode plot over a specified frequency range. Use this approach when you want to focus on the dynamics in a particular range of frequencies.

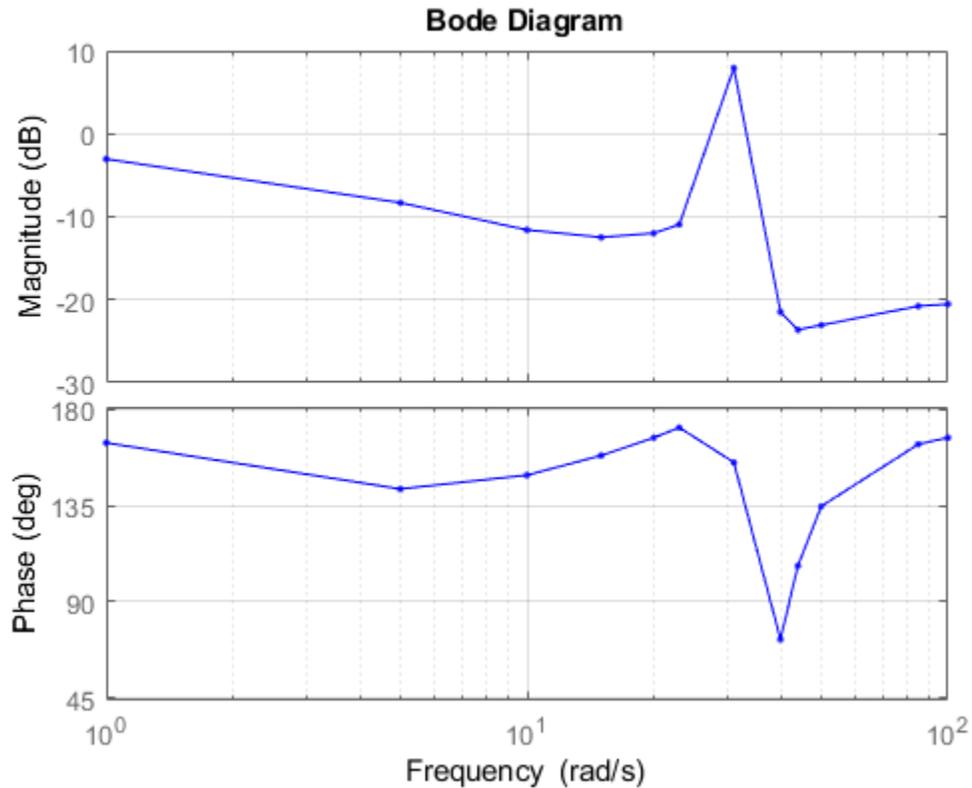
```
H = tf([-0.1, -2.4, -181, -1950], [1, 3.3, 990, 2600]);  
bode(H, {1, 100})  
grid on
```



The cell array `{1, 100}` specifies the minimum and maximum frequency values in the Bode plot. When you provide frequency bounds in this way, the function selects intermediate points for frequency response data.

Alternatively, specify a vector of frequency points to use for evaluating and plotting the frequency response.

```
w = [1 5 10 15 20 23 31 40 44 50 85 100];  
bode(H,w, '.-')  
grid on
```



bode plots the frequency response at the specified frequencies only.

### Compare Bode Plots of Several Dynamic Systems

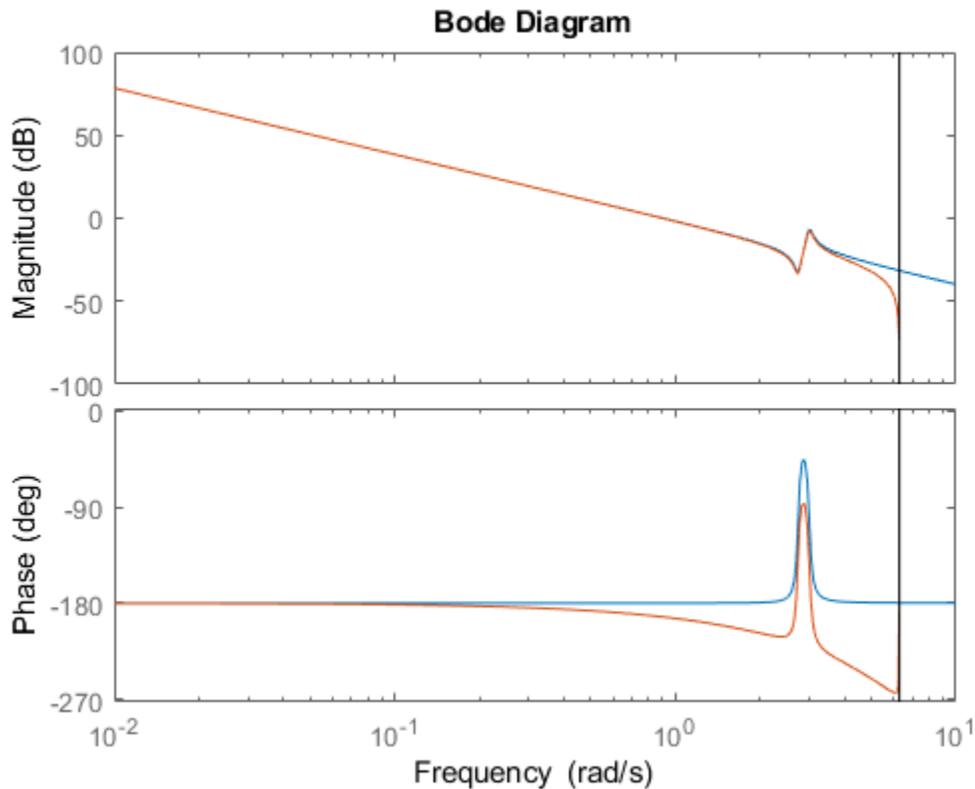
Compare the frequency response of a continuous-time system to an equivalent discretized system on the same Bode plot.

Create continuous-time and discrete-time dynamic systems.

```
H = tf([1 0.1 7.5],[1 0.12 9 0 0]);
Hd = c2d(H,0.5,'zoh');
```

Create a Bode plot that displays both systems.

```
bode(H,Hd)
```

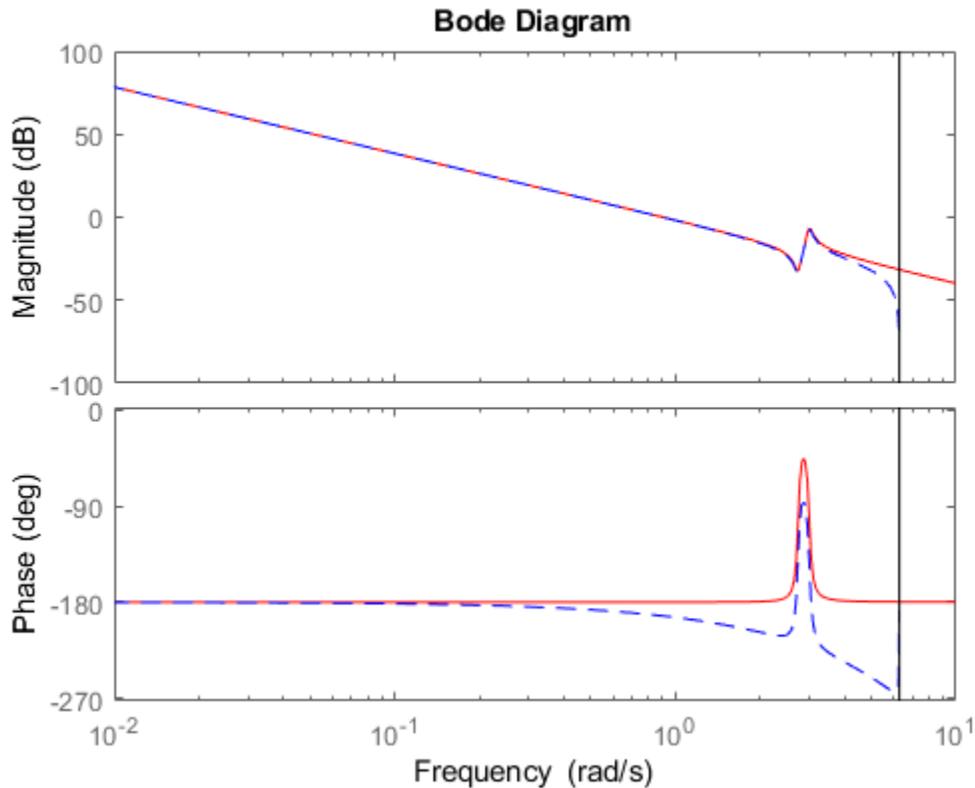


The Bode plot of a discrete-time system includes a vertical line marking the Nyquist frequency of the system.

### Bode Plot with Specified Line Attributes

Specify the line style, color, or marker for each system in a Bode plot using the LineSpec input argument.

```
H = tf([1 0.1 7.5],[1 0.12 9 0 0]);  
Hd = c2d(H,0.5,'zoh');  
bode(H,'r',Hd,'b--')
```



The first LineSpec, 'r', specifies a solid red line for the response of H. The second LineSpec, 'b--', specifies a dashed blue line for the response of Hd.

### Obtain Magnitude and Phase Data

Compute the magnitude and phase of the frequency response of a SISO system.

If you do not specify frequencies, `bode` chooses frequencies based on the system dynamics and returns them in the third output argument.

```
H = tf([1 0.1 7.5],[1 0.12 9 0 0]);  
[mag,phase,wout] = bode(H);
```

Because `H` is a SISO model, the first two dimensions of `mag` and `phase` are both 1. The third dimension is the number of frequencies in `wout`.

```
size(mag)
```

```
ans = 1×3
```

```
     1     1    42
```

```
length(wout)
```

```
ans = 42
```

Thus, each entry along the third dimension of `mag` gives the magnitude of the response at the corresponding frequency in `wout`.

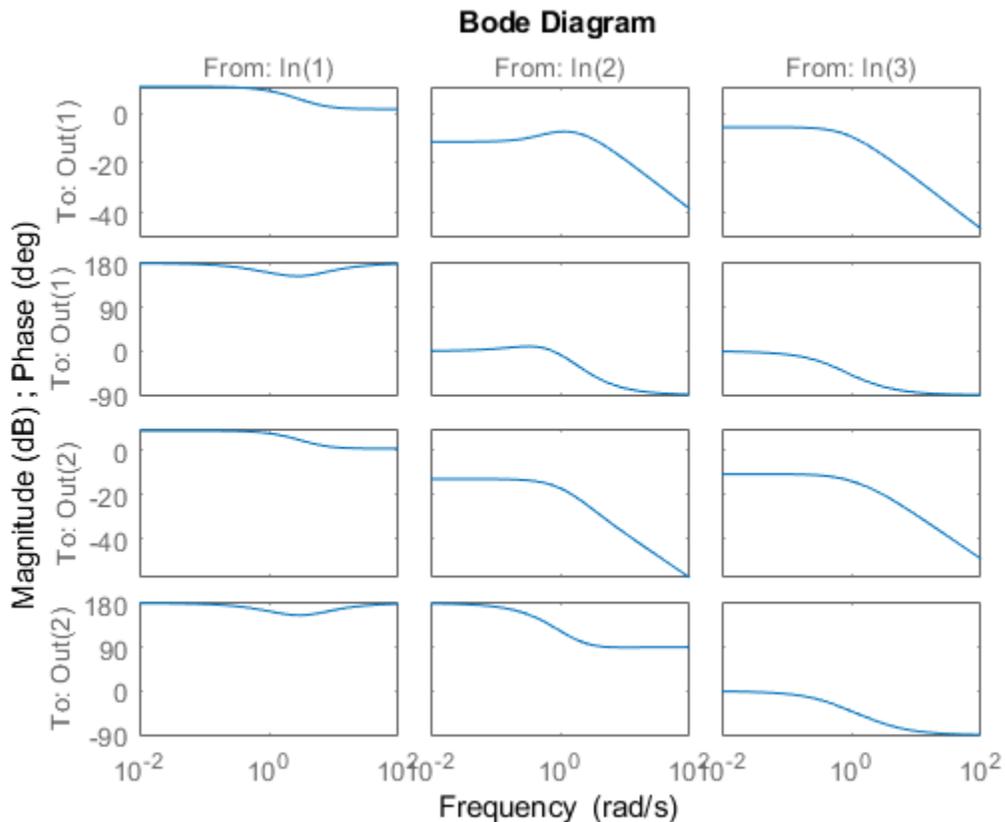
### Magnitude and Phase of MIMO System

For this example, create a 2-output, 3-input system.

```
rng(0, 'twister'); % For reproducibility  
H = rss(4,2,3);
```

For this system, `bode` plots the frequency responses of each I/O channel in a separate plot in a single figure.

```
bode(H)
```



Compute the magnitude and phase of these responses at 20 frequencies between 1 and 10 radians.

```
w = logspace(0,1,20);
[mag,phase] = bode(H,w);
```

mag and phase are three-dimensional arrays, in which the first two dimensions correspond to the output and input dimensions of H, and the third dimension is the number of frequencies. For instance, examine the dimensions of mag.

```
size(mag)
```

```
ans = 1x3
```

2      3      20

Thus, for example, `mag(1,3,10)` is the magnitude of the response from the third input to the first output, computed at the 10th frequency in `w`. Similarly, `phase(1,3,10)` contains the phase of the same response.

## Bode Plot of Identified Model

Compare the frequency response of a parametric model, identified from input/output data, to a nonparametric model identified using the same data.

Identify parametric and nonparametric models based on data.

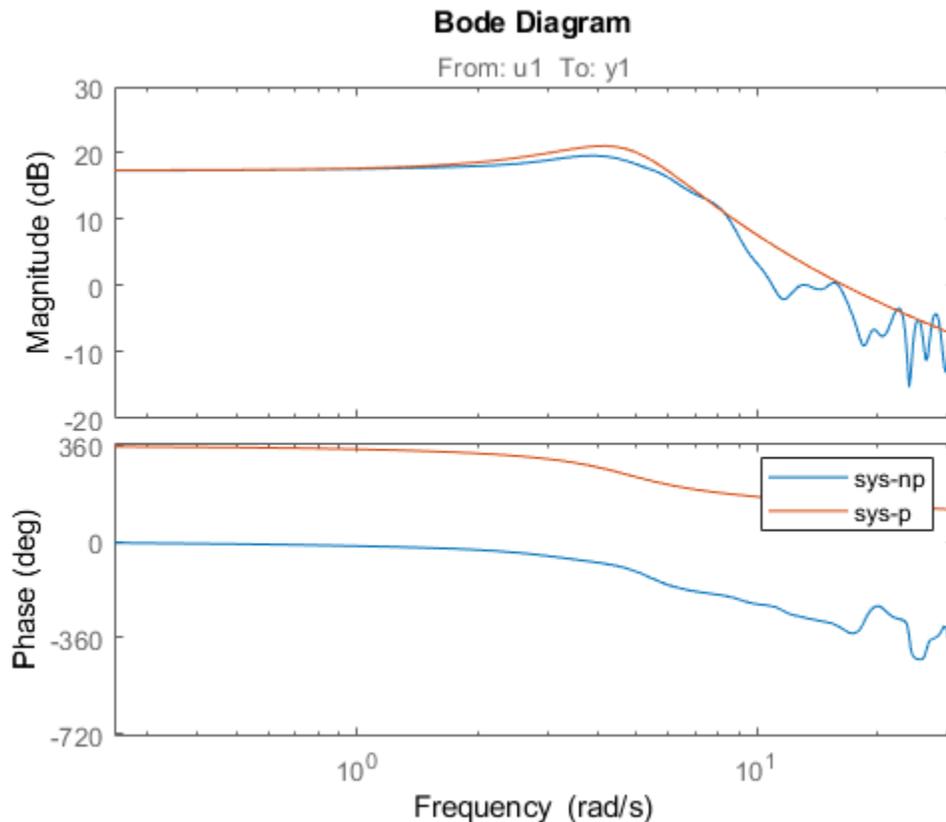
```
load iddata2 z2;  
w = linspace(0,10*pi,128);  
sys_np = spa(z2,[],w);  
sys_p = tfest(z2,2);
```

Using the `spa` and `tfest` commands requires System Identification Toolbox™ software.

`sys_np` is a nonparametric identified model. `sys_p` is a parametric identified model.

Create a Bode plot that includes both systems.

```
bode(sys_np,sys_p,w);  
legend('sys-np','sys-p')
```



You can display the confidence region on the Bode plot by right-clicking the plot and selecting **Characteristics > Confidence Region**.

### Obtain Magnitude and Phase Standard Deviation Data of Identified Model

Compute the standard deviation of the magnitude and phase of an identified model. Use this data to create a  $3\sigma$  plot of the response uncertainty.

Identify a transfer function model based on data. Obtain the standard deviation data for the magnitude and phase of the frequency response.

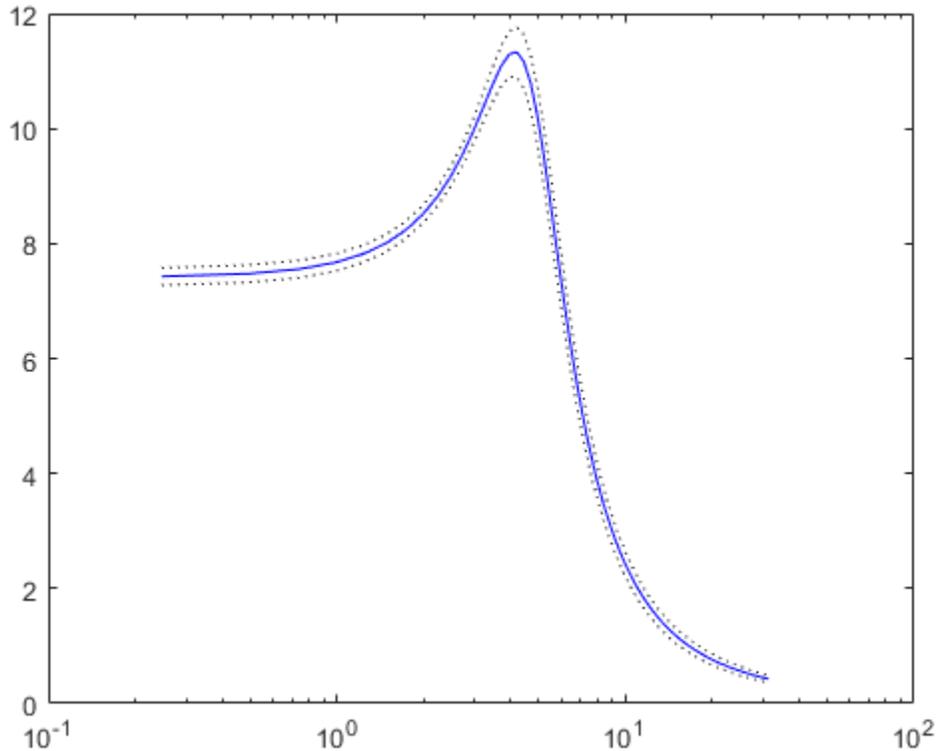
```
load iddata2 z2;  
sys_p = tfest(z2,2);  
w = linspace(0,10*pi,128);  
[mag,ph,w,sdmag,sdphase] = bode(sys_p,w);
```

Using the `tfest` command requires System Identification Toolbox™ software.

`sys_p` is an identified transfer function model. `sdmag` and `sdphase` contain the standard deviation data for the magnitude and phase of the frequency response, respectively.

Use the standard deviation data to create a  $3\sigma$  plot corresponding to the confidence region.

```
mag = squeeze(mag);  
sdmag = squeeze(sdmag);  
semilogx(w,mag,'b',w,mag+3*sdmag,'k:',w,mag-3*sdmag,'k:');
```



## Input Arguments

### **sys** — Dynamic system

dynamic system model | model array

Dynamic system, specified as a SISO or MIMO dynamic system model or array of dynamic system models. Dynamic systems that you can use include:

- Continuous-time or discrete-time numeric LTI models, such as `tf`, `zpk`, or `ss` models.
- Generalized or uncertain LTI models such as `genss` or `uss` models. (Using uncertain models requires Robust Control Toolbox™ software.)

- For tunable control design blocks, the function evaluates the model at its current value for both plotting and returning frequency response data.
- For uncertain control design blocks, the function plots the nominal value and random samples of the model. When you use output arguments, the function returns frequency response data for the nominal model only.
- Frequency-response data models such as `frd` models. For such models, the function plots the response at frequencies defined in the model.
- Identified LTI models, such as `idtf`, `idss`, or `idproc` models. For such models, the function can also plot confidence intervals and return standard deviations of the frequency response. See “Bode Plot of Identified Model” on page 1-142.

If `sys` is an array of models, the function plots the frequency responses of all models in the array on the same axes.

### **LineStyle — Line style, marker, and color**

character vector | string

Line style, marker, and color, specified as a string or vector of one, two, or three characters. The characters can appear in any order. You do not need to specify all three characteristics (line style, marker, and color). For example, if you omit the line style and specify the marker, then the plot shows only the marker and no line. For more information about configuring this argument, see the `LineStyle` input argument of the `plot` function.

Example: `'r--'` specifies a red dashed line

Example: `'*b'` specifies blue asterisk markers

Example: `'y'` specifies a yellow line

### **w — Frequencies**

{`wmin`, `wmax`} | vector

Frequencies at which to compute and plot frequency response, specified as the cell array `{wmin, wmax}` or as a vector of frequency values.

- If `w` is a cell array of the form `{wmin, wmax}`, then the function computes the index at frequencies ranging between `wmin` and `wmax`.
- If `w` is a vector of frequencies, then the function computes the index at each specified frequency. For example, use `logspace` to generate a row vector with logarithmically spaced frequency values.

Specify frequencies in units of rad/TimeUnit, where TimeUnit is the TimeUnit property of the model.

## Output Arguments

### **mag** — Magnitude of system response

3-D array

Magnitude of the system response in absolute units, returned as a 3-D array. The dimensions of this array are (number of system outputs) × (number of system inputs) × (number of frequency points).

- For SISO systems, `mag(1,1,k)` gives the magnitude of the response at the *k*th frequency in `w` or `wout`. For an example, see “Obtain Magnitude and Phase Data” on page 1-139.
- For MIMO systems, `mag(i,j,k)` gives the magnitude of the response at the *k*th frequency from the *j*th input to the *i*th output. For an example, see “Magnitude and Phase of MIMO System” on page 1-140.

To convert the magnitude from absolute units to decibels, use:

$$\text{magdb} = 20 * \log_{10}(\text{mag})$$

### **phase** — Phase of system response

3-D array

Phase of the system response in degrees, returned as a 3-D array. The dimensions of this array are (number of outputs) × (number of inputs) × (number of frequency points).

- For SISO systems, `mag(1,1,k)` gives the phase of the response at the *k*th frequency in `w` or `wout`. For an example, see “Obtain Magnitude and Phase Data” on page 1-139.
- For MIMO systems, `mag(i,j,k)` gives the phase of the response at the *k*th frequency from the *j*th input to the *i*th output. For an example, see “Magnitude and Phase of MIMO System” on page 1-140.

### **wout** — Frequencies

vector

Frequencies at which the function returns the system response, returned as a column vector. The function chooses the frequency values based on the model dynamics, unless you specify frequencies using the input argument `w`.

Frequency values are in radians per `TimeUnit`, where `TimeUnit` is the value of the `TimeUnit` property of `sys`.

### **sdmag — Standard deviation of magnitude**

3-D array | []

Estimated standard deviation of the magnitude of the response at each frequency point, returned as a 3-D array. `sdmag` has the same dimensions as `mag`.

If `sys` is not an identified LTI model, `sdmag` is [].

### **sdphase — Standard deviation of phase**

3-D array | []

Estimated standard deviation of the phase of the response at each frequency point, returned as a 3-D array. `sdphase` has the same dimensions as `phase`.

If `sys` is not an identified LTI model, `sdphase` is [].

## **Tips**

- When you need additional plot customization options, use `bodeplot` instead.

## **Algorithms**

`bode` computes the frequency response as follows:

- 1 Compute the zero-pole-gain (`zpk`) representation of the dynamic system.
- 2 Evaluate the gain and phase of the frequency response based on the zero, pole, and gain data for each input/output channel of the system.
  - For continuous-time systems, `bode` evaluates the frequency response on the imaginary axis  $s = j\omega$  and considers only positive frequencies.
  - For discrete-time systems, `bode` evaluates the frequency response on the unit circle. To facilitate interpretation, the command parameterizes the upper half of the unit circle as:

$$z = e^{j\omega T_s}, \quad 0 \leq \omega \leq \omega_N = \frac{\pi}{T_s},$$

where  $T_s$  is the sample time and  $\omega_N$  is the Nyquist frequency. The equivalent continuous-time frequency  $\omega$  is then used as the x-axis variable. Because

$H(e^{j\omega T_s})$  is periodic with period  $2\omega_N$ , **bode** plots the response only up to the Nyquist frequency  $\omega_N$ . If **sys** is a discrete-time model with unspecified sample time, **bode** uses  $T_s = 1$ .

## See Also

[bodeplot](#) | [freqresp](#) | [nyquist](#) | [spectrum](#) | [step](#)

## Topics

“Plot Bode and Nyquist Plots at the Command Line”

“Dynamic System Models”

**Introduced before R2006a**

## bodemag

Bode magnitude response of LTI models

### Syntax

```
bodemag(sys)
bodemag(sys,{wmin,wmax})
bodemag(sys,w)
bodemag(sys1,sys2,...,sysN,w)
```

### Description

`bodemag(sys)` plots the magnitude of the frequency response of the dynamic system model `sys` (Bode plot without the phase diagram). The frequency range and number of points are chosen automatically.

`bodemag(sys,{wmin,wmax})` draws the magnitude plot for frequencies between `wmin` and `wmax` (in `rad/TimeUnit`, where `TimeUnit` is the time units of the input dynamic system, specified in the `TimeUnit` property of `sys`).

`bodemag(sys,w)` uses the user-supplied vector `W` of frequencies, in `rad/TimeUnit`, at which the frequency response is to be evaluated.

`bodemag(sys1,sys2,...,sysN,w)` shows the frequency response magnitude of several models `sys1,sys2,...,sysN` on a single plot. The frequency vector `w` is optional. You can also specify a color, line style, and marker for each model. For example:

```
bodemag(sys1,'r',sys2,'y--',sys3,'gx')
```

### See Also

**Linear System Analyzer** | `bode`

**Introduced in R2012a**

# bodeoptions

Create list of Bode plot options

## Syntax

```
P = bodeoptions  
P = bodeoptions('cstprefs')
```

## Description

`P = bodeoptions` returns a default set of plot options for use with the `bodeplot`. You can use these options to customize the Bode plot appearance using the command line. This syntax is useful when you want to write a script to generate plots that look the same regardless of the preference settings of the MATLAB session in which you run the script.

`P = bodeoptions('cstprefs')` initializes the plot options with the options you selected in the Control System and System Identification Toolbox Preferences Editor. For more information about the editor, see “Toolbox Preferences Editor”. This syntax is useful when you want to change a few plot options but otherwise use your default preferences. A script that uses this syntax may generate results that look different when run in a session with different preferences.

The following table summarizes the Bode plot options.

Option	Description
Title, XLabel, YLabel	<p>Label text and style, specified as a structure with the following fields:</p> <ul style="list-style-type: none"> <li>• String — Label text, specified as a character vector, for example 'Amplitude'.</li> <li>• FontSize — <b>Default:</b> 8</li> <li>• FontWeight — <b>Default:</b> 'Normal'</li> <li>• Font Angle — <b>Default:</b> 'Normal'</li> <li>• Color — Vector of RGB values ranging from 0 to 1. <b>Default:</b> [0,0,0]</li> <li>• Interpreter — <b>Default:</b> 'tex'</li> </ul>
TickLabel	<p>Tick label style, specified as a structure with the following fields:</p> <ul style="list-style-type: none"> <li>• FontSize <b>Default:</b> 8</li> <li>• FontWeight — <b>Default:</b> 'Normal'</li> <li>• Font Angle — <b>Default:</b> 'Normal'</li> <li>• Color — Vector of RGB values ranging from 0 to 1. <b>Default:</b> [0,0,0]</li> </ul>
Grid	<p>Show or hide the grid            Specified as one of the following values: 'off'   'on'  <b>Default:</b> 'off'</p>
GridColor	<p>Color of the grid lines            Specified as one of the following: Vector of RGB values in the range [0,1]   character vector of color name   'none'. For example, for yellow color, specify as one of the following: [1 1 0], 'yellow', or 'y'.  <b>Default:</b> [0.15,0.15,0.15]</p>
XlimMode, YlimMode	<p>Axis limit modes. <b>Default:</b> 'auto'</p>
Xlim, Ylim	<p>Axes limits, specified as an array of the form [min,max]</p>
IOGrouping	<p>Grouping of input-output pairs            Specified as one of the following values: 'none'   'inputs'   'outputs'   'all'  <b>Default:</b> 'none'</p>

---

<b>Option</b>	<b>Description</b>
InputLabels, OutputLabels	Input and output label styles
InputVisible, OutputVisible	Visibility of input and output channels
ConfidenceRegionNum berSD	Number of standard deviations to use to plotting the response confidence region (identified models only).  <b>Default: 1.</b>

Option	Description
FreqUnits	<p>Frequency units, specified as one of the following values:</p> <ul style="list-style-type: none"> <li>• 'Hz'</li> <li>• 'rad/second'</li> <li>• 'rpm'</li> <li>• 'kHz'</li> <li>• 'MHz'</li> <li>• 'GHz'</li> <li>• 'rad/nanosecond'</li> <li>• 'rad/microsecond'</li> <li>• 'rad/millisecond'</li> <li>• 'rad/minute'</li> <li>• 'rad/hour'</li> <li>• 'rad/day'</li> <li>• 'rad/week'</li> <li>• 'rad/month'</li> <li>• 'rad/year'</li> <li>• 'cycles/nanosecond'</li> <li>• 'cycles/microsecond'</li> <li>• 'cycles/millisecond'</li> <li>• 'cycles/hour'</li> <li>• 'cycles/day'</li> <li>• 'cycles/week'</li> <li>• 'cycles/month'</li> <li>• 'cycles/year'</li> </ul>
FreqScale	<p>Frequency scale            Specified as one of the following values: 'linear'   'log'  <b>Default:</b> 'log'</p>

Option	Description
MagUnits	Magnitude units Specified as one of the following values: 'dB'   'abs' <b>Default:</b> 'dB'
MagScale	Magnitude scale Specified as one of the following values: 'linear'   'log' <b>Default:</b> 'linear'
MagVisible	Magnitude plot visibility Specified as one of the following values: 'on'   'off' <b>Default:</b> 'on'
MagLowerLimMode	Enables a lower magnitude limit Specified as one of the following values: 'auto'   'manual' <b>Default:</b> 'auto'
MagLowerLim	Specifies the lower magnitude limit
PhaseUnits	Phase units Specified as one of the following values: 'deg'   'rad' <b>Default:</b> 'deg'
PhaseVisible	Phase plot visibility Specified as one of the following values: 'on'   'off' <b>Default:</b> 'on'
PhaseWrapping	Enables phase wrapping Specified as one of the following values: 'on'   'off' When you set <code>PhaseWrapping</code> to 'on', the plot wraps accumulated phase at the value specified by the <code>PhaseWrappingBranch</code> property. <b>Default:</b> 'off'
PhaseWrappingBranch	Phase value at which the plot wraps accumulated phase when <code>PhaseWrapping</code> is set to 'on'. <b>Default:</b> -180 (phase wraps into the interval [-180°,180°))
PhaseMatching	Enables phase matching Specified as one of the following values: 'on'   'off' <b>Default:</b> 'off'
PhaseMatchingFreq	Frequency for matching phase
PhaseMatchingValue	The value to which phase responses are matched closely

## Examples

### Create Bode Plot with Custom Settings

Create a Bode plot that suppresses the phase plot and uses frequency units Hz instead of the default radians/second. Otherwise, the plot uses the settings that are saved in the toolbox preferences.

First, create an options set based on the toolbox preferences.

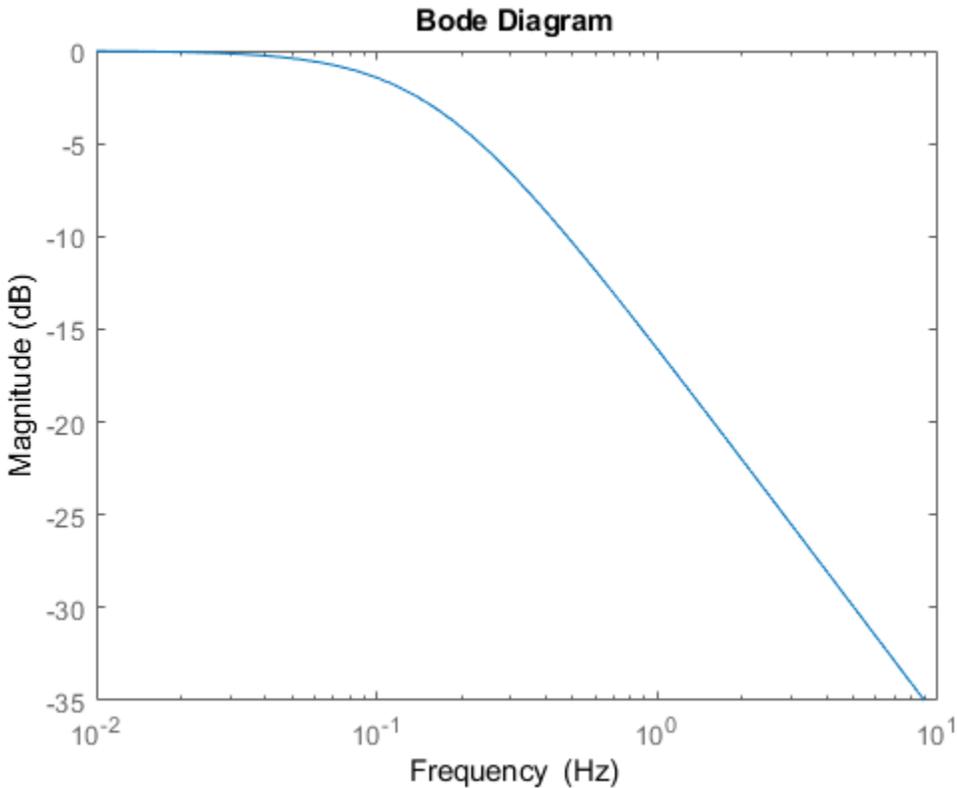
```
opts = bodeoptions('cstprefs');
```

Change properties of the options set.

```
opts.PhaseVisible = 'off';  
opts.FreqUnits = 'Hz';
```

Create a plot using the options.

```
h = bodeplot(tf(1,[1,1]),opts);
```



Depending on your own toolbox preferences, the plot you obtain might look different from this plot. Only the properties that you set explicitly, in this example `PhaseVisible` and `FreqUnits`, override the toolbox preferences.

### Custom Plot Settings Independent of Preferences

Create a Bode plot that uses 14-point red text for the title. This plot should look the same, regardless of the preferences of the MATLAB session in which it is generated.

First, create a default options set.

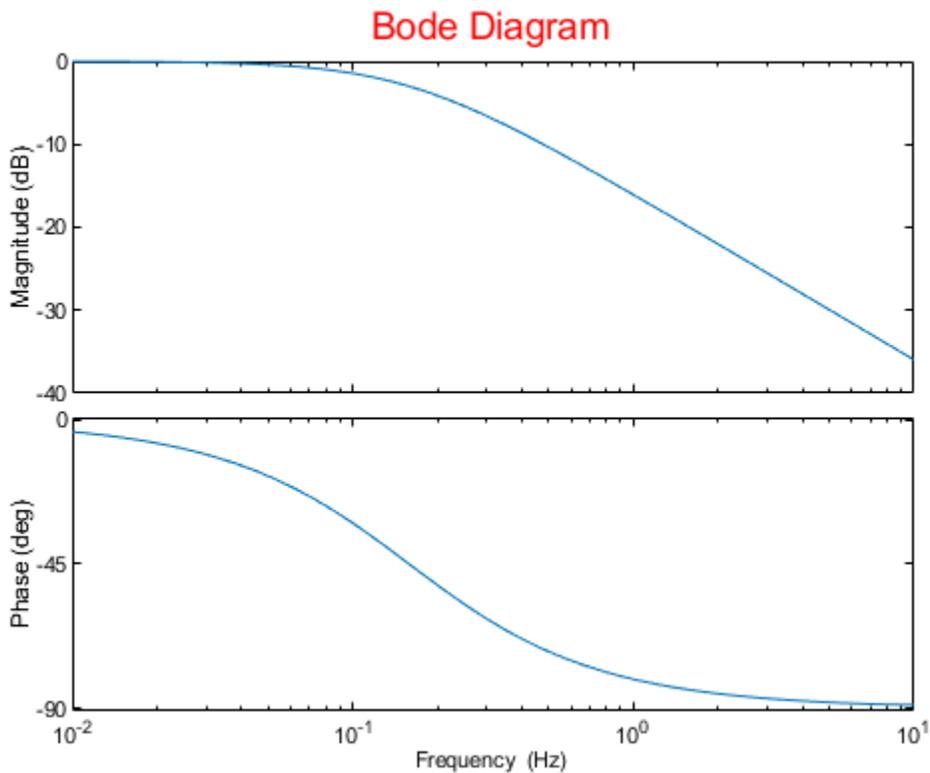
```
opts = bodeoptions;
```

Change properties of the options set.

```
opts.Title.FontSize = 14;  
opts.Title.Color = [1 0 0];  
opts.FreqUnits = 'Hz';
```

Create a plot using the options.

```
h = bodeplot(tf(1,[1,1]),opts);
```



Because `opts` begins with a fixed set of options, the plot result is independent of the toolbox preferences of the MATLAB session.

## **See Also**

bode | bodeplot | getoptions | setoptions | showConfidence

**Introduced in R2012a**

## **bodeplot**

Plot Bode frequency response with additional plot customization options

### **Syntax**

```
h = bodeplot(sys)
bodeplot(sys)
bodeplot(sys1,sys2,...)
bodeplot(AX,...)
bodeplot(..., plotoptions)
bodeplot(sys,w)
```

### **Description**

`h = bodeplot(sys)` plot the Bode magnitude and phase of the dynamic system model `sys` and returns the plot handle `h` to the plot. You can use this handle to customize the plot with the `getoptions` and `setoptions` commands.

`bodeplot(sys)` draws the Bode plot of the model `sys`. The frequency range and number of points are chosen automatically.

`bodeplot(sys1,sys2,...)` graphs the Bode response of multiple models `sys1,sys2,...` on a single plot. You can specify a color, line style, and marker for each model, as in

```
bodeplot(sys1,'r',sys2,'y--',sys3,'gx')
```

`bodeplot(AX,...)` plots into the axes with handle `AX`.

`bodeplot(..., plotoptions)` plots the Bode response with the options specified in `plotoptions`. Type

```
help bodeoptions
```

for a list of available plot options. See “Match Phase at Specified Frequency” on page 1-163 for an example of phase matching using the `PhaseMatchingFreq` and `PhaseMatchingValue` options.

`bodeplot(sys,w)` draws the Bode plot for frequencies specified by `w`. When `w = {wmin,wmax}`, the Bode plot is drawn for frequencies between `wmin` and `wmax` (in `rad/TimeUnit`, where `TimeUnit` is the time units of the input dynamic system, specified in the `TimeUnit` property of `sys`). When `w` is a user-supplied vector `w` of frequencies, in `rad/TimeUnit`, the Bode response is drawn for the specified frequencies.

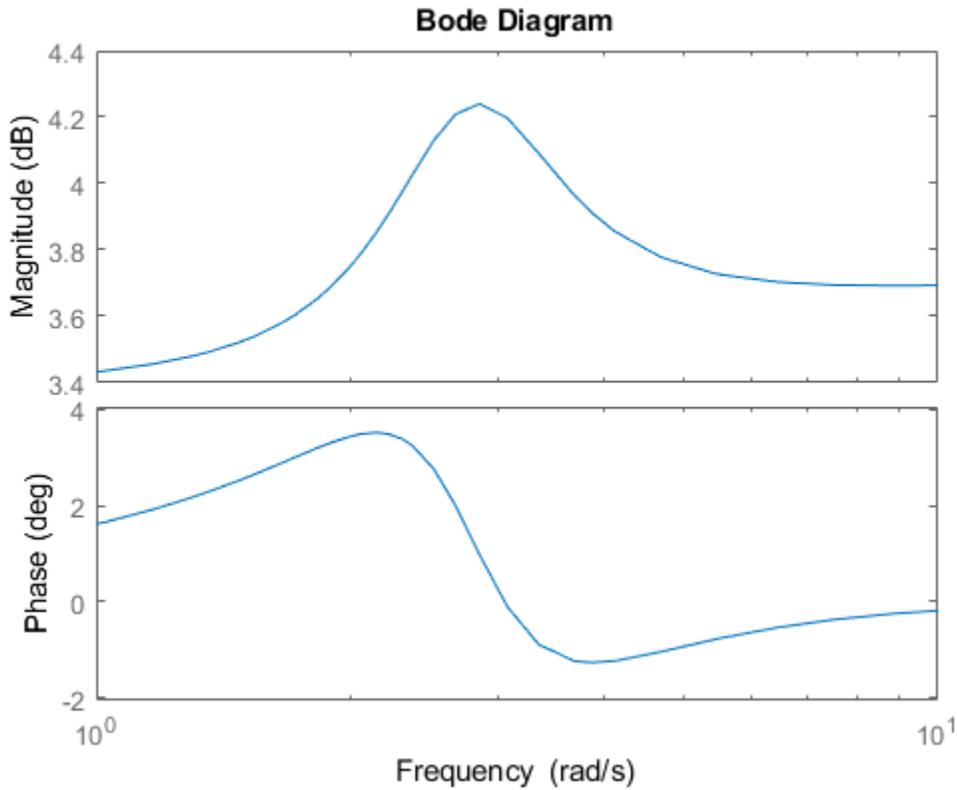
See `logspace` to generate logarithmically spaced frequency vectors.

## Examples

### Change Bode Plot Options with Plot Handle

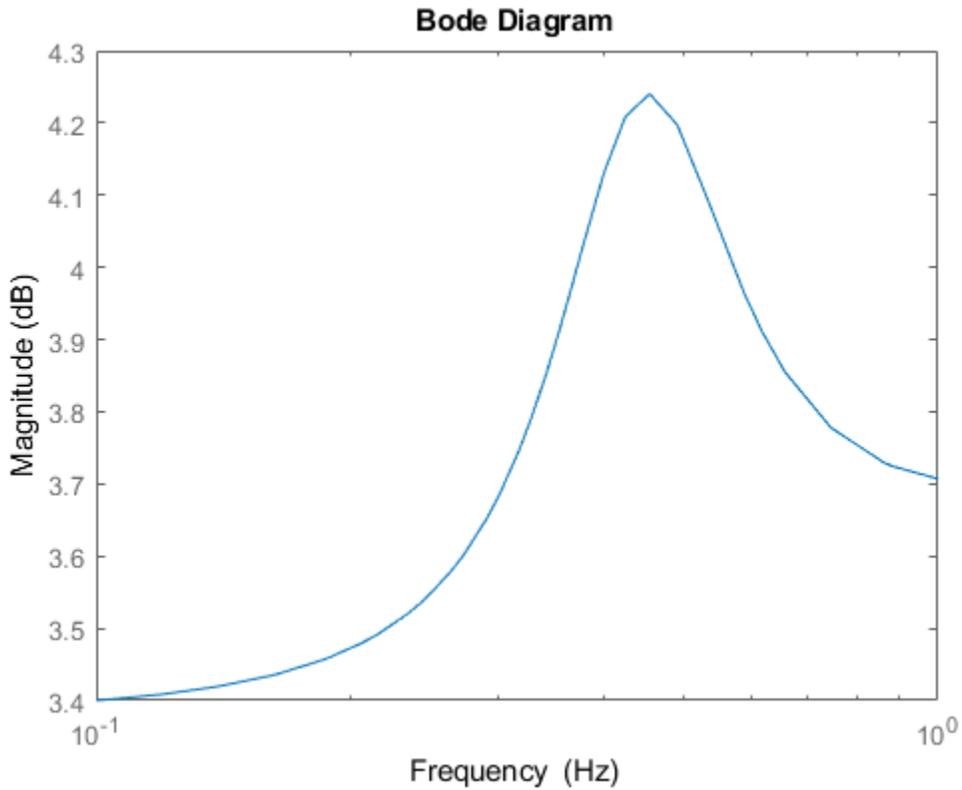
Generate a Bode plot.

```
sys = rss(5);  
h = bodeplot(sys);
```



Change the units to Hz and suppress the phase plot. To do so, edit properties of the plot handle, h.

```
setoptions(h, 'FreqUnits', 'Hz', 'PhaseVisible', 'off');
```

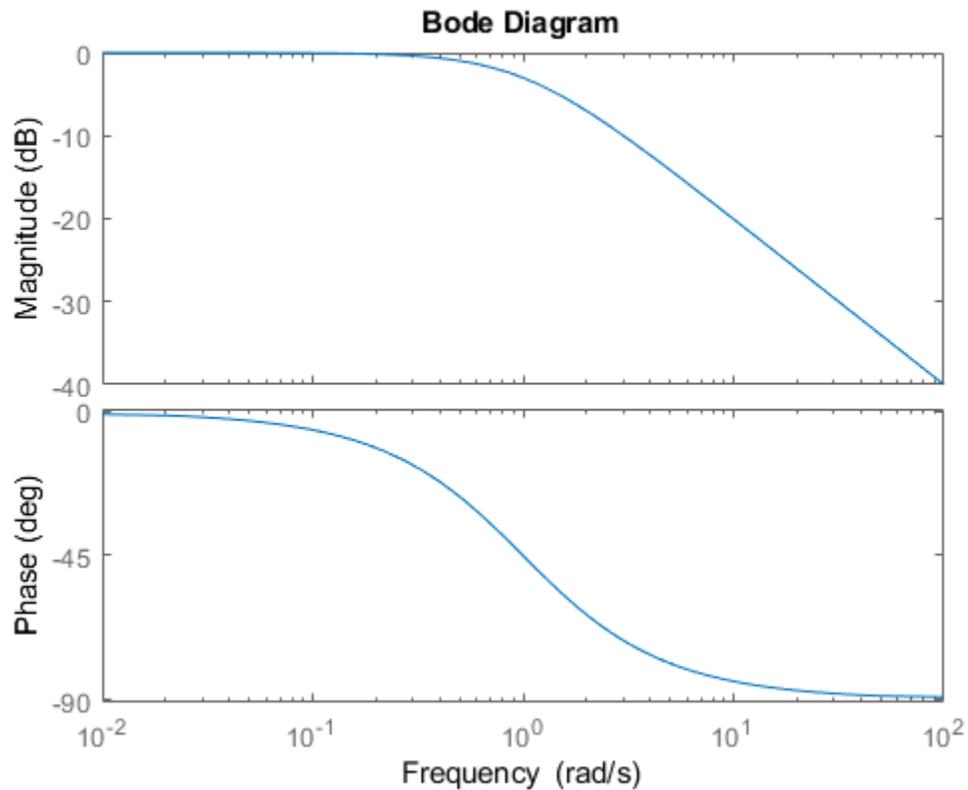


The plot automatically updates when you call `setoptions`.

### Match Phase at Specified Frequency

Create a Bode plot of a dynamic system.

```
sys = tf(1,[1 1]);  
h = bodeplot(sys);
```

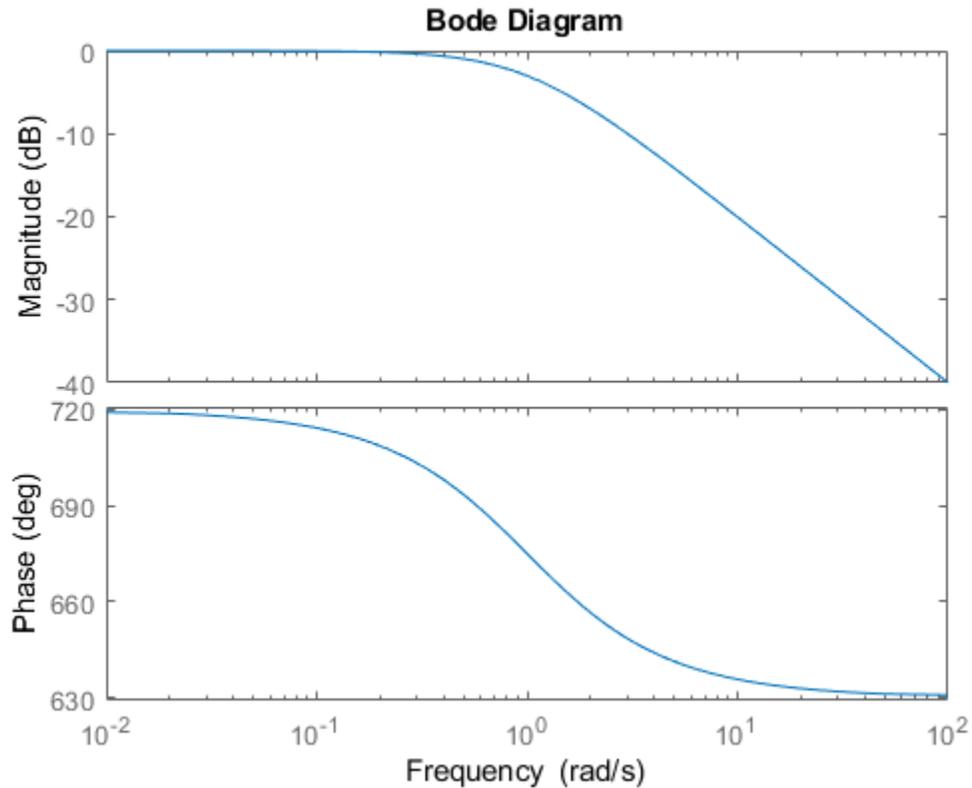


Fix the phase at 1 rad/s to 750 degrees. To do so, get the plot properties. Then alter the properties `PhaseMatchingFreq` and `PhaseMatchingValue` to match a phase to a specified frequency.

```
p = getoptions(h);  
p.PhaseMatching = 'on';  
p.PhaseMatchingFreq = 1;  
p.PhaseMatchingValue = 750;
```

Update the plot.

```
setoptions(h,p);
```



The first bode plot has a phase of -45 degrees at a frequency of 1 rad/s. Setting the phase matching options so that at 1 rad/s the phase is near 750 degrees yields the second Bode plot. Note that, however, the phase can only be  $-45 + N \cdot 360$ , where  $N$  is an integer, and so the plot is set to the nearest allowable phase, namely 675 degrees (or  $2 \cdot 360 - 45 = 675$ ).

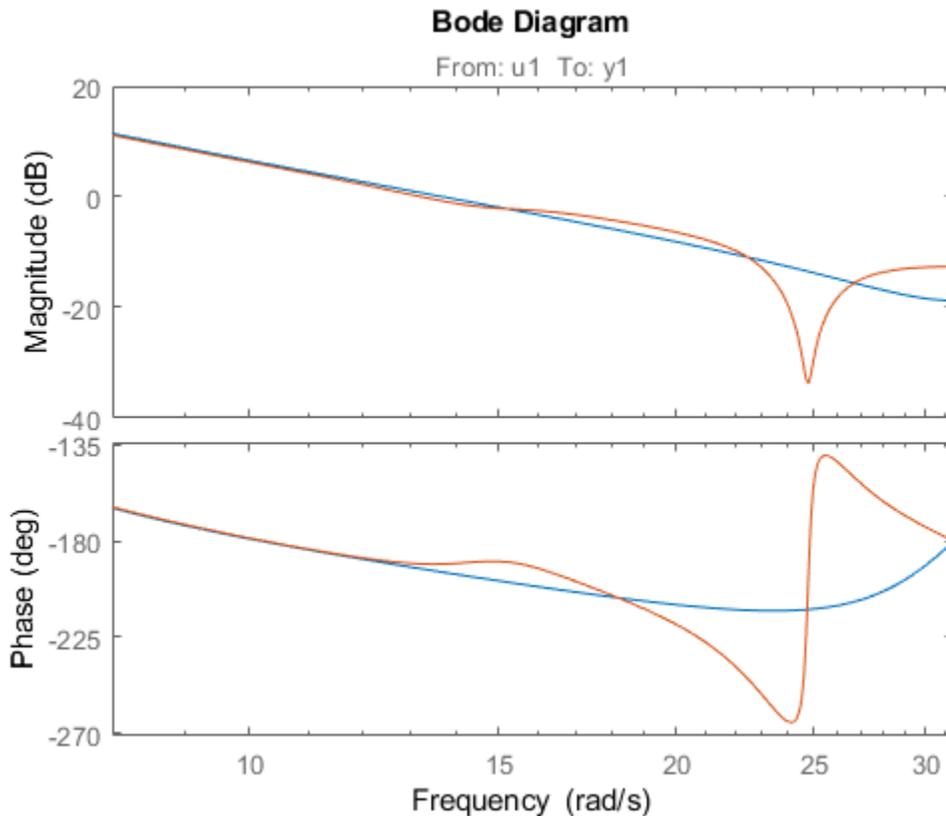
### Display Confidence Regions of Identified Models

Compare the frequency responses of identified state-space models of order 2 and 6 along with their  $2\sigma$  confidence regions.

```
load iddata1
sys1 = n4sid(z1, 2);
sys2 = n4sid(z1, 6);
```

Both models produce about 70% fit to data. However, `sys2` shows higher uncertainty in its frequency response, especially close to Nyquist frequency as shown by the plot:

```
w = linspace(8,10*pi,256);
h = bodeplot(sys1,sys2,w);
setoptions(h, 'PhaseMatching', 'on', 'ConfidenceRegionNumberSD', 2);
```



Right-click the plot and select **Characteristics > Confidence Region** to turn on the confidence region characteristic. Alternatively, type `showConfidence(h)` to plot the confidence region.

## Frequency Response of Identified Parametric and Nonparametric Models

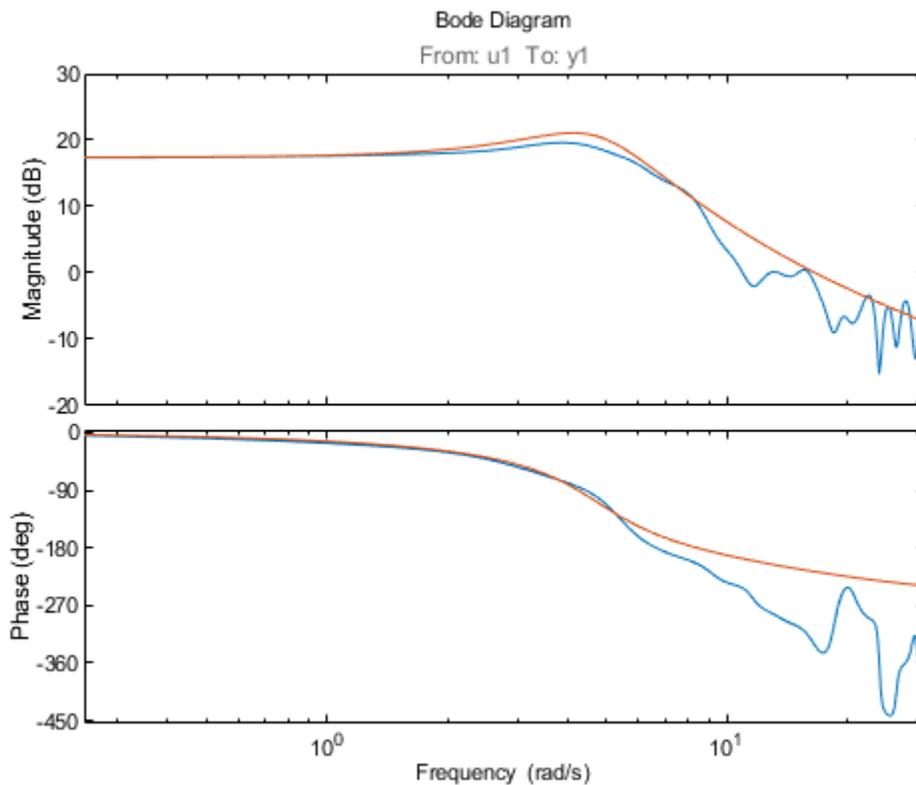
Compare the frequency response of a parametric model, identified from input/output data, to a nonparametric model identified using the same data. Identify parametric and non-parametric models based on data.

```
load iddata2 z2;  
w = linspace(0,10*pi,128);  
sys_np = spa(z2,[],w);  
sys_p = tfest(z2,2);
```

`spa` and `tfest` require System Identification Toolbox™ software. `sys_np` is a nonparametric identified model. `sys_p` is a parametric identified model.

Create a Bode plot that includes both systems.

```
opt = bodeoptions;  
opt.PhaseMatching = 'on';  
bodeplot(sys_np,sys_p,w,opt);
```



## Tips

You can change the properties of your plot, for example the units. For information on the ways to change properties of your plots, see “Ways to Customize Plots” (Control System Toolbox).

## See Also

`bode` | `bodeoptions` | `getoptions` | `setoptions` | `showConfidence`

**Introduced before R2006a**

## c2d

Convert model from continuous to discrete time

### Syntax

```
sysd = c2d(sys,Ts)
sysd = c2d(sys,Ts,method)
sysd = c2d(sys,Ts,opts)
[sysd,G] = c2d(sys,Ts,method)
[sysd,G] = c2d(sys,Ts,opts)
```

### Description

`sysd = c2d(sys,Ts)` discretizes the continuous-time dynamic system model `sys` using zero-order hold on the inputs and a sample time of `Ts` seconds.

`sysd = c2d(sys,Ts,method)` discretizes `sys` using the specified discretization method `method`.

`sysd = c2d(sys,Ts,opts)` discretizes `sys` using the option set `opts`, specified using the `c2dOptions` command.

`[sysd,G] = c2d(sys,Ts,method)` returns a matrix, `G` that maps the continuous initial conditions  $x_0$  and  $u_0$  of the state-space model `sys` to the discrete-time initial state vector `x[0]`. `method` is optional. To specify additional discretization options, use `[sysd,G] = c2d(sys,Ts,opts)`.

### Input Arguments

#### **sys**

Continuous-time dynamic system model (except frequency response data models). `sys` can represent a SISO or MIMO system, except that the 'matched' discretization method supports SISO systems only.

sys can have input/output or internal time delays; however, the 'matched', 'impulse', and least-squares' methods do not support state-space models with internal time delays.

The following identified linear systems cannot be discretized directly:

- idgrey models whose FunctionType is 'c'. Convert to idss model first.
- idproc models. Convert to idtf or idpoly model first.

For the syntax `[sysd,G] = c2d(sys,Ts,opts)`, sys must be a state-space model.

### **Ts**

Sample time.

### **method**

Discretization method, specified as one of the following values:

- 'zoh' — Zero-order hold (default). Assumes the control inputs are piecewise constant over the sample time Ts.
- 'foh' — Triangle approximation (modified first-order hold). Assumes the control inputs are piecewise linear over the sample time Ts.
- 'impulse' — Impulse invariant discretization
- 'tustin' — Bilinear (Tustin) method
- 'matched' — Zero-pole matching method
- 'least-squares' — Least-squares method

For information about the algorithms for each conversion method, see “Continuous-Discrete Conversion Methods”.

### **opts**

Discretization options. Create opts using `c2doptions`.

## **Output Arguments**

### **sysd**

Discrete-time model of the same type as the input system sys.

When `sys` is an identified (IDLTI) model, `sysd`:

- Includes both measured and noise components of `sys`. The innovations variance  $\lambda$  of the continuous-time identified model `sys`, stored in its `NoiseVariance` property, is interpreted as the intensity of the spectral density of the noise spectrum. The noise variance in `sysd` is thus  $\lambda/T_s$ .
- Does not include the estimated parameter covariance of `sys`. If you want to translate the covariance while discretizing the model, use `translatecov`.

## G

Matrix relating continuous-time initial conditions  $x_0$  and  $u_0$  of the state-space model `sys` to the discrete-time initial state vector `x[0]`, as follows:

$$x[0] = G \cdot \begin{bmatrix} x_0 \\ u_0 \end{bmatrix}$$

For state-space models with time delays, `c2d` pads the matrix `G` with zeroes to account for additional states introduced by discretizing those delays. See “Continuous-Discrete Conversion Methods” for a discussion of modeling time delays in discretized systems.

## Examples

### Discretize a Transfer Function

Discretize the following continuous-time transfer function:

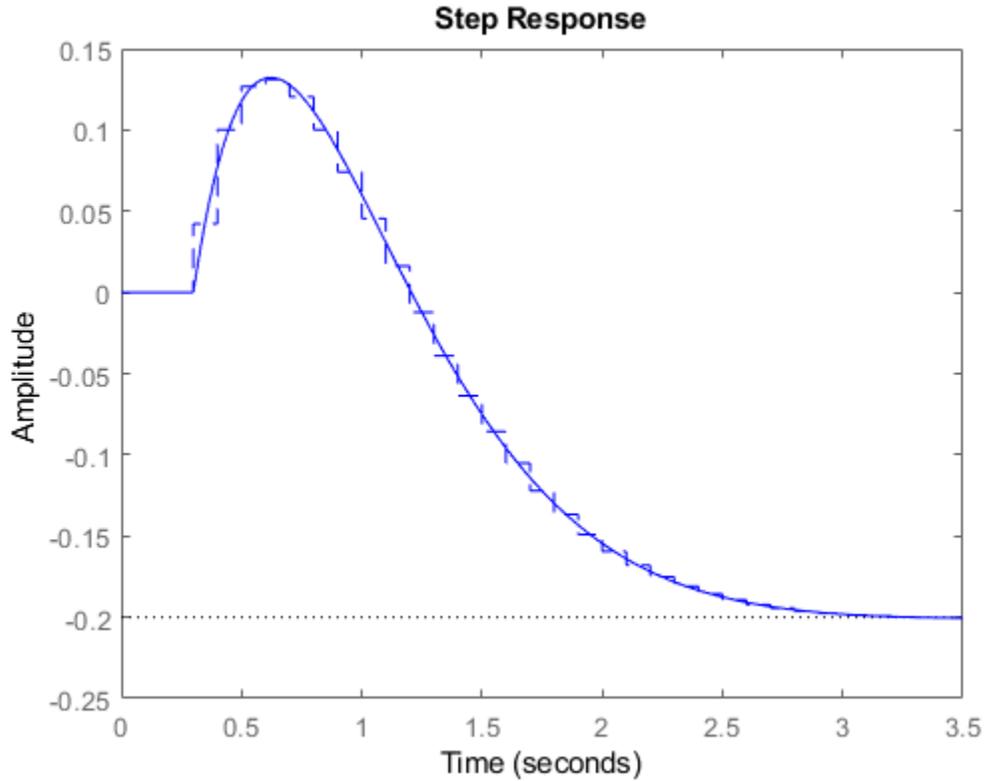
$$H(s) = e^{-0.3s} \frac{s - 1}{s^2 + 4s + 5}$$

This system has an input delay of 0.3 s. Discretize the system using the triangle (first-order-hold) approximation with sample time  $T_s = 0.1$  s.

```
H = tf([1 -1],[1 4 5],'InputDelay', 0.3);  
Hd = c2d(H,0.1,'foh');
```

Compare the step responses of the continuous-time and discretized systems.

```
step(H, '-', Hd, '---')
```



### Discretize Model with Fractional Delay Absorbed into Coefficients

Discretize the following delayed transfer function using zero-order hold on the input, and a 10-Hz sampling rate.

$$H(s) = e^{-0.25s} \frac{10}{s^2 + 3s + 10}$$

```
h = tf(10,[1 3 10], 'IODelay',0.25);
hd = c2d(h,0.1)
```

hd =

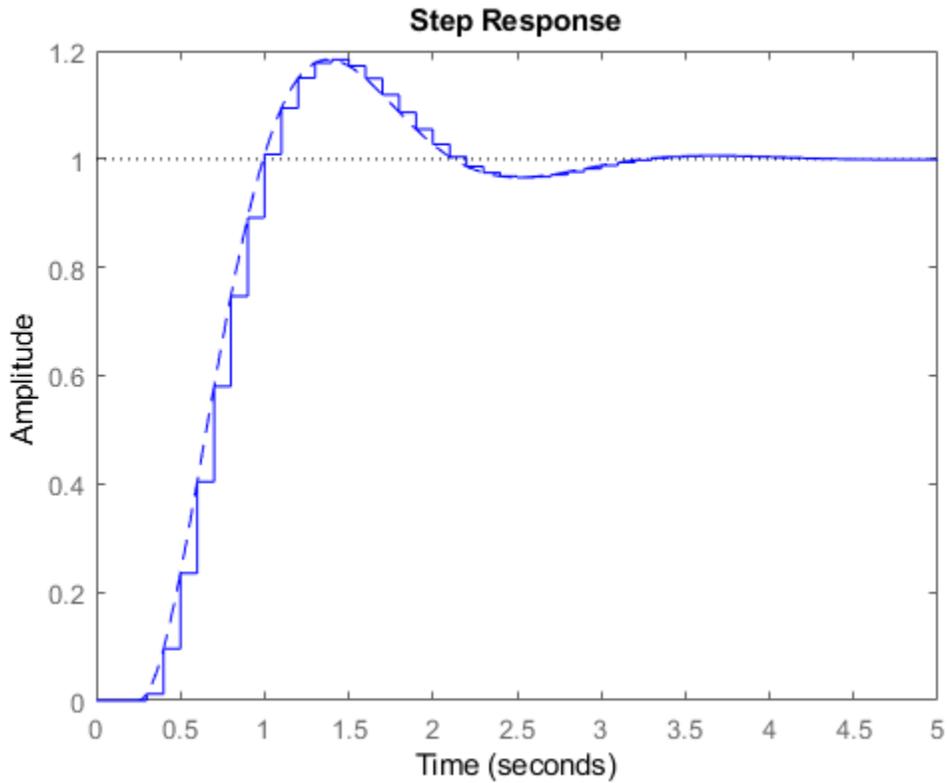
$$z^{-3} * \frac{0.01187 z^2 + 0.06408 z + 0.009721}{z^2 - 1.655 z + 0.7408}$$

Sample time: 0.1 seconds  
Discrete-time transfer function.

In this example, the discretized model `hd` has a delay of three sampling periods. The discretization algorithm absorbs the residual half-period delay into the coefficients of `hd`.

Compare the step responses of the continuous-time and discretized models.

```
step(h, '-',hd, '-')
```



### Discretize Model With Approximated Fractional Delay

Create a continuous-time state-space model with two states and an input delay.

```
sys = ss(tf([1,2],[1,4,2]));  
sys.InputDelay = 2.7
```

```
sys =
```

```
A =  
      x1  x2  
x1  -4  -2
```

```
x2  1  0
B =
      u1
x1  2
x2  0
C =
      x1  x2
y1  0.5  1
D =
      u1
y1  0
```

Input delays (seconds): 2.7

Continuous-time state-space model.

Discretize the model using the Tustin discretization method and a Thiran filter to model fractional delays. The sample time  $T_s = 1$  second.

```
opt = c2dOptions('Method','tustin','FractDelayApproxOrder',3);
sysd1 = c2d(sys,1,opt)
```

```
sysd1 =
```

```
A =
      x1      x2      x3      x4      x5
x1  -0.4286  -0.5714  -0.00265  0.06954  2.286
x2   0.2857   0.7143  -0.001325  0.03477  1.143
x3   0         0      -0.2432   0.1449  -0.1153
x4   0         0         0.25     0         0
x5   0         0         0         0.125    0
```

```
B =
      u1
x1  0.002058
x2  0.001029
x3   8
x4   0
x5   0
```

```
C =
      x1      x2      x3      x4      x5
```

```

y1      0.2857      0.7143  -0.001325      0.03477      1.143
D =
          u1
y1  0.001029

```

Sample time: 1 seconds  
Discrete-time state-space model.

The discretized model now contains three additional states  $x_3$ ,  $x_4$ , and  $x_5$  corresponding to a third-order Thiran filter. Since the time delay divided by the sample time is 2.7, the third-order Thiran filter ('FractDelayApproxOrder' = 3) can approximate the entire time delay.

### Discretize Identified Model

Estimate a continuous-time transfer function, and discretize it.

```

load iddata1
sys1c = tfest(z1,2);
sys1d = c2d(sys1c,0.1,'zoh');

```

Estimate a second order discrete-time transfer function.

```

sys2d = tfest(z1,2,'Ts',0.1);

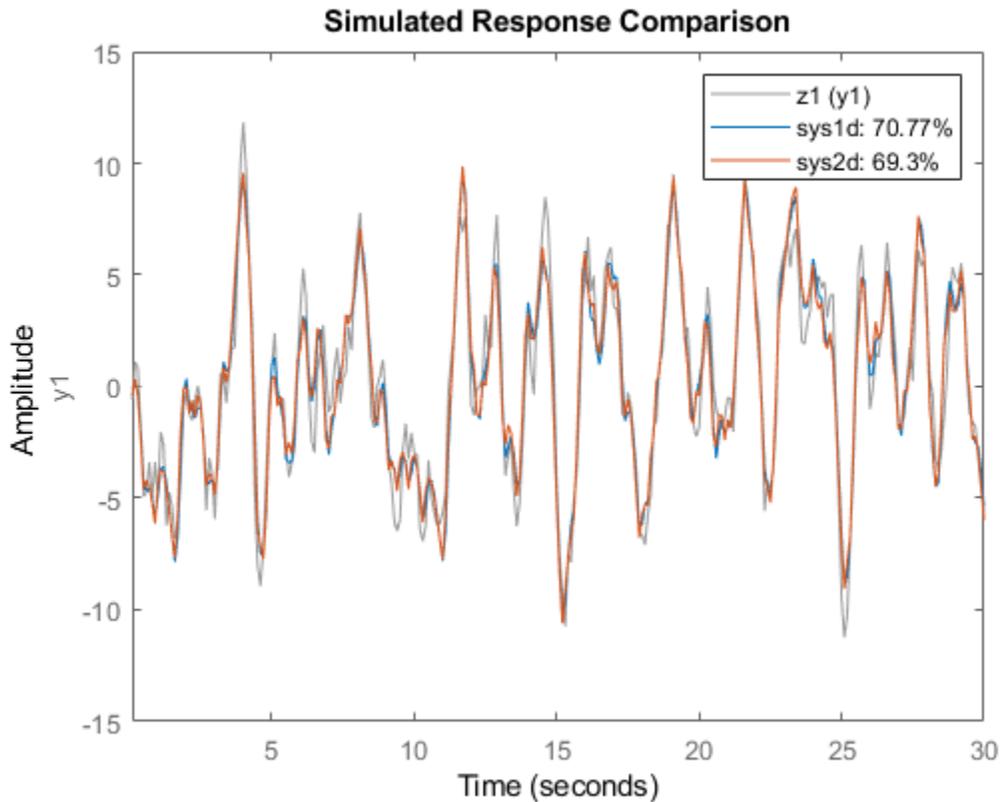
```

Compare the response of the discretized continuous-time transfer function model, `sys1d`, and the directly estimated discrete-time model, `sys2d`.

```

compare(z1,sys1d,sys2d)

```



The two systems are almost identical.

### Build Predictor Model

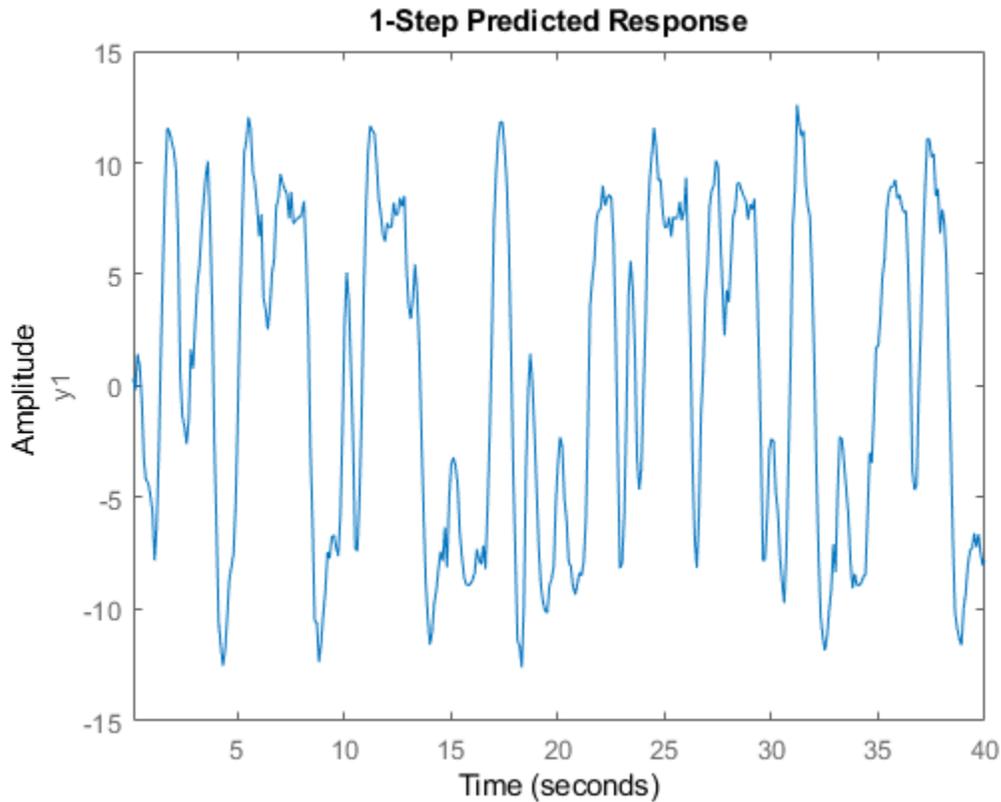
Discretize an identified state-space model to build a one-step ahead predictor of its response.

Create a continuous-time identified state-space model using estimation data.

```
load iddata2  
sysc = sstest(z2,4);
```

Predict the 1-step ahead predicted response of sysc.

```
predict(sysc,z2)
```



Discretize the model.

```
sysd = c2d(sysc,0.1,'zoh');
```

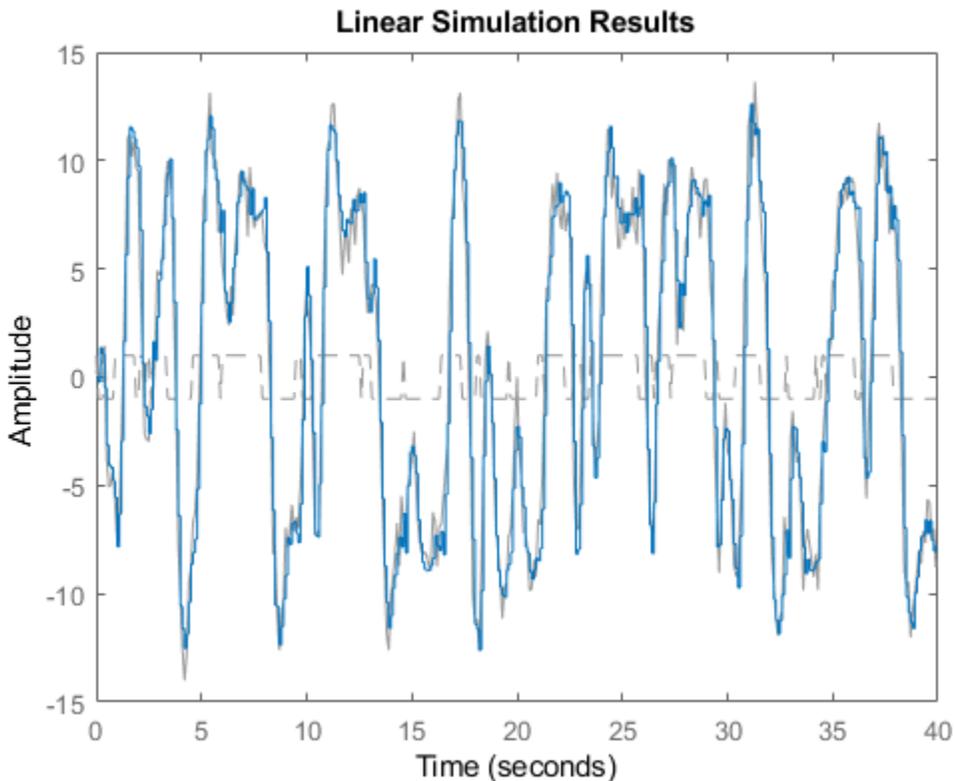
Build a predictor model from the discretized model, sysd.

```
[A,B,C,D,K] = idssdata(sysd);
Predictor = ss(A-K*C,[K B-K*D],C,[0 D],0.1);
```

Predictor is a two-input model which uses the measured output and input signals ( $[z1.y \ z1.u]$ ) to compute the 1-step predicted response of sysc.

Simulate the predictor model to get the same response as the predict command.

```
lsim(Predictor,[z2.y,z2.u])
```



The simulation of the predictor model gives the same response as `predict(sys,z2)`.

## Tips

- Use the syntax `sysd = c2d(sys,Ts,method)` to discretize `sys` using the default options for `method`. To specify additional discretization options, use the syntax `sysd = c2d(sys,Ts,opts)`.

- To specify the `tustin` method with frequency prewarping (formerly known as the 'prewarp' method), use the `PrewarpFrequency` option of `c2dOptions`.

## Algorithms

For information about the algorithms for each c2d conversion method, see “Continuous-Discrete Conversion Methods”.

## See Also

`c2dOptions` | `d2c` | `d2d` | `thiran` | `translatecov`

## Topics

“Dynamic System Models”

“Discretize a Compensator” (Control System Toolbox)

“Continuous-Discrete Conversion Methods” (Control System Toolbox)

**Introduced before R2006a**

## c2dOptions

Create option set for continuous- to discrete-time conversions

### Syntax

```
opts = c2dOptions  
opts = c2dOptions('OptionName', OptionValue)
```

### Description

*opts* = c2dOptions returns the default options for c2d.

*opts* = c2dOptions('OptionName', OptionValue) accepts one or more comma-separated name/value pairs that specify options for the c2d command. Specify *OptionName* inside single quotes.

### Input Arguments

#### Name-Value Pair Arguments

##### Method

Discretization method, specified as one of the following values:

'zoh'	Zero-order hold, where c2d assumes the control inputs are piecewise constant over the sample time <i>Ts</i> .
'foh'	Triangle approximation (modified first-order hold), where c2d assumes the control inputs are piecewise linear over the sample time <i>Ts</i> . (See [1] on page 1-184, p. 228.)
'impulse'	Impulse-invariant discretization.

'tustin'	Bilinear (Tustin) approximation. By default, c2d discretizes with no prewarp and rounds any fractional time delays to the nearest multiple of the sample time. To include prewarp, use the PrewarpFrequency option. To approximate fractional time delays, use theFractDelayApproxOrder option.
'matched'	Zero-pole matching method. (See [1] on page 1-184, p. 224.) By default, c2d rounds any fractional time delays to the nearest multiple of the sample time. To approximate fractional time delays, use the FractDelayApproxOrder option.
'least-squares'	Least-squares method. Minimize the error between the frequency responses of the continuous-time and discrete-time systems up to the Nyquist frequency.

For information about the algorithms for each conversion method, see “Continuous-Discrete Conversion Methods”.

**Default:** 'zoh'

### PrewarpFrequency

Prewarp frequency for 'tustin' method, specified in rad/TimeUnit, where TimeUnit is the time units, specified in the TimeUnit property, of the discretized system. Takes positive scalar values. A value of 0 corresponds to the standard 'tustin' method without prewarp.

**Default:** 0

### FractDelayApproxOrder

Maximum order of the Thiran filter used to approximate fractional delays in the 'tustin' and 'matched' methods. Takes integer values. A value of 0 means that c2d rounds fractional delays to the nearest integer multiple of the sample time.

**Default:** 0

## Examples

### Discretize Two Models Using Tustin Discretization Method

Generate two random continuous-time state-space models.

```
sys1 = rss(3,2,2);  
sys2 = rss(4,4,1);
```

Create an option set for `c2d` to use the Tustin discretization method and 3.4 rad/s prewarp frequency.

```
opt = c2doptions('Method','tustin','PrewarpFrequency',3.4);
```

Discretize the models, `sys1` and `sys2`, using the same option set, but different sample times.

```
dsys1 = c2d(sys1,0.1,opt);  
dsys2 = c2d(sys2,0.2,opt);
```

## References

[1] Franklin, G.F., Powell, D.J., and Workman, M.L., *Digital Control of Dynamic Systems* (3rd Edition), Prentice Hall, 1997.

## See Also

`c2d`

**Introduced in R2012a**

## canon

State-space canonical realization

### Syntax

```
csys = canon(sys,type)
[csys,T]= canon(sys,type)
csys = canon(sys,'modal',condt)
```

### Description

`csys = canon(sys,type)` transforms the linear model `sys` into a canonical state-space model `csys`. The argument `type` specifies whether `csys` is in modal on page 1-189 or companion on page 1-190 form.

`[csys,T]= canon(sys,type)` also returns the state-coordinate transformation `T` that relates the states of the state-space model `sys` to the states of `csys`.

`csys = canon(sys,'modal',condt)` specifies an upper bound `condt` on the condition number of the block-diagonalizing transformation.

### Input Arguments

#### **sys**

Any linear dynamic system model, except for `frd` models.

#### **type**

Canonical form of `csys`, specified as one of the following values:

- 'modal' — convert `sys` to modal form on page 1-189.
- 'companion' — convert `sys` to companion form on page 1-190.

**condt**

Positive scalar value specifying an upper bound on the condition number of the block-diagonalizing transformation that converts `sys` to `csys`. This argument is available only when `type` is 'modal'.

Increase `condt` to reduce the size of the eigenvalue clusters in the  $A$  matrix of `csys`. Setting `condt = Inf` diagonalizes  $A$ .

**Default:** 1e8

## Output Arguments

**csys**

State-space (ss) model. `csys` is a state-space realization of `sys` in the canonical form specified by `type`.

**T**

Matrix specifying the transformation between the state vector  $x$  of the state-space model `sys` and the state vector  $x_c$  of `csys`:

$$x_c = Tx$$

This argument is available only when `sys` is state-space model.

## Examples

### Convert System To Modal Canonical Form

Consider a system with doubled poles and clusters of close poles:

$$G(s) = 100 \frac{(s - 1)(s + 1)}{s(s + 10)(s + 10.0001)(s - (1 + i))^2(s - (1 - i))^2}$$

Create a linear model of this system, and convert it to modal canonical form.

```
G = zpk([1 -1],[0 -10 -10.0001 1+1i 1-1i 1+1i 1-1i],100);
Gc = canon(G, 'modal');
```

The system, G, has a pair of nearby poles at  $s = -10$  and  $s = -10.0001$ . G also has two complex poles of multiplicity 2 at  $s = 1 + i$  and  $s = 1 - i$ . As a result, the modal form has a block of size 2 for the two poles near  $s = -10$ , and a block of size 4 for the complex eigenvalues.

Gc.A

ans = 7x7

0	0	0	0	0	0	0	0
0	1.0000	1.0000	0	0	0	0	0
0	-1.0000	1.0000	2.0548	0	0	0	0
0	0	0	1.0000	1.0000	0	0	0
0	0	0	-1.0000	1.0000	0	0	0
0	0	0	0	0	-10.0000	8.0573	0
0	0	0	0	0	0	-10.0001	0

Separate the two poles near  $s = -10$  by increasing the value of the condition number of the block-diagonalizing transformation. The default value of the condition number is  $1e8$ .

```
Gc2 = canon(G, 'modal', 1e10);
Gc2.A
```

ans = 7x7

0	0	0	0	0	0	0	0
0	1.0000	1.0000	0	0	0	0	0
0	-1.0000	1.0000	2.0548	0	0	0	0
0	0	0	1.0000	1.0000	0	0	0
0	0	0	-1.0000	1.0000	0	0	0
0	0	0	0	0	-10.0000	0	0
0	0	0	0	0	0	-10.0001	0

The A matrix of Gc2 includes separate diagonal elements for the poles near  $s = -10$ . The cost of increasing the condition number of A is that the B matrix includes some large values.

```
format shortE
Gc2.B
```

```
ans = 7×1

    3.2000e-01
   -6.5691e-03
    5.4046e-02
   -1.9502e-01
    1.0637e+00
    3.2533e+05
    3.2533e+05
```

## Convert System to Companion Canonical Form

Estimate a state-space model that is freely parameterized.

```
load icEngine.mat
z = iddata(y,u,0.04);
FreeModel = n4sid(z,4,'InputDelay',2);
```

Convert the estimated model to companion canonical form.

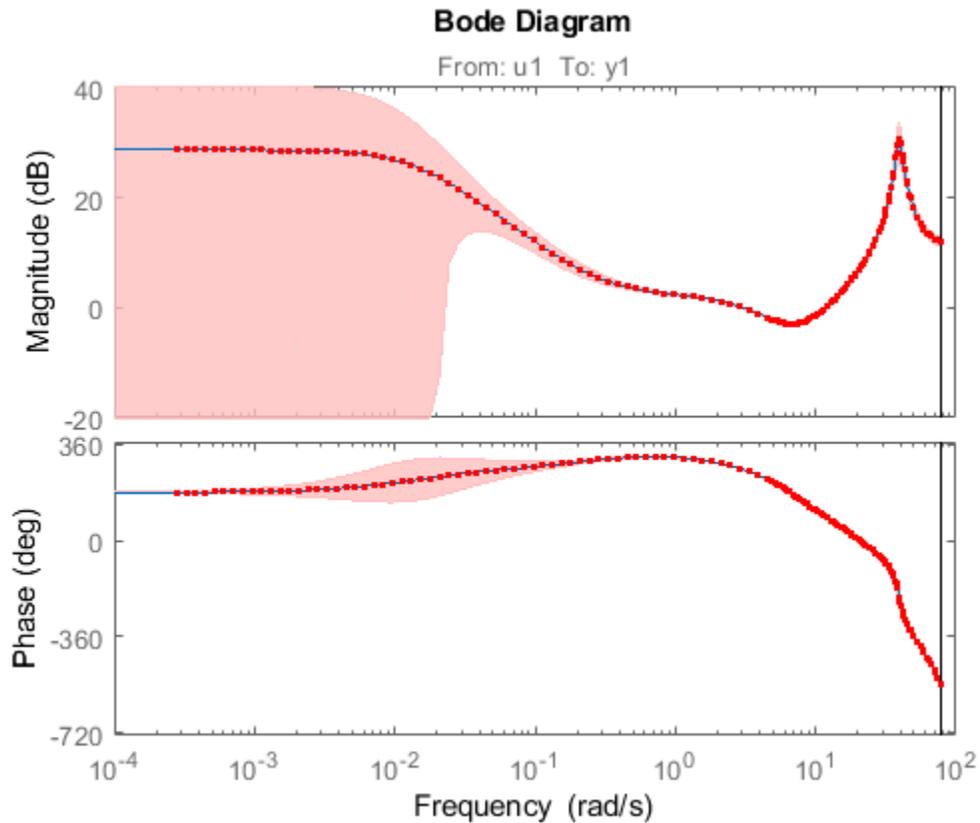
```
CanonicalModel = canon(FreeModel,'companion');
```

Obtain the covariance of the resulting form by running a zero-iteration update to model parameters.

```
opt = ssestOptions;
opt.SearchOptions.MaxIterations = 0;
CanonicalModel = ssest(z,CanonicalModel,opt);
```

Compare frequency response confidence bounds of FreeModel to CanonicalModel.

```
h = bodeplot(FreeModel,CanonicalModel,'r. ');
showConfidence(h)
```



The frequency response confidence bounds are identical.

## Definitions

### Modal Form

In modal form,  $A$  is a block-diagonal matrix. The block size is typically 1-by-1 for real eigenvalues and 2-by-2 for complex eigenvalues. However, if there are repeated eigenvalues or clusters of nearby eigenvalues, the block size can be larger.

For example, for a system with eigenvalues  $(\lambda_1, \sigma \pm j\omega, \lambda_2)$ , the modal  $A$  matrix is of the form

$$\begin{bmatrix} \lambda_1 & 0 & 0 & 0 \\ 0 & \sigma & \omega & 0 \\ 0 & -\omega & \sigma & 0 \\ 0 & 0 & 0 & \lambda_2 \end{bmatrix}$$

## Companion Form

In the companion realization, the characteristic polynomial of the system appears explicitly in the rightmost column of the  $A$  matrix. For a system with characteristic polynomial

$$p(s) = s^n + \alpha_1 s^{n-1} + \dots + \alpha_{n-1} s + \alpha_n$$

the corresponding companion  $A$  matrix is

$$A = \begin{bmatrix} 0 & 0 & \dots & \dots & 0 & -\alpha_n \\ 1 & 0 & 0 & \dots & 0 & -\alpha_{n-1} \\ 0 & 1 & 0 & \dots & \vdots & \vdots \\ \vdots & 0 & \dots & \dots & \vdots & \vdots \\ 0 & \dots & \dots & 1 & 0 & -\alpha_2 \\ 0 & \dots & \dots & 0 & 1 & -\alpha_1 \end{bmatrix}$$

The companion transformation requires that the system be controllable from the first input. The companion form is poorly conditioned for most state-space computations; avoid using it when possible.

## Algorithms

The `canon` command uses the `bdschur` command to convert `sys` into modal form and to compute the transformation `T`. If `sys` is not a state-space model, the algorithm first converts it to state space using `ss`.

The reduction to companion form uses a state similarity transformation based on the controllability matrix [1].

## References

[1] Kailath, T. *Linear Systems*, Prentice-Hall, 1980.

## See Also

`ctrb` | `ctrbf` | `ss2ss`

**Introduced before R2006a**

## chgFreqUnit

Change frequency units of frequency-response data model

### Syntax

```
sys_new = chgFreqUnit(sys,newfrequnits)
```

### Description

`sys_new = chgFreqUnit(sys,newfrequnits)` changes units of the frequency points in `sys` to `newfrequnits`. Both `Frequency` and `FrequencyUnit` properties of `sys` adjust so that the frequency responses of `sys` and `sys_new` match.

### Input Arguments

#### **sys**

Frequency-response data (`frd`, `idfrd`, or `genfrd`) model

#### **newfrequnits**

New units of frequency points, specified as one of the following values:

- 'rad/TimeUnit'
- 'cycles/TimeUnit'
- 'rad/s'
- 'Hz'
- 'kHz'
- 'MHz'
- 'GHz'
- 'rpm'

rad/TimeUnit and cycles/TimeUnit express frequency units relative to the system time units specified in the TimeUnit property.

**Default:** 'rad/TimeUnit'

## Output Arguments

**sys\_new**

Frequency-response data model of the same type as `sys` with new units of frequency points. The frequency response of `sys_new` is same as `sys`.

## Examples

### Change Frequency Units of Frequency-Response Data Model

Create a frequency-response data model.

```
load(fullfile(matlabroot, 'examples', 'controls_id', 'AnalyzerData'));  
sys = frd(resp, freq);
```

The data file `AnalyzerData` has column vectors `freq` and `resp`. These vectors contain 256 test frequencies and corresponding complex-valued frequency response points, respectively. The default frequency units of `sys` is `rad/TimeUnit`, where `TimeUnit` is the system time units.

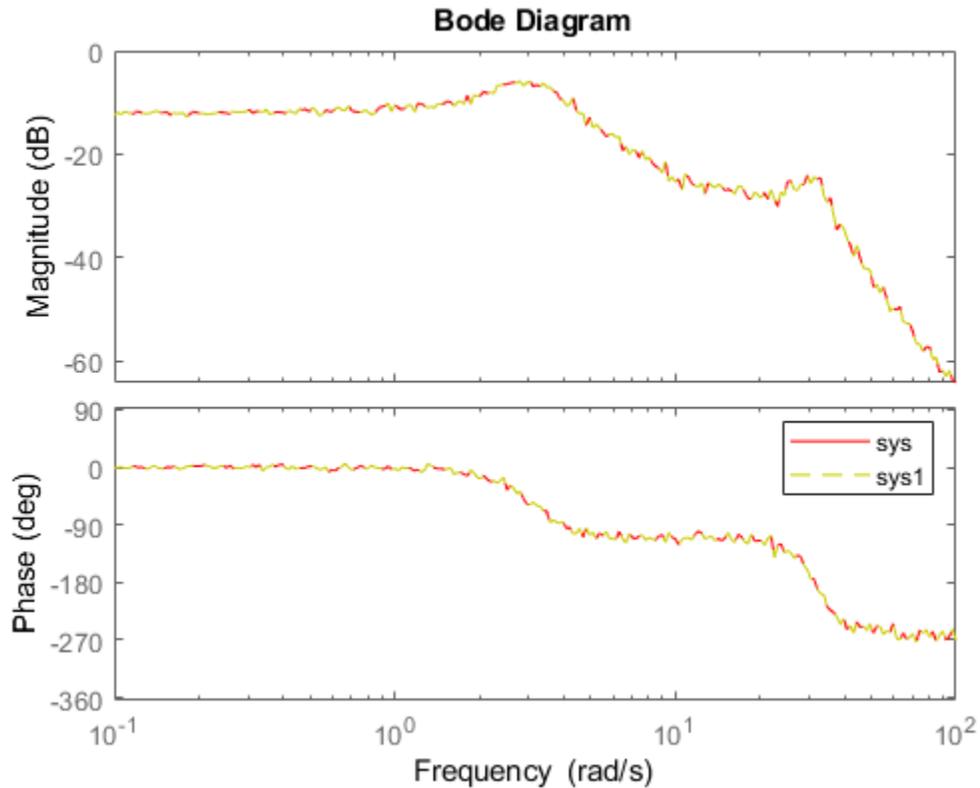
Change the frequency units.

```
sys1 = chgFreqUnit(sys, 'rpm');
```

The `FrequencyUnit` property of `sys1` is `rpm`.

Compare the Bode responses of `sys` and `sys1`.

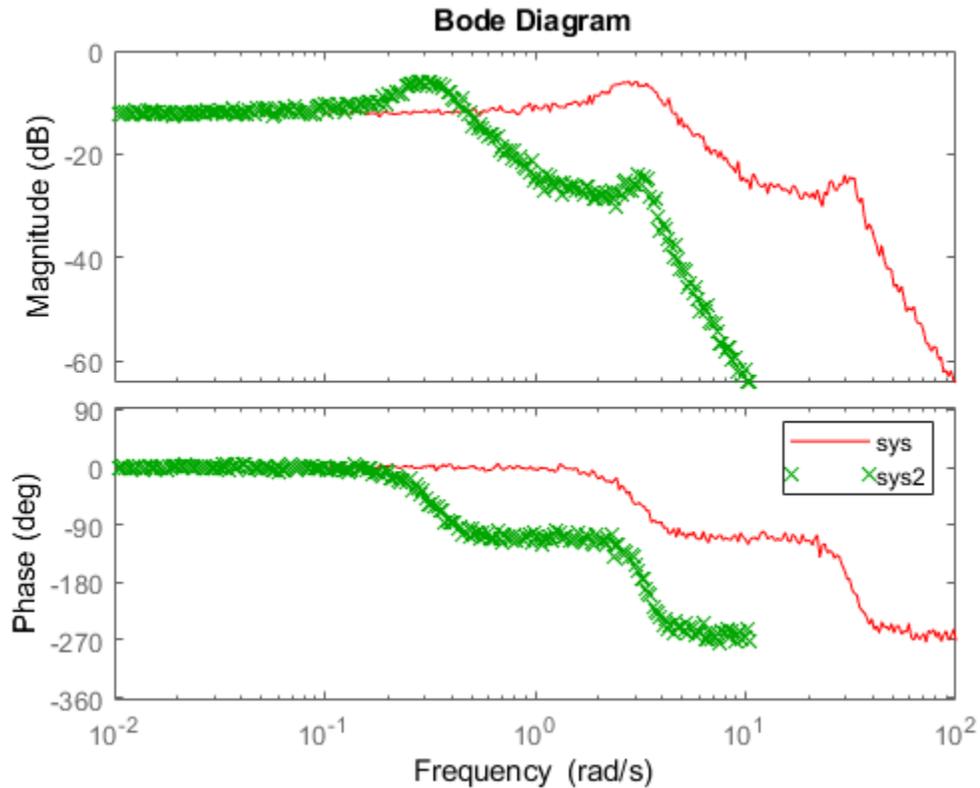
```
bodeplot(sys, 'r', sys1, 'y--');  
legend('sys', 'sys1')
```



The magnitude and phase of `sys` and `sys1` match because `chgFreqUnit` command changes the units of frequency points in `sys` without modifying system behavior.

Change the `FrequencyUnit` property of `sys` to compare the Bode response with the original system.

```
sys2 = sys;  
sys2.FrequencyUnit = 'rpm';  
bodeplot(sys, 'r', sys2, 'gx');  
legend('sys', 'sys2');
```



Changing the `FrequencyUnit` property changes the system behavior. Therefore, the Bode responses of `sys` and `sys2` do not match. For example, the original corner frequency at about 2 rad/s changes to approximately 2 rpm (or 0.2 rad/s).

## Tips

- Use `chgFreqUnit` to change the units of frequency points without modifying system behavior.

## **See Also**

chgTimeUnit | frd | idfrd

## **Topics**

“Specify Frequency Units of Frequency-Response Data Model” (Control System Toolbox)

**Introduced in R2012a**

# chgTimeUnit

Change time units of dynamic system

## Syntax

```
sys_new = chgTimeUnit(sys,newtimeunits)
```

## Description

`sys_new = chgTimeUnit(sys,newtimeunits)` changes the time units of `sys` to `newtimeunits`. The time- and frequency-domain characteristics of `sys` and `sys_new` match.

## Input Arguments

### **sys**

Dynamic system model

### **newtimeunits**

New time units, specified as one of the following values:

- 'nanoseconds'
- 'microseconds'
- 'milliseconds'
- 'seconds'
- 'minutes'
- 'hours'
- 'days'
- 'weeks'

- 'months'
- 'years'

**Default:** 'seconds'

## Output Arguments

### **sys\_new**

Dynamic system model of the same type as `sys` with new time units. The time response of `sys_new` is same as `sys`.

If `sys` is an identified linear model, both the model parameters as and their minimum and maximum bounds are scaled to the new time units.

## Examples

### **Change Time Units of Dynamic System Model**

Create a transfer function model.

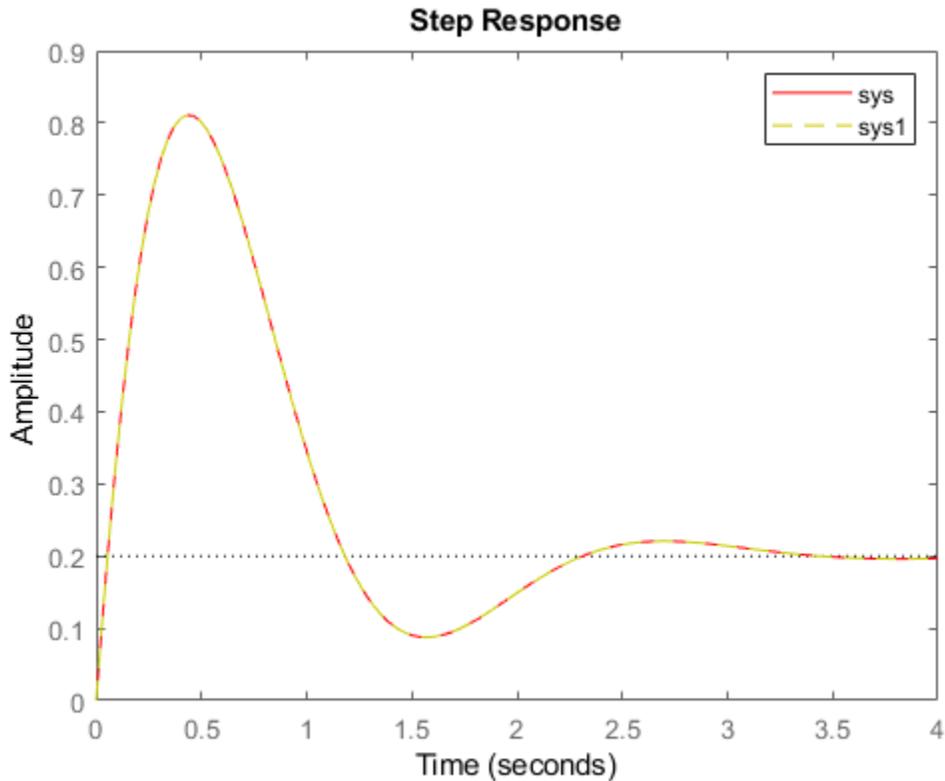
```
num = [4 2];  
den = [1 3 10];  
sys = tf(num,den);
```

By default, the time unit of `sys` is 'seconds'. Create a new model with the time units changed to minutes.

```
sys1 = chgTimeUnit(sys,'minutes');
```

This command sets the `TimeUnit` property of `sys1` to 'minutes', without changing the dynamics. To confirm that the dynamics are unchanged, compare the step responses of `sys` and `sys1`.

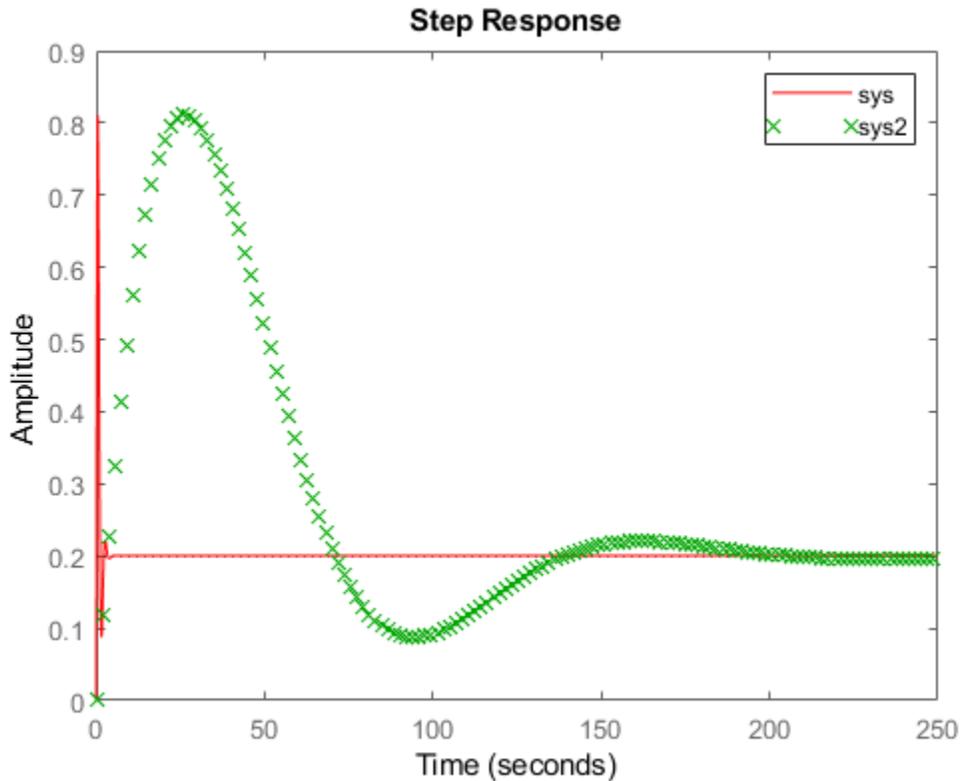
```
stepplot(sys,'r',sys1,'y--');  
legend('sys','sys1');
```



The step responses are the same.

If you change the `TimeUnit` property of the system instead of using `chgTimeUnit`, the dynamics of the system do change. To see this, change the `TimeUnit` property of a copy of `sys` and compare the step response with the original system.

```
sys2 = sys;  
sys2.TimeUnit = 'minutes';  
stepplot(sys, 'r', sys2, 'gx');  
legend('sys', 'sys2');
```



The step responses of `sys` and `sys2` do not match. For example, the original rise time of 0.04 seconds changes to 0.04 minutes.

## Tips

- Use `chgTimeUnit` to change the time units without modifying system behavior.

## See Also

`chgFreqUnit` | `frd` | `idpoly` | `idproc` | `idss` | `idtf` | `pid` | `ss` | `tf` | `zpk`

## **Topics**

“Specify Model Time Units” (Control System Toolbox)

**Introduced in R2012a**

## clone

Copy online parameter estimation System object

## Syntax

```
obj_clone = clone(obj)
```

## Description

`obj_clone = clone(obj)` creates a copy of the online parameter estimation System object™, `obj`, with the same property values. If the object you clone is locked, the new object is also locked.

`clone` is not supported for code generation using MATLAB Coder™.

---

**Note** If you want to copy an existing System object and then modify properties of the copied object, use the `clone` command. Do not create additional objects using syntax `obj2 = obj`. Any changes made to the properties of the new System object created this way (`obj2`) also change the properties of the original System object (`obj`).

---

## Examples

### Clone an Online Estimation System Object

Create a System object™ for online estimation of an ARX model with default properties.

```
obj = recursiveARX
```

```
obj =  
    recursiveARX with properties:
```

```
    A: []  
    B: []
```

```

        InitialA: [1 2.2204e-16]
        InitialB: [0 2.2204e-16]
    ParameterCovariance: []
InitialParameterCovariance: [2x2 double]
    EstimationMethod: 'ForgettingFactor'
    ForgettingFactor: 1
    EnableAdaptation: true
        History: 'Infinite'
    InputProcessing: 'Sample-based'
        DataType: 'double'

```

Use `clone` to generate an object with the same properties as the original object.

```
obj2 = clone(obj)
```

```
obj2 =
```

```
recursiveARX with properties:
```

```

        A: []
        B: []
        InitialA: [1 2.2204e-16]
        InitialB: [0 2.2204e-16]
    ParameterCovariance: []
InitialParameterCovariance: [2x2 double]
    EstimationMethod: 'ForgettingFactor'
    ForgettingFactor: 1
    EnableAdaptation: true
        History: 'Infinite'
    InputProcessing: 'Sample-based'
        DataType: 'double'

```

## Input Arguments

### **obj** — System object for online parameter estimation

recursiveAR object | recursiveARMA object | recursiveARX object |  
 recursiveARMAX object | recursiveOE object | recursiveBJ object | recursiveLS  
 object

System object for online parameter estimation, created using one of the following commands:

- recursiveAR
- recursiveARMA
- recursiveARX
- recursiveARMAX
- recursiveOE
- recursiveBJ
- recursiveLS

## Output Arguments

### **obj\_clone** — Copy of online estimation System object

System object

Copy of online estimation System object, `obj`, returned as a System object with the same properties as `obj`.

## See Also

`isLocked` | `recursiveAR` | `recursiveARMA` | `recursiveARMAX` | `recursiveARX` | `recursiveBJ` | `recursiveLS` | `recursiveOE` | `release` | `reset` | `step`

## Topics

“What Is Online Estimation?”

**Introduced in R2015b**

# clone

Copy online state estimation object

## Syntax

```
obj_clone = clone(obj)
```

## Description

`obj_clone = clone(obj)` creates a copy of the online state estimation object `obj` with the same property values.

If you want to copy an existing object and then modify properties of the copied object, use the `clone` command. Do not create additional objects using syntax `obj2 = obj`. Any changes made to the properties of the new object created in this way (`obj2`) also change the properties of the original object (`obj`).

## Examples

### Clone an Online State Estimation Object

Create an extended Kalman filter object for a van der Pol oscillator with two states and one output. To create the object, use the previously written and saved state transition and measurement functions, `vdpStateFcn.m` and `vdpMeasurementFcn.m`. Specify the initial state values for the two states as `[2;0]`.

```
obj = extendedKalmanFilter(@vdpStateFcn,@vdpMeasurementFcn,[2;0])
```

```
obj =  
    extendedKalmanFilter with properties:  
        HasAdditiveProcessNoise: 1  
        StateTransitionFcn: @vdpStateFcn  
        HasAdditiveMeasurementNoise: 1
```

```
MeasurementFcn: @vdpMeasurementFcn
StateTransitionJacobianFcn: []
MeasurementJacobianFcn: []
State: [2x1 double]
StateCovariance: [2x2 double]
ProcessNoise: [2x2 double]
MeasurementNoise: 1
```

Use `clone` to generate an object with the same properties as the original object.

```
obj2 = clone(obj)
```

```
obj2 =
  extendedKalmanFilter with properties:
```

```
HasAdditiveProcessNoise: 1
StateTransitionFcn: @vdpStateFcn
HasAdditiveMeasurementNoise: 1
MeasurementFcn: @vdpMeasurementFcn
StateTransitionJacobianFcn: []
MeasurementJacobianFcn: []
State: [2x1 double]
StateCovariance: [2x2 double]
ProcessNoise: [2x2 double]
MeasurementNoise: 1
```

Modify the `MeasurementNoise` property of `obj2`.

```
obj2.MeasurementNoise = 2;
```

Verify that `MeasurementNoise` property of original object `obj` remains unchanged and equals 1.

```
obj.MeasurementNoise
```

```
ans = 1
```

## Input Arguments

### **obj** — Object for online state estimation

`extendedKalmanFilter` object | `unscentedKalmanFilter` object | `particleFilter` object

Object for online state estimation of a nonlinear system, created using one of the following commands:

- `extendedKalmanFilter`
- `unscentedKalmanFilter`
- `particleFilter`

## Output Arguments

### **obj\_clone** — Clone of online state estimation object

`extendedKalmanFilter` object | `unscentedKalmanFilter` object | `particleFilter` object

Clone of online state estimation object `obj`, returned as an `extendedKalmanFilter`, `unscentedKalmanFilter` or `particleFilter` object with the same properties as `obj`.

## See Also

`correct` | `extendedKalmanFilter` | `initialize` | `particleFilter` | `predict` | `unscentedKalmanFilter`

## Topics

“What Is Online Estimation?”

**Introduced in R2016b**

## compare

Compare identified model output and measured output

### Syntax

```
compare(data,sys)
compare(data,sys,kstep)
compare(data,sys,LineStyle,kstep)
compare(data,sys1,...,sysN,kstep)
compare(data,sys1,LineStyle1,...,sysN,LineStyleN,kstep)
compare( __ ,opt)
```

```
[y,fit,x0] = compare( __ )
```

### Description

`compare(data,sys)` simulates the response of a dynamic system model, and superimposes that response over plotted measurement data. The plot also displays the normalized root mean square (NRMSE) measure of the goodness of the fit between simulated response and measurement data. Use this function when you want to evaluate a set of candidate models identified from the same measurement data, or when you want to validate a model you have selected. You can use `compare` with time-domain or frequency-domain models and data.

`compare(data,sys,kstep)` also predicts the response of `sys`, using a prediction horizon specified by `kstep`. Prediction uses output measurements as well as input measurements to project a future response. `kstep` represents the number of time samples between the timepoint of each output measurement and the timepoint of the resulting predicted response. For more information on prediction, see “Simulate and Predict Identified Model Output”.

`compare(data,sys,LineStyle,kstep)` also specifies the line type, marker symbol, and color for the model response.

`compare(data,sys1,...,sysN,kstep)` compares the responses of multiple dynamic systems on the same axes. `compare` automatically chooses the line specifications.

`compare(data, sys1, LineSpec1, ..., sysN, LineSpecN, kstep)` also compares the responses of multiple systems on the same axes using the line type, marker symbol, and color specified for each system.

`compare(____, opt)` configures the comparison using an option set. Options include initial condition handling, data offsets, and data selection. You can use this syntax with any of the previous input-argument combinations.

`[y, fit, x0] = compare(____)` returns the model response,  $y$ , goodness of fit value,  $fit$ , and the initial states,  $x_0$ . No plot is generated. You can use this syntax with any of the previous input-argument combinations. However, line specifications are ignored, because there is no plotting.

## Examples

### Compare Response of Estimated Model to Measured Data

Identify a linear model and visualize the simulated model response with the data from which it was generated.

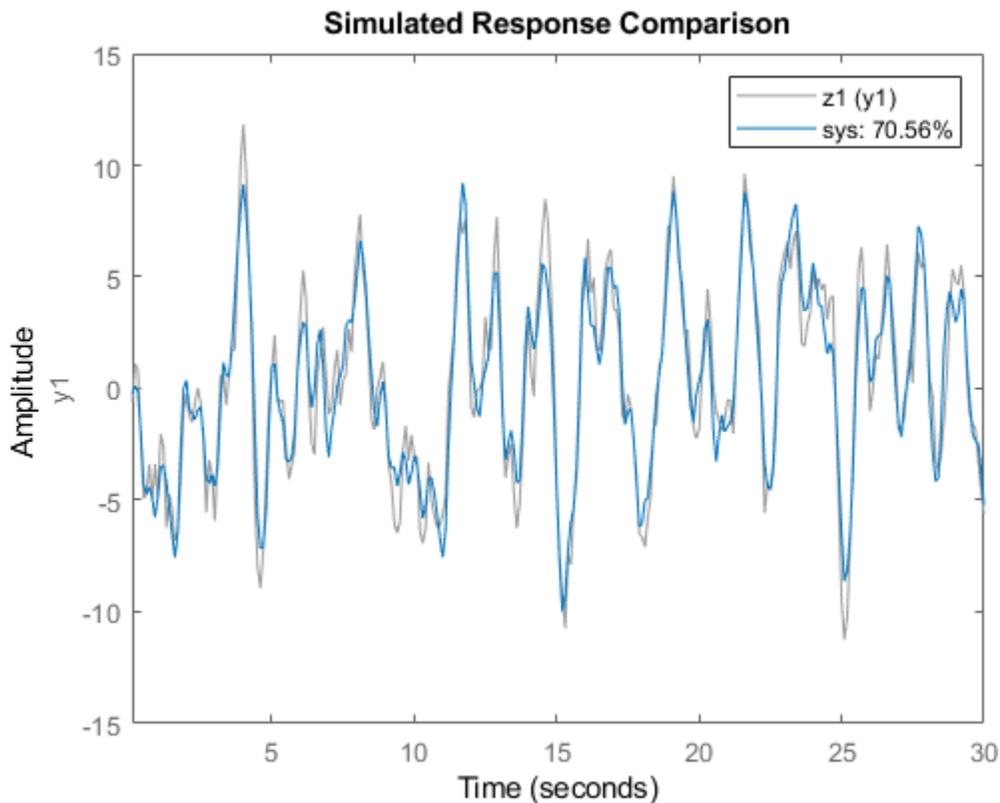
Load input/output measurements `z1`, and identify a third-order state-space model `sys`.

```
load iddata1 z1;  
sys = ssest(z1,3);
```

`sys` is a continuous-time identified state-space (`idss`) model.

Use `compare` to simulate the `sys` response and plot it alongside the data `z1`.

```
figure  
compare(z1,sys)
```



The plot illustrates the differences between the model response and the original data. The percentage shown in the legend is the NRMSE fitness value. It represents how close the predicted model output is to the data.

To change display options in the plot, right-click the plot to access the context menu. For example:

- To plot the error between the predicted output and measured output, select **Error Plot**.
- To view the confidence region for the simulated response, select **Characteristics -> ConfidenceRegion**.
- To specify number of standard deviations to plot, double-click the plot and open the Property Editor dialog box. In the **Options** tab, specify the number of standard

deviations in **Confidence Region for Identified Models**. The default value is 1 standard deviation.

### **Compare Predicted Response of Identified Time-Domain Model to Measured Data**

Identify a linear model and visualize the predicted model response with the data from which it was computed.

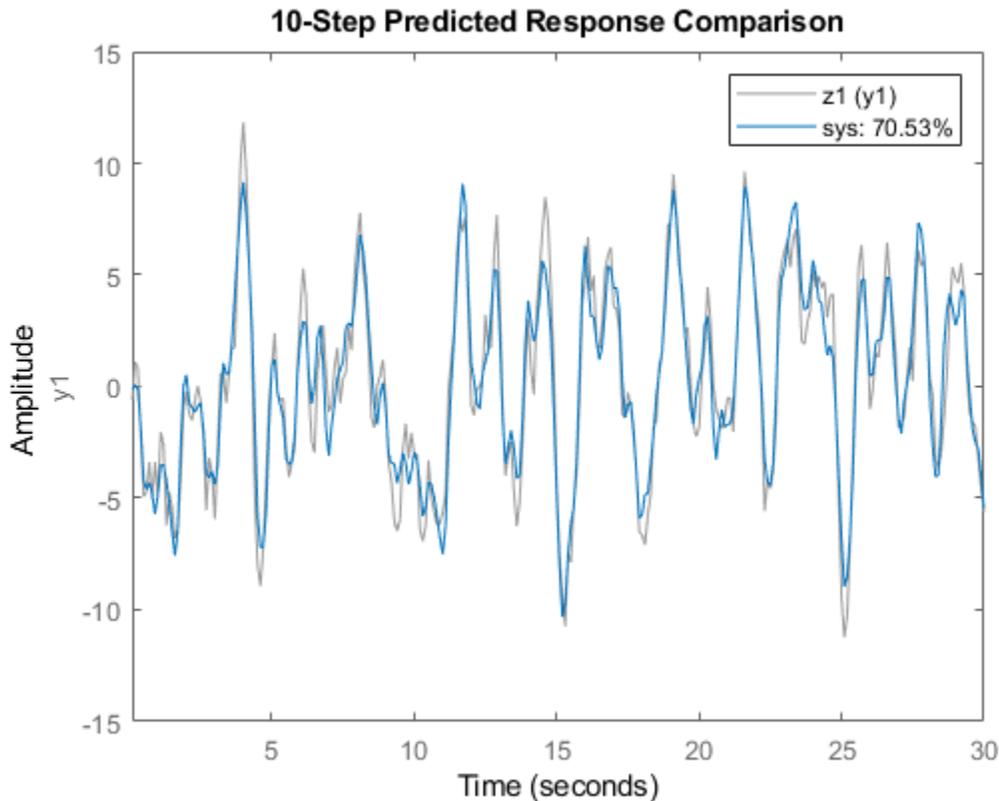
Identify a third-order state-space model using the input/output measurements in `z1`.

```
load iddata1 z1;  
sys = ssest(z1,3);
```

`sys` is a continuous-time identified state-space (`idss`) model.

Now use `compare` to plot the predicted response. Prediction differs from simulation in that it uses both measured input and measured output when computing the system response. The prediction horizon defines how far in the future to predict, relative to your current measured output point. For this example, set the prediction horizon `kstep` to 10 steps, and use `compare` to plot the predicted response against the original measurement data.

```
kstep = 10;  
compare(z1,sys,kstep)
```



In this plot, each sys data point represents the predicted output associated with output measurement data that was taken at least 10 steps earlier. For instance, the point at  $t = 15\text{s}$  is based on output measurements taken at or prior to  $t = 5\text{s}$ . The calculation of this  $t = 15\text{s}$  sys data point also uses input measurements up to  $t = 15\text{s}$ , just as a simulation would.

The plot illustrates the differences between the model response and the original data. The percentage shown in the legend is the NRMSE fitness value. It represents how closely the predicted model output matches the data.

To change display and simulation options in the plot, right-click the plot to access the context menu. For example, to plot the error between the predicted output and measured output, select **Error Plot** from the context menu. To change the prediction horizon value,

or to toggle between simulation and prediction, select **Prediction Horizon** from the context menu.

### Compare Multiple Identified Models to Measured Time-Domain Data

Identify several model types for the same data, and compare the results to see which best fits the data.

Load the data, which contains `iddata` object `z1` with single input and output.

```
load iddata1;
```

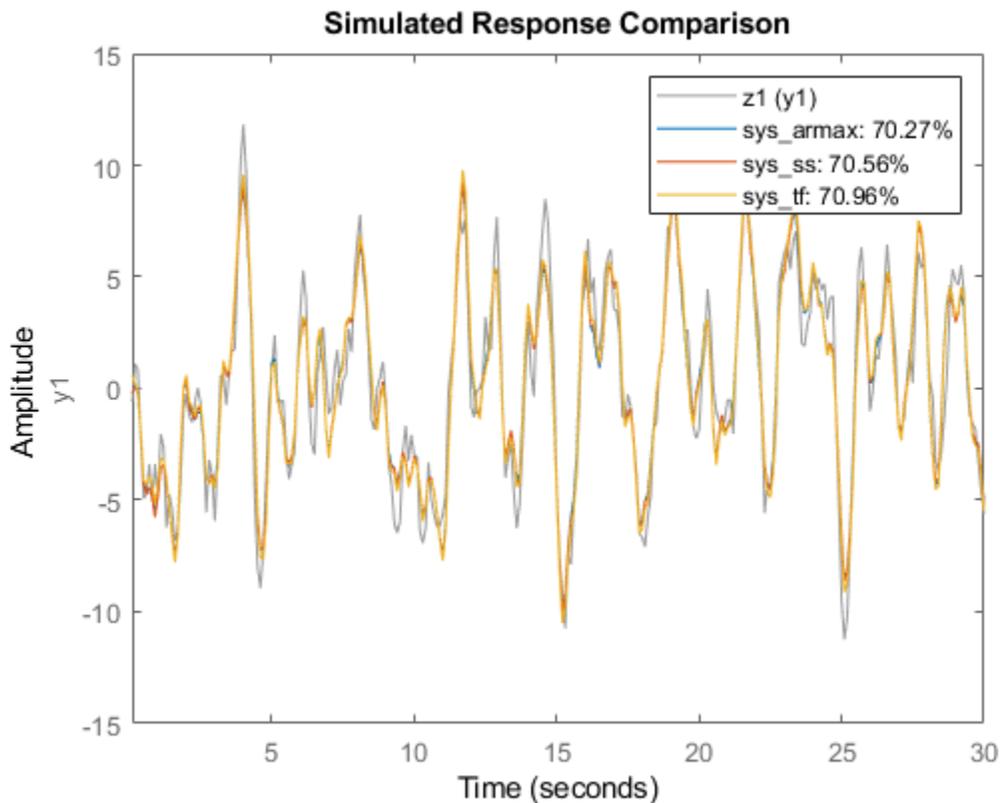
From `z1`, identify a model for each of the following linear forms:

- ARMAX (`idpoly`) of orders 2, 3, and 1, with dead time of 0
- State space (`idss`) with three states
- Transfer function (`idtf`) with three poles

```
sys_armax = armax(z1,[2 3 1 0]);  
sys_ss = ssest(z1,3);  
sys_tf = tfest(z1,3);
```

Using `compare`, plot the simulated responses for the three models with `z1`.

```
compare(z1,sys_armax,sys_ss,sys_tf)
```



For this set of data, along with the default settings for all the models, the transfer-function form has the best NRMSE fit. However, the fits for all models are within about 1% of each other.

You can interactively control which model responses are displayed in the plot by right-clicking on the plot and hovering over **Systems**.

### Compare Multiple Estimated Models to Measured Frequency-Domain Data

Compare the outputs of multiple estimated models of differing types to measured frequency-domain data.

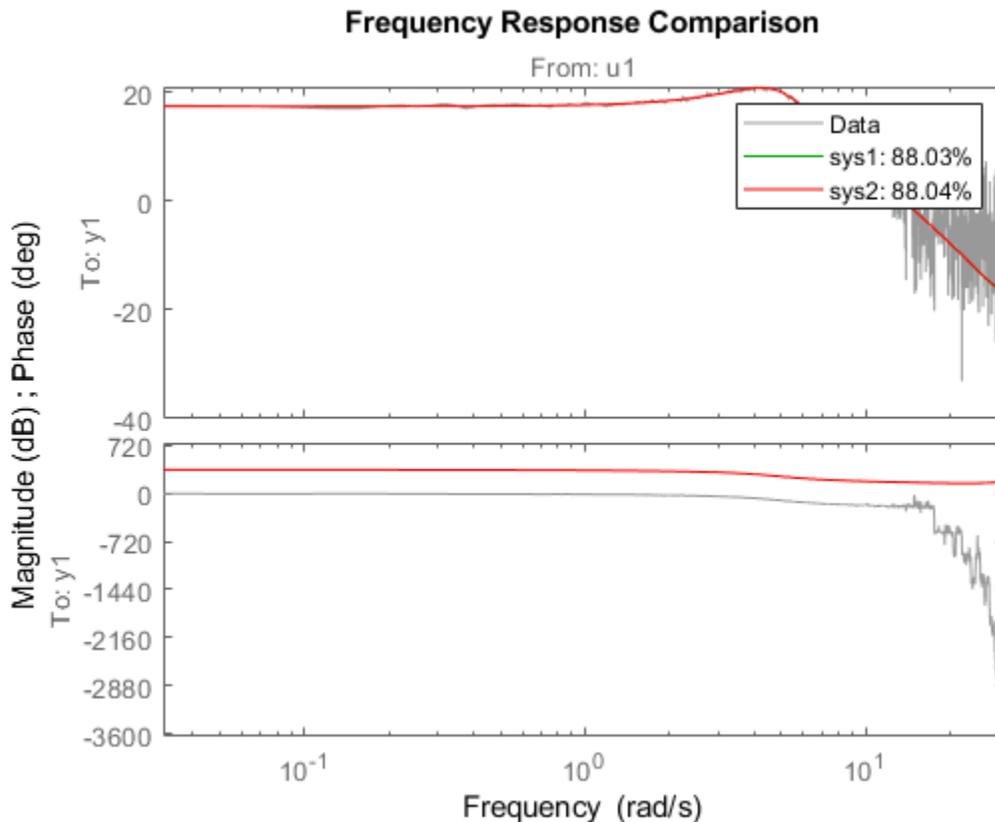
For this example, estimate a process model and an output-error polynomial from frequency response data.

```
load demofr % frequency response data
zfr = AMP.*exp(1i*PHA*pi/180);
Ts = 0.1;
data = idfrd(zfr,W,Ts);
sys1 = procest(data,'P2UDZ');
sys2 = oe(data,[2 2 1]);
```

sys1, an idproc model, is a continuous-time process model. sys2, an idpoly model, is a discrete-time output-error model.

Compare the frequency response of the estimated models to data.

```
compare(data,sys1,'g',sys2,'r');
```



The two models have NRMSE fit values that are nearly equal with respect to the data from which they were calculated.

### Compare Estimated Model to Data and Specify Comparison Options

Modify default behavior when you compare an estimated model to measured data.

Estimate a transfer function for measured data.

```
load iddata1 z1;
sys = tfest(z1,3);
```

`sys` is a continuous-time identified transfer function (`idtf`) model.

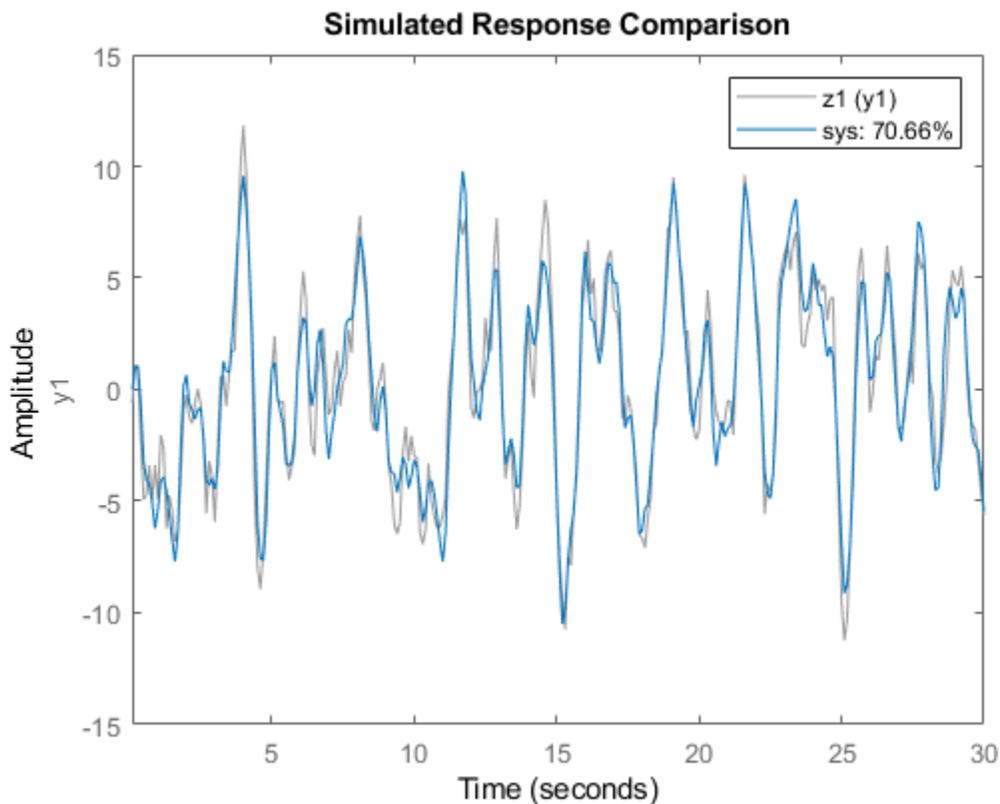
Suppose you want your initial conditions to be zero. The default for `compare` is to estimate initial conditions from the data.

Create an option set to specify the initial condition handling. To use zero for initial conditions, specify `'z'` for the `'InitialCondition'` option.

```
opt = compareOptions('InitialCondition','z');
```

Compare the estimated transfer function model output to the measured data using the comparison option set.

```
compare(z1,sys,opt)
```



## Input Arguments

### **data — Validation data**

`iddata` object | `idfrd` object | `frd` object

Validation data, specified as an `iddata`, `idfrd`, or `frd` object.

If `sys` is:

- An `iddata` object, then `data` must be an `iddata` object with matching domain, number of experiments and time or frequency vectors
- A frequency-response data (FRD) model (defined as either `idfrd` or `frd`), then `data` must also be FRD
- A parametric model (such as `idss`), then `data` can be `iddata` or FRD

`data` can represent either time-domain or frequency-domain data when comparing with linear models. `data` must be time-domain data when comparing with a nonlinear model.

For examples, see:

- “Compare Predicted Response of Identified Time-Domain Model to Measured Data” on page 1-211
- “Compare Multiple Estimated Models to Measured Frequency-Domain Data” on page 1-214

.

### **sys — Identified model**

dynamic system model | `iddata` object | model array

Identified model, specified as a dynamic system model, an `iddata` object, or a model array.

When the time or frequency units of `data` do not match the units of `sys`, `compare` rescales `sys` to match the units of `data`.

### **kstep — Prediction Horizon — steps ahead to predict**

`Inf` (default) | integer

Prediction horizon, specified as one of the following:

- **Inf** — Compare simulated response of `sys` to `data`.
- Positive finite integer — Compare predicted response of `sys` to `data`, where each predicted response point is based not only on measured input data up to that timepoint, but also on measured output data up to `kstep` timepoints earlier.

`compare` ignores `kstep` when `sys` is an `iddata` object, an FRD model, or a dynamic system with no noise component. `compare` also ignores `kstep` when using frequency response validation data.

If you specify `kstep` that is greater than the number of data samples, `compare` sets `kstep` to `Inf` and provides a warning message.

For more information on simulation and prediction, see “Simulate and Predict Identified Model Output”.

For an example, see “Compare Predicted Response of Identified Time-Domain Model to Measured Data” on page 1-211.

### **LineStyle** — Line style, marker, and color

character vector

Line style, marker, and color of both the line and marker, specified as a character vector, such as `'b'` or `'b+:'`.

For more information about configuring `LineStyle`, see the `LineStyle` input argument of `plot`. For an example, see “Compare Multiple Estimated Models to Measured Frequency-Domain Data” on page 1-214.

### **opt** — Comparison options

`compareOptions` object

Comparison options, specified as an option set you create using `compareOptions`.

Available options include:

- Handling of initial conditions
- Sample range for computing fit numbers
- Data offsets
- Output weighting

For examples, see:

- “Compare Estimated Model to Data and Specify Comparison Options” on page 1-216
- “Resolve Fit Value Differences Between Model Identification and compare Command”

## Output Arguments

### **y** — Model response

`iddata` object | `idfrd` object | cell array | array

Model response, returned as an `iddata` object, an `idfrd` object, a cell array, or an array. The output depends on the models and data you provide, as follows:

- For a single model and single-experiment data set, **y** is an `iddata` object or `idfrd` object
- For multimodel comparisons, **y** is a cell array with one `iddata` or `idfrd` object entry for each input model
- For multiexperiment data, **y** is a cell array with one entry for each experiment
- For multimodel comparisons using multiexperiment data, **y** is an  $N_{\text{sys}}$ -by- $N_{\text{exp}}$  cell array, where  $N_{\text{sys}}$  is the number of models, and  $N_{\text{exp}}$  is the number of experiments
- If `sys` is a model array, **y** is an array with an element corresponding to each model in `sys` and experiment in `data`. For more information on model arrays, see `stack`

If `kstep` is not specified or is `Inf`, then `compare` returns the simulated response in **y**.

Otherwise, `compare` returns the predicted response. Measured output values in `data` up to time  $t_{n-k\text{step}}$  are used to predict the output of `sys` at time  $t_n$ . For more information on simulation and prediction, see “Simulate and Predict Identified Model Output”.

The `compare` response computation requires specification of initial condition handling. By default, `compare` estimates the initial conditions to maximize the fit to data. See `compareOptions` for more information on how `compare` determines the initial conditions to use.

### **fit** — NRMSE fitness value

vector | matrix | cell array

NRMSE fitness value indicator of how well the simulated or predicted model response matches the measurement data, returned as a vector, a matrix, or a cell array. The output depends on the models and data you provide, as follows:

- If `data` is an `iddata` object, `fit` is a vector of length  $N_y$ , where  $N_y$  is the number of outputs
- If `data` is an FRD model, `fit` is an  $N_y$ -by- $N_u$  matrix, where  $N_u$  is the number of inputs in `data`
- For a single model and single-experiment data set, `fit` is a vector or matrix
- For multimodel comparisons, `fit` is a cell array with one entry for each input model
- For multiexperiment data, `fit` is a cell array with one entry for each experiment
- For multimodel comparisons using multiexperiment data, `fit` is an  $N_{\text{sys}}$ -by- $N_{\text{exp}}$  cell array, where  $N_{\text{sys}}$  is the number of models, and  $N_{\text{exp}}$  is the number of experiments
- If `sys` is a model array, `fit` is an array with an element corresponding to each model in `sys` and experiment in `data`

`compare` calculates `fit` (in percentage) using:

$$\text{fit} = 100 \left( 1 - \frac{\|y - \hat{y}\|}{\|y - \text{mean}(y)\|} \right),$$

where  $y$  is the validation data output and  $\hat{y}$  is the output of `sys`.

For FRD models — `compare` calculates `fit` by comparing the complex frequency response. The fits of the magnitude and phase curves shown in the `compare` plot are not computed by `compare` separately.

### **x0 — Initial conditions used to compute system response**

[ ] | vector | cell array

Initial conditions used to compute system response, returned as an empty array, a vector, or a cell array. The output depends on the models and data you provide, as follows:

- When `sys` is an `frd` or `iddata` object, `x0` is the empty array [ ], because initial conditions cannot be used with these objects
- For a single model and single-experiment data set, `x0` is a vector
- For multimodel comparisons, `x0` is a cell array, with one vector or matrix entry for each input model
- For multiexperiment data, `x0` is a cell array, with one entry for each experiment

- For multimodel comparisons using multiexperiment data,  $x_0$  is an  $N_{\text{sys}}$ -by- $N_{\text{exp}}$  cell array, where  $N_{\text{sys}}$  is the number of models, and  $N_{\text{exp}}$  is the number of experiments
- If `sys` is a model array,  $x_0$  is an array with an element corresponding to each model in `sys` and experiment in `data`

By default, `compare` uses `findstates` to estimate  $x_0$ . To change this behavior, set the 'InitialCondition' option in `opt` (see `compareOptions`). If you have input/output history that immediately precedes your start point, you can set 'InitialCondition' to that history data. `compare` then uses `data2state` to compute the end state of the history data, and thus the start state for the simulation. Other choices include setting initial conditions to zero, or to specific values that you have determined previously. For more information about finding initial conditions, see “Estimate Initial Conditions for Simulating Identified Models”.

If you are using an estimation model that does not explicitly use states, `compare` first converts the model to its state-space representation and then maps the data to states. For more information, see `compareOptions`.

## Tips

- The NRMSE fit result you obtain with `compare` may not precisely match the fit value reported in model identification. These differences typically arise from mismatches in initial conditions, and in the differences in the prediction horizon defaults for identification and for validation. The differences are generally small, and should not impact your model selection and validation workflow. For more information, see “Resolve Fit Value Differences Between Model Identification and `compare` Command”.
- `compare` matches the input/output channels in `data` and `sys` based on the channel names. Thus, it is possible to evaluate models that do not use all the input channels that are available in `data`. This flexibility allows you to compare multiple models which were each identified independently from different sets of input/output channels.
- The `compare` plot allows you to vary key parameters. For example, you can interactively control:
  - Whether you generate a simulated or predicted response
  - Prediction horizon value
  - Initial condition handling
  - Which experiment data you view

- Which system models you view

To access the controls, right-click the plot to bring up the options menu.

## See Also

`chgFreqUnit` | `chgTimeUnit` | `compareOptions` | `forecast` | `goodnessOfFit` | `plot`  
| `predict` | `sim`

## Topics

“Compare Simulated Output with Measured Validation Data”

“Validating Models After Estimation”

“Model Validation”

“Estimate Initial Conditions for Simulating Identified Models”

**Introduced before R2006a**

## **compareOptions**

Option set for compare

### **Syntax**

```
opt = compareOptions  
opt = compareOptions(Name, Value)
```

### **Description**

`opt = compareOptions` creates the default options set for `compare`.

`opt = compareOptions(Name, Value)` creates an option set with the options specified by one or more `Name, Value` pair arguments.

### **Input Arguments**

#### **Name-Value Pair Arguments**

Specify optional comma-separated pairs of `Name, Value` arguments. `Name` is the argument name and `Value` is the corresponding value. `Name` must appear inside quotes. You can specify several name and value pair arguments in any order as `Name1, Value1, ..., NameN, ValueN`.

#### **Samples**

Data for which `compare` calculates fit values.

Specify `Samples` as a vector containing the data sample indices. For multiexperiment data, use a cell array of  $N_e$  vectors, where  $N_e$  is the number of experiments.

#### **InitialCondition**

Handling of initial conditions.

Specify `InitialCondition` as one of the following:

- 'z' — Zero initial conditions.
- 'e' — Estimate initial conditions such that the prediction error for observed output is minimized.

For nonlinear grey-box models, only those initial states `i` that are designated as free in the model (`sys.InitialStates(i).Fixed = false`) are estimated. To estimate all the states of the model, first specify all the `Nx` states of the `idnlgrey` model `sys` as free.

```
for i = 1:Nx
    sys.InitialStates(i).Fixed = false;
end
```

Similarly, to fix all the initial states to values specified in `sys.InitialStates`, first specify all the states as fixed in the `sys.InitialStates` property of the nonlinear grey-box model.

- 'd' — Similar to 'e', but absorbs nonzero delays into the model coefficients. The delays are first converted to explicit model states, and the initial values of those states are also estimated and returned.

Use this option for linear models only.

- Vector or Matrix — Initial guess for state values, specified as a numerical column vector of length equal to the number of states. For multi-experiment data, specify a matrix with `Ne` columns, where `Ne` is the number of experiments. Otherwise, use a column vector to specify the same initial conditions for all experiments. Use this option for state-space (`idss` and `idgrey`) and nonlinear models (`idnlarx`, `idnlhw`, and `idnlgrey`) only.
- Structure with the following fields, which contain the historical input and output values for a time interval immediately before the start time of the data used by `compare`:

Field	Description
Input	Input history, specified as a matrix with <code>Nu</code> columns, where <code>Nu</code> is the number of input channels. For time series models, use <code>[]</code> . The number of rows must be greater than or equal to the model order.

Field	Description
Output	Output history, specified as a matrix with $N_y$ columns, where $N_y$ is the number of output channels. The number of rows must be greater than or equal to the model order.

For multi-experiment data, configure the initial conditions separately for each experiment by specifying `InitialCondition` as a structure array with  $N_e$  elements. To specify the same initial conditions for all experiments, use a single structure.

The software uses `data2state` to map the historical data to states. If your model is not `idss`, `idgrey`, `idnlgrey`, or `idnlrx`, the software first converts the model to its state-space representation and then maps the data to states. If conversion of your model to `idss` is not possible, the estimated states are returned empty.

- `x0obj` — Specification object created using `idpar`. Use this object for discrete-time state-space (`idss` and `idgrey`) and nonlinear grey-box (`idnlgrey`) models only. Use `x0obj` to impose constraints on the initial states by fixing their value or specifying minimum or maximum bounds.

**Default:** 'e'

### **InputOffset**

Removes offset from time domain input data for model response computation.

Specify as a column vector of length  $N_u$ , where  $N_u$  is the number of inputs.

Use `[]` to indicate no offset.

For multiexperiment data, specify `InputOffset` as a  $N_u$ -by- $N_e$  matrix.  $N_u$  is the number of inputs and  $N_e$  is the number of experiments.

Each entry specified by `InputOffset` is subtracted from the corresponding input data.

**Default:** `[]`

### **OutputOffset**

Removes offset from time-domain output data for model response prediction.

Specify as a column vector of length  $N_y$ , where  $N_y$  is the number of outputs.

Use `[]` to indicate no offset.

For multiexperiment data, specify `OutputOffset` as a  $N_y$ -by- $N_e$  matrix.  $N_y$  is the number of outputs and  $N_e$  is the number of experiments.

Each entry specified by `OutputOffset` is subtracted from the corresponding output data before computing the model response. After computing the model response, the software adds the offset to the response to give the final model response.

**Default:** []

### **OutputWeight**

Weight of output for initial condition estimation.

`OutputWeight` requires one of the following values:

- [] — No weighting is used. This option is the same as using `eye(Ny)` for the output weight.  $N_y$  is the number of outputs.
- 'noise' — Inverse of the noise variance stored with the model.
- Matrix of doubles — A positive semi-definite matrix of dimension  $N_y$ -by- $N_y$ .  $N_y$  is the number of outputs.

**Default:** []

## **Output Arguments**

### **opt**

Option set containing the specified options for `compare`.

## **Examples**

### **Create Default Options Set for Model Comparison**

Create a default options set for `compare`.

```
opt = compareOptions;
```

## Specify Options for Model Comparison

Create an options set for `compare` using zero initial conditions. Set the input offset to 5.

```
opt = compareOptions('InitialCondition','z','InputOffset',5);
```

Alternatively, use dot notation to set the values of `opt`.

```
opt = compareOptions;  
opt.InitialCondition = 'z';  
opt.InputOffset = 5;
```

## See Also

`compare`

**Introduced in R2012a**

## correct

Correct state and state estimation error covariance using extended or unscented Kalman filter, or particle filter and measurements

The `correct` command updates the state and state estimation error covariance of an `extendedKalmanFilter`, `unscentedKalmanFilter` or `particleFilter` object using measured system outputs. To implement extended or unscented Kalman filter, or particle filter, use the `correct` and `predict` commands together. If the current output measurement exists, you can use `correct` and `predict`. If the measurement is missing, you can only use `predict`. For information about the order in which to use the commands, see “Using predict and correct Commands” on page 1-236.

## Syntax

```
[CorrectedState,CorrectedStateCovariance] = correct(obj,y)
[CorrectedState,CorrectedStateCovariance] = correct(obj,y,
Um1,...,Umn)
```

## Description

`[CorrectedState,CorrectedStateCovariance] = correct(obj,y)` corrects the state estimate and state estimation error covariance of an extended or unscented Kalman filter, or particle filter object `obj` using the measured output `y`.

You create `obj` using the `extendedKalmanFilter`, `unscentedKalmanFilter` or `particleFilter` commands. You specify the state transition function and measurement function of your nonlinear system in `obj`. You also specify whether the process and measurement noise terms are additive or nonadditive in these functions. The `State` property of the object stores the latest estimated state value. Assume that at time step `k`,

`obj.State` is  $\hat{x}[k|k-1]$ . This value is the state estimate for time `k`, estimated using measured outputs until time `k-1`. When you use the `correct` command with measured system output `y[k]`, the software returns the corrected state estimate  $\hat{x}[k|k]$  in the `CorrectedState` output. Where  $\hat{x}[k|k]$  is the state estimate at time `k`, estimated using

measured outputs until time  $k$ . The command returns the state estimation error covariance of  $\hat{x}[k|k]$  in the `CorrectedStateCovariance` output. The software also updates the `State` and `StateCovariance` properties of `obj` with these corrected values.

Use this syntax if the measurement function  $h$  that you specified in `obj.MeasurementFcn` has one of the following forms:

- $y(k) = h(x(k))$  — for additive measurement noise.
- $y(k) = h(x(k), v(k))$  — for nonadditive measurement noise.

Where  $y(k)$ ,  $x(k)$ , and  $v(k)$  are the measured output, states, and measurement noise of the system at time step  $k$ . The only inputs to  $h$  are the states and measurement noise.

`[CorrectedState, CorrectedStateCovariance] = correct(obj, y, Um1, ..., Umn)` specifies additional input arguments, if the measurement function of the system requires these inputs. You can specify multiple arguments.

Use this syntax if the measurement function  $h$  has one of the following forms:

- $y(k) = h(x(k), Um1, \dots, Umn)$  — for additive measurement noise.
- $y(k) = h(x(k), v(k), Um1, \dots, Umn)$  — for nonadditive measurement noise.

`correct` command passes these inputs to the measurement function to calculate the estimated outputs.

## Examples

### Estimate States Online Using Extended Kalman Filter

Estimate the states of a van der Pol oscillator using an extended Kalman filter algorithm and measured output data. The oscillator has two states and one output.

Create an extended Kalman filter object for the oscillator. Use previously written and saved state transition and measurement functions, `vdpStateFcn.m` and `vdpMeasurementFcn.m`. These functions describe a discrete-approximation to a van der Pol oscillator with nonlinearity parameter,  $\mu$ , equal to 1. The functions assume additive process and measurement noise in the system. Specify the initial state values for the two

states as  $[1;0]$ . This is the guess for the state value at initial time  $k$ , using knowledge of system outputs until time  $k-1$ ,  $\hat{x}[k|k-1]$ .

```
obj = extendedKalmanFilter(@vdpStateFcn,@vdpMeasurementFcn,[1;0]);
```

Load the measured output data,  $y$ , from the oscillator. In this example, use simulated static data for illustration. The data is stored in the `vdp_data.mat` file.

```
load vdp_data.mat y
```

Specify the process noise and measurement noise covariances of the oscillator.

```
obj.ProcessNoise = 0.01;
obj.MeasurementNoise = 0.16;
```

Implement the extended Kalman filter algorithm to estimate the states of the oscillator by using the `correct` and `predict` commands. You first correct  $\hat{x}[k|k-1]$  using measurements at time  $k$  to get  $\hat{x}[k|k]$ . Then, you predict the state value at next time step,  $\hat{x}[k+1|k]$ , using  $\hat{x}[k|k]$ , the state estimate at time step  $k$  that is estimated using measurements until time  $k$ .

To simulate real-time data measurements, use the measured data one time step at a time.

```
for k = 1:size(y)
    [CorrectedState,CorrectedStateCovariance] = correct(obj,y(k));
    [PredictedState,PredictedStateCovariance] = predict(obj);
end
```

When you use the `correct` command, `obj.State` and `obj.StateCovariance` are updated with the corrected state and state estimation error covariance values for time step  $k$ , `CorrectedState` and `CorrectedStateCovariance`. When you use the `predict` command, `obj.State` and `obj.StateCovariance` are updated with the predicted values for time step  $k+1$ , `PredictedState` and `PredictedStateCovariance`.

In this example, you used `correct` before `predict` because the initial state value was  $\hat{x}[k|k-1]$ , a guess for the state value at initial time  $k$  using system outputs until time  $k-1$ . If your initial state value is  $\hat{x}[k-1|k-1]$ , the value at previous time  $k-1$  using measurement until  $k-1$ , then use the `predict` command first. For more information

about the order of using `predict` and `correct`, see “Using predict and correct Commands” on page 1-236.

### Estimate States Online using Particle Filter

Load the van der Pol ODE data, and specify the sample time.

`vdpODEdata.mat` contains a simulation of the van der Pol ODE with nonlinearity parameter  $\mu=1$ , using `ode45`, with initial conditions  $[2;0]$ . The true state was extracted with sample time  $dt = 0.05$ .

```
addpath(fullfile(matlabroot,'examples','ident','main')) % add example data
```

```
load ('vdpODEdata.mat','xTrue','dt')
tSpan = 0:dt:5;
```

Get the measurements. For this example, a sensor measures the first state with a Gaussian noise with standard deviation  $0.04$ .

```
sqrtr = 0.04;
yMeas = xTrue(:,1) + sqrtr*randn(numel(tSpan),1);
```

Create a particle filter, and set the state transition and measurement likelihood functions.

```
myPF = particleFilter(@vdpParticleFilterStateFcn,@vdpMeasurementLikelihoodFcn);
```

Initialize the particle filter at state  $[2; 0]$  with unit covariance, and use 1000 particles.

```
initialize(myPF,1000,[2;0],eye(2));
```

Pick the mean state estimation and systematic resampling methods.

```
myPF.StateEstimationMethod = 'mean';
myPF.ResamplingMethod = 'systematic';
```

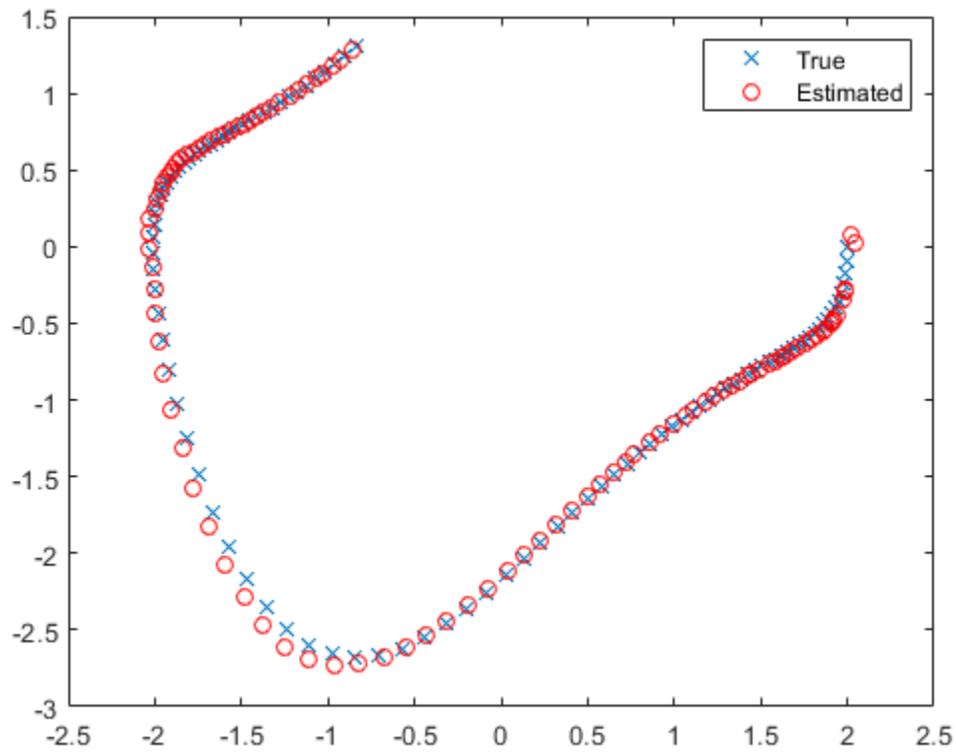
Estimate the states using the `correct` and `predict` commands, and store the estimated states.

```
xEst = zeros(size(xTrue));
for k=1:size(xTrue,1)
    xEst(k,:) = correct(myPF,yMeas(k));
```

```
    predict(myPF);  
end
```

Plot the results, and compare the estimated and true states.

```
figure(1)  
plot(xTrue(:,1),xTrue(:,2),'x',xEst(:,1),xEst(:,2),'ro')  
legend('True','Estimated')
```



```
rmpath(fullfile(matlabroot,'examples','ident','main')) % remove example data
```

### Specify State Transition and Measurement Functions with Additional Inputs

Consider a nonlinear system with input  $u$  whose state  $x$  and measurement  $y$  evolve according to the following state transition and measurement equations:

$$x[k] = \sqrt{x[k-1] + u[k-1]} + w[k-1]$$

$$y[k] = x[k] + 2 * u[k] + v[k]^2$$

The process noise  $w$  of the system is additive while the measurement noise  $v$  is nonadditive.

Create the state transition function and measurement function for the system. Specify the functions with an additional input  $u$ .

```
f = @(x,u)(sqrt(x+u));  
h = @(x,v,u)(x+2*u+v^2);
```

$f$  and  $h$  are function handles to the anonymous functions that store the state transition and measurement functions, respectively. In the measurement function, because the measurement noise is nonadditive,  $v$  is also specified as an input. Note that  $v$  is specified as an input before the additional input  $u$ .

Create an extended Kalman filter object for estimating the state of the nonlinear system using the specified functions. Specify the initial value of the state as 1, and the measurement noise as nonadditive.

```
obj = extendedKalmanFilter(f,h,1,'HasAdditiveMeasurementNoise',false);
```

Specify the measurement noise covariance.

```
obj.MeasurementNoise = 0.01;
```

You can now estimate the state of the system using the `predict` and `correct` commands. You pass the values of  $u$  to `predict` and `correct`, which in turn pass them to the state transition and measurement functions, respectively.

Correct the state estimate with measurement  $y[k]=0.8$  and input  $u[k]=0.2$  at time step  $k$ .

```
correct(obj,0.8,0.2)
```

Predict the state at next time step, given  $u[k]=0.2$ .

```
predict(obj,0.2)
```

## Input Arguments

### **obj** — Extended or unscented Kalman filter, or particle filter object

extendedKalmanFilter object | unscentedKalmanFilter object | particleFilter object

Extended or unscented Kalman filter, or particle filter object for online state estimation, created using one of the following commands:

- `extendedKalmanFilter` — Uses the extended Kalman filter algorithm.
- `unscentedKalmanFilter` — Uses the unscented Kalman filter algorithm.
- `particleFilter` — Uses the particle filter algorithm.

### **y** — Measured system output

vector

Measured system output at the current time step, specified as an  $N$ -element vector, where  $N$  is the number of measurements.

### **Um1, . . . , Umn** — Additional input arguments to measurement function

input arguments of any type

Additional input arguments to the measurement function of the system, specified as input arguments of any type. The measurement function,  $h$ , is specified in the `MeasurementFcn` or `MeasurementLikelihoodFcn` property of `obj`. If the function requires input arguments in addition to the state and measurement noise values, you specify these inputs in the `correct` command syntax. `correct` command passes these inputs to the measurement or the measurement likelihood function to calculate estimated outputs. You can specify multiple arguments.

For example, suppose that your measurement or measurement likelihood function calculates the estimated system output  $y$  using system inputs  $u$  and current time  $k$ , in addition to the state  $x$ :

$$y(k) = h(x(k), u(k), k)$$

Then when you perform online state estimation at time step  $k$ , specify these additional inputs in the `correct` command syntax:

```
[CorrectedState,CorrectedStateCovariance] = correct(obj,y,u(k),k);
```

## Output Arguments

### **CorrectedState — Corrected state estimate**

vector

Corrected state estimate, returned as a vector of size  $M$ , where  $M$  is the number of states of the system. If you specify the initial states of `obj` as a column vector then  $M$  is returned as a column vector, otherwise  $M$  is returned as a row vector.

For information about how to specify the initial states of the object, see the `extendedKalmanFilter`, `unscentedKalmanFilter` and `particleFilter` reference pages.

### **CorrectedStateCovariance — Corrected state estimation error covariance**

matrix

Corrected state estimation error covariance, returned as an  $M$ -by- $M$  matrix, where  $M$  is the number of states of the system.

## Definitions

### Using `predict` and `correct` Commands

After you have created an extended or unscented Kalman filter, or particle filter object, `obj`, to implement the estimation algorithms, use the `correct` and `predict` commands together.

At time step  $k$ , `correct` command returns the corrected value of states and state estimation error covariance using measured system outputs  $y[k]$  at the same time step. If your measurement function has additional input arguments  $U_m$ , you specify these as inputs to the `correct` command. The command passes these values to the measurement function.

```
[CorrectedState,CorrectedCovariance] = correct(obj,y,Um)
```

The `correct` command updates the `State` and `StateCovariance` properties of the object with the estimated values, `CorrectedState` and `CorrectedCovariance`.

The `predict` command returns the prediction of state and state estimation error covariance at the next time step. If your state transition function has additional input arguments  $U_s$ , you specify these as inputs to the `predict` command. The command passes these values to the state transition function.

```
[PredictedState,PredictedCovariance] = predict(obj,Us)
```

The `predict` command updates the `State` and `StateCovariance` properties of the object with the predicted values, `PredictedState` and `PredictedCovariance`.

If the current output measurement exists at a given time step, you can use `correct` and `predict`. If the measurement is missing, you can only use `predict`. For details about how these commands implement the algorithms, see “Extended and Unscented Kalman Filter Algorithms for Online State Estimation”.

The order in which you implement the commands depends on the availability of measured data  $y$ ,  $U_s$ , and  $U_m$  for your system:

- `correct` then `predict` — Assume that at time step  $k$ , the value of `obj.State` is  $\hat{x}[k|k-1]$ . This value is the state of the system at time  $k$ , estimated using measured outputs until time  $k-1$ . You also have the measured output  $y[k]$  and inputs  $U_s[k]$  and  $U_m[k]$  at the same time step.

Then you first execute the `correct` command with measured system data  $y[k]$  and additional inputs  $U_m[k]$ . The command updates the value of `obj.State` to be  $\hat{x}[k|k]$ , the state estimate for time  $k$ , estimated using measured outputs up to time  $k$ . When you then execute the `predict` command with input  $U_s[k]$ , `obj.State` now stores

$\hat{x}[k+1|k]$ . The algorithm uses this state value as an input to the `correct` command in the next time step.

- `predict` then `correct` — Assume that at time step  $k$ , the value of `obj.State` is  $\hat{x}[k-1|k-1]$ . You also have the measured output  $y[k]$  and input  $U_m[k]$  at the same time step but you have  $U_s[k-1]$  from the previous time step.

Then you first execute the `predict` command with input  $U_s[k-1]$ . The command updates the value of `obj.State` to  $\hat{x}[k|k-1]$ . When you then execute the `correct` command with input arguments  $y[k]$  and  $U_m[k]$ , `obj.State` is updated with  $\hat{x}[k|k]$ . The algorithm uses this state value as an input to the `predict` command in the next time step.

Thus, while in both cases the state estimate for time  $k$ ,  $\hat{x}[k|k]$  is the same, if at time  $k$  you do not have access to the current state transition function inputs  $U_s[k]$ , and instead have  $U_s[k-1]$ , then use `predict` first and then `correct`.

For an example of estimating states using the `predict` and `correct` commands, see “Estimate States Online Using Extended Kalman Filter” on page 1-230 or “Estimate States Online using Particle Filter” on page 1-1235.

## See Also

`clone` | `extendedKalmanFilter` | `initialize` | `particleFilter` | `predict` | `unscentedKalmanFilter`

## Topics

“Nonlinear State Estimation Using Unscented Kalman Filter and Particle Filter”

“Generate Code for Online State Estimation in MATLAB”

“What Is Online Estimation?”

“Extended and Unscented Kalman Filter Algorithms for Online State Estimation”

## Introduced in R2016b

## cra

Estimate impulse response using prewhitened-based correlation analysis

## Syntax

```
ir=cra(data)
[ir,R,cl] = cra(data,M,na,plot)
```

## Description

`ir=cra(data)` estimates the impulse response for the time-domain data, `data`.

`[ir,R,cl] = cra(data,M,na,plot)` estimates correlation/covariance information, `R`, and the 99% confidence level for the impulse response, `cl`.

The `cra` command first computes an autoregressive model for the input  $u$  as

$A(q)u(t) = e(t)$ , where  $e$  is uncorrelated (white) noise,  $q$  is the time-shift operator, and  $A(q)$  is a polynomial of order `na`. The command then filters  $u$  and output data  $y$  with  $A(q)$  to obtain the prewhitened data. The command then computes and plots the covariance functions of the prewhitened  $y$  and  $u$  and the cross-correlation function between them. Positive values of the lag variable then correspond to an influence from  $u$  to later values of  $y$ . In other words, significant correlation for negative lags is an indication of feedback from  $y$  to  $u$  in the data. A properly scaled version of this correlation function is also an estimate of the system impulse response. This is also plotted along with 99% confidence levels. The output argument `ir` is this impulse response estimate, so that its first entry corresponds to lag zero. (Negative lags are excluded in `ir`.) In the plot, the impulse response is scaled so that it corresponds to an impulse of height  $1/T$  and duration  $T$ , where  $T$  is the sample time of the data.

## Input Arguments

### **data**

Input-output data.

Specify data as an `iddata` object containing time-domain data only.

data should contain data for a single-input, single-output experiment. For the multivariate case, apply `cra` to two signals at a time, or use `impulse`.

## **M**

Number of lags for which the covariance/correlation functions are computed.

M specifies the number of lags for which the covariance/correlation functions are computed. These are from  $-M$  to  $M$ , so that the length of R is  $2M+1$ . The impulse response is computed from  $0$  to  $M$ .

**Default:** 20

## **na**

Order of the AR model to which the input is fitted.

For the prewhitening, the input is fitted to an AR model of order `na`.

Use `na = 0` to obtain the covariance and correlation functions of the original data sequences.

**Default:** 10

## **plot**

Plot display control.

Specify plot as one of the following integers:

- 0 — No plots are displayed.
- 1 — Plots the estimated impulse response with a 99% confidence region.
- 2 — Plots all the covariance functions.

**Default:** 1

## Output Arguments

### **ir**

Estimated impulse response.

The first entry of `ir` corresponds to lag zero. (Negative lags are excluded in `ir`.)

### **R**

Covariance/correlation information.

- The first column of `R` contains the lag indices.
- The second column contains the covariance function of the (possibly filtered) output.
- The third column contains the covariance function of the (possibly prewhitened) input.
- The fourth column contains the correlation function. The plots can be redisplayed by `cra(R)`.

### **cl**

99 % significance level for the impulse response.

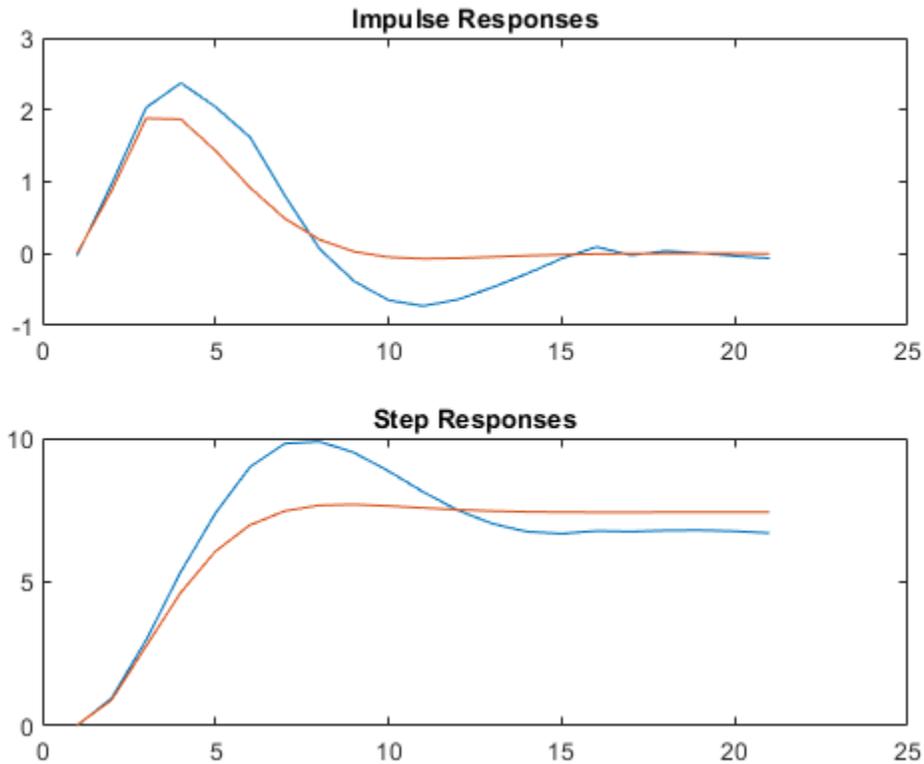
## Examples

### **Estimate the Impulse Response of an ARX Model**

Compare a second-order ARX model's impulse response with the one obtained by correlation analysis.

```
load iddata1
z = z1;
ir = cra(z);
m = arx(z,[2 2 1]);
imp = [1;zeros(20,1)];
irth = sim(m,imp);
subplot(211)
plot([ir irth])
title('Impulse Responses')
subplot(212)
```

```
plot([cumsum(ir),cumsum(irth)])  
title('Step Responses')
```



## Alternatives

An often better alternative to `cra` is `impulseeest`, which use a high-order FIR model to estimate the impulse response.

## See Also

`impulse` | `impulseeest` | `spa` | `step`

**Introduced before R2006a**

## customnet

Custom nonlinearity estimator for nonlinear ARX and Hammerstein-Wiener models

### Syntax

```
C=customnet(H)  
C=customnet(H,PropertyName,PropertyValue)
```

### Description

`customnet` is an object that stores a custom nonlinear estimator with a user-defined unit function. This custom unit function uses a weighted sum of inputs to compute a scalar output.

### Construction

`C=customnet(H)` creates a nonlinearity estimator object with a user-defined unit function using the function handle `H`. `H` must point to a function of the form `[f,g,a] = function_name(x)`, where `f` is the value of the function, `g = df/dx`, and `a` indicates the unit function active range. `g` is significantly nonzero in the interval `[-a a]`. Hammerstein-Wiener models require that your custom nonlinearity have only one input and one output.

`C=customnet(H,PropertyName,PropertyValue)` creates a nonlinearity estimator using property-value pairs defined in “`customnet Properties`” on page 1-244.

### customnet Properties

You can include property-value pairs in the constructor to specify the object.

After creating the object, you can use `get` or dot notation to access the object property values. For example:

```
% List all property values
get(C)
% Get value of NumberOfUnits property
C.NumberOfUnits
```

You can also use the `set` function to set the value of particular properties. For example:

```
set(C, 'LinearTerm', 'on')
```

The first argument to `set` must be the name of a MATLAB variable.

Property Name	Description
NumberOfUnits	<p>Integer specifies the number of nonlinearity units in the expansion. Default=10.</p> <p>For example:</p> <pre>customnet(H, 'NumberOfUnits', 5)</pre>
LinearTerm	<p>Can have the following values:</p> <ul style="list-style-type: none"> <li>'on'—Estimates the vector <math>L</math> in the expansion.</li> <li>'off'—Fixes the vector <math>L</math> to zero.</li> </ul> <p>For example:</p> <pre>customnet(H, 'LinearTerm', 'on')</pre>

Property Name	Description
Parameters	<p>A structure containing the parameters in the nonlinear expansion, as follows:</p> <ul style="list-style-type: none"> <li>• <b>RegressorMean</b>: 1-by-m vector containing the means of <math>x</math> in estimation data, <math>r</math>.</li> <li>• <b>NonLinearSubspace</b>: m-by-q matrix containing <math>Q</math>.</li> <li>• <b>LinearSubspace</b>: m-by-p matrix containing <math>P</math>.</li> <li>• <b>LinearCoef</b>: p-by-1 vector <math>L</math>.</li> <li>• <b>Dilation</b>: q-by-1 matrix containing the values <math>b_n</math>.</li> <li>• <b>Translation</b>: 1-by-n vector containing the values <math>c_n</math>.</li> <li>• <b>OutputCoef</b>: n-by-1 vector containing the values <math>a_n</math>.</li> <li>• <b>OutputOffset</b>: scalar <math>d</math>.</li> </ul> <p>Typically, the values of this structure are set by estimating a model with a customnet nonlinearity.</p>
UnitFcn	Stores the function handle that points to the unit function.

## Examples

Define custom unit function and save it in `gaussunit.m`:

```
function [f, g, a] = GAUSSUNIT(x)
% x: unit function variable
% f: unit function value
% g: df/dx
% a: unit active range (g(x) is significantly
% nonzero in the interval [-a a])

% The unit function must be "vectorized": for
% a vector or matrix x, the output arguments f and g
% must have the same size as x,
% computed element-by-element.

% GAUSSUNIT customnet unit function example
[f, g, a] = gaussunit(x)
f = exp(-x.*x);
if nargin>1
```

```

    g = - 2*x.*f;
    a = 0.2;
end

```

Use custom networks in `nlarx` and `nlhw` model estimation commands:

```

% Define handle to example unit function.
H = @gaussunit;
% Estimate nonlinear ARX model using
% Gauss unit function with 5 units.
m = nlarx(Data,Orders,customnet(H,'NumberOfUnits',5));

```

## Tips

Use `customnet` to define a nonlinear function  $y = F(x)$ , where  $y$  is scalar and  $x$  is an  $m$ -dimensional row vector. The unit function is based on the following function expansion with a possible linear term  $L$ :

$$F(x) = (x-r)PL + a_1 f((x-r)Qb_1 + c_1) + \dots \\ + a_n f((x-r)Qb_n + c_n) + d$$

where  $f$  is a unit function that you define using the function handle  $H$ .

$P$  and  $Q$  are  $m$ -by- $p$  and  $m$ -by- $q$  projection matrices, respectively. The projection matrices  $P$  and  $Q$  are determined by principal component analysis of estimation data. Usually,  $p=m$ . If the components of  $x$  in the estimation data are linearly dependent, then  $p < m$ . The number of columns of  $Q$ ,  $q$ , corresponds to the number of components of  $x$  used in the unit function.

When used to estimate nonlinear ARX models,  $q$  is equal to the size of the `NonlinearRegressors` property of the `idnlarx` object. When used to estimate Hammerstein-Wiener models,  $m=q=1$  and  $Q$  is a scalar.

$r$  is a  $1$ -by- $m$  vector and represents the mean value of the regressor vector computed from estimation data.

$d$ ,  $a$ , and  $c$  are scalars.

$L$  is a  $p$ -by- $1$  vector.

$b$  represents  $q$ -by-1 vectors.

The function handle of the unit function of the custom net must have the form `[f, g, a] = function_name(x)`. This function must be vectorized, which means that for a vector or matrix  $x$ , the output arguments  $f$  and  $g$  must have the same size as  $x$  and be computed element-by-element.

## Algorithms

`customnet` uses an iterative search technique for estimating parameters.

## See Also

`evaluate` | `nlarx` | `nlnhw`

## Topics

“Identifying Nonlinear ARX Models”

“Identifying Hammerstein-Wiener Models”

**Introduced in R2007a**

## customreg

Custom regressor for nonlinear ARX models

### Syntax

```
C=customreg(Function,Variables)
C=customreg(Function,Variables,Delays,Vectorized)
```

### Description

`customreg` class represents arbitrary functions of past inputs and outputs, such as products, powers, and other MATLAB expressions of input and output variables.

You can specify custom regressors in addition to or instead of standard regressors for greater flexibility in modeling your data using nonlinear ARX models. For example, you can define regressors like  $\tan(u(t-1))$ ,  $u(t-1)^2$ , and  $u(t-1)*y(t-3)$ .

For simpler regressor expressions, specify custom regressors directly in the app or in the `nlarx` estimation command. For more complex expressions, create a `customreg` object for each custom regressor and specify these objects as inputs to the estimation. Regardless of how you specify custom regressors, the toolbox represents these regressors as `customreg` objects. Use `getreg` to list the expressions of all standard and custom regressors in your model.

A special case of custom regressors involves polynomial combinations of past inputs and outputs. For example, it is common to capture nonlinearities in the system using polynomial expressions like  $y(t-1)^2$ ,  $u(t-1)^2$ ,  $y(t-2)^2$ ,  $y(t-1)*y(t-2)$ ,  $y(t-1)*u(t-1)$ ,  $y(t-2)*u(t-1)$ . At the command line, use the `polyreg` command to generate polynomial-type regressors automatically by computing all combinations of input and output variables up to a specified degree. `polyreg` produces `customreg` objects that you specify as inputs to the estimation.

The nonlinear ARX model (`idnlarx` object) stores all custom regressors as the `CustomRegressors` property. You can list all custom regressors using `m.CustomRegressors`, where `m` is a nonlinear ARX model. For MIMO models, to retrieve the `r`th custom regressor for output `ky`, use `m.CustomRegressors{ky}(r)`.

Use the `Vectorized` property to specify whether to compute custom regressors using vectorized form during estimation. If you know that your regressor formulas can be vectorized, set `Vectorized` to 1 to achieve better performance. To better understand vectorization, consider the custom regressor function handle `z=@(x,y)x^2*y`. `x` and `y` are vectors and each variable is evaluated over a time grid. Therefore, `z` must be evaluated for each `(xi,yi)` pair, and the results are concatenated to produce a `z` vector:

```
for k = 1:length(x)
    z(k) = x(k)^2*y(k)
end
```

The above expression is a nonvectorized computation and tends to be slow. Specifying a `Vectorized` computation uses MATLAB vectorization rules to evaluate the regressor expression using matrices instead of the FOR-loop and results in faster computation:

```
% "." indicates element-wise operation
z=(x.^2).*y
```

## Construction

`C=customreg(Function,Variables)` specifies a custom regressor for a nonlinear ARX model. `C` is a `customreg` object that stores custom regressor. `Function` is a function of input and output variables. `Variables` represent the names of model inputs and outputs in the function `Function`. Each input and output name must coincide with the `InputName` and `OutputName` properties of the corresponding `idnlarx` object. The size of `Variables` must match the number of `Function` inputs. For multiple-output models with `p` outputs, the custom regressor is a `p`-by-1 cell array or an array of `customreg` object, where the `kyth` entry defines the custom regressor for output `ky`. You must add these regressors to the `model` by assigning the `CustomRegressors model` property or by using `addreg`.

`C=customreg(Function,Variables,Delays,Vectorized)` create a custom regressor that includes the delays corresponding to inputs or outputs in `Arguments`. `Delays` is a vector of positive integers that represent the delays of `Variables` variables (default is 1 for each vector element). The size of `Delays` must match the size of `Variables`. `Vectorized` value of 1 uses MATLAB vectorization rules to evaluate the regressor expression `Function`. By default, `Vectorized` value is 0 (false).

## Properties

After creating the object, you can use `get` or dot notation to access the object property values. For example:

```
% List all property values
get(C)
% Get value of Arguments property
C.Arguments
```

You can also use the `set` function to set the value of particular properties. For example:

```
set(C, 'Vectorized', 1)
```

Property Name	Description
Function	<p>Function handle or character vector representing a function of standards regressors.</p> <p>For example:</p> <pre>cr = @(x,y) x*y</pre>
Variables	<p>Cell array of character vectors that represent the names of model input and output variables in the function <code>Function</code>. Each input and output name must coincide with the <code>InputName</code> and <code>OutputName</code> properties of the <code>idnlarx</code> object—the model for which you define custom regressors. The size of <code>Variables</code> must match the number of <code>Function</code> inputs.</p> <p>For example, <code>Variables</code> correspond to <code>{'y1', 'u1'}</code> in:</p> <pre>C = customreg(cr,{'y1', 'u1'},[2 3])</pre>
Delays	<p>Vector of positive integers representing the delays of <code>Variables</code>. The size of <code>Delays</code> must match the size of <code>Arguments</code>.</p> <p>Default: 1 for each vector element.</p> <p>For example, <code>Delays</code> are <code>[2 3]</code> in:</p> <pre>C = customreg(cr,{'y1', 'u1'},[2 3])</pre>

Property Name	Description
Vectorized	Assignable values: <ul style="list-style-type: none"><li>• 0 (default)—Function is not computed in vectorized form.</li><li>• 1—Function is computed in vectorized form when called with vector arguments.</li></ul>

## Examples

### Define Custom Regressors

Load estimation data.

```
load iddata1
```

Specify the regressors as a cell array of character vectors.

```
C = {'u1(t-1)*sin(y1(t-3))', 'u1(t-2)^3'};
```

u1 and y1 are input and output data, respectively.

Estimate a nonlinear ARX model using the custom regressors.

```
m = nlarx(z1,[2 2 1], 'linear', 'CustomRegressors', C);
```

### Define Custom Regressors During Estimation

Load the estimation data.

```
load iddata1
```

Estimate a nonlinear ARX model with custom regressors.

```
m = nlarx(z1,[2 2 1], 'linear', 'CustomRegressors', ...  
        {'u1(t-1)*sin(y1(t-3))', 'u1(t-2)^3'});
```

## Define Custom Regressors as Array of customreg Objects

Load the estimation data.

```
load iddata1
```

Construct a nonlinear ARX model.

```
m = idnlarx([2 2 1]);
```

Define the custom regressors.

```
cr1 = @(x,y) x*sin(y);  
cr2 = @(x) x^3;  
C = [customreg(cr1,{'u','y'},[1 3]),customreg(cr2,{'u'},2)];
```

Add custom regressors to the model.

```
m2 = addreg(m,C);
```

## Use Vectorization Rules to Evaluate Regressor Expression

Load the estimation data.

```
load iddata1
```

Specify the regressors.

```
C = customreg(@(x,y) x*sin(y),{'u' 'y'},[1 3]);  
set(C,'Vectorized',1);
```

Estimate a nonlinear ARX model with custom regressors.

```
m = nlarx(z1,[2 2 1],'sigmoidnet','CustomReg',C);
```

## See Also

[addreg](#) | [getreg](#) | [idnlarx](#) | [nlarx](#) | [polyreg](#)

**Topics**

“Identifying Nonlinear ARX Models”

**Introduced in R2007a**

## d2c

Convert model from discrete to continuous time

### Syntax

```
sysc = d2c(sysd)
sysc = d2c(sysd,method)
sysc = d2c(sysd,opts)
[sysc,G] = d2c(sysd,method,opts)
```

### Description

`sysc = d2c(sysd)` produces a continuous-time model `sysc` that is equivalent to the discrete-time dynamic system model `sysd` using zero-order hold on the inputs.

`sysc = d2c(sysd,method)` uses the specified conversion method `method`.

`sysc = d2c(sysd,opts)` converts `sysd` using the option set `opts`, specified using the `d2cOptions` command.

`[sysc,G] = d2c(sysd,method,opts)` returns a matrix `G` that maps the states `xd[k]` of the state-space model `sysd` to the states `xc(t)` of `sysc`.

### Input Arguments

#### **sysd**

Discrete-time dynamic system model

You cannot directly use an `idgrey` model whose `FunctionType` is 'd' with `d2c`. Convert the model into `idss` form first.

#### **Default:**

## method

Discrete-to-continuous time conversion method, specified as one of the following values:

- 'zoh' — Zero-order hold on the inputs. Assumes that the control inputs are piecewise constant over the sampling period.
- 'foh' — Linear interpolation of the inputs (modified first-order hold). Assumes that the control inputs are piecewise linear over the sampling period.
- 'tustin' — Bilinear (Tustin) approximation to the derivative
- 'matched' — Zero-pole matching method (for SISO systems only). See [1].

For information about the algorithms for each d2c conversion method, see “Continuous-Discrete Conversion Methods”.

**Default:** 'zoh'

## opts

Discrete-to-continuous time conversion options, created using `d2cOptions`.

## Output Arguments

### sysc

Continuous-time model of the same type as the input system `sysd`.

When `sysd` is an identified (IDLTI) model, `sysc`:

- Includes both the measured and noise components of `sysd`. If the noise variance is  $\lambda$  in `sysd`, then the continuous-time model `sysc` has an indicated level of noise spectral density equal to  $T_s \cdot \lambda$ .
- Does not include the estimated parameter covariance of `sysd`. If you want to translate the covariance while converting the model, use `translatecov`.

### G

Matrix mapping the states `xd[k]` of the state-space model `sysd` to the states `xc(t)` of `sysc`:

$$x_c(kT_s) = G \begin{bmatrix} x_d[k] \\ u[k] \end{bmatrix}.$$

Given an initial condition  $x_0$  for `sysd` and an initial input  $u_0 = u[0]$ , the corresponding initial condition for `sysc` (assuming  $u[k] = 0$  for  $k < 0$ ) is:

$$x_c(0) = G \begin{bmatrix} x_0 \\ u_0 \end{bmatrix}.$$

## Examples

### Convert Discrete-Time Transfer Function to Continuous Time

Create the following discrete-time transfer function:

$$H(z) = \frac{z - 1}{z^2 + z + 0.3}$$

```
H = tf([1 -1],[1 1 0.3],0.1);
```

The sample time of the model is  $T_s = 0.1s$ .

Derive a continuous-time, zero-order-hold equivalent model.

```
Hc = d2c(H)
```

```
Hc =
```

$$\frac{121.7 s + 1.405e-12}{s^2 + 12.04 s + 776.7}$$

Continuous-time transfer function.

Discretize the resulting model, `Hc`, with the default zero-order hold method and sample time 0.1s to return the original discrete model, `H`.

```
c2d(Hc,0.1)
```

```
ans =
```

$$\frac{z - 1}{z^2 + z + 0.3}$$

```
Sample time: 0.1 seconds  
Discrete-time transfer function.
```

Use the Tustin approximation method to convert H to a continuous time model.

```
Hc2 = d2c(H, 'tustin');
```

Discretize the resulting model, Hc2, to get back the original discrete-time model, H.

```
c2d(Hc2,0.1, 'tustin');
```

## Convert Identified Discrete-Time Transfer Function to Continuous Time

Estimate a discrete-time transfer function model, and convert it to a continuous-time model.

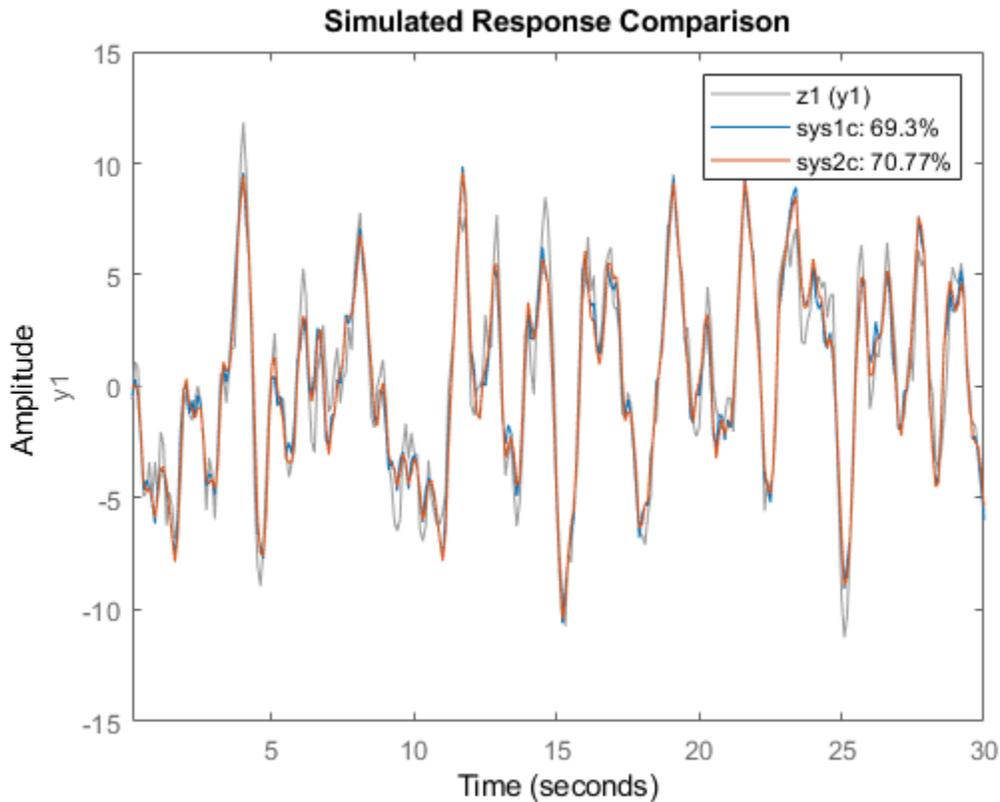
```
load iddata1  
sys1d = tfest(z1,2,'Ts',0.1);  
sys1c = d2c(sys1d,'zoh');
```

Estimate a continuous-time transfer function model.

```
sys2c = tfest(z1,2);
```

Compare the response of sys1c and the directly estimated continuous-time model, sys2c.

```
compare(z1,sys1c,sys2c)
```



The two systems are almost identical.

### Regenerate Covariance Information After Converting to Continuous-Time Model

Convert an identified discrete-time transfer function model to continuous-time.

```
load iddata1
sysd = tfest(z1,2,'Ts',0.1);
sysc = d2c(sysd,'zoh');
```

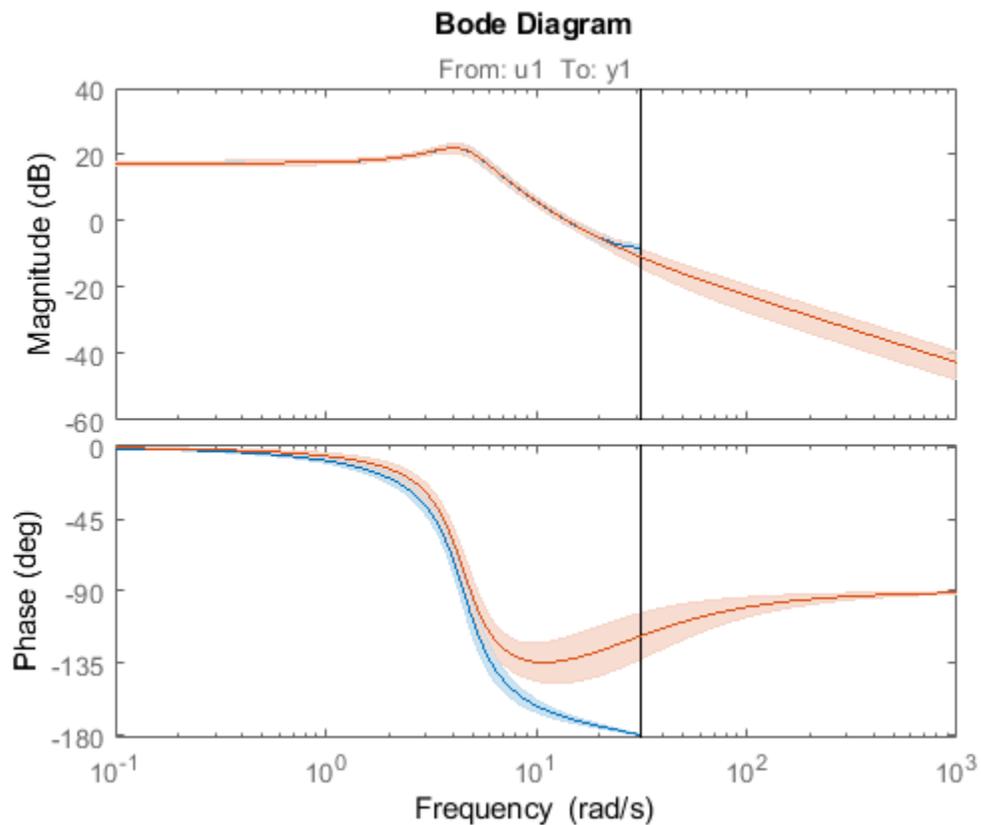
sys1c has no covariance information. The d2c operation leads to loss of covariance data of identified models.

Regenerate the covariance information using a zero iteration update with the same estimation command and estimation data.

```
opt = tfestOptions;  
opt.SearchOptions.MaxIterations = 0;  
sys1c = tfest(z1,sysc,opt);
```

Analyze the effect on frequency-response uncertainty.

```
h = bodeplot(sysd,sys1c);  
showConfidence(h,3)
```



The uncertainties of `sys1c` and `sysd` are comparable up to the Nyquist frequency. However, `sys1c` exhibits large uncertainty in the frequency range for which the estimation data does not provide any information.

If you do not have access to the estimation data, use the `translatecov` command which is a Gauss-approximation formula based translation of covariance across model type conversion operations.

## Limitations

The Tustin approximation is not defined for systems with poles at  $z = -1$  and is ill-conditioned for systems with poles near  $z = -1$ .

The zero-order hold method cannot handle systems with poles at  $z = 0$ . In addition, the 'zoh' conversion increases the model order for systems with negative real poles, [2]. The model order increases because the matrix logarithm maps real negative poles to complex poles. Single complex poles are not physically meaningful because of their complex time response.

Instead, to ensure that all complex poles of the continuous model come in conjugate pairs, `d2c` replaces negative real poles  $z = -\alpha$  with a pair of complex conjugate poles near  $-\alpha$ . The conversion then yields a continuous model with higher order. For example, to convert the discrete-time transfer function

$$H(z) = \frac{z + 0.2}{(z + 0.5)(z^2 + z + 0.4)}$$

type:

```
Ts = 0.1 % sample time 0.1 s
H = zpk(-0.2, -0.5, 1, Ts) * tf(1, [1 1 0.4], Ts)
Hc = d2c(H)
```

These commands produce the following result.

Warning: System order was increased to handle real negative poles.

```
Zero/pole/gain:
-33.6556 (s-6.273) (s^2 + 28.29s + 1041)
-----
(s^2 + 9.163s + 637.3) (s^2 + 13.86s + 1035)
```

To convert  $H_c$  back to discrete time, type:

```
c2d(Hc, Ts)
```

yielding

```
Zero/pole/gain:  
      (z+0.5) (z+0.2)  
-----  
(z+0.5)^2 (z^2 + z + 0.4)
```

```
Sample time: 0.1
```

This discrete model coincides with  $H(z)$  after canceling the pole/zero pair at  $z = -0.5$ .

## Tips

- Use the syntax `sysc = d2c(sysd, 'method')` to convert `sysd` using the default options for 'method'. To specify `tustin` conversion with a frequency prewarp (formerly the 'prewarp' method), use the syntax `sysc = d2c(sysd, opts)`. See the `d2cOptions` reference page for more information.

## Algorithms

`d2c` performs the 'zoh' conversion in state space, and relies on the matrix logarithm (see `logm` in the MATLAB documentation).

See "Continuous-Discrete Conversion Methods" (Control System Toolbox) for more details on the conversion methods.

## References

- [1] Franklin, G.F., Powell, D.J., and Workman, M.L., *Digital Control of Dynamic Systems* (3rd Edition), Prentice Hall, 1997.
- [2] Kollár, I., G.F. Franklin, and R. Pintelon, "On the Equivalence of z-domain and s-domain Models in System Identification," *Proceedings of the IEEE® Instrumentation and*

*Measurement Technology Conference*, Brussels, Belgium, June, 1996, Vol. 1, pp. 14-19.

## **See Also**

c2d | d2cOptions | d2d | logm | translatecov

**Introduced before R2006a**

## d2cOptions

Create option set for discrete- to continuous-time conversions

### Syntax

```
opts = d2cOptions  
opts = d2cOptions(Name,Value)
```

### Description

`opts = d2cOptions` returns the default options for `d2c`.

`opts = d2cOptions(Name,Value)` creates an option set with the options specified by one or more `Name, Value` pair arguments.

### Input Arguments

#### Name-Value Pair Arguments

##### method

Discretization method, specified as one of the following values:

'zoh'	Zero-order hold, where <code>d2c</code> assumes that the control inputs are piecewise constant over the sample time <code>Ts</code> .
'foh'	Linear interpolation of the inputs (modified first-order hold). Assumes that the control inputs are piecewise linear over the sampling period.
'tustin'	Bilinear (Tustin) approximation. By default, <code>d2c</code> converts with no prewarp. To include prewarp, use the <code>PrewarpFrequency</code> option.
'matched'	Zero-pole matching method. (See [1] on page 1-266, p. 224.)

For information about the algorithms for each d2c conversion method, see “Continuous-Discrete Conversion Methods”.

**Default:** 'zoh'

### PrewarpFrequency

Prewarp frequency for 'tustin' method, specified in rad/TimeUnit, where TimeUnit is the time units, specified in the TimeUnit property, of the discrete-time system. Specify the prewarp frequency as a positive scalar value. A value of 0 corresponds to the 'tustin' method without prewarp.

**Default:** 0

## Output Arguments

### opts — Option set for d2c

d2cOptions option set

Option set for d2c, returned as an d2cOptions option set.

## Examples

### Specify Model Discretization Method

Consider the following discrete-time transfer function.

$$H(z) = \frac{z + 1}{z^2 + z + 1}$$

Create the discrete-time transfer function with a sample time of 0.1 seconds.

```
Hd = tf([1 1],[1 1 1],0.1);
```

Specify the discretization method as bilinear (Tustin) approximation and the prewarp frequency as 20 rad/seconds.

```
opts = d2cOptions('Method','tustin','PrewarpFrequency',20);
```

Convert the discrete-time model to continuous-time using the specified discretization method.

```
Hc = d2c(Hd,opts);
```

You can use the discretization option set `opts` to discretize additional models using the same options.

## References

[1] Franklin, G.F., Powell,D.J., and Workman, M.L., *Digital Control of Dynamic Systems* (3rd Edition), Prentice Hall, 1997.

## See Also

d2c

**Introduced in R2012a**

## d2d

Resample discrete-time model

### Syntax

```
sys1 = d2d(sys, Ts)
sys1 = d2d(sys, Ts, 'method')
sys1 = d2d(sys, Ts, opts)
```

### Description

`sys1 = d2d(sys, Ts)` resamples the discrete-time dynamic system model `sys` to produce an equivalent discrete-time model `sys1` with the new sample time `Ts` (in seconds), using zero-order hold on the inputs.

`sys1 = d2d(sys, Ts, 'method')` uses the specified resampling method 'method':

- 'zoh' — Zero-order hold on the inputs
- 'tustin' — Bilinear (Tustin) approximation

For information about the algorithms for each `d2d` conversion method, see “Continuous-Discrete Conversion Methods”.

`sys1 = d2d(sys, Ts, opts)` resamples `sys` using the option set with `d2dOptions`.

### Examples

#### Resample a Discrete-Time Model

Create the following zero-pole-gain-model with sample time 0.1 seconds.

$$H(z) = \frac{z - 0.7}{z - 0.5}$$

```
H = zpk(0.7,0.5,1,0.1);
```

Resample the model at 0.05 s.

```
H2 = d2d(H,0.05)
```

```
H2 =
```

```
(z-0.8243)
-----
(z-0.7071)
```

```
Sample time: 0.05 seconds
Discrete-time zero/pole/gain model.
```

Resample H2 at 0.1 seconds to obtain the original model H.

```
H3 = d2d(H2,0.1)
```

```
H3 =
```

```
(z-0.7)
-----
(z-0.5)
```

```
Sample time: 0.1 seconds
Discrete-time zero/pole/gain model.
```

## Resample an Identified Discrete-Time Model

Suppose that you estimate a discrete-time output-error polynomial model with sample time commensurate with the estimation data (0.1 seconds). However, your deployment application requires a faster sampling frequency (0.01 seconds). You can use `d2d` to resample your estimated model.

Load the estimation data.

```
load iddata1 z1
z1.Ts
```

```
ans = 0.1000
```

`z1` is an `iddata` object containing the estimation input-output data with sample time 0.1 seconds.

Estimate an output-error polynomial model of order `[2 2 1]`.

```
sys = oe(z1,[2 2 1]);  
sys.Ts
```

```
ans = 0.1000
```

Resample the model at sample time 0.01 seconds.

```
sys2 = d2d(sys,0.01);  
sys2.Ts
```

```
ans = 0.0100
```

`d2d` resamples the model using the zero-order hold method.

## Tips

- Use the syntax `sys1 = d2d(sys,Ts,'method')` to resample `sys` using the default options for `'method'`. To specify `tustin` resampling with a frequency prewarp, use the syntax `sys1 = d2d(sys,Ts,opts)`. For more information, see `d2dOptions`.
- When `sys` is an identified (IDLTI) model, `sys1` does not include the estimated parameter covariance of `sys`. If you want to translate the covariance while converting the model, use `translatecov`.

## See Also

`c2d` | `d2c` | `d2dOptions` | `translatecov` | `upsample`

**Introduced before R2006a**

## d2dOptions

Create option set for discrete-time resampling

### Syntax

```
opts = d2dOptions  
opts = d2dOptions('OptionName', OptionValue)
```

### Description

*opts* = d2dOptions returns the default options for d2d.

*opts* = d2dOptions('OptionName', OptionValue) accepts one or more comma-separated name-value pairs that specify options for the d2d command. Specify *OptionName* inside single quotes.

This table summarizes the options that the d2d command supports.

## Input Arguments

### Name-Value Pair Arguments

#### Method

Discretization method, specified as one of the following values:

'zoh'	Zero-order hold, where d2d assumes that the control inputs are piecewise constant over the sample time $T_s$ .
'tustin'	Bilinear (Tustin) approximation. By default, d2d resamples with no prewarp. To include prewarp, use the PrewarpFrequency option.

For information about the algorithms for each d2d conversion method, see “Continuous-Discrete Conversion Methods”.

**Default:** 'zoh'

### PrewarpFrequency

Prewarp frequency for 'tustin' method, specified in rad/TimeUnit, where TimeUnit is the time units, specified in the TimeUnit property, of the resampled system. Takes positive scalar values. The prewarp frequency must be smaller than the Nyquist frequency before and after resampling. A value of 0 corresponds to the standard 'tustin' method without prewarp.

**Default:** 0

## Examples

### Specify Method for Resampling a Discrete-Time Model

Create the following discrete-time transfer function with sample time 0.1 seconds.

$$H(z) = \frac{z + 1}{z^2 + z + 1}$$

```
h1 = tf([1 1],[1 1 1],0.1);
```

Specify the discretization method as bilinear Tustin method with a prewarping frequency of 20 rad/seconds.

```
opts = d2dOptions('Method','tustin','PrewarpFrequency',20);
```

Resample the discrete-time model using the specified options.

```
h2 = d2d(h1,0.05,opts);
```

You can use the option set `opts` to resample additional models using the same options.

## See Also

d2d

**Introduced in R2012a**

# damp

Natural frequency and damping ratio

## Syntax

```
damp(sys)
[Wn,zeta] = damp(sys)
[Wn,zeta,P] = damp(sys)
```

## Description

`damp(sys)` displays a table of the damping ratio (also called damping factor), natural frequency, and time constant of the poles of the linear model `sys`. For a discrete-time model, the table also includes the magnitude of each pole. Frequencies are expressed in units of the reciprocal of the `TimeUnit` property of `sys`. Time constants are expressed in the same units as the `TimeUnit` property of `sys`.

This command requires a Control System Toolbox license.

`[Wn,zeta] = damp(sys)` returns the natural frequencies, `Wn`, and damping ratios, `zeta`, of the poles of `sys`.

`[Wn,zeta,P] = damp(sys)` returns the poles of `sys`.

## Input Arguments

### `sys`

Any linear dynamic system model.

## Output Arguments

### **Wn**

Vector containing the natural frequencies of each pole of `sys`, in order of increasing frequency. Frequencies are expressed in units of the reciprocal of the `TimeUnit` property of `sys`.

If `sys` is a discrete-time model with specified sample time, `Wn` contains the natural frequencies of the equivalent continuous-time poles (see “Algorithms” on page 1-276). If `sys` has an unspecified sample time (`Ts = -1`), then the software uses `Ts = 1` and calculates `Wn` accordingly.

### **zeta**

Vector containing the damping ratios of each pole of `sys`, in the same order as `Wn`.

If `sys` is a discrete-time model with specified sample time, `zeta` contains the damping ratios of the equivalent continuous-time poles (see “Algorithms” on page 1-276). If `sys` has an unspecified sample time (`Ts = -1`), then the software uses `Ts = 1` and calculates `zeta` accordingly.

### **P**

Vector containing the poles of `sys`, in order of increasing natural frequency. `P` is the same as the output of `pole(sys)`, except for the order.

## Examples

### **Display Natural Frequency, Damping Ratio, and Poles of Continuous-Time System**

Create the following continuous-time transfer function:

$$H(s) = \frac{2s^2 + 5s + 1}{s^2 + 2s + 3}.$$

```
H = tf([2 5 1],[1 2 3]);
```

Display the natural frequencies, damping ratios, time constants, and poles of `H`.

```
damp(H)
```

Pole	Damping	Frequency (rad/seconds)	Time Constant (seconds)
-1.00e+00 + 1.41e+00i	5.77e-01	1.73e+00	1.00e+00
-1.00e+00 - 1.41e+00i	5.77e-01	1.73e+00	1.00e+00

Obtain vectors containing the natural frequencies and damping ratios of the poles.

```
[Wn,zeta] = damp(H);
```

Calculate the associated time constants.

```
tau = 1./(zeta.*Wn);
```

### Display Natural Frequency, Damping Ratio, and Poles of Discrete-Time System

Create a discrete-time transfer function.

```
H = tf([5 3 1],[1 6 4 4],0.01);
```

Display information about the poles of  $H$ .

```
damp(H)
```

Pole	Magnitude	Damping	Frequency (rad/seconds)	Time Constant (seconds)
-3.02e-01 + 8.06e-01i	8.61e-01	7.74e-02	1.93e+02	6.68e-02
-3.02e-01 - 8.06e-01i	8.61e-01	7.74e-02	1.93e+02	6.68e-02
-5.40e+00	5.40e+00	-4.73e-01	3.57e+02	-5.93e-03

The **Magnitude** column displays the discrete-time pole magnitudes. The **Damping**, **Frequency**, and **Time Constant** columns display values calculated using the equivalent continuous-time poles.

Obtain vectors containing the natural frequencies and damping ratios of the poles.

```
[Wn,zeta] = damp(H);
```

Calculate the associated time constants.

$$\tau = 1./(\zeta \cdot \omega_n);$$

## Algorithms

The natural frequency, time constant, and damping ratio of the system poles are defined in the following table:

	Continuous Time	Discrete Time with Sample Time Ts
<b>Pole Location</b>	$s$	$z$
<b>Equivalent Continuous-Time Pole</b>	Not applicable	$s = \frac{\ln(z)}{T_s}$
<b>Natural Frequency</b>	$\omega_n =  s $	$\omega_n =  s  = \left  \frac{\ln(z)}{T_s} \right $
<b>Damping Ratio</b>	$\zeta = -\cos(\angle s)$	$\zeta = -\cos(\angle s) = -\cos(\angle \ln(z))$
<b>Time Constant</b>	$\tau = \frac{1}{\omega_n \zeta}$	$\tau = \frac{1}{\omega_n \zeta}$

## See Also

dsort | eig | esort | pole | pzmap | zero

**Introduced before R2006a**

## data2state

Map past data to states of state-space and nonlinear ARX models

### Syntax

```
X = data2state(sys,PastData)
[X,XCov] = data2state(sys,PastData)
```

### Description

`X = data2state(sys,PastData)` maps the past data to the states of a state-space or a nonlinear ARX model `sys`. `X` contains the state values at the time instant immediately after the most recent data sample in `PastData`. The software computes the state estimates by minimizing the 1-step ahead prediction error between predicted response and output signal in `PastData`.

`data2state` is useful for continued model simulation. That is, suppose you have simulated a model up to a certain time instant and would like to then simulate the model for future inputs. Use `data2state` to estimate states of the model at the beginning of the second simulation.

`[X,XCov] = data2state(sys,PastData)` returns the estimated covariance, `XCov`, of the current states.

### Examples

#### Compute Mapped States of Identified Model

Compute the mapped states of an identified model, and use the states as initial state values for model simulation.

Load estimation data.

```
load iddata3 z3
```

Estimate a second-order state-space model using the data.

```
sys = ssest(z3,2);
```

Simulate the model using the entire input signal in z3.

```
Input = z3(:,[],:); % |iddata| object containing only the input signal  
y_all = sim(sys,Input);
```

sim uses zero initial conditions to compute y\_all.

Now simulate the model using only the first-half of the input signal.

```
Input1 = Input(1:150);  
y_1 = sim(sys,Input1);
```

Continue the simulation with the second-half of the input signal such that the results show no discontinuity owing to initial-condition-induced transients. To do so, first construct a past data set comprising of the input and simulated output from the first-half of the input signal. Then calculate the state values corresponding to the start of the second-half of the input signal ( $t = 151$ ).

```
PastData = [y_1,Input1];  
X = data2state(sys,PastData);
```

X contains the state values at the time instant immediately after the most recent data sample in PastData. This time point is also the start of the future data (second-half of the input signal).

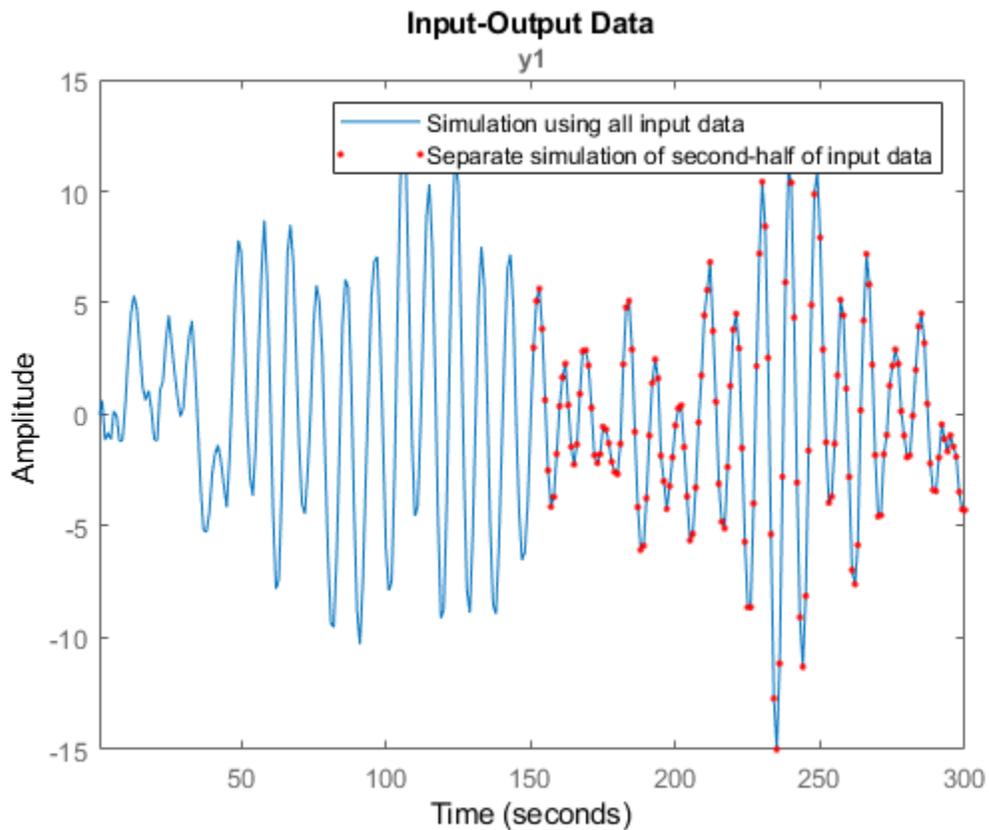
```
FutureData = Input(151:end);
```

Simulate the model using the second-half of the input signal and X as initial conditions.

```
opt = simOptions('InitialCondition',X);  
y_2 = sim(sys,FutureData,opt);
```

Verify that y\_2 matches the second half of y\_all.

```
plot(y_all,y_2,'r.')  
legend('Simulation using all input data',...  
       'Separate simulation of second-half of input data')
```



### Calculate Mapped States and Covariance of States

Load the past data.

```
load iddata1 z1  
PastData = z1;
```

Estimate an ARX model.

```
sys = arx(PastData,[1 1 0]);
```

Convert the model to a state-space model.

```
sys2 = idss(sys);
```

Calculate the mapped states and covariance of states using `PastData`.

```
[X,XCov] = data2state(sys2,PastData);
```

`X` is the state value at the time instant immediately after the most recent data sample in `PastData`.

### Determine Mapped State of a Nonlinear ARX model

Load your data and create a data object.

```
load motorizedcamera;  
z = iddata(y,u,0.02,'Name','Motorized Camera','TimeUnit','s');
```

Estimate a nonlinear ARX model.

```
mw1 = nlarx(z,[ones(2,2),ones(2,6),ones(2,6)],'wavenet');
```

The estimated model has six inputs and two outputs.

Determine the model order, `nx`.

```
nx = order(mw1);
```

Use the first `nx` samples of data to generate initial conditions.

```
PastData = struct('Input', z.u(1:nx,:), 'Output', z.y(1:nx,:));
```

Compute the mapped states of the model.

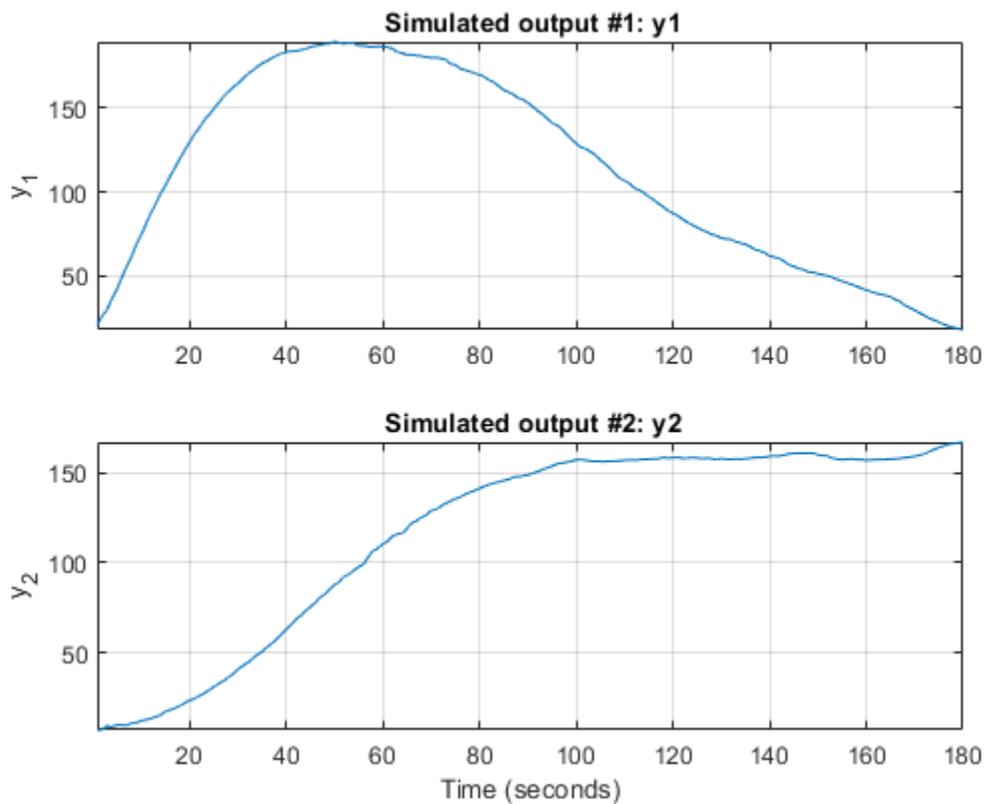
```
X = data2state(mw1,PastData);
```

`X` is the state value at the time instant immediately after the most recent data sample in `PastData`.

Simulate the model using the remaining input data, and specify the initial conditions for simulation.

```
InputSignal = z.u(nx+1:end,:);  
opt = simOptions;
```

```
opt.InitialCondition = X;  
sim(mw1,InputSignal,opt)
```



## Input Arguments

### **sys** — Identified model

idss | idgrey | idnlgrey | idnlrx

Identified model whose current states are estimated, specified as one of the following:

- State-space model (idss, idgrey, or idnlgrey object)

- Nonlinear ARX model (`idnlarx` object) — For a definition of the states of `idnlarx` models, see “Definition of `idnlarx` States” on page 1-631.

## **PastData — Past input-output data**

`iddata` object | structure

Past input-output data, specified as one of the following:

- `iddata` object — The number of samples must be greater than or equal to the model order. To determine model order, use `order`.

$X$  is the value of model states at time `PastData.SamplingInstants(end) + PastData.Ts`.

When `sys` is continuous-time, specify `PastData` as an `iddata` object.  $X$  then corresponds to the discretized (`c2d`) version of the model, where the discretization method is stored in the `InterSample` property of `PastData`.

- Structure — Specified as a structure with the following fields:
  - Input — Past input data, specified as an  $N$ -by- $Nu$  matrix, where  $N$  is great than or equal to the model order.
  - Output — Past output data, specified as an  $N$ -by- $Ny$  matrix, where  $N$  is great than or equal to the model order.

Specify `PastData` as a structure only when `sys` is a discrete-time model.

The data samples in `PastData` should be in the order of increasing time. That is, the last row in `PastData` should correspond to the latest time.

## **Output Arguments**

### **X — Mapped states of model**

row vector

Mapped states of model, returned as a row vector of size equal to the number of states.  $X$  contains the state value at the time instant immediately after the most recent data sample in `PastData`. That is, if `PastData` is an `iddata` object,  $X$  is the state value at time  $t = \text{PastData.SamplingInstants}(end) + \text{PastData.Ts}$ .

### **XCov — Estimated covariance of state values**

matrix

Estimated covariance of state values, returned as a matrix of size  $N_x$ -by- $N_x$ , where  $N_x$  is the number of states.

XCov is empty if sys is a nonlinear ARX model.

## See Also

[findstates](#) | [getDelayInfo](#) | [idnlarx/findop](#) | [order](#) | [sim](#)

**Introduced in R2008a**

## **db2mag**

Convert decibels (dB) to magnitude

### **Syntax**

`y = db2mag(ydb)`

### **Description**

`y = db2mag(ydb)` returns the corresponding magnitude  $y$  for a given decibel (dB) value  $ydb$ . The relationship between magnitude and decibels is  $ydb = 20 * \log_{10}(y)$ .

### **See Also**

`mag2db`

**Introduced in R2008a**

## dcgain

Low-frequency (DC) gain of LTI system

### Syntax

```
k = dcgain(sys)
```

### Description

`k = dcgain(sys)` computes the DC gain `k` of the LTI model `sys`.

#### Continuous Time

The continuous-time DC gain is the transfer function value at the frequency  $s = 0$ . For state-space models with matrices  $(A, B, C, D)$ , this value is

$$K = D - CA^{-1}B$$

#### Discrete Time

The discrete-time DC gain is the transfer function value at  $z = 1$ . For state-space models with matrices  $(A, B, C, D)$ , this value is

$$K = D + C(I - A)^{-1}B$$

## Examples

### Compute the DC Gain of a MIMO Transfer Function

Create the following 2-input 2-output continuous-time transfer function.

$$H(s) = \begin{bmatrix} 1 & \frac{s-1}{s^2+s+3} \\ \frac{1}{s+1} & \frac{s+2}{s-3} \end{bmatrix}$$

```
H = [1 tf([1 -1],[1 1 3]) ; tf(1,[1 1]) tf([1 2],[1 -3])];
```

Compute the DC gain of the transfer function. For continuous-time models, the DC gain is the transfer function value at the frequency  $s = 0$ .

```
K = dcgain(H)
```

```
K = 2x2
```

```
1.0000 -0.3333
1.0000 -0.6667
```

The DC gain for each input-output pair is returned.  $K(i, j)$  is the DC gain from input  $j$  to output  $i$ .

### Compute DC Gain of Identified Model

Load the estimation data.

```
load iddata1 z1
```

`z1` is an `iddata` object containing the input-output estimation data.

Estimate a process model from the data. Specify that the model has one pole and a time delay term.

```
sys = procest(z1, 'PID')
```

```
sys =
Process model with transfer function:
```

$$G(s) = \frac{K_p}{1+T_p1*s} * \exp(-T_d*s)$$

```
Kp = 9.0754
```

```
Tp1 = 0.25655  
Td = 0.068
```

Parameterization:

```
'PID'
```

```
Number of free coefficients: 3
```

```
Use "getpvec", "getcov" for parameters and their uncertainties.
```

Status:

```
Estimated using PROCEST on time domain data "z1".
```

```
Fit to estimation data: 44.85%
```

```
FPE: 6.02, MSE: 5.901
```

Compute the DC gain of the model.

```
K = dcbgain(sys)
```

```
K = 9.0754
```

This DC gain value is stored in the `Kp` property of `sys`.

```
sys.Kp
```

```
ans = 9.0754
```

## Tips

The DC gain is infinite for systems with integrators.

## See Also

`evalfr` | `norm`

**Introduced in R2012a**

## deadzone

Create a dead-zone nonlinearity estimator object

### Syntax

```
NL = deadzone  
NL = deadzone('ZeroInterval',[a,b])
```

### Description

NL = deadzone creates a default dead-zone nonlinearity estimator object for estimating Hammerstein-Wiener models. The interval in which the dead-zone exists (zero interval) is set to [NaN NaN]. The initial value of the zero interval is determined from the estimation data range, during estimation using n1hw. Use dot notation to customize the object properties, if needed.

NL = deadzone('ZeroInterval',[a,b]) creates a dead-zone nonlinearity estimator object initialized with zero interval, [a,b].

Alternatively, use NL = deadzone([a,b]).

### Object Description

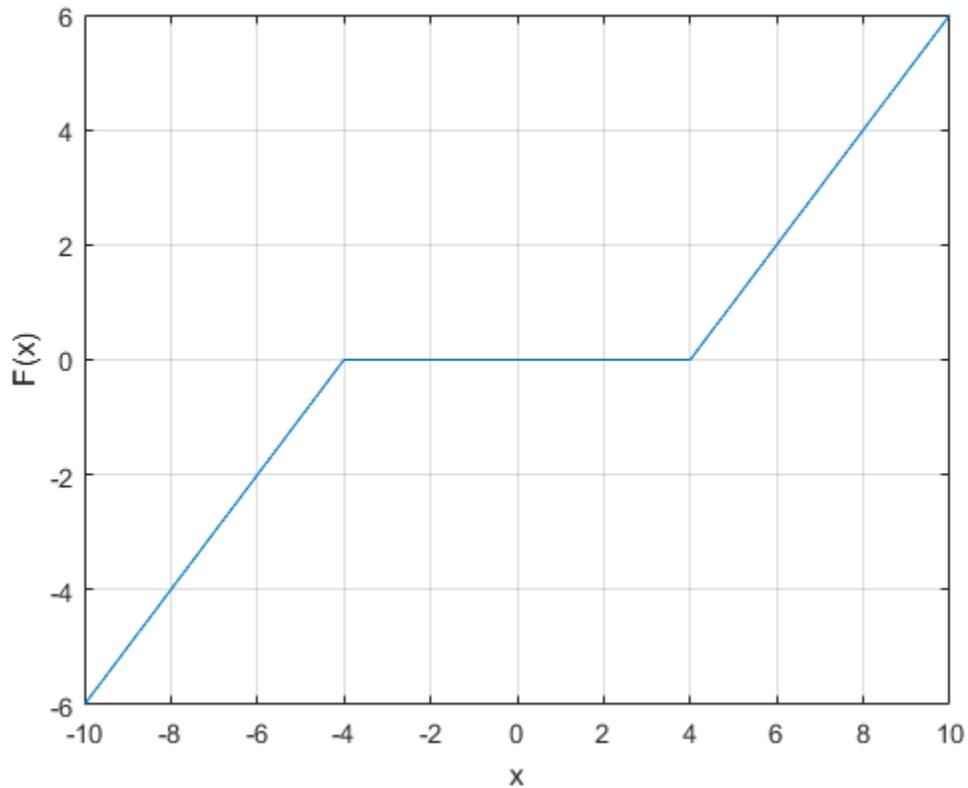
deadzone is an object that stores the dead-zone nonlinearity estimator for estimating Hammerstein-Wiener models.

Use deadzone to define a nonlinear function  $y = F(x, \theta)$ , where  $y$  and  $x$  are scalars, and  $\theta$  represents the parameters  $a$  and  $b$ , which define the zero interval.

The dead-zone nonlinearity function has the following characteristics:

$$\begin{array}{ll} a \leq x < b & F(x) = 0 \\ x < a & F(x) = x - a \\ x \geq b & F(x) = x - b \end{array}$$

For example, in the following plot, the dead-zone is in the interval  $[-4, 4]$ .



The value  $F(x)$  is computed by `evaluate(NL, x)`, where `NL` is the deadzone object.

For deadzone object properties, see “Properties” on page 1-293.

## Examples

### Create a Default Dead-Zone Nonlinearity Estimator

```
NL = deadzone;
```

Specify the zero interval.

```
NL.ZeroInterval = [-4,5];
```

## Estimate a Hammerstein-Wiener Model with Dead-zone Nonlinearity

Load estimation data.

```
load twotankdata;  
z = iddata(y,u,0.2,'Name','Two tank system');  
z1 = z(1:1000);
```

Create a deadzone object, and specify the initial guess for the zero-interval.

```
OutputNL = deadzone('ZeroInterval',[-0.1 0.1]);
```

Estimate model with no input nonlinearity.

```
m = nlhw(z1,[2 3 0],[],OutputNL);
```

## Estimate MIMO Hammerstein-Wiener Model

Load the estimation data.

```
load motorizedcamera;
```

Create an iddata object.

```
z = iddata(y,u,0.02,'Name','Motorized Camera','TimeUnit','s');
```

z is an iddata object with 6 inputs and 2 outputs.

Specify the model orders and delays.

```
Orders = [ones(2,6),ones(2,6),ones(2,6)];
```

Specify the same nonlinearity estimator for each input channel.

```
InputNL = saturation;
```

Specify different nonlinearity estimators for each output channel.

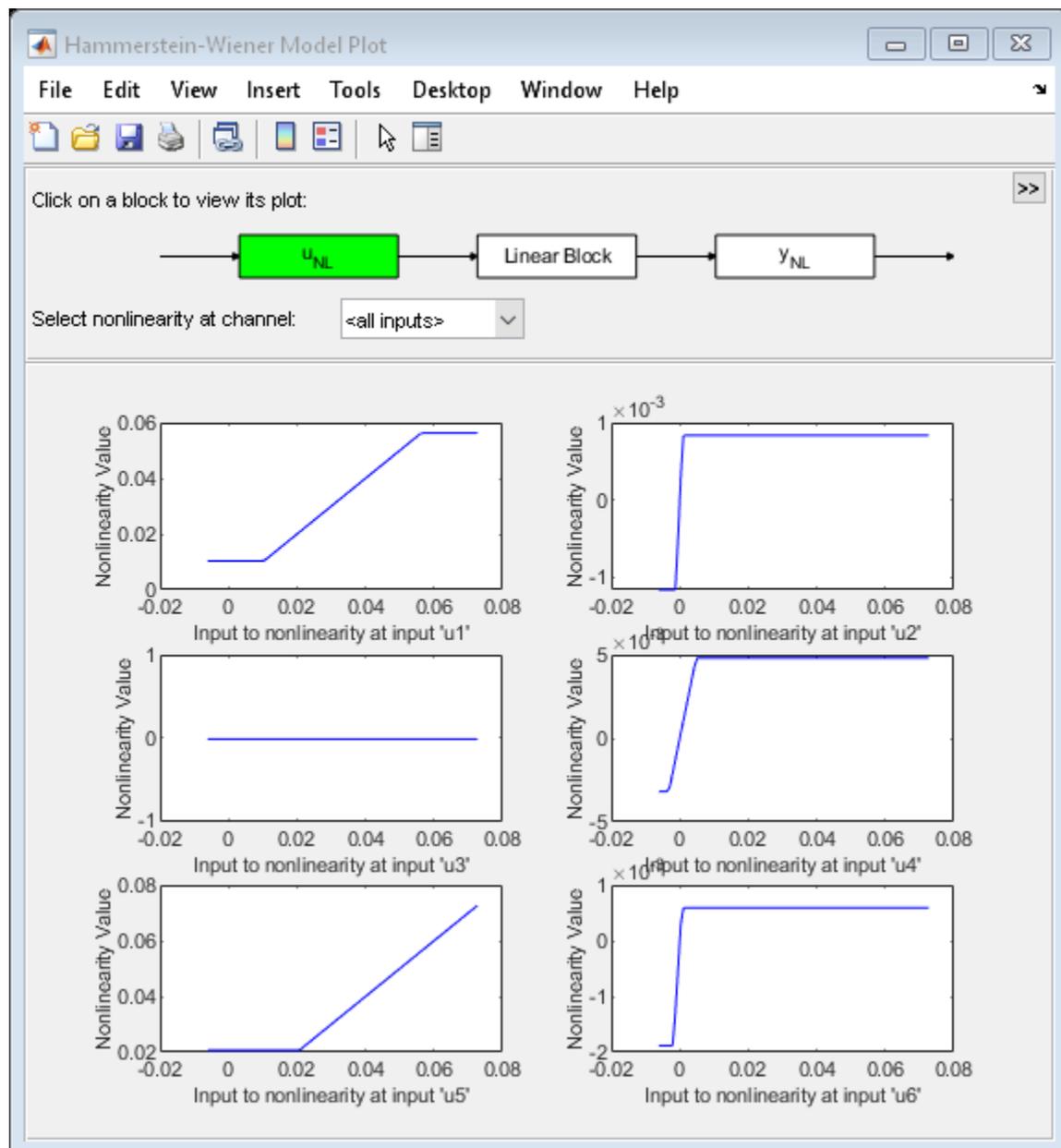
```
OutputNL = [deadzone,wavenet];
```

Estimate the Hammerstein-Wiener model.

```
sys = nlhw(z,Orders,InputNL,OutputNL);
```

To see the shape of the estimated input and output nonlinearities, plot the nonlinearities.

```
plot(sys)
```



Click on the input and output nonlinearity blocks on the top of the plot to see the nonlinearities.

## Input Arguments

### **[a, b] — Zero interval**

[NaN NaN] (default) | 2-element row vector

Zero interval of the dead-zone, specified as a 2-element row vector of doubles.

The dead-zone nonlinearity is initialized at the interval [a, b]. The interval values are adjusted to the estimation data by `n_lhw`. To remove the lower limit, set a to `-Inf`. The lower limit is not adjusted during estimation. To remove the upper limit, set b to `Inf`. The upper limit is not adjusted during estimation.

When the interval is [NaN NaN], the initial value of the zero interval is determined from the estimation data range during estimation using `n_lhw`.

Example: [-2 1]

## Properties

### **ZeroInterval**

Zero interval of the dead-zone, specified as a 2-element row vector of doubles.

**Default:** [NaN NaN]

## Output Arguments

### **NL — Dead-zone nonlinearity estimator object**

deadzone object

Dead-zone nonlinearity estimator object, returned as a `deadzone` object.

## **See Also**

n|hw

**Introduced in R2007a**

# delayest

Estimate time delay (dead time) from data

## Syntax

```
nk = delayest(Data)
nk = delayest(Data, na, nb, nkmin, nkmax, maxtest)
```

## Description

`nk = delayest(Data)` estimates time delay from data. `Data` is an `iddata` object containing the input-output data. It can also be an `idfrd` object defining frequency-response data. Only single-output data can be handled. `nk` is returned as an integer or a row vector of integers, containing the estimated time delay in samples from the input(s) to the output in `Data`.

The estimate is based on a comparison of ARX models with different delays:

$$y(t) + a_1 y(t-1) + \dots + a_{na} y(t-na) = b_1 u(t-nk) + \dots + b_{nb} u(t-nb-nk+1) + e(t)$$

`nk = delayest(Data, na, nb, nkmin, nkmax, maxtest)` specifies additional options. The integer `na` is the order of the A polynomial (default 2). `nb` is a row vector of length equal to the number of inputs, containing the order(s) of the B polynomial(s) (default all 2). `nkmin` and `nkmax` are row vectors of the same length as the number of inputs, containing the smallest and largest delays to be tested. Defaults are `nkmin = 0` and `nkmax = nkmin+20`. If `nb`, `nkmax`, and/or `nkmin` are entered as scalars in the multiple-input case, all inputs will be assigned the same values. `maxtest` is the largest number of tests allowed (default 10,000).

**Introduced before R2006a**

## detrend

Subtract offset or trend from time-domain signals contained in `iddata` objects

`detrend` subtracts offsets or linear trends from time-domain input-output data represented in `iddata` objects. `detrend` either computes the trend data to subtract, or subtracts the trend that you specify.

For a more general detrending function that does not require `iddata` objects, see `detrend`.

## Syntax

```
data_d = detrend(data)
data_d = detrend(data,Type)
[data_d,T_r] = detrend( ___ )

data_d = detrend(data,1,brkpt)
```

## Description

`data_d = detrend(data)` computes and subtracts the mean value from each time-domain signal in `data`. The `iddata` objects `data_d` and `data` each contain input and output data originating from SISO, MIMO, or multiexperiment datasets.

`data_d = detrend(data,Type)` subtracts the trend you specify in `Type`. You can specify a mean-value, linear, or custom trend.

`[data_d,T_r] = detrend( ___ )` also returns the subtracted trend as a `TrendInfo` object, `T_r`. You can obtain `T_r` with any of the input-argument combinations in previous syntaxes.

`data_d = detrend(data,1,brkpt)` computes and subtracts the piecewise-linear trends for data with segmented trends, using the breakpoints that you define with `brkpt`.

The second argument, which corresponds to `Type`, must be `1`.

With this syntax, you cannot retrieve the resulting piecewise-linear trend information as an output.

## Examples

### Remove Biases From Signals

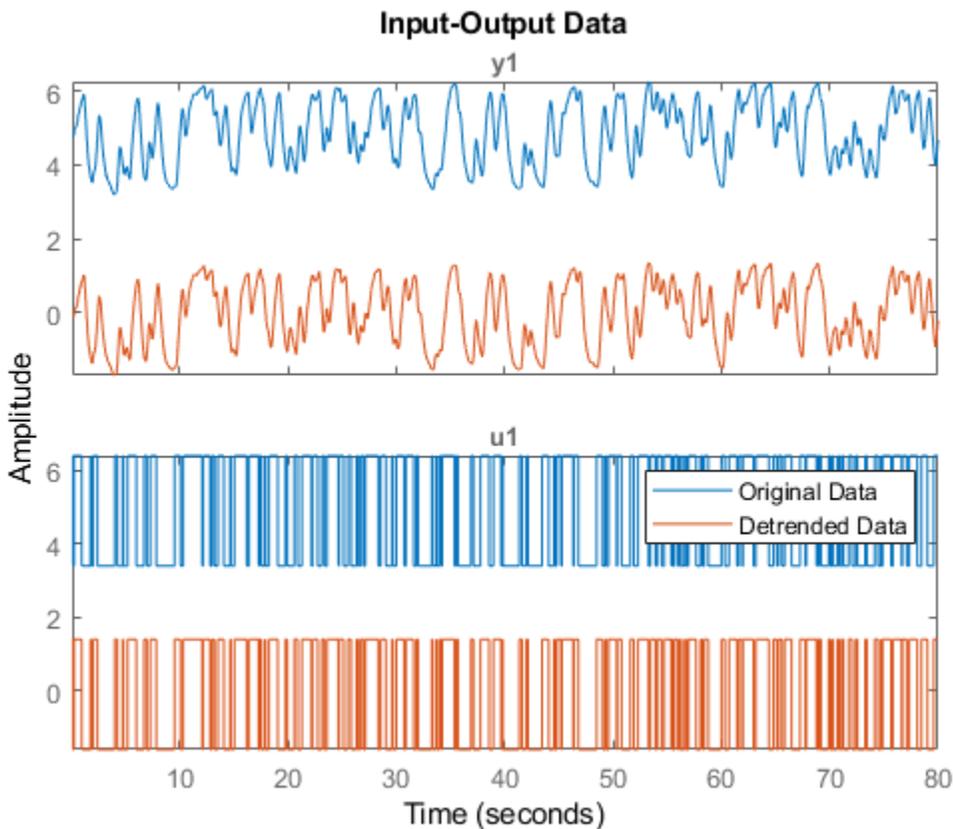
Remove biases from steady-state signals in an `iddata` object by using `detrend` to compute and subtract the mean values of the input and output.

Load the input and output time series data `y2` and `u2`. Construct the `iddata` object `data2`, using the data and a sample time of 0.08 seconds.

```
load dryer2 y2 u2
data2 = iddata(y2,u2,0.08);
```

Use `detrend` to both compute the mean values and subtract them from input and output signals. Use the input argument `Tr` to store the computed trend information. Plot the original data and detrended data together.

```
[data2_d,Tr] = detrend(data2);
plot(data2,data2_d)
legend('Original Data','Detrended Data')
```



The detrended data has shifted by about 5 units. Inspect `Tr` to obtain the precise mean values that `detrend` computed and subtracted. These values are returned in the `InputOffset` and `OutputOffset` properties.

`Tr`

```
Trend specifications for data "data2" with 1 input(s), 1 output(s), 1 experiment(s):
  DataName: 'data2'
  InputOffset: 5.0000
  OutputOffset: 4.8901
  InputSlope: 0
  OutputSlope: 0
```

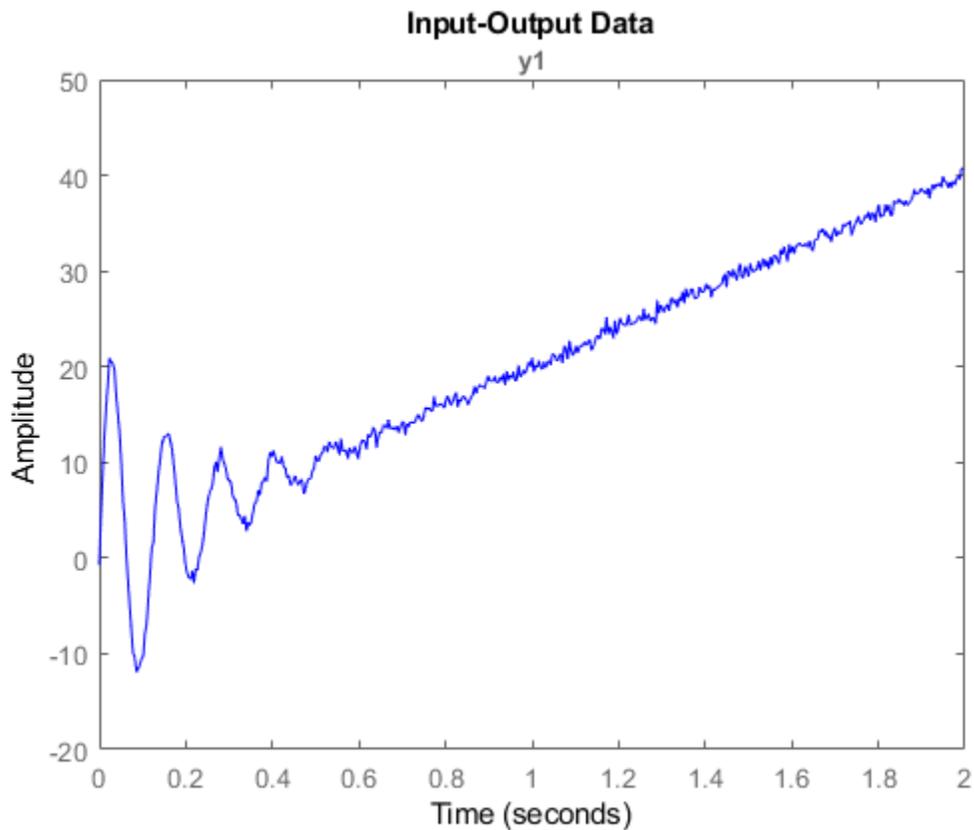
The mean of the original input is higher than the mean of the original output.

## Remove Linear Trend from a Signal

Remove the linear trend from a signal in an `iddata` object, and overlay the trendline on a before-and-after data plot.

Load and plot signal data from the file `lintrend2`. For this example, only output data is provided in `iddata` object `dataL`.

```
load lintrend2 dataL
plot(dataL, 'b')
```

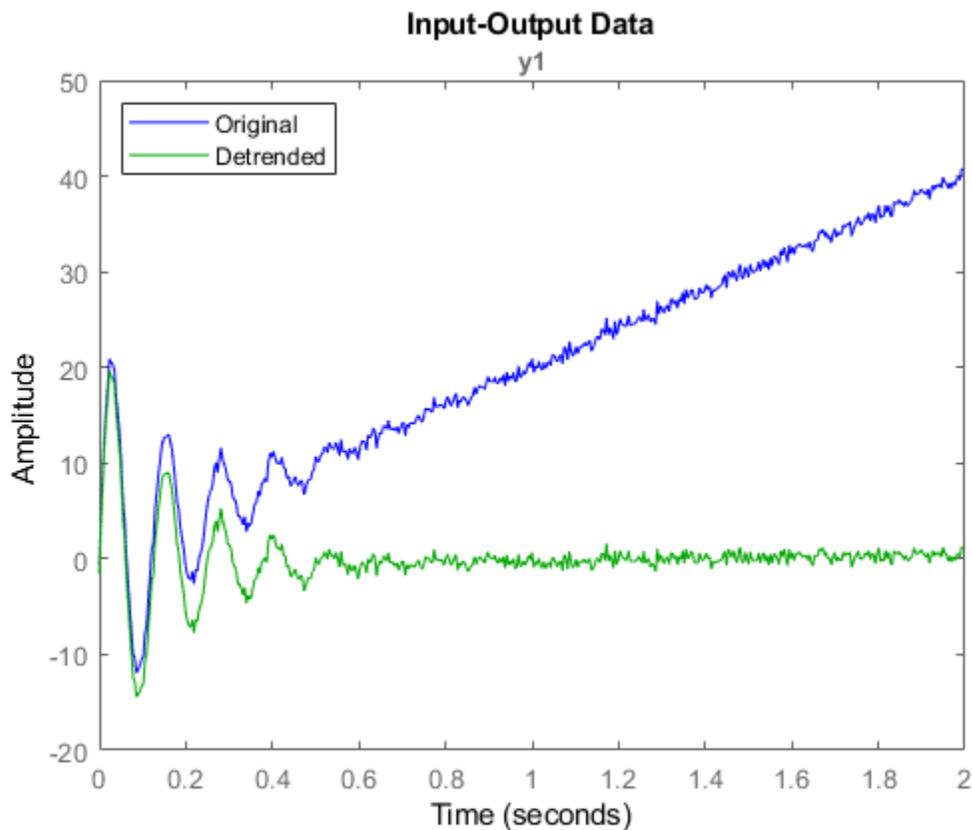


The plot shows a clear linear trend in the data. Use `detrend` linear option (`Type = 1`) to subtract the trend from the data. `detrend` fits the data and determines the linear trend to subtract. Include the `TrendInfo` object `Tr` as an output argument so you can see what `detrend` subtracts.

```
[dataL_d,Tr] = detrend(dataL,1);
```

Plot the detrended data against the original data.

```
hold on  
plot(dataL_d,'g')  
legend('Original','Detrended','Location','northwest')
```



The linear trend has been removed. Inspect `Tr` to get more information about the removed trend.

`Tr`

```
Trend specifications for data "dataL" with 0 input(s), 1 output(s), 1 experiment(s):
  DataName: 'dataL'
  InputOffset: [1x0 double]
  OutputOffset: 0.8888
  InputSlope: [1x0 double]
  OutputSlope: 19.3830
```

The `OutputOffset` and the `OutputSlope` properties provide the parameters of the line that `detrend` removed. You can reconstruct this line, and then overlay it on the before-and-after data plot. The `SamplingInstants` property of `DataL` provides the timepoints associated with the data.

```
m = Tr.OutputSlope
```

```
m = 19.3830
```

```
b = Tr.OutputOffset
```

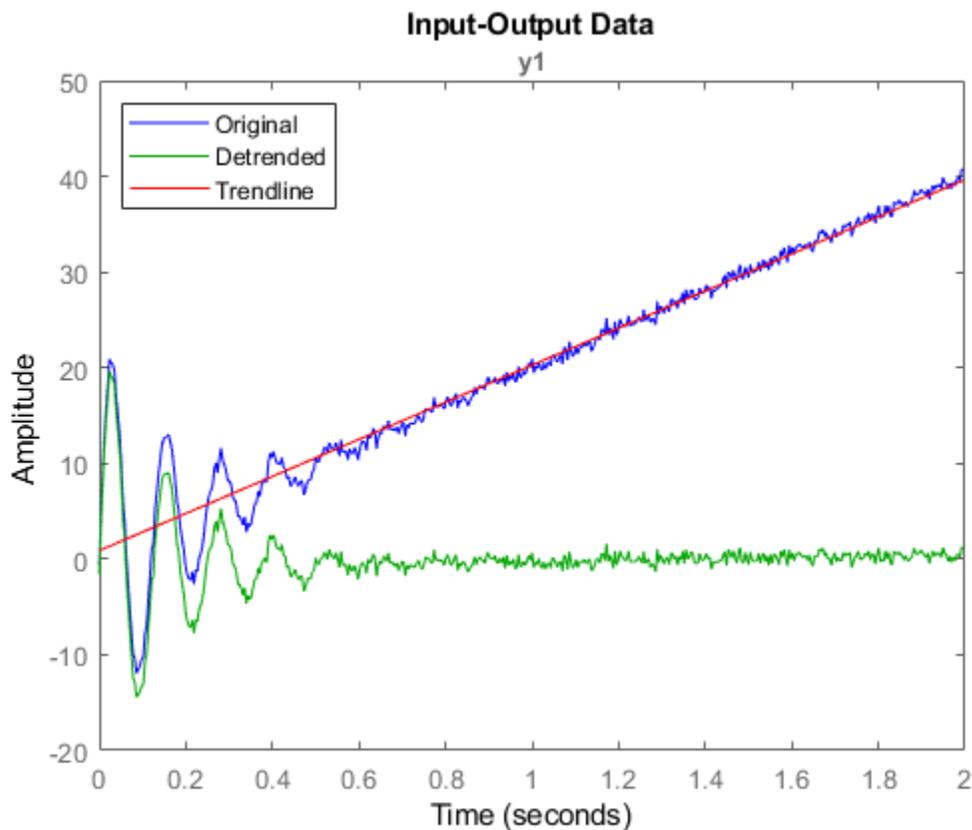
```
b = 0.8888
```

```
t = dataL.SamplingInstants;
```

```
TrLn = m*t+b;
```

```
plot(t,TrLn,'r')
```

```
legend('Original','Detrended','Trendline','Location','northwest')
```



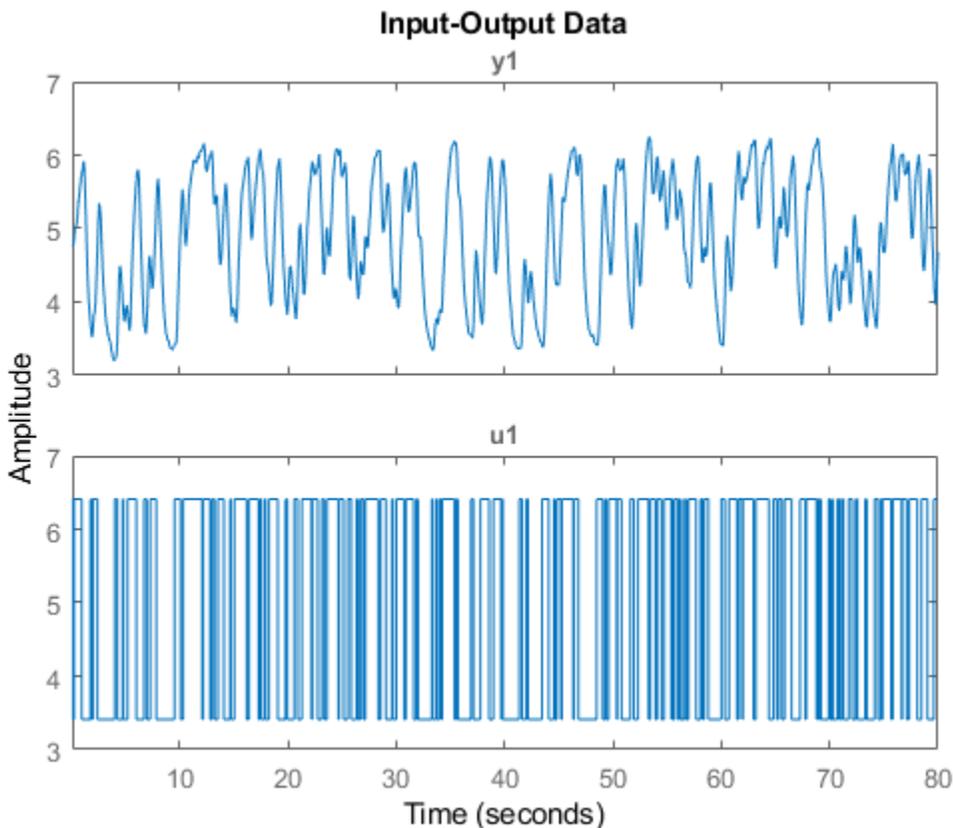
### Remove Specified Offsets from Signals

Remove known offsets from an input-output signal pair contained in an `iddata` object.

`detrend` can compute and subtract the mean values for input and output signals, resulting in zero-mean detrended signals. However, if you already know you have specific data offsets beforehand, you can have `detrend` subtract these from your signals instead. Specifying the offsets also allows you to retain a non-zero operating point in the `detrend` result.

Load SISO data containing vectors `u2` and `y2`. For this example, suppose that you know both signals have an offset of 4 from the expected operating point of 1. Combine these vectors into an `iddata` object, using a sample time of 0.08 seconds, and plot it.

```
load dryer2 u2 y2
data = iddata(y2,u2,0.08);
plot(data)
```



The known offset of 4 (from operating point 1) is visible in the plots. You can construct a `TrendInfo` object to capture this offset, using the function `getTrend`.

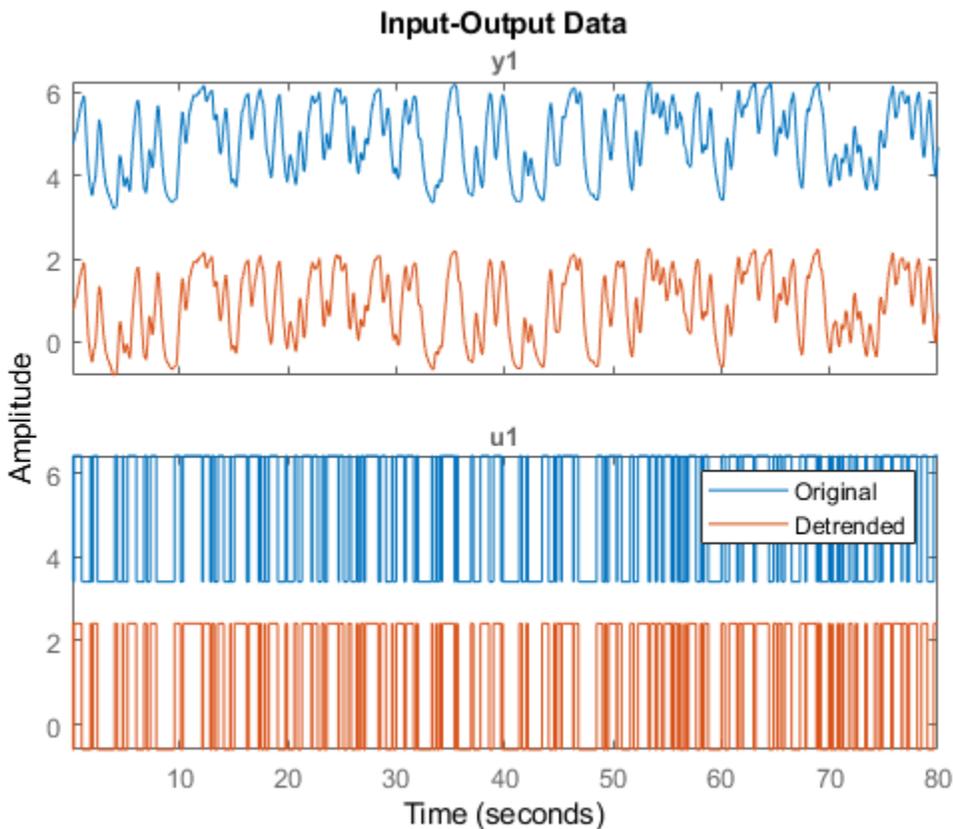
Create the `TrendInfo` object, and then set its offset properties.

```
T = getTrend(data);  
T.InputOffset = 4;  
T.OutputOffset = 4
```

```
Trend specifications for data "data" with 1 input(s), 1 output(s), 1 experiment(s):  
  DataName: 'data'  
  InputOffset: 4  
  OutputOffset: 4  
  InputSlope: 0  
  OutputSlope: 0
```

Use `detrend` to subtract the offset from the data. Plot it alongside the original data.

```
data_d = detrend(data,T);  
hold on  
plot(data_d)  
legend('Original', 'Detrended')
```



The offset of 4 has been removed.

### Remove Segmented Linear Trends from Signals by using Breakpoints

Detrend data with segmented piecewise-linear trends by specifying breakpoints to delimit the segments.

Most of the `detrend` syntaxes assume and compute a single trend for each of the signals. However, in some cases there are discontinuities in the linear trends, caused by test configuration changes, environmental conditions, or other influences. When the signal displays such segmentation, you can have `detrend` operate on the test segments

independently. To do so, specify breakpoints in the `brkpt` input argument. These are the indices to the timepoints in the signal at which linear trends change slope.

You may know these breakpoints up front, based on changes that you know occurred during data collection. Alternatively, you may need to approximate them by inspecting the data itself.

Load the data, inspect its structure and contents, and plot it. This data consists of output data only in the `iddata` object `dataLb2`.

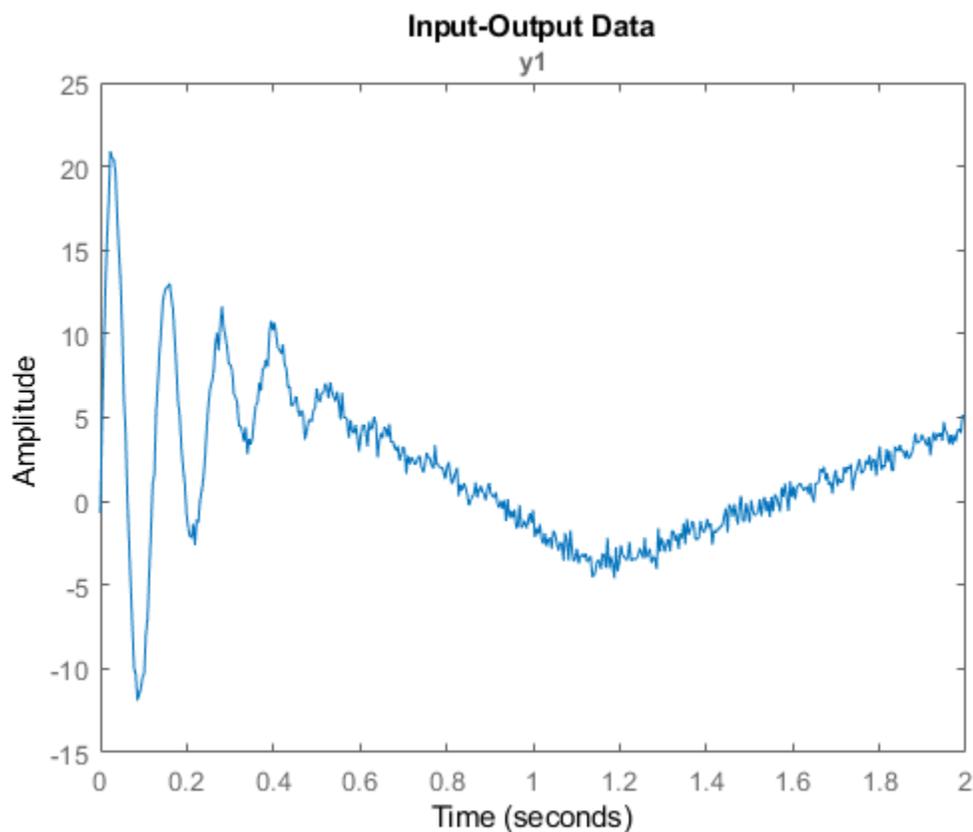
```
load brkTrend dataLb2
dataLb2
```

```
dataLb2 =
```

```
Time domain data set with 512 samples.
Sample time: 0.00390625 seconds
```

```
Outputs      Unit (if specified)
  y1
```

```
plot(dataLb2)
```



For this example, the data has known breakpoints at indices [100 300]. Applying the sample time (property Ts), these breakpoints correspond to the actual timepoints as follows:

```
brkpt=[100 300];
brkpt_time = brkpt*dataLb2.Ts

brkpt_time = 1x2

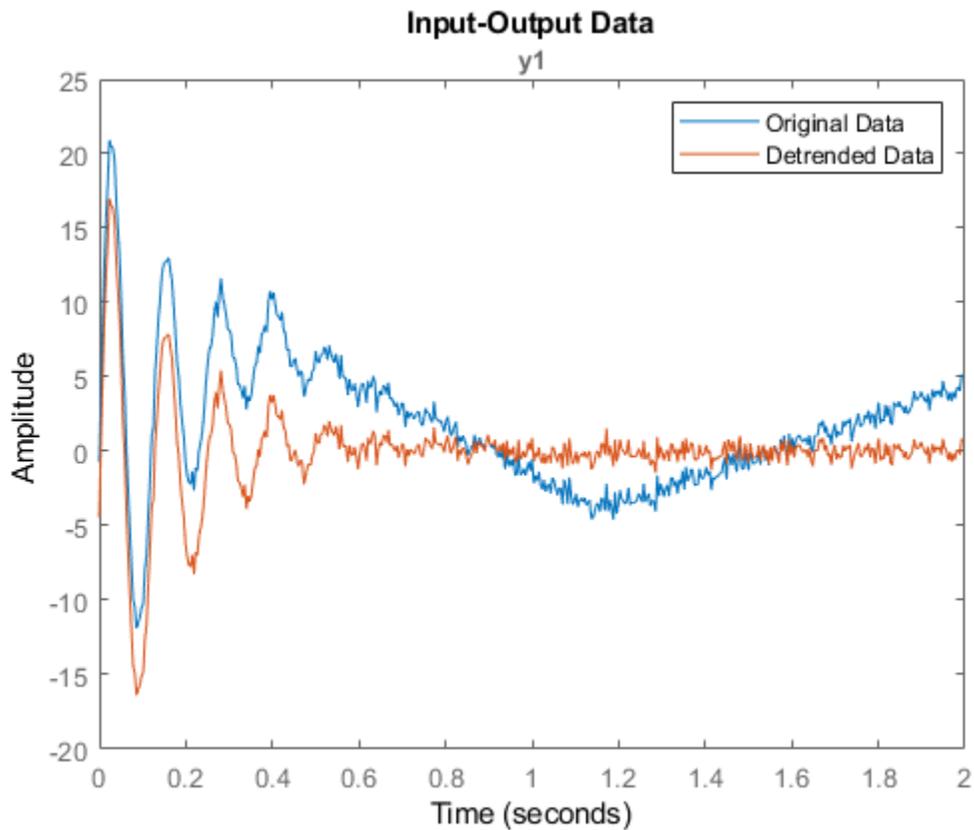
    0.3906    1.1719
```

Detrend the data using brkpt.

```
dataLb2_d = detrend(dataLb2,1,brkpt);
```

Plot the original and detrended data.

```
plot(dataLb2,dataLb2_d)  
legend('Original Data','Detrended Data')
```



The linear trend segments have been removed.

### **Detrend Multiexperiment Signals using Multiple-Breakpoint Sets**

Apply a unique set of breakpoints to each experiment when you detrend a Multiexperiment dataset.

Experiments within a multiexperiment dataset may contain unique linear trending discontinuities. You can apply a unique set of breakpoints to each experiment by expressing them in a cell array.

Load the data, which consists of:

- `datmult`, a multiexperiment `iddata` object containing three experiments (output only)
- `bpn` vectors, which provide known breakpoints for each experiment in the form of indices to timepoints

```
load multiexpdt datmult bp1 bp2 bp3
datmult
```

```
datmult =
Time domain data set containing 3 experiments.
```

Experiment	Samples	Sample Time
exp1	250	1
exp2	320	1
exp3	350	1

Outputs	Unit (if specified)
y1	

```
bp1, bp2, bp3
```

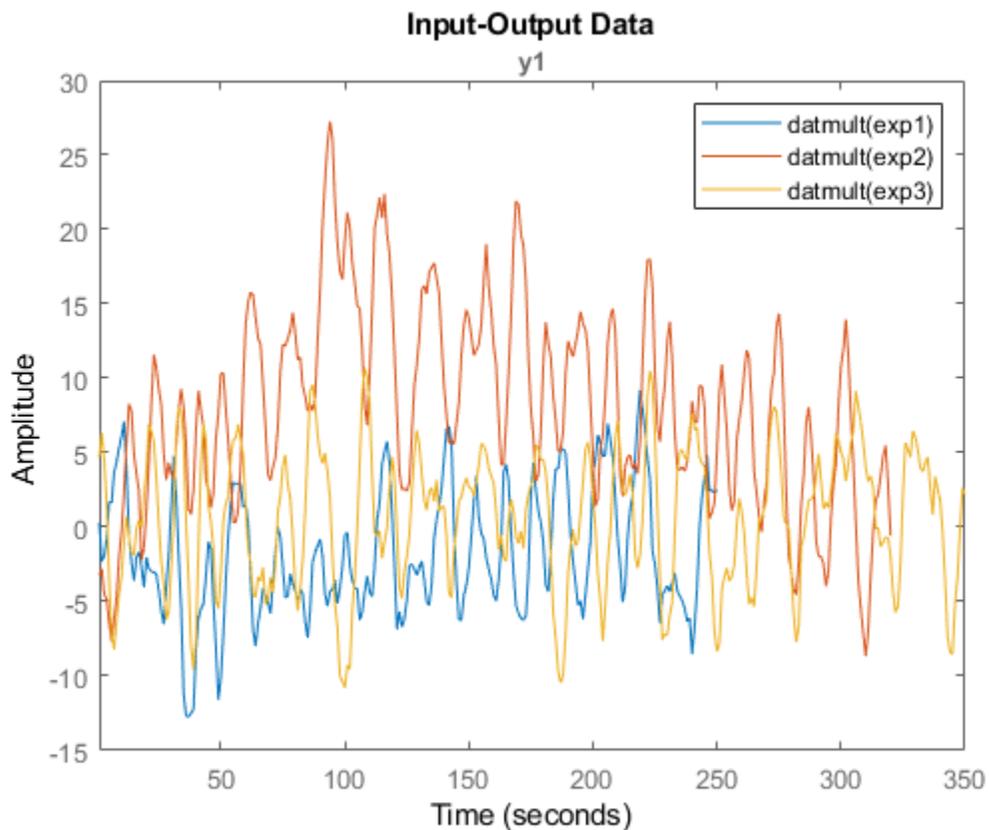
```
bp1 = 1x2
      50 200
```

```
bp2 = 100
```

```
bp3 =
      []
```

Plot the data. There are significant differences among the streams, and they drift at different rates from zero mean.

```
plot(datmult)
legend
```



For this set of experiments, it is known that there is unique trending for each run and unique discontinuities indicated by the `bp` vectors.

`detrend` can incorporate these unique characteristics if the `bp` information is provided as a cell array.

Construct the cell array.

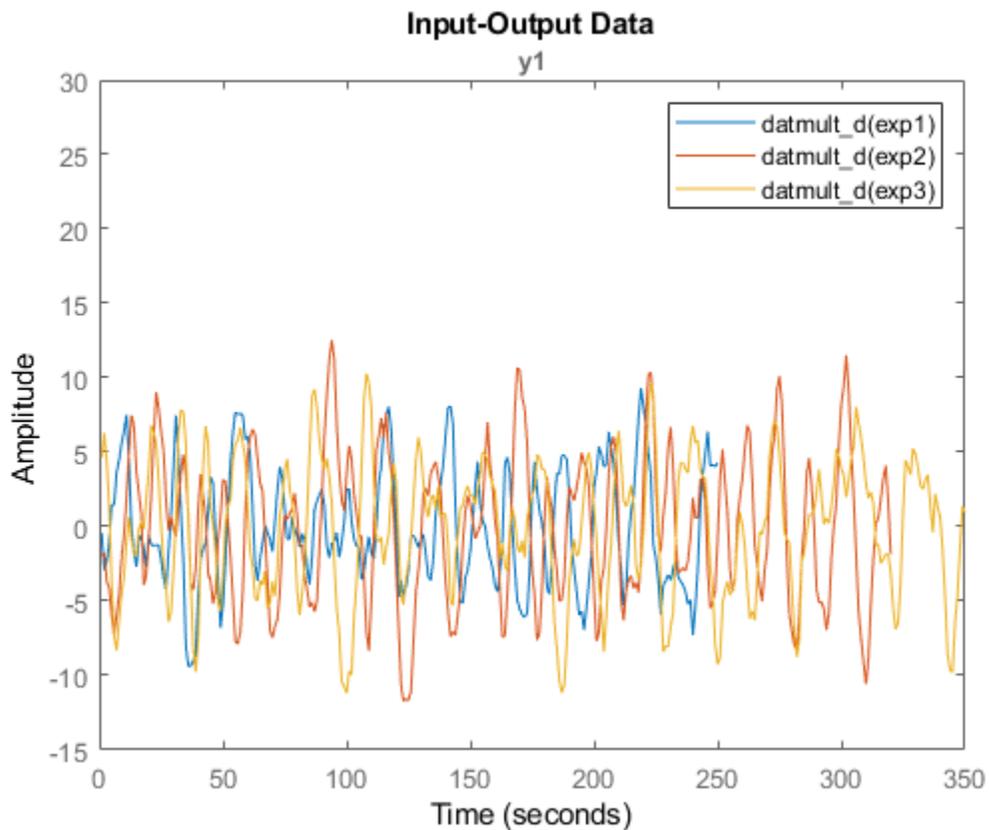
```
bpcell = {bp1;bp2;bp3}
```

```
bpcell = 3x1 cell array  
    {1x2 double}  
    {[    100]}
```

```
{0x0 double}
```

Apply `detrend` and plot the result, using the same scale as the original plot.

```
datmult_d = detrend(datmult,1,bpcell);  
figure  
plot(datmult_d)  
axis([0,inf,-15,30])  
legend
```



The experimental data are now better aligned, and do not drift significantly away from zero mean.

## Detrend Input and Output Signals Separately

Apply different trend types to the input and output signals contained in an `iddata` object.

`Detrend` assumes that the same type of trend applies to both input and output signals. In some cases, there may be a trend type that is present in only one signal. You can perform `detrend` individually on each signal by extracting the signals into separate `iddata` objects. Apply `detrend` to each object using its individual signal trend type. Then reassemble the results back into a single detrended `iddata` object.

Load, examine, and plot the data in `iodatab`.

```
load septrend iodatab;  
iodatab
```

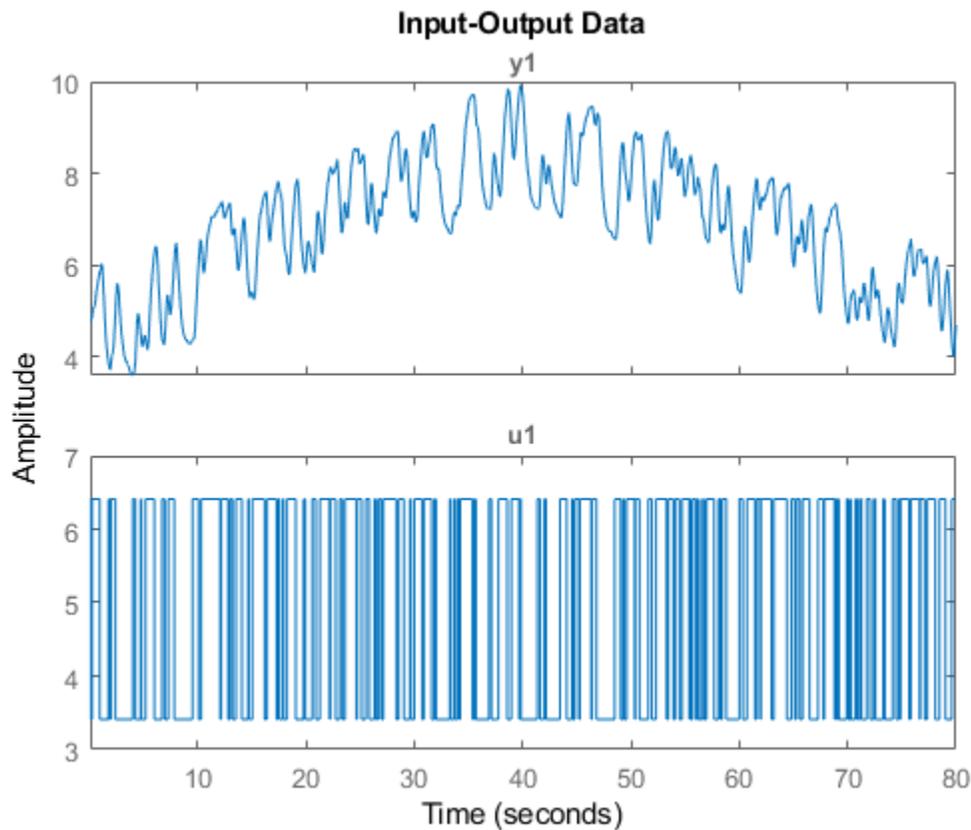
```
iodatab =
```

```
Time domain data set with 1000 samples.  
Sample time: 0.08 seconds
```

```
Outputs      Unit (if specified)  
  y1
```

```
Inputs      Unit (if specified)  
  u1
```

```
plot(iodatab)  
hold on
```



Both input and output plots show a bias. However, the output plot also shows an inverted V-shape trend that is not present in the input data.

Separate the input data and the output data into separate objects for detrending, using the `iddata` general data-selection form (see "Representing Time- and Frequency-Domain Data Using `iddata` Objects"):

```
data(samples,outputchannels,inputchannels)
```

```
idatab = iodatab(:,[],:);
odatab = iodatab(:,[:,[]]);
```

Remove the bias from the input data, using `detrrend` to calculate and subtract the mean.

```
idatab_d = detrend(idatab,0);
```

Remove the bias and the inverted-V trend from the output data, using the midpoint index 500 as a breakpoint.

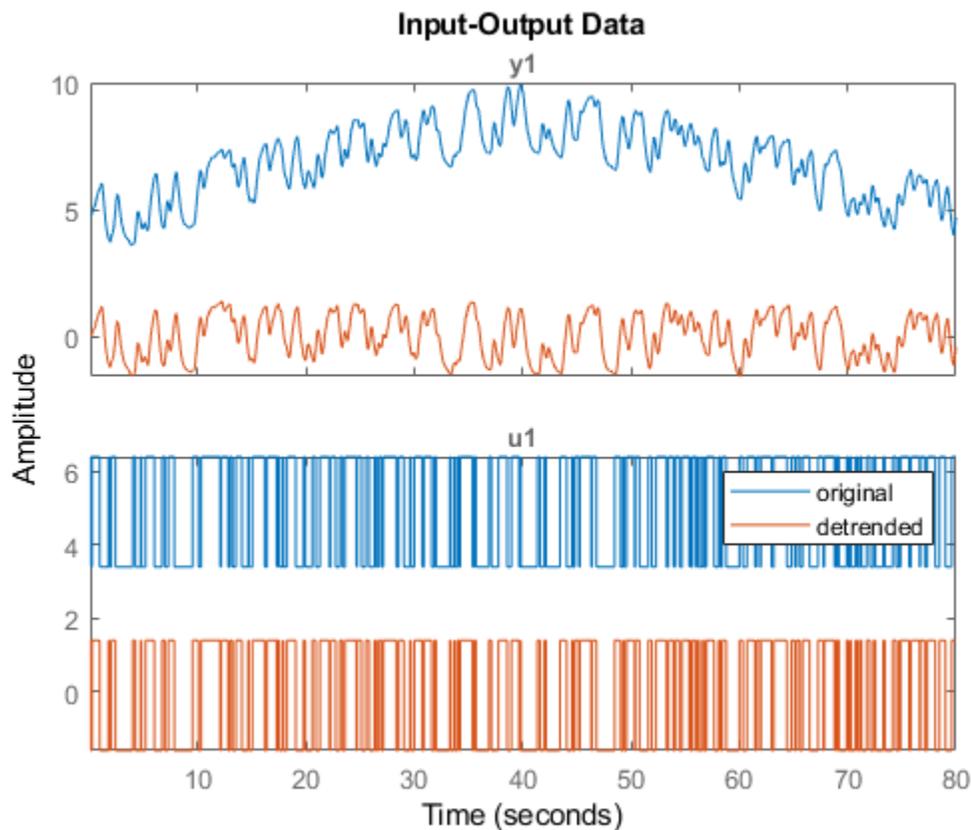
```
odatab_d = detrend(odatab,1,500);
```

Combine the detrended input and output data into a single `iddata` object.

```
iodatab_d = [odatab_d,idatab_d];
```

Overlay the detrended data on the original data.

```
plot(iodatab_d)  
legend('original','detrended')
```



The input and output data now contain neither bias nor V-shape trend.

## Input Arguments

### **data — Time-domain input-output data**

`iddata` object

Specified as an `iddata` object containing one or more sets of time-domain signals. The `iddata` object can contain SISO, MIMO, or multiexperiment data. The signal sets can contain either input and output data, or output data only.

### **Type — Trend type to be subtracted**

0 (default) | 1 | `TrendInfo` object

Specified as one of:

- 0 — Compute and subtract the mean value
- 1 — Compute and subtract the linear trend (least-squares fit)
- `TrendInfo` object — subtract the trend you specify in the `TrendInfo` object. Use `getTrend` to create a `TrendInfo` object. For an example, see “Remove Specified Offsets from Signals” on page 1-302.

### **brkpt — Timepoint locations of trending discontinuities**

integer row vector | cell array of integer vectors

Timepoint locations of trending discontinuities (breakpoints), specified as:

- An integer row vector — For single-experiment SISO and MIMO datasets. Doing so applies a single set of breakpoints to all input and output signals. For an example, see “Remove Segmented Linear Trends from Signals by using Breakpoints” on page 1-305.
- A cell array containing individually-sized integer row vectors — For multiple-experiment datasets. Doing so applies a unique set of breakpoints to the output and input signals for each experiment. For an example, see “Detrend Multiexperiment Signals using Multiple-Breakpoint Sets” on page 1-308.

## Output Arguments

### **data\_d — Detrended signals**

`iddata` object

Detrended signals, returned as an `iddata` object. Dimensions of the contents are the same as dimensions of the contents of `data`.

### **T\_r — Subtracted trend data**

`TrendInfo` object

Trend data subtracted from `data` to produce `data_d`, returned as a `TrendInfo` object .

When you use `brkpt` to specify multiple trends, you cannot retrieve the computed trend data.

## See Also

`getTrend` | `iddata` | `idfilt` | `retrend`

**Introduced before R2006a**

# diff

Difference signals in iddata objects

## Syntax

```
zdi = diff(z)  
zdi = diff(z,n)
```

## Description

`zdi = diff(z)` and `zdi = diff(z,n)` return the difference signals in `iddata` objects. `z` is a time-domain `iddata` object. `diff(z)` and `diff(z,n)` apply this command to each of the input/output signals in `z`.

**Introduced before R2006a**

## etfe

Estimate empirical transfer functions and periodograms

### Syntax

```
g = etfe(data)
g = etfe(data,M)
g = etfe(data,M,N)
```

### Description

`g = etfe(data)` estimates a transfer function of the form:

$$y(t) = G(q)u(t) + v(t)$$

`data` contains time- or frequency-domain input-output data or time-series data:

- If `data` is time-domain input-output signals, `g` is the ratio of the output Fourier transform to the input Fourier transform for the data.

For nonperiodic data, the transfer function is estimated at 128 equally-spaced frequencies  $[1:128]/128*\pi/T_s$ .

For periodic data that contains a whole number of periods (`data.Period = integer`), the response is computed at the frequencies  $k*2*\pi/\text{period}$  for  $k = 0$  up to the Nyquist frequency.

- If `data` is frequency-domain input-output signals, `g` is the ratio of output to input at all frequencies, where the input is nonzero.
- If `data` is time-series data (no input channels), `g` is the periodogram, that is the normed absolute square of the Fourier transform, of the data. The corresponding spectral estimate is normalized, as described in "Spectrum Normalization" and differs from the spectrum normalization in the Signal Processing Toolbox™ product.

`g = etfe(data,M)` applies a smoothing operation on the raw spectral estimates using a Hamming Window that yields a frequency resolution of about  $\pi/M$ . The effect of `M` is

similar to the effect of  $M$  in `spa`.  $M$  is ignored for periodic data. Use this syntax as an alternative to `spa` for narrowband spectra and systems that require large values of  $M$ .

`g = etfe(data,M,N)` specifies the frequency spacing for nonperiodic data.

- For nonperiodic time-domain data,  $N$  specifies the frequency grid  $[1:N]/N*\pi/T_s$  rad/TimeUnit. When not specified,  $N$  is 128.
- For periodic time-domain data,  $N$  is ignored.
- For frequency-domain data, the  $N$  is `fmin:delta_f:fmax`, where `[fmin fmax]` is the range of frequencies in `data`, and `delta_f` is  $(f_{\max} - f_{\min}) / (N - 1)$  rad/TimeUnit. When not specified, the response is computed at the frequencies contained in `data` where input is nonzero.

## Examples

### Compare an Empirical Transfer Function to a Smoothed Spectral Estimate

Load estimation data.

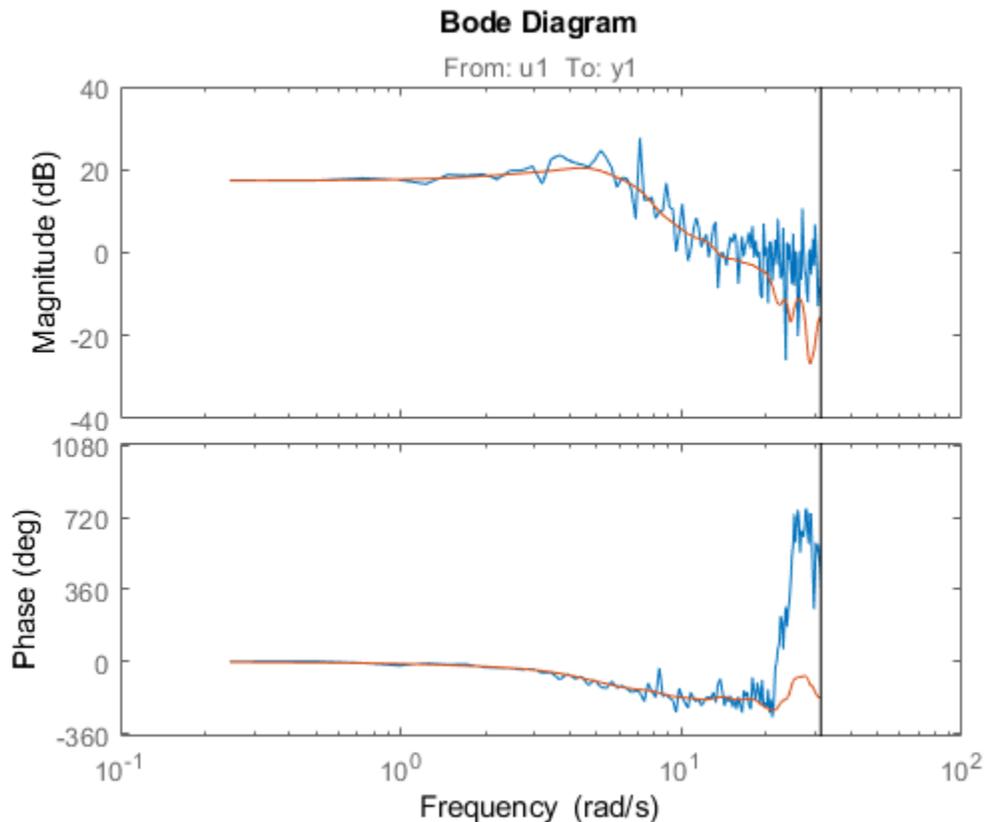
```
load iddata1 z1;
```

Estimate empirical transfer function and smoothed spectral estimate.

```
ge = etfe(z1);  
gs = spa(z1);
```

Compare the two models on a Bode plot.

```
bode(ge,gs)
```



### Generate Empirical Transfer Function Using Periodic Input

Generate a periodic input, simulate a system with it, and compare the frequency response of the estimated model with the original system at the excited frequency points.

Generate a periodic input signal and output signal using simulation.

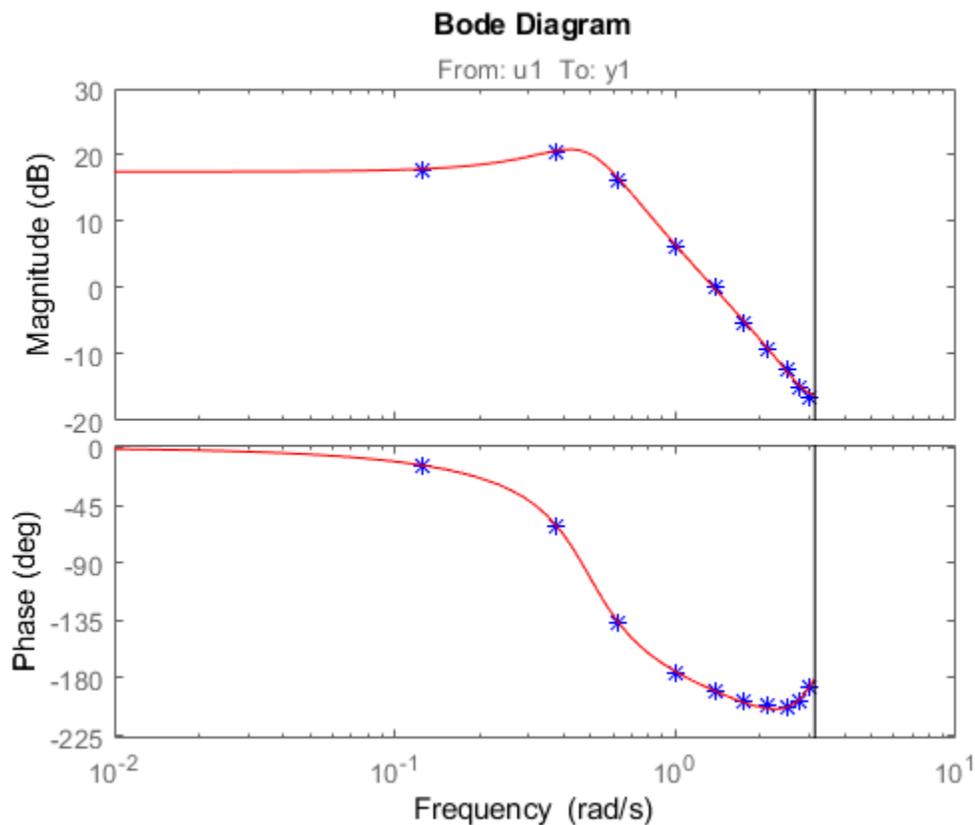
```
m = idpoly([1 -1.5 0.7],[0 1 0.5]);  
u = iddata([],idinput([50,1,10], 'sine'));  
u.Period = 50;  
y = sim(m,u);
```

Estimate an empirical transfer function.

```
me = etfe([y u]);
```

Compare the empirical transfer function with the original model.

```
bode(me, 'b*',m, 'r')
```



### Apply Smoothing Operation on Empirical Transfer Function Estimate

Perform a smoothing operation on raw spectral estimates using a Hamming Window and compare the responses.

Load data.

```
load iddata1
```

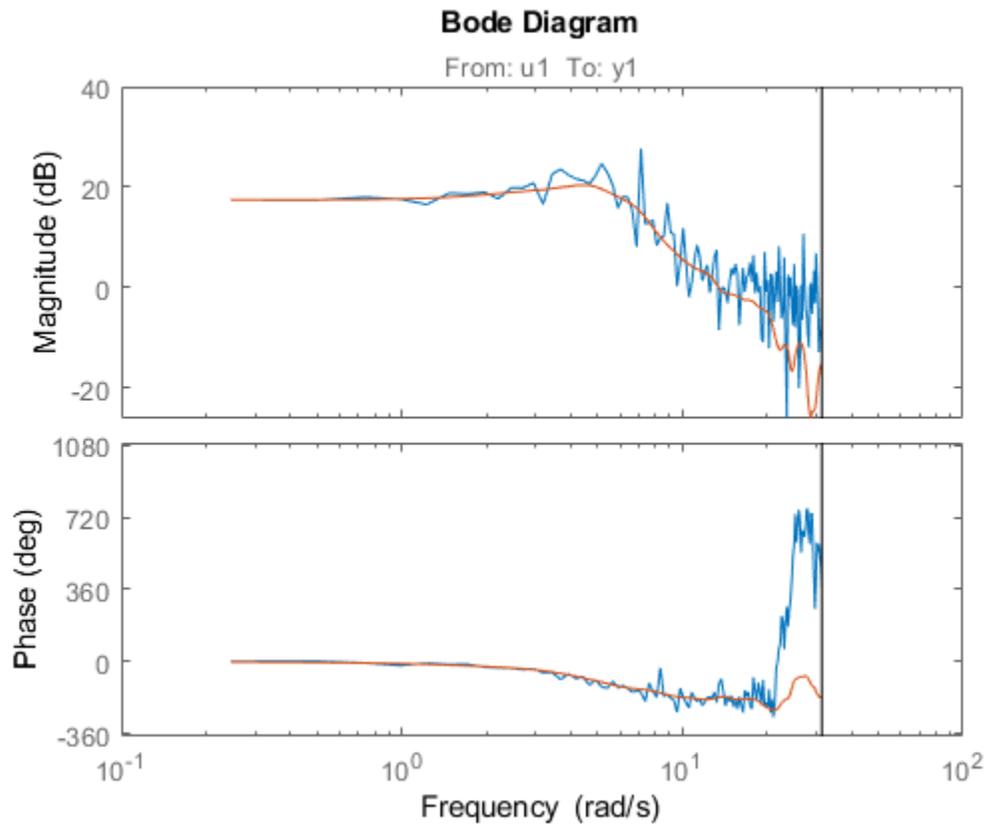
Estimate empirical transfer functions with and without the smoothing operation.

```
ge1 = etfe(z1);  
ge2 = etfe(z1,32);
```

Compare the models on a Bode plot.

ge2 is smoother than ge1 because of the effect of the smoothing operation.

```
bode(ge1,ge2)
```



## Compare Effect of Frequency Spacing on Empirical Transfer Function Estimate

Estimate empirical transfer functions with low- and high-frequency spacings and compare the responses.

Load data.

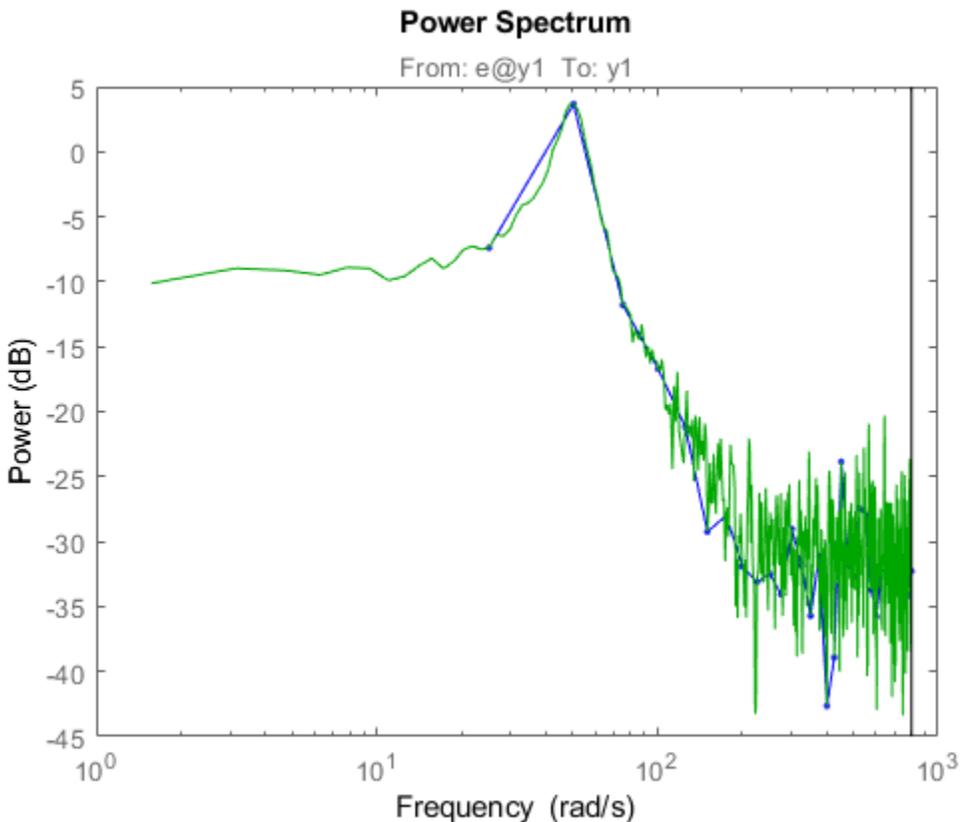
```
load iddata9
```

Estimate empirical transfer functions with low and high frequency spacings.

```
ge1 = etfe(z9, [], 32);  
ge2 = etfe(z9, [], 512);
```

Plot the output power spectrum of the two models.

```
spectrum(ge1, 'b.-', ge2, 'g')
```



## Input Arguments

### **data** — Estimation data

`iddata`

Estimation data, specified as an `iddata` object. The data can be time- or frequency-domain input/output signals or time-series data.

### **M** — Frequency resolution

`[]` (default) | positive scalar

Frequency resolution, specified as a positive scalar.

**N — Frequency spacing**

128 for nonperiodic time-domain data (default) | positive scalar

Frequency spacing, specified as a positive scalar. For frequency-domain data, the default frequency spacing is the spacing inherent in the estimation data.

## Output Arguments

**g — Transfer function estimate**

`idfrd`

Transfer function estimate, returned as an `idfrd` model.

Information about the estimation results and options used is stored in the model's `Report` property. `Report` has the following fields:

Report Field	Description
Status	Summary of the model status, which indicates whether the model was created by construction or obtained by estimation.
Method	Estimation command used.
WindowSize	Size of the Hamming window.

Report Field	Description	
DataUsed	Attributes of the data used for estimation, returned as a structure with the following fields:	
	<b>Field</b> <b>Description</b>	
	Name	Name of the data set.
	Type	Data type.
	Length	Number of data samples.
	Ts	Sample time.
	InterSample	Input intersample behavior, returned as one of the following values: <ul style="list-style-type: none"> <li>'zoh' — Zero-order hold maintains a piecewise-constant input signal between samples.</li> <li>'foh' — First-order hold maintains a piecewise-linear input signal between samples.</li> <li>'bl' — Band-limited behavior specifies that the continuous-time input signal has zero power above the Nyquist frequency.</li> </ul>
	InputOffset	Offset removed from time-domain input data during estimation. For nonlinear models, it is [ ].
OutputOffset	Offset removed from time-domain output data during estimation. For nonlinear models, it is [ ].	

For more information on using Report, see “Estimation Report”.

## See Also

bode | freqresp | idfrd | impulseest | nyquist | spa | spafdr | spectrum

## Topics

“Estimate Frequency-Response Models at the Command Line”

“What is a Frequency-Response Model?”

**Introduced before R2006a**

## evalfr

Evaluate frequency response at given frequency

### Syntax

```
frsp = evalfr(sys,f)
```

### Description

`frsp = evalfr(sys,f)` evaluates the transfer function of the TF, SS, or ZPK model `sys` at the complex number `f`. For state-space models with data  $(A, B, C, D)$ , the result is

$$H(f) = D + C(fI - A)^{-1}B$$

`evalfr` is a simplified version of `freqresp` meant for quick evaluation of the response at a single point. Use `freqresp` to compute the frequency response over a set of frequencies.

## Examples

### Evaluate Discrete-Time Transfer Function

Create the following discrete-time transfer function.

$$H(z) = \frac{z-1}{z^2+z+1}$$

```
H = tf([1 -1],[1 1 1],-1);
```

Evaluate the transfer function at  $z = 1+j$ .

```
z = 1+j;
evalfr(H,z)
```

```
ans = 0.2308 + 0.1538i
```

## Evaluate Frequency Response of Identified Model at Given Frequency

Create the following continuous-time transfer function model:

$$H(s) = \frac{1}{s^2 + 2s + 1}$$

```
sys = idtf(1,[1 2 1]);
```

Evaluate the transfer function at frequency 0.1 rad/second.

```
w = 0.1;  
s = j*w;  
evalfr(sys,s)
```

```
ans = 0.9705 - 0.1961i
```

Alternatively, use the `freqresp` command.

```
freqresp(sys,w)
```

```
ans = 0.9705 - 0.1961i
```

## Limitations

The response is not finite when  $f$  is a pole of `sys`.

## See Also

`bode` | `freqresp` | `sigma`

**Introduced in R2012a**

# extendedKalmanFilter

Create extended Kalman filter object for online state estimation

## Syntax

```
obj = extendedKalmanFilter(StateTransitionFcn,MeasurementFcn,  
InitialState)
```

```
obj = extendedKalmanFilter(StateTransitionFcn,MeasurementFcn,  
InitialState,Name,Value)
```

```
obj = extendedKalmanFilter(StateTransitionFcn,MeasurementFcn)
```

```
obj = extendedKalmanFilter(StateTransitionFcn,MeasurementFcn,  
Name,Value)
```

```
obj = extendedKalmanFilter(Name,Value)
```

## Description

`obj = extendedKalmanFilter(StateTransitionFcn,MeasurementFcn,InitialState)` creates an extended Kalman filter object for online state estimation of a discrete-time nonlinear system. `StateTransitionFcn` is a function that calculates the state of the system at time  $k$ , given the state vector at time  $k-1$ . `MeasurementFcn` is a function that calculates the output measurement of the system at time  $k$ , given the state at time  $k$ . `InitialState` specifies the initial value of the state estimates.

After creating the object, use the `correct` and `predict` commands to update state estimates and state estimation error covariance values using a first-order discrete-time extended Kalman filter algorithm and real-time data.

`obj = extendedKalmanFilter(StateTransitionFcn,MeasurementFcn,InitialState,Name,Value)` specifies additional attributes of the extended Kalman filter object using one or more `Name,Value` pair arguments.

`obj = extendedKalmanFilter(StateTransitionFcn,MeasurementFcn)` creates an extended Kalman filter object using the specified state transition and measurement functions. Before using the `predict` and `correct` commands, specify the initial state

values using dot notation. For example, for a two-state system with initial state values  $[1;0]$ , specify `obj.State = [1;0]`.

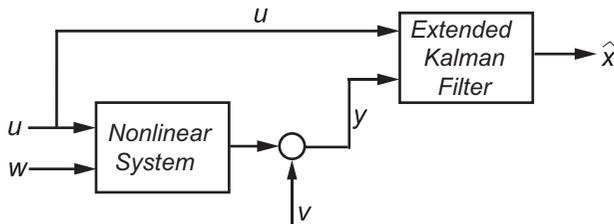
`obj = extendedKalmanFilter(StateTransitionFcn,MeasurementFcn, Name, Value)` specifies additional attributes of the extended Kalman filter object using one or more `Name, Value` pair arguments. Before using the `predict` and `correct` commands, specify the initial state values using `Name, Value` pair arguments or dot notation.

`obj = extendedKalmanFilter(Name, Value)` creates an extended Kalman filter object with properties specified using one or more `Name, Value` pair arguments. Before using the `predict` and `correct` commands, specify the state transition function, measurement function, and initial state values using `Name, Value` pair arguments or dot notation.

## Object Description

`extendedKalmanFilter` creates an object for online state estimation of a discrete-time nonlinear system using the first-order discrete-time extended Kalman filter algorithm.

Consider a plant with states  $x$ , input  $u$ , output  $y$ , process noise  $w$ , and measurement noise  $v$ . Assume that you can represent the plant as a nonlinear system.



The algorithm computes the state estimates  $\hat{x}$  of the nonlinear system using state transition and measurement functions specified by you. The software lets you specify the noise in these functions as additive or nonadditive:

- **Additive Noise Terms** — The state transition and measurements equations have the following form:

$$x[k] = f(x[k-1], u_s[k-1]) + w[k-1]$$

$$y[k] = h(x[k], u_m[k]) + v[k]$$

Here  $f$  is a nonlinear state transition function that describes the evolution of states  $x$  from one time step to the next. The nonlinear measurement function  $h$  relates  $x$  to the measurements  $y$  at time step  $k$ .  $w$  and  $v$  are the zero-mean, uncorrelated process and measurement noises, respectively. These functions can also have additional input arguments that are denoted by  $u_s$  and  $u_m$  in the equations. For example, the additional arguments could be time step  $k$  or the inputs  $u$  to the nonlinear system. There can be multiple such arguments.

Note that the noise terms in both equations are additive. That is,  $x(k)$  is linearly related to the process noise  $w(k-1)$ , and  $y(k)$  is linearly related to the measurement noise  $v(k)$ .

- **Nonadditive Noise Terms** — The software also supports more complex state transition and measurement functions where the state  $x[k]$  and measurement  $y[k]$  are nonlinear functions of the process noise and measurement noise, respectively. When the noise terms are nonadditive, the state transition and measurements equation have the following form:

$$x[k] = f(x[k-1], w[k-1], u_s[k-1])$$

$$y[k] = h(x[k], v[k], u_m[k])$$

When you perform online state estimation, you first create the nonlinear state transition function  $f$  and measurement function  $h$ . You then construct the `extendedKalmanFilter` object using these nonlinear functions, and specify whether the noise terms are additive or nonadditive. You can also specify the Jacobians of the state transition and measurement functions. If you do not specify them, the software numerically computes the Jacobians.

After you create the object, you use the `predict` command to predict state estimate at the next time step, and `correct` to correct state estimates using the algorithm and real-time data. For information about the algorithm, see “Extended and Unscented Kalman Filter Algorithms for Online State Estimation”.

You can use the following commands with `extendedKalmanFilter` objects:

Command	Description
<code>correct</code>	Correct the state and state estimation error covariance at time step $k$ using measured data at time step $k$ .

<b>Command</b>	<b>Description</b>
<code>predict</code>	Predict the state and state estimation error covariance at time the next time step.
<code>clone</code>	Create another object with the same object property values.  Do not create additional objects using syntax <code>obj2 = obj</code> . Any changes made to the properties of the new object created in this way ( <code>obj2</code> ) also change the properties of the original object ( <code>obj</code> ).

For `extendedKalmanFilter` object properties, see “Properties” on page 1-338.

## Examples

### Create Extended Kalman Filter Object for Online State Estimation

To define an extended Kalman filter object for estimating the states of your system, you first write and save the state transition function and measurement function for the system.

In this example, use the previously written and saved state transition and measurement functions, `vdpStateFcn.m` and `vdpMeasurementFcn.m`. These functions describe a discrete-approximation to a van der Pol oscillator with nonlinearity parameter, `mu`, equal to 1. The oscillator has two states.

Specify an initial guess for the two states. You specify the guess as an `M`-element row or column vector, where `M` is the number of states.

```
initialStateGuess = [1;0];
```

Create the extended Kalman filter object. Use function handles to provide the state transition and measurement functions to the object.

```
obj = extendedKalmanFilter(@vdpStateFcn,@vdpMeasurementFcn,initialStateGuess);
```

The object has a default structure where the process and measurement noise are additive.

To estimate the states and state estimation error covariance from the constructed object, use the `correct` and `predict` commands and real-time data.

### Specify Process and Measurement Noise Covariances in Extended Kalman Filter Object

Create an extended Kalman filter object for a van der Pol oscillator with two states and one output. Use the previously written and saved state transition and measurement functions, `vdpStateFcn.m` and `vdpMeasurementFcn.m`. These functions are written for additive process and measurement noise terms. Specify the initial state values for the two states as `[2;0]`.

Since the system has two states and the process noise is additive, the process noise is a 2-element vector and the process noise covariance is a 2-by-2 matrix. Assume there is no cross-correlation between process noise terms, and both the terms have the same variance 0.01. You can specify the process noise covariance as a scalar. The software uses the scalar value to create a 2-by-2 diagonal matrix with 0.01 on the diagonals.

Specify the process noise covariance during object construction.

```
obj = extendedKalmanFilter(@vdpStateFcn,@vdpMeasurementFcn,[2;0],...  
    'ProcessNoise',0.01);
```

Alternatively, you can specify noise covariances after object construction using dot notation. For example, specify the measurement noise covariance as 0.2.

```
obj.MeasurementNoise = 0.2;
```

Since the system has only one output, the measurement noise is a 1-element vector and the `MeasurementNoise` property denotes the variance of the measurement noise.

### Specify Jacobians for State and Measurement Functions

Create an extended Kalman filter object for a van der Pol oscillator with two states and one output. Use the previously written and saved state transition and measurement functions, `vdpStateFcn.m` and `vdpMeasurementFcn.m`. Specify the initial state values for the two states as `[2;0]`.

```
obj = extendedKalmanFilter(@vdpStateFcn,@vdpMeasurementFcn,[2;0]);
```

The extended Kalman filter algorithm uses Jacobians of the state transition and measurement functions for state estimation. You write and save the Jacobian functions and provide them as function handles to the object. In this example, use the previously written and saved functions `vdpStateJacobianFcn.m` and `vdpMeasurementJacobianFcn.m`.

```
obj.StateTransitionJacobianFcn = @vdpStateJacobianFcn;  
obj.MeasurementJacobianFcn = @vdpMeasurementJacobianFcn;
```

Note that if you do not specify the Jacobians of the functions, the software numerically computes the Jacobians. This numerical computation may result in increased processing time and numerical inaccuracy of the state estimation.

### Specify Nonadditive Measurement Noise in Extended Kalman Filter Object

Create an extended Kalman filter object for a van der Pol oscillator with two states and one output. Assume that the process noise terms in the state transition function are additive. That is, there is a linear relation between the state and process noise. Also assume that the measurement noise terms are nonadditive. That is, there is a nonlinear relation between the measurement and measurement noise.

```
obj = extendedKalmanFilter('HasAdditiveMeasurementNoise',false);
```

Specify the state transition function and measurement functions. Use the previously written and saved functions, `vdpStateFcn.m` and `vdpMeasurementNonAdditiveNoiseFcn.m`.

The state transition function is written assuming the process noise is additive. The measurement function is written assuming the measurement noise is nonadditive.

```
obj.StateTransitionFcn = @vdpStateFcn;  
obj.MeasurementFcn = @vdpMeasurementNonAdditiveNoiseFcn;
```

Specify the initial state values for the two states as `[2;0]`.

```
obj.State = [2;0];
```

You can now use the `correct` and `predict` commands to estimate the state and state estimation error covariance values from the constructed object.

### Specify State Transition and Measurement Functions with Additional Inputs

Consider a nonlinear system with input  $u$  whose state  $x$  and measurement  $y$  evolve according to the following state transition and measurement equations:

$$x[k] = \sqrt{x[k-1] + u[k-1]} + w[k-1]$$

$$y[k] = x[k] + 2 * u[k] + v[k]^2$$

The process noise  $w$  of the system is additive while the measurement noise  $v$  is nonadditive.

Create the state transition function and measurement function for the system. Specify the functions with an additional input  $u$ .

```
f = @(x,u)(sqrt(x+u));
h = @(x,v,u)(x+2*u+v^2);
```

$f$  and  $h$  are function handles to the anonymous functions that store the state transition and measurement functions, respectively. In the measurement function, because the measurement noise is nonadditive,  $v$  is also specified as an input. Note that  $v$  is specified as an input before the additional input  $u$ .

Create an extended Kalman filter object for estimating the state of the nonlinear system using the specified functions. Specify the initial value of the state as 1, and the measurement noise as nonadditive.

```
obj = extendedKalmanFilter(f,h,1,'HasAdditiveMeasurementNoise',false);
```

Specify the measurement noise covariance.

```
obj.MeasurementNoise = 0.01;
```

You can now estimate the state of the system using the `predict` and `correct` commands. You pass the values of  $u$  to `predict` and `correct`, which in turn pass them to the state transition and measurement functions, respectively.

Correct the state estimate with measurement  $y[k]=0.8$  and input  $u[k]=0.2$  at time step  $k$ .

```
correct(obj,0.8,0.2)
```

Predict the state at next time step, given  $u[k]=0.2$ .

```
predict(obj,0.2)
```

## Input Arguments

### StateTransitionFcn — State transition function

function handle

State transition function  $f$ , specified as a function handle. The function calculates the  $N_s$ -element state vector of the system at time step  $k$ , given the state vector at time step  $k-1$ .  $N_s$  is the number of states of the nonlinear system.

You write and save the state transition function for your nonlinear system, and use it to construct the object. For example, if `vdpStateFcn.m` is the state transition function, specify `StateTransitionFcn` as `@vdpStateFcn`. You can also specify `StateTransitionFcn` as a function handle to an anonymous function.

The inputs to the function you write depend on whether you specify the process noise as additive or nonadditive in the `HasAdditiveProcessNoise` property of the object:

- `HasAdditiveProcessNoise` is true — The process noise  $w$  is additive, and the state transition function specifies how the states evolve as a function of state values at the previous time step:

$$x(k) = f(x(k-1), Us1, \dots, Usn)$$

Where  $x(k)$  is the estimated state at time  $k$ , and  $Us1, \dots, Usn$  are any additional input arguments required by your state transition function, such as system inputs or the sample time. During estimation, you pass these additional arguments to the `predict` command, which in turn passes them to the state transition function.

- `HasAdditiveProcessNoise` is false — The process noise is nonadditive, and the state transition function also specifies how the states evolve as a function of the process noise:

$$x(k) = f(x(k-1), w(k-1), Us1, \dots, Usn)$$

To see an example of a state transition function with additive process noise, type `edit vdpStateFcn` at the command line.

**MeasurementFcn — Measurement function**

function handle

Measurement function  $h$ , specified as a function handle. The function calculates the  $N$ -element output measurement vector of the nonlinear system at time step  $k$ , given the state vector at time step  $k$ .  $N$  is the number of measurements of the system. You write and save the measurement function, and use it to construct the object. For example, if `vdpMeasurementFcn.m` is the measurement function, specify `MeasurementFcn` as `@vdpMeasurementFcn`. You can also specify `MeasurementFcn` as a function handle to an anonymous function.

The inputs to the function depend on whether you specify the measurement noise as additive or nonadditive in the `HasAdditiveMeasurementNoise` property of the object:

- `HasAdditiveMeasurementNoise` is true — The measurement noise  $v$  is additive, and the measurement function specifies how the measurements evolve as a function of state values:

$$y(k) = h(x(k), Um1, \dots, Umn)$$

Where  $y(k)$  and  $x(k)$  are the estimated output and estimated state at time  $k$ , and  $Um1, \dots, Umn$  are any optional input arguments required by your measurement function. For example, if you are using multiple sensors for tracking an object, an additional input could be the sensor position. During estimation, you pass these additional arguments to the correct command, which in turn passes them to the measurement function.

- `HasAdditiveMeasurementNoise` is false — The measurement noise is nonadditive, and the measurement function also specifies how the output measurement evolves as a function of the measurement noise:

$$y(k) = h(x(k), v(k), Um1, \dots, Umn)$$

To see an example of a measurement function with additive process noise, type `edit vdpMeasurementFcn` at the command line. To see an example of a measurement function with nonadditive process noise, type `edit vdpMeasurementNonAdditiveNoiseFcn`.

**InitialState — Initial state estimate value**

vector

Initial state estimate value, specified as an  $N_s$ -element vector, where  $N_s$  is the number of states in the system. Specify the initial state values based on your knowledge of the system.

The specified value is stored in the `State` property of the object. If you specify `InitialState` as a column vector, then `State` is also a column vector, and the `predict` and `correct` commands return state estimates as a column vector. Otherwise, a row vector is returned.

If you want a filter with single-precision floating-point variables, specify `InitialState` as a single-precision vector variable. For example, for a two-state system with state transition and measurement functions `vdpStateFcn.m` and `vdpMeasurementFcn.m`, create the extended Kalman filter object with initial state estimates `[1;2]` as follows:

```
obj = extendedKalmanFilter(@vdpStateFcn,@vdpMeasurementFcn,single([1;2]))
```

Data Types: `double` | `single`

### Name-Value Pair Arguments

Specify optional comma-separated pairs of `Name`, `Value` arguments. `Name` is the argument name and `Value` is the corresponding value. `Name` must appear inside quotes. You can specify several name and value pair arguments in any order as `Name1,Value1,...,NameN,ValueN`.

Use `Name`, `Value` arguments to specify properties on page 1-338 of `extendedKalmanFilter` object during object creation. For example, to create an extended Kalman filter object and specify the process noise covariance as 0.01:

```
obj = extendedKalmanFilter(StateTransitionFcn,MeasurementFcn,InitialState,'ProcessNoise',0.01);
```

### Properties

`extendedKalmanFilter` object properties are of three types:

- Tunable properties that you can specify multiple times, either during object construction using `Name`, `Value` arguments, or any time afterward during state estimation. After object creation, use dot notation to modify the tunable properties.

```
obj = extendedKalmanFilter(StateTransitionFcn,MeasurementFcn,InitialState);  
obj.ProcessNoise = 0.01;
```

The tunable properties are `State`, `StateCovariance`, `ProcessNoise`, and `MeasurementNoise`.

- Nontunable properties that you can specify once, either during object construction or afterward using dot notation. Specify these properties before state estimation using `correct` and `predict`. The `StateTransitionFcn`, `MeasurementFcn`, `StateTransitionJacobianFcn`, and `MeasurementJacobianFcn` properties belong to this category.
- Nontunable properties that you must specify during object construction. The `HasAdditiveProcessNoise` and `HasAdditiveMeasurementNoise` properties belong to this category.

### **HasAdditiveMeasurementNoise — Measurement noise characteristics**

`true` (default) | `false`

Measurement noise characteristics, specified as one of the following values:

- `true` — Measurement noise  $v$  is additive. The measurement function  $h$  that is specified in `MeasurementFcn` has the following form:

$$y(k) = h(x(k), Um1, \dots, Umn)$$

Where  $y(k)$  and  $x(k)$  are the estimated output and estimated state at time  $k$ , and  $Um1, \dots, Umn$  are any optional input arguments required by your measurement function.

- `false` — Measurement noise is nonadditive. The measurement function specifies how the output measurement evolves as a function of the state *and* measurement noise:

$$y(k) = h(x(k), v(k), Um1, \dots, Umn)$$

`HasAdditiveMeasurementNoise` is a nontunable property, and you can specify it only during object construction. You cannot change it using dot notation.

### **HasAdditiveProcessNoise — Process noise characteristics**

`true` (default) | `false`

Process noise characteristics, specified as one of the following values:

- `true` — Process noise  $w$  is additive. The state transition function  $f$  specified in `StateTransitionFcn` has the following form:

$$x(k) = f(x(k-1), Us1, \dots, Usn)$$

Where  $x(k)$  is the estimated state at time  $k$ , and  $Us1, \dots, Usn$  are any additional input arguments required by your state transition function.

- `false` — Process noise is nonadditive. The state transition function specifies how the states evolve as a function of the state *and* process noise at the previous time step:

$$x(k) = f(x(k-1), w(k-1), Us1, \dots, Usn)$$

`HasAdditiveProcessNoise` is a nontunable property, and you can specify it only during object construction. You cannot change it using dot notation.

### **MeasurementFcn — Measurement function**

function handle

Measurement function  $h$ , specified as a function handle. The function calculates the  $N$ -element output measurement vector of the nonlinear system at time step  $k$ , given the state vector at time step  $k$ .  $N$  is the number of measurements of the system. You write and save the measurement function and use it to construct the object. For example, if `vdpMeasurementFcn.m` is the measurement function, specify `MeasurementFcn` as `@vdpMeasurementFcn`. You can also specify `MeasurementFcn` as a function handle to an anonymous function.

The inputs to the function depend on whether you specify the measurement noise as additive or nonadditive in the `HasAdditiveMeasurementNoise` property of the object:

- `HasAdditiveMeasurementNoise` is true — The measurement noise  $v$  is additive, and the measurement function specifies how the measurements evolve as a function of state values:

$$y(k) = h(x(k), Um1, \dots, Umn)$$

Where  $y(k)$  and  $x(k)$  are the estimated output and estimated state at time  $k$ , and  $Um1, \dots, Umn$  are any optional input arguments required by your measurement function. For example, if you are using multiple sensors for tracking an object, an additional input could be the sensor position. During estimation, you pass these additional arguments to the `correct` command which in turn passes them to the measurement function.

- `HasAdditiveMeasurementNoise` is false — The measurement noise is nonadditive, and the measurement function also specifies how the output measurement evolves as a function of the measurement noise:

$$y(k) = h(x(k), v(k), Um1, \dots, Umn)$$

To see an example of a measurement function with additive process noise, type `edit vdpMeasurementFcn` at the command line. To see an example of a measurement function with nonadditive process noise, type `edit vdpMeasurementNonAdditiveNoiseFcn`.

`MeasurementFcn` is a nontunable property. You can specify it once before using the `correct` command either during object construction or using dot notation after object construction. You cannot change it after using the `correct` command.

### **MeasurementJacobianFcn — Jacobian of measurement function**

`[]` (default) | function handle

Jacobian of measurement function  $h$ , specified as one of the following:

- `[]` — The Jacobian is numerically computed at every call to the `correct` command. This may increase processing time and numerical inaccuracy of the state estimation.
- function handle — You write and save the Jacobian function and specify the handle to the function. For example, if `vdpMeasurementJacobianFcn.m` is the Jacobian function, specify `MeasurementJacobianFcn` as `@vdpMeasurementJacobianFcn`.

The function calculates the partial derivatives of the measurement function with respect to the states and measurement noise. The number of inputs to the Jacobian function must equal the number of inputs to the measurement function and must be specified in the same order in both functions. The number of outputs of the Jacobian function depends on the `HasAdditiveMeasurementNoise` property:

- `HasAdditiveMeasurementNoise` is true — The function calculates the partial derivatives of the measurement function with respect to the states ( $\partial h / \partial x$ ). The output is as an  $N$ -by- $N_s$  Jacobian matrix, where  $N$  is the number of measurements of the system and  $N_s$  is the number of states.
- `HasAdditiveMeasurementNoise` is false — The function also returns a second output that is the partial derivative of the measurement function with respect to the measurement noise terms ( $\partial h / \partial v$ ). The second output is returned as an  $N$ -by- $V$  Jacobian matrix, where  $V$  is the number of measurement noise terms.

To see an example of a Jacobian function for additive measurement noise, type `edit vdpMeasurementJacobianFcn` at the command line.

`MeasurementJacobianFcn` is a nontunable property. You can specify it once before using the `correct` command either during object construction or using dot notation after object construction. You cannot change it after using the `correct` command.

## MeasurementNoise — Measurement noise covariance

1 (default) | scalar | matrix

Measurement noise covariance, specified as a scalar or matrix depending on the value of the `HasAdditiveMeasurementNoise` property:

- `HasAdditiveMeasurementNoise` is true — Specify the covariance as a scalar or an  $N$ -by- $N$  matrix, where  $N$  is the number of measurements of the system. Specify a scalar if there is no cross-correlation between measurement noise terms and all the terms have the same variance. The software uses the scalar value to create an  $N$ -by- $N$  diagonal matrix.
- `HasAdditiveMeasurementNoise` is false — Specify the covariance as a  $V$ -by- $V$  matrix, where  $V$  is the number of measurement noise terms. `MeasurementNoise` must be specified before using `correct`. After you specify `MeasurementNoise` as a matrix for the first time, to then change `MeasurementNoise` you can also specify it as a scalar. Specify as a scalar if there is no cross-correlation between the measurement noise terms and all the terms have the same variance. The software extends the scalar to a  $V$ -by- $V$  diagonal matrix with the scalar on the diagonals.

`MeasurementNoise` is a tunable property. You can change it using dot notation.

## ProcessNoise — Process noise covariance

1 (default) | scalar | matrix

Process noise covariance, specified as a scalar or matrix depending on the value of the `HasAdditiveProcessNoise` property:

- `HasAdditiveProcessNoise` is true — Specify the covariance as a scalar or an  $N_s$ -by- $N_s$  matrix, where  $N_s$  is the number of states of the system. Specify a scalar if there is no cross-correlation between process noise terms, and all the terms have the same variance. The software uses the scalar value to create an  $N_s$ -by- $N_s$  diagonal matrix.
- `HasAdditiveProcessNoise` is false — Specify the covariance as a  $W$ -by- $W$  matrix, where  $W$  is the number of process noise terms. `ProcessNoise` must be specified before using `predict`. After you specify `ProcessNoise` as a matrix for the first time, to then change `ProcessNoise` you can also specify it as a scalar. Specify as a scalar if there is no cross-correlation between the process noise terms and all the terms have the same variance. The software extends the scalar to a  $W$ -by- $W$  diagonal matrix.

`ProcessNoise` is a tunable property. You can change it using dot notation.

## State — State of nonlinear system

[] (default) | vector

State of the nonlinear system, specified as a vector of size  $N_s$ , where  $N_s$  is the number of states of the system.

When you use the `predict` command, `State` is updated with the predicted value at time step  $k$  using the state value at time step  $k-1$ . When you use the `correct` command, `State` is updated with the estimated value at time step  $k$  using measured data at time step  $k$ .

The initial value of `State` is the value you specify in the `InitialState` input argument during object creation. If you specify `InitialState` as a column vector, then `State` is also a column vector, and the `predict` and `correct` commands return state estimates as a column vector. Otherwise, a row vector is returned. If you want a filter with single-precision floating-point variables, you must specify `State` as a single-precision variable during object construction using the `InitialState` input argument.

`State` is a tunable property. You can change it using dot notation.

### **StateCovariance — State estimation error covariance**

1 (default) | scalar | matrix

State estimation error covariance, specified as a scalar or an  $N_s$ -by- $N_s$  matrix, where  $N_s$  is the number of states of the system. If you specify a scalar, the software uses the scalar value to create an  $N_s$ -by- $N_s$  diagonal matrix.

Specify a high value for the covariance when you do not have confidence in the initial state values that you specify in the `InitialState` input argument.

When you use the `predict` command, `StateCovariance` is updated with the predicted value at time step  $k$  using the state value at time step  $k-1$ . When you use the `correct` command, `StateCovariance` is updated with the estimated value at time step  $k$  using measured data at time step  $k$ .

`StateCovariance` is a tunable property. You can change it using dot notation after using the `correct` or `predict` commands.

### **StateTransitionFcn — State transition function**

function handle

State transition function  $f$ , specified as a function handle. The function calculates the  $N_s$ -element state vector of the system at time step  $k$ , given the state vector at time step  $k-1$ .  $N_s$  is the number of states of the nonlinear system.

You write and save the state transition function for your nonlinear system and use it to construct the object. For example, if `vdpStateFcn.m` is the state transition function, specify `StateTransitionFcn` as `@vdpStateFcn`. You can also specify `StateTransitionFcn` as a function handle to an anonymous function.

The inputs to the function you write depend on whether you specify the process noise as additive or nonadditive in the `HasAdditiveProcessNoise` property of the object:

- `HasAdditiveProcessNoise` is true — The process noise  $w$  is additive, and the state transition function specifies how the states evolve as a function of state values at previous time step:

$$x(k) = f(x(k-1), Us1, \dots, Usn)$$

Where  $x(k)$  is the estimated state at time  $k$ , and  $Us1, \dots, Usn$  are any additional input arguments required by your state transition function, such as system inputs or the sample time. During estimation, you pass these additional arguments to the `predict` command, which in turn passes them to the state transition function.

- `HasAdditiveProcessNoise` is false — The process noise is nonadditive, and the state transition function also specifies how the states evolve as a function of the process noise:

$$x(k) = f(x(k-1), w(k-1), Us1, \dots, Usn)$$

To see an example of a state transition function with additive process noise, type `edit vdpStateFcn` at the command line.

`StateTransitionFcn` is a nontunable property. You can specify it once before using the `predict` command either during object construction or using dot notation after object construction. You cannot change it after using the `predict` command.

### **StateTransitionJacobianFcn — Jacobian of state transition function**

`[]` (default) | function handle

Jacobian of state transition function  $f$ , specified as one of the following:

- `[]` — The Jacobian is numerically computed at every call to the `predict` command. This may increase processing time and numerical inaccuracy of the state estimation.
- function handle — You write and save the Jacobian function and specify the handle to the function. For example, if `vdpStateJacobianFcn.m` is the Jacobian function, specify `StateTransitionJacobianFcn` as `@vdpStateJacobianFcn`.

The function calculates the partial derivatives of the state transition function with respect to the states and process noise. The number of inputs to the Jacobian function must equal the number of inputs of the state transition function and must be specified in the same order in both functions. The number of outputs of the function depends on the `HasAdditiveProcessNoise` property:

- `HasAdditiveProcessNoise` is true — The function calculates the partial derivative of the state transition function with respect to the states ( $\partial f / \partial x$ ). The output is an  $N_s$ -by- $N_s$  Jacobian matrix, where  $N_s$  is the number of states.
- `HasAdditiveProcessNoise` is false — The function must also return a second output that is the partial derivative of the state transition function with respect to the process noise terms ( $\partial f / \partial w$ ). The second output is returned as an  $N_s$ -by- $W$  Jacobian matrix, where  $W$  is the number of process noise terms.

The extended Kalman filter algorithm uses the Jacobian to compute the state estimation error covariance.

To see an example of a Jacobian function for additive process noise, type `edit vdpStateJacobianFcn` at the command line.

`StateTransitionJacobianFcn` is a nontunable property. You can specify it once before using the `predict` command either during object construction or using dot notation after object construction. You cannot change it after using the `predict` command.

## Output Arguments

### **obj** — Extended Kalman filter object for online state estimation

`extendedKalmanFilter` object

Extended Kalman filter object for online state estimation, returned as an `extendedKalmanFilter` object. This object is created using the specified properties on page 1-338. Use the `correct` and `predict` commands to estimate the state and state estimation error covariance using the extended Kalman filter algorithm.

When you use `predict`, `obj.State` and `obj.StateCovariance` are updated with the predicted value at time step  $k$  using the state value at time step  $k-1$ . When you use `correct`, `obj.State` and `obj.StateCovariance` are updated with the estimated values at time step  $k$  using measured data at time step  $k$ .

## Extended Capabilities

### C/C++ Code Generation

Generate C and C++ code using MATLAB® Coder™.

Usage notes and limitations:

For more information, see “Generate Code for Online State Estimation in MATLAB”.

Supports MATLAB Function block: No

### See Also

#### Functions

`clone` | `correct` | `predict` | `unscentedKalmanFilter`

#### Blocks

Extended Kalman Filter | Kalman Filter | Unscented Kalman Filter

#### Topics

“Nonlinear State Estimation Using Unscented Kalman Filter and Particle Filter”

“Generate Code for Online State Estimation in MATLAB”

“What Is Online Estimation?”

“Extended and Unscented Kalman Filter Algorithms for Online State Estimation”

“Validate Online State Estimation at the Command Line”

“Troubleshoot Online State Estimation”

#### Introduced in R2016b

# evaluate

Value of nonlinearity estimator at given input

## Syntax

```
value = evaluate(nl,x)
```

## Arguments

`nl`

Nonlinearity estimator object.

`x`

Value at which to evaluate the nonlinearity.

If `nl` is a single nonlinearity estimator, then `x` is a 1-by-`nx` row vector or an `nv`-by-`nx` matrix, where `nx` is the dimension of the regression vector input to `nl` (`size(nl)`) and `nv` is the number of points where `nl` is evaluated.

If `nl` is an array of `ny` nonlinearity estimators, then `x` is a 1-by-`ny` cell array of `nv`-by-`nx` matrices.

## Description

`value = evaluate(nl,x)` computes the value of a nonlinear estimator object of type `customnet`, `deadzone`, `linear`, `neuralnet`, `pwnlinear`, `saturation`, `sigmoidnet`, `treepartition`, or `wavenet`.

## Examples

The following syntax evaluates the nonlinearity of an estimated nonlinear ARX model `m`:

```
value = evaluate(m.Nonlinearity,x)
```

where `m.Nonlinearity` accesses the nonlinearity estimator of the nonlinear ARX model.

## **See Also**

`idnlarx` | `idnlhw`

**Introduced in R2007a**

## fcat

Concatenate FRD models along frequency dimension

### Syntax

```
sys = fcat(sys1,sys2,...)
```

### Description

`sys = fcat(sys1,sys2,...)` takes two or more `frd` models and merges their frequency responses into a single `frd` model `sys`. The resulting frequency vector is sorted by increasing frequency. The frequency vectors of `sys1`, `sys2`, ... should not intersect. If the frequency vectors do intersect, use `fdel` to remove intersecting data from one or more of the models.

### See Also

`fdel` | `frd` | `fselect` | `idfrd` | `interp`

**Introduced before R2006a**

## **fdel**

Delete specified data from frequency response data (FRD) models

### **Syntax**

```
sysout = fdel(sys, freq)
```

### **Description**

*sysout* = fdel(*sys*, *freq*) removes from the frd model *sys* the data nearest to the frequency values specified in the vector *freq*.

### **Input Arguments**

**sys**

frd model.

**freq**

Vector of frequency values.

### **Output Arguments**

**sysout**

frd model containing the data remaining in *sys* after removing the frequency points closest to the entries of *freq*.

### **Examples**

## Delete Specified Data from Frequency Response Data Model

Create a frequency response data (FRD) model at specified frequencies from a transfer function model.

```
w = logspace(0,1,10);
sys = frd(tf([1],[1 1]),w)
```

```
sys =
```

Frequency(rad/s)	Response
-----	-----
1.0000	0.5000 - 0.5000i
1.2915	0.3748 - 0.4841i
1.6681	0.2644 - 0.4410i
2.1544	0.1773 - 0.3819i
2.7826	0.1144 - 0.3183i
3.5938	0.0719 - 0.2583i
4.6416	0.0444 - 0.2059i
5.9948	0.0271 - 0.1623i
7.7426	0.0164 - 0.1270i
10.0000	0.0099 - 0.0990i

Continuous-time frequency response.

w is a logarithmically-spaced grid of 10 frequency points between 1 and 10 rad/second.

Remove the data nearest 2, 3.5, and 6 rad/s from sys.

```
freq = [2, 3.5, 6];
sys2 = fdel(sys,freq)
```

```
sys2 =
```

Frequency(rad/s)	Response
-----	-----
1.0000	0.5000 - 0.5000i
1.2915	0.3748 - 0.4841i
1.6681	0.2644 - 0.4410i
2.7826	0.1144 - 0.3183i
4.6416	0.0444 - 0.2059i
7.7426	0.0164 - 0.1270i
10.0000	0.0099 - 0.0990i

Continuous-time frequency response.

Note that you do not have to specify the exact frequency of the data to remove. The `fdel` command removes the data corresponding to frequencies that are nearest to the specified frequencies.

### Tips

- Use `fdel` to remove unwanted data (for example, outlier points) at specified frequencies.
- Use `fdel` to remove data at intersecting frequencies from `frd` models before merging them with `fcats`. `fcats` produces an error when you attempt to merge `frd` models that have intersecting frequency data.
- To remove data from an `frd` model within a range of frequencies, use `fselect`.

### See Also

`fcats` | `frd` | `fselect` | `idfrd`

**Introduced in R2012a**

# feedback

Identify possible feedback data

## Syntax

```
[fbck,fbck0,nudir] = feedback(Data)
```

## Description

Data is an `iddata` set with  $N_y$  outputs and  $N_u$  inputs.

`fbck` is an  $N_y$ -by- $N_u$  matrix indicating the feedback. The  $ky,ku$  entry is a measure of feedback from output  $ky$  to input  $ku$ . The value is a probability  $P$  in percent. Its interpretation is that if the hypothesis that there is no feedback from output  $ky$  to input  $ku$  were tested at the level  $P$ , it would have been rejected. An intuitive but technically incorrect way of thinking about this is to see  $P$  as “the probability of feedback.” Often only values above 90% are taken as indications of feedback. When `fbck` is calculated, direct dependence at lag zero between  $u(t)$  and  $y(t)$  is not regarded as a feedback effect.

`fbck0`: Same as `fbck`, but direct dependence at lag 0 between  $u(t)$  and  $y(t)$  is viewed as feedback effect.

`nudir`: A vector containing those input numbers that appear to have a direct effect on some outputs, that is, no delay from input to output.

## See Also

`advice` | `iddata`

**Introduced before R2006a**

## fft

Transform `iddata` object to frequency domain data

### Syntax

```
Datf = fft(Data)
Datf = fft(Data,N)
Datf = fft(Data,N,'complex')
```

### Description

`Datf = fft(Data)` transforms time-domain data to frequency domain data. If `Data` is a time-domain `iddata` object with real-valued signals and with constant sample time `Ts`, `Datf` is returned as a frequency-domain `iddata` object with the frequency values equally distributed from frequency 0 to the Nyquist frequency. Whether the Nyquist frequency actually is included or not depends on the signal length (even or odd). Note that the FFTs are normalized by dividing each transform by the square root of the signal length. That is in order to preserve the signal power and noise level.

`Datf = fft(Data,N)` specifies the transformation length. In the default case, the length of the transformation is determined by the signal length. A second argument `N` will force FFT transformations of length `N`, padding with zeros if the signals in `Data` are shorter and truncating otherwise. Thus the number of frequencies in the real signal case will be  $N/2$  or  $(N+1)/2$ . If `Data` contains several experiments, `N` can be a row vector of corresponding length.

`Datf = fft(Data,N,'complex')` specifies to include negative frequencies. For real signals, the default is that `Datf` only contains nonnegative frequencies. For complex-valued signals, negative frequencies are also included. To enforce negative frequencies in the real case, add a last argument, `'Complex'`.

### See Also

`iddata` | `ifft` | `spa`

**Introduced in R2007a**

## idnlarx/findop

Compute operating point for Nonlinear ARX model

### Syntax

```
[X,U] = findop(sys,'steady',InputLevel,OutputLevel)
```

```
[X,U] = findop(sys,spec)
```

```
[X,U] = findop( ____,Options)
```

```
[X,U,Report] = findop( ____,Options)
```

```
[X,U] = findop(sys,'snapshot',T,Uin)
```

```
[X,U] = findop(sys,'snapshot',T,Uin,X0)
```

### Description

`[X,U] = findop(sys,'steady',InputLevel,OutputLevel)` returns the operating-point state values, `X`, and input values, `U`, for the `idnlarx` model, `sys`, using steady-state input and output specifications.

`[X,U] = findop(sys,spec)` returns the steady-state operating point for `sys` using the operating-point specification, `spec`.

`[X,U] = findop( ____,Options)` specifies optimization search options for all of the previous syntaxes.

`[X,U,Report] = findop( ____,Options)` returns a summary report on the optimization search results for all of the previous syntaxes.

`[X,U] = findop(sys,'snapshot',T,Uin)` returns the operating point for `sys` at a simulation snapshot at time, `T`, using the specified input, `Uin`. The initial states of `sys` are assumed to be zero.

`[X,U] = findop(sys,'snapshot',T,Uin,X0)` specifies the initial states of the simulation.

## Examples

### Find Steady-State Nonlinear ARX Operating Point Using Default Specifications

Estimate a nonlinear ARX model.

```
load iddata6;  
M = nlarx(z6,[4 3 1]);
```

Find the steady-state operating point where the input level is fixed to 1 and the output is unknown.

```
[X,U] = findop(M,'steady',1,NaN);
```

### Find Nonlinear ARX Operating Point Using Additional Specifications

Estimate a nonlinear ARX model.

```
load iddata7;  
M = nlarx(z7,[4 3*ones(1,2) 2*ones(1,2)]);
```

Create a default operating point specification object.

```
spec = operspec(M);
```

Set the values for the input signals.

```
spec.Input.Value(1) = -1;  
spec.Input.Value(2) = 1;
```

Set the maximum and minimum values for the output signal.

```
spec.Output.Max = 10;  
spec.Output.Min = -10;
```

Find the steady-state operating point using the given specifications.

```
[X,U] = findop(M,spec);
```

### Find Nonlinear ARX Operating Point Using Custom Options

Estimate a nonlinear ARX model.

```
load iddata6;  
M = nlarx(z6,[4 3 2]);
```

Create a default findopOptions option set.

```
opt = findopOptions(M);
```

Modify the option set to specify a steepest descent gradient search method with a maximum of 50 iterations.

```
opt.SearchMethod = 'grad';  
opt.SearchOptions.MaxIterations = 50;
```

Find the steady-state operating point using the specified options.

```
[X,U] = findop(M,'steady',1,1,opt);
```

### Retrieve Nonlinear ARX Operating Point Search Report

Estimate a nonlinear ARX model.

```
load iddata7;  
M = nlarx(z7,[4 3*ones(1,2) 2*ones(1,2)]);
```

Find the steady-state operating point where input 1 is set to 1 and input 2 is unrestricted. The initial guess for the output value is 2.

```
[X,U,R] = findop(M,'steady',[1 NaN],2);
```

Display the summary report.

```
disp(R);
```

```
SearchMethod: 'auto'  
WhyStop: 'Near (local) minimum, (norm(g) < tol).'Iterations: 11  
FinalCost: 0  
FirstOrderOptimality: 0  
SignalLevels: [1x1 struct]
```

## Find Nonlinear ARX Simulation Snapshot Using Default Initial States

Load the estimation data and estimate a nonlinear ARX model.

```
load twotankdata;  
z = iddata(y,u,1);  
M = nlarx(z,[4 3 1]);
```

Find the simulation snapshot after 10 seconds, assuming initial states of zero.

```
[X,U] = findop(M,'snapshot',10,z);
```

## Find Nonlinear ARX Simulation Snapshot Using Initial State Specifications

Load the estimation data and estimate a nonlinear ARX model.

```
load twotankdata;  
z = iddata(y,u,1);  
M = nlarx(z,[4 3 1]);
```

Create an initial state vector. The first four states correspond to delayed output values and the final three states correspond to delayed inputs.

```
X0 = [2;2;2;2;5;5;5];
```

Find the simulation snapshot after 10 seconds using the specified initial states.

```
[X,U] = findop(M,'snapshot',10,z,X0);
```

## Input Arguments

### sys — Nonlinear ARX model

idnlarx object

Nonlinear ARX model, specified as an idnlarx object.

**InputLevel — Steady-state input level**

vector

Steady-state input level for computing the operating point, specified as a vector. The length of `InputLevel` must equal the number of inputs specified in `sys`.

The optimization algorithm assumes that finite values in `InputLevel` are fixed input values. Use `NaN` to specify unknown input signals with initial guesses of `0`. The minimum and maximum bounds for all inputs have default values of `-Inf` and `+Inf` respectively.

**OutputLevel — Steady-state output level**

vector

Steady-state output level for computing the operating point, specified as a vector. The length of `OutputLevel` must equal the number of outputs specified in `sys`.

The values in `OutputLevel` indicate initial guesses for the optimization algorithm. Use `NaN` to specify unknown output signals with initial guesses of `0`. The minimum and maximum bounds for all outputs have default values of `-Inf` and `+Inf` respectively.

**spec — Operating-point specifications**

operspec object

Operating-point specifications, such as minimum and maximum input/output constraints and known inputs, specified as an `operspec` object.

**T — Operating point snapshot time**

positive scalar

Operating point snapshot time, specified as a positive scalar. The value of `T` must be in the range  $[T_0, N \cdot T_s]$ , where  $N$  is the number of input samples,  $T_s$  is the sample time and  $T_0$  is the input start time (`Uin.Tstart`).

**Uin — Snapshot simulation input**

iddata object | matrix

Snapshot simulation input, specified as one of the following:

- Time-domain `iddata` object with a sample time and input size that matches `sys`.
- Matrix with as many columns as there are input channels. If the matrix has  $N$  rows, the input data is assumed to correspond to the time vector  $(1:N) \cdot \text{sys.Ts}$ .

**X0 — Initial states**

column vector

Initial states of the simulation, specified as a column vector with size equal to the number of states in `sys`. `X0` provides the initial conditions at the time corresponding to the first input sample (`Uin.Start`, if `Uin` is an `iddata` object, or `sys.Ts` if `Uin` is a double matrix).

For more information about the states of an `idnlarx` model, see “Definition of `idnlarx` States” on page 1-631.

**Options — Operating point search options**`findopOptions` option set

Operating point search options, specified as a `findopOptions` option set.

## Output Arguments

**X — Operating point state values**

column vector

Operating point state values, returned as a column vector of length equal to the number of model states.

**U — Operating point input values**

column vector

Operating point input values, returned as a column vector of length equal to the number of inputs.

**Report — Search result summary**

structure

Search result summary report, returned as a structure with the following fields:

Field	Description
<code>SearchMethod</code>	Search method used for iterative parameter estimation. See <code>SearchMethod</code> in <code>findopOptions</code> for more information.
<code>WhyStop</code>	Search algorithm termination condition.

Field	Description
Iterations	Number of estimation iterations performed.
FinalCost	Final value of the minimization objective function (sum of the squared errors).
FirstOrderOptimality	$\infty$ -norm of the search gradient vector when the search algorithm terminates.
SignalLevels	Structure containing the fields <code>Input</code> and <code>Output</code> , which are the operating point input and output signal levels respectively.

## Algorithms

`findop` computes the operating point from steady-state operating point specifications or at a simulation snapshot.

### Computing the Operating Point from Steady-State Specifications

To compute the steady-state operating point, call `findop` using either of the following syntaxes:

```
[X,U] = findop(sys,'steady',InputLevel,OutputLevel)
[X,U] = findop(sys,spec)
```

To compute the states, `X`, and the input, `U`, of the steady-state operating point, `findop` minimizes the norm of the error  $e(t) = y(t) - f(x(t), u(t))$ , where:

- $f$  is the nonlinearity estimator.
- $u(t)$  is the input.
- $x(t)$  is the model state.
- $y(t)$  is the model output.

You can specify the search algorithm and search options using the `findopOptions` option set.

The algorithm uses the following independent variables for minimization:

- Unknown (unspecified) input signal levels
- Output signal levels

Because `idnlarx` model states are delayed samples of the input and output variables, the state values are the constant values of the corresponding steady-state inputs and outputs. For more information about the definition of nonlinear ARX model states, see “Definition of `idnlarx` States” on page 1-631.

## Computing the Operating Point at a Simulation Snapshot

When you use the syntax `[X,U] = findop(sys,'snapshot',T,Uin,X0)`, the algorithm simulates the model output until the snapshot time, `T`. At the snapshot time, the algorithm passes the input and output samples to the `data2state` command to map these values to the current state vector.

---

**Note** For snapshot-based computations, `findop` does not perform numerical optimization.

---

## Extended Capabilities

### Automatic Parallel Support

Accelerate code by automatically running computation in parallel using Parallel Computing Toolbox™.

Parallel computing support is available for estimation using the `lsqnonlin` search method (requires Optimization Toolbox). To enable parallel computing, use `findopOptions`, set `SearchMethod` to `'lsqnonlin'`, and set `SearchOptions.Advanced.UseParallel` to `true`.

For example:

```
opt = findopOptions(idnlarx);  
opt.SearchMethod = 'lsqnonlin';  
opt.SearchOptions.Advanced.UseParallel = true;
```

## **See Also**

`data2state` | `findopOptions` | `idnlarx` | `idnlarx/operspec` | `idnlhw/findop` | `sim`

**Introduced in R2008a**

# idnlhw/findop

Compute operating point for Hammerstein-Wiener model

## Syntax

```
[X,U] = findop(sys,'steady',InputLevel,OutputLevel)
```

```
[X,U] = findop(sys,spec)
```

```
[X,U] = findop( ____,Options)
```

```
[X,U,Report] = findop( ____,Options)
```

```
[X,U] = findop(sys,'snapshot',T,Uin)
```

```
[X,U] = findop(sys,'snapshot',T,Uin,X0)
```

## Description

`[X,U] = findop(sys,'steady',InputLevel,OutputLevel)` returns the operating-point state values, `X`, and input values, `U`, for the `idnlhw` model, `sys`, using steady-state input and output specifications.

`[X,U] = findop(sys,spec)` returns the steady-state operating point for `sys` using the operating point specification in `spec`.

`[X,U] = findop( ____,Options)` specifies optimization search options for all of the previous syntaxes.

`[X,U,Report] = findop( ____,Options)` returns a summary report on the optimization search results for all of the previous syntaxes.

`[X,U] = findop(sys,'snapshot',T,Uin)` returns the operating point for `sys` at a simulation snapshot at time, `T`, using the specified input, `Uin`. The initial states of `sys` are assumed to be zero.

`[X,U] = findop(sys,'snapshot',T,Uin,X0)` specifies the initial states of the simulation.

## Examples

### Find Steady-State Hammerstein-Wiener Operating Point Using Default Specifications

Load the estimation data and estimate a Hammerstein-Wiener model.

```
load twotankdata;  
z = iddata(y,u,1);  
M = nlhw(z,[5 1 3]);
```

Find the steady-state operating point where the input level is set to 1 and the output is unknown.

```
[X,U] = findop(M,'steady',1,NaN);
```

### Find Hammerstein-Wiener Operating Point Using Additional Specifications

Estimate a Hammerstein-Wiener model.

```
load iddata7;  
orders = [4*ones(1,2) 2*ones(1,2) 3*ones(1,2)];  
M = nlhw(z7,orders,'unitgain','pwnlinear');
```

Create a default operating point specification object.

```
spec = operspec(M);
```

Set the values for the input signals.

```
spec.Input.Value(1) = -1;  
spec.Input.Value(2) = 1;
```

Set the maximum and minimum values for the output signal.

```
spec.Output.Max = 10;  
spec.Output.Min = -10;
```

Find the steady-state operating point using the given specifications.

```
[X,U] = findop(M,spec);
```

## Find Hammerstein-Wiener Operating Point Using Custom Options

Load the estimation data and estimate a Hammerstein-Wiener model.

```
load twotankdata;
z = iddata(y,u,1);
M = nlhw(z,[5 1 3]);
```

Create a default findopOptions option set.

```
opt = findopOptions(M);
```

Modify the option set to specify a steepest descent gradient search method with a maximum of 50 iterations.

```
opt.SearchMethod = 'grad';
opt.SearchOptions.MaxIterations = 50;
```

Find the steady-state operating point using the specified options.

```
[X,U] = findop(M,'steady',1,NaN,opt);
```

## Retrieve Hammerstein-Wiener Operating Point Search Report

Load the estimation data and estimate a Hammerstein-Wiener model.

```
load iddata7;
orders = [4*ones(1,2) 2*ones(1,2) 3*ones(1,2)];
M = nlhw(z7,orders,'unitgain','pwlinear');
```

Find the steady-state operating point where input 1 is set to 1 and input 2 is unrestricted. The initial guess for the output value is 2.

```
[X,U,R] = findop(M,'steady',[1 NaN],2);
```

Display the summary report.

```
disp(R);
```

```
SearchMethod: 'auto'
WhyStop: 'Near (local) minimum, (norm(g) < tol).'
```

```
Iterations: 2
FinalCost: 0
FirstOrderOptimality: 0
SignalLevels: [1x1 struct]
```

## Find Hammerstein-Wiener Simulation Snapshot Using Default Initial States

Load the estimation data estimate a Hammerstein-Wiener model.

```
load twotankdata;
z = iddata(y,u,1);
M = nllhw(z,[5 1 3]);
```

Find the simulation snapshot after 10 seconds, assuming initial states of zero.

```
[X,U] = findop(M,'snapshot',10,z);
```

## Find Hammerstein-Wiener Simulation Snapshot Using Initial State Specifications

Load the estimation data and estimate a Hammerstein-Wiener model.

```
load twotankdata
z = iddata(y,u,1);
M = nllhw(z,[5 1 3]);
```

Create an initial state vector.

```
X0 = [10;10;5;5;1;1;0];
```

Find the simulation snapshot after 10 seconds using the specified initial states.

```
[X,U] = findop(M,'snapshot',10,z,X0);
```

## Input Arguments

**sys** — Hammerstein-Wiener model

idnllhw object

Hammerstein-Wiener model, specified as an `idnlhw` object.

### **InputLevel — Steady-state input level**

vector

Steady-state input level for computing the operating point, specified as a vector. The length of `InputLevel` must equal the number of inputs specified in `sys`.

The optimization algorithm assumes that finite values in `InputLevel` are fixed input values. Use `NaN` to specify unknown input signals with initial guesses of  $\emptyset$ . The minimum and maximum bounds for all inputs have default values of `-Inf` and `+Inf` respectively.

### **OutputLevel — Steady-state output level**

vector

Steady-state output level for computing the operating point, specified as a vector. The length of `OutputLevel` must equal the number of outputs specified in `sys`.

The values in `OutputLevel` indicate initial guesses for the optimization algorithm. Use `NaN` to specify unknown output signals with initial guesses of  $\emptyset$ . The minimum and maximum bounds for all outputs have default values of `-Inf` and `+Inf` respectively.

### **spec — Operating-point specifications**

`operspec` object

Operating-point specifications, such as minimum and maximum input/output constraints and known inputs, specified as an `operspec` object.

### **T — Operating point snapshot time**

positive scalar

Operating point snapshot time, specified as a positive scalar. The value of `T` must be in the range  $[T_0, N \cdot T_s]$ , where  $N$  is the number of input samples,  $T_s$  is the sample time and  $T_0$  is the input start time (`Uin.Tstart`).

### **Uin — Snapshot simulation input**

`iddata` object | matrix

Snapshot simulation input, specified as one of the following:

- Time-domain `iddata` object with a sample time and input size that matches `sys`.
- Matrix with as many columns as there are input channels. If the matrix has  $N$  rows, the input data is assumed to correspond to the time vector  $(1:N) \cdot \text{sys}.T_s$ .

**X0 — Initial states**

column vector

Initial states of the simulation, specified as a column vector with length equal to the number of states in `sys`. `X0` provides the initial conditions at the time corresponding to the first input sample (`Uin.Start`, if `Uin` is an `iddata` object, or `sys.Ts` if `Uin` is a double matrix).

For more information about the states of an `idnlhw` model, see “Definition of `idnlhw` States” on page 1-672.

**Options — Operating point search options**`findopOptions` option set

Operating point search options, specified as a `findopOptions` option set.

## Output Arguments

**X — Operating point state values**

column vector

Operating point state values, returned as a column vector of length equal to the number of model states.

**U — Operating point input values**

column vector

Operating point input values, returned as a column vector of length equal to the number of inputs.

**Report — Search result summary**

structure

Search result summary report, returned as a structure with the following fields:

Field	Description
<code>SearchMethod</code>	Search method used for iterative parameter estimation. See <code>SearchMethod</code> in <code>findopOptions</code> for more information.
<code>WhyStop</code>	Search algorithm termination condition.

Field	Description
Iterations	Number of estimation iterations performed.
FinalCost	Final value of the minimization objective function (sum of the squared errors).
FirstOrderOptimality	$\infty$ -norm of the search gradient vector when the search algorithm terminates.
SignalLevels	Structure containing the fields <code>Input</code> and <code>Output</code> , which are the operating point input and output signal levels respectively.

## Algorithms

`findop` computes the operating point from steady-state operating point specifications or at a simulation snapshot.

### Computing the Operating Point from Steady-State Specifications

To compute the steady-state operating point, call `findop` using either of the following syntaxes:

```
[X,U] = findop(sys,'steady',InputLevel,OutputLevel)
[X,U] = findop(sys,spec)
```

`findop` uses a different approach to compute the steady-state operating point depending on how much information you provide for this computation:

- When you specify values for all input levels (no NaN values). For a given input level,  $U$ , the equilibrium state values are  $X = \text{inv}(I-A)*B*f(U)$ , where  $[A, B, C, D] = \text{ssdata}(\text{model.LinearModel})$ , and  $f()$  is the input nonlinearity.
- When you specify known and unknown input levels. `findop` uses numerical optimization to minimize the norm of the error and compute the operating point. The total error is the union of contributions from  $e_1$  and  $e_2$ ,  $e(t) = (e_1(t)e_2(t))$ , such that:
  - $e_1$  applies for known outputs and the algorithm minimizes  $e_1 = y - g(L(x,f(u)))$ , where  $f$  is the input nonlinearity,  $L(x,u)$  is the linear model with states  $x$ , and  $g$  is the output nonlinearity.

- $e_2$  applies for unknown outputs and the error is a measure of whether these outputs are within the specified minimum and maximum bounds. If a variable is within its specified bounds, the corresponding error is zero. Otherwise, the error is equal to the distance from the nearest bound. For example, if a free output variable has a value  $z$  and its minimum and maximum bounds are  $L$  and  $U$ , respectively, then the error is  $e_2 = \max[z-U, L-z, 0]$ .

The independent variables for the minimization problem are the unknown inputs. In the error definition  $e$ , both the input  $u$  and the states  $x$  are free variables. To get an error expression that contains only unknown inputs as free variables, the algorithm `findop` specifies the states as a function of inputs by imposing steady-state conditions:  $x = \text{inv}(I-A)*B*f(U)$ , where  $A$  and  $B$  are state-space parameters corresponding to the linear model  $L(x,u)$ . Thus, substituting  $x = \text{inv}(I-A)*B*f(U)$  into the error function results in an error expression that contains only unknown inputs as free variables computed by the optimization algorithm.

### Computing the Operating Point at a Simulation Snapshot

When you use the syntax `[X,U] = findop(sys, 'snapshot', T, UIN, X0)`, the algorithm simulates the model output until the snapshot time,  $T$ . At the snapshot time, the algorithm computes the inputs for the linear model block of the Hammerstein-Wiener model (`LinearModel` property of the `idnlhw` object) by transforming the given inputs using the input nonlinearity:  $w = f(u)$ . `findop` uses the resulting  $w$  to compute  $x$  until the snapshot time using the following equation:  $x(t+1) = Ax(t) + Bw(t)$ , where `[A,B,C,D] = ssdata(model.LinearModel)`.

---

**Note** For snapshot-based computations, `findop` does not perform numerical optimization.

---

## Extended Capabilities

### Automatic Parallel Support

Accelerate code by automatically running computation in parallel using Parallel Computing Toolbox™.

Parallel computing support is available for estimation using the `lsqnonlin` search method (requires Optimization Toolbox). To enable parallel computing, use `findopOptions`, set `SearchMethod` to `'lsqnonlin'`, and set `SearchOptions.Advanced.UseParallel` to `true`.

For example:

```
opt = findopOptions(idnlhw);  
opt.SearchMethod = 'lsqnonlin';  
opt.SearchOptions.Advanced.UseParallel = true;
```

## See Also

[findopOptions](#) | [idnlarx/findop](#) | [idnlhw](#) | [idnlhw/operspec](#) | [sim](#)

**Introduced in R2008a**

## findopOptions

Option set for findop

### Syntax

```
opt = findopOptions(model)
opt = findopOptions(model,Name,Value)
```

### Description

`opt = findopOptions(model)` creates a default option set for computing the operating point of a specified nonlinear ARX or Hammerstein-Wiener model. Use dot notation to modify this option set for your specific application. Options that you do not modify retain their default values.

`opt = findopOptions(model,Name,Value)` creates an option set with options specified by one or more `Name,Value` pair arguments.

### Examples

#### Create Default Option Set for Operating Point Search

Create a default option set for findop using an `idnlarx` model

```
opt = findopOptions(idnlarx);
```

#### Create and Modify Default Operating Point Search Options

Create a default option set for findop using an `idnlhw` model.

```
opt = findopOptions(idnlhw);
```

Use dot notation to specify a subspace Gauss-Newton least squares search with a maximum of 25 iterations.

```
opt.SearchMethod = 'gn';
opt.SearchOptions.MaxIterations = 25;
```

### Specify Options for Operating Point Search

Create an option set for `findop` using an `idnlarx` model. Specify a steepest descent least squares search with default search options.

```
opt = findopOptions(idnlarx, 'SearchMethod', 'grad');
```

## Input Arguments

### **model** — Estimated nonlinear model

`idnlarx model` | `idnlhw model`

Estimated nonlinear model, specified as one of the following:

- `idnlarx model`
- `idnlhw model`

### Name-Value Pair Arguments

Specify optional comma-separated pairs of `Name`, `Value` arguments. `Name` is the argument name and `Value` is the corresponding value. `Name` must appear inside quotes. You can specify several name and value pair arguments in any order as `Name1, Value1, ..., NameN, ValueN`.

Example: `'SearchMethod', 'grad'` specifies a steepest descent least squares search method

### **SearchMethod** — Numerical search method used for iterative parameter estimation

`'auto'` (default) | `'gn'` | `'gna'` | `'lm'` | `'grad'` | `'lsqnonlin'` | `'fmincon'`

Numerical search method used for iterative parameter estimation, specified as the comma-separated pair consisting of 'SearchMethod' and one of the following:

- 'auto' — A combination of the line search algorithms, 'gn', 'lm', 'gna', and 'grad' methods is tried in sequence at each iteration. The first descent direction leading to a reduction in estimation cost is used.
- 'gn' — Subspace Gauss-Newton least squares search. Singular values of the Jacobian matrix less than  $\text{GnPinvConstant} \cdot \text{eps} \cdot \max(\text{size}(J)) \cdot \text{norm}(J)$  are discarded when computing the search direction.  $J$  is the Jacobian matrix. The Hessian matrix is approximated as  $J^T J$ . If there is no improvement in this direction, the function tries the gradient direction.
- 'gna' — Adaptive subspace Gauss-Newton search. Eigenvalues less than  $\text{gamma} \cdot \max(\text{sv})$  of the Hessian are ignored, where  $\text{sv}$  contains the singular values of the Hessian. The Gauss-Newton direction is computed in the remaining subspace.  $\text{gamma}$  has the initial value `InitialGnaTolerance` (see Advanced in 'SearchOptions' for more information). This value is increased by the factor `LMStep` each time the search fails to find a lower value of the criterion in fewer than five bisections. This value is decreased by the factor  $2 \cdot \text{LMStep}$  each time a search is successful without any bisections.
- 'lm' — Levenberg-Marquardt least squares search, where the next parameter value is  $-\text{pinv}(H+d \cdot I) \cdot \text{grad}$  from the previous one.  $H$  is the Hessian,  $I$  is the identity matrix, and  $\text{grad}$  is the gradient.  $d$  is a number that is increased until a lower value of the criterion is found.
- 'grad' — Steepest descent least squares search.
- 'lsqnonlin' — Trust-region-reflective algorithm of `lsqnonlin`. Requires Optimization Toolbox software.
- 'fmincon' — Constrained nonlinear solvers. You can use the sequential quadratic programming (SQP) and trust-region-reflective algorithms of the `fmincon` solver. If you have Optimization Toolbox software, you can also use the interior-point and active-set algorithms of the `fmincon` solver. Specify the algorithm in the `SearchOptions.Algorithm` option. The `fmincon` algorithms may result in improved estimation results in the following scenarios:
  - Constrained minimization problems when there are bounds imposed on the model parameters.
  - Model structures where the loss function is a nonlinear or non smooth function of the parameters.

- Multi-output model estimation. A determinant loss function is minimized by default for multi-output model estimation. `fmincon` algorithms are able to minimize such loss functions directly. The other search methods such as `'lm'` and `'gn'` minimize the determinant loss function by alternately estimating the noise variance and reducing the loss value for a given noise variance value. Hence, the `fmincon` algorithms can offer better efficiency and accuracy for multi-output model estimations.

**SearchOptions — Option set for the search algorithm**

search option set

Option set for the search algorithm, specified as the comma-separated pair consisting of `'SearchOptions'` and a search option set with fields that depend on the value of `SearchMethod`.

**SearchOptions Structure When SearchMethod is Specified as 'gn', 'gna', 'lm', 'grad', or 'auto'**

Field Name	Description	Default
Tolerance	Minimum percentage difference between the current value of the loss function and its expected improvement after the next iteration, specified as a positive scalar. When the percentage of expected improvement is less than <code>Tolerance</code> , the iterations stop. The estimate of the expected loss-function improvement at the next iteration is based on the Gauss-Newton vector computed for the current parameter value.	0.01
MaxIterations	<p>Maximum number of iterations during loss-function minimization, specified as a positive integer. The iterations stop when <code>MaxIterations</code> is reached or another stopping criterion is satisfied, such as <code>Tolerance</code>.</p> <p>Setting <code>MaxIterations = 0</code> returns the result of the start-up procedure.</p> <p>Use <code>sys.Report.Termination.Iterations</code> to get the actual number of iterations during an estimation, where <code>sys</code> is an <code>idtf</code> model.</p>	20

Field Name	Description	Default
Advanced	Advanced search settings, specified as a structure with the following fields:	
Field Name	Description	Default
GnPinvConstant	Jacobian matrix singular value threshold, specified as a positive scalar. Singular values of the Jacobian matrix that are smaller than $\text{GnPinvConstant} * \max(\text{size}(J) * \text{norm}(J) * \text{eps})$ are discarded when computing the search direction. Applicable when SearchMethod is 'gn'.	10000
InitialGnaTolerance	Initial value of $\gamma$ , specified as a positive scalar. Applicable when SearchMethod is 'gna'.	0.0001
LMStartValue	Starting value of search-direction length $d$ in the Levenberg-Marquardt method, specified as a positive scalar. Applicable when SearchMethod is 'lm'.	0.001
LMStep	Size of the Levenberg-Marquardt step, specified as a positive integer. The next value of the search-direction length $d$ in the Levenberg-Marquardt method is LMStep times the previous one. Applicable when SearchMethod is 'lm'.	2
MaxBisections	Maximum number of bisections used for line search along the search direction, specified as a positive integer.	25
MaxFunctionEvaluations	Maximum number of calls to the model file, specified as a positive integer. Iterations stop if the number of calls to the model file exceeds this value.	Inf
MinParameterChange	Smallest parameter update allowed per iteration, specified as a nonnegative scalar.	0
RelativeImprovement	Relative improvement threshold, specified as a nonnegative scalar. Iterations stop if the relative	0

<b>Field Name</b>	<b>Description</b>		<b>Default</b>
	<b>Field Name</b>	<b>Description</b>	<b>Default</b>
		improvement of the criterion function is less than this value.	
	StepReduction	<p>Step reduction factor, specified as a positive scalar that is greater than 1. The suggested parameter update is reduced by the factor <code>StepReduction</code> after each try. This reduction continues until <code>MaxBisections</code> tries are completed or a lower value of the criterion function is obtained.</p> <p><code>StepReduction</code> is not applicable for SearchMethod 'lm' (Levenberg-Marquardt method).</p>	2

**SearchOptions Structure When SearchMethod is Specified as 'lsqnonlin'**

<b>Field Name</b>	<b>Description</b>	<b>Default</b>
FunctionTolerance	Termination tolerance on the loss function that the software minimizes to determine the estimated parameter values, specified as a positive scalar.  The value of FunctionTolerance is the same as that of <code>opt.SearchOptions.Advanced.TolFun</code> .	1e-5
StepTolerance	Termination tolerance on the estimated parameter values, specified as a positive scalar.  The value of StepTolerance is the same as that of <code>opt.SearchOptions.Advanced.TolX</code> .	1e-6
MaxIterations	Maximum number of iterations during loss-function minimization, specified as a positive integer. The iterations stop when MaxIterations is reached or another stopping criterion is satisfied, such as FunctionTolerance.  The value of MaxIterations is the same as that of <code>opt.SearchOptions.Advanced.MaxIter</code> .	20
Advanced	Advanced search settings, specified as an option set for <code>lsqnonlin</code> .  For more information, see the Optimization Options table in "Optimization Options" (Optimization Toolbox).	Use <code>optimset('lsqnonlin')</code> to create a default option set.

**SearchOptions Structure When SearchMethod is Specified as 'fmincon'**

Field Name	Description	Default
Algorithm	<p>fmincon optimization algorithm, specified as one of the following:</p> <ul style="list-style-type: none"> <li>• 'sqp' — Sequential quadratic programming algorithm. The algorithm satisfies bounds at all iterations, and it can recover from NaN or Inf results. It is not a large-scale algorithm. For more information, see “Large-Scale vs. Medium-Scale Algorithms” (Optimization Toolbox).</li> <li>• 'trust-region-reflective' — Subspace trust-region method based on the interior-reflective Newton method. It is a large-scale algorithm.</li> <li>• 'interior-point' — Large-scale algorithm that requires Optimization Toolbox software. The algorithm satisfies bounds at all iterations, and it can recover from NaN or Inf results.</li> <li>• 'active-set' — Requires Optimization Toolbox software. The</li> </ul>	'sqp'

Field Name	Description	Default
	<p>algorithm can take large steps, which adds speed. It is not a large-scale algorithm.</p> <p>For more information about the algorithms, see “Constrained Nonlinear Optimization Algorithms” (Optimization Toolbox) and “Choosing the Algorithm” (Optimization Toolbox).</p>	
FunctionTolerance	Termination tolerance on the loss function that the software minimizes to determine the estimated parameter values, specified as a positive scalar.	1e-6
StepTolerance	Termination tolerance on the estimated parameter values, specified as a positive scalar.	1e-6
MaxIterations	Maximum number of iterations during loss function minimization, specified as a positive integer. The iterations stop when MaxIterations is reached or another stopping criterion is satisfied, such as FunctionTolerance.	100

To specify field values in `SearchOptions`, create a default `findopOptions` set and modify the fields using dot notation. Any fields that you do not modify retain their default values.

```
opt = findopOptions;  
opt.SearchOptions.MaxIterations = 15;  
opt.SearchOptions.Advanced.RelImprovement = 0.5;
```

## Output Arguments

**opt** — Option set for `findop` command

`findopOptions` object

Option set for `findop` command, returned as a `findopOptions` object.

## See Also

[idnlarx/findop](#) | [idnlhw/findop](#)

**Introduced in R2015a**

# findstates

Estimate initial states of model

## Syntax

```
x0 = findstates(sys,Data)
x0 = findstates(sys,Data,Horizon)
x0 = findstates(sys,Data,Horizon,Options)
```

```
[x0,Report]= findstates( ___ )
```

## Description

`x0 = findstates(sys,Data)` estimates the initial states, `x0`, of an identified model `sys`, to maximize the fit between the model response and the output signal in the estimation data.

`x0 = findstates(sys,Data,Horizon)` specifies the prediction horizon for computing the response of `sys`.

`x0 = findstates(sys,Data,Horizon,Options)` specifies additional options for computation of `x0`.

`[x0,Report]= findstates( ___ )` delivers a report on the initial state estimation. Report is returned with any of the previous syntaxes.

## Examples

### Estimate Initial States of a Model

Create a nonlinear grey-box model. The model is a linear DC motor with one input (voltage), and two outputs (angular position and angular velocity). The structure of the model is specified by `dcmotor_m.m` file.

```
FileName = 'dcmotor_m';
Order = [2 1 2];
Parameters = [0.24365;0.24964];
nlgr = idnlgrey(FileName,Order,Parameters);
nlgr = setinit(nlgr, 'Fixed', false(2,1)); % set initial states free
```

Load data for initial state estimation.

```
load(fullfile(matlabroot,'toolbox','ident',...
    'iddemos','data','dcmotordata'));
z = iddata(y,u,0.1);
```

Estimate the initial states such that the model's response using the estimated states  $X_0$  and measured input  $u$  is as close as possible to the measured output  $y$ .

```
X0 = findstates(nlgr,z,Inf);
```

## Estimate Initial States of State-Space Model

Estimate an `idss` model and simulate it such that the response of the estimated model matches the estimation data's output signal as closely as possible.

Load sample data.

```
load iddata1 z1;
```

Estimate a linear model from the data.

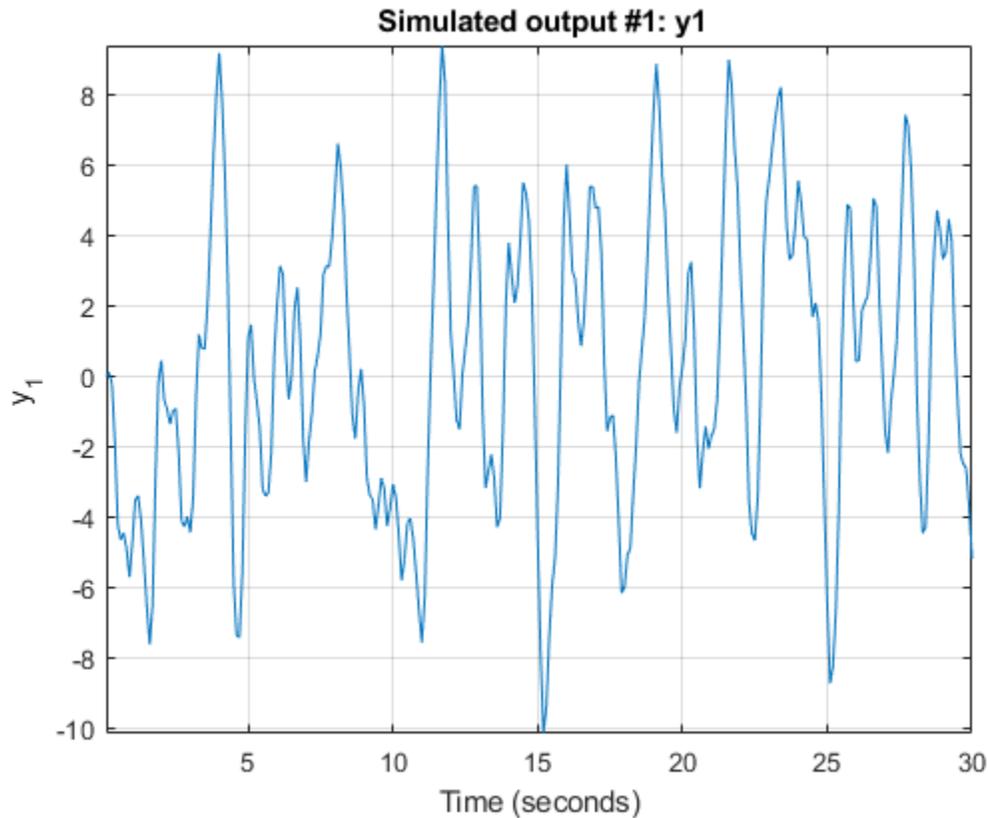
```
model = ssest(z1,2);
```

Estimate the value of the initial states to best fit the measured output  $z1.y$ .

```
x0est = findstates(model,z1,Inf);
```

Simulate the model.

```
opt = simOptions('InitialCondition',x0est);
sim(model,z1(:,[],:),opt);
```



### Selectively Estimate Initial States of a Model

Estimate the initial states of a model selectively by fixing the first state and allowing the second state of the model to be estimated.

Create a nonlinear grey-box model.

```
FileName = 'dcmotor_m';  
Order = [2 1 2];  
Parameters = [0.24365;0.24964];  
nlgr = idnlgrey(FileName,Order,Parameters);
```

The model is a linear DC motor with one input (voltage), and two outputs (angular position and angular velocity). The structure of the model is specified by `dcmotor_m.m` file.

Load the estimation data.

```
load(fullfile(matlabroot, 'toolbox', 'ident', ...  
            'iddemos', 'data', 'dcmotordata'));  
z = iddata(y,u,0.1);
```

Hold the first state fixed at zero, and estimate the value of the second.

```
x0spec = idpar('x0',[0;0]);  
x0spec.Free(1) = false;  
opt = findstatesOptions;  
opt.InitialState = x0spec;  
[X0,Report] = findstates(nlgr,z,Inf,opt)
```

```
X0 = 2x1
```

```
      0  
    0.0061
```

```
Report =
```

```
      Status: 'Estimated by simulation error minimization'  
      Method: 'lsqnonlin'  
      Covariance: [2x2 double]  
      DataUsed: [1x1 struct]  
      Termination: [1x1 struct]
```

## Estimate Initial States by Specifying an Initial State Vector

Create a nonlinear grey-box model.

```
FileName = 'dcmotor_m';  
Order = [2 1 2];  
Parameters = [0.24365;0.24964];  
nlgr = idnlgrey(FileName,Order,Parameters);
```

The model is a linear DC motor with one input (voltage), and two outputs (angular position and angular velocity). The structure of the model is specified by `dcmotor_m.m` file.

Load the estimation data.

```
load(fullfile(matlabroot,'toolbox','ident',...
    'iddemos','data','dcmotordata'));
z = iddata(y,u,0.1);
```

Specify an initial guess for the initial states.

```
x0spec = idpar('x0',[10;10]);
```

`x0spec.Free` is true by default

Estimate the initial states

```
opt = findstatesOptions;
opt.InitialState = x0spec;
x0 = findstates(nlgr,z,Inf,opt)
```

```
x0 = 2×1
```

```
    0.0362
   -0.1322
```

## Estimate Initial States Using Multi-Experiment Data

Create a nonlinear grey-box model.

```
FileName = 'dcmotor_m';
Order = [2 1 2];
Parameters = [0.24365;0.24964];
nlgr = idnlgrey(FileName,Order,Parameters);
set(nlgr, 'InputName','Voltage','OutputName', ...
    {'Angular position','Angular velocity'});
```

The model is a linear DC motor with one input (voltage), and two outputs (angular position and angular velocity). The structure of the model is specified by `dcmotor_m.m` file.

Load the estimation data.

```
load(fullfile(matlabroot,'toolbox','ident',...  
    'iddemos','data','dcmotordata'));  
z = iddata(y,u,0.1,'Name','DC-motor',...  
    'InputName','Voltage','OutputName',...  
    {'Angular position','Angular velocity'});
```

Create a three-experiment data set.

```
z3 = merge(z,z,z);
```

Choose experiment for estimating the initial states:

- Estimate initial state 1 for experiments 1 and 3
- Estimate initial state 2 for experiment 1

The fixed initial states have zero values.

```
x0spec = idpar('x0',zeros(2,3));  
x0spec.Free(1,2) = false;  
x0spec.Free(2,[2 3]) = false;  
opt = findstatesOptions;  
opt.InitialState = x0spec;
```

Estimate the initial states

```
[X0,EstInfo] = findstates(nlgr,z3,Inf,opt);
```

## Input Arguments

### **sys** — Identified model

idss object | idgrey object | idnlarx object | idnlhw object | idnlgrey object

Identified model whose initial states are estimated, represented as a linear state-space (idss or idgrey) or nonlinear model (idnlarx, idnlhw, or idnlgrey).

### **Data** — Estimation data

iddata object

Estimation data, specified as an `iddata` object with input/output dimensions that match `sys`.

If `sys` is a linear model, `Data` can be a frequency-domain `iddata` object. For easier interpretation of initial conditions, make the frequency vector of `Data` be symmetric about the origin. For converting time-domain data into frequency-domain data, use `fft` with `'compl'` input argument, and ensure that there is sufficient zero padding. Scale your data appropriately when you compare  $x_0$  between the time-domain and frequency-domain. Since for an  $N$ -point `fft`, the input/output signals are scaled by  $1/\sqrt{N}$ , the estimated  $x_0$  vector is also scaled by this factor.

### Horizon — Prediction horizon for computing model response

1 (default) | positive integer between 1 and `Inf`

Prediction horizon for computing the response of `sys`, specified as a positive integer between 1 and `Inf`. The most common values used are:

- `Horizon = 1` — Minimizes the 1-step prediction error. The 1-step ahead prediction response of `sys` is compared to the output signals in `Data` to determine  $x_0$ . See `predict` for more information.
- `Horizon = Inf` — Minimizes the simulation error. The difference between measured output, `Data.y`, and simulated response of `sys` to the measured input data, `Data.u` is minimized. See `sim` for more information.

Specify `Horizon` as any positive integer between 1 and `Inf`, with the following restrictions:

Scenario	Horizon
Continuous-time model with time-domain data	1 or <code>Inf</code>
Continuous-time frequency-domain data ( <code>data.Ts = 0</code> )	<code>Inf</code>
Output Error models (trivial noise component):	Irrelevant
<ul style="list-style-type: none"> <li>• Nonlinear grey-box (<code>idnlgrey</code>)</li> <li>• Hammerstein-Wiener (<code>idnlhw</code>)</li> <li>• Linear state-space with disturbance matrix, <math>K = 0</math></li> </ul>	Any value of <code>Horizon</code> returns the same answer for $x_0$
Nonlinear ARX ( <code>idnlarx</code> )	1 or <code>Inf</code>

### Options — Estimation options for findstates

`findstates` Option set

Estimation options for `findstates`, specified as an option set created using `findstatesOptions`

## Output Arguments

### **x0** — Estimated initial states

vector | matrix

Estimated initial states of model `sys`, returned as a vector or matrix. For multi-experiment data, `x0` is a matrix with one column for each experiment.

### **Report** — Initial state estimation information

structure

Initial state estimation information, returned as a structure. `Report` contains information about the data used, state covariance, and results of any numerical optimization performed to search for the initial states. `Report` has the following fields:

<b>Report Field</b>	<b>Description</b>
Status	Summary of how the initial state were estimated.
Method	Search method used.
Covariance	Covariance of state estimates, returned as a $N_s$ -by- $N_s$ matrix, where $N_s$ is the number of states.

Report Field	Description																
DataUsed	Attributes of the data used for estimation, returned as a structure with the following fields:																
	<table border="1"> <thead> <tr> <th data-bbox="385 425 516 465">Field</th> <th data-bbox="516 425 1335 465">Description</th> </tr> </thead> <tbody> <tr> <td data-bbox="385 465 516 512">Name</td> <td data-bbox="516 465 1335 512">Name of the data set.</td> </tr> <tr> <td data-bbox="385 512 516 558">Type</td> <td data-bbox="516 512 1335 558">Data type.</td> </tr> <tr> <td data-bbox="385 558 516 605">Length</td> <td data-bbox="516 558 1335 605">Number of data samples.</td> </tr> <tr> <td data-bbox="385 605 516 652">Ts</td> <td data-bbox="516 605 1335 652">Sample time.</td> </tr> <tr> <td data-bbox="385 652 516 956">InterSample</td> <td data-bbox="516 652 1335 956">           Input intersample behavior, returned as one of the following values:           <ul style="list-style-type: none"> <li>• 'zoh' — Zero-order hold maintains a piecewise-constant input signal between samples.</li> <li>• 'foh' — First-order hold maintains a piecewise-linear input signal between samples.</li> <li>• 'bl' — Band-limited behavior specifies that the continuous-time input signal has zero power above the Nyquist frequency.</li> </ul> </td> </tr> <tr> <td data-bbox="385 956 516 1038">InputOffset</td> <td data-bbox="516 956 1335 1038">Offset removed from time-domain input data during estimation. For nonlinear models, it is [ ].</td> </tr> <tr> <td data-bbox="385 1038 516 1111">OutputOffset</td> <td data-bbox="516 1038 1335 1111">Offset removed from time-domain output data during estimation. For nonlinear models, it is [ ].</td> </tr> </tbody> </table>	Field	Description	Name	Name of the data set.	Type	Data type.	Length	Number of data samples.	Ts	Sample time.	InterSample	Input intersample behavior, returned as one of the following values: <ul style="list-style-type: none"> <li>• 'zoh' — Zero-order hold maintains a piecewise-constant input signal between samples.</li> <li>• 'foh' — First-order hold maintains a piecewise-linear input signal between samples.</li> <li>• 'bl' — Band-limited behavior specifies that the continuous-time input signal has zero power above the Nyquist frequency.</li> </ul>	InputOffset	Offset removed from time-domain input data during estimation. For nonlinear models, it is [ ].	OutputOffset	Offset removed from time-domain output data during estimation. For nonlinear models, it is [ ].
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InputOffset	Offset removed from time-domain input data during estimation. For nonlinear models, it is [ ].																
OutputOffset	Offset removed from time-domain output data during estimation. For nonlinear models, it is [ ].																

Report Field	Description	
Termination	Termination conditions for the iterative search used for initial state estimation of nonlinear models. Structure with the following fields:	
	<b>Field</b>	<b>Description</b>
	WhyStop	Reason for terminating the numerical search.
	Iterations	Number of search iterations performed by the estimation algorithm.
	FirstOrderOptimality	$\infty$ -norm of the gradient search vector when the search algorithm terminates.
	FcnCount	Number of times the objective function was called.
	UpdateNorm	Norm of the gradient search vector in the last iteration. Omitted when the search method is 'lsqnonlin' or 'fmincon'.
	LastImprovement	Criterion improvement in the last iteration, expressed as a percentage. Omitted when the search method is 'lsqnonlin' or 'fmincon'.
	Algorithm	Algorithm used by 'lsqnonlin' or 'fmincon' search method. Omitted when other search methods are used.
Termination is empty for linear models.		

## Extended Capabilities

### Automatic Parallel Support

Accelerate code by automatically running computation in parallel using Parallel Computing Toolbox™.

Parallel computing support is available for estimation using the `lsqnonlin` search method (requires Optimization Toolbox). To enable parallel computing, use `findstatesOptions`, set `SearchMethod` to 'lsqnonlin', and set `SearchOptions.Advanced.UseParallel` to true.

For example:

```
opt = findstatesOptions;  
opt.SearchMethod = 'lsqnonlin';  
opt.SearchOptions.Advanced.UseParallel = true;
```

## See Also

[findstatesOptions](#) | [predict](#) | [sim](#)

**Introduced in R2015a**

## findstatesOptions

Option set for findstates

### Syntax

```
opt = findstatesOptions  
opt = findstatesOptions(Name,Value)
```

### Description

`opt = findstatesOptions` creates the default option set for `findstates`. Use dot notation to customize the option set, if needed.

`opt = findstatesOptions(Name,Value)` creates an option set with options specified by one or more `Name,Value` pair arguments. The options that you do not specify retain their default value.

### Examples

#### Identify Initial States Using Option Set

Create an option set for `findstates` by configuring a specification object for the initial states.

Identify a fourth-order state-space model from data.

```
load iddata8 z8;  
sys = ssest(z8,4);
```

`z8` is an `iddata` object containing time-domain system response data. `sys` is a fourth-order `idss` model that is identified from the data.

Configure a specification object for the initial states of the model.

```
x0obj = idpar([1;nan(3,1)]);
x0obj.Free(1) = false;
x0obj.Minimum(2) = 0;
x0obj.Maximum(2) = 1;
```

`x0obj` specifies estimation constraints on the initial conditions. The value of the first state is specified as 1 when `x0obj` is created. `x0obj.Free(1) = false` specifies the first initial state as a fixed estimation parameter. The second state is unknown. But, `x0obj.Minimum(2) = 0` and `x0obj.Maximum(2) = 1` specify the lower and upper bounds of the second state as 0 and 1, respectively.

Create an option set for `findstates` to identify the initial states of the model.

```
opt = findstatesOptions;
opt.InitialState = x0obj;
```

Identify the initial states of the model.

```
x0_estimated = findstates(sys,z8,Inf,opt);
```

## Specify Option Set for Initial States Estimation

Create an option set for `findstates` where:

- Initial states are estimated such that the norm of prediction error is minimized. The initial values of the states corresponding to nonzero delays are also estimated.
- Adaptive subspace Gauss-Newton search is used for estimation.

```
opt = findstatesOptions('InitialState','d','SearchMethod','gna');
```

## Input Arguments

### Name-Value Pair Arguments

Specify optional comma-separated pairs of `Name,Value` arguments. `Name` is the argument name and `Value` is the corresponding value. `Name` must appear inside quotes. You can specify several name and value pair arguments in any order as `Name1,Value1,...,NameN,ValueN`.

Example: `findstatesOptions('InitialState','d')`

### **InitialState — Estimation of initial states**

'e' (default) | 'd' | vector or matrix | idpar object x0obj

Estimation of initial states, specified as the comma-separated pair consisting of 'InitialState' and one of the following:

- 'e' — The initial states are estimated such that the norm of prediction error is minimized.

For nonlinear grey-box models, only those initial states `i` that are designated as free in the model (`sys.InitialStates(i).Fixed = false`) are estimated. To estimate all the states of the model, first specify all the `Nx` states of the `idnlgrey` model `sys` as free.

```
for i = 1:Nx
    sys.InitialStates(i).Fixed = false;
end
```

Similarly, to fix all the initial states to values specified in `sys.InitialStates`, first specify all the states as fixed in the `sys.InitialStates` property of the nonlinear grey-box model.

- 'd' — Similar to 'e', but absorbs nonzero delays into the model coefficients. The delays are first converted to explicit model states, and the initial values of those states are also estimated and returned.

Use this option for discrete-time linear models only.

- **Vector or Matrix** — Initial guess for state values, when using nonlinear models. Specify a column vector of length equal to the number of states. For multi-experiment data, use a matrix with `Ne` columns, where `Ne` is the number of experiments.

Use this option for nonlinear models only.

- `x0obj` — Specification object created using `idpar`. Use `x0obj` to impose constraints on the initial states by fixing their value or specifying minimum or maximum bounds.

Use `x0obj` only for nonlinear grey-box models and linear state-space models (`idss` or `idgrey`). This option is applicable only for prediction horizon equal to 1 or `Inf`. See `findstates` for more details about the prediction horizon.

### **InputOffset — Removal of offset from time-domain input data during estimation**

[ ] (default) | vector of positive integers | matrix

Removal of offset from time-domain input data during estimation, specified as the comma-separated pair consisting of 'InputOffset' and one of the following:

- A column vector of positive integers of length  $Nu$ , where  $Nu$  is the number of inputs.
- [ ] — Indicates no offset.
- $Nu$ -by- $Ne$  matrix — For multi-experiment data, specify InputOffset as an  $Nu$ -by- $Ne$  matrix.  $Nu$  is the number of inputs, and  $Ne$  is the number of experiments.

Each entry specified by InputOffset is subtracted from the corresponding input data.

### **OutputOffset — Removal of offset from time-domain output data during estimation**

[ ] (default) | vector | matrix

Removal of offset from time-domain output data during estimation, specified as the comma-separated pair consisting of 'OutputOffset' and one of the following:

- A column vector of length  $Ny$ , where  $Ny$  is the number of outputs.
- [ ] — Indicates no offset.
- $Ny$ -by- $Ne$  matrix — For multi-experiment data, specify OutputOffset as a  $Ny$ -by- $Ne$  matrix.  $Ny$  is the number of outputs, and  $Ne$  is the number of experiments.

Each entry specified by OutputOffset is subtracted from the corresponding output data.

### **OutputWeight — Weighting of prediction errors when using multi-output data**

[ ] (default) | 'noise' | matrix

Weighting of prediction errors when using multi-output data, specified as the comma-separated pair consisting of 'OutputWeight' and one of the following:

- [ ] — No weighting is used. Specifying as [ ] is the same as  $\text{eye}(Ny)$ , where  $Ny$  is the number of outputs.
- 'noise' — Inverse of the noise variance stored with the model is used for weighting during estimation of initial states.

- Positive semidefinite matrix,  $W$ , of size  $N_y$ -by- $N_y$  — This weighting minimizes  $\text{trace}(E' * E * W)$  for estimation of initial states, where  $E$  is the matrix of prediction errors.

**SearchMethod — Numerical search method used for iterative parameter estimation**

'auto' (default) | 'gn' | 'gna' | 'lm' | 'grad' | 'lsqnonlin' | 'fmincon'

Numerical search method used for iterative parameter estimation, specified as the comma-separated pair consisting of 'SearchMethod' and one of the following:

- 'auto' — A combination of the line search algorithms, 'gn', 'lm', 'gna', and 'grad' methods is tried in sequence at each iteration. The first descent direction leading to a reduction in estimation cost is used.
- 'gn' — Subspace Gauss-Newton least squares search. Singular values of the Jacobian matrix less than  $GnPinvConstant * \epsilon * \max(\text{size}(J)) * \text{norm}(J)$  are discarded when computing the search direction.  $J$  is the Jacobian matrix. The Hessian matrix is approximated as  $J'J$ . If there is no improvement in this direction, the function tries the gradient direction.
- 'gna' — Adaptive subspace Gauss-Newton search. Eigenvalues less than  $\gamma * \max(sv)$  of the Hessian are ignored, where  $sv$  contains the singular values of the Hessian. The Gauss-Newton direction is computed in the remaining subspace.  $\gamma$  has the initial value `InitialGnaTolerance` (see Advanced in 'SearchOptions' for more information). This value is increased by the factor `LMStep` each time the search fails to find a lower value of the criterion in fewer than five bisections. This value is decreased by the factor  $2 * LMStep$  each time a search is successful without any bisections.
- 'lm' — Levenberg-Marquardt least squares search, where the next parameter value is  $-\text{pinv}(H + d * I) * \text{grad}$  from the previous one.  $H$  is the Hessian,  $I$  is the identity matrix, and  $\text{grad}$  is the gradient.  $d$  is a number that is increased until a lower value of the criterion is found.
- 'grad' — Steepest descent least squares search.
- 'lsqnonlin' — Trust-region-reflective algorithm of `lsqnonlin`. Requires Optimization Toolbox software.
- 'fmincon' — Constrained nonlinear solvers. You can use the sequential quadratic programming (SQP) and trust-region-reflective algorithms of the `fmincon` solver. If you have Optimization Toolbox software, you can also use the interior-point and active-set algorithms of the `fmincon` solver. Specify the algorithm in the

`SearchOptions.Algorithm` option. The `fmincon` algorithms may result in improved estimation results in the following scenarios:

- Constrained minimization problems when there are bounds imposed on the model parameters.
- Model structures where the loss function is a nonlinear or non smooth function of the parameters.
- Multi-output model estimation. A determinant loss function is minimized by default for multi-output model estimation. `fmincon` algorithms are able to minimize such loss functions directly. The other search methods such as `'lm'` and `'gn'` minimize the determinant loss function by alternately estimating the noise variance and reducing the loss value for a given noise variance value. Hence, the `fmincon` algorithms can offer better efficiency and accuracy for multi-output model estimations.

### **SearchOptions — Option set for the search algorithm**

search option set

Option set for the search algorithm, specified as the comma-separated pair consisting of `'SearchOptions '` and a search option set with fields that depend on the value of `SearchMethod`.

**SearchOptions Structure When SearchMethod is Specified as 'gn', 'gna', 'lm', 'grad', or 'auto'**

Field Name	Description	Default
Tolerance	Minimum percentage difference between the current value of the loss function and its expected improvement after the next iteration, specified as a positive scalar. When the percentage of expected improvement is less than <code>Tolerance</code> , the iterations stop. The estimate of the expected loss-function improvement at the next iteration is based on the Gauss-Newton vector computed for the current parameter value.	0.01
MaxIterations	<p>Maximum number of iterations during loss-function minimization, specified as a positive integer. The iterations stop when <code>MaxIterations</code> is reached or another stopping criterion is satisfied, such as <code>Tolerance</code>.</p> <p>Setting <code>MaxIterations = 0</code> returns the result of the start-up procedure.</p> <p>Use <code>sys.Report.Termination.Iterations</code> to get the actual number of iterations during an estimation, where <code>sys</code> is an <code>idtf</code> model.</p>	20

Field Name	Description	Default
Advanced	Advanced search settings, specified as a structure with the following fields:	
Field Name	Description	Default
GnPinvConstant	Jacobian matrix singular value threshold, specified as a positive scalar. Singular values of the Jacobian matrix that are smaller than $\text{GnPinvConstant} * \max(\text{size}(J) * \text{norm}(J) * \text{eps})$ are discarded when computing the search direction. Applicable when SearchMethod is 'gn'.	10000
InitialGnaTolerance	Initial value of $\gamma$ , specified as a positive scalar. Applicable when SearchMethod is 'gna'.	0.0001
LMStartValue	Starting value of search-direction length $d$ in the Levenberg-Marquardt method, specified as a positive scalar. Applicable when SearchMethod is 'lm'.	0.001
LMStep	Size of the Levenberg-Marquardt step, specified as a positive integer. The next value of the search-direction length $d$ in the Levenberg-Marquardt method is LMStep times the previous one. Applicable when SearchMethod is 'lm'.	2
MaxBisections	Maximum number of bisections used for line search along the search direction, specified as a positive integer.	25
MaxFunctionEvaluations	Maximum number of calls to the model file, specified as a positive integer. Iterations stop if the number of calls to the model file exceeds this value.	Inf
MinParameterChange	Smallest parameter update allowed per iteration, specified as a nonnegative scalar.	0
RelativeImprovement	Relative improvement threshold, specified as a nonnegative scalar. Iterations stop if the relative	0

Field Name	Description		Default
	Field Name	Description	Default
		improvement of the criterion function is less than this value.	
	StepReduction	<p>Step reduction factor, specified as a positive scalar that is greater than 1. The suggested parameter update is reduced by the factor <code>StepReduction</code> after each try. This reduction continues until <code>MaxBisections</code> tries are completed or a lower value of the criterion function is obtained.</p> <p><code>StepReduction</code> is not applicable for SearchMethod 'lm' (Levenberg-Marquardt method).</p>	2

**SearchOptions Structure When SearchMethod is Specified as 'lsqnonlin'**

<b>Field Name</b>	<b>Description</b>	<b>Default</b>
FunctionTolerance	Termination tolerance on the loss function that the software minimizes to determine the estimated parameter values, specified as a positive scalar.  The value of FunctionTolerance is the same as that of <code>opt.SearchOptions.Advanced.TolFun</code> .	1e-5
StepTolerance	Termination tolerance on the estimated parameter values, specified as a positive scalar.  The value of StepTolerance is the same as that of <code>opt.SearchOptions.Advanced.TolX</code> .	1e-6
MaxIterations	Maximum number of iterations during loss-function minimization, specified as a positive integer. The iterations stop when MaxIterations is reached or another stopping criterion is satisfied, such as FunctionTolerance.  The value of MaxIterations is the same as that of <code>opt.SearchOptions.Advanced.MaxIter</code> .	20
Advanced	Advanced search settings, specified as an option set for <code>lsqnonlin</code> .  For more information, see the Optimization Options table in "Optimization Options" (Optimization Toolbox).	Use <code>optimset('lsqnonlin')</code> to create a default option set.

**SearchOptions Structure When SearchMethod is Specified as 'fmincon'**

Field Name	Description	Default
Algorithm	<p>fmincon optimization algorithm, specified as one of the following:</p> <ul style="list-style-type: none"> <li>• 'sqp' — Sequential quadratic programming algorithm. The algorithm satisfies bounds at all iterations, and it can recover from NaN or Inf results. It is not a large-scale algorithm. For more information, see “Large-Scale vs. Medium-Scale Algorithms” (Optimization Toolbox).</li> <li>• 'trust-region-reflective' — Subspace trust-region method based on the interior-reflective Newton method. It is a large-scale algorithm.</li> <li>• 'interior-point' — Large-scale algorithm that requires Optimization Toolbox software. The algorithm satisfies bounds at all iterations, and it can recover from NaN or Inf results.</li> <li>• 'active-set' — Requires Optimization Toolbox software. The</li> </ul>	'sqp'

Field Name	Description	Default
	<p>algorithm can take large steps, which adds speed. It is not a large-scale algorithm.</p> <p>For more information about the algorithms, see “Constrained Nonlinear Optimization Algorithms” (Optimization Toolbox) and “Choosing the Algorithm” (Optimization Toolbox).</p>	
FunctionTolerance	Termination tolerance on the loss function that the software minimizes to determine the estimated parameter values, specified as a positive scalar.	1e-6
StepTolerance	Termination tolerance on the estimated parameter values, specified as a positive scalar.	1e-6
MaxIterations	Maximum number of iterations during loss function minimization, specified as a positive integer. The iterations stop when MaxIterations is reached or another stopping criterion is satisfied, such as FunctionTolerance.	100

To specify field values in `SearchOptions`, create a default `findstatesOptions` set and modify the fields using dot notation. Any fields that you do not modify retain their default values.

```
opt = findstatesOptions;  
opt.SearchOptions.Tolerance = 0.02;  
opt.SearchOptions.Advanced.MaxBisections = 30;
```

## Output Arguments

**opt** — Option set for `findstates`

`findstatesOptions` option set

Option set for `findstates`, returned as an `findstatesOptions` option set.

## See Also

`findstates` | `idpar`

**Introduced in R2012a**

# fnorm

Pointwise peak gain of FRD model

## Syntax

```
fnrm = fnorm(sys)  
fnrm = fnorm(sys, ntype)
```

## Description

`fnrm = fnorm(sys)` computes the pointwise 2-norm of the frequency response contained in the FRD model `sys`, that is, the peak gain at each frequency point. The output `fnrm` is an FRD object containing the peak gain across frequencies.

`fnrm = fnorm(sys, ntype)` computes the frequency response gains using the matrix norm specified by `ntype`. See `norm` for valid matrix norms and corresponding `NTYPE` values.

## See Also

`abs` | `norm`

**Introduced in R2006a**

## forecast

Forecast identified model output

### Syntax

```
yf = forecast(sys,PastData,K)
yf = forecast(sys,PastData,K,FutureInputs)
```

```
yf = forecast( ____,opts)
```

```
[yf,x0,sysf] = forecast( ____)
[yf,x0,sysf,yf_sd,x,x_sd] = forecast( ____)
```

```
forecast(sys,PastData,K, ____)
forecast(sys,Linespec,PastData,K, ____)
forecast(sys1,...,sysN,PastData,K, ____)
forecast(sys1,Linespec1,...,sysN,LinespecN,PastData,K, ____)
```

### Description

`yf = forecast(sys,PastData,K)` forecasts the output of an identified time series model `sys`, `K` steps into the future using past measured data, `PastData`.

`forecast` performs prediction into the future, in a time range beyond the last instant of measured data. In contrast, the `predict` command predicts the response of an identified model over the time span of measured data. Use `predict` to determine if the predicted result matches the observed response of an estimated model. If `sys` is a good prediction model, consider using it with `forecast`.

`yf = forecast(sys,PastData,K,FutureInputs)` uses the future values of the inputs, `FutureInputs`, to forecast the response of an identified model with input channels.

`yf = forecast( ____,opts)` uses the option set, `opts`, to specify additional forecast options. Use `opts` with any of the previous input argument combinations.

`[yf,x0,sysf] = forecast(____)` also returns the estimated values for initial states, `x0`, and a forecasting model, `sysf`, and can include any of the previous input argument combinations.

`[yf,x0,sysf,yf_sd,x,x_sd] = forecast(____)` also returns estimated standard deviation of the output, `yf_sd`, state trajectory, `x`, and standard deviation of the trajectory, `x_sd`. Use with any of the previous input argument combinations.

`forecast(sys,PastData,K,____)` plots the forecasted output. Use with any of the previous input argument combinations.

To change display options, right-click the plot to access the context menu. For example, to view the estimated standard deviation of the forecasted output, select **Confidence Region** from the context menu. For more details about the menu, see “Tips” on page 1-427.

`forecast(sys,Linespec,PastData,K,____)` uses `Linespec` to specify the line type, marker symbol, and color.

`forecast(sys1,...,sysN,PastData,K,____)` plots the forecasted outputs for multiple identified models. `forecast` automatically chooses colors and line styles.

`forecast(sys1,Linespec1,...,sysN,LinespecN,PastData,K,____)` uses the line type, marker symbol, and color specified for each system.

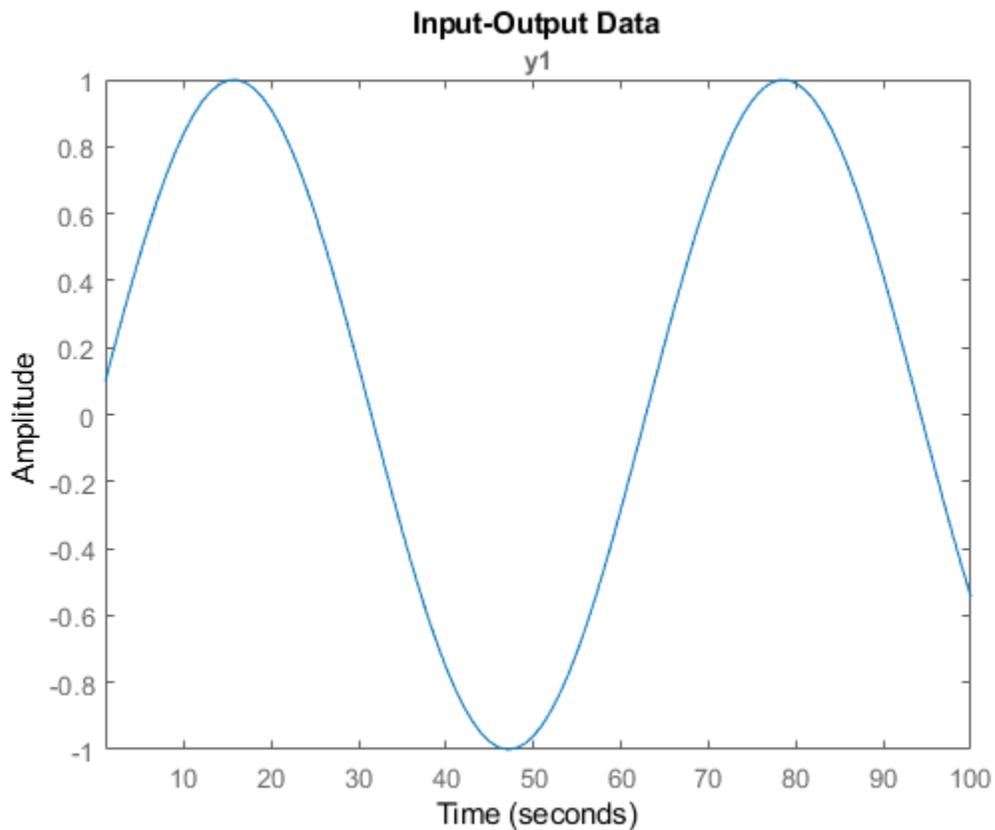
## Examples

### Forecast Future Values of a Sinusoidal Signal

Forecast the values of a sinusoidal signal using an AR model.

Generate and plot data.

```
data = iddata(sin(0.1*[1:100])',[]);
plot(data)
```



Fit an AR model to the sine wave.

```
sys = ar(data,2);
```

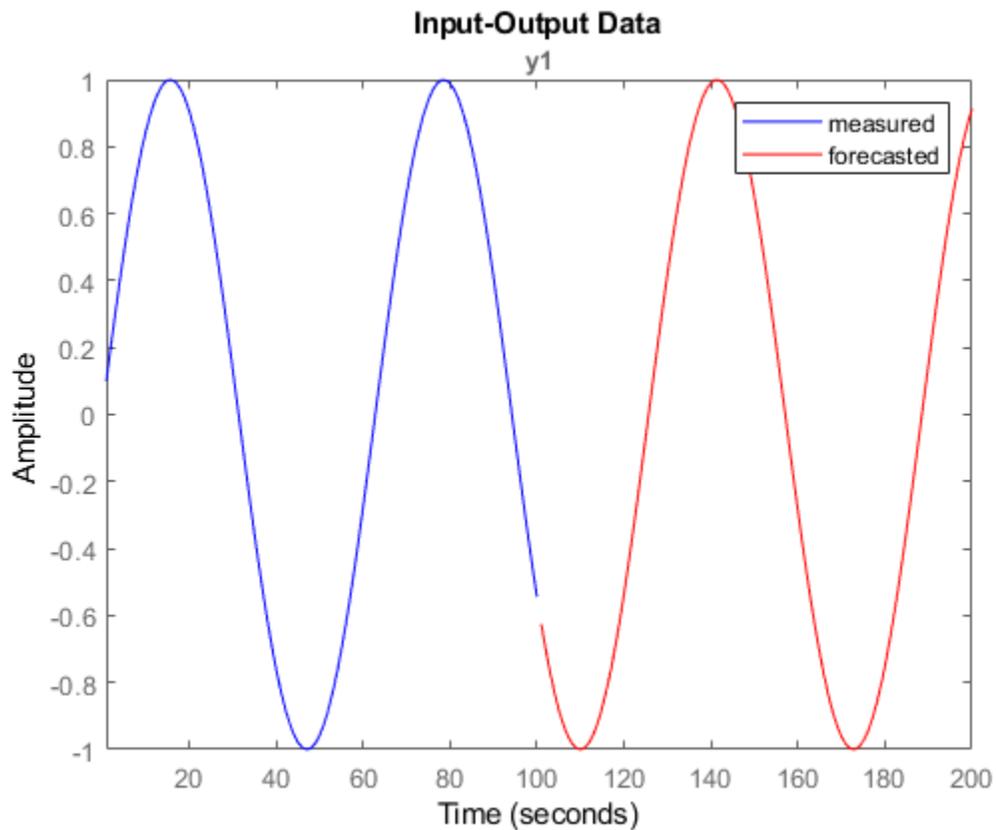
Forecast the values into the future for a given time horizon.

```
K = 100;  
p = forecast(sys,data,K);
```

K specifies the forecasting time horizon as 100 samples. p is the forecasted model response.

Plot the forecasted data.

```
plot(data,'b',p,'r'), legend('measured','forecasted')
```



Alternatively, plot the forecasted output using the syntax `forecast(sys,data,K)`.

### Forecast Response of Time Series Model

Obtain past data, and identify a time series model.

```
load iddata9 z9
past_data = z9.OutputData(1:50);
model = ar(z9,4);
```

`z9` is an `iddata` object that contains measured output only.

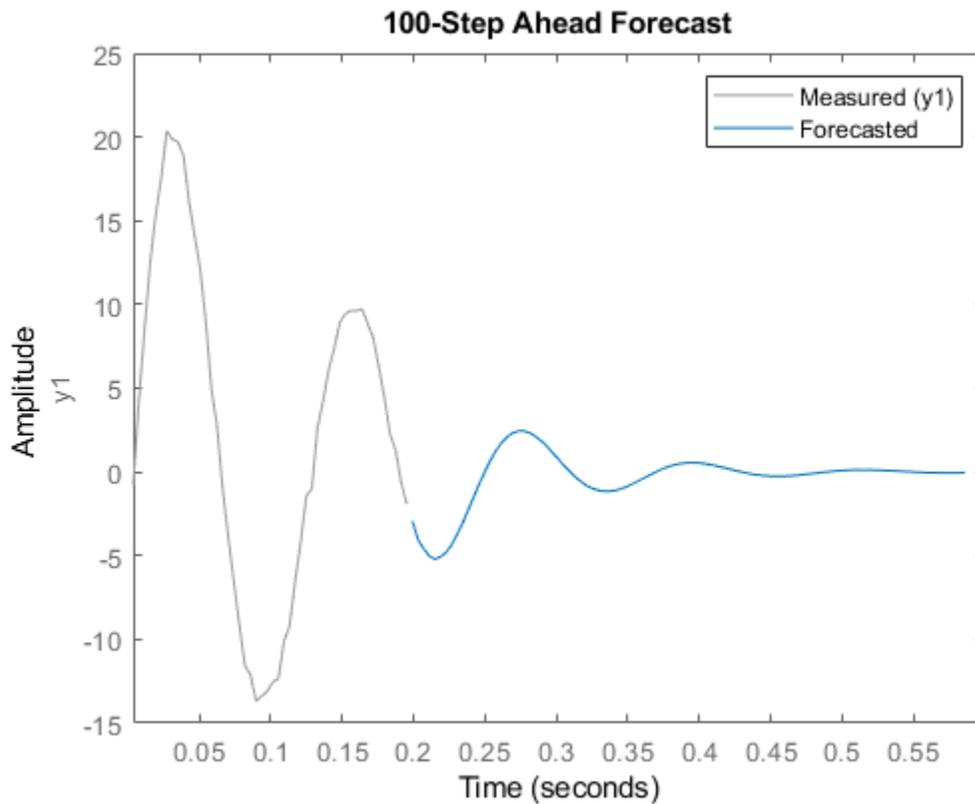
model is an idpoly time series model.

Specify initial conditions for forecasting.

```
opt = forecastOptions('InitialCondition','e');
```

Plot the forecasted system response for a given time horizon.

```
K = 100;  
forecast(model,past_data,K,opt);  
legend('Measured','Forecasted')
```



### Plot Forecasted Output With Specified Line Type

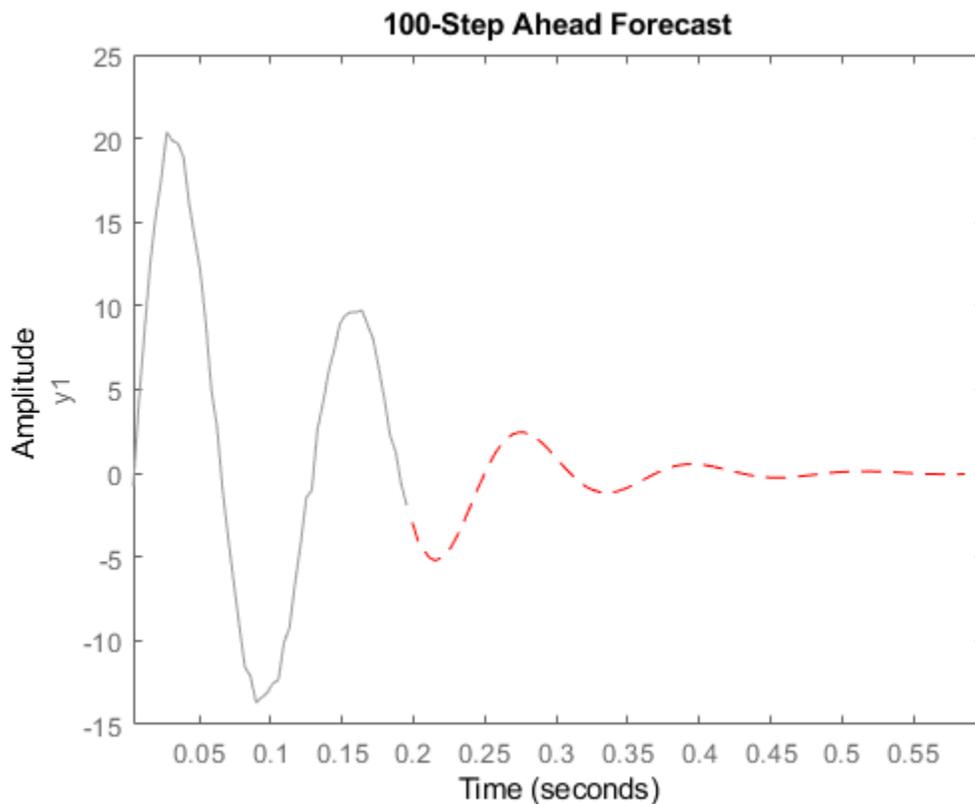
Obtain past data, and identify a time series model.

```
load iddata9 z9
past_data = z9.OutputData(1:50);
model = ar(z9,4);
```

z9 is an `iddata` object that contains measured output only.

Plot the forecasted system response for a given time horizon as a red dashed line.

```
K = 100;
forecast(model, 'r--', past_data, K);
```



The plot also displays the past data by default. To change display options, right-click the plot to access the context menu. For example, to view the estimated standard deviation of the forecasted output, select **ConfidenceRegion** from the context menu. To specify number of standard deviations to plot, double-click the plot and open the Property Editor dialog box. In the dialog box, in the **Options** tab, specify the number of standard deviations in **Confidence Region for Identified Models**. The default value is 1 standard deviation.

### Forecast Model Response for Known Future Inputs

Obtain past data, future inputs, and an identified linear model.

```
load iddata1 z1
z1 = iddata(cumsum(z1.y),cumsum(z1.u),z1.Ts,'InterSample','foh');
past_data = z1(1:100);
future_inputs = z1.u(101:end);
sys = polyest(z1,[2 2 2 0 0 1],'IntegrateNoise',true);
```

`z1` is an `iddata` object that contains integrated data. `sys` is an `idpoly` model. `past_data` contains the first 100 data points of `z1`.

`future_inputs` contains the last 200 data points of `z1`.

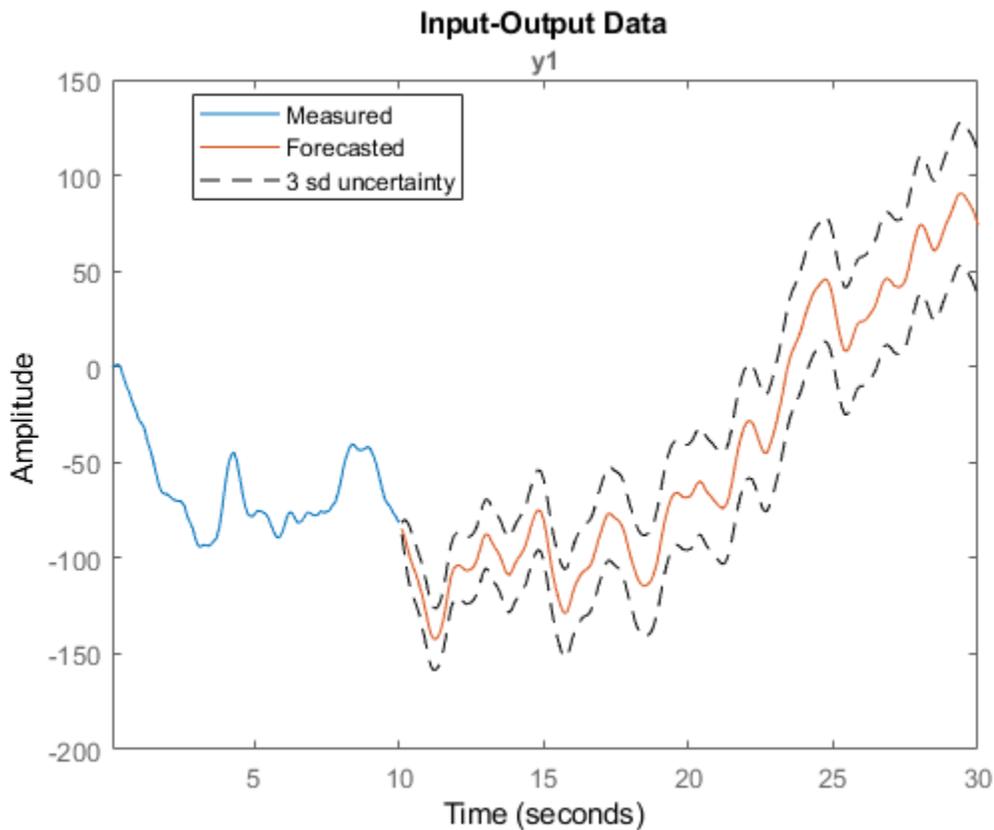
Forecast the system response into the future for a given time horizon and future inputs.

```
K = 200;
[yf,x0,sysf,yf_sd,x,x_sd] = forecast(sys,past_data,K,future_inputs);
```

`yf` is the forecasted model response, and `yf_sd` is the standard deviation of the output. `x0` is the estimated value for initial states, and `sysf` is the forecasting state-space model. Also returned are the state trajectory, `x`, and standard deviation of the trajectory, `x_sd`.

Plot the forecasted response.

```
UpperBound = iddata(yf.OutputData+3*yf_sd,[],yf.Ts,'Tstart',yf.Tstart);
LowerBound = iddata(yf.OutputData-3*yf_sd,[],yf.Ts,'Tstart',yf.Tstart);
plot(past_data(:,:,[]),yf(:,:,[]),UpperBound,'k--',LowerBound,'k--')
legend({'Measured','Forecasted','3 sd uncertainty'},'Location','best')
```

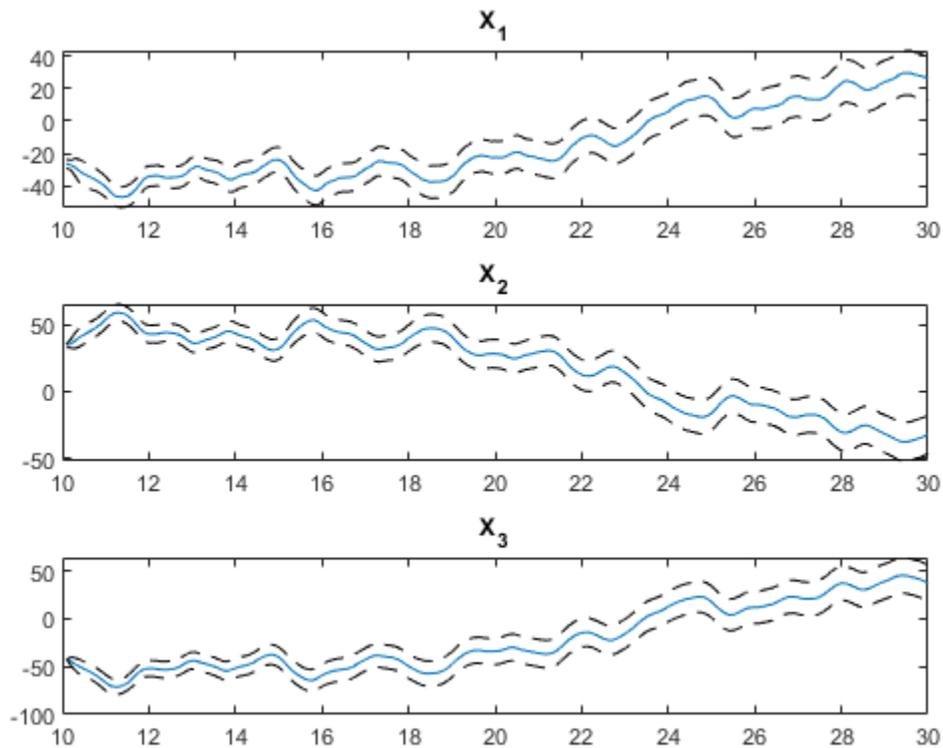


Plot the state trajectory.

```
t = z1.SamplingInstants(101:end);
subplot(3,1,1)
plot(t,x(:,1),t,x(:,1)+3*x_sd(:,1), 'k--',t,x(:,1)-3*x_sd(:,1), 'k--')
title('X_1')

subplot(3,1,2)
plot(t, x(:,2),t,x(:,2)+3*x_sd(:,2), 'k--',t, x(:,2)-3*x_sd(:,2), 'k--')
title('X_2')

subplot(3,1,3)
plot(t,x(:,3),t,x(:,3)+3*x_sd(:,3), 'k--',t, x(:,3)-3*x_sd(:,3), 'k--')
title('X_3')
```



The response uncertainty does not grow over the forecasting time span because of the specification of future inputs.

## Forecast Response of Multi-Output Nonlinear Time Series Model

Load data.

```
load(fullfile(matlabroot,'toolbox','ident','iddemos','data','predprey2data'));  
z = iddata(y,[],0.1);  
set(z,'Tstart',0,'OutputUnit',{'Population (in thousands)',...  
    'Population (in thousands)'},'TimeUnit','Years');
```

$z$  is a two output time-series data set (no inputs) from a 1-predator 1-prey population. The population exhibits a decline in predator population due to crowding. The data set contains 201 data samples covering 20 years of evolution.

The changes in the predator ( $y_1$ ) and prey ( $y_2$ ) population can be represented as:

$$y_1(t) = p_1 * y_1(t - 1) + p_2 * y_1(t - 1) * y_2(t - 1)$$

$$y_2(t) = p_3 * y_2(t - 1) - p_4 * y_1(t - 1) * y_2(t - 1) - p_5 * y_2(t - 1)^2$$

The nonlinearity in the predator and prey populations can be fit using a nonlinear ARX model with custom regressors.

Use part of the data as past data.

```
past_data = z(1:100);
```

Specify the standard regressors.

```
na = [1 0; 0 1];
```

```
nb = [];
```

```
nk = [];
```

Specify the custom regressors.

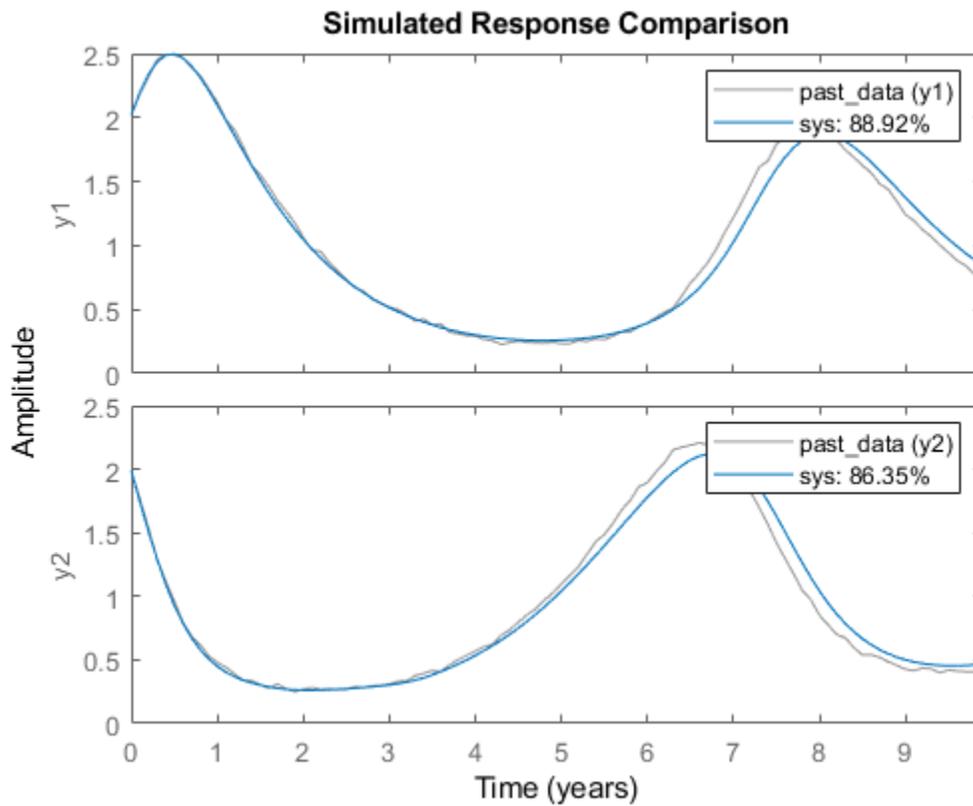
```
C = {'y1(t-1)*y2(t-1)'};{'y1(t-1)*y2(t-1)', 'y2(t-1)^2'};
```

Estimate a nonlinear ARX model using `past_data` as estimation data.

```
sys = nlarx(past_data,[na nb nk], 'wavenet', 'CustomRegressors',C);
```

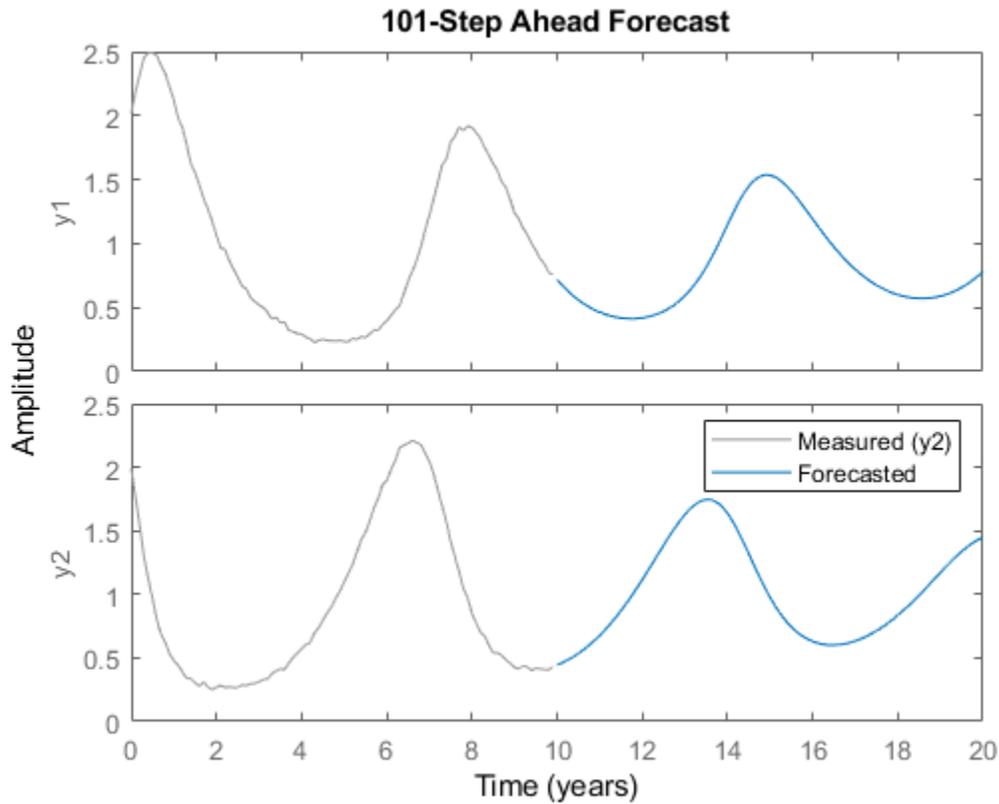
Compare the simulated output of `sys` with measured data to ensure it is a good fit.

```
compare(past_data,sys);
```



Plot the forecasted output of sys.

```
forecast(sys,past_data,101);  
legend('Measured','Forecasted');
```



### Reproduce Forecasting Results by Simulation

Obtain past data, future inputs, and identified linear model.

```
load iddata3 z3
past_data = z3(1:100);
future_inputs = z3.u(101:end);
sys = polyest(z3,[2 2 2 0 0 1]);
```

Forecast the system response into the future for a given time horizon and future inputs.

```
K = size(future_inputs,1);  
[yf,x0,sysf] = forecast(sys,past_data,K,future_inputs);
```

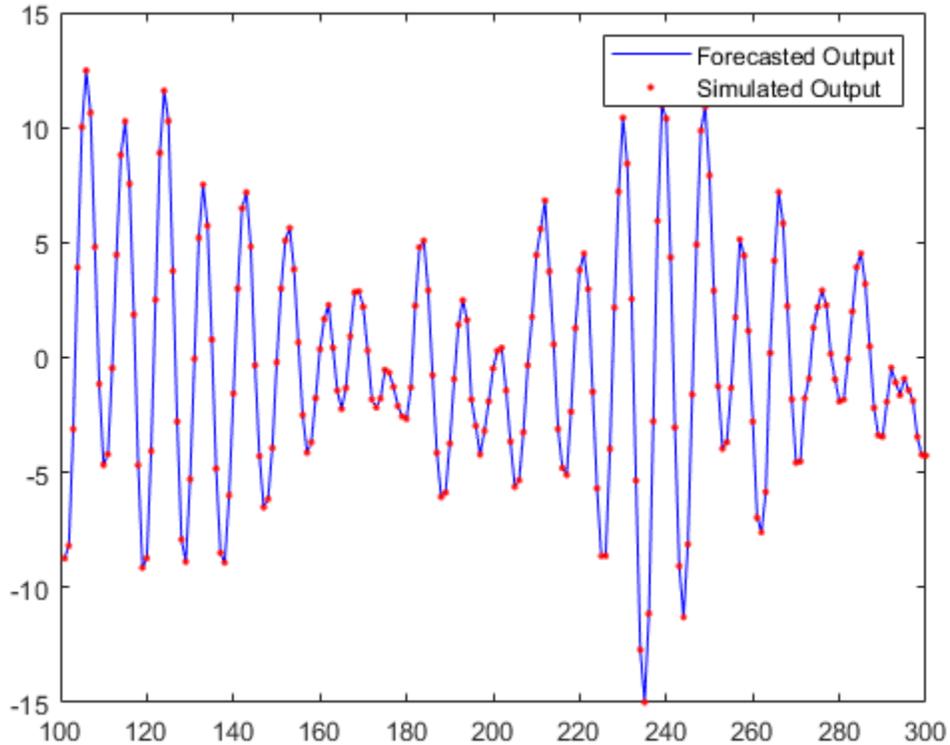
`yf` is the forecasted model response, `x0` is the estimated value for initial states, and `sysf` is the forecasting state-space model.

Simulate the forecasting state-space model with inputs, `future_inputs`, and initial conditions, `x0`.

```
opt = simOptions;  
opt.InitialCondition = x0;  
ys = sim(sysf,future_inputs(1:K),opt);
```

Plot the forecasted and simulated outputs.

```
t = yf.SamplingInstants;  
plot(t,yf.OutputData,'b',t,ys,'.r');  
legend('Forecasted Output','Simulated Output')
```



Simulation of forecasting model, `sysf`, with inputs, `future_inputs`, and initial conditions, `x0`, yields the forecasted output, `yf`.

## Input Arguments

### **sys** — Identified model

linear model | nonlinear model

Identified model whose output is to be forecasted, specified as one of the following:

- Linear model — `idpoly`, `idproc`, `idss`, `idtf`, or `idgrey`

- Nonlinear model — `idnlgrey`, `idnlhw`, or `idnlarx`

If a model is unavailable, estimate `sys` from `PastData` using commands such as `ar`, `arx`, `armax`, `nlarx`, and `ssest`.

## **PastData — Past input-output time-domain data**

`iddata` object | matrix of doubles

Past input-output time-domain data, specified as one of the following:

- `iddata` object — Use observed input and output signals to create an `iddata` object. For time-series data (no inputs), specify as an `iddata` object with no inputs `iddata(output, [])`.
- Matrix of doubles — For discrete-time models only. Specify as an  $N$ -by- $N_y$  matrix for time-series data. Here,  $N$  is the number of observations and  $N_y$  is the number of outputs.

For models with  $N_u$  inputs, specify `PastData` as an  $N$ -by- $(N_y+N_u)$  matrix.

## **K — Time horizon of forecasting**

positive integer

Time horizon of forecasting, specified as a positive integer. The output, `yf`, is calculated  $K$  steps into the future, such that the prediction time horizon is  $T_s * K$ .

## **FutureInputs — Future input values**

`[]` | matrix of doubles | `iddata` object | cell array of matrices

Future input values, specified as one of the following:

- `[]` — Future input values are assumed to be zero, or equal to input offset levels (if they are specified in `opts`). For time series models, specify as `[]`.
- `iddata` object — Specify as an `iddata` object with no outputs.
- $K$ -by- $N_u$  matrix of doubles —  $K$  is the forecast horizon, and  $N_u$  is the number of inputs.

If you have data from multiple experiments, you can specify a cell array of matrices, one for each experiment in `PastData`.

## **opts — Forecast options**

`forecastOptions` option set

Forecast options, specified as a `forecastOptions` option set.

**Linespec — Line style, marker, and color**

character vector

Line style, marker, and color, specified as a character vector. For example, 'b' or 'b+:'.

For more information about configuring Linespec, see the Linespec argument of `plot`.

## Output Arguments

**yf — Forecasted response**

iddata object

Forecasted response, returned as an `iddata` object. `yf` is the forecasted response at times after the last sample time in `PastData`. `yf` contains data for the time interval  $T_0 + (N+1:N+K)*T_1$ , where  $T_0 = \text{PastData.Tstart}$  and  $T_1 = \text{PastData.Ts}$ .  $N$  is the number of samples in `PastData`.

**x0 — Estimated initial states**

column vector | cell array

Estimated initial states at the start of forecasting, returned as a column vector of size equal to the number of states. Use `x0` with the forecasting model `sysf` to reproduce the result of forecasting by pure simulation.

If `PastData` is multi-experiment, `x0` is a cell array of size  $N_e$ , where  $N_e$  is the number of experiments.

When `sys` is not a state-space model (`idss`, `idgrey`, or `idnlgrey`), the definition of states depends on if `sys` is linear or nonlinear:

- Linear model (`idpoly`, `idproc`, `idtf`) - `sys` is converted to a discrete-time state-space model, and `x0` is returned as the states of the converted model at a time-point beyond the last data in `PastData`.

If conversion of `sys` to `idss` is not possible, `x0` is returned empty. For example, if `sys` is a MIMO continuous-time model with irreducible internal delays.

- Nonlinear model (`idnlhw` or `idnlarx`) — For a definition of the states of `idnlarx` and `idnlhw` models, see “Definition of `idnlarx` States” on page 1-631, and “Definition of `idnlhw` States” on page 1-672.

**sysf — Forecasting model**discrete-time `idss` | `idnlarx` | `idnlhw` | `idnlgrey` | cell array of models

Forecasting model, returned as one of the following:

- Discrete-time `idss` — If `sys` is a discrete-time `idss` model, `sysf` is the same as `sys`. If `sys` is a linear model that is not a state-space model (`idpoly`, `idproc`, `idtf`), or is a continuous-time state-space model (`idss`, `idgrey`), `sys` is converted to a discrete-time `idss` model. The converted model is returned in `sysf`.
- `idnlarx`, `idnlhw`, or `idnlgrey`— If `sys` is a nonlinear model, `sysf` is the same as `sys`.
- Cell array of models — If `PastData` is multiexperiment, `sysf` is an array of  $N_e$  models, where  $N_e$  is the number of experiments.

Simulation of `sysf` using `sim`, with inputs, `FutureInputs`, and initial conditions, `x0`, yields `yf` as the output. For time-series models, `FutureInputs` is empty.

**yf\_sd — Estimated standard deviations of forecasted response**

matrix | cell array

Estimated standard deviations of forecasted response, returned as a  $K$ -by- $N_y$  matrix, where  $K$  is the forecast horizon, and  $N_y$  is the number of outputs. The software computes the standard deviation by taking into account the model parameter covariance, initial state covariance, and additive noise covariance. The additive noise covariance is stored in the `NoiseVariance` property of the model.

If `PastData` is multiexperiment, `yf_sd` is a cell array of size  $N_e$ , where  $N_e$  is the number of experiments.

`yf_sd` is empty if `sys` is a nonlinear ARX (`idnlarx`) or Hammerstein-Wiener model (`idnlhw`). `yf_sd` is also empty if `sys` does not contain parameter covariance information, that is if `getcov(sys)` is empty. For more information, see `getcov`.

**x — Forecasted state trajectory**

matrix | cell array

Forecasted state trajectory, returned as a  $K$ -by- $N_x$  matrix, where  $K$ , the forecast horizon and  $N_x$  is the number of states. `x` are the states of the forecasting model.

If `PastData` is multiexperiment, `x` is a cell array of size  $N_e$ , where  $N_e$  is the number of experiments.

If `sys` is linear model other than a state-space model (not `idss` or `idgrey`), then it is converted to a discrete-time state-space model, and the states of the converted model are calculated. If conversion of `sys` to `idss` is not possible, `x` is returned empty. For example, if `sys` is a MIMO continuous-time model with irreducible internal delays.

`x` is empty if `sys` is a nonlinear ARX (`idnlarx`) or Hammerstein-Wiener model (`idnlhw`).

### **`x_sd` — Estimated standard deviations of forecasted states**

matrix | cell array

Estimated standard deviations of forecasted states `x`, returned as a  $K$ -by- $N_s$  matrix, where  $K$ , the forecast horizon and  $N_s$  is the number of states. The software computes the standard deviation by taking into account the model parameter covariance, initial state covariance, and additive noise covariance. The additive noise covariance is stored in the `NoiseVariance` property of the model.

If `PastData` is multiexperiment, `x_sd` is a cell array of size  $N_e$ , where  $N_e$  is the number of experiments.

If `sys` is linear model other than a state-space model (not `idss` or `idgrey`), then it is converted to a discrete-time state-space model, and the states and standard deviations of the converted model are calculated. If conversion of `sys` to `idss` is not possible, `x_sd` is returned empty. For example, if `sys` is a MIMO continuous-time model with irreducible internal delays.

`x_sd` is empty if `sys` is a nonlinear ARX (`idnlarx`) or Hammerstein-Wiener model (`idnlhw`).

## Tips

- Right-clicking the plot opens the context menu, where you can access the following options:
  - **Systems** — Select systems to view forecasted output. By default, the forecasted output of all systems is plotted.
  - **Data Experiment** — For multi-experiment data only. Toggle between data from different experiments.
  - **Characteristics** — View the following data characteristics:

- **Peak Value** — View peak value of the data.
- **Mean Value** — View mean value of the data.
- **Confidence Region** — View the estimated standard deviation of the forecasted output. To specify number of standard deviations to plot, double-click the plot and open the Property Editor dialog box. Specify the number of standard deviations in the **Options** tab, in **Confidence Region for Identified Models**. The default value is 1 standard deviation.

The confidence region is not generated for nonlinear ARX and Hammerstein-Wiener models and models that do not contain parameter covariance information.

- **Show Past Data** — Plot the past output data used for forecasting. By default, the past output data is plotted.
- **I/O Grouping** — For datasets containing more than one input or output channel. Select grouping of input and output channels on the plot.
  - **None** — Plot input-output channels in their own separate axes.
  - **All** — Group all input channels together and all output channels together.
- **I/O Selector** — For datasets containing more than one input or output channel. Select a subset of the input and output channels to plot. By default, all output channels are plotted.
- **Grid** — Add grids to the plot.
- **Normalize** — Normalize the y-scale of all data in the plot.
- **Full View** — Return to full view. By default, the plot is scaled to full view.
- **Properties** — Open the Property Editor dialog box to customize plot attributes.

## See Also

[ar](#) | [arx](#) | [compare](#) | [forecastOptions](#) | [iddata](#) | [predict](#) | [sim](#) | [ssest](#)

## Topics

“Forecast the Output of a Dynamic System”

“Forecast Multivariate Time Series”

“Time Series Prediction and Forecasting for Prognosis”

“Introduction to Forecasting of Dynamic System Response”

**Introduced in R2012a**

## forecastOptions

Option set for forecast

### Syntax

```
opt = forecastOptions  
opt = forecastOptions(Name,Value)
```

### Description

`opt = forecastOptions` creates the default option set for `forecast`. Use dot notation to modify this option set. Any options that you do not modify retain their default values.

`opt = forecastOptions(Name,Value)` creates an option set with the options specified by one or more `Name,Value` pair arguments.

### Examples

#### Specify Input Signal Offset for Model Forecasting

Create a default option set for `forecast`.

```
opt = forecastOptions;
```

Specify the input offset for a single-input data set as 5.

```
opt.InputOffset = 5;
```

You can now use this option set for forecasting. Before forecasting model response, the `forecast` command subtracts this offset value from the past input data signal.

### Specify Handling of Initial Conditions During Model Forecasting

Create an option set for `forecast` using zero initial conditions.

```
opt = forecastOptions('InitialCondition','z');
```

### Specify Output Offset for Forecasting Multi-Experiment Data

Load past measured data from two experiments.

```
load iddata1
load iddata2
```

`z1` and `z2` are `iddata` objects that store SISO input-output data. Create a two-experiment data set from `z1` and `z2`.

```
z = merge(z1,z2);
```

Estimate a transfer function model with 2 poles using the multi-experiment data.

```
sys = tfest(z,2);
```

Specify the offset as -1 and 1 for the output signals of the two experiments.

```
opt = forecastOptions('OutputOffset',[-1 1]);
```

`OutputOffset` is specified as an  $N_y$ -by- $N_e$  matrix where  $N_y$  is the number of outputs in each experiment, and  $N_e$  is the number of experiments. In this example,  $N_y$  is 1 and  $N_e$  is 2.

Using the option set `opt`, forecast the response of the model 10 time steps into the future. The software subtracts the offset value `OutputOffset(i,j)` from the output signal  $i$  of experiment  $j$  before using the data in the forecasting algorithm. The removed offsets are added back to generate the final result.

```
y = forecast(sys,z,10,opt)
```

```
y =
Time domain data set containing 2 experiments.
```

Experiment	Samples	Sample Time
Exp1	10	0.1

```
    Exp2          10          0.1
Outputs
  y1          Unit (if specified)
Inputs
  u1          Unit (if specified)
```

`y` is an `iddata` object that returns the forecasted response corresponding to each set of past experimental data.

## Input Arguments

### Name-Value Pair Arguments

Specify optional comma-separated pairs of `Name`, `Value` arguments. `Name` is the argument name and `Value` is the corresponding value. `Name` must appear inside quotes. You can specify several name and value pair arguments in any order as `Name1, Value1, ..., NameN, ValueN`.

Example: `forecastOptions('InitialCondition', 'e')` specifies that the software estimates the initial conditions of the measured input-output data such that the 1-step prediction error for observed output is minimized.

#### **InitialCondition — Handling of initial conditions**

'e' (default) | 'z' | `idpar` object `x00bj`

Handling of initial conditions, specified as the comma-separated pair consisting of 'InitialCondition' and one of the following values:

- 'z' — Zero initial conditions.
- 'e' — Estimate initial conditions such that the 1-step prediction error is minimized for the observed output.

For nonlinear grey-box models, only those initial states `i` that are designated as free in the model (`sys.InitialStates(i).Fixed = false`) are estimated. To estimate all the states of the model, first specify all the `Nx` states of the `idnlgrey` model `sys` as free.

```

for i = 1:Nx
sys.InitialStates(i).Fixed = false;
end

```

Similarly, to fix all the initial states to values specified in `sys.InitialStates`, first specify all the states as fixed in the `sys.InitialStates` property of the nonlinear grey-box model.

- `x0obj` — Specification object created using `idpar`. Use this object for discrete-time state-space models only (`idss`, `idgrey`, and `idnlgrey`). Use `x0obj` to impose constraints on the initial states by fixing their value or specifying minimum or maximum bounds.

### InputOffset — Input signal offset

`[]` (default) | column vector | matrix

Input signal offset for time-domain data, specified as the comma-separated pair consisting of 'InputOffset' and one of the following values:

- `[]` — No input offsets.
- A column vector of length  $Nu$ , where  $Nu$  is the number of inputs. When you use the `forecast` command, the software subtracts the offset value `InputOffset(i)` from the  $i$ th input signals in the past and future input values. You specify these values in the `PastData` and `FutureInputs` arguments of `forecast`. The software then uses the offset subtracted inputs to forecast the model response.
- $Nu$ -by- $Ne$  matrix — For multi-experiment data, specify `InputOffset` as an  $Nu$ -by- $Ne$  matrix, where  $Ne$  is the number of experiments. The software subtracts the offset value `InputOffset(i, j)` from the  $i$ th input signal of the  $j$ th experiment in the `PastData` and `FutureInputs` arguments of `forecast` before forecasting.

### OutputOffset — Output signal offset

`[]` (default) | column vector | matrix

Output signal offset for time-domain data, specified as the comma-separated pair consisting of 'OutputOffset' and one of the following values:

- `[]` — No output offsets.
- A column vector of length  $Ny$ , where  $Ny$  is the number of outputs. When you use the `forecast` command, the software subtracts the offset value `OutputOffset(i)` from the  $i$ th past output signal that you specify in the `PastData` argument of `forecast`. The software then uses the offset subtracted output to compute the detrended

forecasts. The removed offsets are added back to the detrended forecasts to generate the final result.

- *N<sub>y</sub>*-by-*N<sub>e</sub>* matrix — For multi-experiment data, specify `OutputOffset` as an *N<sub>y</sub>*-by-*N<sub>e</sub>* matrix, where *N<sub>e</sub>* is the number of experiments. Before forecasting, the software subtracts the offset value `OutputOffset(i, j)` from the *i*th output signal of the *j*th experiment in the `PastData` argument of `forecast`. For an example, see “Specify Output Offset for Forecasting Multi-Experiment Data” on page 1-431.

## Output Arguments

### **opt** — Option set for forecast

`forecastOptions` option set

Option set for forecast, returned as a `forecastOptions` option set.

## See Also

`forecast` | `idpar`

## Topics

“Introduction to Forecasting of Dynamic System Response”

**Introduced in R2012a**

## fpe

Akaike's Final Prediction Error for estimated model

### Syntax

```
value = fpe(model)
value = fpe(model1, ..., modeln)
```

### Description

`value = fpe(model)` returns the Final Prediction Error (FPE) value for the estimated model.

`value = fpe(model1, ..., modeln)` returns the FPE value for multiple estimated models.

### Examples

#### Compute Final Prediction Error of Estimated Model

Estimate a transfer function model.

```
load iddata1 z1;
np = 2;
sys = tfest(z1,np);
```

Compute the Final Prediction Error (FPE) value.

```
value = fpe(sys)
value = 1.7252
```

Alternatively, use the `Report` property of the model to access the value.

```
sys.Report.Fit.FPE
```

```
ans = 1.7252
```

## Pick Model with Optimal Tradeoff Between Accuracy and Complexity Using FPE Criterion

Estimate multiple Output-Error (OE) models and use Akaike's Final Prediction Error (FPE) value to pick the one with optimal tradeoff between accuracy and complexity.

Load the estimation data.

```
load iddata2
```

Specify model orders varying in 1:4 range.

```
nf = 1:4;  
nb = 1:4;  
nk = 0:4;
```

Estimate OE models with all possible combinations of chosen order ranges.

```
NN = struc(nf,nb,nk);  
models = cell(size(NN,1),1);  
for ct = 1:size(NN,1)  
    models{ct} = oe(z2, NN(ct,:));  
end
```

Compute the small sample-size corrected AIC values for the models, and return the smallest value.

```
V = fpe(models{:});  
[Vmin, I] = min(V);
```

Return the optimal model that has the smallest AICc value.

```
models{I}
```

```
ans =  
Discrete-time OE model:  $y(t) = [B(z)/F(z)]u(t) + e(t)$   
     $B(z) = 1.067 z^{-2}$   
  
     $F(z) = 1 - 1.824 z^{-1} + 1.195 z^{-2} - 0.2307 z^{-3}$ 
```

Sample time: 0.1 seconds

Parameterization:

Polynomial orders: nb=1 nf=3 nk=2

Number of free coefficients: 4

Use "polydata", "getpvec", "getcov" for parameters and their uncertainties.

Status:

Estimated using OE on time domain data "z2".

Fit to estimation data: 86.53%

FPE: 0.9809, MSE: 0.9615

## Input Arguments

**model** — Identified model

idtf | idgrey | idpoly | idproc | idss | idnlarx, | idnlhw | idnlgrey

Identified model, specified as one of the following model objects:

- idtf
- idgrey
- idpoly
- idproc
- idss
- idnlarx, except nonlinear ARX model that includes a binary-tree or neural network nonlinearity estimator
- idnlhw
- idnlgrey

## Output Arguments

**value** — Final Prediction Error (FPE) value

scalar | vector

Final Prediction Error (FPE) value, returned as a scalar or vector. For multiple models, value is a row vector where value(k) corresponds to the kth estimated model modelk.

## Definitions

### Akaike's Final Prediction Error (FPE)

Akaike's Final Prediction Error (FPE) criterion provides a measure of model quality by simulating the situation where the model is tested on a different data set. After computing several different models, you can compare them using this criterion. According to Akaike's theory, the most accurate model has the smallest FPE.

If you use the same data set for both model estimation and validation, the fit always improves as you increase the model order and, therefore, the flexibility of the model structure.

Akaike's Final Prediction Error (FPE) is defined by the following equation:

$$FPE = \det \left( \frac{1}{N} \sum_{t=1}^N e(t, \theta_N) (e(t, \theta_N))^T \right) \begin{pmatrix} 1 + d/N \\ 1 - d/N \end{pmatrix}$$

where:

- $N$  is the number of values in the estimation data set.
- $e(t)$  is a  $ny$ -by-1 vector of prediction errors.
- $\theta_N$  represents the estimated parameters.
- $d$  is the number of estimated parameters.

If number of parameters exceeds the number of samples, FPE is not computed when model estimation is performed (`model.Report.FPE` is empty). The `fpe` command returns NaN.

## Tips

- The software computes and stores the FPE value during model estimation. If you want to access this value, see the `Report.Fit.FPE` property of the model.

## References

- [1] Ljung, L. *System Identification: Theory for the User*, Upper Saddle River, NJ, Prentice-Hall PTR, 1999. See sections 7.4 and 16.4.

## See Also

aic | goodnessOfFit

## Topics

“Loss Function and Model Quality Metrics”

**Introduced before R2006a**

## frdata

Access data for frequency response data (FRD) object

### Syntax

```
[response, freq] = frdata(sys)
[response, freq, covresp] = frdata(sys)
[response, freq, Ts, covresp] = frdata(sys, 'v')
[response, freq, Ts] = frdata(sys)
```

### Description

`[response, freq] = frdata(sys)` returns the response data and frequency samples of the FRD model `sys`. For an FRD model with `Ny` outputs and `Nu` inputs at `Nf` frequencies:

- `response` is an `Ny-by-Nu-by-Nf` multidimensional array where the  $(i, j)$  entry specifies the response from input  $j$  to output  $i$ .
- `freq` is a column vector of length `Nf` that contains the frequency samples of the FRD model.

See the `frd` reference page for more information on the data format for FRD response data.

`[response, freq, covresp] = frdata(sys)` also returns the covariance `covresp` of the response data `resp` for `idfrd` model `sys`. The covariance `covresp` is a 5D-array where `covH(i, j, k, :, :)` contains the 2-by-2 covariance matrix of the response `resp(i, j, k)`. The  $(1, 1)$  element is the variance of the real part, the  $(2, 2)$  element the variance of the imaginary part and the  $(1, 2)$  and  $(2, 1)$  elements the covariance between the real and imaginary parts.

For SISO FRD models, the syntax

```
[response, freq] = frdata(sys, 'v')
```

forces `frdata` to return the response data as a column vector rather than a 3-dimensional array (see example below). Similarly

```
[response, freq, Ts, covresp] = frdata(sys, 'v')
```

 for an IDFRD model `sys` returns `covresp` as a 3-dimensional rather than a 5-dimensional array.

```
[response, freq, Ts] = frdata(sys)
```

 also returns the sample time `Ts`.

Other properties of `sys` can be accessed with `get` or by direct structure-like referencing (e.g., `sys.Frequency`).

## Arguments

The input argument `sys` to `frdata` must be an FRD model.

## Examples

### Extract Data from Frequency Response Data Model

Create a frequency response data model by computing the response of a transfer function on a grid of frequencies.

```
H = tf([-1.2, -2.4, -1.5], [1, 20, 9.1]);  
w = logspace(-2, 3, 101);  
sys = frd(H, w);
```

`sys` is a SISO frequency response data (`frd`) model containing the frequency response at 101 frequencies.

Extract the frequency response data from `sys`.

```
[response, freq] = frdata(sys);
```

`response` is a 1-by-1-by-101 array. `response(1, 1, k)` is the complex frequency response at the frequency `freq(k)`.

## **See Also**

frd | freqresp | get | idfrd | set | spectrum

**Introduced before R2006a**

# freqresp

Frequency response over grid

## Syntax

```
[H,wout] = freqresp(sys)
H = freqresp(sys,w)
H = freqresp(sys,w,units)
[H,wout,covH] = freqresp(idsys,...)
```

## Description

`[H,wout] = freqresp(sys)` returns the frequency response on page 1-447 of the dynamic system model `sys` at frequencies `wout`. The `freqresp` command automatically determines the frequencies based on the dynamics of `sys`.

`H = freqresp(sys,w)` returns the frequency response on page 1-447 on the real frequency grid specified by the vector `w`.

`H = freqresp(sys,w,units)` explicitly specifies the frequency units of `w` with `units`.

`[H,wout,covH] = freqresp(idsys,...)` also returns the covariance `covH` of the frequency response of the identified model `idsys`.

## Input Arguments

### `sys`

Any dynamic system model or model array.

### `w`

Vector of real frequencies at which to evaluate the frequency response. Specify frequencies in units of `rad/TimeUnit`, where `TimeUnit` is the time units specified in the `TimeUnit` property of `sys`.

## units

Units of the frequencies in the input frequency vector  $w$ , specified as one of the following values:

- 'rad/TimeUnit' — radians per the time unit specified in the TimeUnit property of `sys`
- 'cycles/TimeUnit' — cycles per the time unit specified in the TimeUnit property of `sys`
- 'rad/s'
- 'Hz'
- 'kHz'
- 'MHz'
- 'GHz'
- 'rpm'

**Default:** 'rad/TimeUnit'

## idsys

Any identified model.

## Output Arguments

### H

Array containing the frequency response values.

If `sys` is an individual dynamic system model having  $N_y$  outputs and  $N_u$  inputs, `H` is a 3D array with dimensions  $N_y$ -by- $N_u$ -by- $N_w$ , where  $N_w$  is the number of frequency points. Thus, `H(:, :, k)` is the response at the frequency  $w(k)$  or `wout(k)`.

If `sys` is a model array of size  $[N_y N_u S_1 \dots S_n]$ , `H` is an array with dimensions  $N_y$ -by- $N_u$ -by- $N_w$ -by- $S_1$ -by-...-by- $S_n$  array.

If `sys` is a frequency response data model (such as `frd`, `genfrd`, or `idfrd`), `freqresp(sys, w)` evaluates to `NaN` for values of  $w$  falling outside the frequency interval defined by `sys.frequency`. The `freqresp` command can interpolate between

frequencies in `sys.frequency`. However, `freqresp` cannot extrapolate beyond the frequency interval defined by `sys.frequency`.

### **wout**

Vector of frequencies corresponding to the frequency response values in `H`. If you omit `w` from the inputs to `freqresp`, the command automatically determines the frequencies of `wout` based on the system dynamics. If you specify `w`, then `wout = w`

### **covH**

Covariance of the response `H`. The covariance is a 5D array where `covH(i, j, k, :, :)` contains the 2-by-2 covariance matrix of the response from the  $i$ th input to the  $j$ th output at frequency  $w(k)$ . The (1,1) element of this 2-by-2 matrix is the variance of the real part of the response. The (2,2) element is the variance of the imaginary part. The (1,2) and (2,1) elements are the covariance between the real and imaginary parts of the response.

## Examples

### Compute Frequency Response of System

Create the following 2-input, 2-output system:

$$\mathit{sys} = \begin{bmatrix} 0 & \frac{1}{s+1} \\ \frac{s-1}{s+2} & 1 \end{bmatrix}$$

```
sys11 = 0;
sys22 = 1;
sys12 = tf(1,[1 1]);
sys21 = tf([1 -1],[1 2]);
sys = [sys11,sys12;sys21,sys22];
```

Compute the frequency response of the system.

```
[H,wout] = freqresp(sys);
```

`H` is a 2-by-2-by-45 array. Each entry `H(:, :, k)` in `H` is a 2-by-2 matrix giving the complex frequency response of all input-output pairs of `sys` at the corresponding frequency

wout(k). The 45 frequencies in wout are automatically selected based on the dynamics of sys.

### Compute Frequency Response on Specified Frequency Grid

Create the following 2-input, 2-output system:

$$\text{sys} = \begin{bmatrix} 0 & \frac{1}{s+1} \\ \frac{s-1}{s+2} & 1 \end{bmatrix}$$

```
sys11 = 0;  
sys22 = 1;  
sys12 = tf(1,[1 1]);  
sys21 = tf([1 -1],[1 2]);  
sys = [sys11,sys12;sys21,sys22];
```

Create a logarithmically-spaced grid of 200 frequency points between 10 and 100 radians per second.

```
w = logspace(1,2,200);
```

Compute the frequency response of the system on the specified frequency grid.

```
H = freqresp(sys,w);
```

H is a 2-by-2-by-200 array. Each entry  $H(:, :, k)$  in H is a 2-by-2 matrix giving the complex frequency response of all input-output pairs of sys at the corresponding frequency  $w(k)$ .

### Compute Frequency Response and Associated Covariance

Compute the frequency response and associated covariance for an identified process model at its peak response frequency.

Load estimation data z1.

```
load iddata1 z1
```

Estimate a SISO process model using the data.

```
model = procest(z1, 'P2UZ');
```

Compute the frequency at which the model achieves the peak frequency response gain. To get a more accurate result, specify a tolerance value of  $1e-6$ .

```
[gpeak, fpeak] = getPeakGain(model, 1e-6);
```

Compute the frequency response and associated covariance for `model` at its peak response frequency.

```
[H, wout, covH] = freqresp(model, fpeak);
```

`H` is the response value at `fpeak` frequency, and `wout` is the same as `fpeak`.

`covH` is a 5-dimensional array that contains the covariance matrix of the response from the input to the output at frequency `fpeak`. Here `covH(1, 1, 1, 1, 1)` is the variance of the real part of the response, and `covH(1, 1, 1, 2, 2)` is the variance of the imaginary part. The `covH(1, 1, 1, 1, 2)` and `covH(1, 1, 1, 2, 1)` elements are the covariance between the real and imaginary parts of the response.

## Definitions

### Frequency Response

In continuous time, the frequency response at a frequency  $\omega$  is the transfer function value at  $s = j\omega$ . For state-space models, this value is given by

$$H(j\omega) = D + C(j\omega I - A)^{-1}B$$

In discrete time, the frequency response is the transfer function evaluated at points on the unit circle that correspond to the real frequencies. `freqresp` maps the real

frequencies  $w(1), \dots, w(N)$  to points on the unit circle using the transformation  $z = e^{j\omega T_s}$ .  $T_s$  is the sample time. The function returns the values of the transfer function at the resulting  $z$  values. For models with unspecified sample time, `freqresp` uses  $T_s = 1$ .

## Algorithms

For transfer functions or zero-pole-gain models, `freqresp` evaluates the numerator(s) and denominator(s) at the specified frequency points. For continuous-time state-space models  $(A, B, C, D)$ , the frequency response is

$$D + C(j\omega - A)^{-1}B, \quad \omega = \omega_1, \dots, \omega_N$$

For efficiency,  $A$  is reduced to upper Hessenberg form and the linear equation  $(j\omega - A)X = B$  is solved at each frequency point, taking advantage of the Hessenberg structure. The reduction to Hessenberg form provides a good compromise between efficiency and reliability. See [1] for more details on this technique.

## Alternatives

Use `evalfr` to evaluate the frequency response at individual frequencies or small numbers of frequencies. `freqresp` is optimized for medium-to-large vectors of frequencies.

## References

- [1] Laub, A.J., "Efficient Multivariable Frequency Response Computations," *IEEE Transactions on Automatic Control*, AC-26 (1981), pp. 407-408.

## See Also

`bode` | `evalfr` | `interp` | `nichols` | `nyquist` | `sigma` | `spectrum`

**Introduced before R2006a**

## fselect

Select frequency points or range in FRD model

### Syntax

```
subsys = fselect(sys, fmin, fmax)
subsys = fselect(sys, index)
```

### Description

`subsys = fselect(sys, fmin, fmax)` takes an FRD model `sys` and selects the portion of the frequency response between the frequencies `fmin` and `fmax`. The selected range `[fmin, fmax]` should be expressed in the FRD model units. For an IDFRD model, the `SpectrumData`, `CovarianceData` and `NoiseCovariance` values, if non-empty, are also selected in the chosen range.

`subsys = fselect(sys, index)` selects the frequency points specified by the vector of indices `index`. The resulting frequency grid is

```
sys.Frequency(index)
```

### See Also

`fcats` | `fdel` | `frd` | `idfrd` | `interp`

**Introduced before R2006a**

## get

Access model property values

## Syntax

```
Value = get(sys, 'PropertyName')  
Struct = get(sys)
```

## Description

`Value = get(sys, 'PropertyName')` returns the current value of the property `PropertyName` of the model object `sys`. `'PropertyName'` can be the full property name (for example, `'UserData'`) or any unambiguous case-insensitive abbreviation (for example, `'user'`). See reference pages for the individual model object types for a list of properties available for that model.

`Struct = get(sys)` converts the TF, SS, or ZPK object `sys` into a standard MATLAB structure with the property names as field names and the property values as field values.

Without left-side argument,

```
get(sys)
```

displays all properties of `sys` and their values.

## Examples

### Display Model Property Values

Create the following discrete-time SISO transfer function model:

$$H(z) = \frac{1}{z+2}$$

Specify the sample time as 0.1 seconds and input channel name as Voltage.

```
h = tf(1,[1 2],0.1,'InputName','Voltage')
```

```
h =
```

```
From input "Voltage" to output:
```

```
  1
-----
z + 2
```

```
Sample time: 0.1 seconds
Discrete-time transfer function.
```

Display all the properties of the transfer function.

```
get(h)
```

```

    Numerator: {[0 1]}
  Denominator: {[1 2]}
      Variable: 'z'
      IODelay: 0
    InputDelay: 0
    OutputDelay: 0
           Ts: 0.1000
    TimeUnit: 'seconds'
    InputName: {'Voltage'}
    InputUnit: {''}
    InputGroup: [1x1 struct]
    OutputName: {''}
    OutputUnit: {''}
    OutputGroup: [1x1 struct]
           Notes: [0x1 string]
    UserData: []
           Name: ''
    SamplingGrid: [1x1 struct]
```

Display the numerator of the transfer function.

```
num = get(h,'Numerator')
```

```
num = 1x1 cell array
      {1x2 double}
```

The numerator data is stored as a cell array, thus the `Numerator` property is a cell array containing the row vector `[0 1]`.

```
num{1}
ans = 1x2
      0     1
```

Display the sample time `Ts` of the transfer function.

```
get(h, 'Ts')
ans = 0.1000
```

Alternatively, use dot notation to access the property value.

```
h.Ts
ans = 0.1000
```

## Tips

An alternative to the syntax

```
Value = get(sys, 'PropertyName')
```

is the structure-like referencing

```
Value = sys.PropertyName
```

For example,

```
sys.Ts
sys.A
sys.user
```

return the values of the sample time, `A` matrix, and `UserData` property of the (state-space) model `sys`.

## See Also

frdata | getcov | getpvec | idssdata | polydata | set | ssdata | tfdata

**Introduced before R2006a**

## getcov

Parameter covariance of identified model

### Syntax

```
cov_data = getcov(sys)
cov_data = getcov(sys,cov_type)
cov_data = getcov(sys,cov_type,'free')
```

### Description

`cov_data = getcov(sys)` returns the raw covariance of the parameters of an identified model.

- If `sys` is a single model, then `cov_data` is an  $np$ -by- $np$  matrix.  $np$  is the number of parameters of `sys`.
- If `sys` is a model array, then `cov_data` is a cell array of size equal to the array size of `sys`.

`cov_data(i,j,k,...)` contains the covariance data for `sys(:, :, i, j, k, ...)`.

`cov_data = getcov(sys,cov_type)` returns the parameter covariance as either a matrix or a structure, depending on the covariance type that is specified.

`cov_data = getcov(sys,cov_type,'free')` returns the covariance data of only the free model parameters.

### Examples

#### Obtain Raw Parameter Covariance for Identified Model

Obtain the identified model.

```
load iddata1 z1
sys = tfest(z1,2);
```

Get the raw parameter covariance for the model.

```
cov_data = getcov(sys)
```

```
cov_data = 5x5
```

```
    1.2131   -4.3949   -0.0309   -0.5531     0
   -4.3949   115.0838    1.8598   10.6660     0
   -0.0309    1.8598    0.0636    0.1672     0
   -0.5531   10.6660    0.1672    1.2433     0
         0         0         0         0         0
```

`cov_data` contains the covariance matrix for the parameter vector [`sys.Numerator`, `sys.Denominator(2:end)`, `sys.IODelay`].

`sys.Denominator(1)` is fixed to 1 and not treated as a parameter. The covariance matrix entries corresponding to the delay parameter (fifth row and column) are zero because the delay was not estimated.

### Obtain Raw Parameter Covariance for Identified Model Array

Obtain the identified model array.

```
load iddata1 z1;
sys1 = tfest(z1,2);
sys2 = tfest(z1,3);
sysarr = stack(1,sys1,sys2);
```

`sysarr` is a 2-by-1 array of continuous-time, identified transfer functions.

Get the raw parameter covariance for the models in the array.

```
cov_data = getcov(sysarr)
```

```
cov_data = 2x1 cell array
    {5x5 double}
    {7x7 double}
```

cov\_data is a 2-by-1 cell array. cov\_data{1} and cov\_data{2} are the raw parameter covariance matrices for sys1 and sys2.

### Obtain Raw Covariance of Estimated Parameters of Identified Model

Load the estimation data.

```
load iddata1 z1
z1.y = cumsum(z1.y);
```

Estimate the model.

```
init_sys = idtf([100 1500],[1 10 10 0]);
init_sys.Structure.Numerator.Minimum = eps;
init_sys.Structure.Denominator.Minimum = eps;
init_sys.Structure.Denominator.Free(end) = false;
opt = tfestOptions('SearchMethod','lm');
sys = tfest(z1,init_sys,opt);
```

sys is an idtf model with six parameters, four of which are estimated.

Get the covariance matrix for the estimated parameters.

```
cov_type = 'value';
cov_data = getcov(sys,cov_type,'free')
```

```
cov_data = 4x4
105 x
```

```
    0.0269    -0.1237    -0.0001    -0.0017
   -0.1237    1.0221     0.0016     0.0133
   -0.0001     0.0016     0.0000     0.0000
   -0.0017     0.0133     0.0000     0.0002
```

cov\_data is a 4x4 covariance matrix, with entries corresponding to the four estimated parameters.

### Obtain Factored Parameter Covariance for Identified Model

Obtain the identified model.

```
load iddata1 z1
sys = tfest(z1,2);
```

Get the factored parameter covariance for the model.

```
cov_type = 'factors';
cov_data = getcov(sys,cov_type);
```

### Obtain Factored Parameter Covariance for Identified Model Array

Obtain the identified model array.

```
load iddata1 z1
sys1 = tfest(z1,2);
sys2 = tfest(z1,3);
sysarr = stack(1,sys1,sys2);
```

sysarr is a 2-by-1 array of continuous-time, identified transfer functions.

Get the factored parameter covariance for the models in the array.

```
cov_type = 'factors';
cov_data = getcov(sysarr,cov_type)
```

```
cov_data = 2x1 struct array with fields:
```

```
    R
    T
    Free
```

cov\_data is a 2-by-1 structure array. cov\_data(1) and cov\_data(2) are the factored covariance structures for sys1 and sys2.

### Obtain Factored Covariance of Estimated Parameters of Identified Model

Load the estimation data.

```
load iddata1 z1
z1.y = cumsum(z1.y);
```

Estimate the model.

```
init_sys = idtf([100 1500],[1 10 10 0]);
init_sys.Structure.Numerator.Minimum = eps;
init_sys.Structure.Denominator.Minimum = eps;
init_sys.Structure.Denominator.Free(end) = false;
opt = tfestOptions('SearchMethod','lm');
sys = tfest(z1,init_sys,opt);
```

sys, an idtf model, has six parameters, four of which are estimated.

Get the factored covariance for the estimated parameters.

```
cov_type = 'factors';
cov_data = getcov(sys,cov_type,'free');
```

## Input Arguments

### sys — Identified model

idtf, idss, idgrey, idpoly, idproc, idnlarx, idnlhw, or idnlgrey object | model array

Identified model, specified as an idtf, idss, idgrey, idpoly, idproc, idnlarx, idnlhw, or idnlgrey model or an array of such models.

The getcov command returns cov\_data as [] for idnlarx and idnlhw models because these models do not store parameter covariance data.

### cov\_type — Covariance type

'value' (default) | 'factors'

Covariance return type, specified as either 'value' or 'factors'.

- If cov\_type is 'value', then cov\_data is returned as a matrix (raw covariance).
- If cov\_type is 'factors', then cov\_data is returned as a structure containing the factors of the covariance matrix.

Use this option for fetching the covariance data if the covariance matrix contains nonfinite values, is not positive definite, or is ill conditioned. You can calculate the

response uncertainty using the covariance factors instead of the numerically disadvantageous covariance matrix.

This option does not offer a numerical advantage in the following cases:

- `sys` is estimated using certain instrument variable methods, such as `iv4`.
- You have explicitly specified the parameter covariance of `sys` using the deprecated `CovarianceMatrix` model property.

Data Types: `char`

## Output Arguments

### **cov\_data** — Parameter covariance of `sys`

matrix or cell array of matrices | structure or cell array of structures

Parameter covariance of `sys`, returned as a matrix, cell array of matrices, structure, or cell array of structures. `cov_data` is [] for `idnlarx` and `idnlhw` models.

- If `sys` is a single model and `cov_type` is 'value', then `cov_data` is an  $np$ -by- $np$  matrix.  $np$  is the number of parameters of `sys`.

The value of the nonzero elements of this matrix is equal to `sys.Report.Parameters.FreeParCovariance` when `sys` is obtained via estimation. The row and column entries that correspond to fixed parameters are zero.

- If `sys` is a single model and `cov_type` is 'factors', then `cov_data` is a structure with fields:
  - `R` — Usually an upper triangular matrix.
  - `T` — Transformation matrix.
  - `Free` — Logical vector of length  $np$ , indicating if a model parameter is free (estimated) or not.  $np$  is the number of parameters of `sys`.

To obtain the covariance matrix using the factored form, enter:

```
Free = cov_factored.Free;
T = cov_factored.T;
R = cov_factored.R;
np = nparams(sys);
```

```
cov_matrix = zeros(np);  
cov_matrix(Free, Free) = T*inv(R'*R)*T';
```

For numerical accuracy, calculate  $T*inv(R'*R)*T'$  as  $X*X'$ , where  $X = T/R$ .

- If `sys` is a model array, then `cov_data` is a cell array of size equal to the array size of `sys`.

`cov_data(i,j,k,...)` contains the covariance data for `sys(:, :, i, j, k, ...)`.

### See Also

[getpvec](#) | [nparams](#) | [rsample](#) | [setcov](#) | [sim](#) | [simsd](#)

### Topics

“What Is Model Covariance?”

“Types of Model Uncertainty Information”

**Introduced in R2012a**

# getDelayInfo

Get input/output delay information for `idnlarx` model structure

## Syntax

```
DELAYS = getDelayInfo(MODEL)
DELAYS = getDelayInfo(MODEL,TYPE)
```

## Description

`DELAYS = getDelayInfo(MODEL)` obtains the maximum delay in each input and output variable of an `idnlarx` model.

`DELAYS = getDelayInfo(MODEL,TYPE)` lets you choose between obtaining maximum delays across all input and output variables or maximum delays for each output variable individually. When delays are obtained for each output variable individually a matrix is returned, where each row is a vector containing  $n_y+n_u$  maximum delays for each output variable, and:

- $n_y$  is the number of outputs of `MODEL`.
- $n_u$  is the number of inputs of `MODEL`.

Delay information is useful for determining the number of states in the model. For nonlinear ARX models, the states are related to the set of delayed input and output variables that define the model structure (regressors). For example, if an input or output variable  $p$  has a maximum delay of  $D$  samples, then it contributes  $D$  elements to the state vector:

$$p(t-1), p(t-2), \dots, p(t-D)$$

The number of states of a nonlinear ARX model equals the sum of the maximum delays of each input and output variable. For more information about the definition of states for `idnlarx` models, see “Definition of `idnlarx` States” on page 1-631

## Input Arguments

getDelayInfo accepts the following arguments:

- **MODEL**: idnlarx model.
- **TYPE**: (Optional) Specifies whether to obtain channel delays 'channelwise' or 'all' as follows:
  - 'all': Default value. DELAYS contains the maximum delays across each output (vector of  $n_y+n_u$  entries, where  $[n_y, n_u] = \text{size}(\text{MODEL})$ ).
  - 'channelwise': DELAYS contains delay values separated for each output ( $n_y$ -by- $(n_y+n_u)$  matrix).

## Output Arguments

- **DELAYS**: Contains delay information in a vector of length  $n_y+n_u$  arranged with output channels preceding the input channels, i.e.,  $[y1, y2, \dots, u1, u2, \dots]$ .

## Examples

### Get Input and Output Delay Information for Nonlinear ARX Model

Create a two-output, three-input nonlinear ARX model.

```
M = idnlarx([2 0 2 2 1 1 0 0; 1 0 1 5 0 1 1 0], 'linear');
```

Compute the maximum delays for each output variable individually.

```
Del = getDelayInfo(M, 'channelwise')
```

```
Del = 2x5
```

```
     2     0     2     1     0  
     1     0     1     5     0
```

The matrix `Del` contains the maximum delays for the first and second output of model `M`. You can interpret the contents of matrix `Del` as follows:

- In the dynamics for output 1 ( $y_1$ ), the maximum delays in channels  $y_1$ ,  $y_2$ ,  $u_1$ ,  $u_2$ ,  $u_3$  are 2, 0, 2, 1, and 0 respectively.
- Similarly, in the dynamics for output 2 ( $y_2$ ) of the model, the maximum delays in channels  $y_1$ ,  $y_2$ ,  $u_1$ ,  $u_2$ ,  $u_3$  are 1, 0, 1, 5, and 0 respectively.

Find maximum delays for all the input and output variables in the order  $y_1$ ,  $y_2$ ,  $u_1$ ,  $u_2$ ,  $u_3$ .

```
Del = getDelayInfo(M, 'all')
```

```
Del = 1x5
```

```
     2     0     2     5     0
```

Note, The maximum delay across all output equations can be obtained by executing

$\text{MaxDel} = \max(\text{Del}, [], 1)$ . Since input  $u_2$  has 5 delays (the fourth entry in  $\text{Del}$ ), there are 5 terms corresponding to  $u_2$  in the state vector. Applying this definition to all I/O channels, the complete state vector for model  $M$  becomes:

$$X(t) = [y_1(t-1), y_1(t-2), u_1(t-1), u_1(t-2), u_2(t-1), u_2(t-2), u_2(t-3), u_2(t-4), u_2(t-5)]$$

## See Also

[data2state](#) | [getreg](#) | [idnlarx](#)

**Introduced in R2008b**

## getexp

Specific experiments from multiple-experiment data set

### Syntax

```
d1 = getexp(data,ExperimentNumber)
d1 = getexp(data,ExperimentName)
```

### Description

`d1 = getexp(data,ExperimentNumber)` retrieves specific experiments from multiple-experiment data set. `data` is an `iddata` object that contains several experiments. `d1` is another `iddata` object containing the indicated experiment(s). `ExperimentNumber` is the experiment number as in `d1 = getexp(data,3)` or `d1 = getexp(data,[4 2])`.

`d1 = getexp(data,ExperimentName)` specifies the experiment name as in `d1 = getexp(data,'Period1')` or `d1 = getexp(data,{'Day1','Day3'})`.

See `merge (iddata)` and `iddata` for how to create multiple-experiment data objects.

You can also retrieve the experiments using a fourth subscript, as in `d1 = data(:, :, :, ExperimentNumber)`. Type `help iddata/subsref` for details on this.

**Introduced before R2006a**

# getinit

Values of idnlgrey model initial states

## Syntax

```
getinit(model)  
getinit(model,prop)
```

## Arguments

model

Name of the idnlgrey model object.

Property

Name of the InitialStates model property field, such as 'Name', 'Unit', 'Value', 'Minimum', 'Maximum', and 'Fixed'.

Default: 'Value'.

## Description

getinit(model) gets the initial-state values in the 'Value' field of the InitialStates model property.

getinit(model,prop) gets the initial-state values of the prop field of the InitialStates model property. prop can be 'Name', 'Unit', 'Value', 'Minimum', 'Maximum', and 'Fixed'.

The returned values are an Nx-by-1 cell array of values, where Nx is the number of states.

## See Also

getpar | idnlgrey | setinit | setpar

**Introduced in R2007a**

# getoptions

Return @PlotOptions handle or plot options property

## Syntax

```
p = getoptions(h)
p = getoptions(h,propertyname)
```

## Description

`p = getoptions(h)` returns the plot options handle associated with plot handle `h`. `p` contains all the settable options for a given response plot.

`p = getoptions(h,propertyname)` returns the specified options property, `propertyname`, for the plot with handle `h`. You can use this to interrogate a plot handle. For example,

```
p = getoptions(h,'Grid')
```

returns 'on' if a grid is visible, and 'off' when it is not.

For a list of the properties and values available for each plot type, see “Properties and Values Reference” (Control System Toolbox).

## See Also

setoptions

**Introduced in R2012a**

## idParametric/getpar

Obtain attributes such as values and bounds of linear model parameters

### Syntax

```
value = getpar(sys, 'value')  
free = getpar(sys, 'free')  
bounds = getpar(sys, 'bounds')  
label = getpar(sys, 'label')  
getpar(sys)
```

### Description

`value = getpar(sys, 'value')` returns the parameter values of the model `sys`. If `sys` is a model array, the returned value is a cell array of size equal to the model array.

`free = getpar(sys, 'free')` returns the free or fixed status of the parameters.

`bounds = getpar(sys, 'bounds')` returns the minimum and maximum bounds on the parameters.

`label = getpar(sys, 'label')` returns the labels for the parameters.

`getpar(sys)` prints a table of parameter values, labels, free status and minimum and maximum bounds.

### Examples

#### Get Parameter Values

Get the parameter values of an estimated ARMAX model.

Estimate an ARMAX model.

```
load iddata8
init_data = z8(1:100);
na = 1;
nb = [1 1 1];
nc = 1;
nk = [0 0 0];
sys = armax(init_data,[na nb nc nk]);
```

Get the parameter values.

```
val = getpar(sys, 'value')
```

```
val = 5×1
```

```
-0.7519
-0.4341
 0.4442
 0.0119
 0.3431
```

To set parameter values, use `sys = setpar(sys, 'value', value)`.

### Get Free Parameters and Their Bounds

Get the free parameters and their bounds for a process model.

Construct a process model, and set its parameter values and free status.

```
m = idproc('P2DUZI');
m.Kp = 1;
m.Tw = 100;
m.Zeta = .3;
m.Tz = 10;
m.Td = 0.4;
m.Structure.Td.Free = 0;
```

Here, the value of Td is fixed.

Get the parameter values.

```
Val = getpar(m, 'Value')
```

```
Val = 5×1
```

```
1.0000  
100.0000  
0.3000  
0.4000  
10.0000
```

Get the free statuses of the parameters.

```
Free = getpar(m, 'Free')
```

```
Free = 5×1 logical array
```

```
1  
1  
1  
0  
1
```

The output indicates that Td is a fixed parameter and the remaining parameters are free.

Get the default bounds on the parameters.

```
MinMax = getpar(m, 'bounds')
```

```
MinMax = 5×2
```

```
-Inf   Inf  
0     Inf  
0     Inf  
0     Inf  
-Inf  Inf
```

Extract the values of the free parameters.

```
FreeValues = Val(Free)
```

```
FreeValues = 4×1
```

```
1.0000  
100.0000  
0.3000
```

```
10.0000
```

Extract the bounds on the free parameters.

```
FreeValBounds = MinMax(Free,:)
```

```
FreeValBounds = 4x2
```

```
-Inf   Inf
      0   Inf
      0   Inf
-Inf   Inf
```

### Get Parameter Labels

Get the parameter labels of an estimated ARMAX model.

Estimate an ARMAX model.

```
load iddata8;
init_data = z8(1:100);
na = 1;
nb = [1 1 1];
nc = 1;
nk = [0 0 0];
sys = armax(init_data,[na nb nc nk]);
```

Assign parameter labels.

```
sys.Structure.A.Info(2).Label = 'a2';
```

Get the parameter labels.

```
label = getpar(sys,'label')
```

```
label = 5x1 cell array
      {'a2'      }
      {0x0 char}
      {0x0 char}
      {0x0 char}
```

```
{0x0 char}
```

### Obtain a Table of Model Parameter Attributes

Obtain a table of all model parameter attributes of an ARMAX model.

Estimate an ARMAX model.

```
load iddata8;  
init_data = z8(1:100);  
na = 4;  
nb = [3 2 3];  
nc = 2;  
nk = [0 0 0];  
sys = armax(init_data,[na nb nc nk]);
```

Get all parameter attributes.

```
getpar(sys)
```

#	Label	Value	Free	Min.	Max.
1.		-1.4328	1	-Inf	Inf
2.		0.497	1	-Inf	Inf
3.		0.22904	1	-Inf	Inf
4.		-0.09849	1	-Inf	Inf
5.		-0.10246	1	-Inf	Inf
6.		1.1671	1	-Inf	Inf
7.		0.39579	1	-Inf	Inf
8.		0.97219	1	-Inf	Inf
9.		0.026995	1	-Inf	Inf
10.		-0.17113	1	-Inf	Inf
11.		0.16155	1	-Inf	Inf
12.		0.48468	1	-Inf	Inf
13.		-1.8871	1	-Inf	Inf
14.		0.97391	1	-Inf	Inf

## Input Arguments

### **sys** — Identified linear model

idss | idproc | idgrey | idtf | idpoly | array of model objects

Identified linear model, specified as an `idss`, `idpoly`, `idgrey`, `idtf`, or `idfrd` model object or an array of model objects.

## Output Arguments

### **value** — Parameter values

vector of doubles

Parameter values, returned as a double vector of length `nparams(sys)`.

### **free** — Free or fixed status of parameters

vector of logical values

Free or fixed status of parameters, returned as a logical vector of length `nparams(sys)`.

### **bounds** — Minimum and maximum bounds on parameters

matrix of doubles

Minimum and maximum bounds on parameters, returned as a double matrix of size `nparams(sys)-by-2`. The first column contains the minimum bound, and the second column the maximum bound.

### **label** — Parameter labels

cell array of character vectors

Parameter labels, returned as a cell array of character vectors of length `nparams(sys)`. For example, `{'a2', 'a3'}`, if `nparams(sys)` is two.

## See Also

getcov | getpvec | idssdata | polydata | setpar | tfdata

**Introduced in R2013b**

## getpar

Parameter values and properties of idnlgrey model parameters

### Syntax

```
getpar(model)  
getpar(model,prop)
```

### Arguments

model

Name of the idnlgrey model object.

Property

Name of the Parameters model property field, such as 'Name', 'Unit', 'Value', 'Minimum', 'Maximum', or 'Fixed'.

Default: 'Value'.

### Description

getpar(model) gets the model parameter values in the 'Value' field of the Parameters model property.

getpar(model,prop) gets the model parameter values in the prop field of the Parameters model property. prop can be 'Name', 'Unit', 'Value', 'Minimum', and 'Maximum'.

The returned values are an Np-by-1 cell array of values, where Np is the number of parameters.

### See Also

getinit | getpvec | idnlgrey | setinit | setpar

**Introduced in R2007a**

## getpvec

Model parameters and associated uncertainty data

### Syntax

```
pvec = getpvec(sys)
[pvec,pvec_sd] = getpvec(sys)
[___] = getpvec(sys,'free')
```

### Description

`pvec = getpvec(sys)` returns a vector, `pvec`, containing the values of all the parameters of the identified model `sys`.

`[pvec,pvec_sd] = getpvec(sys)` also returns the 1 standard deviation value of the uncertainty associated with the parameters of `sys`. If the model covariance information for `sys` is not available, `pvec_sd` is `[]`.

`[___] = getpvec(sys,'free')` returns data for only the free parameters of `sys`, using any of the output arguments in previous syntaxes. For `idnlarx` and `idnlhw` models, all parameters are treated as free.

### Input Arguments

#### **sys**

Identified model.

Identified model, specified as an `idtf`, `idss`, `idgrey`, `idpoly`, `idproc`, `idnlarx`, `idnlhw`, or `idnlgrey` model or an array of such models.

## Output Arguments

### **pvec**

Values of the parameters of `sys`.

If `sys` is an array of models, then `pvec` is a cell array with parameter value vectors corresponding to each model in `sys`. `pvec` is `[]` for `idnlarx` and `idnlhw` models that have not been estimated.

### **pvec\_sd**

1 standard deviation value of the parameters of `sys`.

If the model covariance information for `sys` is not available, `pvec_sd` is `[]`. Thus, `pvec_sd` is always `[]` for `idnlarx` and `idnlhw` models because these models do not store parameter covariance information.

If `sys` is an array of models, then `pvec_sd` is a cell array with standard deviation vectors corresponding to each model in `sys`.

## Examples

### **Retrieve Parameter Values from Estimated Model**

Load the estimation data.

```
load iddata1 z1;
```

Estimate a transfer function model.

```
sys = tfest(z1,3);
```

Retrieve the parameter values from the estimated model.

```
pvec = getpvec(sys);
```

## Retrieve Parameter Values and Standard Deviations from Estimated Model

Load the estimation data

```
load iddata2 z2;
```

Estimate a state-space model.

```
sys = ssest(z2,3);
```

Retrieve the model parameters, `pvec`, and associated standard deviations, `pvec_sd`, from the estimated model.

```
[pvec,pvec_sd] = getpvec(sys);
```

## Retrieve Values of Free Parameters from Estimated Model

Load the estimation data.

```
load iddata2 z2;
```

Estimate a state-space model.

```
sys = ssest(z2,3);
```

Retrieve the values of the free parameters from the estimated model.

```
pvec = getpvec(sys,'free');
```

## See Also

[getcov](#) | [idssdata](#) | [setpvec](#) | [tfdata](#) | [zpkdata](#)

**Introduced in R2012a**

## getreg

Regressor expressions and numerical values in nonlinear ARX model

### Syntax

```
Rs = getreg(model)
Rs = getreg(model,subset)
Rm = getreg(model,subset,data)
Rm = getreg(model,subset,data,init)
```

### Description

`Rs = getreg(model)` returns expressions for computing regressors in the nonlinear ARX model. `model` is an `idnlarx` object.

`Rs = getreg(model,subset)` returns regressor expressions for a specified subset of regressors.

`Rm = getreg(model,subset,data)` returns regressor values as a matrix for a specified subset of regressors.

`Rm = getreg(model,subset,data,init)` returns regressor values as matrices for a specified subset of regressors. The first `N` rows of each regressor matrix depend on the initial states `init`, where `N` is the maximum delay in the regressors (see `getDelayInfo`). For multiple-output models, `Rm` is a cell array of cell arrays.

### Input Arguments

`data`

`iddata` object containing measured data.

`init`

    Initial conditions of your data:

- 'z' (default) specifies zero initial state.
- Real column vector containing the initial state values. input and output data values at a time instant before the first sample in `data`. To create the initial state vector from the input-output data, use the `data2state` command. For multiple-experiment data, this is a matrix where each column specifies the initial state of the model corresponding to that experiment.
- `iddata` object containing input and output samples at time instants before to the first sample in `data`. When the `iddata` object contains more samples than the maximum delay in the model, only the most recent samples are used. The minimum number of samples required is equal to `max(getDelayInfo(model))`.

`model`

`iddata` object representing nonlinear ARX model.

`subset`

Subset of all regressors, specified as one of the following values:

- (Default) 'all' — All regressors.
- 'custom'—Only custom regressors.
- 'input'—Only standard regressors computed from input data.
- 'linear'—Only regressors not used in the nonlinear block.
- 'nonlinear'—Only regressors used in the nonlinear block.

---

**Note** You can use 'nl' as an abbreviation of 'nonlinear'.

---

- 'output'—Only regressors computed from output data.
- 'standard'—Only standard regressors (excluding any custom regressors).

## Output Arguments

`Rm`

Matrix of regressor values for all or a specified subset of regressors. Each matrix in `Rm` contains as many rows as there are data samples. For a model with `ny` outputs, `Rm` is an `ny`-by-1 cell array of matrices. When `data` contains multiple experiments, `Rm` is a cell array where each element corresponds to a matrix of regressor values for an experiment.

**Rs**

Regressor expressions represented as a cell array of character vectors. For a model with  $n_y$  outputs, **Rs** is an  $n_y$ -by-1 cell array of cell array of character vectors. For example, the expression 'u1(t-2)' computes the regressor by delaying the input signal **u1** by two time samples. Similarly, the expression 'y2(t-1)' computes the regressor by delaying the output signal **y2** by one time sample.

The order of regressors in **Rs** corresponds to regressor indices in the `idnlarx` object property `model.NonlinearRegressors`.

## Examples

### Get Regressor Expressions and Values, and Evaluate Predicted Model Output

Load sample data **u** and **y**.

```
load twotankdata;  
Ts = 0.2;
```

Sample time is 0.2 min

Create data object and use first 1000 samples for estimation.

```
z = iddata(y,u,Ts);  
ze = z(1:1000);
```

Estimate nonlinear ARX model.

```
model = nlarx(ze,[3 2 1]);
```

Get regressor expressions.

```
Rs = getreg(model);
```

Get regressor values.

```
Rm = getreg(model,'all',ze);
```

Evaluate model output for one-step-prediction.

```
Y = evaluate(model.Nonlinearity,Rm);
```

The previous result is equivalent to.

```
Y_p = predict(model, ze, 1, 'z');
```

### See Also

[addreg](#) | [customreg](#) | [evaluate](#) | [polyreg](#)

### Topics

“Identifying Nonlinear ARX Models”

**Introduced in R2007a**

# getStateEstimate

Extract best state estimate and covariance from particles

## Syntax

```
State = getStateEstimate(pf)
[State,StateCovariance] = getStateEstimate(pf)
```

## Description

`State = getStateEstimate(pf)` returns the best state estimate based on the current set of particles. The estimate is extracted based on the `StateEstimationMethod` property from the `particleFilter` object, `pf`.

`[State,StateCovariance] = getStateEstimate(pf)` also returns the covariance of the state estimate. The covariance is a measure of the uncertainty of the state estimate. Not all state estimation methods support covariance output. In this case, `getStateEstimate` returns `StateCovariance` as `[]`.

The `State` and `StateCovariance` information can directly be accessed as properties of the particle filter object `pf`, as `pf.State` and `pf.StateCovariance`. However, when both these quantities are needed, using the `getStateEstimation` method with two output arguments is more computationally efficient.

## Examples

### State Estimation using Particle Filter

Create a particle filter, and set the state transition and measurement likelihood functions.

```
myPF = particleFilter(@vdpParticleFilterStateFcn,@vdpMeasurementLikelihoodFcn);
```

Initialize the particle filter at state `[2; 0]` with unit covariance, and use 1000 particles.

```
initialize(myPF, 1000, [2;0], eye(2));
```

Pick the mean state estimation and systematic resampling methods.

```
myPF.StateEstimationMethod = 'mean';  
myPF.ResamplingMethod = 'systematic';  
myPF
```

```
myPF =  
  particleFilter with properties:  
  
    NumStateVariables: 2  
      NumParticles: 1000  
    StateTransitionFcn: @vdpParticleFilterStateFcn  
    MeasurementLikelihoodFcn: @vdpMeasurementLikelihoodFcn  
    IsStateVariableCircular: [0 0]  
      ResamplingPolicy: [1x1 particleResamplingPolicy]  
      ResamplingMethod: 'systematic'  
    StateEstimationMethod: 'mean'  
      StateOrientation: 'column'  
        Particles: [2x1000 double]  
          Weights: [1x1000 double]  
            State: 'Use the getStateEstimate function to see the value.'  
              StateCovariance: 'Use the getStateEstimate function to see the value.'
```

Assuming a measurement 2.1, run one predict and correct step.

```
[PredictedState, PredictedStateCovariance] = predict(myPF);  
[CorrectedState, CorrectedStateCovariance] = correct(myPF,2.1);
```

Get the best state estimate and covariance based on the StateEstimationMethod property.

```
[State, StateCovariance] = getStateEstimate(myPF)
```

```
State = 2x1
```

```
    2.1063  
   -0.1162
```

```
StateCovariance = 2x2
```

```
    0.0169    0.0091
```

0.0091 0.7130

## Input Arguments

### **pf** — Particle filter

particleFilter object

Particle filter, specified as an object. See `particleFilter` for more information.

## Output Arguments

### **State** — Best state estimate

[ ] (default) | vector

Best state estimate, defined as a vector based on the condition of the `StateOrientation` property:

- If `StateOrientation` is 'row' then `State` is a 1-by-`NumStateVariables` vector
- If `StateOrientation` is 'column' then `State` is a `NumStateVariables`-by-1 vector

### **StateCovariance** — Current estimate of state estimation error covariance

`NumStateVariables`-by-`NumStateVariables` array (default) | [ ] | array

Current estimate of state estimation error covariance, defined as an `NumStateVariables`-by-`NumStateVariables` array. `StateCovariance` is calculated based on the `StateEstimationMethod`. If you specify a state estimation method that does not support covariance, then the function returns `StateCovariance` as [ ].

## See Also

`correct` | `initialize` | `particleFilter` | `predict`

## Topics

“Nonlinear State Estimation Using Unscented Kalman Filter and Particle Filter”

“What Is Online Estimation?”

**Introduced in R2017b**

# getTrend

Data offset and trend information

## Syntax

```
T = getTrend(data)
T = getTrend(data,0)
T = getTrend(data,1)
```

## Description

`T = getTrend(data)` constructs a `TrendInfo` object to store offset, mean, or linear trend information for detrending or retrending data. You can assign specific offset and slope values to `T`.

`T = getTrend(data,0)` computes the means of input and output signals and stores them as `InputOffset` and `OutputOffset` properties of `T`, respectively.

`T = getTrend(data,1)` computes a best-fit straight line for both input and output signals and stores them as properties of `T`.

## Examples

### Remove Offsets From Data

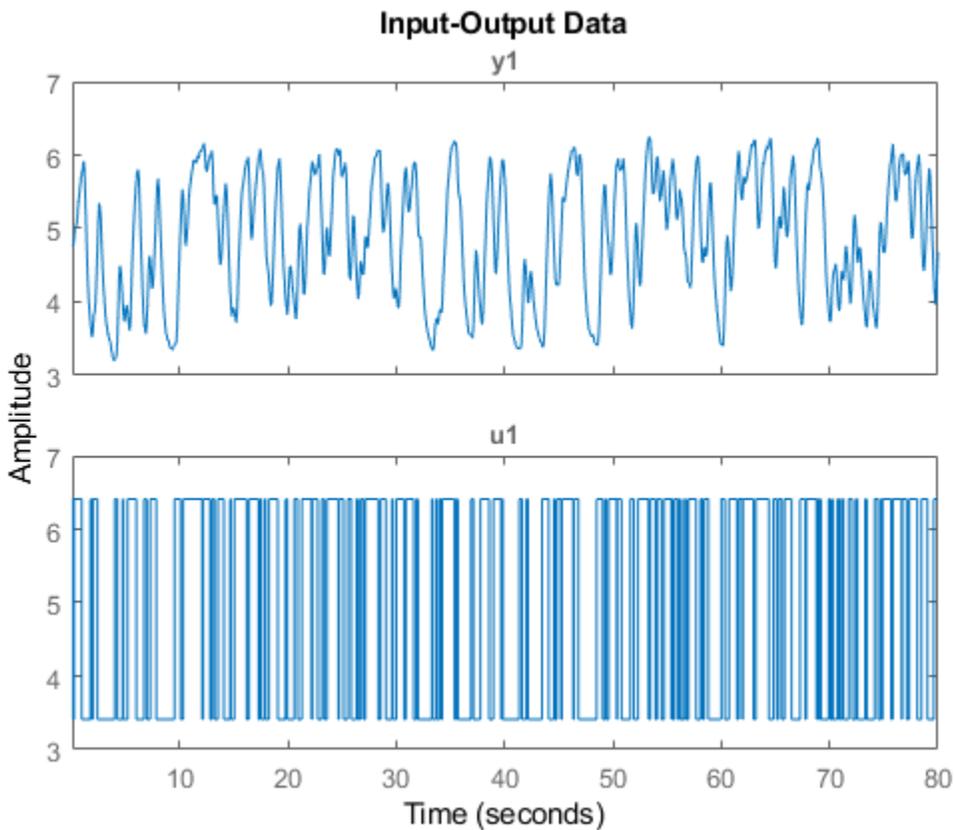
Remove specified offset from input and output signals.

Load SISO data containing vectors `u2` and `y2`.

```
load dryer2
```

Create a data object with sample time of 0.08 seconds and plot it.

```
data = iddata(y2,u2,0.08);
plot(data)
```



The data has a nonzero mean value.

Store the data offset and trend information in a `TrendInfo` object.

```
T = getTrend(data);
```

Assign offset values to the `TrendInfo` object.

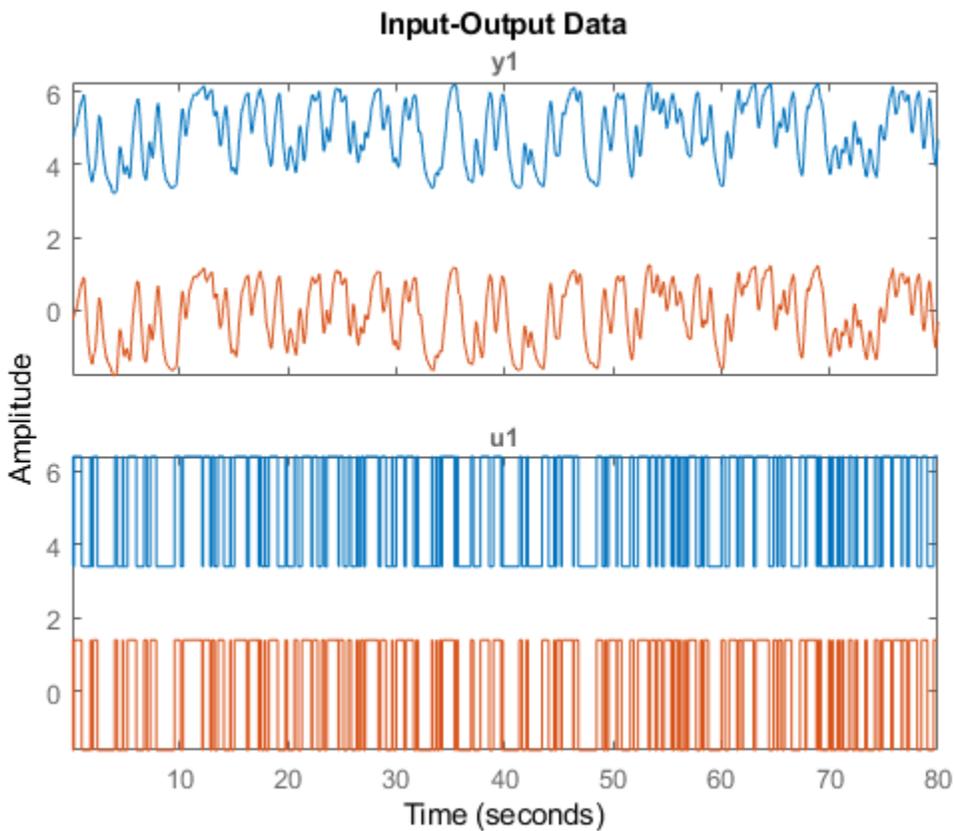
```
T.InputOffset = 5;  
T.OutputOffset = 5;
```

Subtract offset from the data.

```
data_d = detrend(data,T);
```

Plot the detrended data on the same plot.

```
hold on  
plot(data_d)
```



View the mean value removed from the data.

```
get(T)
```

```
ans = struct with fields:  
    DataName: 'data'  
    InputOffset: 5  
    OutputOffset: 5  
    InputSlope: 0
```

OutputSlope: 0

### **Compute and Store Means of Input and Output Signals**

Compute input-output signal means, store them, and detrend the data.

Load SISO data containing vectors `u2` and `y2`.

```
load dryer2
```

Create a data object with sample time of 0.08 seconds.

```
data = iddata(y2,u2,0.08);
```

Compute the mean of the data.

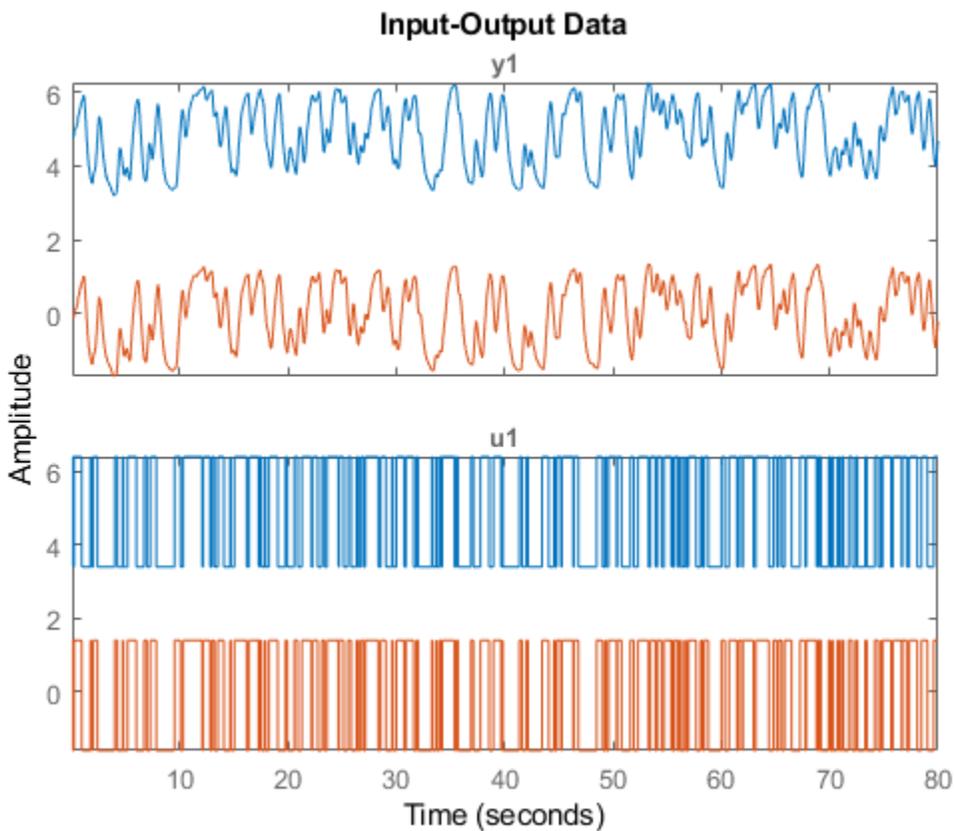
```
T = getTrend(data,0);
```

Remove the mean from the data.

```
data_d = detrend(data,T);
```

Plot the original and detrended data on the same plot.

```
plot(data,data_d)
```



## See Also

TrendInfo | detrend | retrend

## Topics

“Handling Offsets and Trends in Data”

**Introduced in R2009a**

## **goodnessOfFit**

Goodness of fit between test and reference data

### **Syntax**

```
fit = goodnessOfFit(x,xref,cost_func)
```

### **Description**

`fit = goodnessOfFit(x,xref,cost_func)` returns the goodness of fit between the data, `x`, and the reference, `xref` using a cost function specified by `cost_func`.

### **Input Arguments**

#### **x**

Test data.

`x` is an `Ns`-by-`N` matrix, where `Ns` is the number of samples and `N` is the number of channels.

`x` can also be a cell array of multiple test data sets.

`x` must not contain any `NaN` or `Inf` values.

#### **xref**

Reference data.

`xref` must be of the same size as `x`.

`xref` can also be a cell array of multiple reference sets. In this case, each individual reference set must be of the same size as the corresponding test data set.

`xref` must not contain any `NaN` or `Inf` values.

**cost\_func**

Cost function to determine goodness of fit.

cost\_func is specified as one of the following values:

- 'MSE' — Mean square error:

$$fit = \frac{\|x - xref\|^2}{N_s}$$

where,  $N_s$  is the number of samples, and  $\|$  indicates the 2-norm of a vector. *fit* is a scalar value.

- 'NRMSE' — Normalized root mean square error:

$$fit(i) = 1 - \frac{\|xref(:,i) - x(:,i)\|}{\|xref(:,i) - mean(xref(:,i))\|}$$

where,  $\|$  indicates the 2-norm of a vector. *fit* is a row vector of length  $N$  and  $i = 1, \dots, N$ , where  $N$  is the number of channels.

NRMSE costs vary between -Inf (bad fit) to 1 (perfect fit). If the cost function is equal to zero, then  $x$  is no better than a straight line at matching *xref*.

- 'NMSE' — Normalized mean square error:

$$fit(i) = 1 - \frac{\|xref(:,i) - x(:,i)\|^2}{\|xref(:,i) - mean(xref(:,i))\|^2}$$

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NMSE costs vary between -Inf (bad fit) to 1 (perfect fit). If the cost function is equal to zero, then  $x$  is no better than a straight line at matching *xref*.

## Output Arguments

### **fit**

Goodness of fit between test and reference data.

For a single test data set and reference pair, `fit` is returned as a:

- Scalar if `cost_func` is MSE.
- Row vector of length N if `cost_func` is NRMSE or NMSE. N is the number of channels.

If `x` and/or `xref` are cell arrays, then `fit` is an array containing the goodness of fit values for each test data and reference pair.

## Examples

### **Calculate Goodness of Fit of Between Estimated and Measured Data**

Obtain the measured output.

```
load iddata1 z1
yref = z1.y;
```

`z1` is an `iddata` object containing measured input/output data. `z1.y` is the measured output.

Obtain the estimated output.

```
sys = tfest(z1,2);
y_sim = sim(sys,z1(:,[],:));
```

`sys` is a second-order transfer function estimated using the measured input/output data. `y` is the output estimated using `sys` and the measured input.

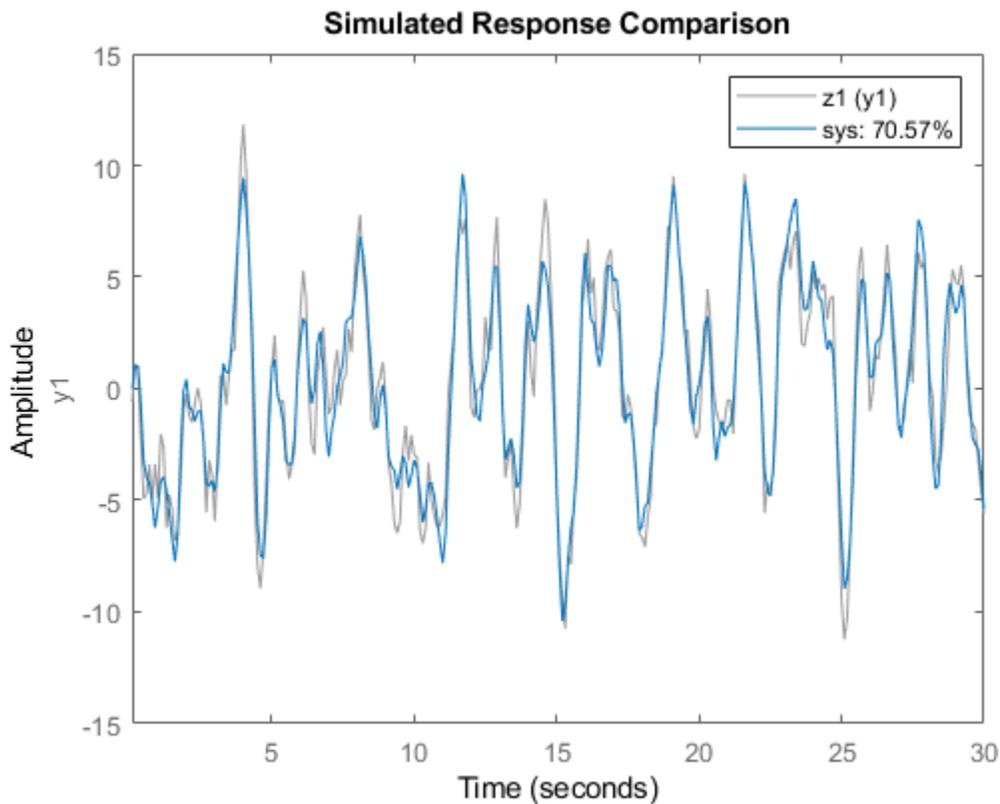
Calculate the goodness of the fit between the measured and estimated outputs.

```
cost_func = 'NRMSE';
y = y_sim.y;
fit = goodnessOfFit(y,yref,cost_func);
```

The goodness of fit is calculated using the normalized root mean square error as the cost function.

Alternatively, you can use `compare` to calculate the goodness of fit:

```
opt = compareOptions('InitialCondition','z');  
compare(z1,sys,opt);
```



## See Also

[aic](#) | [compare](#) | [fpe](#) | [pe](#) | [resid](#)

**Introduced in R2012a**

# greyest

Linear grey-box model estimation

## Syntax

```
sys = greyest(data,init_sys)
sys = greyest(data,init_sys,opt)
```

## Description

`sys = greyest(data,init_sys)` estimates a linear grey-box model, `sys`, using time or frequency domain data, `data`. The dimensions of the inputs and outputs of `data` and `init_sys`, an `idgrey` model, must match. `sys` is an identified `idgrey` model that has the same structure as `init_sys`.

`sys = greyest(data,init_sys,opt)` estimates a linear grey-box model using the option set, `opt`, to configure the estimation options.

## Input Arguments

### **data**

Estimation data.

The dimensions of the inputs and outputs of `data` and `init_sys` must match.

For time-domain estimation, `data` is an `iddata` object containing the input and output signal values.

For frequency domain estimation, `data` can be one of the following:

- Recorded frequency response data (`frd` or `idfrd`)
- `iddata` object with its `Domain` property set to 'Frequency'

**init\_sys**

Identified linear grey-box model that configures the initial parameterization of `sys`.

`init_sys`, an `idgrey` model, must have the same input and output dimensions as `data`.

**opt**

Estimation options.

`opt` is an option set, created using `greyestOptions`, which specifies options including:

- Estimation objective
- Initialization choice
- Disturbance model handling
- Numerical search method to be used in estimation

## Output Arguments

**sys**

Estimated grey-box model, returned as an `idgrey` model. This model is created using the specified initial system, and estimation options.

Information about the estimation results and options used is stored in the `Report` property of the model. `Report` has the following fields:

<b>Report Field</b>	<b>Description</b>
Status	Summary of the model status, which indicates whether the model was created by construction or obtained by estimation.
Method	Estimation command used.

Report Field	Description
InitialState	<p>Handling of initial states during estimation, returned as one of the following:</p> <ul style="list-style-type: none"> <li>• 'model' — The initial state is parameterized by the ODE file used by the idgrey model.</li> <li>• 'zero' — The initial state is set to zero.</li> <li>• 'estimate' — The initial state is treated as an independent estimation parameter.</li> <li>• 'backcast' — The initial state is estimated using the best least squares fit.</li> <li>• Vector of doubles of length <math>N_x</math>, where <math>N_x</math> is the number of states. For multiexperiment data, a matrix with <math>N_e</math> columns, where <math>N_e</math> is the number of experiments.</li> </ul> <p>This field is especially useful to view how the initial states were handled when the InitialState option in the estimation option set is 'auto'.</p>
DisturbanceModel	<p>Handling of the disturbance component (<math>K</math>) during estimation, returned as one of the following values:</p> <ul style="list-style-type: none"> <li>• 'model' — <math>K</math> values are parameterized by the ODE file used by the idgrey model.</li> <li>• 'fixed' — The value of the <math>K</math> property of the idgrey model is fixed to its original value.</li> <li>• 'none' — <math>K</math> is fixed to zero.</li> <li>• 'estimate' — <math>K</math> is treated as an independent estimation parameter.</li> </ul> <p>This field is especially useful to view the how the disturbance component was handled when the DisturbanceModel option in the estimation option set is 'auto'.</p>

Report Field	Description																		
Fit	<p>Quantitative assessment of the estimation, returned as a structure. See “Loss Function and Model Quality Metrics” for more information on these quality metrics. The structure has the following fields:</p> <table border="1" data-bbox="387 456 1340 951"> <thead> <tr> <th data-bbox="387 456 520 499">Field</th> <th data-bbox="520 456 1340 499">Description</th> </tr> </thead> <tbody> <tr> <td data-bbox="387 499 520 609">FitPercent</td> <td data-bbox="520 499 1340 609">Normalized root mean squared error (NRMSE) measure of how well the response of the model fits the estimation data, expressed as a percentage.</td> </tr> <tr> <td data-bbox="387 609 520 652">LossFcn</td> <td data-bbox="520 609 1340 652">Value of the loss function when the estimation completes.</td> </tr> <tr> <td data-bbox="387 652 520 730">MSE</td> <td data-bbox="520 652 1340 730">Mean squared error (MSE) measure of how well the response of the model fits the estimation data.</td> </tr> <tr> <td data-bbox="387 730 520 774">FPE</td> <td data-bbox="520 730 1340 774">Final prediction error for the model.</td> </tr> <tr> <td data-bbox="387 774 520 817">AIC</td> <td data-bbox="520 774 1340 817">Raw Akaike Information Criteria (AIC) measure of model quality.</td> </tr> <tr> <td data-bbox="387 817 520 861">AICc</td> <td data-bbox="520 817 1340 861">Small sample-size corrected AIC.</td> </tr> <tr> <td data-bbox="387 861 520 904">nAIC</td> <td data-bbox="520 861 1340 904">Normalized AIC.</td> </tr> <tr> <td data-bbox="387 904 520 951">BIC</td> <td data-bbox="520 904 1340 951">Bayesian Information Criteria (BIC).</td> </tr> </tbody> </table>	Field	Description	FitPercent	Normalized root mean squared error (NRMSE) measure of how well the response of the model fits the estimation data, expressed as a percentage.	LossFcn	Value of the loss function when the estimation completes.	MSE	Mean squared error (MSE) measure of how well the response of the model fits the estimation data.	FPE	Final prediction error for the model.	AIC	Raw Akaike Information Criteria (AIC) measure of model quality.	AICc	Small sample-size corrected AIC.	nAIC	Normalized AIC.	BIC	Bayesian Information Criteria (BIC).
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BIC	Bayesian Information Criteria (BIC).																		
Parameters	Estimated values of model parameters.																		
OptionsUsed	Option set used for estimation. If no custom options were configured, this is a set of default options. See <code>greyestOptions</code> for more information.																		
RandState	State of the random number stream at the start of estimation. Empty, [], if randomization was not used during estimation. For more information, see <code>rng</code> in the MATLAB documentation.																		

Report Field	Description																
DataUsed	Attributes of the data used for estimation, returned as a structure with the following fields:																
	<table border="1"> <thead> <tr> <th data-bbox="385 425 516 465">Field</th> <th data-bbox="516 425 1328 465">Description</th> </tr> </thead> <tbody> <tr> <td data-bbox="385 465 516 512">Name</td> <td data-bbox="516 465 1328 512">Name of the data set.</td> </tr> <tr> <td data-bbox="385 512 516 558">Type</td> <td data-bbox="516 512 1328 558">Data type.</td> </tr> <tr> <td data-bbox="385 558 516 605">Length</td> <td data-bbox="516 558 1328 605">Number of data samples.</td> </tr> <tr> <td data-bbox="385 605 516 652">Ts</td> <td data-bbox="516 605 1328 652">Sample time.</td> </tr> <tr> <td data-bbox="385 652 516 956">InterSample</td> <td data-bbox="516 652 1328 956">           Input intersample behavior, returned as one of the following values:           <ul style="list-style-type: none"> <li>• 'zoh' — Zero-order hold maintains a piecewise-constant input signal between samples.</li> <li>• 'foh' — First-order hold maintains a piecewise-linear input signal between samples.</li> <li>• 'bl' — Band-limited behavior specifies that the continuous-time input signal has zero power above the Nyquist frequency.</li> </ul> </td> </tr> <tr> <td data-bbox="385 956 516 1038">InputOffset</td> <td data-bbox="516 956 1328 1038">Offset removed from time-domain input data during estimation. For nonlinear models, it is [ ].</td> </tr> <tr> <td data-bbox="385 1038 516 1104">OutputOffset</td> <td data-bbox="516 1038 1328 1104">Offset removed from time-domain output data during estimation. For nonlinear models, it is [ ].</td> </tr> </tbody> </table>	Field	Description	Name	Name of the data set.	Type	Data type.	Length	Number of data samples.	Ts	Sample time.	InterSample	Input intersample behavior, returned as one of the following values: <ul style="list-style-type: none"> <li>• 'zoh' — Zero-order hold maintains a piecewise-constant input signal between samples.</li> <li>• 'foh' — First-order hold maintains a piecewise-linear input signal between samples.</li> <li>• 'bl' — Band-limited behavior specifies that the continuous-time input signal has zero power above the Nyquist frequency.</li> </ul>	InputOffset	Offset removed from time-domain input data during estimation. For nonlinear models, it is [ ].	OutputOffset	Offset removed from time-domain output data during estimation. For nonlinear models, it is [ ].
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<b>Report Field</b>	<b>Description</b>	
Termination	Termination conditions for the iterative search used for prediction error minimization. Structure with the following fields:	
	<b>Field</b>	<b>Description</b>
	WhyStop	Reason for terminating the numerical search.
	Iterations	Number of search iterations performed by the estimation algorithm.
	FirstOrderOptimality	$\infty$ -norm of the gradient search vector when the search algorithm terminates.
	FcnCount	Number of times the objective function was called.
	UpdateNorm	Norm of the gradient search vector in the last iteration. Omitted when the search method is 'lsqnonlin' or 'fmincon'.
	LastImprovement	Criterion improvement in the last iteration, expressed as a percentage. Omitted when the search method is 'lsqnonlin' or 'fmincon'.
	Algorithm	Algorithm used by 'lsqnonlin' or 'fmincon' search method. Omitted when other search methods are used.
For estimation methods that do not require numerical search optimization, the Termination field is omitted.		

For more information on using Report, see “Estimation Report”.

## Examples

### Estimate Grey-Box Model

Estimate the parameters of a DC motor using the linear grey-box framework.

Load the measured data.

```
load(fullfile(matlabroot, 'toolbox', 'ident', 'iddemos', 'data', 'dcmotordata'));
data = iddata(y, u, 0.1, 'Name', 'DC-motor');
data.InputName = 'Voltage';
data.InputUnit = 'V';
data.OutputName = {'Angular position', 'Angular velocity'};
data.OutputUnit = {'rad', 'rad/s'};
data.Tstart = 0;
data.TimeUnit = 's';
```

`data` is an `iddata` object containing the measured data for the outputs, the angular position, the angular velocity. It also contains the input, the driving voltage.

Create a grey-box model representing the system dynamics.

For the DC motor, choose the angular position (rad) and the angular velocity (rad/s) as the outputs and the driving voltage (V) as the input. Set up a linear state-space structure of the following form:

$$\dot{x}(t) = \begin{bmatrix} 0 & 1 \\ 0 & -\frac{1}{\tau} \end{bmatrix} x(t) + \begin{bmatrix} 0 \\ \frac{G}{\tau} \end{bmatrix} u(t)$$

$$y(t) = \begin{bmatrix} 1 & 0 \\ 0 & 1 \end{bmatrix} x(t).$$

$\tau$  is the time constant of the motor in seconds, and  $G$  is the static gain from the input to the angular velocity in rad/(V\*s) .

```
G = 0.25;
tau = 1;
```

```
init_sys = idgrey('motorDynamics',tau,'cd',G,0);
```

The governing equations in state-space form are represented in the MATLAB® file `motorDynamics.m`. To view the contents of this file, enter `edit motorDynamics.m` at the MATLAB command prompt.

$G$  is a known quantity that is provided to `motorDynamics.m` as an optional argument.

$\tau$  is a free estimation parameter.

`init_sys` is an `idgrey` model associated with `motor.m`.

Estimate  $\tau$ .

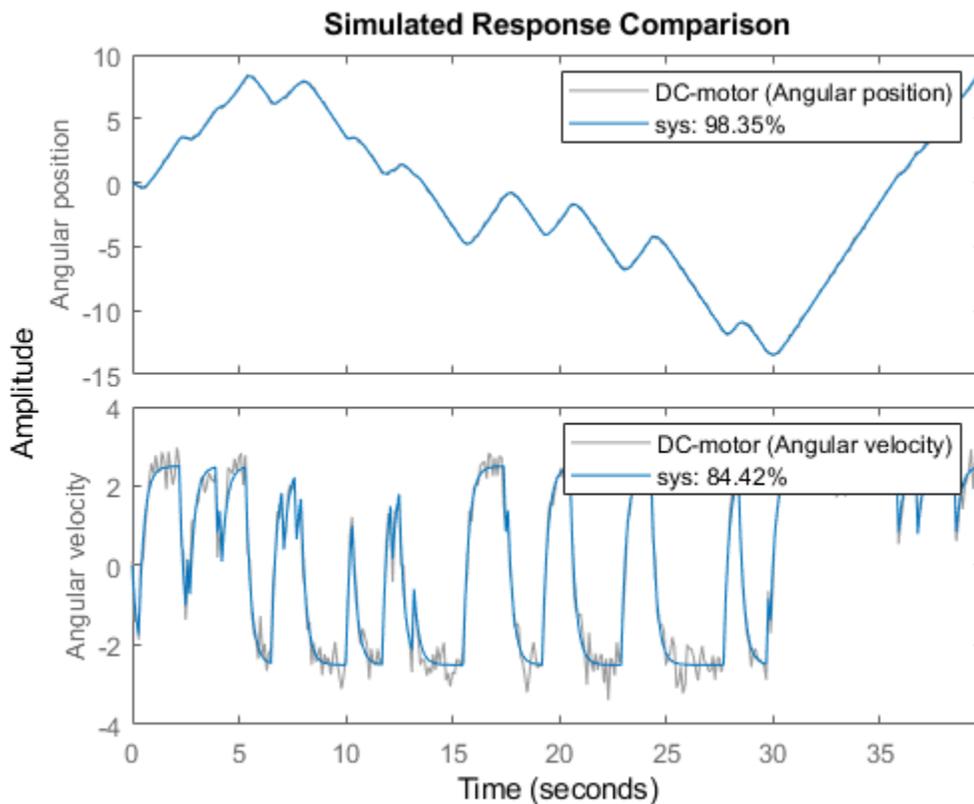
```
sys = greyest(data,init_sys);
```

`sys` is an `idgrey` model containing the estimated value of  $\tau$ .

To obtain the estimated parameter values associated with `sys`, use `getpvec(sys)`.

Analyze the result.

```
opt = compareOptions('InitialCondition','zero');  
compare(data,sys,Inf,opt)
```



sys provides a 98.35% fit for the angular position and an 84.42% fit for the angular velocity.

## Estimate Grey-Box Model Using Regularization

Estimate the parameters of a DC motor by incorporating prior information about the parameters when using regularization constants.

The model is parameterized by static gain  $G$  and time constant  $\tau$ . From prior knowledge, it is known that  $G$  is about 4 and  $\tau$  is about 1. Also, you have more confidence in the value of  $\tau$  than  $G$  and would like to guide the estimation to remain close to the initial guess.

Load estimation data.

```
load regularizationExampleData.mat motorData
```

The data contains measurements of motor's angular position and velocity at given input voltages.

Create an `idgrey` model for DC motor dynamics. Use the function `DCMotorODE` that represents the structure of the grey-box model.

```
mi = idgrey(@DCMotorODE,{'G', 4; 'Tau', 1},'cd',{}, 0);
mi = setpar(mi, 'label', 'default');
```

If you want to view the `DCMotorODE` function, type:

```
type DCMotorODE.m
```

```
function [A,B,C,D] = DCMotorODE(G,Tau,Ts)
%DCMOTORODE ODE file representing the dynamics of a DC motor parameterized
%by gain G and time constant Tau.
%
% [A,B,C,D,K,X0] = DCMOTORODE(G,Tau,Ts) returns the state space matrices
% of the DC-motor with time-constant Tau and static gain G. The sample
% time is Ts.
%
% This file returns continuous-time representation if input argument Ts
% is zero. If Ts>0, a discrete-time representation is returned.
%
```

```
% See also IDGREY, GREYEST.  
  
% Copyright 2013 The MathWorks, Inc.  
  
A = [0 1;0 -1/Tau];  
B = [0; G/Tau];  
C = eye(2);  
D = [0;0];  
if Ts>0 % Sample the model with sample time Ts  
    s = expm([[A B]*Ts; zeros(1,3)]);  
    A = s(1:2,1:2);  
    B = s(1:2,3);  
end
```

Specify regularization options Lambda.

```
opt = greyestOptions;  
opt.Regularization.Lambda = 100;
```

Specify regularization options R.

```
opt.Regularization.R = [1, 1000];
```

You specify more weighting on the second parameter because you have more confidence in the value of  $\tau$  than G.

Specify the initial values of the parameters as regularization option  $\theta^*$ .

```
opt.Regularization.Nominal = 'model';
```

Estimate the regularized grey-box model.

```
sys = greyest(motorData, mi, opt);
```

## Extended Capabilities

### Automatic Parallel Support

Accelerate code by automatically running computation in parallel using Parallel Computing Toolbox™.

Parallel computing support is available for estimation using the `lsqnonlin` search method (requires Optimization Toolbox). To enable parallel computing, use `greyestOptions`, set `SearchMethod` to `'lsqnonlin'`, and set `SearchOptions.Advanced.UseParallel` to `true`.

For example:

```
opt = greyestOptions;  
opt.SearchMethod = 'lsqnonlin';  
opt.SearchOptions.Advanced.UseParallel = true;
```

## See Also

`greyestOptions` | `iddata` | `idfrd` | `idgrey` | `idnlgrey` | `pem` | `ssest`

## Topics

“Estimate Model Using Zero/Pole/Gain Parameters”

“Regularized Estimates of Model Parameters”

**Introduced in R2012a**

## greyestOptions

Option set for greyest

### Syntax

```
opt = greyestOptions  
opt = greyestOptions(Name,Value)
```

### Description

`opt = greyestOptions` creates the default options set for `greyest`.

`opt = greyestOptions(Name,Value)` creates an option set with the options specified by one or more `Name,Value` pair arguments.

### Input Arguments

#### Name-Value Pair Arguments

Specify optional comma-separated pairs of `Name,Value` arguments. `Name` is the argument name and `Value` is the corresponding value. `Name` must appear inside quotes. You can specify several name and value pair arguments in any order as `Name1,Value1,...,NameN,ValueN`.

#### InitialState — Handling of initial states

'auto' (default) | 'model' | 'zero' | 'estimate' | 'backcast'

Handling of initial states during estimation, specified as one of the following values:

- 'model' — The initial state is parameterized by the ODE file used by the `idgrey` model. The ODE file must return 6 or more output arguments.
- 'zero' — The initial state is set to zero. Any values returned by the ODE file are ignored.

- `'estimate'` — The initial state is treated as an independent estimation parameter.
- `'backcast'` — The initial state is estimated using the best least squares fit.
- `'auto'` — The software chooses the method to handle initial states based on the estimation data.
- Vector of doubles — Specify a column vector of length  $N_x$ , where  $N_x$  is the number of states. For multiexperiment data, specify a matrix with  $N_e$  columns, where  $N_e$  is the number of experiments. The specified values are treated as fixed values during the estimation process.

### **DisturbanceModel — Handling of disturbance component**

`'auto'` (default) | `'model'` | `'fixed'` | `'none'` | `'estimate'`

Handling of the disturbance component ( $K$ ) during estimation, specified as one of the following values:

- `'model'` —  $K$  values are parameterized by the ODE file used by the `idgrey` model. The ODE file must return 5 or more output arguments.
- `'fixed'` — The value of the  $K$  property of the `idgrey` model is fixed to its original value.
- `'none'` —  $K$  is fixed to zero. Any values returned by the ODE file are ignored.
- `'estimate'` —  $K$  is treated as an independent estimation parameter.
- `'auto'` — The software chooses the method to handle how the disturbance component is handled during estimation. The software uses the `'model'` method if the ODE file returns 5 or more output arguments with a finite value for  $K$ . Else, the software uses the `'fixed'` method.

---

**Note** Noise model cannot be estimated using frequency domain data.

---

### **Focus — Error to be minimized**

`'prediction'` (default) | `'simulation'`

Error to be minimized in the loss function during estimation, specified as the comma-separated pair consisting of `'Focus'` and one of the following values:

- `'prediction'` — The one-step ahead prediction error between measured and predicted outputs is minimized during estimation. As a result, the estimation focuses on producing a good predictor model.

- 'simulation' — The simulation error between measured and simulated outputs is minimized during estimation. As a result, the estimation focuses on making a good fit for simulation of model response with the current inputs.

The `Focus` option can be interpreted as a weighting filter in the loss function. For more information, see “Loss Function and Model Quality Metrics”.

### **WeightingFilter — Weighting prefilter**

[ ] (default) | vector | matrix | cell array | linear system

Weighting prefilter applied to the loss function to be minimized during estimation. To understand the effect of `WeightingFilter` on the loss function, see “Loss Function and Model Quality Metrics”.

Specify `WeightingFilter` as one of the following values:

- [ ] — No weighting prefilter is used.
- Passbands — Specify a row vector or matrix containing frequency values that define desired passbands. You select a frequency band where the fit between estimated model and estimation data is optimized. For example, `[wl,wh]` where `wl` and `wh` represent lower and upper limits of a passband. For a matrix with several rows defining frequency passbands, `[w1l,w1h;w2l,w2h;w3l,w3h;...]`, the estimation algorithm uses the union of the frequency ranges to define the estimation passband.

Passbands are expressed in `rad/TimeUnit` for time-domain data and in `FrequencyUnit` for frequency-domain data, where `TimeUnit` and `FrequencyUnit` are the time and frequency units of the estimation data.

- SISO filter — Specify a single-input-single-output (SISO) linear filter in one of the following ways:
  - A SISO LTI model
  - `{A,B,C,D}` format, which specifies the state-space matrices of a filter with the same sample time as estimation data.
  - `{numerator,denominator}` format, which specifies the numerator and denominator of the filter as a transfer function with same sample time as estimation data.

This option calculates the weighting function as a product of the filter and the input spectrum to estimate the transfer function.

- Weighting vector — Applicable for frequency-domain data only. Specify a column vector of weights. This vector must have the same length as the frequency vector of

the data set, `Data.Frequency`. Each input and output response in the data is multiplied by the corresponding weight at that frequency.

### **EnforceStability** — Control whether to enforce stability of model

`false` (default) | `true`

Control whether to enforce stability of estimated model, specified as the comma-separated pair consisting of 'EnforceStability' and either `true` or `false`.

Data Types: `logical`

### **EstimateCovariance** — Control whether to generate parameter covariance data

`true` (default) | `false`

Controls whether parameter covariance data is generated, specified as `true` or `false`.

If `EstimateCovariance` is `true`, then use `getcov` to fetch the covariance matrix from the estimated model.

### **Display** — Specify whether to display the estimation progress

'off' (default) | 'on'

Specify whether to display the estimation progress, specified as one of the following values:

- 'on' — Information on model structure and estimation results are displayed in a progress-viewer window.
- 'off' — No progress or results information is displayed.

### **InputOffset** — Removal of offset from time-domain input data during estimation

`[]` (default) | vector of positive integers | matrix

Removal of offset from time-domain input data during estimation, specified as the comma-separated pair consisting of 'InputOffset' and one of the following:

- A column vector of positive integers of length  $Nu$ , where  $Nu$  is the number of inputs.
- `[]` — Indicates no offset.
- $Nu$ -by- $Ne$  matrix — For multi-experiment data, specify `InputOffset` as an  $Nu$ -by- $Ne$  matrix.  $Nu$  is the number of inputs, and  $Ne$  is the number of experiments.

Each entry specified by `InputOffset` is subtracted from the corresponding input data.

**OutputOffset — Removal of offset from time-domain output data during estimation**

[] (default) | vector | matrix

Removal of offset from time-domain output data during estimation, specified as the comma-separated pair consisting of 'OutputOffset' and one of the following:

- A column vector of length  $N_y$ , where  $N_y$  is the number of outputs.
- [] — Indicates no offset.
- $N_y$ -by- $N_e$  matrix — For multi-experiment data, specify OutputOffset as a  $N_y$ -by- $N_e$  matrix.  $N_y$  is the number of outputs, and  $N_e$  is the number of experiments.

Each entry specified by OutputOffset is subtracted from the corresponding output data.

**OutputWeight — Weighting of prediction errors in multi-output estimations**

[] (default) | 'noise' | positive semidefinite symmetric matrix

Weighting of prediction errors in multi-output estimations, specified as one of the following values:

- 'noise' — Minimize  $\det(E^*E/N)$ , where  $E$  represents the prediction error and  $N$  is the number of data samples. This choice is optimal in a statistical sense and leads to maximum likelihood estimates if nothing is known about the variance of the noise. It uses the inverse of the estimated noise variance as the weighting function.

---

**Note** OutputWeight must not be 'noise' if SearchMethod is 'lsqnonlin'.

---

- Positive semidefinite symmetric matrix ( $W$ ) — Minimize the trace of the weighted prediction error matrix  $\text{trace}(E^*E*W/N)$  where:
  - $E$  is the matrix of prediction errors, with one column for each output, and  $W$  is the positive semidefinite symmetric matrix of size equal to the number of outputs. Use  $W$  to specify the relative importance of outputs in multiple-output models, or the reliability of corresponding data.
  - $N$  is the number of data samples.
- [] — The software chooses between the 'noise' or using the identity matrix for  $W$ .

This option is relevant for only multi-output models.

## Regularization — Options for regularized estimation of model parameters

structure

Options for regularized estimation of model parameters. For more information on regularization, see “Regularized Estimates of Model Parameters”.

Regularization is a structure with the following fields:

- **Lambda** — Constant that determines the bias versus variance tradeoff.

Specify a positive scalar to add the regularization term to the estimation cost.

The default value of zero implies no regularization.

**Default:** 0

- **R** — Weighting matrix.

Specify a vector of nonnegative numbers or a square positive semi-definite matrix. The length must be equal to the number of free parameters of the model.

For black-box models, using the default value is recommended. For structured and grey-box models, you can also specify a vector of `np` positive numbers such that each entry denotes the confidence in the value of the associated parameter.

The default value of 1 implies a value of `eye(npfree)`, where `npfree` is the number of free parameters.

**Default:** 1

- **Nominal** — The nominal value towards which the free parameters are pulled during estimation.

The default value of zero implies that the parameter values are pulled towards zero. If you are refining a model, you can set the value to `'model'` to pull the parameters towards the parameter values of the initial model. The initial parameter values must be finite for this setting to work.

**Default:** 0

## SearchMethod — Numerical search method used for iterative parameter estimation

`'auto'` (default) | `'gn'` | `'gna'` | `'lm'` | `'grad'` | `'lsqnonlin'` | `'fmincon'`

Numerical search method used for iterative parameter estimation, specified as the comma-separated pair consisting of 'SearchMethod' and one of the following:

- 'auto' — A combination of the line search algorithms, 'gn', 'lm', 'gna', and 'grad' methods is tried in sequence at each iteration. The first descent direction leading to a reduction in estimation cost is used.
- 'gn' — Subspace Gauss-Newton least squares search. Singular values of the Jacobian matrix less than  $\text{GnPinvConstant} \cdot \text{eps} \cdot \max(\text{size}(J)) \cdot \text{norm}(J)$  are discarded when computing the search direction.  $J$  is the Jacobian matrix. The Hessian matrix is approximated as  $J^T J$ . If there is no improvement in this direction, the function tries the gradient direction.
- 'gna' — Adaptive subspace Gauss-Newton search. Eigenvalues less than  $\text{gamma} \cdot \max(\text{sv})$  of the Hessian are ignored, where  $\text{sv}$  contains the singular values of the Hessian. The Gauss-Newton direction is computed in the remaining subspace.  $\text{gamma}$  has the initial value `InitialGnaTolerance` (see Advanced in 'SearchOptions' for more information). This value is increased by the factor `LMStep` each time the search fails to find a lower value of the criterion in fewer than five bisections. This value is decreased by the factor  $2 \cdot \text{LMStep}$  each time a search is successful without any bisections.
- 'lm' — Levenberg-Marquardt least squares search, where the next parameter value is  $-\text{pinv}(H+d \cdot I) \cdot \text{grad}$  from the previous one.  $H$  is the Hessian,  $I$  is the identity matrix, and  $\text{grad}$  is the gradient.  $d$  is a number that is increased until a lower value of the criterion is found.
- 'grad' — Steepest descent least squares search.
- 'lsqnonlin' — Trust-region-reflective algorithm of `lsqnonlin`. Requires Optimization Toolbox software.
- 'fmincon' — Constrained nonlinear solvers. You can use the sequential quadratic programming (SQP) and trust-region-reflective algorithms of the `fmincon` solver. If you have Optimization Toolbox software, you can also use the interior-point and active-set algorithms of the `fmincon` solver. Specify the algorithm in the `SearchOptions.Algorithm` option. The `fmincon` algorithms may result in improved estimation results in the following scenarios:
  - Constrained minimization problems when there are bounds imposed on the model parameters.
  - Model structures where the loss function is a nonlinear or non smooth function of the parameters.

- Multi-output model estimation. A determinant loss function is minimized by default for multi-output model estimation. `fmincon` algorithms are able to minimize such loss functions directly. The other search methods such as `'lm'` and `'gn'` minimize the determinant loss function by alternately estimating the noise variance and reducing the loss value for a given noise variance value. Hence, the `fmincon` algorithms can offer better efficiency and accuracy for multi-output model estimations.

**SearchOptions — Option set for the search algorithm**

search option set

Option set for the search algorithm, specified as the comma-separated pair consisting of `'SearchOptions'` and a search option set with fields that depend on the value of `SearchMethod`.

**SearchOptions Structure When SearchMethod is Specified as 'gn', 'gna', 'lm', 'grad', or 'auto'**

Field Name	Description	Default
Tolerance	Minimum percentage difference between the current value of the loss function and its expected improvement after the next iteration, specified as a positive scalar. When the percentage of expected improvement is less than <code>Tolerance</code> , the iterations stop. The estimate of the expected loss-function improvement at the next iteration is based on the Gauss-Newton vector computed for the current parameter value.	0.01
MaxIterations	<p>Maximum number of iterations during loss-function minimization, specified as a positive integer. The iterations stop when <code>MaxIterations</code> is reached or another stopping criterion is satisfied, such as <code>Tolerance</code>.</p> <p>Setting <code>MaxIterations = 0</code> returns the result of the start-up procedure.</p> <p>Use <code>sys.Report.Termination.Iterations</code> to get the actual number of iterations during an estimation, where <code>sys</code> is an <code>idtf</code> model.</p>	20

Field Name	Description	Default
Advanced	Advanced search settings, specified as a structure with the following fields:	
Field Name	Description	Default
GnPinvConstant	Jacobian matrix singular value threshold, specified as a positive scalar. Singular values of the Jacobian matrix that are smaller than $\text{GnPinvConstant} * \max(\text{size}(J) * \text{norm}(J) * \text{eps})$ are discarded when computing the search direction. Applicable when SearchMethod is 'gn'.	10000
InitialGnaTolerance	Initial value of $\gamma$ , specified as a positive scalar. Applicable when SearchMethod is 'gna'.	0.0001
LMStartValue	Starting value of search-direction length $d$ in the Levenberg-Marquardt method, specified as a positive scalar. Applicable when SearchMethod is 'lm'.	0.001
LMStep	Size of the Levenberg-Marquardt step, specified as a positive integer. The next value of the search-direction length $d$ in the Levenberg-Marquardt method is LMStep times the previous one. Applicable when SearchMethod is 'lm'.	2
MaxBisections	Maximum number of bisections used for line search along the search direction, specified as a positive integer.	25
MaxFunctionEvaluations	Maximum number of calls to the model file, specified as a positive integer. Iterations stop if the number of calls to the model file exceeds this value.	Inf
MinParameterChange	Smallest parameter update allowed per iteration, specified as a nonnegative scalar.	0
RelativeImprovement	Relative improvement threshold, specified as a nonnegative scalar. Iterations stop if the relative	0

Field Name	Description		Default
	Field Name	Description	Default
		improvement of the criterion function is less than this value.	
	StepReduction	<p>Step reduction factor, specified as a positive scalar that is greater than 1. The suggested parameter update is reduced by the factor <code>StepReduction</code> after each try. This reduction continues until <code>MaxBisections</code> tries are completed or a lower value of the criterion function is obtained.</p> <p><code>StepReduction</code> is not applicable for SearchMethod 'lm' (Levenberg-Marquardt method).</p>	2

**SearchOptions Structure When SearchMethod is Specified as 'lsqnonlin'**

<b>Field Name</b>	<b>Description</b>	<b>Default</b>
FunctionTolerance	Termination tolerance on the loss function that the software minimizes to determine the estimated parameter values, specified as a positive scalar.  The value of FunctionTolerance is the same as that of <code>opt.SearchOptions.Advanced.TolFun</code> .	1e-5
StepTolerance	Termination tolerance on the estimated parameter values, specified as a positive scalar.  The value of StepTolerance is the same as that of <code>opt.SearchOptions.Advanced.TolX</code> .	1e-6
MaxIterations	Maximum number of iterations during loss-function minimization, specified as a positive integer. The iterations stop when MaxIterations is reached or another stopping criterion is satisfied, such as FunctionTolerance.  The value of MaxIterations is the same as that of <code>opt.SearchOptions.Advanced.MaxIter</code> .	20
Advanced	Advanced search settings, specified as an option set for <code>lsqnonlin</code> .  For more information, see the Optimization Options table in "Optimization Options" (Optimization Toolbox).	Use <code>optimset('lsqnonlin')</code> to create a default option set.

**SearchOptions Structure When SearchMethod is Specified as 'fmincon'**

Field Name	Description	Default
Algorithm	<p>fmincon optimization algorithm, specified as one of the following:</p> <ul style="list-style-type: none"> <li>• 'sqp' — Sequential quadratic programming algorithm. The algorithm satisfies bounds at all iterations, and it can recover from NaN or Inf results. It is not a large-scale algorithm. For more information, see “Large-Scale vs. Medium-Scale Algorithms” (Optimization Toolbox).</li> <li>• 'trust-region-reflective' — Subspace trust-region method based on the interior-reflective Newton method. It is a large-scale algorithm.</li> <li>• 'interior-point' — Large-scale algorithm that requires Optimization Toolbox software. The algorithm satisfies bounds at all iterations, and it can recover from NaN or Inf results.</li> <li>• 'active-set' — Requires Optimization Toolbox software. The</li> </ul>	'sqp'

Field Name	Description	Default
	<p>algorithm can take large steps, which adds speed. It is not a large-scale algorithm.</p> <p>For more information about the algorithms, see “Constrained Nonlinear Optimization Algorithms” (Optimization Toolbox) and “Choosing the Algorithm” (Optimization Toolbox).</p>	
FunctionTolerance	Termination tolerance on the loss function that the software minimizes to determine the estimated parameter values, specified as a positive scalar.	1e-6
StepTolerance	Termination tolerance on the estimated parameter values, specified as a positive scalar.	1e-6
MaxIterations	Maximum number of iterations during loss function minimization, specified as a positive integer. The iterations stop when MaxIterations is reached or another stopping criterion is satisfied, such as FunctionTolerance.	100

### Advanced – Additional advanced options structure

Additional advanced options, specified as a structure with the following fields:

- `ErrorThreshold` — Specifies when to adjust the weight of large errors from quadratic to linear.

Errors larger than `ErrorThreshold` times the estimated standard deviation have a linear weight in the loss function. The standard deviation is estimated robustly as the median of the absolute deviations from the median of the prediction errors, divided by 0.7. For more information on robust norm choices, see section 15.2 of [2].

`ErrorThreshold = 0` disables robustification and leads to a purely quadratic loss function. When estimating with frequency-domain data, the software sets `ErrorThreshold` to zero. For time-domain data that contains outliers, try setting `ErrorThreshold` to 1.6.

**Default:** 0

- `MaxSize` — Specifies the maximum number of elements in a segment when input-output data is split into segments.

`MaxSize` must be a positive integer.

**Default:** 250000

- `StabilityThreshold` — Specifies thresholds for stability tests.

`StabilityThreshold` is a structure with the following fields:

- `s` — Specifies the location of the right-most pole to test the stability of continuous-time models. A model is considered stable when its right-most pole is to the left of `s`.

**Default:** 0

- `z` — Specifies the maximum distance of all poles from the origin to test stability of discrete-time models. A model is considered stable if all poles are within the distance `z` from the origin.

**Default:**  $1 + \sqrt{\text{eps}}$

- `AutoInitThreshold` — Specifies when to automatically estimate the initial state.

The initial state is estimated when

$$\frac{\|y_{p,z} - y_{meas}\|}{\|y_{p,e} - y_{meas}\|} > \text{AutoInitThreshold}$$

- $y_{meas}$  is the measured output.
- $y_{p,z}$  is the predicted output of a model estimated using zero initial states.
- $y_{p,e}$  is the predicted output of a model estimated using estimated initial states.

Applicable when `InitialState` is 'auto'.

**Default:** 1.05

## Output Arguments

**opt** — Options set for `greyest`

`greyestOptions` option set

Option set for `greyest`, returned as an `greyestOptions` option set.

## Examples

### Create Default Options Set for Linear Grey Box Estimation

```
opt = greyestOptions;
```

### Specify Options for Linear Grey Box Estimation

Create an options set for `greyest` using the 'backcast' algorithm to initialize the state. Specify `Display` as 'on'.

```
opt = greyestOptions('InitialState','backcast','Display','on');
```

Alternatively, use dot notation to set the values of `opt`.

```
opt = greyestOptions;  
opt.InitialState = 'backcast';  
opt.Display = 'on';
```

## References

- [1] Wills, Adrian, B. Ninness, and S. Gibson. "On Gradient-Based Search for Multivariable System Estimates". *Proceedings of the 16th IFAC World Congress, Prague, Czech Republic, July 3-8, 2005*. Oxford, UK: Elsevier Ltd., 2005.
- [2] Ljung, L. *System Identification: Theory for the User*. Upper Saddle River, NJ: Prentice-Hall PTR, 1999.

## See Also

greyest | idgrey | idnlgrey | pem | ssest

## Topics

"Loss Function and Model Quality Metrics"

**Introduced in R2012a**

# hasdelay

True for linear model with time delays

## Syntax

```
B = hasdelay(sys)
B = hasdelay(sys, 'elem')
```

## Description

`B = hasdelay(sys)` returns 1 (true) if the model `sys` has input delays, output delays, I/O delays, or internal delays, and 0 (false) otherwise. If `sys` is a model array, then `B` is true if least one model in `sys` has delays.

`B = hasdelay(sys, 'elem')` returns a logical array of the same size as the model array `sys`. The logical array indicates which models in `sys` have delays.

## See Also

`absorbDelay` | `totaldelay`

**Introduced in R2012a**

## iddata

Time- or frequency-domain data

### Syntax

```
data = iddata(y,[],Ts)
data = iddata(y,u,Ts)
data = iddata(y,u,Ts,'Frequency',W)
data = iddata(y,u,Ts,'P1',V1,...,'PN',VN)
data = iddata(idfrd_object)
```

### Description

`data = iddata(y,[],Ts)` creates an `iddata` object for time-series data, containing a time-domain output signal `y` and an empty input signal `[]`, respectively. `Ts` specifies the sample time of the experimental data.

`data = iddata(y,u,Ts)` creates an `iddata` object containing a time-domain output signal `y` and input signal `u`, respectively. `Ts` specifies the sample time of the experimental data.

`data = iddata(y,u,Ts,'Frequency',W)` creates an `iddata` object containing a frequency-domain output signal `y` and input signal `u`, respectively. `Ts` specifies the sample time of the experimental data. `W` specifies the `iddata` property `'frequency'` as a vector of frequencies.

`data = iddata(y,u,Ts,'P1',V1,...,'PN',VN)` creates an `iddata` object containing a time-domain or frequency-domain output signal `y` and input signal `u`, respectively. `Ts` specifies the sample time of the experimental data. `'P1',V1,...,'PN',VN` are property-value pairs, as described in “Properties” on page 1-529.

`data = iddata(idfrd_object)` transforms an `idfrd` object to a frequency-domain `iddata` object.

## Arguments

*y*

Output signal from a system, specified as a column vector of length  $N$  for a single-output system, where  $N$  is the number of observations. For a multiple-output system with  $N_y$  output channels, specify *y* as an  $N$ -by- $N_y$  matrix. *y* sets the `OutputData` property of the `iddata` object.

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**Note** Output data must be in the same domain as input data.

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*u*

Input signal to a system, specified as a column vector of length  $N$  for a single-input system, where  $N$  is the number of observations. For a multiple-input system with  $N_u$  input channels, specify *u* as an  $N$ -by- $N_u$  matrix. *u* sets the `InputData` property of the `iddata` object.

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**Note** Input data must be in the same domain as output data.

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*Ts*

Time interval between successive data samples in seconds. Default value is 1. For continuous-time data in the frequency domain, set *Ts* to 0.

'P1', *V1*, ..., 'PN', *VN*

Pairs of `iddata` property names and property values.

*idfrd\_object*

Name of `idfrd` data object.

## Constructor

### Requirements for Constructing an `iddata` Object

To construct an `iddata` object, you must have already imported data into the MATLAB workspace, as described in “Time-Domain Data Representation”.

## Constructing an `iddata` Object for Time-Domain Data

Use the following syntax to create a time-domain `iddata` object data:

```
data = iddata(y,u,Ts)
```

You can also specify additional properties, as follows:

```
data = iddata(y,u,Ts,'Property1',Value1,...,'PropertyN',ValueN)
```

For more information about accessing object properties, see “Properties” on page 1-529.

Here, `Ts` is the sample time, or the time interval, between successive data samples:

- For uniformly sampled data, `Ts` is a scalar value equal to the sample time of your experiment.
- For nonuniformly sampled data, `Ts` is `[]`, and the value of the `SamplingInstants` property is a column vector containing individual time values. For example:

```
data = iddata(y,u,[],'SamplingInstants',TimeVector)
```

where `TimeVector` represents a vector of time values.

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**Note** You can modify the property `SamplingInstants` by setting it to a new vector with the length equal to the number of data samples.

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The default time unit is seconds, but you can specify any unit using the `TimeUnit` property. For more information about `iddata` time properties, see “Modifying Time and Frequency Vectors”.

To represent time-series data, use the following syntax:

```
ts_data = iddata(y,[],Ts)
```

where `y` is the output data, `[]` indicates empty input data, and `Ts` is the sample time.

## Constructing an `iddata` Object for Frequency-Domain Data

Frequency-domain data is the Fourier transform of the input and output signals at specific frequency values. To represent frequency-domain data, use the following syntax to create the `iddata` object:

```
data = iddata(y,u,Ts,'Frequency',w)
```

'Frequency' is an `iddata` property that specifies the frequency values  $w$ , where  $w$  is the frequency column vector that defines the frequencies at which the Fourier transform values of  $y$  and  $u$  are computed.  $T_s$  is the time interval between successive data samples in seconds for the original time-domain data.  $w$ ,  $y$ , and  $u$  have the same number of rows.

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**Note** You must specify the frequency vector for frequency-domain data.

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For more information about `iddata` time and frequency properties, see “Modifying Time and Frequency Vectors”.

To specify a continuous-time system, set  $T_s$  to  $0$ .

You can specify additional properties when you create the `iddata` object, as follows:

```
data = iddata(y,u,Ts,'Property1',Value1,...,'PropertyN',ValueN)
```

For more information about accessing object properties, see “Properties” on page 1-529.

## Properties

After creating the object, you can use `get` or dot notation to access the object property values.

Use `set` or dot notation to set a property of an existing object.

The following table describes `iddata` object properties and their values. These properties are specified as property-value arguments 'P1', V1, ..., 'PN', VN' in the `iddata` constructor, or you can set them using the `set` command or dot notation. In the list below,  $N$  denotes the number of data samples in the input and output signals,  $n_y$  is the number of output channels,  $n_u$  is the number of input channels, and  $N_e$  is the number of experiments.

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**Tip** Property names are not case sensitive. You do not need to type the entire property name. However, the portion you enter must be enough to uniquely identify the property.

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Property Name	Description	Value
Domain	Specifies whether the data is in the time domain or frequency domain.	<ul style="list-style-type: none"> <li>'Frequency' — Frequency-domain data.</li> <li>'Time' (Default) — Time-domain data.</li> </ul>
ExperimentName	Name of each data set contained in the <code>iddata</code> object.	For $N_e$ experiments, a 1-by- $N_e$ cell array of character vectors. Each cell contains the name of the corresponding experiment. Default names are <code>{'Exp1', 'Exp2', ...}</code> .
Frequency	(Frequency-domain data only)  Frequency values for defining the Fourier Transforms of the signals.	For a single experiment, this is an $N$ -by-1 vector. For $N_e$ experiments, a 1-by- $N_e$ cell array and each cell contains the frequencies of the corresponding experiment.
InputData	Name of MATLAB variable that stores the input signal to a system.	For $nu$ input channels and $N$ data samples, this is an $N$ -by- $nu$ matrix.
InputName	Specifies the names of individual input channels.	Cell array of length $nu$ -by-1 contains the name of each input channel. Default names are <code>{'u1'; 'u2'; ...}</code> .
InputUnit	Specifies the units of each input channel.	Cell array of character vectors of length $nu$ -by-1. Each cell contains the units of the corresponding input channel, for example <code>{'rad', 'rad/s'}</code> .

Property Name	Description	Value
InterSample	Specifies the behavior of the input signals between samples for transformations between discrete-time and continuous-time.	<p>For a single experiment:</p> <ul style="list-style-type: none"> <li>• <code>zoh</code>— (Default) Zero-order hold maintains a piecewise-constant input signal between samples.</li> <li>• <code>foh</code>— First-order hold maintains a piecewise-linear input signal between samples.</li> <li>• <code>bl</code>— Band-limited behavior specifies that the continuous-time input signal has zero power above the Nyquist frequency.</li> </ul> <p>For <code>Ne</code> experiments, <code>InterSample</code> is an <code>nu-by-Ne</code> cell array. Each cell contains one of these values corresponding to each experiment.</p>
Name	Name of the data set.	Character vector that contains the name of the data set. For example, <code>'dryer data'</code> . The default value is <code>''</code> .
Notes	Comments about the data set.	Character vector or cell array of character vectors that contain optional details about the data set. For example, <code>'data from experiment 2'</code> .
OutputData	Name of MATLAB variable that stores the output signal from a system.	For <code>ny</code> output channels and <code>N</code> samples, this is an <code>N-by-ny</code> matrix.

Property Name	Description	Value
OutputName	For a multiple-output system, specifies the names of individual output channels.	Cell array of length ny-by-1 contains the name of each output channel. Default names are {'y1'; 'y2'; ...}.
OutputUnit	Specifies the units of each output channel.	A cell array of character vectors of length ny-by-1. Each cell contains the units of the corresponding output channel, for example {'rad', 'rad/s'}.
Period	Period of the input signal.	(Default) For a nonperiodic signal, set to <code>inf</code> . For a multiple-input signal, this is an nu-by-1 vector and the kth entry contains the period of the kth input. For Ne experiments, this is a 1-by-Ne cell array and each cell contains a scalar or vector of periods for the corresponding experiment.
SamplingInstants	(Time-domain data only) The time values in the time vector calculated from the properties Tstart and Ts.	For a single experiment, this is an N-by-1 vector. For Ne experiments, this is a 1-by-Ne cell array and each cell contains the sampling instants of the corresponding experiment.
TimeUnit	(Time-domain data only) Time unit.	Specify TimeUnit as one of the following values: 'nanoseconds', 'microseconds', 'milliseconds', 'seconds' (default), 'minutes', 'hours', 'days', 'weeks', 'months' or 'years'.

Property Name	Description	Value
Ts	<p>Time interval between successive data samples in seconds. Must be specified for both time- and frequency-domain data. For frequency-domain, it is used to compute Fourier transforms of the signals as discrete-time Fourier transforms (DTFT) with the indicated sample time.</p> <hr/> <p><b>Note</b> Your data must be uniformly sampled.</p> <hr/>	<p>Default value is 1. For continuous-time data in the frequency domain, set to 0; the inputs and outputs are interpreted as continuous-time Fourier transforms of the signals. Note that Ts is essential also for frequency-domain data, for proper interpretation of how the Fourier transforms were computed: They are interpreted as discrete-time Fourier transforms (DTFT) with the indicated sample time. For multiple-experiment data, Ts is a 1-by-Ne cell array and each cell contains the sample time of the corresponding experiment.</p>
Tstart	<p>(Time-domain data only) Specifies the start time of the time vector.</p>	<p>For a single experiment, this is a scalar. For Ne experiments, Tstart is a 1-by-Ne cell array and each cell contains the starting time of the corresponding experiment.</p>

Property Name	Description	Value
FrequencyUnit	(Frequency-domain data only) Frequency unit.	Specifies the units of the frequency vector (see Frequency). Specify as one of the following: 'rad/TimeUnit', 'cycles/TimeUnit', 'rad/s', 'Hz', 'kHz', 'MHz', 'GHz', or 'rpm'. The units 'rad/TimeUnit' and 'cycles/TimeUnit' are relative to the time units specified in the TimeUnit property. Setting FrequencyUnit does not change the frequency vector. To convert the units and automatically scale frequency points, use chgFreqUnit.
UserData	Additional comments.	Any MATLAB data type. The default value is [].

## Examples

### Create an iddata Object for Time-Domain Data

Create an `iddata` object using single-input/single-output (SISO) data. The input and output each contain 1000 samples with the sample time of 0.08 second.

```
load dryer2
data = iddata(y2,u2,0.08)

data =

Time domain data set with 1000 samples.
Sample time: 0.08 seconds

Outputs      Unit (if specified)
  y1
```

Inputs	Unit (if specified)
u1	

The default channel name 'y1' is assigned to the first and only output channel. When y2 contains several channels, the channels are assigned default names 'y1', 'y2', 'y2', ..., 'yn'. Similarly, the default channel name 'u1' is assigned to the first and only input channel. For more information about naming channels, see “Naming, Adding, and Removing Data Channels”.

### View and Modify Properties of iddata Object

To view and modify a property of an `iddata` object, use dot notation.

Load input u2 and output y2 of the data.

```
load dryer2
```

Create an `iddata` object.

```
data = iddata(y2,u2,0.08);
```

You can use `get(data)` to view all properties of the `iddata` object. You can specify properties when you create an `iddata` object using the constructor syntax. For example, `data = iddata(y,u,Ts,'Property1',Value1,...,'PropertyN',ValueN)`.

Use dot notation to change property values for an existing `iddata` object.

```
data.ts = 0.05;
```

Property names are not case sensitive. You do not need to type the entire property name if the first few letters uniquely identify the property.

You can use `data.y` as an alternative to `data.OutputData` to access the output values, or use `data.u` as an alternative to `data.InputData` to access the input values.

## Examine `iddata` Object that Contains Frequency-Domain Data

An `iddata` object containing frequency-domain data includes frequency-specific properties, such as `Frequency` for the frequency vector and `Units` for frequency units (instead of `Tstart` and `SamplingInstants` for time-domain data).

Load input `u2` and output `y2` of the data.

```
load dryer2;
```

Create an `iddata` object.

```
data = iddata(y2,u2,0.08);
```

Transform the data to frequency domain using the Fourier transform.

```
data = fft(data);
```

Get the frequency vector of the data.

```
data.Frequency;
```

Get the frequency units of the data.

```
data.Units;
```

## See Also

[advice](#) | [detrend](#) | [fcad](#) | [getexp](#) | [idfilt](#) | [idfrd](#) | [plot](#) | [resample](#) | [size](#)

**Introduced before R2006a**

# iddataPlotOptions

Option set for iddata/plot

## Syntax

```
opt = iddataPlotOptions('time')
opt = iddataPlotOptions('frequency')
opt = iddataPlotOptions( ____, 'identpref')
```

## Description

`opt = iddataPlotOptions('time')` creates the default option set for plotting time-domain data. Use dot notation to customize the option set, if needed.

`opt = iddataPlotOptions('frequency')` creates a default option set for plotting frequency-domain data. Use dot notation to customize the option set, if needed.

`opt = iddataPlotOptions( ____, 'identpref')` initializes the plot options with the System Identification Toolbox preferences. This syntax can include any of the input argument combinations in the previous syntaxes. Use this syntax to change a few plot options but otherwise use your toolbox preferences.

## Examples

### Create Option Set for Plotting Time-Domain Data

Create an options set with default options for time-domain data.

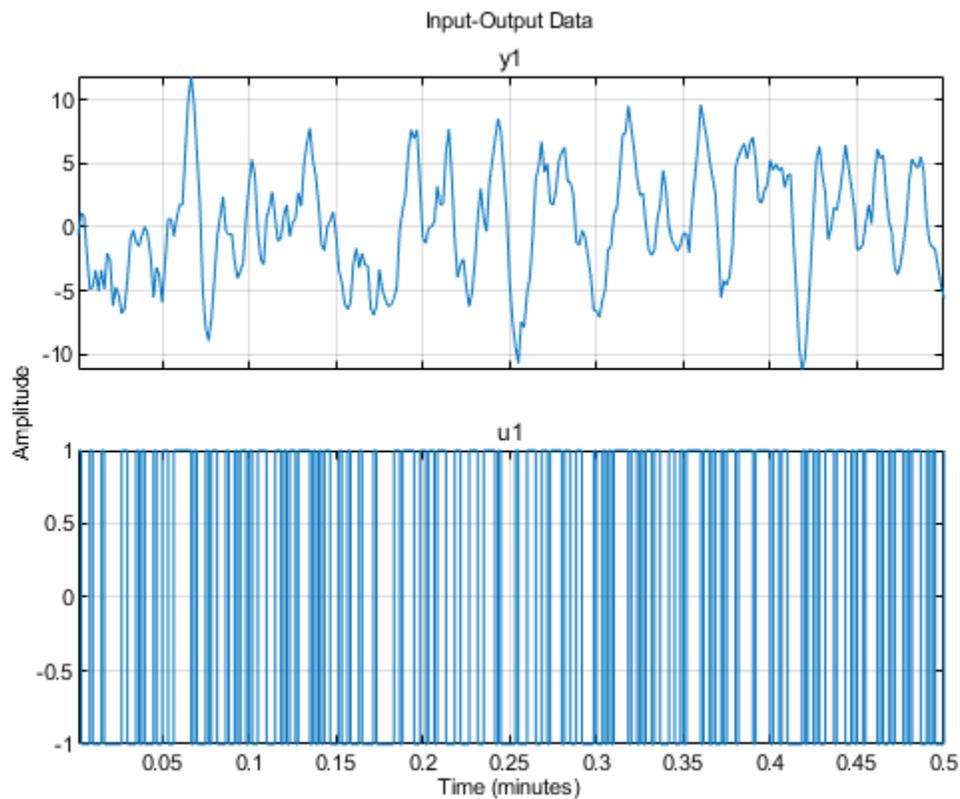
```
opt = iddataPlotOptions('time');
```

Specify plot properties, such as time units and grid. View the plot in minutes

```
opt.TimeUnits = 'minutes';
% Turn grid on
opt.Grid = 'on';
```

Create a plot using the specified options.

```
load iddata1 z1  
h = plot(z1, opt);
```



### Change Orientation of Input-Output Data Axes

Generate data with two inputs and one output.

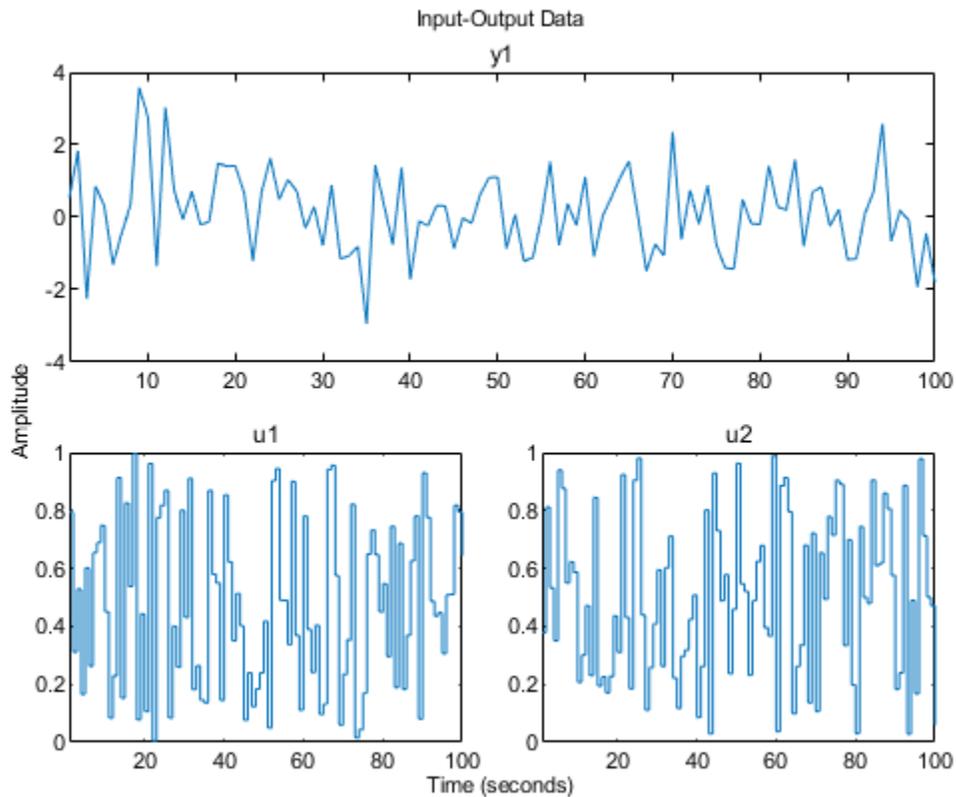
```
z = iddata(randn(100,1),rand(100,2));
```

Configure a time plot.

```
opt = iddataPlotOptions('time');
```

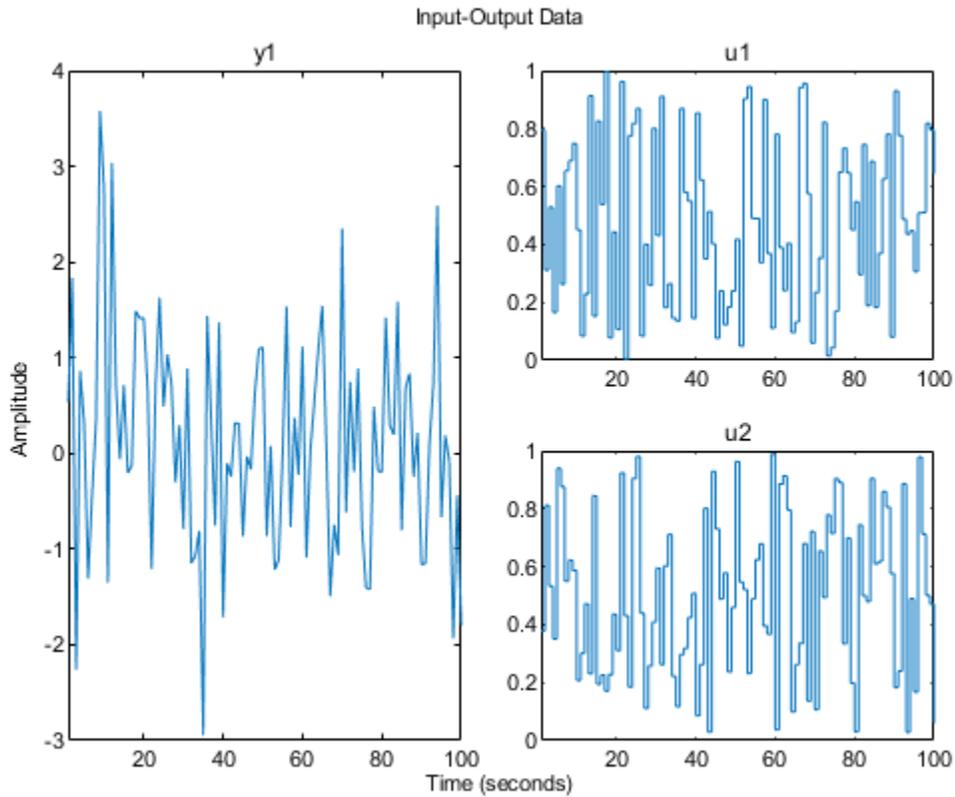
Plot the data.

```
h = plot(z,opt);
```



Change the orientation of the plots such that all inputs are plotted in one column, and all outputs are in a second column.

```
opt.Orientation = 'two-column';  
h = plot(z,opt);
```



Alternatively, use `setoptions`.

```
setoptions(h, 'Orientation', 'two-column')
```

You can also change the orientation by right-clicking the plot and choosing `Orientation` in the context menu.

### Create Option Set for Plotting Frequency-Domain Data

Create an option set with default options for frequency-domain data.

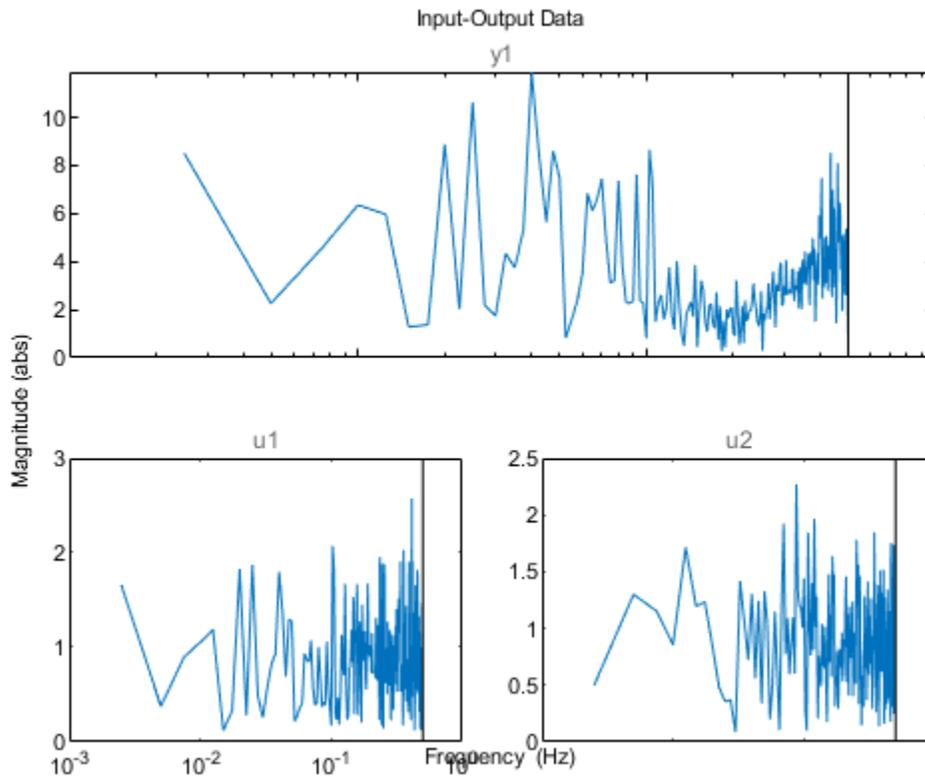
```
opt = iddataPlotOptions('frequency');
```

Specify plot properties, such as phase visibility and frequency units.

```
opt.PhaseVisible = 'off';  
opt.FreqUnits = 'Hz';
```

Create a plot with the specified options.

```
load iddata7 z7  
zf = fft(z7);  
h = plot(zf,opt);
```



## Initialize a Plot Using Toolbox Preferences

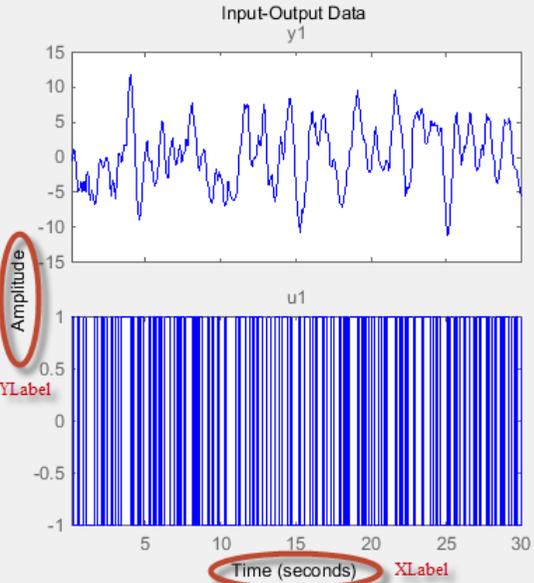
```
opt = iddataPlotOptions('time','identpref');
```

## Output Arguments

### **opt** — Option set for `iddata/plot`

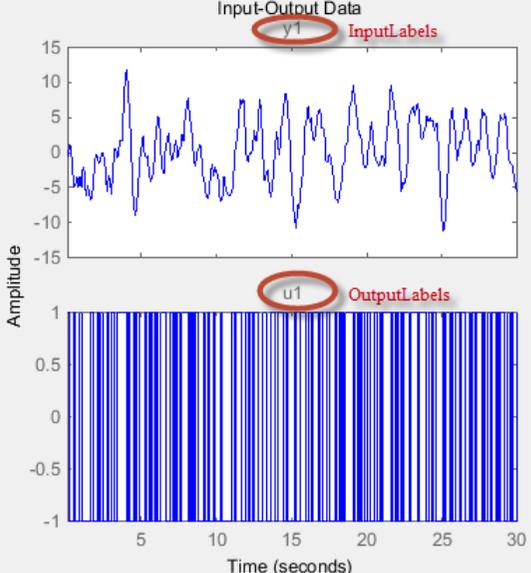
`iddataPlotOptions` option set

Option set containing the specified options for `iddata/plot`. The structure has the following fields:

Field	Description
Title, XLabel, YLabel	<p data-bbox="782 302 1307 395">Text and style for axes labels and plot title, specified as a structure array with the following fields:</p> <div data-bbox="782 423 1337 1020" style="border: 1px solid gray; padding: 5px;">  </div> <ul data-bbox="782 1052 1307 1529" style="list-style-type: none"> <li>• <b>String</b> — Title and axes label text, specified as a character vector. <b>Default Title:</b> 'Input-Output Data' <b>Default XLabel:</b> 'Time' <b>Default YLabel:</b> 'Amplitude'</li> <li>• <b>FontSize</b> — Font size, specified as scalar value greater than 0. <b>Default:</b> 8</li> <li>• <b>FontWeight</b> — Thickness of text, specified as one of the following values: 'Normal'   'Bold' <b>Default:</b> 'Normal'</li> </ul>

Field	Description
	<ul style="list-style-type: none"> <li>• <b>Font Angle</b> — Text character angle, specified as one of the following values: 'Normal'   'Italic' <b>Default:</b> 'Normal'</li> <li>• <b>Color</b> — Color of text, specified as vector of RGB values between 0 to 1. <b>Default:</b> [0,0,0]</li> <li>• <b>Interpreter</b> — Interpretation of text characters, specified as one of the following values: 'tex'   'latex'   'none' <b>Default:</b> 'tex'</li> </ul>
TickLabel	<p>Tick label style, specified as a structure array with the following fields:</p> <ul style="list-style-type: none"> <li>• <b>FontSize</b> — Font size, specified as scalar value greater than 0. <b>Default:</b> 8</li> <li>• <b>FontWeight</b> — Thickness of text, specified as one of the following values: 'Normal'   'Bold' <b>Default:</b> 'Normal'</li> <li>• <b>Font Angle</b> — Text character angle, specified as one of the following values: 'Normal'   'Italic' <b>Default:</b> 'Normal'</li> <li>• <b>Color</b> — Color of text, specified as vector of RGB values between 0 to 1   character vector of color name   'none'. For example, for yellow color, specify as one of the following: [1 1 0], 'yellow', or 'y'. <b>Default:</b> [0,0,0]</li> </ul>

Field	Description
Grid	<p>Show or hide the grid, specified as one of the following values: 'off'   'on'</p> <p><b>Default:</b> 'off'</p>
GridColor	<p>Color of the grid lines, specified as one of the following values: vector of RGB values in the range [0,1]   character vector of color name   'none'. For example, for yellow color, specify as one of the following: [1 1 0], 'yellow', or 'y'.</p> <p><b>Default:</b> [0.15,0.15,0.15]</p>
XlimMode, YlimMode	<p>Axes limit modes, specified as one of the following values:</p> <ul style="list-style-type: none"> <li>• 'auto' — The axis limits are based on the data plotted</li> <li>• 'manual' — The values explicitly set with Xlim, Ylim</li> </ul> <p><b>Default:</b> 'auto'</p>
Xlim, Ylim	<p>Axes limits, specified as maximum and minimum values.</p> <p><b>Default:</b> [0 1]</p>
IOGrouping	<p>Grouping of input-output pairs, specified as one of the following values: 'none'   'inputs'   'outputs'   'all'</p> <p><b>Default:</b> 'none'</p>

Field	Description
InputLabels, OutputLabels	<p data-bbox="783 302 1332 395">Input and output label styles on individual plot axes, specified as a structure array with the following fields:</p> <div data-bbox="783 427 1332 1017" style="border: 1px solid #ccc; padding: 5px;">  </div> <ul data-bbox="786 1052 1332 1538" style="list-style-type: none"> <li>• <b>FontSize</b> — Font size, specified as data type scalar. <b>Default:</b> 8</li> <li>• <b>FontWeight</b> — Thickness of text, specified as one of the following values: 'Normal'   'Bold' <b>Default:</b> 'Normal'</li> <li>• <b>Font Angle</b> — Text character angle, specified as one of the following values: 'Normal'   'Italic' <b>Default:</b> 'Normal'</li> <li>• <b>Color</b> — Color of text, specified as a vector of RGB values between 0 to 1   character vector of color name   'none'.</li> </ul>

Field	Description
	<p>For example, for yellow color, specify as one of the following: [1 1 0], 'yellow', or 'y'.  <b>Default:</b> [0.4,0.4,0.4]</p> <ul style="list-style-type: none"> <li>• <b>Interpreter</b> — Interpretation of text characters, specified as one of the following values: 'tex'   'latex'   'none'  <b>Default:</b> 'tex'</li> </ul>
InputVisible, OutputVisible	<p>Visibility of input and output channels, specified as one of the following values: 'off'   'on'  <b>Default:</b> 'on'</p>
Orientation	<p>Orientation of the input and output data plots, specified as one of the following values:</p> <ul style="list-style-type: none"> <li>• 'two-row' — Plot all outputs in one row and all inputs in a second row</li> <li>• 'two-column' — Plot all outputs in one column and all inputs in a second column</li> <li>• 'single-row' — Plot all inputs and outputs in one row</li> <li>• 'single-column' — Plot all inputs and outputs in one column</li> </ul> <p><b>Default:</b> 'two-row'.</p>

<b>Field</b>	<b>Description</b>
For time-domain data plots only:	
<b>Field</b>	<b>Description</b>
TimeUnits	<p>Time units, specified as one of the following values:</p> <ul style="list-style-type: none"><li>• 'nanoseconds'</li><li>• 'microseconds'</li><li>• 'milliseconds'</li><li>• 'seconds'</li><li>• 'minutes'</li><li>• 'hours'</li><li>• 'days'</li><li>• 'weeks'</li><li>• 'months'</li><li>• 'years'</li></ul> <p>You can also specify 'auto' which uses time units specified in the <code>TimeUnit</code> property of the data. For multiple systems with different time units, the units of the first system is used.</p>
Normalize	<p>Normalize responses, specified as one of the following values: 'on'   'off'</p> <p><b>Default:</b> 'off'</p>

Field	Description
For frequency-domain data plots only:	
Field	Description
FreqUnits	<p>Frequency units, specified as one of the following values:</p> <ul style="list-style-type: none"> <li>• 'Hz'</li> <li>• 'rad/second'</li> <li>• 'rpm'</li> <li>• 'kHz'</li> <li>• 'MHz'</li> <li>• 'GHz'</li> <li>• 'rad/nanosecond'</li> <li>• 'rad/microsecond'</li> <li>• 'rad/millisecond'</li> <li>• 'rad/minute'</li> <li>• 'rad/hour'</li> <li>• 'rad/day'</li> <li>• 'rad/week'</li> <li>• 'rad/month'</li> <li>• 'rad/year'</li> <li>• 'cycles/nanosecond'</li> <li>• 'cycles/microsecond'</li> <li>• 'cycles/millisecond'</li> <li>• 'cycles/hour'</li> <li>• 'cycles/day'</li> <li>• 'cycles/week'</li> <li>• 'cycles/month'</li> <li>• 'cycles/year'</li> </ul> <p><b>Default:</b> 'rad/s'</p>

Field	Description
<b>Field</b>	<b>Description</b>
	You can also specify 'auto' which uses frequency units rad/TimeUnit relative to system time units specified in the TimeUnit property. For multiple systems with different time units, the units of the first system are used.
FreqScale	Frequency scale, specified as one of the following values: 'linear'   'log'  <b>Default:</b> 'log'
MagUnits	Magnitude units, specified as one of the following values: 'dB'   'abs'  <b>Default:</b> 'dB'
MagScale	Magnitude scale, specified as one of the following values: 'linear'   'log'  <b>Default:</b> 'linear'
MagVisible	Magnitude plot visibility, specified as one of the following values: 'on'   'off'  <b>Default:</b> 'on'
MagLowerLimMode	Enables a lower magnitude limit, specified as one of the following values: 'auto'   'manual'  <b>Default:</b> 'auto'
MagLowerLim	Lower magnitude limit, , specified as data type double. It is typically decided by the range of the amplitudes the plotted data takes.
PhaseUnits	Phase units, specified as one of the following values: 'deg'   'rad'  <b>Default:</b> 'deg'

Field	Description
<b>Field</b>	<b>Description</b>
PhaseVisible	Phase plot visibility, specified as one of the following values: 'on'   'off'  <b>Default:</b> 'on'
PhaseWrapping	Enable phase wrapping, specified as one of the following values: 'on'   'off'  <b>Default:</b> 'off'
PhaseWrappingBranch	Phase value at which the plot wraps accumulated phase when PhaseWrapping is set to 'on'.  <b>Default:</b> -180 (phase wraps into the interval [-180°,180°))
PhaseMatching	Enable phase matching, specified as one of the following values: 'on'   'off'  <b>Default:</b> 'off'
PhaseMatchingFreq	Frequency for matching phase, specified as data type double.
PhaseMatchingValue	The value to which phase responses are matched closely, specified as a real number representing the desired phase value PhaseMatchingFreq.

## See Also

iddata/plot | identpref

**Introduced in R2014a**

# identpref

Set System Identification Toolbox preferences

## Syntax

identpref

## Description

identpref opens a Graphical User Interface (GUI) which allows you to change the System Identification Toolbox preferences. Preferences set in this GUI affect future plots only (existing plots are not altered).

Your preferences are stored to disk (in a system-dependent location) and will be automatically reloaded in future MATLAB sessions using the System Identification Toolbox software.

## See Also

### Topics

“Toolbox Preferences Editor”

**Introduced in R2012a**

## idfilt

Filter data using user-defined passbands, general filters, or Butterworth filters

### Syntax

```
Zf = idfilt(Z,filter)
Zf = idfilt(Z,filter,causality)
Zf = idfilt(Z,filter,'FilterOrder',NF)
```

### Description

`Zf = idfilt(Z,filter)` filters data using user-defined passbands, general filters, or Butterworth filters. `Z` is the data, defined as an `iddata` object. `Zf` contains the filtered data as an `iddata` object. The filter can be defined in three ways:

- As an explicit system that defines the filter.

```
filter = idm or filter = {num,den} or filter = {A,B,C,D}
```

`idm` can be any SISO identified linear model or LTI model object. Alternatively the filter can be defined as a cell array `{A,B,C,D}` of SISO state-space matrices or as a cell array `{num,den}` of numerator/denominator filter coefficients.

- As a vector or matrix that defines one or several passbands.

```
filter=[[wp1l,wp1h];[ wp2l,wp2h]; ... ;[wpnl,wpnh]]
```

The matrix is  $n$ -by-2, where each row defines a passband. A filter is constructed that gives the union of these passbands. For time-domain data, it is computed as cascaded Butterworth filters or order `NF`. The default value of `NF` is 5.

- For time-domain data — The passbands are in units of `rad/TimeUnit`, where `TimeUnit` is the time units of the estimation data.
- For frequency-domain data — The passbands are in the frequency units (`FrequencyUnit` property) of the estimation data.

For example, to define a stopband between `ws1` and `ws2`, use

```
filter = [0 ws1; ws2,Nyqf]
```

where `Nyqf` is the Nyquist frequency.

- For frequency-domain data, only the frequency response of the filter can be specified.

```
filter = Wf
```

Here `Wf` is a vector of possibly complex values that define the filter's frequency response, so that the inputs and outputs at frequency `Z.Frequency(kf)` are multiplied by `Wf(kf)`. `Wf` is a column vector of length = number of frequencies in `Z`. If the data object has several experiments, `Wf` is a cell array of length = # of experiments in `Z`.

`Zf = idfilt(Z,filter,causality)` specifies causality. For time-domain data, the filtering is carried out in the time domain as causal filtering as default. This corresponds to a last argument `causality = 'causal'`. With `causality = 'noncausal'`, a noncausal, zero-phase filter is used for the filtering (corresponding to `filtfilt` in the Signal Processing Toolbox product).

For frequency-domain data, the signals are multiplied by the frequency response of the filter. With the filters defined as passband, this gives ideal, zero-phase filtering (“brickwall filters”). Frequencies that have been assigned zero weight by the filter (outside the passband, or via the frequency response) are removed from the `iddata` object `Zf`.

`Zf = idfilt(Z,filter,'FilterOrder',NF)` specifies the filter order. The time domain filters in the pass-band case are calculated as cascaded Butterworth pass-band and stop-band filters. The orders of these filters are 5 by default, which can be changed to an arbitrary integer `NF`.

It is common practice in identification to select a frequency band where the fit between model and data is concentrated. Often this corresponds to bandpass filtering with a passband over the interesting breakpoints in a Bode diagram. For identification where a disturbance model is also estimated, it is better to achieve the desired estimation result by using the `'WeightingFilter'` option of the estimation command than just to prefilter the data. The values for `'WeightingFilter'` are the same as the argument `filter` in `idfilt`.

## Algorithms

The Butterworth filter is the same as `butter` in the Signal Processing Toolbox product. Also, the zero-phase filter is equivalent to `filtfilt` in that toolbox.

## References

Ljung (1999), Chapter 14.

## See Also

`iddata` | `resample`

**Introduced before R2006a**

## idfrd

Frequency-response data or model

### Syntax

```
h = idfrd(Response,Freq,Ts)
h =
idfrd(Response,Freq,Ts,'CovarianceData',Covariance,'SpectrumData',Spec,
'NoiseCovariance',Speccov)
h = idfrd(Response,Freq,Ts,...
    'P1',V1,'PN',VN)
h = idfrd(mod)
h = idfrd(mod,Freqs)
```

### Description

`h = idfrd(Response,Freq,Ts)` constructs an `idfrd` object that stores the frequency response, `Response`, of a linear system at frequency values, `Freq`. `Ts` is the sample time. For a continuous-time system, set `Ts=0`.

`h = idfrd(Response,Freq,Ts,'CovarianceData',Covariance,'SpectrumData',Spec,'NoiseCovariance',Speccov)` also stores the uncertainty of the response, `Covariance`, the spectrum of the additive disturbance (noise), `Spec`, and the covariance of the noise, `Speccov`.

`h = idfrd(Response,Freq,Ts,... 'P1',V1,'PN',VN)` constructs an `idfrd` object that stores a frequency-response model with properties specified by the `idfrd` model property-value pairs.

`h = idfrd(mod)` converts a System Identification Toolbox or Control System Toolbox linear model to frequency-response data at default frequencies, including the output noise spectra and their covariance.

`h = idfrd(mod,Freqs)` converts a System Identification Toolbox or Control System Toolbox linear model to frequency-response data at frequencies `Freqs`.

For a model

$$y(t) = G(q)u(t) + H(q)e(t)$$

`idfrd` object stores the transfer function estimate  $G(e^{i\omega})$ , as well as the spectrum of the additive noise ( $\Phi_v$ ) at the output.

$$\Phi_v(\omega) = \lambda T \left| H(e^{i\omega T}) \right|^2$$

where  $\lambda$  is the estimated variance of  $e(t)$ , and  $T$  is the sample time.

For a continuous-time system, the noise spectrum is given by:

$$\Phi_v(\omega) = \lambda \left| H(e^{i\omega}) \right|^2$$

## Creating `idfrd` from Given Responses

`Response` is a 3-D array of dimension `ny-by-nu-by-Nf`, with `ny` being the number of outputs, `nu` the number of inputs, and `Nf` the number of frequencies (that is, the length of `Freqs`). `Response(ky, ku, kf)` is thus the complex-valued frequency response from input `ku` to output `ky` at frequency  $\omega = \text{Freqs}(kf)$ . When defining the response of a SISO system, `Response` can be given as a vector.

`Freqs` is a column vector of length `Nf` containing the frequencies of the response.

`Ts` is the sample time. `Ts = 0` means a continuous-time model.

Intersample behavior: For discrete-time frequency response data (`Ts > 0`), you can also specify the intersample behavior of the input signal that was in effect when the samples were collected originally from an experiment. To specify the intersample behavior, use:

```
mf = idfrd(Response, Freq, Ts, 'InterSample', 'zoh');
```

For multi-input systems, specify the intersample behavior using an `Nu-by-1` cell array, where `Nu` is the number of inputs. The `InterSample` property is irrelevant for continuous-time data.

Covariance is a 5-D array containing the covariance of the frequency response. It has dimension `ny-by-nu-by-Nf-by-2-by-2`. The structure is such that `Covariance(ky, ku, kf, :, :)` is the 2-by-2 covariance matrix of the response `Response(ky, ku, kf)`. The 1-1 element is the variance of the real part, the 2-2 element is the variance of the imaginary part, and the 1-2 and 2-1 elements are the covariance between the real and imaginary parts. `squeeze(Covariance(ky, ku, kf, :, :))` thus gives the covariance matrix of the corresponding response.

The format for spectrum information is as follows:

`spec` is a 3-D array of dimension `ny-by-ny-by-Nf`, such that `spec(ky1, ky2, kf)` is the cross spectrum between the noise at output `ky1` and the noise at output `ky2`, at frequency `Freqs(kf)`. When `ky1 = ky2` the (power) spectrum of the noise at output `ky1` is thus obtained. For a single-output model, `spec` can be given as a vector.

`speccov` is a 3-D array of dimension `ny-by-ny-by-Nf`, such that `speccov(ky1, ky1, kf)` is the variance of the corresponding power spectrum.

If only `SpectrumData` is to be packaged in the `idfrd` object, set `Response = []`.

## Converting to `idfrd`

An `idfrd` object can also be computed from a given linear identified model, `mod`.

If the frequencies `Freqs` are not specified, a default choice is made based on the dynamics of the model `mod`.

Estimated covariance:

- If you obtain `mod` by identification, the software computes the estimated covariance for the `idfrd` object from the uncertainty information in `mod`. The software uses the Gauss approximation formula for this calculation for all model types, except grey-box models. For grey-box models (`idgrey`), the software applies numerical differentiation. The step sizes for the numerical derivatives are determined by `nuderst`.
- If you create `mod` by using commands such as `idss`, `idtf`, `idproc`, `idgrey`, or `idpoly`, then the software sets `CovarianceData` to `[]`.

Delay treatment: If `mod` contains delays, then the software assigns the delays of the `idfrd` object, `h`, as follows:

- `h.InputDelay = mod.InputDelay`

- `h.IODelay = mod.IODelay+repmat(mod.OutputDelay,[1,nu])`

The expression `repmat(mod.OutputDelay,[1,nu])` returns a matrix containing the output delay for each input/output pair.

Frequency responses for submodels can be obtained by the standard subreferencing, `h = idfrd(m(2,3))`. `h = idfrd(m(:,[]))` gives an `h` that just contains `SpectrumData`.

The `idfrd` models can be graphed with `bode`, `spectrum`, and `nyquist`, which accept mixtures of parametric models, such as `idtf` and `idfrd` models as arguments. Note that `spa`, `spafdr`, and `etfe` return their estimation results as `idfrd` objects.

## Constructor

The `idfrd` represents complex frequency-response data. Before you can create an `idfrd` object, you must import your data as described in “Frequency-Response Data Representation”.

---

**Note** The `idfrd` object can only encapsulate one frequency-response data set. It does not support the `iddata` equivalent of multiexperiment data.

---

Use the following syntax to create the data object `fr_data`:

```
fr_data = idfrd(response,f,Ts)
```

Suppose that `ny` is the number of output channels, `nu` is the number of input channels, and `nf` is a vector of frequency values. `response` is an `ny-by-nu-by-nf` 3-D array. `f` is the frequency vector that contains the frequencies of the response. `Ts` is the sample time, which is used when measuring or computing the frequency response. If you are working with a continuous-time system, set `Ts` to `0`.

`response(ky,ku,kf)`, where `ky`, `ku`, and `kf` reference the `k`th output, input, and frequency value, respectively, is interpreted as the complex-valued frequency response from input `ku` to output `ky` at frequency `f(kf)`.

You can specify object properties when you create the `idfrd` object using the constructor syntax:

```
fr_data = idfrd(response,f,Ts,
    'Property1',Value1,...,'PropertyN',ValueN)
```

## Properties

idfrd object properties include:

### ResponseData

Frequency response data.

The 'ResponseData' property stores the frequency response data as a 3-D array of complex numbers. For SISO systems, 'ResponseData' is a vector of frequency response values at the frequency points specified in the 'Frequency' property. For MIMO systems with  $N_u$  inputs and  $N_y$  outputs, 'ResponseData' is an array of size  $[N_y \ N_u \ N_w]$ , where  $N_w$  is the number of frequency points.

### Frequency

Frequency points of the frequency response data. Specify Frequency values in the units specified by the FrequencyUnit property.

### FrequencyUnit

Frequency units of the model.

Units of the frequency vector in the Frequency property, specified as one of the following values:

- 'rad/TimeUnit'
- 'cycles/TimeUnit'
- 'rad/s'
- 'Hz'
- 'kHz'
- 'MHz'
- 'GHz'
- 'rpm'

The units 'rad/TimeUnit' and 'cycles/TimeUnit' are relative to the time units specified in the TimeUnit property.

Changing this property changes the overall system behavior. Use chgFreqUnit to convert between frequency units without modifying system behavior.

**Default:** 'rad/TimeUnit'

### **SpectrumData**

Power spectra and cross spectra of the system output disturbances (noise). Specify `SpectrumData` as a 3-D array of complex numbers.

Specify `SpectrumData` as a 3-D array with dimension `ny-by-ny-by-Nf`. Here, `ny` is the number of outputs and `Nf` is the number of frequency points.

`SpectrumData(ky1, ky2, kf)` is the cross spectrum between the noise at output `ky1` and the noise at output `ky2`, at frequency `Freqs(kf)`. When `ky1 = ky2` the (power) spectrum of the noise at output `ky1` is thus obtained.

For a single-output model, specify `SpectrumData` as a vector.

### **CovarianceData**

Response data covariance matrices.

Specify `CovarianceData` as a 5-D array with dimension `ny-by-nu-by-Nf-by-2-by-2`. Here, `ny`, `nu`, and `Nf` are the number of outputs, inputs and frequency points, respectively. `CovarianceData(ky, ku, kf, :, :)` is the 2-by-2 covariance matrix of the response data `ResponseData(ky, ku, kf)`. The 1-1 element is the variance of the real part, the 2-2 element is the variance of the imaginary part, and the 1-2 and 2-1 elements are the covariance between the real and imaginary parts.

```
squeeze(Covariance(ky, ku, kf, :, :))
```

### **NoiseCovariance**

Power spectra variance.

Specify `NoiseCovariance` as a 3-D array with dimension `ny-by-ny-by-Nf`. Here, `ny` is the number of outputs and `Nf` is the number of frequency points.

`NoiseCovariance(ky1, ky1, kf)` is the variance of the corresponding power spectrum. To eliminate the influence of the noise component from the model, specify `NoiseVariance` as `0`. Zero variance makes the predicted output the same as the simulated output.

## Report

Summary report that contains information about the estimation options and results when the frequency-response model is obtained using estimation commands, such as `spa`, `spafdr`, and `etfe`. Use `Report` to query a model for how it was estimated, including its:

- Estimation method
- Estimation options
- Search termination conditions
- Estimation data fit and other quality metrics

The contents of `Report` are irrelevant if the model was created by construction.

```
f = logspace(-1,1,100);  
[mag,phase] = bode(idtf([1 .2],[1 2 1 1]),f);  
response = mag.*exp(1j*phase*pi/180);  
m = idfrd(response,f,0.08);  
m.Report.Method
```

```
ans =
```

```
''
```

If you obtain the frequency-response model using estimation commands, the fields of `Report` contain information on the estimation data, options, and results.

```
load iddata3;  
m = spa(z3);  
m.Report.Method
```

```
ans =
```

```
SPA
```

`Report` is a read-only property.

For more information on this property and how to use it, see the Output Arguments section of the corresponding estimation command reference page and “Estimation Report”.

## InterSample

Input intersample behavior.

Specifies the behavior of the input signals between samples for transformations between discrete-time and continuous-time. This property is meaningful for discrete-time `idfrd` models only.

Set `InterSample` to one of the following:

- `'zoh'` — The input signal used for construction/estimation of the frequency response data was subject to a zero-order-hold filter.
- `'foh'` — The input signal was subject to a first-order-hold filter.
- `'bl'` — The input signal has no power above the Nyquist frequency ( $\pi/\text{sys.Ts}$  rad/s). This is typically the case when the input signal is measured experimentally using an anti-aliasing filter and a sampler. Ideally, treat the data as continuous-time. That is, if the signals used for the estimation of the frequency response were subject to anti-aliasing filters, set `sys.Ts` to zero.

For multi-input data, specify `InterSample` as an  $Nu$ -by-1 cell array, where  $Nu$  is the number of inputs.

### **IODElay**

Transport delays. `IODElay` is a numeric array specifying a separate transport delay for each input/output pair.

For continuous-time systems, specify transport delays in the time unit stored in the `TimeUnit` property. For discrete-time systems, specify transport delays in integer multiples of the sample time, `Ts`.

For a MIMO system with  $Ny$  outputs and  $Nu$  inputs, set `IODElay` to a  $Ny$ -by- $Nu$  array. Each entry of this array is a numerical value that represents the transport delay for the corresponding input/output pair. You can also set `IODElay` to a scalar value to apply the same delay to all input/output pairs.

**Default:** 0 for all input/output pairs

### **InputDelay**

Input delay for each input channel, specified as a scalar value or numeric vector. For continuous-time systems, specify input delays in the time unit stored in the `TimeUnit` property. For discrete-time systems, specify input delays in integer multiples of the sample time `Ts`. For example, `InputDelay = 3` means a delay of three sample times.

For a system with `Nu` inputs, set `InputDelay` to an `Nu-by-1` vector. Each entry of this vector is a numerical value that represents the input delay for the corresponding input channel.

You can also set `InputDelay` to a scalar value to apply the same delay to all channels.

**Default:** 0

### **OutputDelay**

Output delays.

For identified systems, like `idfrd`, `OutputDelay` is fixed to zero.

### **Ts**

Sample time. For continuous-time models,  $T_s = 0$ . For discrete-time models, `Ts` is a positive scalar representing the sample time expressed in the unit specified by the `TimeUnit` property of the model. To denote a discrete-time model with unspecified sample time, set  $T_s = -1$ .

Changing this property does not discretize or resample the model.

**Default:** 1

### **TimeUnit**

Units for the time variable, the sample time `Ts`, and any time delays in the model, specified as one of the following values:

- 'nanoseconds'
- 'microseconds'
- 'milliseconds'
- 'seconds'
- 'minutes'
- 'hours'
- 'days'
- 'weeks'
- 'months'

- 'years'

Changing this property has no effect on other properties, and therefore changes the overall system behavior. Use `chgTimeUnit` to convert between time units without modifying system behavior.

**Default:** 'seconds'

### **InputName**

Input channel names, specified as one of the following:

- Character vector — For single-input models, for example, 'controls'.
- Cell array of character vectors — For multi-input models.

Alternatively, use automatic vector expansion to assign input names for multi-input models. For example, if `sys` is a two-input model, enter:

```
sys.InputName = 'controls';
```

The input names automatically expand to {'controls(1)'; 'controls(2)'}

When you estimate a model using an `iddata` object, `data`, the software automatically sets `InputName` to `data.InputName`.

You can use the shorthand notation `u` to refer to the `InputName` property. For example, `sys.u` is equivalent to `sys.InputName`.

Input channel names have several uses, including:

- Identifying channels on model display and plots
- Extracting subsystems of MIMO systems
- Specifying connection points when interconnecting models

**Default:** '' for all input channels

### **InputUnit**

Input channel units, specified as one of the following:

- Character vector — For single-input models, for example, 'seconds'.

- Cell array of character vectors — For multi-input models.

Use `InputUnit` to keep track of input signal units. `InputUnit` has no effect on system behavior.

**Default:** '' for all input channels

### **InputGroup**

Input channel groups. The `InputGroup` property lets you assign the input channels of MIMO systems into groups and refer to each group by name. Specify input groups as a structure. In this structure, field names are the group names, and field values are the input channels belonging to each group. For example:

```
sys.InputGroup.controls = [1 2];  
sys.InputGroup.noise = [3 5];
```

creates input groups named `controls` and `noise` that include input channels 1, 2 and 3, 5, respectively. You can then extract the subsystem from the `controls` inputs to all outputs using:

```
sys(:, 'controls')
```

**Default:** Struct with no fields

### **OutputName**

Output channel names, specified as one of the following:

- Character vector — For single-output models. For example, `'measurements'`.
- Cell array of character vectors — For multi-output models.

Alternatively, use automatic vector expansion to assign output names for multi-output models. For example, if `sys` is a two-output model, enter:

```
sys.OutputName = 'measurements';
```

The output names automatically expand to `{'measurements(1)'; 'measurements(2)'}`.

When you estimate a model using an `iddata` object, `data`, the software automatically sets `OutputName` to `data.OutputName`.

You can use the shorthand notation `y` to refer to the `OutputName` property. For example, `sys.y` is equivalent to `sys.OutputName`.

Output channel names have several uses, including:

- Identifying channels on model display and plots
- Extracting subsystems of MIMO systems
- Specifying connection points when interconnecting models

**Default:** '' for all output channels

### OutputUnit

Output channel units, specified as one of the following:

- Character vector — For single-output models. For example, 'seconds'.
- Cell array of character vectors — For multi-output models.

Use `OutputUnit` to keep track of output signal units. `OutputUnit` has no effect on system behavior.

**Default:** '' for all output channels

### OutputGroup

Output channel groups. The `OutputGroup` property lets you assign the output channels of MIMO systems into groups and refer to each group by name. Specify output groups as a structure. In this structure, field names are the group names, and field values are the output channels belonging to each group. For example:

```
sys.OutputGroup.temperature = [1];  
sys.InputGroup.measurement = [3 5];
```

creates output groups named `temperature` and `measurement` that include output channels 1, and 3, 5, respectively. You can then extract the subsystem from all inputs to the measurement outputs using:

```
sys('measurement', :)
```

**Default:** Struct with no fields

## Name

System name, specified as a character vector. For example, 'system\_1'.

**Default:** ''

## Notes

Any text that you want to associate with the system, stored as a string or a cell array of character vectors. The property stores whichever data type you provide. For instance, if `sys1` and `sys2` are dynamic system models, you can set their `Notes` properties as follows:

```
sys1.Notes = "sys1 has a string.";
sys2.Notes = 'sys2 has a character vector.';
sys1.Notes
sys2.Notes
```

```
ans =
```

```
    "sys1 has a string."
```

```
ans =
```

```
    'sys2 has a character vector.'
```

**Default:** [0×1 string]

## UserData

Any type of data you want to associate with system, specified as any MATLAB data type.

**Default:** []

## SamplingGrid

Sampling grid for model arrays, specified as a data structure.

For arrays of identified linear (IDLTI) models that are derived by sampling one or more independent variables, this property tracks the variable values associated with each

model. This information appears when you display or plot the model array. Use this information to trace results back to the independent variables.

Set the field names of the data structure to the names of the sampling variables. Set the field values to the sampled variable values associated with each model in the array. All sampling variables should be numeric and scalar valued, and all arrays of sampled values should match the dimensions of the model array.

For example, if you collect data at various operating points of a system, you can identify a model for each operating point separately and then stack the results together into a single system array. You can tag the individual models in the array with information regarding the operating point:

```
nominal_engine_rpm = [1000 5000 10000];
sys.SamplingGrid = struct('rpm', nominal_engine_rpm)
```

where `sys` is an array containing three identified models obtained at rpms 1000, 5000 and 10000, respectively.

For model arrays generated by linearizing a Simulink® model at multiple parameter values or operating points, the software populates `SamplingGrid` automatically with the variable values that correspond to each entry in the array. For example, the Simulink Control Design™ commands `linearize` and `sLinearizer` populate `SamplingGrid` in this way.

**Default:** []

## Subreferencing

The different channels of the `idfrd` are retrieved by subreferencing.

`h(outputs,inputs)`

`h(2,3)` thus contains the response data from input channel 3 to output channel 2, and, if applicable, the output spectrum data for output channel 2. The channels can also be referred to by their names, as in `h('power',{ 'voltage', 'speed' })`.

## Horizontal Concatenation

Adding input channels,

```
h = [h1,h2,...,hN]
```

creates an `idfrd` model `h`, with `ResponseData` containing all the input channels in `h1, ..., hN`. The output channels of `hk` must be the same, as well as the frequency vectors. `SpectrumData` is ignored.

## Vertical Concatenation

Adding output channels,

```
h = [h1;h2;... ;hN]
```

creates an `idfrd` model `h` with `ResponseData` containing all the output channels in `h1, h2, ..., hN`. The input channels of `hk` must all be the same, as well as the frequency vectors. `SpectrumData` is also appended for the new outputs. The cross spectrum between output channels of `h1, h2, ..., hN` is then set to zero.

## Converting to iddata

You can convert an `idfrd` object to a frequency-domain `iddata` object by

```
Data = iddata(Idfrdmodel)
```

See `iddata`.

## Examples

### View and Modify Properties of idfrd Object

To view and modify a property of an `idfrd` object, use dot notation.

The following example shows how to create an `idfrd` object that contains 100 frequency-response values with a sample time of 0.08s and get its properties.

Create an `idfrd` object.

```
f = logspace(-1,1,100);  
[mag, phase] = bode(idtf([1 .2],[1 2 1 1]),f);
```

```
response = mag.*exp(1j*phase*pi/180);  
fr_data = idfrd(response, f, 0.08);
```

response and f are variables in the MATLAB Workspace browser, representing the frequency-response data and frequency values, respectively.

You can use `get(fr_data)` to view all properties of the `idfrd` object. You can specify properties when you create an `idfrd` object using the constructor syntax. For example, `fr_data = idfrd(y,u,Ts, 'Property1', Value1, ..., 'PropertyN', ValueN)`.

Use dot notation to change property values for an existing `idfrd` object. For example, change the name of the `idfrd` object.

```
fr_data.Name = 'DC_Converter';
```

If you import `fr_data` into the System Identification app, this data is named `DC_Converter` in the app, and not the variable name `fr_data`.

## See Also

[bode](#) | [etfe](#) | [freqresp](#) | [nyquist](#) | [spa](#) | [spafdr](#) | [tfest](#)

**Introduced before R2006a**

## idgrey

Linear ODE (grey-box model) with identifiable parameters

### Syntax

```
sys = idgrey(odefun,parameters,fcn_type)
sys = idgrey(odefun,parameters,fcn_type,optional_args)
sys = idgrey(odefun,parameters,fcn_type,optional_args,Ts)
sys = idgrey(odefun,parameters,fcn_type,optional_args,Ts,Name,Value)
```

### Description

`sys = idgrey(odefun,parameters,fcn_type)` creates a linear grey-box model with identifiable parameters, `sys`. `odefun` specifies the user-defined function that relates the model parameters, `parameters`, to its state-space representation.

`sys = idgrey(odefun,parameters,fcn_type,optional_args)` creates a linear grey-box model with identifiable parameters using the optional arguments required by `odefun`.

`sys = idgrey(odefun,parameters,fcn_type,optional_args,Ts)` creates a linear grey-box model with identifiable parameters with the specified sample time, `Ts`.

`sys = idgrey(odefun,parameters,fcn_type,optional_args,Ts,Name,Value)` creates a linear grey-box model with identifiable parameters with additional options specified by one or more `Name, Value` pair arguments.

### Object Description

An `idgrey` model represents a system as a continuous-time or discrete-time state-space model with identifiable (estimable) coefficients.

A state-space model of a system with input vector,  $u$ , output vector,  $y$ , and disturbance,  $e$ , takes the following form in continuous time:

$$\begin{aligned} \dot{x}(t) &= Ax(t) + Bu(t) + Ke(t) \\ y(t) &= Cx(t) + Du(t) + e(t) \end{aligned}$$

In discrete time, the state-space model takes the form:

$$\begin{aligned} x[k+1] &= Ax[k] + Bu[k] + Ke[k] \\ y[k] &= Cx[k] + Du[k] + e[k] \end{aligned}$$

For `idgrey` models, the state-space matrices  $A$ ,  $B$ ,  $C$ , and  $D$  are expressed as a function of user-defined parameters using a MATLAB function. You access estimated parameters using `sys.Structures.Parameters`, where `sys` is an `idgrey` model.

Use an `idgrey` model when you know the system of equations governing the system dynamics explicitly. You should be able to express these dynamics in the form of ordinary differential or difference equations. You specify complex relationships and constraints among the parameters that cannot be done through structured state-space models (`idss`).

You can create an `idgrey` model using the `idgrey` command. To do so, write a MATLAB function that returns the  $A$ ,  $B$ ,  $C$ , and  $D$  matrices for given values of the estimable parameters and sample time. The MATLAB function can also return the  $K$  matrix and accept optional input arguments. The matrices returned may represent a continuous-time or discrete-time model, as indicated by the sample time.

Use the estimating functions `pem` or `greyest` to obtain estimated values for the unknown parameters of an `idgrey` model.

You can convert an `idgrey` model into other dynamic systems, such as `idpoly`, `idss`, `tf`, `ss` etc. You cannot convert a dynamic system into an `idgrey` model.

## Examples

### Create Grey-Box Model with Estimable Parameters

Create an `idgrey` model to represent a DC motor. Specify the motor time-constant as an estimable parameter and that the ODE function can return continuous- or discrete-time state-space matrices.

Create the `idgrey` model.

```
odefun = 'motorDynamics';  
parameters = 1;  
fcn_type = 'cd';  
optional_args = 0.25;  
Ts = 0;  
sys = idgrey(odefun,parameters,fcn_type,optional_args,Ts);
```

`sys` is an `idgrey` model that is configured to use the shipped file `motorDynamics.m` to return the  $A$ ,  $B$ ,  $C$ ,  $D$ , and  $K$  matrices. `motorDynamics.m` also returns the initial conditions,  $X0$ . The motor constant,  $\tau$ , is defined in `motorDynamics.m` as an estimable parameter, and `parameters = 1` specifies its initial value as 1.

You can use `pem` or `greyest` to refine the estimate for  $\tau$ .

## Configure Estimable Parameter of Grey-Box Model

Specify the known parameters of a grey-box model as fixed for estimation. Also specify a minimum bound for an estimable parameter.

Create an ODE file that relates the pendulum model coefficients to its state-space representation. Save this function as `LinearPendulum.m` such that it is in the MATLAB® search path.

```
function [A,B,C,D] = LinearPendulum(m,g,l,b,Ts)  
A = [0 1; -g/l, -b/m/l^2];  
B = zeros(2,0);  
C = [1 0];  
D = zeros(1,0);  
end
```

In this function:

- $m$  is the pendulum mass.
- $g$  is the gravitational acceleration.

- $l$  is the pendulum length.
- $b$  is the viscous friction coefficient.
- $T_s$  is the model sample time.

Create a linear grey-box model associated with the ODE function.

```
odefun = 'LinearPendulum';

m = 1;
g = 9.81;
l = 1;
b = 0.2;
parameters = {'mass',m;'gravity',g;'length',l;'friction',b};

fcn_type = 'c';

sys = idgrey(odefun,parameters,fcn_type);
```

`sys` has four parameters.

Specify the known parameters,  $m$ ,  $g$ , and  $l$ , as fixed for estimation.

```
sys.Structure.Parameters(1).Free = false;
sys.Structure.Parameters(2).Free = false;
sys.Structure.Parameters(3).Free = false;
```

$m$ ,  $g$ , and  $l$  are the first three parameters of `sys`.

Specify a zero lower bound for  $b$ , the fourth parameter of `sys`.

```
sys.Structure.Parameters(4).Minimum = 0;
```

Similarly, to specify an upper bound for an estimable parameter, use the `Maximum` field of the parameter.

### Specify Additional Attributes of Grey-Box Model

Create a grey-box model with identifiable parameters. Name the input and output channels of the model, and specify seconds for the model time units.

Use `Name`, `Value` pair arguments to specify additional model properties on model creation.

```
odefun = 'motorDynamics';
parameters = 1;
fcn_type = 'cd';
optional_args = 0.25;
Ts = 0;
sys = idgrey(odefun,parameters,fcn_type,optional_args,Ts,'InputName','Voltage',...
            'OutputName',{'Angular Position','Angular Velocity'});
```

To change or specify more attributes of an existing model, you can use dot notation. For example:

```
sys.TimeUnit = 'seconds';
```

### Create Array of Grey-Box Models

Use the `stack` command to create an array of linear grey-box models.

```
odefun1 = @motorDynamics;
parameters1 = [1 2];
fcn_type = 'cd';
optional_args1 = 1;
sys1 = idgrey(odefun1,parameters1,fcn_type,optional_args1);

odefun2 = 'motorDynamics';
parameters2 = {[1 2]};
optional_args2 = 0.5;
sys2 = idgrey(odefun2,parameters2,fcn_type,optional_args2);

sysarr = stack(1,sys1,sys2);
```

`stack` creates a 2-by-1 array of `idgrey` models, `sysarr`.

## Input Arguments

### **odefun**

MATLAB function that relates the model parameters to its state-space representation.

`odefun` specifies the name of a MATLAB function (.m, .p, a function handle or .mex\* file). This function establishes the relationship between the model parameters, `parameters`,

and its state-space representation. The function may optionally relate the model parameters to the disturbance matrix and initial states.

If the function is not on the MATLAB path, then specify the full file name, including the path.

The syntax for `odefun` must be as follows:

```
[A,B,C,D] = odefun(par1,par2,...,parN,Ts,optional_arg1,optional_arg2,...)
```

The function outputs describe the model in the following linear state-space innovations form:

$$\begin{aligned}x_n(t) &= Ax(t) + Bu(t) + Ke(t); x(0) = x_0 \\ y(t) &= Cx(t) + Du(t) + e(t)\end{aligned}$$

In discrete time  $x_n(t)=x(t+Ts)$  and in continuous time,  $x_n(t) = \dot{x}(t)$ .

`par1, par2, . . . , parN` are model parameters. Each entry may be a scalar, vector or matrix.

`Ts` is the sample time.

`optional_arg1, optional_arg2, . . .` are the optional inputs that `odefun` may require. The values of the optional input arguments are unchanged through the estimation process. However, the values of `par1, par2, . . . , parN` are updated during estimation to fit the data. Use optional input arguments to vary the constants and coefficients used by your model without editing `odefun`.

The disturbance matrix,  $K$ , and the initial state values,  $x_0$ , are not parametrized. Instead, these values are determined separately, using the `DisturbanceModel` and `InitialState` estimation options, respectively. For more information regarding the estimation options, see `greyestOptions`.

A good choice for achieving the best simulation results is to set the `DisturbanceModel` option to 'none', which fixes  $K$  to zero.

(Optional) Parameterizing Disturbance: `odefun` can also return the disturbance component,  $K$ , using the syntax:

```
[A,B,C,D,K] = odefun(par1,par2,...,parN,Ts,optional_arg1,optional_arg2,...)
```

If `odefun` returns a value for  $K$  that contains NaN values, then the estimating function assumes that  $K$  is not parameterized. In this case, the value of the `DisturbanceModel` estimation option determines how  $K$  is handled.

(Optional) Parameterizing Initial State Values: To make the model initial states,  $X0$ , dependent on the model parameters, use the following syntax for `odefun`:

```
[A,B,C,D,K,X0] = odefun(par1,par2,...,parN,Ts,optional_arg1,optional_arg2,...)
```

If `odefun` returns a value for  $X0$  that contains NaN values, then the estimating function assumes that  $X0$  is not parameterized. In this case,  $X0$  may be fixed to zero or estimated separately, using the `InitialStates` estimation option.

## parameters

Initial values of the parameters required by `odefun`.

Specify `parameters` as a cell array containing the parameter initial values. If your model requires only one parameter, which may itself be a vector or a matrix, you may specify `parameters` as a matrix.

You may also specify parameter names using an  $N$ -by-2 cell array, where  $N$  is the number of parameters. The first column specifies the names, and the second column specifies the values of the parameters.

For example:

```
parameters = {'mass',par1;'stiffness',par2;'damping',par3}
```

## fcn\_type

Indicates whether the model is parameterized in continuous-time, discrete-time, or both.

`fcn_type` requires one of the following values:

- 'c' — `odefun` returns matrices corresponding to a continuous-time system, regardless of the value of  $Ts$ .
- 'd' — `odefun` returns matrices corresponding to a discrete-time system, whose values may or may not depend on the value of  $Ts$ .
- 'cd' — `odefun` returns matrices corresponding to a continuous-time system, if  $Ts=0$ .

Otherwise, if  $T_s > 0$ , `odefun` returns matrices corresponding to a discrete-time system. Select this option to sample your model using the values returned by `odefun`, rather than using the software's internal sample time conversion routines.

### **optional\_args**

Optional input arguments required by `odefun`.

Specify `optional_args` as a cell array.

If `odefun` does not require optional input arguments, specify `optional_args` as `{}`.

### **Ts**

Model sample time.

If  $T_s$  is unspecified, it is assumed to be:

- $-1$  — If `fcn_type` is 'd' or 'cd'.

$T_s = -1$  indicates a discrete-time model with unknown sample time.

- $0$  — If `fcn_type` is 'c'.

$T_s = 0$  indicates a continuous-time model.

### **Name, Value**

Specify optional comma-separated pairs of `Name, Value` arguments, where `Name` is the argument name and `Value` is the corresponding value. `Name` must appear inside single quotes (' '). You can specify several name and value pair arguments in any order as `Name1, Value1, ..., NameN, ValueN`.

Use `Name, Value` arguments to specify additional properties of `idgrey` models during model creation. For example, `idgrey(odefun, parameters, fcn_type, 'InputName', 'Voltage')` creates an `idgrey` model with the `InputName` property set to `Voltage`.

## **Properties**

`idgrey` object properties include:

**A, B, C, D**

Values of state-space matrices.

- **A** — State matrix  $A$ , an  $N_x$ -by- $N_x$  matrix, as returned by the ODE function associated with the `idgrey` model.  $N_x$  is the number of states.
- **B** — Input-to-state matrix  $B$ , an  $N_x$ -by- $N_u$  matrix, as returned by the ODE function associated with the `idgrey` model.  $N_u$  is the number of inputs and  $N_x$  is the number of states.
- **C** — State-to-output matrix  $C$ , an  $N_y$ -by- $N_x$  matrix, as returned by the ODE function associated with the `idgrey` model.  $N_x$  is the number of states and  $N_y$  is the number of outputs.
- **D** — Feedthrough matrix  $D$ , an  $N_y$ -by- $N_u$  matrix, as returned by the ODE function associated with the `idgrey` model.  $N_y$  is the number of outputs and  $N_u$  is the number of inputs.

The values  $A, B, C, D$  are returned by the ODE function associated with the `idgrey` model. Thus, you can only read these matrices; you cannot set their values.

**K**

Value of state disturbance matrix,  $K$

$K$  is  $N_x$ -by- $N_y$  matrix, where  $N_x$  is the number of states and  $N_y$  is the number of outputs.

- If `odefun` parameterizes the  $K$  matrix, then  $K$  has the value returned by `odefun`. `odefun` parameterizes the  $K$  matrix if it returns at least five outputs and the value of the fifth output does not contain NaN values.
- If `odefun` does not parameterize the  $K$  matrix, then  $K$  is a zero matrix of size  $N_x$ -by- $N_y$ .  $N_x$  is the number of states and  $N_y$  is the number of outputs. The value is treated as a fixed value of the  $K$  matrix during estimation. To make the value estimable, use the `DisturbanceModel` estimation option.
- Regardless of whether the  $K$  matrix is parameterized by `odefun` or not, you can set the value of the  $K$  property explicitly as an  $N_x$ -by- $N_y$  matrix.  $N_x$  is the number of states and  $N_y$  is the number of outputs. The specified value is treated as a fixed value of the  $K$  matrix during estimation. To make the value estimable, use the `DisturbanceModel` estimation option.

To create an estimation option set for `idgrey` models, use `greyestOptions`.

## StateName

State names, specified as one of the following:

- Character vector — For first-order models, for example, 'velocity'.
- Cell array of character vectors — For models with two or more states
- '' — For unnamed states.

**Default:** '' for all states

## StateUnit

State units, specified as one of the following:

- Character vector — For first-order models, for example, 'velocity'
- Cell array of character vectors — For models with two or more states
- '' — For states without specified units

Use `StateUnit` to keep track of the units each state is expressed in. `StateUnit` has no effect on system behavior.

**Default:** '' for all states

## Structure

Information about the estimable parameters of the `idgrey` model.

`Structure` stores information regarding the MATLAB function that parameterizes the `idgrey` model.

- `Structure.Function` — Name or function handle of the MATLAB function used to create the `idgrey` model.
- `Structure.FunctionType` — Indicates whether the model is parameterized in continuous-time, discrete-time, or both.
- `Structure.Parameters` — Information about the estimated parameters. `Structure.Parameters` contains the following fields:
  - `Value` — Parameter values. For example, `sys.Structure.Parameters(2).Value` contains the initial or estimated values of the second parameter.

NaN represents unknown parameter values.

- **Minimum** — Minimum value that the parameter can assume during estimation. For example, `sys.Structure.Parameters(1).Minimum = 0` constrains the first parameter to be greater than or equal to zero.
- **Maximum** — Maximum value that the parameter can assume during estimation.
- **Free** — Boolean value specifying whether the parameter is estimable. If you want to fix the value of a parameter during estimation, set `Free = false` for the corresponding entry.
- **Scale** — Scale of the parameter's value. `Scale` is not used in estimation.
- **Info** — Structure array for storing parameter units and labels. The structure has `Label` and `Unit` fields.

Specify parameter units and labels as character vectors. For example, `'Time'`.

- `Structure.ExtraArguments` — Optional input arguments required by the ODE function.
- `Structure.StateName` — Names of the model states.
- `Structure.StateUnit` — Units of the model states.

### NoiseVariance

The variance (covariance matrix) of the model innovations,  $e$ .

An identified model includes a white, Gaussian noise component,  $e(t)$ . `NoiseVariance` is the variance of this noise component. Typically, the model estimation function (such as `greyest` or `pem`) determines this variance.

For SISO models, `NoiseVariance` is a scalar. For MIMO models, `NoiseVariance` is a  $N_y$ -by- $N_y$  matrix, where  $N_y$  is the number of outputs in the system.

### Report

Summary report that contains information about the estimation options and results when the grey-box model is obtained using the `greyest` estimation command. Use `Report` to query a model for how it was estimated, including its:

- Estimation method
- Estimation options
- Search termination conditions

- Estimation data fit and other quality metrics

The contents of `Report` are irrelevant if the model was created by construction.

```
odefun = 'motorDynamics';
m = idgrey(odefun,1,'cd',0.25,0);
m.Report.OptionsUsed
```

```
ans =
```

```
[]
```

If you obtain the grey-box model using estimation commands, the fields of `Report` contain information on the estimation data, options, and results.

```
load(fullfile(matlabroot,'toolbox','ident','iddemos','data','dcmotordata'));
data = iddata(y,u,0.1,'Name','DC-motor');
odefun = 'motorDynamics';
init_sys = idgrey('motorDynamics',1,'cd',0.25,0);
m = greyest(data,init_sys);
m.Report.OptionsUsed
```

```
InitialState: 'auto'
  DisturbanceModel: 'auto'
    Focus: 'prediction'
  EstimateCovariance: 1
    Display: 'off'
  InputOffset: []
  OutputOffset: []
  Regularization: [1x1 struct]
  OutputWeight: []
  SearchMethod: 'auto'
  SearchOptions: [1x1 idoptions.search.identsolver]
  Advanced: [1x1 struct]
```

`Report` is a read-only property.

For more information on this property and how to use it, see the Output Arguments section of the corresponding estimation command reference page and “Estimation Report”.

### InputDelay

Input delay for each input channel, specified as a scalar value or numeric vector. For continuous-time systems, specify input delays in the time unit stored in the `TimeUnit`

property. For discrete-time systems, specify input delays in integer multiples of the sample time  $T_s$ . For example, `InputDelay = 3` means a delay of three sample times.

For a system with  $N_u$  inputs, set `InputDelay` to an  $N_u$ -by-1 vector. Each entry of this vector is a numerical value that represents the input delay for the corresponding input channel.

You can also set `InputDelay` to a scalar value to apply the same delay to all channels.

**Default:** 0

### **OutputDelay**

Output delays.

For identified systems, like `idgrey`, `OutputDelay` is fixed to zero.

### **Ts**

Sample time.

For continuous-time models,  $T_s = 0$ . For discrete-time models,  $T_s$  is a positive scalar representing the sample time expressed in the unit specified by the `TimeUnit` property of the model. To denote a discrete-time model with unspecified sample time, set  $T_s = -1$ .

Changing this property does not discretize or resample the model.

For `idgrey` models, there is no unique default value for  $T_s$ .  $T_s$  depends on the value of `fcn_type`.

### **TimeUnit**

Units for the time variable, the sample time  $T_s$ , and any time delays in the model, specified as one of the following values:

- 'nanoseconds'
- 'microseconds'
- 'milliseconds'
- 'seconds'
- 'minutes'

- 'hours'
- 'days'
- 'weeks'
- 'months'
- 'years'

Changing this property has no effect on other properties, and therefore changes the overall system behavior. Use `chgTimeUnit` to convert between time units without modifying system behavior.

**Default:** 'seconds'

### **InputName**

Input channel names, specified as one of the following:

- Character vector — For single-input models, for example, 'controls'.
- Cell array of character vectors — For multi-input models.

Alternatively, use automatic vector expansion to assign input names for multi-input models. For example, if `sys` is a two-input model, enter:

```
sys.InputName = 'controls';
```

The input names automatically expand to {'controls(1)'; 'controls(2)'}

When you estimate a model using an `iddata` object, `data`, the software automatically sets `InputName` to `data.InputName`.

You can use the shorthand notation `u` to refer to the `InputName` property. For example, `sys.u` is equivalent to `sys.InputName`.

Input channel names have several uses, including:

- Identifying channels on model display and plots
- Extracting subsystems of MIMO systems
- Specifying connection points when interconnecting models

**Default:** '' for all input channels

## InputUnit

Input channel units, specified as one of the following:

- Character vector — For single-input models, for example, 'seconds'.
- Cell array of character vectors — For multi-input models.

Use `InputUnit` to keep track of input signal units. `InputUnit` has no effect on system behavior.

**Default:** '' for all input channels

## InputGroup

Input channel groups. The `InputGroup` property lets you assign the input channels of MIMO systems into groups and refer to each group by name. Specify input groups as a structure. In this structure, field names are the group names, and field values are the input channels belonging to each group. For example:

```
sys.InputGroup.controls = [1 2];  
sys.InputGroup.noise = [3 5];
```

creates input groups named `controls` and `noise` that include input channels 1, 2 and 3, 5, respectively. You can then extract the subsystem from the `controls` inputs to all outputs using:

```
sys(:, 'controls')
```

**Default:** Struct with no fields

## OutputName

Output channel names, specified as one of the following:

- Character vector — For single-output models. For example, 'measurements'.
- Cell array of character vectors — For multi-output models.

Alternatively, use automatic vector expansion to assign output names for multi-output models. For example, if `sys` is a two-output model, enter:

```
sys.OutputName = 'measurements';
```

The output names automatically expand to `{'measurements(1)'; 'measurements(2)'}`.

When you estimate a model using an `iddata` object, `data`, the software automatically sets `OutputName` to `data.OutputName`.

You can use the shorthand notation `y` to refer to the `OutputName` property. For example, `sys.y` is equivalent to `sys.OutputName`.

Output channel names have several uses, including:

- Identifying channels on model display and plots
- Extracting subsystems of MIMO systems
- Specifying connection points when interconnecting models

**Default:** `''` for all output channels

### OutputUnit

Output channel units, specified as one of the following:

- Character vector — For single-output models. For example, `'seconds'`.
- Cell array of character vectors — For multi-output models.

Use `OutputUnit` to keep track of output signal units. `OutputUnit` has no effect on system behavior.

**Default:** `''` for all output channels

### OutputGroup

Output channel groups. The `OutputGroup` property lets you assign the output channels of MIMO systems into groups and refer to each group by name. Specify output groups as a structure. In this structure, field names are the group names, and field values are the output channels belonging to each group. For example:

```
sys.OutputGroup.temperature = [1];
sys.InputGroup.measurement = [3 5];
```

creates output groups named `temperature` and `measurement` that include output channels 1, and 3, 5, respectively. You can then extract the subsystem from all inputs to the measurement outputs using:

```
sys('measurement', :)
```

**Default:** Struct with no fields

## Name

System name, specified as a character vector. For example, 'system\_1'.

**Default:** ''

## Notes

Any text that you want to associate with the system, stored as a string or a cell array of character vectors. The property stores whichever data type you provide. For instance, if `sys1` and `sys2` are dynamic system models, you can set their `Notes` properties as follows:

```
sys1.Notes = "sys1 has a string.";
sys2.Notes = 'sys2 has a character vector.';
sys1.Notes
sys2.Notes
```

```
ans =
```

```
    "sys1 has a string."
```

```
ans =
```

```
    'sys2 has a character vector.'
```

**Default:** [0×1 string]

## UserData

Any type of data you want to associate with system, specified as any MATLAB data type.

**Default:** []

## SamplingGrid

Sampling grid for model arrays, specified as a data structure.

For arrays of identified linear (IDLTI) models that are derived by sampling one or more independent variables, this property tracks the variable values associated with each

model. This information appears when you display or plot the model array. Use this information to trace results back to the independent variables.

Set the field names of the data structure to the names of the sampling variables. Set the field values to the sampled variable values associated with each model in the array. All sampling variables should be numeric and scalar valued, and all arrays of sampled values should match the dimensions of the model array.

For example, if you collect data at various operating points of a system, you can identify a model for each operating point separately and then stack the results together into a single system array. You can tag the individual models in the array with information regarding the operating point:

```
nominal_engine_rpm = [1000 5000 10000];  
sys.SamplingGrid = struct('rpm', nominal_engine_rpm)
```

where `sys` is an array containing three identified models obtained at rpms 1000, 5000 and 10000, respectively.

For model arrays generated by linearizing a Simulink model at multiple parameter values or operating points, the software populates `SamplingGrid` automatically with the variable values that correspond to each entry in the array. For example, the Simulink Control Design commands `linearize` and `sLinearizer` populate `SamplingGrid` in this way.

**Default:** []

## See Also

[getpvec](#) | [greyest](#) | [greyestOptions](#) | [idnlgrey](#) | [idss](#) | [pem](#) | [setpvec](#) | [ssest](#)

## Topics

“Estimate Coefficients of ODEs to Fit Given Solution”

“Estimate Model Using Zero/Pole/Gain Parameters”

“Estimate Discrete-Time Grey-Box Model with Parameterized Disturbance”

“Specifying the Linear Grey-Box Model Structure”

**Introduced before R2006a**

## idinput

Generate input signals

The `idinput` command generates an input signal with specified characteristics for your system. You can use the generated input, and simulate the response of your system to study system behavior. For example, you can study the system response to periodic inputs. The system can be an actual physical system or a model such as a Simulink model. You can also design optimal experiments. For example, you can determine which input signals isolate faults or nonlinearities in your system. You can also use `idinput` to design an input that has sufficient bandwidth to excite the dynamic range of your system.

### Syntax

```
u = idinput(N)
u = idinput([N,Nu])
u = idinput([Period,Nu,NumPeriod])

u = idinput(___,Type)
u = idinput(___,Type,Band)
u = idinput(___,Type,Band,Range)
[u,freq] = idinput(___, 'sine',Band,Range,SineData)
```

### Description

`u = idinput(N)` returns a single-channel random binary input signal `u` of length `N`. The generated signal values are either -1 or 1.

`u = idinput([N,Nu])` returns an `Nu`-channel random binary input signal, where each channel signal has length `N`. The signals in each channel differ from each other.

`u = idinput([Period,Nu,NumPeriod])` returns an `Nu`-channel periodic random binary input signal with specified period and number of periods. Each input channel signal is of length `NumPeriod*Period`.

`u = idinput(___,Type)` specifies the type of input to be generated as one of the following:

- 'rbs' — Random binary signal
- 'rgs' — Random Gaussian signal
- 'prbs' — Pseudorandom binary signal
- 'sine' — Sum-of-sinusoids signal

Use with any of the previous input argument combinations.

`u = idinput( ____, Type, Band)` specifies the frequency band of the signal. For pseudorandom binary signals (PRBS), `Band` specifies the inverse of the clock period of the signal.

`u = idinput( ____, Type, Band, Range)` specifies the amplitude-range of the signal.

`[u, freq] = idinput( ____, 'sine', Band, Range, SineData)` specifies the `Type` as a sum-of-sinusoids signal and specifies the characteristics of the sine waves used to generate the signal in `SineData`. You can specify characteristics such as the number of sine waves and their frequency separation. The frequencies of the sine waves are returned in `freq`.

## Examples

### Generate a Random Binary Input Signal

Generate a single-channel random binary input signal with 200 samples.

```
N = 200;  
u = idinput(N);
```

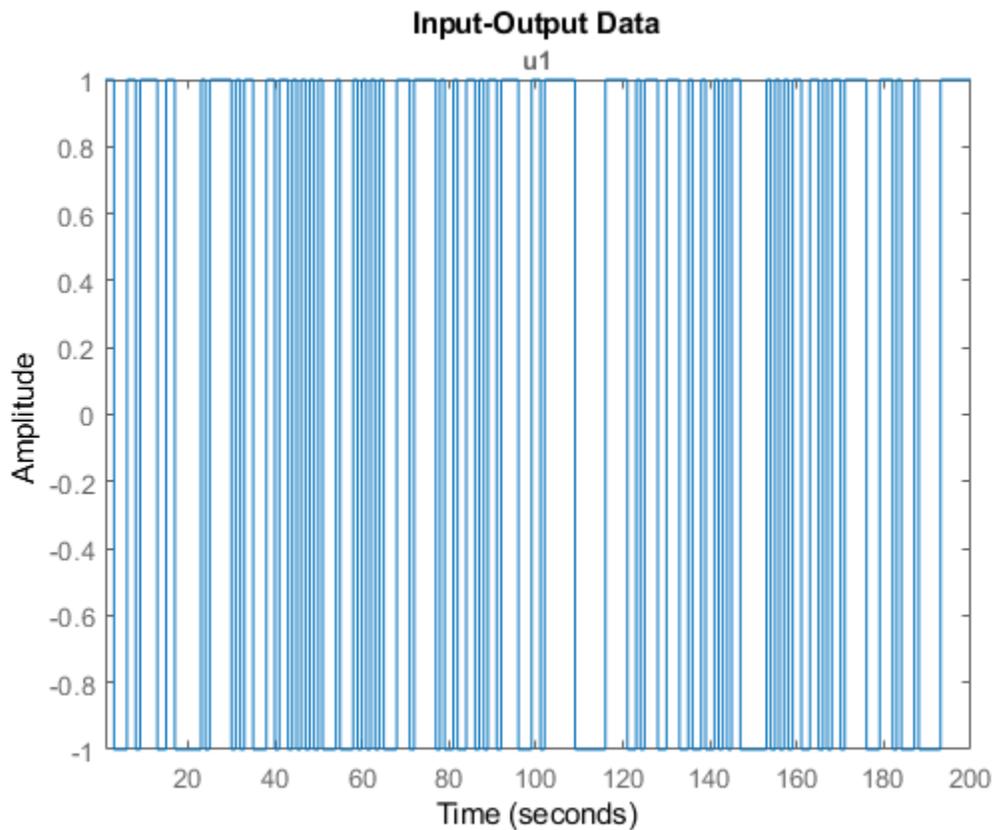
`u` is a column vector of length 200. The values in `u` are either -1 or 1.

Create an `iddata` object from the generated signal. For this example, specify the sample time as 1 second.

```
u = iddata([], u, 1);
```

To examine the signal, plot it.

```
plot(u)
```



The generated signal is a random binary input signal with values -1 or 1. You can use the generated input signal to simulate the output of your system using the `sim` command.

### Generate a Multichannel Random Binary Input Signal

Generate a two-channel random binary input signal with 200 samples.

```
N = 200;  
u = idinput([N,2]);
```

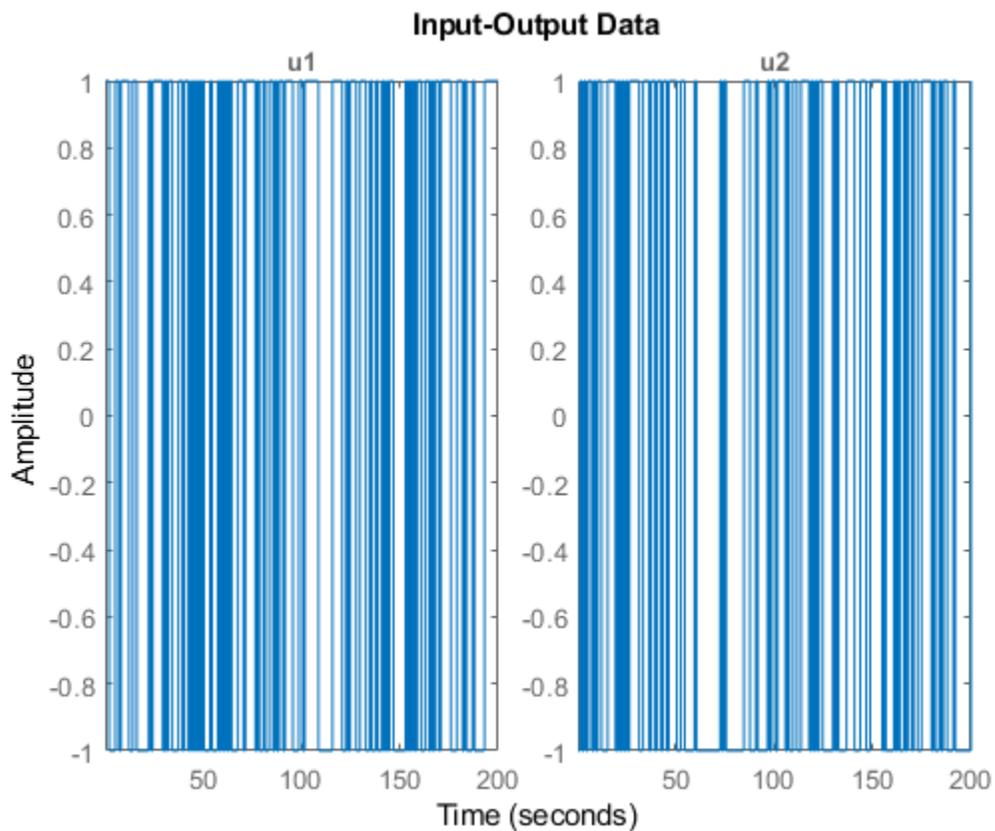
`u` is a 200-by-2 matrix with values -1 or 1.

Create an `iddata` object from the generated signal. For this example, specify the sample time as 1 second.

```
u = iddata([],u,1);
```

Plot the signals for the two channels, and examine the signals.

```
plot(u)
```



The plot shows the two generated random binary signals with values -1 or 1.

## Generate a Periodic Random Binary Input Signal

Generate a single-channel periodic random binary input signal with a period of 10 samples and 5 periods in the signal.

```
NumChannel = 1;  
Period = 10;  
NumPeriod = 5;  
u = idinput([Period,NumChannel,NumPeriod]);
```

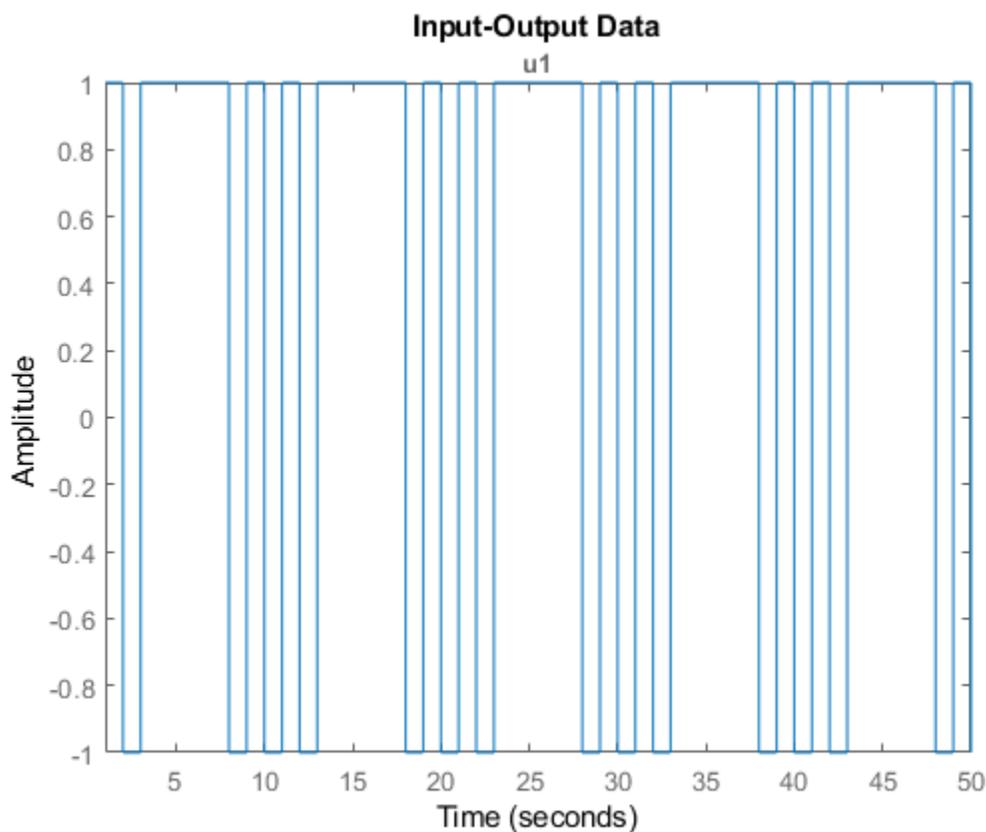
`u` is a column vector of length 50 (= `Period*NumPeriod`). The values in `u` are either -1 or 1.

Create an `iddata` object from the generated signal. Specify the sample time as 1 second.

```
u = iddata([],u,1);
```

Plot the signal.

```
plot(u)
```



As specified, the generated single-channel periodic random binary input signal has a period of 10 seconds, and there are 5 whole periods in the signal.

### **Generate a Periodic Random Gaussian Input Signal in Specified Frequency Range**

Generate a single-channel periodic random Gaussian input signal with a period of 50 samples and 5 periods in the signal. First generate the signal using the entire frequency range, then specify a passband.

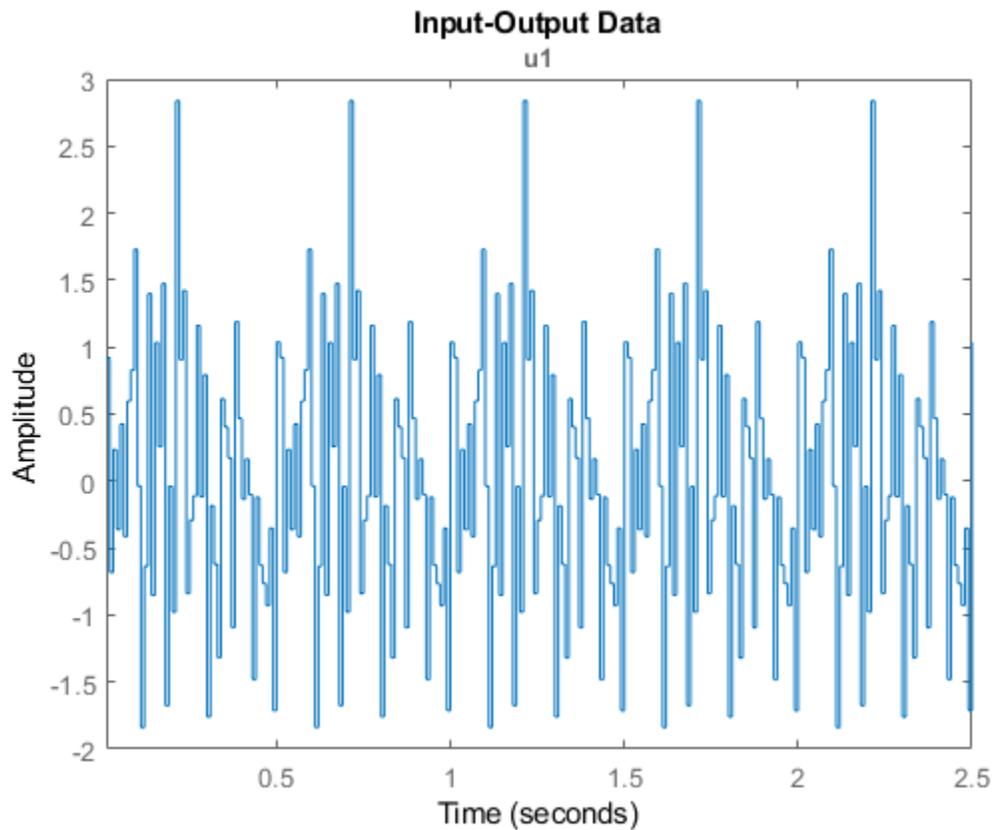
```
NumChannel = 1;  
Period = 50;
```

```
NumPeriod = 5;  
u = idinput([Period,NumChannel,NumPeriod],'rgs');
```

u is a column vector of length 250 (= Period\*NumPeriod).

Create an `iddata` object from the generated signal, and plot the signal. For this example, specify the sample time as 0.01 seconds.

```
u = iddata([],u,0.01);  
plot(u)
```



The plot shows that `u` contains a random segment of 50 samples, repeated 5 times. The signal is a Gaussian white noise signal with zero mean and variance one.

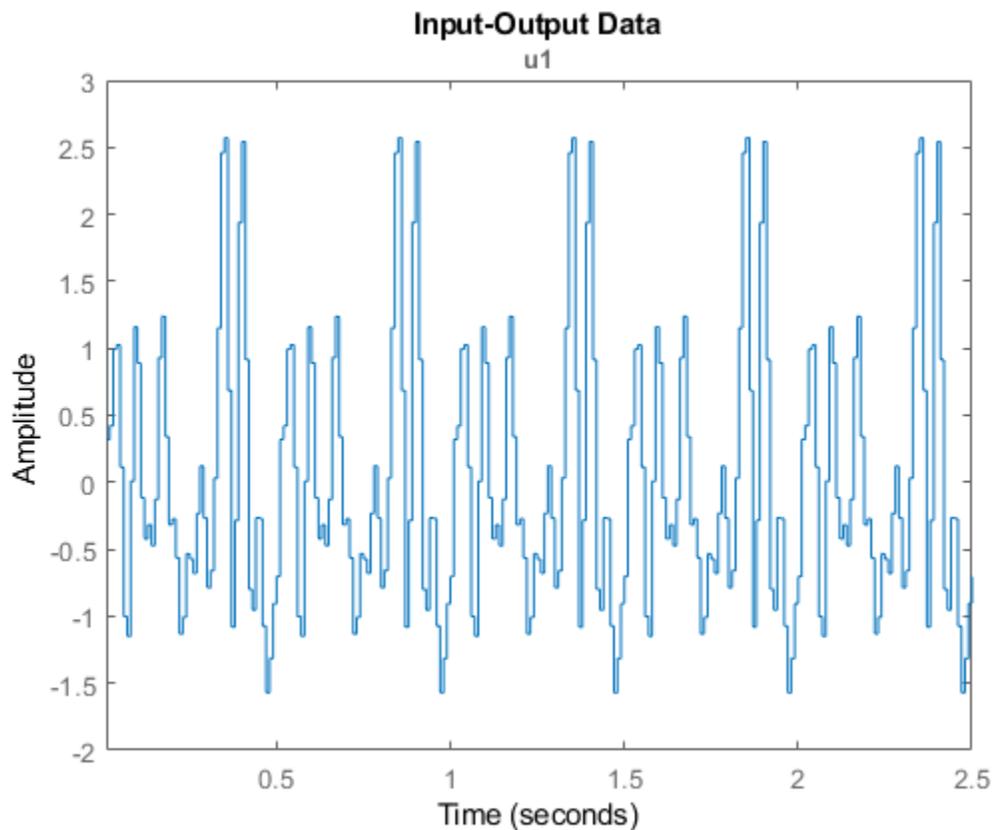
Since the sample time is 0.01 seconds, the generated signal has a period of 0.5 seconds. The frequency content of the signal spans the entire available range (0-50 Hz).

Now specify a passband between 0 and 25 Hz (= 0.5 times the Nyquist frequency).

```
Band = [0 0.5];  
u2 = idinput([Period,NumChannel,NumPeriod], 'rgs', Band);
```

Create an iddata object, and plot the signal.

```
u2 = iddata([],u2,0.01);  
plot(u2)
```



The frequency content of the generated signal u2 is limited to 0-25 Hz.

### Generate a Nonperiodic Pseudorandom Binary Input Signal

A pseudorandom binary input signal (PRBS) is a deterministic signal whose frequency properties mimic white noise. A PRBS is inherently periodic with a maximum period length of  $2^n - 1$ , where integer  $n$  is the order of the PRBS. For more information, see “Pseudorandom Binary Signals” on page 1-613.

Specify that the single-channel PRBS value switches between -2 and 2.

```
Range = [-2,2];
```

Specify the clock period of the signal as 1 sample. That is, the signal value can change at each time step. For PRBS signals, the clock period is specified in `Band = [0 B]`, where `B` is the inverse of the required clock period.

```
Band = [0 1];
```

Generate a nonperiodic PRBS of length 100 samples.

```
u = idinput(100,'prbs',Band,Range);
```

Warning: The PRBS signal delivered is the 100 first values of a full sequence of length

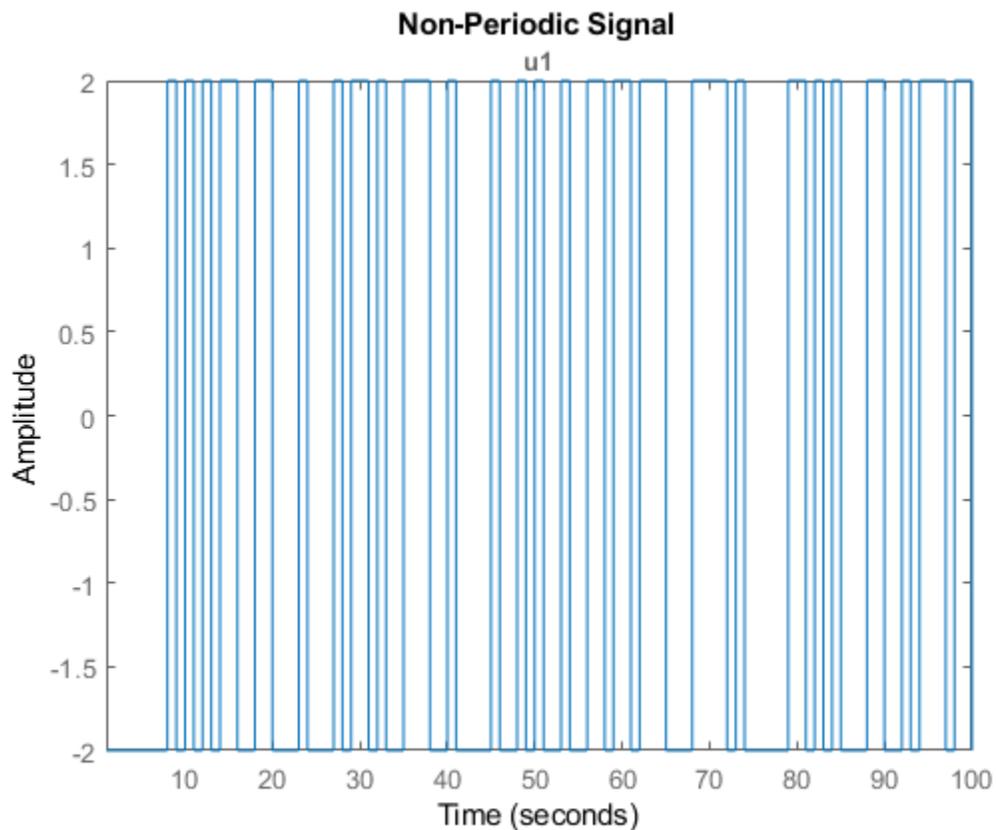
A PRBS is inherently periodic. To generate a nonperiodic signal, the software generates a maximum length PRBS of length 127 that has a period greater than the required number of samples, 100. The software returns the first 100 samples of the generated PRBS. This action ensures that the generated signal is not periodic, as indicated in the generated warning.

Create an `iddata` object from the generated signal. For this example, specify the sample time as 1 second.

```
u = iddata([],u,1);
```

Plot, and examine the generated signal.

```
plot(u);  
title('Non-Periodic Signal')
```



The generated signal is a nonperiodic PRBS of length 100 that switches between -2 and 2.

### Generate a Periodic Pseudorandom Binary Input Signal

Specify that the pseudorandom binary input signal (PRBS) switches between -2 and 2.

Range = [-2,2];

Specify the clock period of the signal as 1 sample. That is, the signal value can change at each time step. For PRBS signals, the clock period is specified in Band = [0 B], where B is the inverse of the required clock period.

```
Band = [0 1];
```

Generate a single-channel, periodic PRBS with a period of 100 samples and 3 periods in the signal.

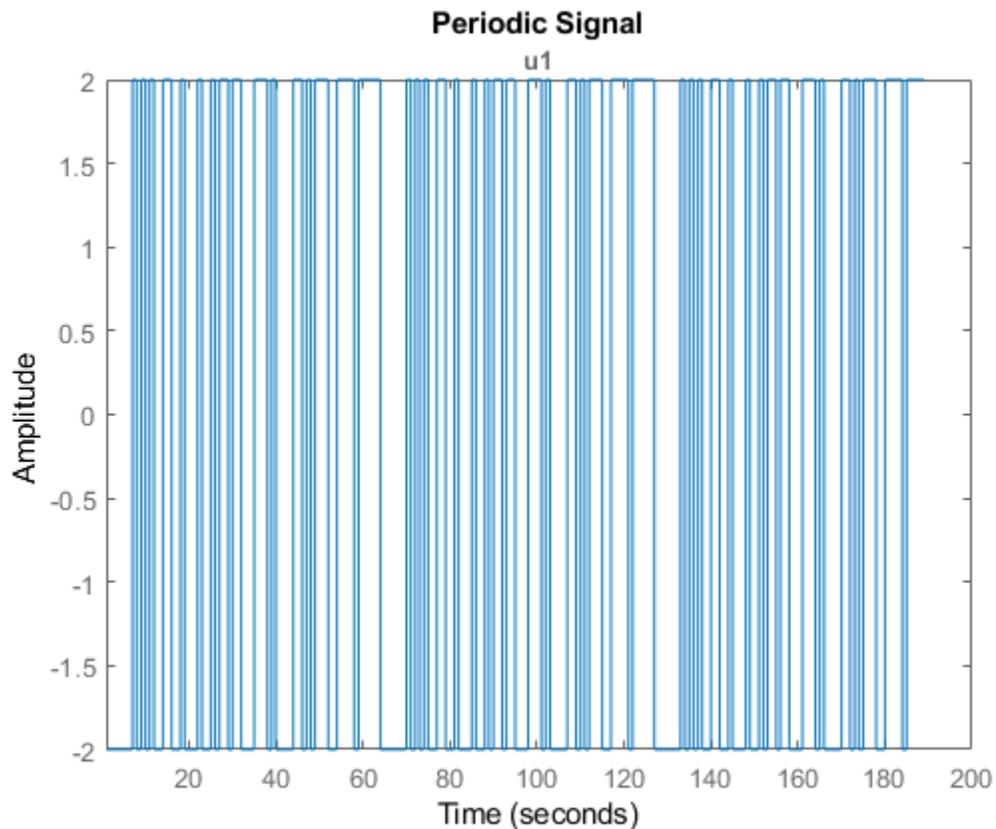
```
u1 = idinput([100,1,3], 'prbs', Band, Range);
```

Warning: The period of the PRBS signal was changed to 63. Accordingly, the length of the

A PRBS is inherently periodic with a maximum period length of  $2^n - 1$ , where integer  $n$  is the order of the PRBS. If the period you specify is not equal to a maximum length PRBS, the software adjusts the period of the generated signal to obtain an integer number of maximum length PRBS, and issues a warning. For more information about maximum length PRBS, see “Pseudorandom Binary Signals” on page 1-613. In this example, the desired period, 100, is not equal to a maximum length PRBS, thus the software instead generates a maximum length PRBS of order  $n = \text{floor}(\log_2(\text{Period})) = 6$ . Thus, the period of the PRBS signal is 63 ( $= 2^6 - 1$ ), and the length of the generated signal is 189 ( $= \text{NumPeriod} * 63$ ). This result is indicated in the generated warning.

Create an `iddata` object from the generated signal, and plot the signal. Specify the period of the signal as 63 samples.

```
u1 = iddata([],u1,1, 'Period', 63);  
plot(u1)  
title('Periodic Signal')
```



The generated signal is a periodic PRBS with three periods.

### **Generate Pseudorandom Binary Input Signal with Specified Clock Period**

Generate periodic and nonperiodic pseudorandom binary input signals (PRBS) with specified clock period.

Generate a single-channel PRBS that switches between -2 and 2. Specify the clock period of the signal as 4 samples. That is, the signal has to stay constant for at least 4 consecutive samples before it can change. For PRBS signals, the clock period is specified in  $\text{Band} = [0 B]$ , where  $B$  is the inverse of the required clock period.

```
Range = [-2,2];  
Band = [0 1/4];
```

First generate a nonperiodic signal of length 100.

```
u1 = idinput(100, 'prbs', Band, Range);
```

Warning: The PRBS signal delivered is the 100 first values of a full sequence of length

To understand the generated warning, first note that the code is equivalent to generating a single-channel PRBS with a 100-sample period and 1 period.

```
u1 = idinput([100,1,1], 'prbs', Band, Range);
```

The generated PRBS signal has to remain constant for at least 4 samples before the value can change. To satisfy this requirement, the software first computes the order of the smallest possible maximum length PRBS as  $n = \text{floor}(\log_2(\text{Period} \cdot B)) = 4$  and period  $2^n - 1 = 15$ . For information about maximum length PRBS, see “Pseudorandom Binary Signals” on page 1-613. The software then stretches this PRBS such that the period of the stretched signal is  $P = (1/B)(2^n - 1) = 60$ .

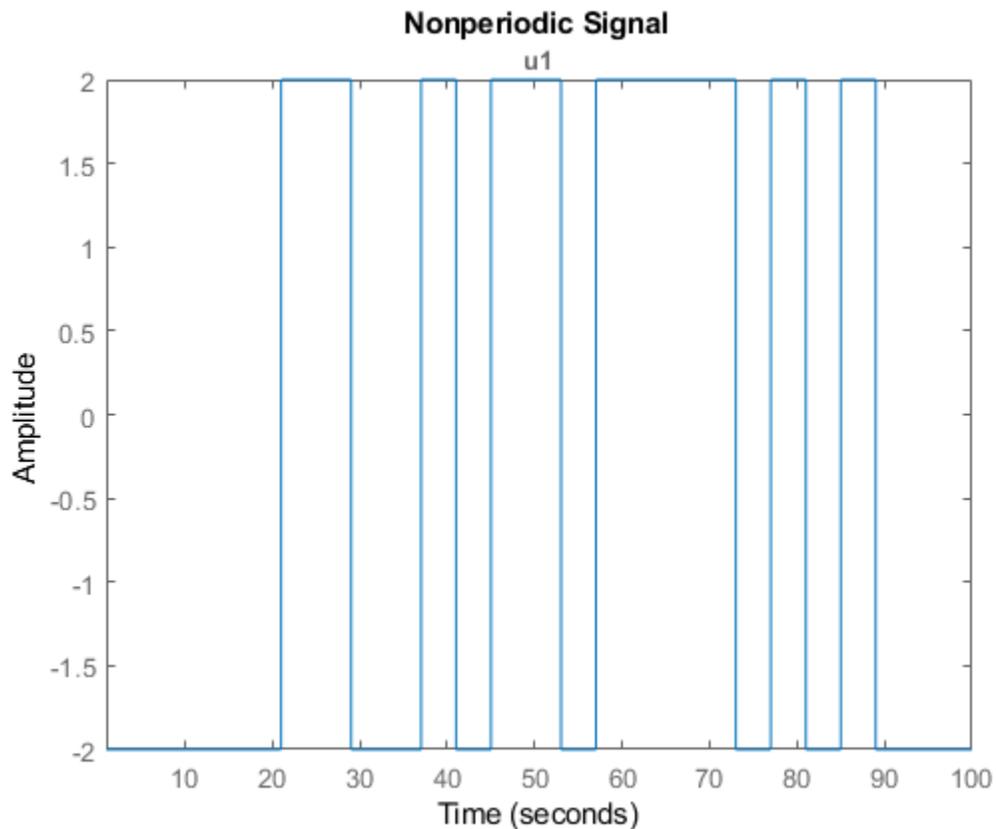
However, since this period is less than the specified length, 100, the software computes instead a maximum length PRBS of order  $m = n+1 = 5$ . The software then stretches this PRBS such that the period is now  $P2 = (1/B)(2^m - 1) = 124$ . The software returns the first 100 samples of this signal as `u1`. This result ensures that the generated signal is not periodic but is constant for every 4 samples.

Create an `iddata` object from the generated signal. For this example, specify the sample time as 1 second.

```
u1 = iddata([], u1, 1);
```

Plot, and examine the signal.

```
plot(u1);  
title('Nonperiodic Signal')
```



The generated signal is a nonperiodic PRBS of length 100. The signal remains constant for at least 4 samples before each change in value. Thus, the signal satisfies the clock period specified in Band.

Now generate a periodic signal with a 100-sample period and 3 periods.

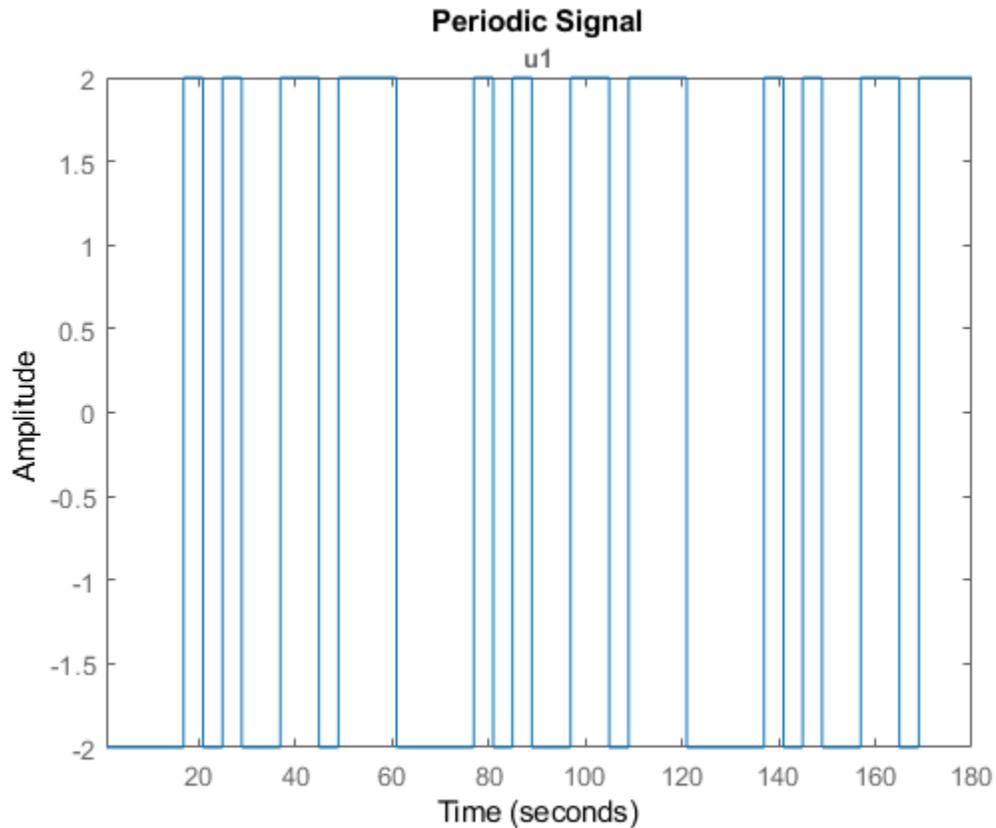
```
u2 = idinput([100,1,3], 'prbs', Band, Range);
```

Warning: The period of the PRBS signal was changed to 60. Accordingly, the length of the

To generate a periodic signal with specified clock period, the software generates u2 as 3 repetitions of the original stretched signal of period  $P = 60$ . Thus, the length of u2 is  $P \cdot \text{NumPeriod} = 60 \cdot 3 = 180$ . This change in period and length of the generated signal is indicated in the generated warning.

Create an `iddata` object from the generated signal, and plot the signal. Specify the period of the signal as 60 seconds.

```
u2 = iddata([],u2,1,'Period',60);  
plot(u2)  
title('Periodic Signal')
```



The generated signal is a periodic PRBS with a 60-second period and 3 periods. The signal remains constant for at least 4 samples before each change in value. Thus, the signal satisfies the specified clock period.

## Generate a Sum-of-Sinusoids Signal

You can generate a sum-of-sinusoids signal using default characteristics for the sine waves. Alternatively, you configure the number of sine waves, and the frequencies and phases of the sine waves. This example shows both approaches.

Specify that the signal has 50 samples in each period and 3 periods. Also specify that the signal amplitude range is between -1 and 1.

```
Period = 50;  
NumPeriod = 3;  
Range = [-1 1];
```

Specify the frequency range of the signal. For a sum-of-sinusoids signal, you specify the lower and upper frequencies of the passband in fractions of the Nyquist frequency. In this example, use the entire frequency range between 0 and Nyquist frequency.

```
Band = [0 1];
```

First generate the signal using default characteristics for the sine waves. By default, the software uses 10 sine waves to generate the signal. The software assigns a random phase to each sinusoid, and then changes these phases 10 times to get the smallest signal spread. The signal spread is the difference between the minimum and the maximum value of the signal over all samples.

```
[u,freq] = idinput([Period 1 NumPeriod], 'sine',Band,Range);
```

The software returns the sum-of-sinusoids signal in `u` and the frequencies of the sinusoids in `freq`. The values in `freq` are scaled assuming that the sample time is 1 time unit. Suppose that the sample time is 0.01 hours. To retrieve the actual frequencies in rad/hours, divide the values by the sample time.

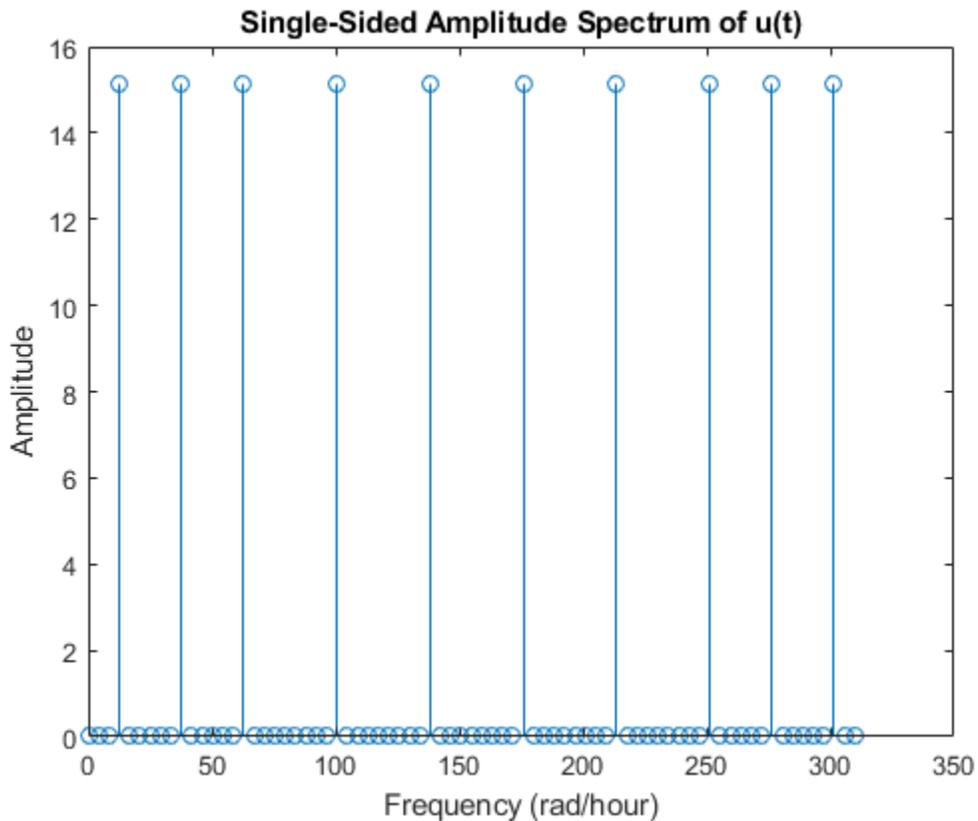
```
Ts = 0.01; % Sample time in hours  
freq = freq/Ts;  
freq(1)
```

```
ans = 12.5664
```

`freq(1)` is the frequency of the first sine wave. To see how the software chooses the frequencies, see the `SineData` argument description on the `idinput` reference page.

To verify that 10 sine waves were used to generate the signal, you can view the frequency content of the signal. Perform a Fourier transform of the signal, and plot the single-sided amplitude spectrum of the signal.

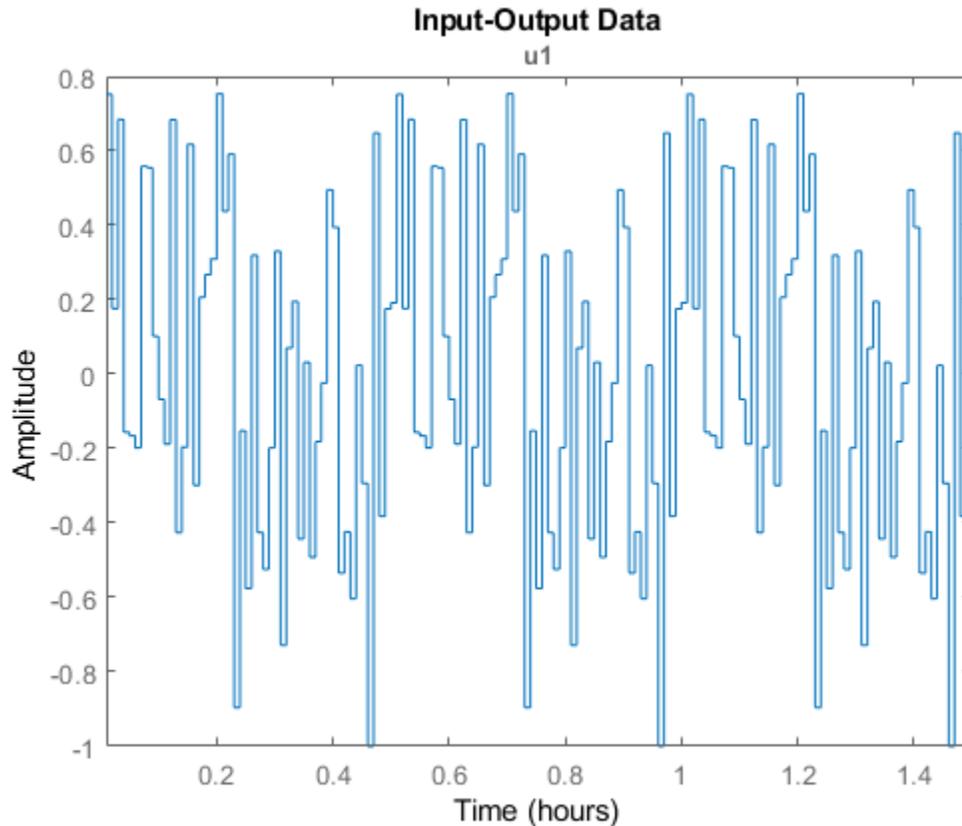
```
ufft = fft(u);  
Fs = 2*pi/Ts; % Sampling frequency in rad/hour  
L = length(u);  
w = (0:L-1)*Fs/L;  
stem(w(1:L/2),abs(ufft(1:L/2))) % Plot until Nyquist frequency  
title('Single-Sided Amplitude Spectrum of u(t)')  
xlabel('Frequency (rad/hour)')  
ylabel('Amplitude')
```



The generated plot shows the frequencies of the 10 sine waves used to generate the signal. For example, the plot shows that the first sine wave has a frequency of 12.57 rad/hour, the same as `freq(1)`.

Convert the generated signal into an iddata object, and plot the signal. Specify the sample time as 0.01 hours.

```
u = iddata([],u,Ts,'TimeUnit','hours');
plot(u)
```



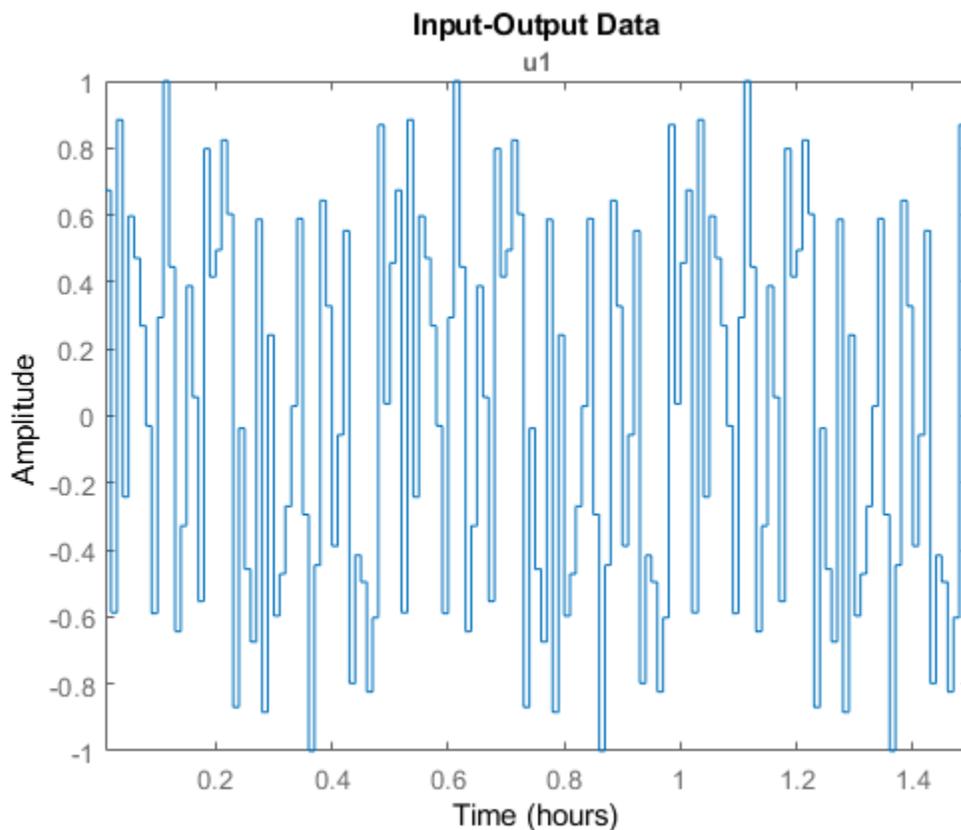
The signal  $u$  is generated using 10 sinusoids and has a period of 0.5 hours and 3 periods.

Now modify the number, frequency, and phase of the sinusoids that are used to generate the sum-of-sinusoids signal. Use 12 sinusoids and try 15 different sets of phases. To set the frequencies of the sinusoids, specify `GridSkip = 2`. The software selects the frequencies of the sinusoids from the intersection of the frequency grid  $2\pi*[1:\text{GridSkip}:\text{fix}(\text{Period}/2)]/\text{Period}$  and the passband  $\pi*\text{Band}$ .

```
NumSinusoids = 12;  
NumTrials = 15;  
GridSkip = 2;  
SineData = [NumSinusoids,NumTrials,GridSkip];  
u2 = idinput([Period 1 NumPeriod],'sine',Band,Range,SineData);
```

Convert the generated signal into an iddata object, and plot the signal.

```
u2 = iddata([],u2,Ts,'TimeUnit','hours');  
plot(u2)
```



The signal  $u_2$  is generated using 12 sinusoids and has a period of 0.5 hours and 3 periods.

## Input Arguments

### **N — Number of generated input data samples**

real positive integer

Number of generated input data samples, specified as a real positive integer. For a single-channel input data, the generated input *u* has *N* rows. For an *Nu*-channel input data, *u* is returned as an *N*-by-*Nu* matrix, where each channel signal has length *N*.

### **Nu — Number of input channels**

1 (default) | real positive integer

Number of input channels in generated signal, specified as a real positive integer.

### **Period — Number of samples in each period**

real positive integer

Number of samples in each period of generated signal, specified as a real positive integer. Use this input to specify a periodic signal. Also specify the number of periods in *NumPeriod*. Each generated input channel signal has *NumPeriod*\**Period* samples.

### **NumPeriod — Number of periods in generated signal**

1 (default) | real positive integer

Number of periods in generated signal, specified as a real positive integer. Use this input to specify a periodic signal. Also specify the signal *Period*. Each generated input channel signal has *NumPeriod*\**Period* samples.

### **Type — Type of generated signal**

'rbs' (default) | 'rgs' | 'prbs' | 'sine'

Type of generated signal, specified as one of the following values:

- 'rbs' — Generates a random binary signal. A random binary signal is a random process that assumes only two values. You can specify these values using *Range*. To generate a band-limited signal, specify the passband in *Band*. To generate a periodic signal, specify *Period* and *NumPeriod*.
- 'rgs' — Generates a random Gaussian signal. The generated Gaussian signal has mean  $\mu$  and standard deviation  $\sigma$  such that  $[\mu-\sigma, \mu+\sigma]$  equals *Range*. To generate a band-limited Gaussian signal, specify the passband in *Band*. To generate a periodic Gaussian signal with an *n* samples period that repeats itself *m* times, specify *Period* as *n* and *NumPeriod* as *m*.

- 'prbs' — Generates a pseudorandom binary signal (PRBS). A PRBS is a periodic, deterministic signal with white-noise-like properties that shifts between two values. You can specify these two values using `Range`. You can also specify the clock period, the minimum number of sampling intervals for which the value of the signal does not change. You specify the inverse of the clock period in `Band`.

The length of the generated signal is not always the same as what you specify. The length depends on whether you require a periodic or nonperiodic signal and also on the clock period you specify. For more information, see “Pseudorandom Binary Signals” on page 1-613.

- 'sine' — Generates a signal that is a sum-of-sinusoids. The software selects the frequencies of the sinusoids to be equally spread over a chosen grid and assigns each sinusoid a random phase. The software then tries several random phases for each sinusoid and selects the phases that give the smallest signal spread. The signal spread is the difference between the minimum and the maximum value of the signal over all samples. The amplitude of the generated sum-of-sinusoids signal is scaled to satisfy the `Range` you specify.

You can specify the characteristics of the sine waves used to generate the signal, such as the number of sine waves and their frequency separation, in the `SineData` argument.

## **Band — Frequency range of generated signal**

[0 1] (default) | 1-by-2 row vector

Frequency range of generated signal, specified as a 1-by-2 row vector containing minimum and maximum frequency values.

- If `Type` is 'rgs', 'rbs', or 'sine' — Specify `Band` as a passband [`wlow` `whigh`]. Where, `wlow` and `whigh` are the lower and upper frequencies of the passband, expressed in fractions of the Nyquist frequency. For example, to generate an input with white noise characteristics, use `Band = [0 1]`.

The software achieves the frequency contents for a random Gaussian signal ('rgs') using `idfilt` with an eighth-order Butterworth, noncausal filter. For generating a random binary signal ('rbs'), the software uses the same filter and then makes the signal binary. Thus, the frequency content in the generated random binary signal may not match the specified passband.

For 'sine' signals, the frequencies of the sinusoids are selected to be equally spread over a chosen grid in the specified passband. For more information, see the `SineData` argument description.

- If Type is 'prbs' — Specify Band as  $[0 \ B]$ , where B is the inverse of the clock period of the signal. The clock period is the minimum number of sampling intervals for which the value of the signal does not change. Thus, the generated signal is constant over intervals of length  $1/B$  samples. If  $1/B$  is not an integer, the software uses  $\text{floor}(1/B)$  as the clock period.

### Range — Generated input signal range

`[-1, 1]` (default) | two-element row vector

Generated input signal range, specified as a two-element row vector of the form `[umin, umax]`.

- If Type is 'rbs' or 'prbs' — The generated signal  $u$  has values  $u_{\min}$  or  $u_{\max}$ .
- If Type is 'sine' — The generated signal  $u$  has values between  $u_{\min}$  and  $u_{\max}$ .
- If Type is 'rgs' — The generated Gaussian signal has mean  $\mu$  and standard deviation  $\sigma$  such that  $u_{\min}$  and  $u_{\max}$  are equal to  $\mu - \sigma$  and  $\mu + \sigma$ , respectively. For example, Range = `[-1, 1]` returns a Gaussian white noise signal with zero mean and variance one.

### SineData — Characterization of sinusoids

`[10, 10, 1]` (default) | three-element row vector

`[NumSinusoids, NumTrials, GridSkip]`

Characterization of sinusoids used to generate a sum-of-sinusoids signal, specified as a three-element row vector `[NumSinusoids, NumTrials, GridSkip]`. Where,

- NumSinusoids is the number of sinusoids used to generate the signal. The default value is 10.
- NumTrials is the number of different random relative phases of the sinusoids that the software tries to find the lowest signal spread. The signal spread is the difference between the minimum and the maximum value of the signal over all samples.

The maximum amplitude of the sum-of-sinusoids signal depends on the relative phases of the different sinusoids. To find the phases that give the smallest signal spread, the software tries NumTrials different random choices of phases to find the best phase values. For example, suppose that NumSinusoids is 20 and NumTrials is 5. The software tries 5 different sets of relative phases for the 20 sinusoids, and selects the phases that give the smallest signal spread. The default value for NumTrials is 10.

- GridSkip is used to characterize the frequency of the sinusoids. The software selects the frequency of the sinusoids from the intersection of the frequency grid

$2\pi * [1:\text{GridSkip}:\text{fix}(\text{Period}/2)]/\text{Period}$  and the pass band  $\pi * [\text{Band}(1) \text{Band}(2)]$ . For multichannel input signals, the software uses different frequencies from this frequency grid to generate the different input channels. You can use `GridSkip` for controlling odd and even frequency multiples, for example, to detect nonlinearities of different kinds.

To extract the frequencies `freq` that are selected by the software to generate the signal, use the following syntax.

```
[u,freq] = idinput(__)
```

## Output Arguments

### **u** — Generated input signal

column vector | matrix

Generated input signal, returned as a column vector of length `N` for a single-channel input or an `N`-by-`Nu` matrix for an `Nu`-channel signal. You use the generated signal to simulate the response of your system using `sim`.

You can create an `iddata` object from `u` by specifying output data as `[]`.

```
u = iddata([],u);
```

In the `iddata` object, you can also specify the properties of the signal such as sample time, input names, and periodicity.

### **freq** — Frequencies of sine waves

column vector | matrix

Frequencies of sine waves used for sum-of-sinusoids signal, returned as a column vector of length equal to the number of sinusoids, `NumSinusoids`. You specify `NumSinusoids` in the `SineData` argument. The frequency values are scaled assuming the sample time is 1 time unit. To retrieve the actual frequencies, divide the values by the sample time. For an example, see “Generate a Sum-of-Sinusoids Signal” on page 1-604.

For multichannel input signals, `freq` is an `Nu`-by-`NumSinusoids` matrix where the  $k$ th row contains the frequencies corresponding to the  $k$ th channel. For information about how the software selects the frequencies, see the `SineData` argument description.

## Definitions

### Pseudorandom Binary Signals

A pseudorandom binary signal (PRBS) is a periodic, deterministic signal with white-noise-like properties that shifts between two values.

A PRBS is generated as:

$$u(t) = \text{rem}(a_1 u(t-1) + \dots + a_n u(t-n), 2)$$

Here,  $u(t-1), \dots, u(t-n)$  is the vector of past inputs,  $n$  is the PRBS order, and  $\text{rem}$  denotes the remainder when  $(a_1 u(t-1) + \dots + a_n u(t-n))$  is divided by 2. Thus, a PRBS can only take the values 0 and 1. The software scales these values according to the Range you specify.

In addition, the vector of past inputs  $u(t-1), \dots, u(t-n)$  can only take  $2^n$  values. Out of these values, the state with all zeros is ignored because it will result in future signals equal to zero. Thus, a PRBS is an inherently periodic signal with a maximum period length of  $2^n - 1$ . The following table lists the maximum length possible for different orders  $n$  of the PRBS.

Order $n$	Maximum length PRBS ( $2^n - 1$ )
2	3
3	7
4	15
5	31
6	63
7	127
⋮	⋮
32	4294967295

---

**Note** The software does not generate signals with period greater than  $2^{32} - 1$ .

---

Since PRBS are inherently periodic, the length and period of the generated signal depends on the clock period that you specify and whether you require a periodic or nonperiodic signal. The clock period is the minimum number of sampling intervals for which the value of the signal does not change. You specify the clock period in *Band*.

*Clock period = 1 sample (Band = [0 B] = [0 1]):*

- To generate a *nonperiodic signal* of length *N*, (*NumPeriod* = 1), the software first computes a maximum length PRBS with a period greater than *N*. The software then returns the first *N* samples of the PRBS as *u*. This action ensures that *u* is not periodic. For example, if *N* is 100, the software creates a maximum length PRBS of period 127 (order 7), and returns the first 100 samples as *u*.

For an example, see “Generate a Nonperiodic Pseudorandom Binary Input Signal” on page 1-598.

- To generate a *periodic signal* (*NumPeriod* > 1), the software adjusts the period of the signal to obtain an integer number of maximum length PRBS. To do so, the software computes a PRBS of order  $n = \text{floor}(\log_2(\text{Period}))$  and period  $P = 2^n - 1$ . The signal *u* is then generated as *NumPeriod* repetitions of this PRBS signal of period *P*. Thus, the length of *u* is  $P * \text{NumPeriod}$ .

For an example, see “Generate a Periodic Pseudorandom Binary Input Signal” on page 1-599.

In the multiple-input channel case, the signals are maximally shifted. That is, the overlap between the different inputs is minimized. This means  $\text{Period}/\text{NumPeriod}$  is an upper bound for the model orders that you can estimate using such a signal.

*Clock period > 1 sample (Band = [0 B], where  $B < 1$ ):*

The generated signal has to remain constant for at least  $1/B$  samples. To satisfy this requirement, the software first computes the order of the smallest possible maximum length PRBS as  $n = \text{floor}(\log_2(\text{Period} * B))$  and period  $2^n - 1$ . The software then stretches the PRBS such that period of the stretched signal is  $P = B^{-1} (2^n - 1)$ .

- To generate a *nonperiodic signal* of length *N*, if the period *P* of the stretched signal is greater than or equal to *N*, the software returns the first *N* samples of the stretched signal as *u*. This ensures that *u* is nonperiodic but constant for every  $1/B$  samples. Note that for a nonperiodic signal, *Period* is equal to *N*.

If the period  $P$  is less than  $N$ , the software computes instead a maximum length PRBS of order  $n_2 = n+1$ . The software then stretches this PRBS such that the period is now  $P_2 = B^{-1}(2^{n_2} - 1)$ . The software then returns the first  $N$  samples of this signal as  $u$ .

- To generate a *periodic signal*, the software generates  $u$  as `NumPeriod` repetitions of the stretched signal of period  $P$ . Thus, the length of  $u$  is  $P * \text{NumPeriod}$ .

For an example, see “Generate Pseudorandom Binary Input Signal with Specified Clock Period” on page 1-601.

## References

- [1] Söderström, T. and P. Stoica., Chapter C5.3 in *System Identification*, Prentice Hall, 1989.
- [2] Ljung, L., Section 13.3 in *System Identification: Theory for the User*, Prentice Hall PTR, 1999.

## See Also

`iddata` | `sim`

## Topics

“Ways to Obtain Identification Data”

“Generate Data Using Simulation”

“Simulate and Predict Identified Model Output”

**Introduced before R2006a**

## idnlarx

Nonlinear ARX model

### Syntax

```
sys = idnlarx(Orders)
sys = idnlarx(Orders,Nonlinearity)
sys = idnlarx(Orders,Nonlinearity,Name,Value)
sys = idnlarx(LinModel)
sys = idnlarx(LinModel,Nonlinearity)
sys = idnlarx(LinModel,Nonlinearity,Name,Value)
```

### Description

`sys = idnlarx(Orders)` creates a nonlinear ARX model with the specified orders using a default wavelet network nonlinearity estimator.

`sys = idnlarx(Orders,Nonlinearity)` specifies a nonlinearity estimator for the model.

`sys = idnlarx(Orders,Nonlinearity,Name,Value)` specifies additional attributes of the `idnlarx` model structure using one or more `Name, Value` pair arguments.

`sys = idnlarx(LinModel)` uses a linear ARX model `LinModel` to specify the model orders and the initial values of the linear coefficients of the model.

`sys = idnlarx(LinModel,Nonlinearity)` specifies a nonlinearity estimator for the model.

`sys = idnlarx(LinModel,Nonlinearity,Name,Value)` specifies additional attributes of the `idnlarx` model structure using one or more `Name, Value` pair arguments.

## Object Description

`idnlarx` represents a nonlinear ARX model, which is an extension of the linear ARX structure and contains linear and nonlinear functions. For more information, see “Nonlinear ARX Model Extends the Linear ARX Structure”.

Use the `nlarx` command to both construct an `idnlarx` object and estimate the model parameters.

You can also use the `idnlarx` constructor to create the nonlinear ARX model and then estimate the model parameters using `nlarx` or `pem`.

For `idnlarx` object properties, see “Properties” on page 1-621.

## Examples

### Create Nonlinear ARX Model with Default Wavelet Network Nonlinearity

```
m = idnlarx([2 2 1]);
```

### Create and Configure Nonlinear ARX Model

Create a nonlinear ARX model with specific orders.

```
M = idnlarx([3 2 1]);
```

Configure the model with the following property settings:

- Sigmoid network nonlinearity with default settings
- Use inputs only as nonlinear regressors

```
M.Nonlinearity = 'sigmoidnet';  
M.NonlinearRegressors = 'input';
```

## Create Nonlinear ARX Model with Sigmoid Network Nonlinearity

```
m = idnlarx([2 3 1],sigmoidnet('NumberOfUnits',15));
```

## Create Nonlinear ARX Model Without Nonlinear Function in Nonlinearity Estimator

```
m = idnlarx([2 2 1],[,]);
```

## Create Nonlinear ARX Model with Custom Regressors

Specify two custom regressors.

```
C = {'y1(t-1)^2', 'y1(t-2)*u1(t-3)'};
```

Create a nonlinear ARX model with custom regressors and no standard regressors.

```
sys = idnlarx([0 0 0], 'wavenet', 'CustomRegressors', C);
```

## Create Nonlinear ARX Model Using Linear ARX Model

Construct a linear ARX model.

```
A = [1 -1.2 0.5];  
B = [0.8 1];  
LinearModel = idpoly(A, B, 'Ts', 0.1);
```

Construct nonlinear ARX model using the linear ARX model.

```
m1 = idnlarx(LinearModel);
```

## Input Arguments

### Orders — Model orders and delays

1-by-3 vector of positive integers | 1-by-3 vector of matrices

Model orders and delays for defining the regressor configuration, specified as a 1-by-3 vector, `[na nb nk]`.

For a model with  $n_y$  output channels and  $n_u$  input channels:

- `na` is an  $n_y$ -by- $n_y$  matrix, where `na(i, j)` specifies the number of regressors from the  $j$ th output used to predict the  $i$ th output.
- `nb` is an  $n_y$ -by- $n_u$  matrix, where `nb(i, j)` specifies the number of regressors from the  $j$ th input used to predict the  $i$ th output.
- `nk` is an  $n_y$ -by- $n_u$  matrix, where `nk(i, j)` specifies the lag in the  $j$ th input used to predict the  $i$ th output.

```
na = [1 2; 2 3]
nb = [1 2 3; 2 3 1];
nk = [2 0 3; 1 0 5];
```

The estimation data for this system has three inputs (`u1`, `u2`, `u3`) and two outputs (`y1`, `y2`). Consider the regressors used to predict output, `y2(t)`:

- Since `na(2, :)` is `[2 3]`, the contributing regressors from the outputs are:
  - `y1(t-1)` and `y1(t-2)`
  - `y2(t-1)`, `y2(t-2)`, and `y2(t-3)`
- Since `nb(2, :)` is `[2 3 1]` and `nk(2, :)` is `[1 0 5]`, the contributing regressors from the inputs are:
  - `u1(t-1)` and `u1(t-2)`
  - `u2(t)`, `u2(t-1)`, and `u2(t-2)`
  - `u3(t-5)`

---

**Note** The minimum lag for regressors based on output variables is always 1, while the minimum lag for regressors based on input variables is dictated by `nk`. Use `getreg` to view the complete set of regressors used by the nonlinear ARX model.

---

### Nonlinearity – Nonlinearity estimator

'wavenet' (default) | 'sigmoidnet' | 'treepartition' | 'linear' | nonlinearity estimator object | array of nonlinearity estimator objects

Nonlinearity estimator, specified as one of the following:

'wavenet' or wavenet object	Wavelet network
'sigmoidnet' or sigmoidnet object	Sigmoid network
'treepartition' or treepartition object	Binary-tree
'linear' or [] or linear object	Linear function
neuralnet object	Neural network — Requires Deep Learning Toolbox™.
customnet object	Custom network — Similar to sigmoidnet, but with a user-defined replacement for the sigmoid function.

For more information, see “Available Nonlinearity Estimators for Nonlinear ARX Models”.

Specifying a character vector, for example 'sigmoidnet', creates a nonlinearity estimator object with default settings. Alternatively, you can specify nonlinearity estimator settings in two ways:

- Use the associated nonlinearity estimator function with Name-Value pair arguments.

```
NL = sigmoidnet('NumberOfUnits',10);
```

- Create and modify a default nonlinearity estimator object.

```
NL = sigmoidnet;  
NL.NumberOfUnits = 10;
```

For  $n_y$  output channels, you can specify nonlinear estimators individually for each channel by setting `Nonlinearity` to an  $n_y$ -by-1 array of nonlinearity estimator objects. To specify the same nonlinearity for all outputs, specify `Nonlinearity` as a character vector or a single nonlinearity estimator object.

Example: 'sigmoidnet' specifies a sigmoid network nonlinearity with a default configuration.

Example: `treepartition('NumberOfUnits',5)` specifies a binary-tree nonlinearity with 5 terms in the binary tree expansion.

Example: `[wavenet('NumberOfUnits',10);sigmoidnet]` specifies different nonlinearity estimators for two output channels.

**LinModel** — Discrete time input-output polynomial model of ARX structure  
idpoly model

Discrete time input-output polynomial model of ARX structure, specified as an `idpoly` model. Create this object using the `idpoly` constructor or estimate it using the `arx` command.

## Name-Value Pair Arguments

Specify optional comma-separated pairs of `Name`, `Value` arguments. `Name` is the argument name and `Value` is the corresponding value. `Name` must appear inside quotes. You can specify several name and value pair arguments in any order as `Name1, Value1, ..., NameN, ValueN`.

Use `Name`, `Value` arguments to specify additional properties on page 1-621 of `idnlarx` models during model creation. For example, `m = idnlarx([2 3 1], 'treepartition', 'InputName', 'Pressure', 'Ts', 0.1)` creates an `idnlarx` model with the input name `Pressure`, and a sample time of 0.1 seconds

## Properties

### `na`, `nb`, `nk`

Model orders and delays for defining the regressor configuration, specified as nonnegative integers.

For a model with  $n_y$  output channels and  $n_u$  input channels:

- `na` is an  $n_y$ -by- $n_y$  matrix, where `na(i, j)` specifies the number of regressors from the  $j$ th output used to predict the  $i$ th output.
- `nb` is an  $n_y$ -by- $n_u$  matrix, where `nb(i, j)` specifies the number of regressors from the  $j$ th input used to predict the  $i$ th output.
- `nk` is an  $n_y$ -by- $n_u$  matrix, where `nk(i, j)` specifies the lag in the  $j$ th input used to predict the  $i$ th output.

```
na = [1 2; 2 3]
nb = [1 2 3; 2 3 1];
nk = [2 0 3; 1 0 5];
```

The estimation data for this system has three inputs (`u1`, `u2`, `u3`) and two outputs (`y1`, `y2`). Consider the regressors used to predict output,  $y_2(t)$ :

- Since  $na(2, :)$  is  $[2 \ 3]$ , the contributing regressors from the outputs are:
  - $y1(t-1)$  and  $y1(t-2)$
  - $y2(t-1)$ ,  $y2(t-2)$ , and  $y2(t-3)$
- Since  $nb(2, :)$  is  $[2 \ 3 \ 1]$  and  $nk(2, :)$  is  $[1 \ 0 \ 5]$ , the contributing regressors from the inputs are:
  - $u1(t-1)$  and  $u1(t-2)$
  - $u2(t)$ ,  $u2(t-1)$ , and  $u2(t-2)$
  - $u3(t-5)$

---

**Note** The minimum lag for regressors based on output variables is always 1, while the minimum lag for regressors based on input variables is dictated by  $nk$ . Use `getreg` to view the complete set of regressors used by the nonlinear ARX model.

---

## CustomRegressors

Regressors constructed from combinations of inputs and outputs, specified as one of the following:

- Cell array of character vectors. For example:
  - `{'y1(t-3)^3', 'y2(t-1)*u1(t-3)', 'sin(u3(t-2))'}`

Each character vector must represent a valid formula for a regressor contributing towards the prediction of the model output. The formula must be written using the input and output names and the time-variable name as variables.

- Array of custom regressor objects, created using `customreg` or `polyreg`.

For a model with  $n_y$  outputs, specify an  $n_y$ -by-1 cell array of `customreg` objects arrays or an  $n_y$ -by-1 cell array of cell array of character vectors.

These regressors are in addition to the standard regressors based on  $na$ ,  $nb$ , and  $nk$ .

**Default:** `{}`

## NonlinearRegressors

Subset of regressors that enter as inputs to the nonlinear block of the model, specified as one of the following:

- 'all' — All regressors
- 'output' — Regressors containing output variables
- 'input' — Regressors containing input variables
- 'standard' — Standard regressors
- 'custom' — Custom regressors
- 'search' — The estimation algorithm performs a search for the best regressor combination. This option must be applied to all output models simultaneously.
- [] — No regressors
- Vector of regressor indices. To determine the number and order of regressors, use `getreg`.

For a model with multiple outputs, specify a cell array of  $n_y$  elements, where  $n_y$  is the number of output channels. For each output, specify one of the preceding options. Alternatively, to apply the same regressor subset to all model outputs, specify [] or any of the character vector options alone, for example 'standard'.

**Default:** 'all'

### Nonlinearity

Nonlinearity estimator, specified as one of the following:

'wavenet' or wavenet object	Wavelet network
'sigmoidnet' or sigmoidnet object	Sigmoid network
'treepartition' or treepartition object	Binary-tree
'linear' or [] or linear object	Linear function
neuralnet object	Neural network — Requires Deep Learning Toolbox.
customnet object	Custom network — Similar to <code>sigmoidnet</code> , but with a user-defined replacement for the sigmoid function.

For more information, see “Available Nonlinearity Estimators for Nonlinear ARX Models”.

Specifying a character vector, for example 'sigmoidnet', creates a nonlinearity estimator object with default settings. Alternatively, you can specify nonlinearity estimator settings in two ways:

- Use the associated nonlinearity estimator function with Name-Value pair arguments:

```
NL = sigmoidnet('NumberOfUnits',10);
```

- Create and modify a default nonlinearity estimator object:

```
NL = sigmoidnet;  
NL.NumberOfUnits = 10;
```

For  $n_y$  output channels, you can specify nonlinear estimators individually for each channel by setting `Nonlinearity` to an array of  $n_y$  nonlinearity estimator objects, where  $n_y$  is the number of outputs. To specify the same nonlinearity for all outputs, specify `Nonlinearity` as a character vector or a single nonlinearity estimator object.

**Default:** 'wavenet'

### Report

Summary report that contains information about the estimation options and results when the model is estimated using the `nlarx` command. Use `Report` to query a model for how it was estimated, including its:

- Estimation method
- Estimation options
- Search termination conditions
- Estimation data fit

The contents of `Report` are irrelevant if the model was constructed.

```
m = idnlarx([2 2 1]);  
m.Report.OptionsUsed
```

```
ans =
```

```
 []
```

If you use `nlarx` to estimate the model, the fields of `Report` contain information on the estimation data, options, and results.

```
load iddata1;  
m = nlarx(z1, [2 2 1]);  
m.Report.OptionsUsed
```

```
Option set for the nlarx command:
```

```
IterativeWavenet: 'auto'  
    Focus: 'prediction'  
    Display: 'off'  
Regularization: [1x1 struct]  
    SearchMethod: 'auto'  
SearchOptions: [1x1 idoptions.search.identsolver]  
    OutputWeight: 'noise'  
    Advanced: [1x1 struct]
```

Report is a read-only property.

For more information on this property and how to use it, see “Output Arguments” on page 1-956 in the `nlrx` reference page and “Estimation Report”.

### **TimeVariable**

Independent variable for the inputs, outputs, and—when available—internal states, specified as a character vector.

**Default:** 't' (time)

### **NoiseVariance**

Noise variance (covariance matrix) of the model innovations  $e$ . Assignable value is an  $ny$ -by- $ny$  matrix. Typically set automatically by the estimation algorithm.

### **Ts**

Sample time.  $T_s$  is a positive scalar representing the sampling period. This value is expressed in the unit specified by the `TimeUnit` property of the model.

**Default:** 1

### **TimeUnit**

Units for the time variable, the sample time  $T_s$ , and any time delays in the model, specified as one of the following values:

- 'nanoseconds'
- 'microseconds'
- 'milliseconds'

- 'seconds'
- 'minutes'
- 'hours'
- 'days'
- 'weeks'
- 'months'
- 'years'

Changing this property has no effect on other properties, and therefore changes the overall system behavior. Use `chgTimeUnit` to convert between time units without modifying system behavior.

**Default:** 'seconds'

### **InputName**

Input channel names, specified as one of the following:

- Character vector — For single-input models, for example, 'controls'.
- Cell array of character vectors — For multi-input models.

Alternatively, use automatic vector expansion to assign input names for multi-input models. For example, if `sys` is a two-input model, enter:

```
sys.InputName = 'controls';
```

The input names automatically expand to {'controls(1)'; 'controls(2)'}

When you estimate a model using an `iddata` object, `data`, the software automatically sets `InputName` to `data.InputName`.

You can use the shorthand notation `u` to refer to the `InputName` property. For example, `sys.u` is equivalent to `sys.InputName`.

Input channel names have several uses, including:

- Identifying channels on model display and plots
- Extracting subsystems of MIMO systems
- Specifying connection points when interconnecting models

**Default:** '' for all input channels

### InputUnit

Input channel units, specified as one of the following:

- Character vector — For single-input models, for example, 'seconds'.
- Cell array of character vectors — For multi-input models.

Use `InputUnit` to keep track of input signal units. `InputUnit` has no effect on system behavior.

**Default:** '' for all input channels

### InputGroup

Input channel groups. The `InputGroup` property lets you assign the input channels of MIMO systems into groups and refer to each group by name. Specify input groups as a structure. In this structure, field names are the group names, and field values are the input channels belonging to each group. For example:

```
sys.InputGroup.controls = [1 2];  
sys.InputGroup.noise = [3 5];
```

creates input groups named `controls` and `noise` that include input channels 1, 2 and 3, 5, respectively. You can then extract the subsystem from the `controls` inputs to all outputs using:

```
sys(:, 'controls')
```

**Default:** Struct with no fields

### OutputName

Output channel names, specified as one of the following:

- Character vector — For single-output models. For example, 'measurements'.
- Cell array of character vectors — For multi-output models.

Alternatively, use automatic vector expansion to assign output names for multi-output models. For example, if `sys` is a two-output model, enter:

```
sys.OutputName = 'measurements';
```

The output names automatically expand to `{'measurements(1)'; 'measurements(2)'}.`

When you estimate a model using an `iddata` object, `data`, the software automatically sets `OutputName` to `data.OutputName`.

You can use the shorthand notation `y` to refer to the `OutputName` property. For example, `sys.y` is equivalent to `sys.OutputName`.

Output channel names have several uses, including:

- Identifying channels on model display and plots
- Extracting subsystems of MIMO systems
- Specifying connection points when interconnecting models

**Default:** `''` for all output channels

### **OutputUnit**

Output channel units, specified as one of the following:

- Character vector — For single-output models. For example, `'seconds'`.
- Cell array of character vectors — For multi-output models.

Use `OutputUnit` to keep track of output signal units. `OutputUnit` has no effect on system behavior.

**Default:** `''` for all output channels

### **OutputGroup**

Output channel groups. The `OutputGroup` property lets you assign the output channels of MIMO systems into groups and refer to each group by name. Specify output groups as a structure. In this structure, field names are the group names, and field values are the output channels belonging to each group. For example:

```
sys.OutputGroup.temperature = [1];  
sys.InputGroup.measurement = [3 5];
```

creates output groups named `temperature` and `measurement` that include output channels 1, and 3, 5, respectively. You can then extract the subsystem from all inputs to the `measurement` outputs using:

```
sys('measurement',:)
```

**Default:** Struct with no fields

### Name

System name, specified as a character vector. For example, 'system\_1'.

**Default:** ''

### Notes

Any text that you want to associate with the system, stored as a string or a cell array of character vectors. The property stores whichever data type you provide. For instance, if `sys1` and `sys2` are dynamic system models, you can set their `Notes` properties as follows:

```
sys1.Notes = "sys1 has a string.";
sys2.Notes = 'sys2 has a character vector.';
sys1.Notes
sys2.Notes
```

```
ans =
```

```
    "sys1 has a string."
```

```
ans =
```

```
    'sys2 has a character vector.'
```

**Default:** [0×1 string]

### UserData

Any type of data you want to associate with system, specified as any MATLAB data type.

**Default:** []

## Output Arguments

### sys — Nonlinear ARX model

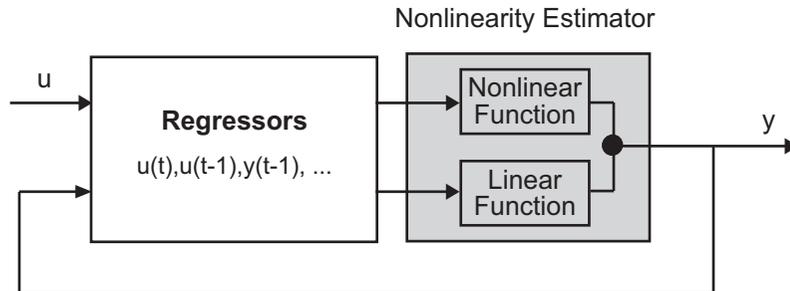
idnlarx object

Nonlinear ARX model, returned as an `idnlarx` object. This model is created using the specified model orders, nonlinearity estimator, and properties.

## Definitions

### Nonlinear ARX Model Structure

A nonlinear ARX model consists of model regressors and a nonlinearity estimator. The nonlinearity estimator comprises both linear and nonlinear functions that act on the model regressors to give the model output. This block diagram represents the structure of a nonlinear ARX model in a simulation scenario.



The software computes the nonlinear ARX model output  $y$  in two stages:

- 1 It computes regressor values from the current and past input values and past output data.

In the simplest case, regressors are delayed inputs and outputs, such as  $u(t-1)$  and  $y(t-3)$ . These kind of regressors are called *standard regressors*. You specify the standard regressors using the model orders and delay. For more information, see “Nonlinear ARX Model Orders and Delay”. You can also specify *custom* regressors, which are nonlinear functions of delayed inputs and outputs. For example,  $u(t-1)*y(t-3)$ . To create a set of polynomial type regressors, use `polyreg`.

By default, all regressors are inputs to both the linear and the nonlinear function blocks of the nonlinearity estimator. You can choose a subset of regressors as inputs to the nonlinear function block.

- 2 It maps the regressors to the model output using the nonlinearity estimator block. The nonlinearity estimator block can include linear and nonlinear blocks in parallel. For example:

$$F(x) = L^T(x - r) + d + g(Q(x - r))$$

Here,  $x$  is a vector of the regressors, and  $r$  is the mean of the regressors  $x$ .  $L^T(x) + d$  is the output of the linear function block and is affine when  $d \neq 0$ .  $d$  is a scalar offset.

$g(Q(x - r))$  represents the output of the nonlinear function block.  $Q$  is a projection matrix that makes the calculations well conditioned. The exact form of  $F(x)$  depends on your choice of the nonlinearity estimator. You can select from available nonlinearity estimators, such as tree-partition networks, wavelet networks, and multilayer neural networks. You can also exclude either the linear or the nonlinear function block from the nonlinearity estimator.

When estimating a nonlinear ARX model, the software computes the model parameter values, such as  $L$ ,  $r$ ,  $d$ ,  $Q$ , and other parameters specifying  $g$ .

Resulting nonlinear ARX models are `idnlarx` objects that store all model data, including model regressors and parameters of the nonlinearity estimator. For more information about these objects, see “Nonlinear Model Structures”.

## Definition of `idnlarx` States

The states of an `idnlarx` object are an ordered list of delayed input and output variables that define the structure of the model. The toolbox uses this definition of states for creating the initial state vector that is used during simulation and prediction with `sim`, `predict`, and `compare`. This definition is also used for linearization of nonlinear ARX models using `linearize`.

This toolbox provides several options to facilitate how you specify the initial states. For example, you can use `findstates` and `data2state` to search for state values in simulation and prediction applications. For linearization, use `findop`. You can also specify the states manually.

The states of an `idnlarx` model depend on the maximum delay in each input and output variable used by the regressors. If a variable  $p$  has a maximum delay of  $D$  samples, then it contributes  $D$  elements to the state vector at time  $t$ :  $p(t-1)$ ,  $p(t-2)$ , ...,  $p(t-D)$ .

For example, if you have a single-input, single-output `idnlarx` model.

```
m = idnlarx([2 3 0], 'wavenet', 'CustomRegressors', {'y1(t-10)*u1(t-1)'});
```

This model has these regressors.

```
getreg(m)
```

```
Regressors:
```

```
  y1(t-1)
  y1(t-2)
  u1(t)
  u1(t-1)
  u1(t-2)
  y1(t-10)*u1(t-1)
```

The regressors show that the maximum delay in the output variable `y1` is 10 samples and the maximum delay in the input `u1` is two samples. Thus, this model has a total of 12 states:

$$X(t) = [y1(t-1), y2(t-2), \dots, y1(t-10), u1(t-1), u1(t-2)]$$

---

**Note** The state vector includes the output variables first, followed by input variables.

---

As another example, consider the 2-output and 3-input model.

```
m = idnlarx([2 0 2 2 1 1 0 0; 1 0 1 5 0 1 1 0],[wavenet; linear]);
```

This model has these regressors.

```
getreg(m)
```

```
Regressors:
```

```
  For output 1:
  y1(t-1)
  y1(t-2)
  u1(t-1)
  u1(t-2)
  u2(t)
```

```

    u2(t-1)
    u3(t)
For output 2:
    y1(t-1)
    u1(t-1)
    u2(t-1)
    u2(t-2)
    u2(t-3)
    u2(t-4)
    u2(t-5)

```

The maximum delay in output variable  $y_1$  is two samples. This delay occurs in the regressor set for output 1. The maximum delays in the three input variables are 2, 5, and 0, respectively. Thus, the state vector is:

$$X(t) = [y_1(t-1), y_1(t-2), u_1(t-1), u_1(t-2), u_2(t-1), u_2(t-2), u_2(t-3), u_2(t-4), u_2(t-5)]$$

Variables  $y_2$  and  $u_3$  do not contribute to the state vector because the maximum delay in these variables is zero.

A simpler way to determine states by inspecting regressors is to use `getDelayInfo`, which returns the maximum delays in all I/O variables across all model outputs. For the multiple-input multiple-output model  $m$ , `getDelayInfo` returns:

```
maxDel = getDelayInfo(m)
```

```
maxDel = 1x5
```

```
     2     0     2     5     0
```

`maxDel` contains the maximum delays for all input and output variables in the order ( $y_1$ ,  $y_2$ ,  $u_1$ ,  $u_2$ ,  $u_3$ ). The total number of model states is `sum(maxDel) = 9`.

The set of states for an `idnlarx` model is not required to be minimal.

## See Also

`addreg` | `customnet` | `customreg` | `getreg` | `idnlarx/findop` | `linear` | `linearize` | `nlarx` | `pem` | `polyreg` | `sigmoidnet` | `wavenet`

## **Topics**

“Use nlarx to Estimate Nonlinear ARX Models”

“Estimate Nonlinear ARX Models Initialized Using Linear ARX Models”

“Initialize Nonlinear ARX Estimation Using Linear Model”

“Identifying Nonlinear ARX Models”

“Nonlinear Model Structures”

## **Introduced in R2007a**

# idnlgrey

Nonlinear grey-box model

## Syntax

```
sys = idnlgrey(FileName,Order,Parameters)
sys = idnlgrey(FileName,Order,Parameters,InitialStates)
sys = idnlgrey(FileName,Order,Parameters,InitialStates,Ts)
sys = idnlgrey(FileName,Order,Parameters,InitialStates,Ts,
Name,Value)
```

## Description

`sys = idnlgrey(FileName,Order,Parameters)` creates a nonlinear grey-box model using the specified model structure in `FileName`, number of outputs, inputs, and states in `Order`, and the model parameters.

`sys = idnlgrey(FileName,Order,Parameters,InitialStates)` specifies the initial states of the model.

`sys = idnlgrey(FileName,Order,Parameters,InitialStates,Ts)` specifies the sample time of a discrete-time model.

`sys = idnlgrey(FileName,Order,Parameters,InitialStates,Ts,Name,Value)` specifies additional attributes of the `idnlgrey` model structure using one or more `Name,Value` pair arguments.

## Object Description

`idnlgrey` represents a nonlinear grey-box model. For information about the nonlinear grey-box models, see “Estimate Nonlinear Grey-Box Models”.

Use the `idnlgrey` constructor to create the nonlinear grey-box model and then estimate the model parameters using `nlgreyest`.

For `idnlgrey` object properties, see “Properties” on page 1-642.

## Examples

### Create a Nonlinear Grey-Box Model

Load data.

```
load(fullfile(matlabroot, 'toolbox', 'ident', 'iddemos', 'data', 'dcmotordata'));  
z = iddata(y,u,0.1, 'Name', 'DC-motor');
```

The data is from a linear DC motor with one input (voltage), and two outputs (angular position and angular velocity). The structure of the model is specified by `dcmotor_m.m` file.

Create a nonlinear grey-box model.

```
file_name = 'dcmotor_m';  
Order = [2 1 2];  
Parameters = [1;0.28];  
InitialStates = [0;0];  
  
sys = idnlgrey(file_name,Order,Parameters,InitialStates,0, ...  
    'Name', 'DC-motor');
```

### Selectively Estimate Parameters of Nonlinear Grey-Box Model

Load data.

```
load(fullfile(matlabroot, 'toolbox', 'ident', 'iddemos', 'data', 'twotankdata'));  
z = iddata(y,u,0.2, 'Name', 'Two tanks');
```

The data contains 3000 input-output data samples of a two tank system. The input is the voltage applied to a pump, and the output is the liquid level of the lower tank.

Specify file describing the model structure for a two-tank system. The file specifies the state derivatives and model outputs as a function of time, states, inputs, and model parameters.

```
FileName = 'twotanks_c';  
Specify model orders [ny nu nx].  
Order = [1 1 2];  
Specify initial parameters (Np = 6).  
Parameters = {0.5;0.0035;0.019; ...  
             9.81;0.25;0.016};  
Specify initial initial states.  
InitialStates = [0;0.1];  
Specify as continuous system.  
Ts = 0;  
Create idnlgrey model object.  
nlgr = idnlgrey(FileName,Order,Parameters,InitialStates,Ts, ...  
              'Name','Two tanks');  
Set some parameters as constant.  
nlgr.Parameters(1).Fixed = true;  
nlgr.Parameters(4).Fixed = true;  
nlgr.Parameters(5).Fixed = true;  
Estimate the model parameters.  
nlgr = nlgreyest(z,nlgr);
```

## Input Arguments

**FileName** — Name of the function or MEX-file that stores the model structure  
character vector | function handle

Name of the function or MEX-file storing the model structure, specified as a character vector (without the file extension) or a function handle for computing the states and the outputs. If `FileName` is a character vector, for example `'twotanks_c'`, then it must

point to a MATLAB file, P-code file, or MEX-file. For more information about the file variables, see “Specifying the Nonlinear Grey-Box Model Structure”.

**Order — Number of outputs, inputs, and states of the model**

vector | structure

Number of outputs, inputs, and states of the model, specified as one of the following:

- Vector  $[N_y \ N_u \ N_x]$ , specifying the number of model outputs  $N_y$ , inputs  $N_u$ , and states  $N_x$ .
- Structure with fields 'Ny', 'Nu', and 'Nx'.

For time series,  $N_u$  is set to 0, and for static model structures,  $N_x$  is set to 0.

**Parameters — Parameters of the model**

structure | vector | cell array

Parameters of the model, specified as one of the following:

- $N_p$ -by-1 structure array, where  $N_p$  is the number of parameters. The structure contains the following fields:

Field	Description	Default
Name	Name of the parameter, specified as a character vector. For example, 'pressure'.	'pi', where i is an integer in $[1, N_p]$
Unit	Unit of the parameter, specified as a character vector.	''
Value	Initial value of the parameter, specified as: <ul style="list-style-type: none"> <li>• Finite real scalar</li> <li>• Finite real column vector</li> <li>• Two-dimensional real matrix</li> </ul>	

Field	Description	Default
Minimum	Minimum value of the parameter, specified as a real scalar, column vector, or matrix of the same size as <code>Value</code> .  Minimum $\geq$ Value for all components.	<code>-Inf(size(Value))</code>
Maximum	Maximum value of the parameter, specified as a real scalar, column vector, or matrix of the same size as <code>Value</code> .  Value $\leq$ Maximum for all components.	<code>Inf(size(Value))</code>
Fixed	Specifies whether parameter is fixed to their initial values, specified as a boolean scalar, column vector, or matrix of the same size as <code>Value</code> .	<code>false(size(Value))</code>  Implies, estimate all parameters

Use dot notation to access the subfields of the  $i$ th parameter. For example, for `idnlgrey` model `M`, the  $i$ th parameter is accessed through `M.Parameters(i)` and its subfield `Fixed` by `M.Parameters(i).Fixed`.

- $N_p$ -by-1 vector of real finite initial values, `InParameters`.

The data is converted into a structure with default values for the fields `Name`, `Unit`, `Minimum`, `Maximum`, and `Fixed`.

`Value` is assigned the value `InParameters(i)`, where  $i$  is an integer in  $[1, N_p]$

- $N_p$ -by-1 cell array containing finite real scalars, finite real vectors, or finite real two-dimensional matrices of initial values.

Default values are used for the fields `Name`, `Unit`, `Minimum`, `Maximum`, and `Fixed`.

### **InitialStates – Initial states of the model**

structure | [] | cell array | {}

Initial states of the model parameters specified as one of the following:

- $N_x$ -by-1 structure array, where  $N_x$  is the number of states. The structure contains the following fields:

Field	Description	Default
Name	Name of the states, specified as a character vector.	'xi', where i is an integer in [1, $N_x$ ]
Unit	Unit of the states, specified as a character vector.	' '
Value	Initial value of the initial states, specified as: <ul style="list-style-type: none"> <li>• A finite real scalar</li> <li>• A finite real 1-by-<math>N_e</math> vector, where <math>N_e</math> is the number of experiments in the data set to be used for estimation</li> </ul>	
Minimum	Minimum value of the initial states, specified as a real scalar or 1-by- $N_e$ vector of the same size as Value.  Minimum $\geq$ Value for all components.	-Inf(size(Value))
Maximum	Maximum value of the parameters, specified as a real scalar or 1-by- $N_e$ vector of the same size as Value.  Value $\leq$ Maximum for all components.	Inf(size(Value))

Field	Description	Default
Fixed	Specifies whether initial states are fixed to their initial values, specified as boolean scalar or 1-by-Ne vector of the same size as Value	true(size(Value)) Implies, do not estimate the initial states.

Use dot notation to access the subfields of the *i*th initial state. For example, for idnlgrey model *M*, the *i*th initial state is accessed through `M.InitialStates(i)` and its subfield `Fixed` by `M.InitialStates(i).Fixed`.

- `[]`.

A structure is created with default values for the fields `Name`, `Unit`, `Minimum`, `Maximum`, and `Fixed`.

`Value` is assigned the value `0`.

- A real finite *N<sub>x</sub>*-by-*N<sub>e</sub>* matrix (`InitStates`).

`Value` of the *i*th structure array element is `InitStates(i,Ne)`, a row vector with *N<sub>e</sub>* elements. `Minimum`, `Maximum`, and `Fixed` will be `-Inf`, `Inf` and `true` row vectors of the same size as `InitStates(i,Ne)`.

- Cell array with finite real vectors of size 1-by-*N<sub>e</sub>* or `{}` (same as `[]`).

### Ts — Sample time

0 (default) | scalar

Sample time, specified as a positive scalar representing the sampling period. The value is expressed in the unit specified by the `TimeUnit` property of the model. For a continuous time model `Ts` is equal to 0 (default).

### Name-Value Pair Arguments

Specify optional comma-separated pairs of `Name`, `Value` arguments. `Name` is the argument name and `Value` is the corresponding value. `Name` must appear inside quotes. You can specify several name and value pair arguments in any order as `Name1,Value1,...,NameN,ValueN`.

Use `Name`, `Value` arguments to specify additional properties on page 1-642 of idnlgrey models during model creation.

## Properties

idnlgrey object properties include:

### FileName

Name of the function or MEX-file storing the model structure, specified as a character vector (without extension) or a function handle for computing the states and the outputs. If `FileName` is a character vector, for example `'twotanks_c'`, then it must point to a MATLAB file, P-code file, or MEX-file. For more information about the file variables, see “Specifying the Nonlinear Grey-Box Model Structure”.

### Order

Number of outputs, inputs, and states of the model, specified as one of the following:

- Vector `[Ny Nu Nx]`, specifying the number of model outputs `Ny`, inputs `Nu`, and states `Nx`.
- Structure with fields `'Ny'`, `'Nu'`, and `'Nx'`.

For time series, `Nu` is set to `0`, and for static model structures, `Nx` is set to `0`.

### Parameters

Parameters of the model, specified as one of the following:

- `Np`-by-1 structure array, where `Np` is the number of parameters. The structure contains the following fields:

Field	Description	Default
Name	Name of the parameter, specified as a character vector. For example, <code>'pressure'</code> .	<code>'pi'</code> , where <code>i</code> is an integer in <code>[1,Np]</code>
Unit	Unit of the parameter, specified as a character vector.	<code>''</code>

Field	Description	Default
Value	Initial value of the parameter, specified as: <ul style="list-style-type: none"> <li>• Finite real scalar</li> <li>• Finite real column vector</li> <li>• Two-dimensional real matrix</li> </ul>	
Minimum	Minimum value of the parameter, specified as a real scalar, column vector, or matrix of the same size as Value.  Minimum $\geq$ Value for all components.	<code>-Inf(size(Value))</code>
Maximum	Maximum value of the parameter, specified as a real scalar, column vector, or matrix of the same size as Value.  Value $\leq$ Maximum for all components.	<code>Inf(size(Value))</code>
Fixed	Specifies whether parameter is fixed to their initial values, specified as a boolean scalar, column vector, or matrix of the same size as Value.	<code>false(size(Value))</code>  Implies, estimate all parameters

- $N_p$ -by-1 vector of real finite initial values, `InParameters`.

The data is converted into a structure with default values for the fields `Name`, `Unit`, `Minimum`, `Maximum`, and `Fixed`.

`Value` is assigned the value `InParameters(i)`, where `i` is an integer in `[1,  $N_p$ ]`

- $N_p$ -by-1 cell array containing finite real scalars, finite real vectors, or finite real two-dimensional matrices of initial values.

A structure is created with default values for the fields **Name**, **Unit**, **Minimum**, **Maximum**, and **Fixed**.

Use dot notation to access the subfields of the  $i$ th parameter. For example, for `idnlgrey` model  $M$ , the  $i$ th parameter is accessed through `M.Parameters(i)` and its subfield `Fixed` by `M.Parameters(i).Fixed`.

### **InitialStates**

Initial states of the model parameters specified as one of the following:

- $N_x$ -by-1 structure array, where  $N_x$  is the number of states. The structure contains the following fields:

<b>Field</b>	<b>Description</b>	<b>Default</b>
<b>Name</b>	Name of the states, specified as a character vector.	'xi', where $i$ is an integer in $[1, N_x]$
<b>Unit</b>	Unit of the states, specified as a character vector.	''
<b>Value</b>	Initial value of the initial states, specified as: <ul style="list-style-type: none"> <li>• A finite real scalar</li> <li>• A finite real 1-by-<math>N_e</math> vector, where <math>N_e</math> is the number of experiments in the data set to be used for estimation</li> </ul>	

Field	Description	Default
Minimum	Minimum value of the initial states, specified as a real scalar or 1-by-Ne vector of the same size as Value.  Minimum $\geq$ Value for all components.	-Inf(size(Value))
Maximum	Maximum value of the parameters, specified as a real scalar or 1-by-Ne vector of the same size as Value.  Value $\leq$ Maximum for all components.	Inf(size(Value))
Fixed	Specifies whether initial states are fixed to their initial values, specified as boolean scalar or 1-by-Ne vector of the same size as Value	true(size(Value))  Implies, do not estimate the initial states.

- `[]`.

A structure is created with default values for the fields `Name`, `Unit`, `Minimum`, `Maximum`, and `Fixed`.

`Value` is assigned the value `0`.

- A real finite  $N_x$ -by- $N_e$  matrix (`InitStates`).

Value of the  $i$ th structure array element is `InitStates(i,Ne)`, a row vector with  $N_e$  elements. `Minimum`, `Maximum`, and `Fixed` will be `-Inf`, `Inf` and `true` row vectors of the same size as `InitStates(i,Ne)`.

- Cell array with finite real vectors of size 1-by- $N_e$  or `{}` (same as `[]`).

A structure is created with default values for the fields `Name`, `Unit`, `Minimum`, `Maximum`, and `Fixed`.

Use dot notation to access the subfields of the *i*th initial state. For example, for `idnlgrey` model *M*, the *i*th initial state is accessed through `M.InitialStates(i)` and its subfield `Fixed` by `M.InitialStates(i).Fixed`.

### **FileArgument**

Contains auxiliary variables passed to the ODE file (function or MEX-file) specified in `FileName`, specified as a cell array. These variables are used as extra inputs for specifying the state and/or output equations.

Default: `{}`.

### **SimulationOptions**

A structure that specifies the simulation method and related options, containing the following fields:

<b>Field</b>	<b>Description</b>	<b>Default</b>
<code>AbsTol</code>	Absolute error tolerance. This scalar applies to all components of the state vector.  Applicable to: Variable step solvers.  Assignable value: A positive real value.	<code>1e-6</code>
<code>FixedStep</code>	Step size used by the solver.  Applicable to: Fixed-step time-continuous solvers.  Assignable values: <ul style="list-style-type: none"><li>• <code>'Auto'</code> — Automatically chooses the initial step.</li><li>• A real value such that <math>0 &lt; \text{FixedStep} \leq 1</math>.</li></ul>	<code>'Auto'</code>  Automatically chooses the initial step.

Field	Description	Default
InitialStep	<p>Specifies the initial step at which the ODE solver starts.</p> <p>Applicable to: Variable-step, time-continuous solvers.</p> <p>Assignable values:</p> <ul style="list-style-type: none"> <li>• 'Auto' — Automatically chooses the initial step.</li> <li>• A positive real value such that <math>\text{MinStep} \leq \text{InitialStep} \leq \text{MaxStep}</math>.</li> </ul>	<p>'Auto'</p> <p>Automatically chooses the initial step.</p>
MaxOrder	<p>Specifies the order of the Numerical Differentiation Formulas (NDF).</p> <p>Applicable to: ode15s.</p> <p>Assignable values: 1, 2, 3, 4 or 5.</p>	5
MaxStep	<p>Specifies the largest time step of the ODE solver.</p> <p>Applicable to: Variable-step, time-continuous solvers.</p> <p>Assignable values:</p> <ul style="list-style-type: none"> <li>• 'Auto' — Automatically chooses the time step.</li> <li>• A positive real value <math>&gt; \text{MinStep}</math>.</li> </ul>	<p>'Auto'</p> <p>Automatically chooses the time step.</p>

<b>Field</b>	<b>Description</b>	<b>Default</b>
MinStep	<p>Specifies the smallest time step of the ODE solver.</p> <p>Applicable to: Variable-step, time-continuous solvers.</p> <p>Assignable values:</p> <ul style="list-style-type: none"><li>• 'Auto' — Automatically chooses the time step.</li><li>• A positive real value &lt; MaxStep.</li></ul>	<p>'Auto'</p> <p>Automatically chooses the time step.</p>
RelTol	<p>Relative error tolerance that applies to all components of the state vector. The estimated error in each integration step satisfies <math> e(i)  \leq \max(\text{RelTol} * \text{abs}(x(i)), \text{AbsTol}(i))</math>.</p> <p>Applicable to: Variable-step, time-continuous solvers.</p> <p>Assignable value: A positive real value.</p>	<p>1e-3</p> <p>(0.1% accuracy).</p>

Field	Description	Default
Solver	<p>ODE (Ordinary Differential/Difference Equation) solver for solving state space equations.</p> <ul style="list-style-type: none"> <li>• Variable-step solvers for time-continuous idnlgrey models: <ul style="list-style-type: none"> <li>• 'ode45' — Runge-Kutta (4,5) solver for nonstiff problems.</li> <li>• 'ode23' — Runge-Kutta (2,3) solver for nonstiff problems.</li> <li>• 'ode113' — Adams-Bashforth-Moulton solver for nonstiff problems.</li> <li>• 'ode15s' — Numerical Differential Formula solver for stiff problems.</li> <li>• 'ode23s' — Modified Rosenbrock solver for stiff problems.</li> <li>• 'ode23t' — Trapezoidal solver for moderately stiff problems.</li> <li>• 'ode23tb' — Implicit Runge-Kutta solver for stiff problems.</li> </ul> </li> <li>• Fixed-step solvers for time-continuous idnlgrey models: <ul style="list-style-type: none"> <li>• 'ode5' — Dormand-Prince solver.</li> <li>• 'ode4' — Fourth-order Runge-Kutta solver.</li> <li>• 'ode3' — Bogacki-Shampine solver.</li> <li>• 'ode2' — Heun or improved Euler solver.</li> <li>• 'ode1' — Euler solver.</li> </ul> </li> <li>• Fixed-step solvers for time-discrete idnlgrey models: 'FixedStepDiscrete'</li> <li>• General: 'Auto' — Automatically chooses one of the previous solvers.</li> </ul>	<p>'Auto'</p> <p>Automatically chooses one of the solvers.</p>

## Report

Summary report that contains information about the estimation options and results when the model is estimated using the `nlgreyest` command. Use `Report` to query a model for how it was estimated, including:

- Estimation method
- Estimation options
- Search termination conditions
- Estimation data fit

The contents of `Report` are irrelevant if the model was created by construction.

```
nlgr = idnlgrey('dcmotor_m',[2,1,2],[1;0.28],[0;0],0,'Name','DC-motor');  
nlgr.Report.OptionsUsed
```

```
ans =
```

```
 []
```

If you use `nlgreyest` to estimate the model, the fields of `Report` contain information on the estimation data, options, and results.

```
load(fullfile(matlabroot,'toolbox','ident','iddemos','data','dcmotordata'));  
z = iddata(y,u,0.1,'Name','DC-motor');  
nlgr = idnlgrey('dcmotor_m',[2,1,2],[1;0.28],[0;0],0,'Name','DC-motor');  
nlgr = nlgreyest(z,nlgr);  
nlgr.Report.OptionsUsed
```

Option set for the `nlgreyest` command:

```
GradientOptions: [1x1 struct]  
EstimateCovariance: 1  
    Display: 'off'  
Regularization: [1x1 struct]  
  SearchMethod: 'auto'  
SearchOptions: [1x1 idoptions.search.lsqnonlin]  
  OutputWeight: []  
    Advanced: [1x1 struct]
```

`Report` is a read-only property.

For more information on this property and how to use it, see “Output Arguments” on page 1-979 in the `nlgreyest` reference page and “Estimation Report”.

**TimeVariable**

Independent variable for the inputs, outputs, and—when available—internal states, specified as a character vector.

**Default:** 't'

**NoiseVariance**

Noise variance (covariance matrix) of the model innovations  $e$ . Assignable value is an  $n_y$ -by- $n_y$  matrix. Typically set automatically by the estimation algorithm.

**Ts**

Sample time.  $T_s$  is a positive scalar representing the sampling period. This value is expressed in the unit specified by the `TimeUnit` property of the model. For a continuous time model,  $T_s$  is equal to 0 (default).

Changing this property does not discretize or resample the model.

**Default:** 0

**TimeUnit**

Units for the time variable, the sample time  $T_s$ , and any time delays in the model, specified as one of the following values:

- 'nanoseconds'
- 'microseconds'
- 'milliseconds'
- 'seconds'
- 'minutes'
- 'hours'
- 'days'
- 'weeks'
- 'months'
- 'years'

Changing this property has no effect on other properties, and therefore changes the overall system behavior. Use `chgTimeUnit` to convert between time units without modifying system behavior.

**Default:** `'seconds'`

### **InputName**

Input channel names, specified as one of the following:

- Character vector — For single-input models, for example, `'controls'`.
- Cell array of character vectors — For multi-input models.

Alternatively, use automatic vector expansion to assign input names for multi-input models. For example, if `sys` is a two-input model, enter:

```
sys.InputName = 'controls';
```

The input names automatically expand to `{'controls(1)'; 'controls(2)'}`.

When you estimate a model using an `iddata` object, `data`, the software automatically sets `InputName` to `data.InputName`.

You can use the shorthand notation `u` to refer to the `InputName` property. For example, `sys.u` is equivalent to `sys.InputName`.

Input channel names have several uses, including:

- Identifying channels on model display and plots
- Extracting subsystems of MIMO systems
- Specifying connection points when interconnecting models

**Default:** `''` for all input channels

### **InputUnit**

Input channel units, specified as one of the following:

- Character vector — For single-input models, for example, `'seconds'`.
- Cell array of character vectors — For multi-input models.

Use `InputUnit` to keep track of input signal units. `InputUnit` has no effect on system behavior.

**Default:** '' for all input channels

### InputGroup

Input channel groups. The `InputGroup` property lets you assign the input channels of MIMO systems into groups and refer to each group by name. Specify input groups as a structure. In this structure, field names are the group names, and field values are the input channels belonging to each group. For example:

```
sys.InputGroup.controls = [1 2];  
sys.InputGroup.noise = [3 5];
```

creates input groups named `controls` and `noise` that include input channels 1, 2 and 3, 5, respectively. You can then extract the subsystem from the `controls` inputs to all outputs using:

```
sys(:, 'controls')
```

**Default:** Struct with no fields

### OutputName

Output channel names, specified as one of the following:

- Character vector — For single-output models. For example, 'measurements'.
- Cell array of character vectors — For multi-output models.

Alternatively, use automatic vector expansion to assign output names for multi-output models. For example, if `sys` is a two-output model, enter:

```
sys.OutputName = 'measurements';
```

The output names automatically expand to `{'measurements(1)'; 'measurements(2)'}`.

When you estimate a model using an `iddata` object, `data`, the software automatically sets `OutputName` to `data.OutputName`.

You can use the shorthand notation `y` to refer to the `OutputName` property. For example, `sys.y` is equivalent to `sys.OutputName`.

Output channel names have several uses, including:

- Identifying channels on model display and plots
- Extracting subsystems of MIMO systems
- Specifying connection points when interconnecting models

**Default:** '' for all output channels

### **OutputUnit**

Output channel units, specified as one of the following:

- Character vector — For single-output models. For example, 'seconds'.
- Cell array of character vectors — For multi-output models.

Use `OutputUnit` to keep track of output signal units. `OutputUnit` has no effect on system behavior.

**Default:** '' for all output channels

### **OutputGroup**

Output channel groups. The `OutputGroup` property lets you assign the output channels of MIMO systems into groups and refer to each group by name. Specify output groups as a structure. In this structure, field names are the group names, and field values are the output channels belonging to each group. For example:

```
sys.OutputGroup.temperature = [1];  
sys.InputGroup.measurement = [3 5];
```

creates output groups named `temperature` and `measurement` that include output channels 1, and 3, 5, respectively. You can then extract the subsystem from all inputs to the `measurement` outputs using:

```
sys('measurement', :)
```

**Default:** Struct with no fields

### **Name**

System name, specified as a character vector. For example, 'system\_1'.

**Default:** ''

## Notes

Any text that you want to associate with the system, stored as a string or a cell array of character vectors. The property stores whichever data type you provide. For instance, if `sys1` and `sys2` are dynamic system models, you can set their `Notes` properties as follows:

```
sys1.Notes = "sys1 has a string.";
sys2.Notes = 'sys2 has a character vector.';
sys1.Notes
sys2.Notes
```

```
ans =
```

```
    "sys1 has a string."
```

```
ans =
```

```
    'sys2 has a character vector.'
```

**Default:** [0×1 string]

## UserData

Any type of data you want to associate with system, specified as any MATLAB data type.

**Default:** []

## Output Arguments

**sys** — Nonlinear grey-box model

idnlgrey object

Nonlinear grey-box model, returned as an idnlgrey object.

## Definitions

### Definition of idnlgrey States

The states of an idnlgrey model are defined explicitly in the function or MEX-file storing the model structure. States are required for simulation and prediction of nonlinear grey-box models. Use `findstates` to search for state values for simulation and prediction with `sim`, `predict`, and `compare`.

---

**Note** The initial values of the states are configured by the `InitialStates` property of the idnlgrey model.

---

### See Also

`get` | `getinit` | `getpar` | `nlgreyest` | `pem` | `set` | `setinit` | `setpar`

### Topics

“Represent Nonlinear Dynamics Using MATLAB File for Grey-Box Estimation”

“Creating IDNLGREY Model Files”

“Estimate Nonlinear Grey-Box Models”

**Introduced in R2007a**

# idnlhw

Hammerstein-Wiener model

## Syntax

```
sys = idnlhw(Orders)
sys = idnlhw(Orders,InputNL,OutputNL)
sys = idnlhw(Orders,InputNL,OutputNL,Name,Value)
sys = idnlhw(LinModel)
sys = idnlhw(LinModel,InputNL,OutputNL)
sys = idnlhw(LinModel,InputNL,OutputNL,Name,Value)
```

## Description

`sys = idnlhw(Orders)` creates a Hammerstein-Wiener model with the specified orders, and using piecewise linear functions as input and output nonlinearity estimators.

`sys = idnlhw(Orders,InputNL,OutputNL)` uses `InputNL` and `OutputNL` as the input and output nonlinearity estimators, respectively.

`sys = idnlhw(Orders,InputNL,OutputNL,Name,Value)` specifies additional attributes of the `idnlhw` model structure using one or more `Name,Value` pair arguments.

`sys = idnlhw(LinModel)` uses a linear model `LinModel` to specify the model orders and default piecewise linear functions for the input and output nonlinearity estimators.

`sys = idnlhw(LinModel,InputNL,OutputNL)` specifies input and output nonlinearity estimators for the model.

`sys = idnlhw(LinModel,InputNL,OutputNL,Name,Value)` specifies additional attributes of the `idnlhw` model structure using one or more `Name,Value` pair arguments.

## Object Description

`idnlhw` represents a Hammerstein-Wiener model. The Hammerstein-Wiener structure on page 1-671 represents a linear model with input-output nonlinearities.

Use the `nlhw` command to both construct an `idnlhw` object and estimate the model parameters.

You can also use the `idnlhw` constructor to create the Hammerstein-Wiener model and then estimate the model parameters using `nlhw`.

For `idnlhw` object properties, see “Properties” on page 1-662.

## Examples

### Create a Hammerstein-Wiener Model Structure with Default Nonlinearities

Create a Hammerstein-Wiener model with `nb` and `nf` = 2 and `nk` = 1.

```
m = idnlhw([2 2 1]);
```

`m` has piecewise linear input and output nonlinearity.

### Create Hammerstein-Wiener Model with Specific Input-Output Nonlinearities

```
m = idnlhw([2 2 1], 'sigmoidnet', 'deadzone');
```

The above is equivalent to:

```
m = idnlhw([2 2 1], 'sig', 'dead');
```

The specified nonlinearities have a default configuration.

### Create Hammerstein-Wiener Model and Configure the Nonlinearities

```
m = idnlhw([2 2 1], sigmoidnet('num',5), deadzone([-1,2]), 'InputName', 'Volts', 'OutputName');
```

## Create a Wiener Model and Estimate Model Parameters

Create a Wiener model (no input nonlinearity).

```
m = idnlhw([2 2 1],[], 'saturation');
```

Estimate the model.

```
load iddata1;
m = nllhw(z1,m);
```

## Create Hammerstein-Wiener Model Using Input-Output Polynomial Model of Output-Error Structure

Construct an input-output polynomial model of OE structure.

```
B = [0.8 1];
F = [1 -1.2 0.5];
LinearModel = idpoly(1,B,1,1,F, 'Ts',0.1);
```

Construct Hammerstein-Wiener model using OE model as its linear component.

```
m1 = idnlhw(LinearModel, 'saturation', [], 'InputName', 'Control');
```

## Input Arguments

### Orders — Order and delays of the linear subsystem transfer function

[nb nf nk] vector of positive integers | [nb nf nk] vector of matrices

Order and delays of the linear subsystem transfer function, specified as a [nb nf nk] vector.

Dimensions of Orders:

- For a SISO transfer function, Orders is a vector of positive integers.
  - nb is the number of zeros plus 1, nf is the number of poles, and nk is the input delay.

- For a MIMO transfer function with  $n_u$  inputs and  $n_y$  outputs, `Orders` is a vector of matrices.

`nb`, `nf`, and `nk` are  $n_y$ -by- $n_u$  matrices whose  $i$ - $j$ th entry specifies the orders and delay of the transfer function from the  $j$ th input to the  $i$ th output.

### **InputNL — Input static nonlinearity**

'pwlinear' (default) | 'sigmoidnet' | 'wavenet' | 'saturation' | 'deadzone' | 'poly1d' | 'unitgain' | nonlinearity estimator object | array of nonlinearity estimators

Input static nonlinearity estimator, specified as one of the following.

'pwlinear' or pwlinear object (default)	Piecewise linear function
'sigmoidnet' or sigmoidnet object	Sigmoid network
'wavenet' or wavenet object	Wavelet network
'saturation' or saturation object	Saturation
'deadzone' or deadzone object	Dead zone
'poly1d' or poly1d object	One-dimensional polynomial
'unitgain' or [] or unitgain object	Unit gain
customnet object	Custom network — Similar to <code>sigmoidnet</code> , but with a user-defined replacement for the sigmoid function.

Specifying a character vector, for example `'sigmoidnet'`, creates a nonlinearity estimator object with default settings. Use object representation instead to configure the properties of a nonlinearity estimator.

```
InputNL = wavenet;  
InputNL.NumberOfUnits = 10;
```

Alternatively, use the associated input nonlinearity estimator function with Name-Value pair arguments.

```
InputNL = wavenet('NumberOfUnits',10);
```

For  $n_u$  input channels, you can specify nonlinear estimators individually for each input channel by setting `InputNL` to an  $n_u$ -by-1 array of nonlinearity estimators.

```
InputNL = [sigmoidnet('NumberOfUnits',5); deadzone([-1,2])]
```

To specify the same nonlinearity for all inputs, specify a single input nonlinearity estimator.

### OutputNL — Output static nonlinearity

'pwlinear' (default) | 'sigmoidnet' | 'wavenet' | 'saturation' | 'deadzone' | 'polyld' | 'unitgain' | nonlinearity estimator object | array of nonlinearity estimators

Output static nonlinearity estimator, specified as one of the following:

'pwlinear' or pwlinear object (default)	Piecewise linear function
'sigmoidnet' or sigmoidnet object	Sigmoid network
'wavenet' or wavenet object	Wavelet network
'saturation' or saturation object	Saturation
'deadzone' or deadzone object	Dead zone
'polyld' or polyld object	One-dimensional polynomial
'unitgain' or [] or unitgain object	Unit gain
customnet object	Custom network — Similar to <b>sigmoidnet</b> , but with a user-defined replacement for the sigmoid function.

Specifying a character vector creates a nonlinearity estimator object with default settings. Use object representation instead to configure the properties of a nonlinearity estimator.

```
OutputNL = sigmoidnet;
OutputNL.NumberOfUnits = 10;
```

Alternatively, use the associated input nonlinearity estimator function with Name-Value pair arguments.

```
OutputNL = sigmoidnet('NumberOfUnits',10);
```

For  $n_y$  output channels, you can specify nonlinear estimators individually for each output channel by setting **OutputNL** to an  $n_y$ -by-1 array of nonlinearity estimators. To specify the same nonlinearity for all outputs, specify a single output nonlinearity estimator.

## LinModel — Discrete time linear model

`idpoly` | `idss` with `K = 0` | `idtf`

Discrete-time linear model used to specify the linear subsystem, specified as one of the following:

- Input-output polynomial model of Output-Error (OE) structure (`idpoly`)
- State-space model with no disturbance component (`idss` with `K = 0`)
- Transfer function model (`idtf`)

Typically, you estimate the model using `oe`, `n4sid`, or `tfest`.

## Name-Value Pair Arguments

Specify optional comma-separated pairs of `Name`, `Value` arguments. `Name` is the argument name and `Value` is the corresponding value. `Name` must appear inside quotes. You can specify several name and value pair arguments in any order as `Name1, Value1, ..., NameN, ValueN`.

Use `Name`, `Value` arguments to specify additional properties on page 1-662 of `idnlnhw` models during model creation. For example, `m = idnlnhw([2 3 1], 'pwlinear', 'wavenet', 'InputName', 'Volts', 'Ts', 0.1)` creates an `idnlnhw` model object with input nonlinearity estimator 'pwlinear', output nonlinearity estimator 'wavenet', input name `Volts`, and a sample time of 0.1 seconds.

## Properties

`idnlnhw` object properties include:

**`nb`, `nf`, `nk`**

Model orders and delays of the linear subsystem transfer function, where `nb` is the number of zeros plus 1, `nf` is the number of poles, and `nk` is the input delay.

For a MIMO transfer function with  $n_u$  inputs and  $n_y$  outputs, `nb`, `nf`, and `nk` are  $n_y$ -by- $n_u$  matrices whose  $i$ - $j$ th entry specifies the orders and delay of the transfer function from the  $j$ th input to the  $i$ th output.

**B**

$B$  polynomial of the linear block in the model structure on page 1-671, specified as a cell array of  $n_y$ -by- $n_u$  elements, where  $n_y$  is the number of outputs and  $n_u$  is the number of inputs. An element  $B\{i, j\}$  is a row vector representing the numerator polynomial for the  $j$ th input to  $i$ th output transfer function. The element contains  $n_k$  leading zeros, where  $n_k$  is the number of input delays.

**F**

$F$  polynomial of the linear block in the model structure on page 1-671, specified as a cell array of  $n_y$ -by- $n_u$  elements, where  $n_y$  is the number of outputs and  $n_u$  is the number of inputs. An element  $F\{i, j\}$  is a row vector representing the denominator polynomial for the  $j$ th input to  $i$ th output transfer function.

**InputNonlinearity**

Input nonlinearity estimator, specified as one of the following:

'pwllinear' or pwllinear object (default)	Piecewise linear function
'sigmoidnet' or sigmoidnet object	Sigmoid network
'wavenet' or wavenet object	Wavelet network
'saturation' or saturation object	Saturation
'deadzone' or deadzone object	Dead zone
'poly1d' or poly1d object	One- dimensional polynomial
'unitgain' or [] or unitgain object	Unit gain
customnet object	Custom network

Specifying a character vector creates a nonlinearity estimator object with default settings. Use object representation instead to configure the properties of a nonlinearity estimator.

```
InputNonlinearity = wavenet;
InputNonlinearity.NumberOfUnits = 10;
```

Alternatively, use the associated input nonlinearity estimator function with Name-Value pair arguments.

```
InputNonlinearity = wavenet('NumberOfUnits',10);
```

For  $n_u$  input channels, you can specify nonlinear estimators individually for each input channel by setting `InputNL` to an  $n_u$ -by-1 array of nonlinearity estimators. To specify the same nonlinearity for all inputs, specify a single input nonlinearity estimator.

**Default:** 'pwnlinear'

### **OutputNonlinearity**

Output nonlinearity estimator, specified as one of the following:

'pwnlinear' or pwnlinear object (default)	Piecewise linear function
'sigmoidnet' or sigmoidnet object	Sigmoid network
'wavenet' or wavenet object	Wavelet network
'saturation' or saturation object	Saturation
'deadzone' or deadzone object	Dead zone
'poly1d' or poly1d object	One- dimensional polynomial
'unitgain' or [] or unitgain object	Unit gain
customnet object	Custom network

Specifying a character vector creates a nonlinearity estimator object with default settings. Use object representation instead to configure the properties of a nonlinearity estimator.

```
OutputNonlinearity = sigmoidnet;  
OutputNonlinearity.NumberOfUnits = 10;
```

Alternatively, use the associated input nonlinearity estimator function with Name-Value pair arguments.

```
OutputNonlinearity = sigmoidnet('NumberOfUnits',10);
```

For  $n_y$  output channels, you can specify nonlinear estimators individually for each output channel by setting `OutputNL` to an  $n_y$ -by-1 array of nonlinearity estimators. To specify the same nonlinearity for all outputs, specify a single output nonlinearity estimator.

**Default:** 'pwnlinear'

## LinearModel

The linear model in the linear block of the model structure, represented as an `idpoly` object. This property is read only.

## Report

Summary report that contains information about the estimation options and results when the model is estimated using the `nlhw` command. Use `Report` to query a model for how it was estimated, including:

- Estimation method
- Estimation options
- Search termination conditions
- Estimation data fit

The contents of `Report` are irrelevant if the model was created by construction.

```
m = idnlhw([2 2 1]);
m.Report.OptionsUsed
```

```
ans =
```

```
    []
```

If you use `nlhw` to estimate the model, the fields of `Report` contain information on the estimation data, options, and results.

```
load iddata1;
m = nlhw(z1,[2 2 1],[],'pwlinear');
m.Report.OptionsUsed
```

Option set for the `nlhw` command:

```
InitialCondition: 'zero'
    Display: 'off'
Regularization: [1x1 struct]
SearchMethod: 'auto'
SearchOption: [1x1 idoptions.search.identsolver]
OutputWeight: 'noise'
    Advanced: [1x1 struct]
```

`Report` is a read-only property.

For more information on this property and how to use it, see “Output Arguments” on page 1-1007 in the *nIhw* reference page and “Estimation Report”.

## **TimeVariable**

Independent variable for the inputs, outputs, and—when available—internal states, specified as a character vector.

**Default:** 't'

## **NoiseVariance**

Noise variance (covariance matrix) of the model innovations  $e$ . Assignable value is an  $n_y$ -by- $n_y$  matrix. Typically set automatically by the estimation algorithm.

## **Ts**

Sample time.  $T_s$  is a positive scalar representing the sampling period. This value is expressed in the unit specified by the `TimeUnit` property of the model.

Changing this property does not discretize or resample the model.

**Default:** 1

## **TimeUnit**

Units for the time variable, the sample time  $T_s$ , and any time delays in the model, specified as one of the following values:

- 'nanoseconds'
- 'microseconds'
- 'milliseconds'
- 'seconds'
- 'minutes'
- 'hours'
- 'days'
- 'weeks'
- 'months'

- 'years'

Changing this property has no effect on other properties, and therefore changes the overall system behavior. Use `chgTimeUnit` to convert between time units without modifying system behavior.

**Default:** 'seconds'

### **InputName**

Input channel names, specified as one of the following:

- Character vector — For single-input models, for example, 'controls'.
- Cell array of character vectors — For multi-input models.

Alternatively, use automatic vector expansion to assign input names for multi-input models. For example, if `sys` is a two-input model, enter:

```
sys.InputName = 'controls';
```

The input names automatically expand to {'controls(1)'; 'controls(2)'}

When you estimate a model using an `iddata` object, `data`, the software automatically sets `InputName` to `data.InputName`.

You can use the shorthand notation `u` to refer to the `InputName` property. For example, `sys.u` is equivalent to `sys.InputName`.

Input channel names have several uses, including:

- Identifying channels on model display and plots
- Extracting subsystems of MIMO systems
- Specifying connection points when interconnecting models

**Default:** '' for all input channels

### **InputUnit**

Input channel units, specified as one of the following:

- Character vector — For single-input models, for example, 'seconds'.

- Cell array of character vectors — For multi-input models.

Use `InputUnit` to keep track of input signal units. `InputUnit` has no effect on system behavior.

**Default:** '' for all input channels

### **InputGroup**

Input channel groups. The `InputGroup` property lets you assign the input channels of MIMO systems into groups and refer to each group by name. Specify input groups as a structure. In this structure, field names are the group names, and field values are the input channels belonging to each group. For example:

```
sys.InputGroup.controls = [1 2];  
sys.InputGroup.noise = [3 5];
```

creates input groups named `controls` and `noise` that include input channels 1, 2 and 3, 5, respectively. You can then extract the subsystem from the `controls` inputs to all outputs using:

```
sys(:, 'controls')
```

**Default:** Struct with no fields

### **OutputName**

Output channel names, specified as one of the following:

- Character vector — For single-output models. For example, 'measurements'.
- Cell array of character vectors — For multi-output models.

Alternatively, use automatic vector expansion to assign output names for multi-output models. For example, if `sys` is a two-output model, enter:

```
sys.OutputName = 'measurements';
```

The output names automatically expand to `{'measurements(1)'; 'measurements(2)'}`.

When you estimate a model using an `iddata` object, `data`, the software automatically sets `OutputName` to `data.OutputName`.

You can use the shorthand notation `y` to refer to the `OutputName` property. For example, `sys.y` is equivalent to `sys.OutputName`.

Output channel names have several uses, including:

- Identifying channels on model display and plots
- Extracting subsystems of MIMO systems
- Specifying connection points when interconnecting models

**Default:** '' for all output channels

### OutputUnit

Output channel units, specified as one of the following:

- Character vector — For single-output models. For example, 'seconds'.
- Cell array of character vectors — For multi-output models.

Use `OutputUnit` to keep track of output signal units. `OutputUnit` has no effect on system behavior.

**Default:** '' for all output channels

### OutputGroup

Output channel groups. The `OutputGroup` property lets you assign the output channels of MIMO systems into groups and refer to each group by name. Specify output groups as a structure. In this structure, field names are the group names, and field values are the output channels belonging to each group. For example:

```
sys.OutputGroup.temperature = [1];  
sys.InputGroup.measurement = [3 5];
```

creates output groups named `temperature` and `measurement` that include output channels 1, and 3, 5, respectively. You can then extract the subsystem from all inputs to the measurement outputs using:

```
sys('measurement', :)
```

**Default:** Struct with no fields

## Name

System name, specified as a character vector. For example, 'system\_1'.

**Default:** ''

## Notes

Any text that you want to associate with the system, stored as a string or a cell array of character vectors. The property stores whichever data type you provide. For instance, if `sys1` and `sys2` are dynamic system models, you can set their `Notes` properties as follows:

```
sys1.Notes = "sys1 has a string.";
sys2.Notes = 'sys2 has a character vector.';
sys1.Notes
sys2.Notes
```

```
ans =
```

```
    "sys1 has a string."
```

```
ans =
```

```
    'sys2 has a character vector.'
```

**Default:** [0×1 string]

## UserData

Any type of data you want to associate with system, specified as any MATLAB data type.

**Default:** []

## Output Arguments

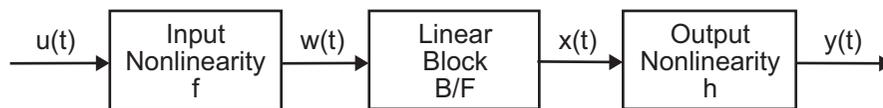
**sys** — Hammerstein-Wiener model  
idnlhw object

Hammerstein-Wiener model, returned as an `idnlhw` object. This model is created using the specified model orders and delays, input and output nonlinearity estimators, and properties.

## Definitions

### Hammerstein-Wiener Model Structure

This block diagram represents the structure of a Hammerstein-Wiener model:



Where,

- $f$  is a nonlinear function that transforms input data  $u(t)$  as  $w(t) = f(u(t))$ .

$w(t)$ , an internal variable, is the output of the Input Nonlinearity block and has the same dimension as  $u(t)$ .

- $B/F$  is a linear transfer function that transforms  $w(t)$  as  $x(t) = (B/F)w(t)$ .

$x(t)$ , an internal variable, is the output of the Linear block and has the same dimension as  $y(t)$ .

$B$  and  $F$  are similar to polynomials in a linear Output-Error model. For more information about Output-Error models, see “What Are Polynomial Models?”.

For  $n_y$  outputs and  $n_u$  inputs, the linear block is a transfer function matrix containing entries:

$$\frac{B_{j,i}(q)}{F_{j,i}(q)}$$

where  $j = 1, 2, \dots, n_y$  and  $i = 1, 2, \dots, n_u$ .

- $h$  is a nonlinear function that maps the output of the linear block  $x(t)$  to the system output  $y(t)$  as  $y(t) = h(x(t))$ .

Because  $f$  acts on the input port of the linear block, this function is called the *input nonlinearity*. Similarly, because  $h$  acts on the output port of the linear block, this function is called the *output nonlinearity*. If your system contains several inputs and outputs, you must define the functions  $f$  and  $h$  for each input and output signal. You do not have to include both the input and the output nonlinearity in the model structure. When a model contains only the input nonlinearity  $f$ , it is called a Hammerstein model. Similarly, when the model contains only the output nonlinearity  $h$ , it is called a *Wiener* model.

The software computes the Hammerstein-Wiener model output  $y$  in three stages:

- 1 Compute  $w(t) = f(u(t))$  from the input data.

$w(t)$  is an input to the linear transfer function  $B/F$ .

The input nonlinearity is a static (*memoryless*) function, where the value of the output a given time  $t$  depends only on the input value at time  $t$ .

You can configure the input nonlinearity as a sigmoid network, wavelet network, saturation, dead zone, piecewise linear function, one-dimensional polynomial, or a custom network. You can also remove the input nonlinearity.

- 2 Compute the output of the linear block using  $w(t)$  and initial conditions:  $x(t) = (B/F)w(t)$ .

You can configure the linear block by specifying the orders of numerator  $B$  and denominator  $F$ .

- 3 Compute the model output by transforming the output of the linear block  $x(t)$  using the nonlinear function  $h$  as  $y(t) = h(x(t))$ .

Similar to the input nonlinearity, the output nonlinearity is a static function. You can configure the output nonlinearity in the same way as the input nonlinearity. You can also remove the output nonlinearity, such that  $y(t) = x(t)$ .

Resulting models are `idnlhw` objects that store all model data, including model parameters and nonlinearity estimators. For more information about these objects, see “Nonlinear Model Structures”.

### Definition of `idnlhw` States

The states of a Hammerstein-Wiener model correspond to the states of the linear block in the model structure. The linear block contains all the dynamic elements of the model. If

the linear block is not a state-space structure, the states are defined as those of model `Mss`, where `Mss = idss(Model.LinearModel)` and `Model` is the `idnlhw` object.

States are required for simulation, prediction, and linearization of Hammerstein-Wiener models. To specify the initial states:

- Use `findstates` to search for state values for simulation and prediction with `sim`, `predict`, and `compare`.
- Use `findop` when linearizing the model with `linearize`.
- Alternatively, specify the states manually.

## See Also

`customnet` | `findop` | `linear` | `linearize` | `nlhw` | `pem` | `poly1d` | `saturation` | `saturation` | `sigmoidnet` | `wavenet`

## Topics

“Estimate Multiple Hammerstein-Wiener Models”

“Estimate Hammerstein-Wiener Models Initialized Using Linear OE Models”

“Identifying Hammerstein-Wiener Models”

“Initialize Hammerstein-Wiener Estimation Using Linear Model”

**Introduced in R2007a**

## idpar

Create parameter for initial states and input level estimation

### Syntax

```
p = idpar(paramvalue)
p = idpar(paramname,paramvalue)
```

### Description

`p = idpar(paramvalue)` creates an estimable parameter with initial value `paramvalue`. The parameter, `p`, is either scalar or array-valued, with the same dimensions as `paramvalue`. You can configure attributes of the parameter, such as which elements are fixed and which are estimated, and lower and upper bounds.

`p = idpar(paramname,paramvalue)` sets the `Name` property of `p` to `paramname`.

### Input Arguments

#### **paramvalue**

Initial parameter value.

`paramvalue` is a numeric scalar or array that determines both the dimensions and initial values of the estimable parameter `p`. For example, `p = idpar(eye(3))` creates a 3-by-3 parameter whose initial value is the identity matrix.

`paramvalue` should be:

- A column vector of length  $N_x$ , the number of states to estimate, if you are using `p` for initial state estimation.
- An  $N_x$ -by- $N_e$  array, if you are using `p` for initial state estimation with multi-experiment data.  $N_e$  is the number of experiments.

- A column vector of length  $N_u$ , the number of inputs to estimate, if you are using `p` for input level estimation.
- An  $N_u$ -by- $N_e$  array, if you are using `p` for input level estimation with multi-experiment data.

If the initial value of a parameter is unknown, use NaN.

### **paramname**

Name property of `p`, specified as a character vector. For example, you can assign 'x0' as the name of a parameter created for initial state estimation.

The `Name` property is not used in state estimation or input level estimation. You can optionally assign a name for convenience.

**Default:** 'par'

## **Output Arguments**

### **p**

Estimable parameter, specified as a `param.Continuous` object.

`p` can be either scalar- or array-valued. `p` takes its dimensions and initial value from `paramvalue`.

`p` contains the following fields:

- `Value` — Scalar or array value of the parameter.

The dimension and initial value of `p.Value` are taken from `paramvalue` when `p` is created.

- `Minimum` — Lower bound for the parameter value. When you use `p` in state estimation or input value estimation, the estimated value of the parameter does not drop below `p.Minimum`.

The dimensions of `p.Minimum` must match the dimensions of `p.Value`.

For array-valued parameters, you can:

- Specify lower bounds on individual array elements. For example, `p.Minimum([1 4]) = -5`.
- Use scalar expansion to set the lower bound for all array elements. For example, `p.Minimum = -5`

**Default:** `-Inf`

- **Maximum** — Upper bound for the parameter value. When you use `p` in state estimation or input value estimation, the estimated value of the parameter does not exceed `p.Maximum`.

The dimensions of `p.Maximum` must match the dimensions of `p.Value`.

For array-valued parameters, you can:

- Specify upper bounds on individual array elements. For example, `p.Maximum([1 4]) = 5`.
- Use scalar expansion to set the upper bound for all array elements. For example, `p.Maximum = 5`

**Default:** `Inf`

- **Free** — Boolean specifying whether the parameter is a free estimation variable.

The dimensions of `p.Free` must match the dimensions of `p.Value`. By default, all values are free (`p.Free = true`).

If you want to estimate `p.Value(k)`, set `p.Free(k) = true`. To fix `p.Value(k)`, set `p.Free(k) = false`. Doing so allows you to control which states or input values are estimated and which are not.

For array-valued parameters, you can:

- Fix individual array elements. For example, `p.Free([1 4]) = false;`  
`p.Free = [1 0; 0 1]`.
- Use scalar expansion to fix all array elements. For example, `p.Free = false`.

**Default:** `true (1)`

- **Scale** — Scaling factor for normalizing the parameter value.

`p.Scale` is not used in initial state estimation or input value estimation.

**Default:** `1`

- **Info** — Structure array for storing parameter units and labels. The structure has `Label` and `Unit` fields.

Use these fields for your convenience, to store parameter units and labels. For example, `p.Info(1,1).Unit = 'rad/m';`  
`p.Info(1,1).Label = 'engine speed'.`

The dimensions of `p.Info` must match the dimensions of `p.Value`.

**Default:** '' for both `Label` and `Unit` fields

- **Name** — Parameter name.

This property is read-only. It is set to the `paramname` input argument when you create the parameter.

**Default:** ''

## Examples

### Create and Configure Parameter for State Estimation

Create and configure a parameter for estimating the initial state values of a 4-state system. Fix the first state value to 1. Limit the second and third states to values between 0 and 1.

```
paramvalue = [1; nan(3,1)];
p = idpar('x0',paramvalue);
p.Free(1) = 0;
p.Minimum([2 3]) = 0;
p.Maximum([2 3]) = 1;
```

The column vector `paramvalue` specifies an initial value of 1 for the first state. `paramvalue` further specifies unknown values for the remaining 3 states.

Setting `p.Free(1)` to false fixes `p.Value(1)` to 1. Estimation using `p` does not alter that value.

Setting `p.Minimum` and `p.Maximum` for the second and third entries in `p` limits the range that those values can take when `p` is used in estimation.

You can now use `p` in initial state estimation, such as with the `findstates` command. For example, use `opt = findstatesOptions('InitialState',p)` to create a `findstates` options set that uses `p`. Then, call `findstates` with that options set.

### Tips

Use `idpar` to create estimable parameters for:

- Initial state estimation for state-space model estimation (`ssest`), prediction (`predict`), and forecasting (`forecast`)
- Explicit initial state estimation with `findstates`
- Input level estimation for process model estimation with `pem`

Specifying estimable state values or input levels gives you explicit control over the behavior of individual state values during estimation.

### See Also

`findstates` | `findstatesOptions` | `forecast` | `pem` | `predict` | `ssest`

**Introduced in R2012a**

# idpoly

Polynomial model with identifiable parameters

## Syntax

```
sys = idpoly(A,B,C,D,F,NoiseVariance,Ts)
sys = idpoly(A,B,C,D,F,NoiseVariance,Ts,Name,Value)

sys = idpoly(A)
sys = idpoly(A,[],C,D,[],NoiseVariance,Ts)
sys = idpoly(A,[],C,D,[],NoiseVariance,Ts,Name,Value)

sys = idpoly(sys0)
sys = idpoly(sys0,'split')
```

## Description

`sys = idpoly(A,B,C,D,F,NoiseVariance,Ts)` creates a polynomial model with identifiable coefficients. `A`, `B`, `C`, `D`, and `F` specify the initial values of the coefficients. `NoiseVariance` specifies the initial value of the variance of the white noise source. `Ts` is the model sample time.

`sys = idpoly(A,B,C,D,F,NoiseVariance,Ts,Name,Value)` creates a polynomial model using additional options specified by one or more `Name,Value` pair arguments.

`sys = idpoly(A)` creates a time-series model with only an autoregressive term. In this case, `sys` represents the AR model given by  $A(q^{-1})y(t) = e(t)$ . The noise  $e(t)$  has variance 1. `A` specifies the initial values of the estimable coefficients.

`sys = idpoly(A,[],C,D,[],NoiseVariance,Ts)` creates a time-series model with an autoregressive and a moving average term. The inputs `A`, `C`, and `D`, specify the initial values of the estimable coefficients. `NoiseVariance` specifies the initial value of the noise  $e(t)$ . `Ts` is the model sample time. (Omit `NoiseVariance` and `Ts` to use their default values.)

If `D = []`, then `sys` represents the ARMA model given by:

$$A(q^{-1})y(t) = C(q^{-1})e(t).$$

`sys = idpoly(A, [], C, D, [], NoiseVariance, Ts, Name, Value)` creates a time-series model using additional options specified by one or more `Name, Value` pair arguments.

`sys = idpoly(sys0)` converts any dynamic system model, `sys0`, to `idpoly` model form.

`sys = idpoly(sys0, 'split')` converts `sys0` to `idpoly` model form, and treats the last  $N_y$  input channels of `sys0` as noise channels in the returned model. `sys0` must be a numeric (nonidentified) `tf`, `zpk`, or `ss` model object. Also, `sys0` must have at least as many inputs as outputs.

## Object Description

An `idpoly` model represents a system as a continuous-time or discrete-time polynomial model with identifiable (estimable) coefficients.

A polynomial model of a system with input vector  $u$ , output vector  $y$ , and disturbance  $e$  takes the following form in discrete time:

$$A(q)y(t) = \frac{B(q)}{F(q)}u(t) + \frac{C(q)}{D(q)}e(t)$$

In continuous time, a polynomial model takes the following form:

$$A(s)Y(s) = \frac{B(s)}{F(s)}U(s) + \frac{C(s)}{D(s)}E(s)$$

$U(s)$  are the Laplace transformed inputs to `sys`.  $Y(s)$  are the Laplace transformed outputs.  $E(s)$  is the Laplace transform of the disturbance.

For `idpoly` models, the coefficients of the polynomials  $A$ ,  $B$ ,  $C$ ,  $D$ , and  $F$  can be estimable parameters. The `idpoly` model stores the values of these matrix elements in the `A`, `B`, `C`, `D`, and `F` properties of the model.

Time-series models are special cases of polynomial models for systems without measured inputs. For AR models, B and F are empty, and C and D are 1 for all outputs. For ARMA models, B and F are empty, while D is 1.

There are three ways to obtain an `idpoly` model:

- Estimate the `idpoly` model based on output or input-output measurements of a system, using commands such as `polyest`, `arx`, `armax`, `oe`, `bj`, `iv4`, or `ivar`. These commands estimate the values of the free polynomial coefficients. The estimated values are stored in the A, B, C, D, and F properties of the resulting `idpoly` model. The `Report` property of the resulting model stores information about the estimation, such as handling of initial conditions and options used in estimation.

When you obtain an `idpoly` model by estimation, you can extract estimated coefficients and their uncertainties from the model using commands such as `polydata`, `getpar`, or `getcov`.

- Create an `idpoly` model using the `idpoly` command. You can create an `idpoly` model to configure an initial parameterization for estimation of a polynomial model to fit measured response data. When you do so, you can specify constraints on the polynomial coefficients. For example, you can fix the values of some coefficients, or specify minimum or maximum values for the free coefficients. You can then use the configured model as an input argument to `polyest` to estimate parameter values with those constraints.
- Convert an existing dynamic system model to an `idpoly` model using the `idpoly` command.

## Examples

### Multi-Output ARMAX Model

Create an `idpoly` model representing the one-input, two-output ARMAX model described by the following equations:

$$\begin{aligned}
 y_1(t) + 0.5y_1(t-1) + 0.9y_2(t-1) + 0.1y_2(t-2) &= \\
 &u(t) + 5u(t-1) + 2u(t-2) + e_1(t) + 0.01e_1(t-1) \\
 y_2(t) + 0.05y_2(t-1) + 0.3y_2(t-2) &= \\
 &10u(t-2) + e_2(t) + 0.1e_2(t-1) + 0.02e_2(t-2).
 \end{aligned}$$

$y_1$  and  $y_2$  are the two outputs, and  $u$  is the input.  $e_1$  and  $e_2$  are the white noise disturbances on the outputs  $y_1$  and  $y_2$  respectively.

To create the `idpoly` model, define the A, B, and C polynomials that describe the relationships between the outputs, inputs, and noise values. (Because there are no denominator terms in the system equations, B and F are 1.)

Define the cell array containing the coefficients of the A polynomials.

```
A = cell(2,2);  
A{1,1} = [1 0.5];  
A{1,2} = [0 0.9 0.1];  
A{2,1} = [0];  
A{2,2} = [1 0.05 0.3];
```

You can read the values of each entry in the A cell array from the left side of the equations describing the system. For example, A{1,1} describes the polynomial that gives the dependence of  $y_1$  on itself. This polynomial is  $A_{11} = 1 + 0.5q^{-1}$ , because each factor of  $q^{-1}$  corresponds to a unit time decrement. Therefore, A{1,1} = [1 0.5], giving the coefficients of  $A_{11}$  in increasing exponents of  $q^{-1}$ .

Similarly, A{1,2} describes the polynomial that gives the dependence of  $y_1$  on  $y_2$ . From the equations,  $A_{12} = 0 + 0.9q^{-1} + 0.1q^{-2}$ . Thus, A{1,2} = [0 0.9 0.1].

The remaining entries in A are similarly constructed.

Define the cell array containing the coefficients of the B polynomials.

```
B = cell(2,1);  
B{1,1} = [1 5 2];  
B{2,1} = [0 0 10];
```

B describes the polynomials that give the dependence of the outputs  $y_1$  and  $y_2$  on the input  $u$ . From the equations,  $B_{11} = 1 + 5q^{-1} + 2q^{-2}$ . Therefore, B{1,1} = [1 5 2].

Similarly, from the equations,  $B_{21} = 0 + 0q^{-1} + 10q^{-2}$ . Therefore, B{2,1} = [0 0 10].

Define the cell array containing the coefficients of the C polynomials.

```
C = cell(2,1);
C{1,1} = [1 0.01];
C{2,1} = [1 0.1 0.02];
```

C describes the polynomials that give the dependence of the outputs  $y_1$  and  $y_2$  on the noise terms  $e_1$  and  $e_2$ . The entries of C can be read from the equations similarly to those of A and B.

Create an idpoly model with the specified coefficients.

```
sys = idpoly(A,B,C)
```

```
sys =
Discrete-time ARMAX model:
  Model for output number 1:  $A(z)y_1(t) = -A_i(z)y_i(t) + B(z)u(t) + C(z)e_1(t)$ 
     $A(z) = 1 + 0.5 z^{-1}$ 

     $A_2(z) = 0.9 z^{-1} + 0.1 z^{-2}$ 

     $B(z) = 1 + 5 z^{-1} + 2 z^{-2}$ 

     $C(z) = 1 + 0.01 z^{-1}$ 

  Model for output number 2:  $A(z)y_2(t) = B(z)u(t) + C(z)e_2(t)$ 
     $A(z) = 1 + 0.05 z^{-1} + 0.3 z^{-2}$ 

     $B(z) = 10 z^{-2}$ 

     $C(z) = 1 + 0.1 z^{-1} + 0.02 z^{-2}$ 

Sample time: unspecified

Parameterization:
  Polynomial orders:  na=[1 2;0 2]  nb=[3;1]  nc=[1;2]
  nk=[0;2]
  Number of free coefficients: 12
  Use "polydata", "getpvec", "getcov" for parameters and their uncertainties.

Status:
Created by direct construction or transformation. Not estimated.
```

The display shows all the polynomials and allows you to verify them. The display also states that there are 12 free coefficients. Leading terms of diagonal entries in A are always fixed to 1. Leading terms of all other entries in A are always fixed to 0.

You can use `sys` to specify an initial parametrization for estimation with such commands as `polyest` or `armax`.

## Input Arguments

### A, B, C, D, F

Initial values of polynomial coefficients.

For SISO models, specify the initial values of the polynomial coefficients as row vectors. Specify the coefficients in order of:

- Ascending powers of  $z^{-1}$  or  $q^{-1}$  (for discrete-time polynomial models).
- Descending powers of  $s$  or  $p$  (for continuous-time polynomial models).

The leading coefficients of A, C, D, and F must be 1. Use NaN for any coefficient whose initial value is not known.

For MIMO models with  $N_y$  outputs and  $N_u$  inputs, A, B, C, D, and F are cell arrays of row vectors. Each entry in the cell array contains the coefficients of a particular polynomial that relates input, output, and noise values.

Polynomial	Dimension	Relation Described
A	$N_y$ -by- $N_y$ array of row vectors	A{ <i>i</i> , <i>j</i> } contains coefficients of relation between output $y_i$ and output $y_j$
B, F	$N_y$ -by- $N_u$ array of row vectors	B{ <i>i</i> , <i>j</i> } and F{ <i>i</i> , <i>j</i> } contain coefficients of relations between output $y_i$ and input $u_j$
C, D	$N_y$ -by-1 array of row vectors	C{ <i>i</i> } and D{ <i>i</i> } contain coefficients of relations between output $y_i$ and noise $e_i$

The leading coefficients of the diagonal entries of  $A$  ( $A\{i, i\}$ ,  $i=1:N_y$ ) must be 1. The leading coefficients of the off-diagonal entries of  $A$  must be zero, for causality. The leading coefficients of all entries of  $C$ ,  $D$ , and  $F$ , must be 1.

Use `[]` for any polynomial that is not present in the wanted model structure. For example, to create an ARX model, use `[]` for  $C$ ,  $D$ , and  $F$ . For an ARMA time series, use `[]` for  $B$  and  $F$ .

**Default:**  $B = []$ ;  $C = 1$  for all outputs;  $D = 1$  for all outputs;  $F = []$

### **Ts**

Sample time. For continuous-time models,  $T_s = 0$ . For discrete-time models,  $T_s$  is a positive scalar representing the sample time expressed in the unit specified by the `TimeUnit` property of the model. To denote a discrete-time model with unspecified sample time, set  $T_s = -1$ .

**Default:** -1 (discrete-time model with unspecified sample time)

### **NoiseVariance**

The variance (covariance matrix) of the model innovations  $e$ .

An identified model includes a white, Gaussian noise component  $e(t)$ . `NoiseVariance` is the variance of this noise component. Typically, a model estimation function (such as `polyest`) determines this variance. Use this input to specify an initial value for the noise variance when you create an `idpoly` model.

For SISO models, `NoiseVariance` is a scalar. For MIMO models, `NoiseVariance` is a  $N_y$ -by- $N_y$  matrix, where  $N_y$  is the number of outputs in the system.

**Default:**  $N_y$ -by- $N_y$  identity matrix

### **sys0**

Dynamic system.

Any dynamic system to be converted into an `idpoly` object.

When `sys0` is an identified model, its estimated parameter covariance is lost during conversion. If you want to translate the estimated parameter covariance during the conversion, use `translatecov`.

For the syntax `sys = idpoly(sys0, 'split')`, `sys0` must be a numeric (non-identified) `tf`, `zpk`, or `ss` model object. Also, `sys0` must have at least as many inputs as outputs. Finally, the subsystem `sys0(:, Ny+1:Nu)` must be biproper.

### Name-Value Pair Arguments

Specify optional comma-separated pairs of `Name`, `Value` arguments. `Name` is the argument name and `Value` is the corresponding value. `Name` must appear inside quotes. You can specify several name and value pair arguments in any order as `Name1, Value1, . . . , NameN, ValueN`.

Use `Name`, `Value` arguments to specify additional properties on page 1-686 of `idpoly` models during model creation. For example, `idpoly(A,B,C,D,F, 1,0, 'InputName', 'Voltage')` creates an `idpoly` model with the `InputName` property set to `Voltage`.

### Properties

`idpoly` object properties include:

#### **A, B, C, D, F**

Values of polynomial coefficients.

If you create an `idpoly` model `sys` using the `idpoly` command, `sys.A`, `sys.B`, `sys.C`, `sys.D`, and `sys.F` contain the initial coefficient values that you specify with the `A`, `B`, `C`, `D`, and `F` input arguments, respectively.

If you obtain an `idpoly` model by identification, then `sys.A`, `sys.B`, `sys.C`, `sys.D`, and `sys.F` contain the estimated values of the coefficients.

For an `idpoly` model `sys`, each property `sys.A`, `sys.B`, `sys.C`, `sys.D`, and `sys.F` is an alias to the corresponding `Value` entry in the `Structure` property of `sys`. For example, `sys.A` is an alias to the value of the property `sys.Structure.A.Value`.

For SISO polynomial models, the values of the numerator coefficients are stored as a row vector in order of:

- Ascending powers of  $z^{-1}$  or  $q^{-1}$  (for discrete-time transfer functions).

- Descending powers of  $s$  or  $p$  (for continuous-time transfer functions).

The leading coefficients of A, C, and D are fixed to 1. Any coefficient whose initial value is not known is stored as NaN.

For MIMO models with  $N_y$  outputs and  $N_u$  inputs, A, B, C, D, and F are cell arrays of row vectors. Each entry in the cell array contains the coefficients of a particular polynomial that relates input, output, and noise values.

Polynomial	Dimension	Relation Described
A	$N_y$ -by- $N_y$ array of row vectors	A{ $i, j$ } contains coefficients of relation between output $y_i$ and output $y_j$
B, F	$N_y$ -by- $N_u$ array of row vectors	B{ $i, j$ } and F{ $i, j$ } contain coefficients of relations between output $y_i$ and input $u_j$
C, D	$N_y$ -by-1 array of row vectors	C{ $i$ } and D{ $i$ } contain coefficients of relations between output $y_i$ and noise $e_i$

The leading coefficients of the diagonal entries of A (A{ $i, i$ },  $i=1:N_y$ ) are fixed to 1. The leading coefficients of the off-diagonal entries of A are fixed to zero. The leading coefficients of all entries of C, D, and F, are fixed to 1.

For a time series (a model with no measured inputs), B = [] and F = [].

**Default:** B = []; C = 1 for all outputs; D = 1 for all outputs; F = []

### Variable

Polynomial model display variable, specified as one of the following values:

- 'z^-1' — Default for discrete-time models
- 'q^-1' — Equivalent to 'z^-1'
- 's' — Default for continuous-time models
- 'p' — Equivalent to 's'

The value of Variable is reflected in the display, and also affects the interpretation of the A, B, C, D, and F coefficient vectors for discrete-time models. For

Variable = 'z<sup>-1</sup>' or 'q<sup>-1</sup>', the coefficient vectors are ordered as ascending powers of the variable.

## **IODelay**

Transport delays. `IODelay` is a numeric array specifying a separate transport delay for each input/output pair.

If you create an `idpoly` model `sys` using the `idpoly` command, `sys.IODelay` contains the initial values of the transport delay that you specify with a `Name, Value` argument pair.

For an `idpoly` model `sys`, the property `sys.IODelay` is an alias to the value of the property `sys.Structure.IODelay.Value`.

For continuous-time systems, transport delays are expressed in the time unit stored in the `TimeUnit` property. For discrete-time systems, transport delays are expressed as integers denoting delay of a multiple of the sample time `Ts`.

For a MIMO system with `Ny` outputs and `Nu` inputs, set `IODelay` is a `Ny-by-Nu` array, where each entry is a numerical value representing the transport delay for the corresponding input/output pair. You can set `IODelay` to a scalar value to apply the same delay to all input/output pairs.

**Default:** 0 for all input/output pairs

## **IntegrateNoise**

Logical vector, denoting presence or absence of integration on noise channels.

Specify `IntegrateNoise` as a logical vector of length equal to the number of outputs.

`IntegrateNoise(i) = true` indicates that the noise channel for the *i*th output contains an integrator. In this case, the corresponding *D* polynomial contains an additional term which is not represented in the property `sys.D`. This integrator term is equal to `[1 0]` for continuous-time systems, and equal to `[1 -1]` for discrete-time systems.

**Default:** 0 for all output channels

## **Structure**

Information about the estimable parameters of the `idpoly` model. `sys.Structure.A`, `sys.Structure.B`, `sys.Structure.C`, `sys.Structure.D`, and `sys.Structure.F`

contain information about the polynomial coefficients. `sys.Structure.IODelay` contains information about the transport delay. `sys.Structure.IntegrateNoise` contain information about the integration terms on the noise. Each contains the following fields:

- **Value** — Parameter values. For example, `sys.Structure.A.Value` contains the initial or estimated values of the  $A$  coefficients.

NaN represents unknown parameter values.

For SISO models, each property `sys.A`, `sys.B`, `sys.C`, `sys.D`, `sys.F`, and `sys.IODelay` is an alias to the corresponding `Value` entry in the `Structure` property of `sys`. For example, `sys.A` is an alias to the value of the property `sys.Structure.A.Value`

For MIMO models, `sys.A{i,j}` is an alias to `sys.Structure.A(i,j).Value`, and similarly for the other identifiable coefficient values.

- **Minimum** — Minimum value that the parameter can assume during estimation. For example, `sys.Structure.IODelay.Minimum = 0.1` constrains the transport delay to values greater than or equal to 0.1.

`sys.Structure.IODelay.Minimum` must be greater than or equal to zero.

- **Maximum** — Maximum value that the parameter can assume during estimation.
- **Free** — Logical value specifying whether the parameter is a free estimation variable. If you want to fix the value of a parameter during estimation, set the corresponding `Free = false`. For example, if  $B$  is a 3-by-3 matrix, `sys.Structure.B.Free = eyes(3)` fixes all of the off-diagonal entries in  $B$  to the values specified in `sys.Structure.B.Value`. In this case, only the diagonal entries in  $B$  are estimable.

For fixed values, such as the leading coefficients in `sys.Structure.B.Value`, the corresponding value of `Free` is always `false`.

- **Scale** — Scale of the parameter's value. `Scale` is not used in estimation.
- **Info** — Structure array for storing parameter units and labels. The structure has `Label` and `Unit` fields.

Specify parameter units and labels as character vectors. For example, `'Time'`.

For a MIMO model with  $N_y$  outputs and  $N_u$  inputs, the dimensions of the `Structure` elements are as follows:

- `sys.Structure.A` — Ny-by-Ny
- `sys.Structure.B` — Ny-by-Nu
- `sys.Structure.C` — Ny-by-1
- `sys.Structure.D` — Ny-by-1
- `sys.Structure.F` — Ny-by-Nu

An inactive polynomial, such as the **B** polynomial in a time-series model, is not available as a parameter in the `Structure` property. For example, `sys = idpoly([1 -0.2 0.5])` creates an AR model. `sys.Structure` contains the fields `sys.Structure.A`, `sys.Structure.IODelay`, and `sys.Structure.IntegrateNoise`. However, there is no field in `sys.Structure` corresponding to **B**, **C**, **D**, or **F**.

## NoiseVariance

The variance (covariance matrix) of the model innovations  $e$ .

An identified model includes a white Gaussian noise component  $e(t)$ . `NoiseVariance` is the variance of this noise component. Typically, the model estimation function (such as `arx`) determines this variance.

For SISO models, `NoiseVariance` is a scalar. For MIMO models, `NoiseVariance` is a  $N_y$ -by- $N_y$  matrix, where  $N_y$  is the number of outputs in the system.

## Report

Summary report that contains information about the estimation options and results when the polynomial model is obtained using estimation commands, such as `polyest`, `armax`, `oe`, and `bj`. Use `Report` to query a model for how it was estimated, including its:

- Estimation method
- Estimation options
- Search termination conditions
- Estimation data fit and other quality metrics

The contents of `Report` are irrelevant if the model was created by construction.

```
m = idpoly({[1 0.5]}, {[1 5]}, {[1 0.01]});  
m.Report.OptionsUsed
```

```
ans =  
  
    []
```

If you obtain the polynomial model using estimation commands, the fields of `Report` contain information on the estimation data, options, and results.

```
load iddata2 z2;  
m = polyest(z2,[2 2 3 3 2 1]);  
m.Report.OptionsUsed
```

Option set for the `polyest` command:

```
    InitialCondition: 'auto'  
                Focus: 'prediction'  
EstimateCovariance: 1  
                Display: 'off'  
                InputOffset: []  
                OutputOffset: []  
Regularization: [1x1 struct]  
SearchMethod: 'auto'  
SearchOptions: [1x1 idoptions.search.identsolver]  
Advanced: [1x1 struct]
```

`Report` is a read-only property.

For more information on this property and how to use it, see the `Output Arguments` section of the corresponding estimation command reference page and “`Estimation Report`”.

### **InputDelay**

Input delay for each input channel, specified as a scalar value or numeric vector. For continuous-time systems, specify input delays in the time unit stored in the `TimeUnit` property. For discrete-time systems, specify input delays in integer multiples of the sample time `Ts`. For example, `InputDelay = 3` means a delay of three sample times.

For a system with `Nu` inputs, set `InputDelay` to an `Nu`-by-1 vector. Each entry of this vector is a numerical value that represents the input delay for the corresponding input channel.

You can also set `InputDelay` to a scalar value to apply the same delay to all channels.

**Default:** 0

## OutputDelay

Output delays.

For identified systems, such as `idpoly`, `OutputDelay` is fixed to zero.

## Ts

Sample time. For continuous-time models,  $T_s = 0$ . For discrete-time models,  $T_s$  is a positive scalar representing the sample time expressed in the unit specified by the `TimeUnit` property of the model. To denote a discrete-time model with unspecified sample time, set  $T_s = -1$ .

Changing this property does not discretize or resample the model. Use `c2d` and `d2c` to convert between continuous- and discrete-time representations. Use `d2d` to change the sample time of a discrete-time system.

**Default:** -1 (discrete-time model with unspecified sample time)

## TimeUnit

Units for the time variable, the sample time  $T_s$ , and any time delays in the model, specified as one of the following values:

- 'nanoseconds'
- 'microseconds'
- 'milliseconds'
- 'seconds'
- 'minutes'
- 'hours'
- 'days'
- 'weeks'
- 'months'
- 'years'

Changing this property has no effect on other properties, and therefore changes the overall system behavior. Use `chgTimeUnit` to convert between time units without modifying system behavior.

**Default:** 'seconds'

## InputName

Input channel names, specified as one of the following:

- Character vector — For single-input models, for example, 'controls'.
- Cell array of character vectors — For multi-input models.

Alternatively, use automatic vector expansion to assign input names for multi-input models. For example, if `sys` is a two-input model, enter:

```
sys.InputName = 'controls';
```

The input names automatically expand to `{'controls(1)'; 'controls(2)'}`.

When you estimate a model using an `iddata` object, `data`, the software automatically sets `InputName` to `data.InputName`.

You can use the shorthand notation `u` to refer to the `InputName` property. For example, `sys.u` is equivalent to `sys.InputName`.

Input channel names have several uses, including:

- Identifying channels on model display and plots
- Extracting subsystems of MIMO systems
- Specifying connection points when interconnecting models

**Default:** '' for all input channels

## InputUnit

Input channel units, specified as one of the following:

- Character vector — For single-input models, for example, 'seconds'.
- Cell array of character vectors — For multi-input models.

Use `InputUnit` to keep track of input signal units. `InputUnit` has no effect on system behavior.

**Default:** '' for all input channels

## InputGroup

Input channel groups. The `InputGroup` property lets you assign the input channels of MIMO systems into groups and refer to each group by name. Specify input groups as a

structure. In this structure, field names are the group names, and field values are the input channels belonging to each group. For example:

```
sys.InputGroup.controls = [1 2];  
sys.InputGroup.noise = [3 5];
```

creates input groups named `controls` and `noise` that include input channels 1, 2 and 3, 5, respectively. You can then extract the subsystem from the `controls` inputs to all outputs using:

```
sys(:, 'controls')
```

**Default:** Struct with no fields

### OutputName

Output channel names, specified as one of the following:

- Character vector — For single-output models. For example, `'measurements'`.
- Cell array of character vectors — For multi-output models.

Alternatively, use automatic vector expansion to assign output names for multi-output models. For example, if `sys` is a two-output model, enter:

```
sys.OutputName = 'measurements';
```

The output names automatically expand to `{'measurements(1)'; 'measurements(2)'}`.

When you estimate a model using an `iddata` object, `data`, the software automatically sets `OutputName` to `data.OutputName`.

You can use the shorthand notation `y` to refer to the `OutputName` property. For example, `sys.y` is equivalent to `sys.OutputName`.

Output channel names have several uses, including:

- Identifying channels on model display and plots
- Extracting subsystems of MIMO systems
- Specifying connection points when interconnecting models

**Default:** `''` for all output channels

## OutputUnit

Output channel units, specified as one of the following:

- Character vector — For single-output models. For example, 'seconds'.
- Cell array of character vectors — For multi-output models.

Use `OutputUnit` to keep track of output signal units. `OutputUnit` has no effect on system behavior.

**Default:** '' for all output channels

## OutputGroup

Output channel groups. The `OutputGroup` property lets you assign the output channels of MIMO systems into groups and refer to each group by name. Specify output groups as a structure. In this structure, field names are the group names, and field values are the output channels belonging to each group. For example:

```
sys.OutputGroup.temperature = [1];
sys.InputGroup.measurement = [3 5];
```

creates output groups named `temperature` and `measurement` that include output channels 1, and 3, 5, respectively. You can then extract the subsystem from all inputs to the `measurement` outputs using:

```
sys('measurement', :)
```

**Default:** Struct with no fields

### Name

System name, specified as a character vector. For example, 'system\_1'.

**Default:** ''

### Notes

Any text that you want to associate with the system, stored as a string or a cell array of character vectors. The property stores whichever data type you provide. For instance, if `sys1` and `sys2` are dynamic system models, you can set their `Notes` properties as follows:

```
sys1.Notes = "sys1 has a string.";
sys2.Notes = 'sys2 has a character vector.';
sys1.Notes
sys2.Notes
```

```
ans =

    "sys1 has a string."
```

```
ans =

    'sys2 has a character vector.'
```

**Default:** [0×1 string]

### **UserData**

Any type of data you want to associate with system, specified as any MATLAB data type.

**Default:** []

### **SamplingGrid**

Sampling grid for model arrays, specified as a data structure.

For arrays of identified linear (IDLTI) models that are derived by sampling one or more independent variables, this property tracks the variable values associated with each model. This information appears when you display or plot the model array. Use this information to trace results back to the independent variables.

Set the field names of the data structure to the names of the sampling variables. Set the field values to the sampled variable values associated with each model in the array. All sampling variables should be numeric and scalar valued, and all arrays of sampled values should match the dimensions of the model array.

For example, if you collect data at various operating points of a system, you can identify a model for each operating point separately and then stack the results together into a single system array. You can tag the individual models in the array with information regarding the operating point:

```
nominal_engine_rpm = [1000 5000 10000];  
sys.SamplingGrid = struct('rpm', nominal_engine_rpm)
```

where `sys` is an array containing three identified models obtained at rpms 1000, 5000 and 10000, respectively.

For model arrays generated by linearizing a Simulink model at multiple parameter values or operating points, the software populates `SamplingGrid` automatically with the variable values that correspond to each entry in the array. For example, the Simulink Control Design commands `linearize` and `sLinearizer` populate `SamplingGrid` in this way.

**Default:** []

## Tips

- Although `idpoly` supports continuous-time models, `idtf` and `idproc` enable more choices for estimation of continuous-time models. Therefore, for some continuous-time applications, these model types are preferable.

## See Also

[ar](#) | [armax](#) | [arx](#) | [bj](#) | [idproc](#) | [idss](#) | [idtf](#) | [iv4](#) | [ivar](#) | [oe](#) | [polydata](#) | [polyest](#) | [setPolyFormat](#) | [translatecov](#)

## Topics

[“Estimate Polynomial Models in the App”](#)

[“Estimate Polynomial Models at the Command Line”](#)

[“Polynomial Sizes and Orders of Multi-Output Polynomial Models”](#)

[“What Are Polynomial Models?”](#)

[“Dynamic System Models”](#)

**Introduced before R2006a**

## idproc

Continuous-time process model with identifiable parameters

### Syntax

```
sys = idproc(type)
sys = idproc(type,Name,Value)
```

### Description

`sys = idproc(type)` creates a continuous-time process model with identifiable parameters. `type` specifies aspects of the model structures, such as the number of poles in the model, whether the model includes an integrator, and whether the model includes a time delay.

`sys = idproc(type,Name,Value)` creates a process model with additional attributes specified by one or more `Name, Value` pair arguments.

### Object Description

An `idproc` model represents a system as a continuous-time process model with identifiable (estimable) coefficients.

A simple SISO process model has a gain, a time constant, and a delay:

$$sys = \frac{K_p}{1 + T_{p1}s} e^{-T_d s}.$$

$K_p$  is a proportional gain.  $K_{p1}$  is the time constant of the real pole, and  $T_d$  is the transport delay (dead time).

More generally, `idproc` can represent process models with up to three poles and a zero:

$$\text{sys} = K_p \frac{1 + T_z s}{(1 + T_{p1} s)(1 + T_{p2} s)(1 + T_{p3} s)} e^{-T_d s}.$$

Two of the poles can be a complex conjugate (underdamped) pair. In that case, the general form of the process model is:

$$\text{sys} = K_p \frac{1 + T_z s}{(1 + 2\zeta T_\omega s + (T_\omega s)^2)(1 + T_{p3} s)} e^{-T_d s}.$$

$T_\omega$  is the time constant of the complex pair of poles, and  $\zeta$  is the associated damping constant.

In addition, any `idproc` model can have an integrator. For example, the following is a process model that you can represent with `idproc`:

$$\text{sys} = K_p \frac{1}{s(1 + 2\zeta T_\omega s + (T_\omega s)^2)} e^{-T_d s}.$$

This model has no zero ( $T_z = 0$ ). The model has a complex pair of poles. The model also has an integrator, represented by the  $1/s$  term.

For `idproc` models, all the time constants, the delay, the proportional gain, and the damping coefficient can be estimable parameters. The `idproc` model stores the values of these parameters in properties of the model such as `Kp`, `Tp1`, and `Zeta`. (See “Properties” on page 1-706 for more information.)

A MIMO process model contains a SISO process model corresponding to each input-output pair in the system. For `idproc` models, the form of each input-output pair can be independently specified. For example, a two-input, one-output process can have one channel with two poles and no zero, and another channel with a zero, a pole, and an integrator. All the coefficients are independently estimable parameters.

There are two ways to obtain an `idproc` model:

- Estimate the `idproc` model based on output or input-output measurements of a system, using the `procest` command. `procest` estimates the values of the free parameters such as gain, time constants, and time delay. The estimated values are stored as properties of the resulting `idproc` model. For example, the properties

`sys.Tz` and `sys.Kp` of an `idproc` model `sys` store the zero time constant and the proportional gain, respectively. (See “Properties” on page 1-706 for more information.) The `Report` property of the resulting model stores information about the estimation, such as handling of initial conditions and options used in estimation.

When you obtain an `idproc` model by estimation, you can extract estimated coefficients and their uncertainties from the model using commands such as `getpar` and `getcov`.

- Create an `idproc` model using the `idproc` command.

You can create an `idproc` model to configure an initial parameterization for estimation of a process model. When you do so, you can specify constraints on the parameters. For example, you can fix the values of some coefficients, or specify minimum or maximum values for the free coefficients. You can then use the configured model as an input argument to `procest` to estimate parameter values with those constraints.

## Examples

### Create SISO Process Model with Complex Poles and Time Delay

Create a process model with a pair of complex poles and a time delay. Set the initial value of the model to the following:

$$sys = \frac{0.01}{1 + 2(0.1)(10)s + (10s)^2} e^{-5s}$$

Create a process model with the specified structure.

```
sys = idproc('P2DU')
```

```
sys =
```

```
Process model with transfer function:
```

$$G(s) = \frac{Kp}{1+2*Zeta*Tw*s+(Tw*s)^2} * \exp(-Td*s)$$

```
Kp = NaN
```

```
Tw = NaN
```

```
Zeta = NaN
Td = NaN
```

Parameterization:

```
'P2DU'
Number of free coefficients: 4
Use "getpvec", "getcov" for parameters and their uncertainties.
```

Status:

Created by direct construction or transformation. Not estimated.

The input 'P2DU' specifies an underdamped pair of poles and a time delay. The display shows that `sys` has the desired structure. The display also shows that the four free parameters, `Kp`, `Tw`, `Zeta`, and `Td` are all initialized to `NaN`.

Set the initial values of all parameters to the desired values.

```
sys.Kp = 0.01;
sys.Tw = 10;
sys.Zeta = 0.1;
sys.Td = 5;
```

You can use `sys` to specify this parametrization and these initial guesses for process model estimation with `procest`.

## Create a MIMO Process Model

Create a one-input, three-output process model, where each channel has two real poles and a zero, but only the first channel has a time delay, and only the first and third channels have an integrator.

```
type = {'P2ZDI'; 'P2Z'; 'P2ZI'};
sys = idproc(type)
```

```
sys =
```

Process model with 3 outputs:  $y_k = G_k(s)u$

From input 1 to output 1:

$$G_1(s) = K_p * \frac{1+T_z*s}{s(1+T_{p1}*s)(1+T_{p2}*s)} * \exp(-T_d*s)$$

```
Kp = NaN
```

Tp1 = NaN  
Tp2 = NaN  
Td = NaN  
Tz = NaN

From input 1 to output 2:

$$G1(s) = Kp * \frac{1+Tz*s}{(1+Tp1*s)(1+Tp2*s)}$$

Kp = NaN  
Tp1 = NaN  
Tp2 = NaN  
Tz = NaN

From input 1 to output 3:

$$G1(s) = Kp * \frac{1+Tz*s}{s(1+Tp1*s)(1+Tp2*s)}$$

Kp = NaN  
Tp1 = NaN  
Tp2 = NaN  
Tz = NaN

Parameterization:

'P2DIZ'  
'P2Z'  
'P2IZ'

Number of free coefficients: 13

Use "getpvec", "getcov" for parameters and their uncertainties.

Status:

Created by direct construction or transformation. Not estimated.

`idproc` creates a MIMO model where each character vector in the `type` array defines the structure of the corresponding I/O pair. Since `type` is a column vector of character vectors, `sys` is a one-input, three-output model having the specified parametrization structure. `type{k, 1}` specifies the structure of the subsystem `sys(k, 1)`. All identifiable parameters are initialized to `NaN`.

## Create Array of Process Models

Create a 3-by-1 array of process models, each containing one output and two input channels.

Specify the structure for each model in the array of process models.

```
type1 = {'P1D', 'P2DZ'};
type2 = {'P0', 'P3UI'};
type3 = {'P2D', 'P2DI'};
type = cat(3,type1,type2,type3);
size(type)
```

```
ans = 1×3
```

```
    1    2    3
```

Use `type` to create the array.

```
sysarr = idproc(type);
```

The first two dimensions of the cell array `type` set the output and input dimensions of each model in the array of process models. The remaining dimensions of the cell array set the array dimensions. Thus, `sysarr` is a 3-model array of 2-input, one-output process models.

Select a model from the array.

```
sysarr(:,:,2)
```

```
ans =
```

```
Process model with 2 inputs: y = G11(s)u1 + G12(s)u2
```

```
From input 1 to output 1:
```

```
G11(s) = Kp
```

```
Kp = NaN
```

```
From input 2 to output 1:
```

```
G12(s) = -----
                Kp
            s(1+2*Zeta*Tw*s+(Tw*s)^2)(1+Tp3*s)
```

```
Kp = NaN
```

```
Tw = NaN
```

```
Zeta = NaN  
Tp3 = NaN
```

```
Parameterization:  
'P0' 'P3IU'
```

```
Number of free coefficients: 5
```

```
Use "getpvec", "getcov" for parameters and their uncertainties.
```

```
Status:
```

```
Created by direct construction or transformation. Not estimated.
```

This two-input, one-output model corresponds to the `type2` entry in the `type` cell array.

## Input Arguments

### **type**

Model structure, specified as a character vector or cell array of character vectors.

For SISO models, `type` is a character vector made up of one or more of the following characters that specify aspects of the model structure:

Characters	Meaning
Pk	A process model with $k$ poles (not including an integrator). $k$ must be 0, 1, 2, or 3.
Z	The process model includes a zero ( $T_z \neq 0$ ). A <code>type</code> with P0 cannot include Z (a process model with no poles cannot include a zero).
D	The process model includes a time delay (deadtime) ( $T_d \neq 0$ ).
I	The process model includes an integrator (1/s).
U	The process model is underdamped. In this case, the process model includes a complex pair of poles

Every `type` character vector must begin with one of P0, P1, P2, or P3. All other components are optional. For example:

- 'P1D' specifies a process model with one pole and a time delay (deadtime) term:

$$\text{sys} = \frac{K_p}{1 + T_{p1}s} e^{-T_d s}.$$

$K_p$ ,  $T_{p1}$ , and  $T_d$  are the identifiable parameters of this model.

- 'P2U' creates a process model with a pair of complex poles:

$$\text{sys} = \frac{K_p}{\left(1 + 2\zeta T_\omega s + (T_\omega s)^2\right)}.$$

$K_p$ ,  $T_\omega$ , and  $\zeta$  are the identifiable parameters of this model.

- 'P3ZDI' creates a process model with three poles. All poles are real, because  $U$  is not included. The model also includes a zero, a time delay, and an integrator:

$$\text{sys} = K_p \frac{1 + T_z s}{s(1 + T_{p1}s)(1 + T_{p2}s)(1 + T_{p3}s)} e^{-T_d s}.$$

The identifiable parameters of this model are  $K_p$ ,  $T_z$ ,  $T_{p1}$ ,  $T_{p2}$ ,  $T_{p3}$ , and  $T_d$ .

The values of all parameters in a particular model structure are initialized to NaN. You can change them to finite values by setting the values of the corresponding `idproc` model properties after you create the model. For example, `sys.Td = 5` sets the initial value of the time delay of `sys` to 5.

For a MIMO process model with  $N_y$  outputs and  $N_u$  inputs, `type` is an  $N_y$ -by- $N_u$  cell array of character vectors specifying the structure of each input/output pair in the model. For example, `type{i, j}` specifies the type of the subsystem `sys(i, j)` from the  $j$ th input to the  $i$ th output.

## Name-Value Pair Arguments

Specify optional comma-separated pairs of `Name`, `Value` arguments. `Name` is the argument name and `Value` is the corresponding value. `Name` must appear inside quotes. You can specify several name and value pair arguments in any order as `Name1, Value1, ..., NameN, ValueN`.

Use `Name`, `Value` arguments to specify parameter initial values and additional properties on page 1-706 of `idproc` models during model creation. For example, `sys =`

`idproc('p2z','InputName','Voltage','Kp',10,'Tz',0);` creates an `idproc` model with the `InputName` property set to `Voltage`. The command also initializes the parameter `Kp` to a value of 10, and `Tz` to 0.

## Properties

`idproc` object properties include:

### Type

Model structure, specified as a character vector or cell array of character vectors.

For a SISO model `sys`, the property `sys.Type` contains a character vector specifying the structure of the system. For example, `'PID'`.

For a MIMO model with `Ny` outputs and `Nu` inputs, `sys.Type` is an `Ny`-by-`Nu` cell array of character vectors specifying the structure of each input/output pair in the model. For example, `type{i,j}` specifies the structure of the subsystem `sys(i,j)` from the `j`th input to the `i`th output.

The character vectors are made up of one or more of the following characters that specify aspects of the model structure:

Characters	Meaning
Pk	A process model with $k$ poles (not including an integrator). $k$ is 0, 1, 2, or 3.
Z	The process model includes a zero ( $T_z \neq 0$ ).
D	The process model includes a time delay (deadtime) ( $T_d \neq 0$ ).
I	The process model includes an integrator (1/s).
U	The process model is underdamped. In this case, the process model includes a complex pair of poles

If you create an `idproc` model `sys` using the `idproc` command, `sys.Type` contains the model structure that you specify with the `type` input argument.

If you obtain an `idproc` model by identification using `procest`, then `sys.Type` contains the model structures that you specified for that identification.

In general, you cannot change the type of an existing model. However, you can change whether the model contains an integrator using the property `sys.Integration`.

### **Kp, Tp1, Tp2, Tp3, Tz, Tw, Zeta, Td**

Values of process model parameters.

If you create an `idproc` model using the `idproc` command, the values of all parameters present in the model structure initialize by default to `NaN`. The values of parameters not present in the model structure are fixed to `0`. For example, if you create a model, `sys`, of type 'PID', then `Kp`, `Tp1`, and `Td` are initialized to `NaN` and are identifiable (free) parameters. All remaining parameters, such as `Tp2` and `Tz`, are inactive in the model. The values of inactive parameters are fixed to zero and cannot be changed.

For a MIMO model with `Ny` outputs and `Nu` inputs, each parameter value is an `Ny-by-Nu` cell array of character vectors specifying the corresponding parameter value for each input/output pair in the model. For example, `sys.Kp(i,j)` specifies the `Kp` value of the subsystem `sys(i,j)` from the `j`th input to the `i`th output.

For an `idproc` model `sys`, each parameter value property such as `sys.Kp`, `sys.Tp1`, `sys.Tz`, and the others is an alias to the corresponding `Value` entry in the `Structure` property of `sys`. For example, `sys.Tp3` is an alias to the value of the property `sys.Structure.Tp3.Value`.

**Default:** For each parameter value, `NaN` if the process model structure includes the particular parameter; `0` if the structure does not include the parameter.

### **Integration**

Logical value or matrix denoting the presence or absence of an integrator in the transfer function of the process model.

For a SISO model `sys`, `sys.Integration = true` if the model contains an integrator.

For a MIMO model, `sys.Integration(i,j) = true` if the transfer function from the `j`th input to the `i`th output contains an integrator.

When you create a process model using the `idproc` command, the value of `sys.Integration` is determined by whether the corresponding `type` contains `I`.

### **NoiseTF**

Coefficients of the noise transfer function.

`sys.NoiseTF` stores the coefficients of the numerator and the denominator polynomials for the noise transfer function  $H(s) = N(s)/D(s)$ .

`sys.NoiseTF` is a structure with fields `num` and `den`. Each field is a cell array of  $N_y$  row vectors, where  $N_y$  is the number of outputs of `sys`. These row vectors specify the coefficients of the noise transfer function numerator and denominator in order of decreasing powers of  $s$ .

Typically, the noise transfer function is automatically computed by the estimation function `procest`. You can specify a noise transfer function that `procest` uses as an initial value. For example:

```
NoiseNum = {[1 2.2]; [1 0.54]};  
NoiseDen = {[1 1.3]; [1 2]};  
NoiseTF = struct('num', {NoiseNum}, 'den', {NoiseDen});  
sys = idproc({'p2'; 'pldi'}); % 2-output, 1-input process model  
sys.NoiseTF = NoiseTF;
```

Each vector in `sys.NoiseTF.num` and `sys.NoiseTF.den` must be of length 3 or less (second-order in  $s$  or less). Each vector must start with 1. The length of a numerator vector must be equal to that of the corresponding denominator vector, so that  $H(s)$  is always biproper.

**Default:** `struct('num', {num2cell(ones(Ny,1))}, 'den', {num2cell(ones(Ny,1))})`

## Structure

Information about the estimable parameters of the `idproc` model.

`sys.Structure` includes one entry for each parameter in the model structure of `sys`. For example, if `sys` is of type 'PID', then `sys` includes identifiable parameters `Kp`, `Tp1`, and `Td`. Correspondingly, `sys.Structure.Kp`, `sys.Structure.Tp1`, and `sys.Structure.Td` contain information about each of these parameters, respectively.

Each of these parameter entries in `sys.Structure` contains the following fields:

- **Value** — Parameter values. For example, `sys.Structure.Kp.Value` contains the initial or estimated values of the  $K_p$  parameter.

`NaN` represents unknown parameter values.

For SISO models, each parameter value property such as `sys.Kp`, `sys.Tp1`, `sys.Tz`, and the others is an alias to the corresponding `Value` entry in the `Structure`

property of `sys`. For example, `sys.Tp3` is an alias to the value of the property `sys.Structure.Tp3.Value`.

For MIMO models, `sys.Kp{i,j}` is an alias to `sys.Structure(i,j).Kp.Value`, and similarly for the other identifiable coefficient values.

- **Minimum** — Minimum value that the parameter can assume during estimation. For example, `sys.Structure.Kp.Minimum = 1` constrains the proportional gain to values greater than or equal to 1.
- **Maximum** — Maximum value that the parameter can assume during estimation.
- **Free** — Logical value specifying whether the parameter is a free estimation variable. If you want to fix the value of a parameter during estimation, set the corresponding `Free = false`. For example, to fix the dead time to 5:

```
sys.Td = 5;
sys.Structure.Td.Free = false;
```

- **Scale** — Scale of the parameter's value. `Scale` is not used in estimation.
- **Info** — Structure array for storing parameter units and labels. The structure has `Label` and `Unit` fields.

Specify parameter units and labels as character vectors. For example, `'Time'`.

`Structure` also includes a field `Integration` that stores a logical array indicating whether each corresponding process model has an integrator. `sys.Structure.Integration` is an alias to `sys.Integration`.

For a MIMO model with  $N_y$  outputs and  $N_u$  input, `Structure` is an  $N_y$ -by- $N_u$  array. The element `Structure(i,j)` contains information corresponding to the process model for the  $(i,j)$  input-output pair.

## NoiseVariance

The variance (covariance matrix) of the model innovations  $e$ .

An identified model includes a white, Gaussian noise component  $e(t)$ . `NoiseVariance` is the variance of this noise component. Typically, the model estimation function (such as `procest`) determines this variance.

For SISO models, `NoiseVariance` is a scalar. For MIMO models, `NoiseVariance` is a  $N_y$ -by- $N_y$  matrix, where  $N_y$  is the number of outputs in the system.

## Report

Summary report that contains information about the estimation options and results when the process model is obtained using the `procest` estimation command. Use `Report` to query a model for how it was estimated, including its:

- Estimation method
- Estimation options
- Search termination conditions
- Estimation data fit and other quality metrics

The contents of `Report` are irrelevant if the model was created by construction.

```
m = idproc('P2DU');  
m.Report.OptionsUsed
```

```
ans =
```

```
 []
```

If you obtain the process model using estimation commands, the fields of `Report` contain information on the estimation data, options, and results.

```
load iddata2 z2;  
m = procest(z2, 'P2DU');  
m.Report.OptionsUsed
```

```
DisturbanceModel: 'estimate'  
  InitialCondition: 'auto'  
        Focus: 'prediction'  
EstimateCovariance: 1  
  Display: 'off'  
  InputOffset: [1x1 param.Continuous]  
  OutputOffset: []  
Regularization: [1x1 struct]  
SearchMethod: 'auto'  
SearchOptions: [1x1 idoptions.search.identsolver]  
OutputWeight: []  
Advanced: [1x1 struct]
```

`Report` is a read-only property.

For more information on this property and how to use it, see the Output Arguments section of the corresponding estimation command reference page and “Estimation Report”.

### **InputDelay**

Input delays. **InputDelay** is a numeric vector specifying a time delay for each input channel. Specify input delays in the time unit stored in the **TimeUnit** property.

For a system with **Nu** inputs, set **InputDelay** to an **Nu**-by-1 vector, where each entry is a numerical value representing the input delay for the corresponding input channel. You can also set **InputDelay** to a scalar value to apply the same delay to all channels.

**Default:** 0 for all input channels

### **OutputDelay**

Output delays.

For identified systems, like **idproc**, **OutputDelay** is fixed to zero.

### **Ts**

Sample time. For **idproc**, **Ts** is fixed to zero because all **idproc** models are continuous time.

### **TimeUnit**

Units for the time variable, the sample time **Ts**, and any time delays in the model, specified as one of the following values:

- 'nanoseconds'
- 'microseconds'
- 'milliseconds'
- 'seconds'
- 'minutes'
- 'hours'
- 'days'
- 'weeks'

- 'months'
- 'years'

Changing this property has no effect on other properties, and therefore changes the overall system behavior. Use `chgTimeUnit` to convert between time units without modifying system behavior.

**Default:** 'seconds'

### **InputName**

Input channel names, specified as one of the following:

- Character vector — For single-input models, for example, 'controls'.
- Cell array of character vectors — For multi-input models.

Alternatively, use automatic vector expansion to assign input names for multi-input models. For example, if `sys` is a two-input model, enter:

```
sys.InputName = 'controls';
```

The input names automatically expand to {'controls(1)'; 'controls(2)'}

When you estimate a model using an `iddata` object, `data`, the software automatically sets `InputName` to `data.InputName`.

You can use the shorthand notation `u` to refer to the `InputName` property. For example, `sys.u` is equivalent to `sys.InputName`.

Input channel names have several uses, including:

- Identifying channels on model display and plots
- Extracting subsystems of MIMO systems
- Specifying connection points when interconnecting models

**Default:** '' for all input channels

### **InputUnit**

Input channel units, specified as one of the following:

- Character vector — For single-input models, for example, 'seconds'.
- Cell array of character vectors — For multi-input models.

Use `InputUnit` to keep track of input signal units. `InputUnit` has no effect on system behavior.

**Default:** '' for all input channels

### InputGroup

Input channel groups. The `InputGroup` property lets you assign the input channels of MIMO systems into groups and refer to each group by name. Specify input groups as a structure. In this structure, field names are the group names, and field values are the input channels belonging to each group. For example:

```
sys.InputGroup.controls = [1 2];
sys.InputGroup.noise = [3 5];
```

creates input groups named `controls` and `noise` that include input channels 1, 2 and 3, 5, respectively. You can then extract the subsystem from the `controls` inputs to all outputs using:

```
sys(:, 'controls')
```

**Default:** Struct with no fields

### OutputName

Output channel names, specified as one of the following:

- Character vector — For single-output models. For example, 'measurements'.
- Cell array of character vectors — For multi-output models.

Alternatively, use automatic vector expansion to assign output names for multi-output models. For example, if `sys` is a two-output model, enter:

```
sys.OutputName = 'measurements';
```

The output names automatically expand to `{'measurements(1)'; 'measurements(2)'}`.

When you estimate a model using an `iddata` object, `data`, the software automatically sets `OutputName` to `data.OutputName`.

You can use the shorthand notation `y` to refer to the `OutputName` property. For example, `sys.y` is equivalent to `sys.OutputName`.

Output channel names have several uses, including:

- Identifying channels on model display and plots
- Extracting subsystems of MIMO systems
- Specifying connection points when interconnecting models

**Default:** '' for all output channels

### OutputUnit

Output channel units, specified as one of the following:

- Character vector — For single-output models. For example, 'seconds'.
- Cell array of character vectors — For multi-output models.

Use `OutputUnit` to keep track of output signal units. `OutputUnit` has no effect on system behavior.

**Default:** '' for all output channels

### OutputGroup

Output channel groups. The `OutputGroup` property lets you assign the output channels of MIMO systems into groups and refer to each group by name. Specify output groups as a structure. In this structure, field names are the group names, and field values are the output channels belonging to each group. For example:

```
sys.OutputGroup.temperature = [1];  
sys.InputGroup.measurement = [3 5];
```

creates output groups named `temperature` and `measurement` that include output channels 1, and 3, 5, respectively. You can then extract the subsystem from all inputs to the measurement outputs using:

```
sys('measurement', :)
```

**Default:** Struct with no fields

**Name**

System name, specified as a character vector. For example, 'system\_1'.

**Default:** ''

**Notes**

Any text that you want to associate with the system, stored as a string or a cell array of character vectors. The property stores whichever data type you provide. For instance, if `sys1` and `sys2` are dynamic system models, you can set their `Notes` properties as follows:

```
sys1.Notes = "sys1 has a string.";
sys2.Notes = 'sys2 has a character vector.';
sys1.Notes
sys2.Notes
```

```
ans =
```

```
    "sys1 has a string."
```

```
ans =
```

```
    'sys2 has a character vector.'
```

**Default:** [0×1 string]

**UserData**

Any type of data you want to associate with system, specified as any MATLAB data type.

**Default:** []

**SamplingGrid**

Sampling grid for model arrays, specified as a data structure.

For arrays of identified linear (IDLTI) models that are derived by sampling one or more independent variables, this property tracks the variable values associated with each

model. This information appears when you display or plot the model array. Use this information to trace results back to the independent variables.

Set the field names of the data structure to the names of the sampling variables. Set the field values to the sampled variable values associated with each model in the array. All sampling variables should be numeric and scalar valued, and all arrays of sampled values should match the dimensions of the model array.

For example, if you collect data at various operating points of a system, you can identify a model for each operating point separately and then stack the results together into a single system array. You can tag the individual models in the array with information regarding the operating point:

```
nominal_engine_rpm = [1000 5000 10000];  
sys.SamplingGrid = struct('rpm', nominal_engine_rpm)
```

where `sys` is an array containing three identified models obtained at rpms 1000, 5000 and 10000, respectively.

For model arrays generated by linearizing a Simulink model at multiple parameter values or operating points, the software populates `SamplingGrid` automatically with the variable values that correspond to each entry in the array. For example, the Simulink Control Design commands `linearize` and `sLinearizer` populate `SamplingGrid` in this way.

**Default:** []

## See Also

`idss` | `idtf` | `pem` | `procest` | `ssest` | `tfest`

**Introduced before R2006a**

# idresamp

Resample time-domain data by decimation or interpolation

## Syntax

```
datar = idresamp(data,R)
datar = idresamp(data,R,order,tol)
[datar,res_fact] = idresamp(data,R,order,tol)
```

## Description

`datar = idresamp(data,R)` resamples data on a new sample interval  $R$  and stores the resampled data as `datar`.

`datar = idresamp(data,R,order,tol)` filters the data by applying a filter of specified `order` before interpolation and decimation. Replaces  $R$  by a rational approximation that is accurate to a tolerance `tol`.

`[datar,res_fact] = idresamp(data,R,order,tol)` returns `res_fact`, which corresponds to the value of  $R$  approximated by a rational expression.

## Input Arguments

`data`

Name of time-domain `iddata` object or a matrix of data. Can be input-output or time-series data.

Data must be sampled at equal time intervals.

`R`

Resampling factor, such that  $R > 1$  results in decimation and  $R < 1$  results in interpolation.

Any positive number you specify is replaced by the rational approximation,  $Q/P$ .

`order`

Order of the filters applied before interpolation and decimation.

Default: 8

`tol`

Tolerance of the rational approximation for the resampling factor R.

Smaller tolerance might result in larger P and Q values, which produces more accurate answers at the expense of slower computation.

Default: 0.1

## Output Arguments

`datar`

Name of the resampled data variable. `datar` class matches the `data` class, as specified.

`res_fact`

Rational approximation for the specified resampling factor R and tolerance `tol`.

Any positive number you specify is replaced by the rational approximation, Q/P, where the data is interpolated by a factor P and then decimated by a factor Q.

## See Also

`resample`

**Introduced in R2007a**

## idss

State-space model with identifiable parameters

### Syntax

```
sys = idss(A,B,C,D)
sys = idss(A,B,C,D,K)
sys = idss(A,B,C,D,K,x0)
sys = idss(A,B,C,D,K,x0,Ts)
sys = idss( ____,Name,Value)

sys = idss(sys0)
sys = idss(sys0,'split')
```

### Description

`sys = idss(A,B,C,D)` creates a state-space model with identifiable parameters. *A*, *B*, *C*, and *D* are the initial values of the state-space matrices. By default, `sys` is discrete-time model with unspecified sample time and no state disturbance element.

`sys = idss(A,B,C,D,K)` creates a state-space model with a disturbance element given by the matrix *K*.

`sys = idss(A,B,C,D,K,x0)` creates a state-space model with initial state values given by the vector *x0*.

`sys = idss(A,B,C,D,K,x0,Ts)` creates a state-space model with sample time *Ts*. Use *Ts* = 0 to create a continuous-time model.

`sys = idss( ____,Name,Value)` creates a state-space model using additional options specified by one or more *Name,Value* pair arguments.

`sys = idss(sys0)` converts any dynamic system model, `sys0`, to `idss` model form.

`sys = idss(sys0,'split')` converts `sys0` to `idss` model form, and treats the last *Ny* input channels of `sys0` as noise channels in the returned model. `sys0` must be a numeric

(non-identified) `tf`, `zpk`, or `ss` model object. Also, `sys0` must have at least as many inputs as outputs.

## Object Description

An `idss` model represents a system as a continuous-time or discrete-time state-space model with identifiable (estimable) coefficients.

A state-space model of a system with input vector  $u$ , output vector  $y$ , and disturbance  $e$  takes the following form in continuous time:

$$\begin{aligned}\frac{dx(t)}{dt} &= Ax(t) + Bu(t) + Ke(t) \\ y(t) &= Cx(t) + Du(t) + e(t).\end{aligned}$$

In discrete time, the state-space model takes the form:

$$\begin{aligned}x[k+1] &= Ax[k] + Bu[k] + Ke[k] \\ y[k] &= Cx[k] + Du[k] + e[k].\end{aligned}$$

For `idss` models, the elements of the state-space matrices  $A$ ,  $B$ ,  $C$ , and  $D$  can be estimable parameters. The elements of the state disturbance  $K$  can also be estimable parameters. The `idss` model stores the values of these matrix elements in the `A`, `B`, `C`, `D`, and `K` properties of the model.

There are three ways to obtain an `idss` model.

- Estimate the `idss` model based on input-output measurements of a system, using `n4sid` or `ssest`. These estimation commands estimate the values of the estimable elements of the state-space matrices. The estimated values are stored in the `A`, `B`, `C`, `D`, and `K` properties of the resulting `idss` model. The `Report` property of the resulting model stores information about the estimation, such as handling of initial state values and options used in estimation.

When you obtain an `idss` model by estimation, you can extract estimated coefficients and their uncertainties from the model using commands such as `idssdata`, `getpar`, or `getcov`.

- Create an `idss` model using the `idss` command.

You can create an `idss` model to configure an initial parameterization for estimation of a state-space model to fit measured response data. When you do so, you can specify constraints on one or more of the state-space matrix elements. For example, you can fix the values of some elements, or specify minimum or maximum values for the free elements. You can then use the configured model as an input argument to an estimation command (`n4sid` or `ssest`) to estimate parameter values with those constraints.

- Convert an existing dynamic system model to an `idss` model using the `idss` command.

To configure an `idss` model in a desired form, such as a companion or modal form, use state transformation commands such as `canon` and `ss2ss`.

## Examples

### Create State-Space Model with Identifiable Parameters

Create a 4th-order SISO state-space model with identifiable parameters. Initialize the initial state values to 0.1 for all entries. Set the sample time to 0.1 s as well.

```
A = blkdiag([-0.1 0.4; -0.4 -0.1],[-1 5; -5 -1]);
B = [1; zeros(3,1)];
C = [1 0 1 0];
D = 0;
K = zeros(4,1);
x0 = [0.1,0.1,0.1,0.1];
Ts = 0.1;
```

```
sys = idss(A,B,C,D,K,x0,Ts);
```

`sys` is a 4th-order, SISO `idss` model. The number of states and input-output dimensions are determined by the dimensions of the state-space matrices. By default, all entries in the matrices `A`, `B`, `C`, `D`, and `K` are identifiable parameters.

You can use `sys` to specify an initial parametrization for state-space model estimation with `ssest` or `n4sid`.

### Specify Additional Attributes of State-Space Model

Create a 4th-order SISO state-space model with identifiable parameters. Name the input and output channels of the model, and specify minutes for the model time units.

You can use `Name`, `Value` pair arguments to specify additional model properties on model creation.

```
A = blkdiag([-0.1 0.4; -0.4 -0.1],[-1 5; -5 -1]);  
B = [1; zeros(3,1)];  
C = [1 0 1 0];  
D = 0;
```

```
sys = idss(A,B,C,D,'InputName','Drive','TimeUnit','minutes');
```

To change or specify most attributes of an existing model, you can use dot notation. For example:

```
sys.OutputName = 'Torque';
```

### Configure Identifiable Parameters of State-Space Model

Configure an `idss` model so that it has no state disturbance element and only the non-zero entries of the `A` matrix are estimable. Additionally, fix the values of the `B` matrix.

You can configure individual parameters of an `idss` model to specify constraints for state-space model estimation with `ssest` or `n4sid`.

Create an `idss` model.

```
A = blkdiag([-0.1 0.4; -0.4 -0.1],[-1 5; -5 -1]);  
B = [1; zeros(3,1)];  
C = [1 0 1 0];  
D = 0;  
K = zeros(4,1);  
x0 = [0.1,0.1,0.1,0.1];
```

```
sys = idss(A,B,C,D,K,x0,0);
```

Setting all entries of `K = 0` creates an `idss` model with no state disturbance element.

Use the `Structure` property of the model to fix the values of some of the parameters.

```
sys.Structure.A.Free = (A~=0);
sys.Structure.B.Free = false;
sys.Structure.K.Free = false;
```

The entries in `sys.Structure.A.Free` determine whether the corresponding entries in `sys.A` are free (identifiable) or fixed. The first line sets `sys.Structure.A.Free` to a logical matrix that is `true` wherever `A` is non-zero, and `false` everywhere else. Doing so fixes the value of the zero entries in `sys.A`.

The remaining lines fix all the values in `sys.B` and `sys.K` to the values you specified when you created the model.

## Array of State-Space Models

Create an array of state-space models.

There are several ways to create arrays of state-space models:

- Direct array construction using  $n$ -dimensional state-space arrays
- Array-building by indexed assignment
- Array-building using the `stack` command
- Sampling an identified model using the `rsample` command

Create an array by providing  $n$ -dimensional arrays as an input argument to `idss`, instead of 2-dimensional matrices.

```
A = rand(2,2,3,4);
sysarr = idss(A,[2;1],[1 1],0);
```

When you provide a multi-dimensional array to `idss` in place of one of the state-space matrices, the first two dimensions specify the numbers of states, inputs, or outputs of each model in the array. The remaining dimensions specify the dimensions of the array itself. `A` is a 2-by-2-by-3-by-4 array. Therefore, `sysarr` is a 3-by-4 array of `idss` models. Each model in `sysarr` has two states, specified by the first two dimensions of `A`. Further, each model in `sysarr` has the same `B`, `C`, and `D` values.

Create an array by indexed assignment.

```
sysarr = idss(zeros(1,1,2));  
sysarr(:,:,1) = idss([4 -3; -2 0],[2;1],[1 1],0);  
sysarr(:,:,2) = idss(rand(2),rand(2,1),rand(1,2),1);
```

The first command preallocates the array. The first two dimensions of the array are the I/O dimensions of each model in the array. Therefore, `sysarr` is a 2-element vector of SISO models.

The remaining commands assign an `idss` model to each position in `sysarr`. Each model in an array must have the same I/O dimensions.

Add another model to `sysarr` using `stack`.

`stack` is an alternative to building an array by indexing.

```
sysarr = stack(1,sysarr,idss([1 -2; -4 9],[0;-1],[1 1],0));
```

This command adds another `idss` model along the first array dimension of `sysarr`. `sysarr` is now a 3-by-1 array of SISO `idss` models

## Input Arguments

### A, B, C, D

Initial values of the state-space matrices.

For a system with  $N_y$  outputs,  $N_u$  inputs, and  $N_x$  states, specify initial values of the state-space matrix elements as follows:

- A —  $N_x$ -by- $N_x$  matrix.
- B —  $N_x$ -by- $N_u$  matrix.
- C —  $N_y$ -by- $N_x$  matrix.
- D —  $N_y$ -by- $N_u$  matrix.

Use NaN for any matrix element whose initial value is not known.

### K

Initial value of the state disturbance matrix.

Specify  $K$  as an  $N_x$ -by- $N_y$  matrix.

Use NaN for any matrix element whose initial value is not known.

**Default:**  $N_x$ -by- $N_y$  zero matrix.

### **x0**

Initial state values.

Specify the initial condition as a column vector of  $N_x$  values.

**Default:**  $N_x$  column vector of zeros.

### **Ts**

Sample time. For continuous-time models,  $T_s = 0$ . For discrete-time models,  $T_s$  is a positive scalar representing the sampling period expressed in the unit specified by the `TimeUnit` property of the model. To denote a discrete-time model with unspecified sample time, set  $T_s = -1$ .

**Default:** -1 (discrete-time model with unspecified sample time)

### **sys0**

Dynamic system.

Any dynamic system to convert to an `idss` model:

- When `sys0` is an identified model, its estimated parameter covariance is lost during conversion. If you want to translate the estimated parameter covariance during the conversion, use `translatecov`.
- When `sys0` is a numeric (non-identified) model, the state-space data of `sys0` define the  $A$ ,  $B$ ,  $C$ , and  $D$  matrices of the converted model. The disturbance matrix  $K$  is fixed to zero. The `NoiseVariance` value defaults to `eye(Ny)`, where  $N_y$  is the number of outputs of `sys`.

For the syntax `sys = idss(sys0, 'split')`, `sys0` must be a numeric (non-identified) `tf`, `zpk`, or `ss` model object. Also, `sys0` must have at least as many inputs as outputs. Finally, the subsystem `sys0(:, Ny+1:Ny+Nu)` must contain a non-zero feedthrough term (the subsystem must be biproper).

## Name-Value Pair Arguments

Specify optional comma-separated pairs of `Name`, `Value` arguments. `Name` is the argument name and `Value` is the corresponding value. `Name` must appear inside quotes. You can specify several name and value pair arguments in any order as `Name1, Value1, . . . , NameN, ValueN`.

Use `Name`, `Value` arguments to specify additional properties on page 1-726 of `idss` models during model creation. For example, `idss(A,B,C,D, 'InputName', 'Voltage')` creates an `idss` model with the `InputName` property set to `Voltage`.

## Properties

Specify optional comma-separated pairs of `Name`, `Value` arguments. `Name` is the argument name and `Value` is the corresponding value. `Name` must appear inside quotes. You can specify several name and value pair arguments in any order as `Name1, Value1, . . . , NameN, ValueN`.

`idss` object properties include:

### **A, B, C, D**

Values of state-space matrices.

- `A` — State matrix `A`, an  $N_x$ -by- $N_x$  matrix.
- `B` —  $N_x$ -by- $N_u$  matrix.
- `C` —  $N_y$ -by- $N_x$  matrix.
- `D` —  $N_y$ -by- $N_u$  matrix.

If you create an `idss` model `sys` using the `idss` command, `sys.A`, `sys.B`, `sys.C`, and `sys.D` contain the initial values of the state-space matrices that you specify with the `A, B, C, D` input arguments.

If you obtain an `idss` model `sys` by identification using `ssest` or `n4sid`, then `sys.A`, `sys.B`, `sys.C`, and `sys.D` contain the estimated values of the matrix elements.

For an `idss` model `sys`, each property `sys.A`, `sys.B`, `sys.C`, and `sys.D` is an alias to the corresponding `Value` entry in the `Structure` property of `sys`. For example, `sys.A` is an alias to the value of the property `sys.Structure.A.Value`.

## K

Value of state disturbance matrix  $K$ , an  $N_x$ -by- $N_y$  matrix.

If you create an `idss` model `sys` using the `idss` command, `sys.K` contains the initial values of the state-space matrices that you specify with the `K` input argument.

If you obtain an `idss` model `sys` by identification using `ssest` or `n4sid`, then `sys.K` contains the estimated values of the matrix elements.

For an `idss` model `sys`, `sys.K` is an alias to the value of the property `sys.Structure.K.Value`.

**Default:**  $N_x$ -by- $N_y$  zero matrix.

## StateName

State names, specified as one of the following:

- Character vector — For first-order models, for example, 'velocity'.
- Cell array of character vectors — For models with two or more states
- '' — For unnamed states.

**Default:** '' for all states

## StateUnit

State units, specified as one of the following:

- Character vector — For first-order models, for example, 'velocity'
- Cell array of character vectors — For models with two or more states
- '' — For states without specified units

Use `StateUnit` to keep track of the units each state is expressed in. `StateUnit` has no effect on system behavior.

**Default:** '' for all states

## Structure

Information about the estimable parameters of the `idss` model. `Structure.A`, `Structure.B`, `Structure.C`, `Structure.D`, and `Structure.K` contain information about the  $A$ ,  $B$ ,  $C$ ,  $D$ , and  $K$  matrices, respectively. Each contains the following fields:

- **Value** — Parameter values. For example, `sys.Structure.A.Value` contains the initial or estimated values of the  $A$  matrix.

NaN represents unknown parameter values.

Each property `sys.A`, `sys.B`, `sys.C`, and `sys.D` is an alias to the corresponding `Value` entry in the `Structure` property of `sys`. For example, `sys.A` is an alias to the value of the property `sys.Structure.A.Value`

- **Minimum** — Minimum value that the parameter can assume during estimation. For example, `sys.Structure.K.Minimum = 0` constrains all entries in the  $K$  matrix to be greater than or equal to zero.
- **Maximum** — Maximum value that the parameter can assume during estimation.
- **Free** — Boolean specifying whether the parameter is a free estimation variable. If you want to fix the value of a parameter during estimation, set the corresponding `Free = false`. For example, if  $A$  is a 3-by-3 matrix, `sys.Structure.A.Free = eyes(3)` fixes all of the off-diagonal entries in  $A$ , to the values specified in `sys.Structure.A.Value`. In this case, only the diagonal entries in  $A$  are estimable.
- **Scale** — Scale of the parameter's value. `Scale` is not used in estimation.
- **Info** — Structure array for storing parameter units and labels. The structure has `Label` and `Unit` fields.

Specify parameter units and labels as character vectors. For example, `'Time'`.

## NoiseVariance

The variance (covariance matrix) of the model innovations  $e$ .

An identified model includes a white, Gaussian noise component  $e(t)$ . `NoiseVariance` is the variance of this noise component. Typically, the model estimation function (such as `ssest`) determines this variance.

For SISO models, `NoiseVariance` is a scalar. For MIMO models, `NoiseVariance` is a  $N_y$ -by- $N_y$  matrix, where  $N_y$  is the number of outputs in the system.

## Report

Summary report that contains information about the estimation options and results when the state-space model is obtained using estimation commands, such as `ssest`, `ssregest`, and `n4sid`. Use `Report` to query a model for how it was estimated, including its:

- Estimation method
- Estimation options
- Search termination conditions
- Estimation data fit and other quality metrics

The contents of `Report` are irrelevant if the model was created by construction.

```
A = [-0.1 0.4; -0.4 -0.1];
B = [1; 0];
C = [1 0];
D = 0;
m = idss(A,B,C,D);
m.Report.OptionsUsed
```

```
ans =
```

```
 []
```

If you obtain the state-space model using estimation commands, the fields of `Report` contain information on the estimation data, options, and results.

```
load iddata2 z2;
m = ssest(z2,3);
m.Report.OptionsUsed

InitialState: 'auto'
    N4Weight: 'auto'
    N4Horizon: 'auto'
    Focus: 'prediction'
EstimateCovariance: 1
    Display: 'off'
    InputOffset: []
    OutputOffset: []
    OutputWeight: []
    SearchMethod: 'auto'
    SearchOptions: [1x1 idoptions.search.identsolver]
Regularization: [1x1 struct]
    Advanced: [1x1 struct]
```

Report is a read-only property.

For more information on this property and how to use it, see the Output Arguments section of the corresponding estimation command reference page and “Estimation Report”.

### **InputDelay**

Input delay for each input channel, specified as a scalar value or numeric vector. For continuous-time systems, specify input delays in the time unit stored in the `TimeUnit` property. For discrete-time systems, specify input delays in integer multiples of the sample time `Ts`. For example, `InputDelay = 3` means a delay of three sample times.

For a system with `Nu` inputs, set `InputDelay` to an `Nu`-by-1 vector. Each entry of this vector is a numerical value that represents the input delay for the corresponding input channel.

You can also set `InputDelay` to a scalar value to apply the same delay to all channels.

**Default:** 0

### **OutputDelay**

Output delays.

For identified systems, like `idss`, `OutputDelay` is fixed to zero.

### **Ts**

Sample time. For continuous-time models, `Ts = 0`. For discrete-time models, `Ts` is a positive scalar representing the sampling period expressed in the unit specified by the `TimeUnit` property of the model. To denote a discrete-time model with unspecified sample time, set `Ts = -1`.

Changing this property does not discretize or resample the model. Use `c2d` and `d2c` to convert between continuous- and discrete-time representations. Use `d2d` to change the sample time of a discrete-time system.

**Default:** -1 (discrete-time model with unspecified sample time)

## TimeUnit

Units for the time variable, the sample time  $T_s$ , and any time delays in the model, specified as one of the following values:

- 'nanoseconds'
- 'microseconds'
- 'milliseconds'
- 'seconds'
- 'minutes'
- 'hours'
- 'days'
- 'weeks'
- 'months'
- 'years'

Changing this property has no effect on other properties, and therefore changes the overall system behavior. Use `chgTimeUnit` to convert between time units without modifying system behavior.

**Default:** 'seconds'

## InputName

Input channel names, specified as one of the following:

- Character vector — For single-input models, for example, 'controls'.
- Cell array of character vectors — For multi-input models.

Alternatively, use automatic vector expansion to assign input names for multi-input models. For example, if `sys` is a two-input model, enter:

```
sys.InputName = 'controls';
```

The input names automatically expand to {'controls(1)'; 'controls(2)'}

When you estimate a model using an `iddata` object, `data`, the software automatically sets `InputName` to `data.InputName`.

You can use the shorthand notation `u` to refer to the `InputName` property. For example, `sys.u` is equivalent to `sys.InputName`.

Input channel names have several uses, including:

- Identifying channels on model display and plots
- Extracting subsystems of MIMO systems
- Specifying connection points when interconnecting models

**Default:** '' for all input channels

### **InputUnit**

Input channel units, specified as one of the following:

- Character vector — For single-input models, for example, 'seconds'.
- Cell array of character vectors — For multi-input models.

Use `InputUnit` to keep track of input signal units. `InputUnit` has no effect on system behavior.

**Default:** '' for all input channels

### **InputGroup**

Input channel groups. The `InputGroup` property lets you assign the input channels of MIMO systems into groups and refer to each group by name. Specify input groups as a structure. In this structure, field names are the group names, and field values are the input channels belonging to each group. For example:

```
sys.InputGroup.controls = [1 2];  
sys.InputGroup.noise = [3 5];
```

creates input groups named `controls` and `noise` that include input channels 1, 2 and 3, 5, respectively. You can then extract the subsystem from the `controls` inputs to all outputs using:

```
sys(:, 'controls')
```

**Default:** Struct with no fields

## OutputName

Output channel names, specified as one of the following:

- Character vector — For single-output models. For example, 'measurements'.
- Cell array of character vectors — For multi-output models.

Alternatively, use automatic vector expansion to assign output names for multi-output models. For example, if `sys` is a two-output model, enter:

```
sys.OutputName = 'measurements';
```

The output names automatically expand to `{'measurements(1)'; 'measurements(2)'}`.

When you estimate a model using an `iddata` object, `data`, the software automatically sets `OutputName` to `data.OutputName`.

You can use the shorthand notation `y` to refer to the `OutputName` property. For example, `sys.y` is equivalent to `sys.OutputName`.

Output channel names have several uses, including:

- Identifying channels on model display and plots
- Extracting subsystems of MIMO systems
- Specifying connection points when interconnecting models

**Default:** '' for all output channels

## OutputUnit

Output channel units, specified as one of the following:

- Character vector — For single-output models. For example, 'seconds'.
- Cell array of character vectors — For multi-output models.

Use `OutputUnit` to keep track of output signal units. `OutputUnit` has no effect on system behavior.

**Default:** '' for all output channels

## OutputGroup

Output channel groups. The `OutputGroup` property lets you assign the output channels of MIMO systems into groups and refer to each group by name. Specify output groups as a structure. In this structure, field names are the group names, and field values are the output channels belonging to each group. For example:

```
sys.OutputGroup.temperature = [1];  
sys.InputGroup.measurement = [3 5];
```

creates output groups named `temperature` and `measurement` that include output channels 1, and 3, 5, respectively. You can then extract the subsystem from all inputs to the measurement outputs using:

```
sys('measurement', :)
```

**Default:** Struct with no fields

### Name

System name, specified as a character vector. For example, `'system_1'`.

**Default:** `''`

### Notes

Any text that you want to associate with the system, stored as a string or a cell array of character vectors. The property stores whichever data type you provide. For instance, if `sys1` and `sys2` are dynamic system models, you can set their `Notes` properties as follows:

```
sys1.Notes = "sys1 has a string."  
sys2.Notes = 'sys2 has a character vector.'  
sys1.Notes  
sys2.Notes
```

```
ans =
```

```
    "sys1 has a string."
```

```
ans =
```

---

'sys2 has a character vector.'

**Default:** [0×1 string]

### **UserData**

Any type of data you want to associate with system, specified as any MATLAB data type.

**Default:** []

### **SamplingGrid**

Sampling grid for model arrays, specified as a data structure.

For arrays of identified linear (IDLTI) models that are derived by sampling one or more independent variables, this property tracks the variable values associated with each model. This information appears when you display or plot the model array. Use this information to trace results back to the independent variables.

Set the field names of the data structure to the names of the sampling variables. Set the field values to the sampled variable values associated with each model in the array. All sampling variables should be numeric and scalar valued, and all arrays of sampled values should match the dimensions of the model array.

For example, if you collect data at various operating points of a system, you can identify a model for each operating point separately and then stack the results together into a single system array. You can tag the individual models in the array with information regarding the operating point:

```
nominal_engine_rpm = [1000 5000 10000];  
sys.SamplingGrid = struct('rpm', nominal_engine_rpm)
```

where `sys` is an array containing three identified models obtained at rpms 1000, 5000 and 10000, respectively.

For model arrays generated by linearizing a Simulink model at multiple parameter values or operating points, the software populates `SamplingGrid` automatically with the variable values that correspond to each entry in the array. For example, the Simulink Control Design commands `linearize` and `sLinearizer` populate `SamplingGrid` in this way.

**Default:** []

## **See Also**

idgrey | idpoly | idproc | idssdata | idtf | n4sid | pem | ssest | ssestOptions |  
translatecov

## **Topics**

“Dynamic System Models”

**Introduced before R2006a**

# idssdata

State-space data of identified system

## Syntax

```
[A,B,C,D,K] = idssdata(sys)
[A,B,C,D,K,x0] = idssdata(sys)
[A,B,C,D,K,x0,dA,dB,dC,dD,dK,dx0] = idssdata(sys)
[A,B,C,D,K, ___] = idssdata(sys,j1,...,jN)
[A,B,C,D,K, ___] = idssdata(sys,'cell')
```

## Description

`[A,B,C,D,K] = idssdata(sys)` returns the A,B,C,D and K matrices of the identified state-space model `sys`.

`[A,B,C,D,K,x0] = idssdata(sys)` returns the initial state values, `x0`.

`[A,B,C,D,K,x0,dA,dB,dC,dD,dK,dx0] = idssdata(sys)` returns the uncertainties in the system matrices for `sys`.

`[A,B,C,D,K, ___] = idssdata(sys,j1,...,jN)` returns data for the `j1, ..., jN` entries in the model array `sys`.

`[A,B,C,D,K, ___] = idssdata(sys,'cell')` returns data for all the entries in the model array `sys` as separate cells in cell arrays.

## Input Arguments

### `sys`

Identified model.

If `sys` is not an identified state-space model (`idss` or `idgrey`), then it is first converted to an `idss` model. This conversion results in a loss of the model uncertainty information.

`sys` can be an array of identified models.

**`j1, ..., jN`**

Integer indices of `N` entries in the array `sys` of identified systems.

## Output Arguments

**`A, B, C, D, K`**

State-space matrices that represent `sys` as:

$$\begin{aligned}x[k+1] &= Ax[k] + Bu[k] + Ke[k]; x[0] = x0; \\ y[k] &= Cx[k] + Du[k] + e[k];\end{aligned}$$

If `sys` is an array of identified models, then `A, B, C, D, K` are multi-dimension arrays. To access the state-space matrix, say `A`, for the  $k$ -th entry of `sys`, use `A(:, :, k)`.

**`x0`**

Initial state.

If `sys` is an `idss` or `idgrey` model, then `x0` is the value obtained during estimation. It is also stored using the `Report.Parameters` property of `sys`.

For other model types, `x0` is zero.

If `sys` is an array of identified models, then `x0` contains a column for each entry in `sys`.

**`dA, dB, dC, dD, dK`**

Uncertainties associated with the state-space matrices `A, B, C, D, K`.

The uncertainty matrices represents 1 standard deviation of uncertainty.

If `sys` is an array of identified models, then `dA, dB, dC, dD, dK` are multi-dimension arrays. To access the state-space matrix, say `A`, for the  $k$ -th entry of `sys`, use `A(:, :, k)`.

**`dx0`**

Uncertainty associated with the initial state.

$dx_0$  represents 1 standard deviation of uncertainty.

If `sys` is an array of identified models, then `dx0` contains a column for each entry in `sys`.

## Examples

### Obtain Identified State-Space Matrices

Obtain the identified state-space matrices for a model estimated from data.

Identify a model using data.

```
load icEngine.mat
data = iddata(y,u,0.04);
sys = n4sid(data,4,'InputDelay',2);
```

`data` is an `iddata` object representing data sampled at a sampling rate of 0.04 seconds.

`sys` is an `idss` model representing the identified system.

Obtain identified state-space matrices of `sys`.

```
[A,B,C,D,K] = idssdata(sys);
```

### Obtain Initial State of Identified Model

Obtain the initial state associated with an identified model.

Identify a model using data.

```
load icEngine.mat
data = iddata(y,u,0.04);
sys = n4sid(data,4,'InputDelay',2);
```

`data` is an `iddata` object representing data sampled at a sampling rate of 0.04 seconds.

`sys` is an `idss` model representing the identified system.

Obtain the initial state associated with `sys`.

```
[A,B,C,D,K,x0] = idssdata(sys);
```

A, B, C, D and K represent the state-space matrices of the identified model `sys`. `x0` is the initial state identified for `sys`.

### Obtain Uncertainty Data of State-Space Matrices of Identified Model

Obtain the uncertainty matrices of the state-space matrices of an identified model.

Identify a model using data.

```
load icEngine.mat
data = iddata(y,u,0.04);
sys = n4sid(data,4,'InputDelay',2);
```

`data` is an `iddata` object representing data sampled at a sampling rate of 0.04 seconds.

`sys` is an `idss` model representing the identified system.

Obtain the uncertainty matrices associated with the state-space matrices of `sys`.

```
[A,B,C,D,K,x0,dA,dB,dC,dD,dx0] = idssdata(sys);
```

`dA`, `dB`, `dC`, `dD` and `dK` represent the uncertainty associated with the state-space matrices of the identified model `sys`. `dx0` represents the uncertainty associated with the estimated initial state.

### Obtain State-Space Matrices for Multiple Identified Models

Obtain the state-space matrices for multiple models from an array of identified models.

Identify multiple models using data.

```
load icEngine.mat
data = iddata(y,u,0.04);
sys2 = n4sid(data,2,'InputDelay',2);
sys3 = n4sid(data,3,'InputDelay',2);
sys4 = n4sid(data,4,'InputDelay',2);
sys = stack(1,sys2,sys3,sys4);
```

`data` is an `iddata` object representing data sampled at a sampling rate of 0.04 seconds.

`sys` is an array of `idss` models. The first entry of `sys` is a second-order identified system. The second and third entries of `sys` are third- and fourth-order identified systems, respectively.

Obtain the state-space matrices for the first and third entries of `sys`.

```
[A,B,C,D,K,x0] = idssdata(sys,1);  
[A,B,C,D,K,x0] = idssdata(sys,3);
```

### Obtain State-Space Matrices for Identified Model as Cell Array

Obtain the state-space matrices of an array of identified models in cell arrays.

Identify multiple models using `data`.

```
load icEngine.mat  
data = iddata(y,u,0.04);  
sys3 = n4sid(data,3,'InputDelay',2);  
sys4 = n4sid(data,4,'InputDelay',2);  
sys = stack(1,sys3,sys4);
```

`data` is an `iddata` object representing data sampled at a sampling rate of 0.04 seconds.

`sys` is an array of `idss` models. The first entry of `sys` is a third-order identified system and the second entry is a fourth-order identified system.

Obtain the state-space matrices of `sys` in cell arrays.

```
[A,B,C,D,K,x0] = idssdata(sys,'cell');
```

`A`, `B`, `C`, `D` and `K` are cell arrays containing the state-space matrices of the individual entries of the identified model array `sys`. `x0` is a cell array containing the estimated initial state of the individual entries of the identified model array `sys`.

## See Also

`idss` | `polydata` | `ssdata` | `tfdata` | `zpkdata`

**Introduced in R2012a**

# idtf

Transfer function model with identifiable parameters

## Syntax

```
sys = idtf(num,den)
sys = idtf(num,den,Ts)
sys = idtf( ____,Name,Value)

sys = idtf(sys0)
```

## Description

`sys = idtf(num,den)` creates a continuous-time transfer function with identifiable parameters (an `idtf` model). `num` specifies the current values of the transfer function numerator coefficients. `den` specifies the current values of the transfer function denominator coefficients.

`sys = idtf(num,den,Ts)` creates a discrete-time transfer function with identifiable parameters. `Ts` is the sample time.

`sys = idtf( ____,Name,Value)` creates a transfer function with properties on page 1-753 specified by one or more `Name,Value` pair arguments.

`sys = idtf(sys0)` converts any dynamic system model, `sys0`, to `idtf` model form.

## Object Description

An `idtf` model represents a system as a continuous-time or discrete-time transfer function with identifiable (estimable) coefficients.

A SISO transfer function is a ratio of polynomials with an exponential term. In continuous time,

$$G(s) = e^{-\tau s} \frac{b_n s^n + b_{n-1} s^{n-1} + \dots + b_0}{s^m + a_{m-1} s^{m-1} + \dots + a_0}.$$

In discrete time,

$$G(z^{-1}) = z^{-k} \frac{b_n z^{-n} + b_{n-1} z^{-n+1} + \dots + b_0}{z^{-m} + a_{m-1} z^{-m+1} + \dots + a_0}.$$

In discrete time,  $z^{-k}$  represents a time delay of  $kT_s$ , where  $T_s$  is the sample time.

For `idtf` models, the denominator coefficients  $a_0, \dots, a_{m-1}$  and the numerator coefficients  $b_0, \dots, b_n$  can be estimable parameters. (The leading denominator coefficient is always fixed to 1.) The time delay  $\tau$  (or  $k$  in discrete time) can also be an estimable parameter. The `idtf` model stores the polynomial coefficients  $a_0, \dots, a_{m-1}$  and  $b_0, \dots, b_n$  in the `Denominator` and `Numerator` properties of the model, respectively. The time delay  $\tau$  or  $k$  is stored in the `IODelay` property of the model.

A MIMO transfer function contains a SISO transfer function corresponding to each input-output pair in the system. For `idtf` models, the polynomial coefficients and transport delays of each input-output pair are independently estimable parameters.

There are three ways to obtain an `idtf` model.

- Estimate the `idtf` model based on input-output measurements of a system, using `tfest`. The `tfest` command estimates the values of the transfer function coefficients and transport delays. The estimated values are stored in the `Numerator`, `Denominator`, and `IODelay` properties of the resulting `idtf` model. The `Report` property of the resulting model stores information about the estimation, such as handling of initial conditions and options used in estimation.

When you obtain an `idtf` model by estimation, you can extract estimated coefficients and their uncertainties from the model. To do so, use commands such as `tfdata`, `getpar`, or `getcov`.

- Create an `idtf` model using the `idtf` command.

You can create an `idtf` model to configure an initial parameterization for estimation of a transfer function to fit measured response data. When you do so, you can specify constraints on such values as the numerator and denominator coefficients and transport delays. For example, you can fix the values of some parameters, or specify

minimum or maximum values for the free parameters. You can then use the configured model as an input argument to `tfest` to estimate parameter values with those constraints.

- Convert an existing dynamic system model to an `idtf` model using the `idtf` command.

---

**Note** Unlike `idss` and `idpoly`, `idtf` uses a trivial noise model and does not parameterize the noise.

So,  $H = 1$  in  $y = Gu + He$ .

---

## Examples

### Create a Continuous-Time Transfer Function Model

Specify a continuous-time, single-input, single-output (SISO) transfer function with estimable parameters. The initial values of the transfer function are:

$$G(s) = \frac{s + 4}{s^2 + 20s + 5}$$

```
num = [1 4];  
den = [1 20 5];  
G = idtf(num,den);
```

`G` is an `idtf` model. `num` and `den` specify the initial values of the numerator and denominator polynomial coefficients in descending powers of  $s$ . The numerator coefficients having initial values 1 and 4 are estimable parameters. The denominator coefficient having initial values 20 and 5 are also estimable parameters. The leading denominator coefficient is always fixed to 1.

You can use `G` to specify an initial parametrization for estimation with `tfest`.

### Create Transfer Function with Known Input Delay and Specified Attributes

Specify a continuous-time, SISO transfer function with known input delay. The transfer function initial values are given by:

$$G(s) = e^{-5.8s} \frac{5}{s + 5}$$

Label the input of the transfer function with the name 'Voltage' and specify the input units as volt.

Use Name, Value input pairs to specify the delay, input name, and input unit.

```
num = 5;  
den = [1 5];  
input_delay = 5.8;  
input_name = 'Voltage';  
input_unit = 'volt';  
G = idtf(num,den,'InputDelay',input_delay,...  
         'InputName',input_name,'InputUnit',input_unit);
```

$G$  is an `idtf` model. You can use  $G$  to specify an initial parametrization for estimation with `tfest`. If you do so, model properties such as `InputDelay`, `InputName`, and `InputUnit` are applied to the estimated model. The estimation process treats `InputDelay` as a fixed value. If you want to estimate the delay and specify an initial value of 5.8 s, use the `IODelay` property instead.

### Create Discrete-Time Transfer Function

Specify a discrete-time SISO transfer function with estimable parameters. The initial values of the transfer function are:

$$H(z) = \frac{z - 0.1}{z + 0.8}$$

Specify the sample time as 0.2 seconds.

```
num = [1 -0.1];  
den = [1 0.8];
```

```
Ts = 0.2;
H = idtf(num,den,Ts);
```

num and den are the initial values of the numerator and denominator polynomial coefficients. For discrete-time systems, specify the coefficients in ascending powers of  $z^{-1}$ .

Ts specifies the sample time for the transfer function as 0.2 seconds.

H is an idtf model. The numerator and denominator coefficients are estimable parameters (except for the leading denominator coefficient, which is fixed to 1).

### Create MIMO Discrete-Time Transfer Function

Specify a discrete-time, two-input, two-output transfer function. The initial values of the MIMO transfer function are:

$$H(z) = \begin{bmatrix} \frac{1}{z+0.2} & \frac{z}{z+0.7} \\ \frac{-z+2}{z-0.3} & \frac{3}{z+0.3} \end{bmatrix}$$

Specify the sample time as 0.2 seconds.

```
nums = {1,[1,0];[-1,2],3};
dens = {[1,0.2],[1,0.7]};[1,-0.3],[1,0.3]};
Ts = 0.2;
H = idtf(nums,dens,Ts);
```

nums and dens specify the initial values of the coefficients in cell arrays. Each entry in the cell array corresponds to the numerator or denominator of the transfer function of one input-output pair. For example, the first row of nums is {1, [1, 0]}. This cell array specifies the numerators across the first row of transfer functions in H. Likewise, the first row of dens, {[1, 0.2], [1, 0.7]}, specifies the denominators across the first row of H.

Ts specifies the sample time for the transfer function as 0.2 seconds.

H is an `idtf` model. All of the polynomial coefficients are estimable parameters, except for the leading coefficient of each denominator polynomial. These coefficients are always fixed to 1.

### Specify Transfer Function Display Variable

Specify the following discrete-time transfer function in terms of  $q^{-1}$ :

$$H(q^{-1}) = \frac{1 + 0.4q^{-1}}{1 + 0.1q^{-1} - 0.3q^{-2}}$$

Specify the sample time as 0.1 seconds.

```
num = [1 0.4];  
den = [1 0.1 -0.3];  
Ts = 0.1;  
convention_variable = 'q^-1';  
H = idtf(num,den,Ts,'Variable',convention_variable);
```

Use a `Name, Value` pair argument to specify the variable  $q^{-1}$ .

`num` and `den` are the numerator and denominator polynomial coefficients in ascending powers of  $q^{-1}$ .

`Ts` specifies the sample time for the transfer function as 0.1 seconds.

H is an `idtf` model.

### Gain Matrix Transfer Function

Specify a transfer function with estimable coefficients whose initial value is the static gain matrix:

$$H(s) = \begin{bmatrix} 1 & 0 & 1 \\ 1 & 1 & 0 \\ 3 & 0 & 2 \end{bmatrix}$$

```
M = [1 0 1; 1 1 0; 3 0 2];
H = idtf(M);
```

H is an `idtf` model that describes a three input ( $N_u=3$ ), three output ( $N_y=3$ ) transfer function. Each input/output channel is an estimable static gain. The initial values of the gains are given by the values in the matrix M.

### Convert Identifiable State-Space Model to Identifiable Transfer Function

Convert a state-space model with identifiable parameters to a transfer function with identifiable parameters.

Convert the following identifiable state-space model to an identifiable transfer function.

$$\begin{aligned}\tilde{x}(t) &= \begin{bmatrix} -0.2 & 0 \\ 0 & -0.3 \end{bmatrix} x(t) + \begin{bmatrix} -2 \\ 4 \end{bmatrix} u(t) + \begin{bmatrix} 0.1 \\ 0.2 \end{bmatrix} e(t) \\ y(t) &= [1 \quad 1]x(t)\end{aligned}$$

```
A = [-0.2, 0; 0, -0.3];
B = [2;4];
C = [1, 1];
D = 0;
K = [0.1; 0.2];
sys0 = idss(A,B,C,D,K,'NoiseVariance',0.1);
sys = idtf(sys0);
```

A, B, C, D and K are matrices that specify `sys0`, an identifiable state-space model with a noise variance of 0.1.

`sys = idtf(sys0)` creates an `idtf` model, `sys`.

### Estimate Transfer Function Model By Specifying Number of Poles

Load time-domain system response data and use it to estimate a transfer function for the system.

```
load iddata1 z1;  
np = 2;  
sys = tfest(z1,np);
```

`z1` is an `iddata` object that contains time-domain, input-output data.

`np` specifies the number of poles in the estimated transfer function.

`sys` is an `idtf` model containing the estimated transfer function.

To see the numerator and denominator coefficients of the resulting estimated model `sys`, enter:

```
sys.Numerator
```

```
ans = 1×2
```

```
2.4554 176.9856
```

```
sys.Denominator
```

```
ans = 1×3
```

```
1.0000 3.1625 23.1631
```

To view the uncertainty in the estimates of the numerator and denominator and other information, use `tfdata`.

## Create Array of Transfer Function Models

Create an array of transfer function models with identifiable coefficients. Each transfer function in the array is of the form:

$$H(s) = \frac{a}{s + a}.$$

The initial value of the coefficient  $a$  varies across the array, from 0.1 to 1.0, in increments of 0.1.

```
H = idtf(zeros(1,1,10));
for k = 1:10
    num = k/10;
    den = [1 k/10];
    H(:,:,k) = idtf(num,den);
end
```

The first command preallocates a one-dimensional, 10-element array, H, and fills it with empty `idtf` models.

The first two dimensions of a model array are the output and input dimensions. The remaining dimensions are the array dimensions. `H(:, :, k)` represents the  $k^{\text{th}}$  model in the array. Thus, the `for` loop replaces the  $k^{\text{th}}$  entry in the array with a transfer function whose coefficients are initialized with  $a = k/10$ .

## Input Arguments

### **num**

Initial values of transfer function numerator coefficients.

For SISO transfer functions, specify the initial values of the numerator coefficients `num` as a row vector. Specify the coefficients in order of:

- Descending powers of  $s$  or  $p$  (for continuous-time transfer functions)
- Ascending powers of  $z^{-1}$  or  $q^{-1}$  (for discrete-time transfer functions)

Use `NaN` for any coefficient whose initial value is not known.

For MIMO transfer functions with  $N_y$  outputs and  $N_u$  inputs, `num` is a  $N_y$ -by- $N_u$  cell array of numerator coefficients for each input/output pair.

### **den**

Initial values of transfer function denominator coefficients.

For SISO transfer functions, specify the initial values of the denominator coefficients `den` as a row vector. Specify the coefficients in order of:

- Descending powers of  $s$  or  $p$  (for continuous-time transfer functions)
- Ascending powers of  $z^{-1}$  or  $q^{-1}$  (for discrete-time transfer functions)

The leading coefficient in `den` must be 1. Use `NaN` for any coefficient whose initial value is not known.

For MIMO transfer functions with `Ny` outputs and `Nu` inputs, `den` is a `Ny-by-Nu` cell array of denominator coefficients for each input/output pair.

## **Ts**

Sample time. For continuous-time models, `Ts = 0`. For discrete-time models, `Ts` is a positive scalar representing the sampling period. This value is expressed in the unit specified by the `TimeUnit` property of the model. To denote a discrete-time model with unspecified sample time, set `Ts = -1`.

Changing this property does not discretize or resample the model. Use `c2d` and `d2c` to convert between continuous- and discrete-time representations. Use `d2d` to change the sample time of a discrete-time system.

**Default:** 0 (continuous time)

## **sys0**

Dynamic system.

Any dynamic system to convert to an `idtf` model.

When `sys0` is an identified model, its estimated parameter covariance is lost during conversion. If you want to translate the estimated parameter covariance during the conversion, use `translatecov`.

## **Name-Value Pair Arguments**

Specify optional comma-separated pairs of `Name`, `Value` arguments. `Name` is the argument name and `Value` is the corresponding value. `Name` must appear inside quotes. You can specify several name and value pair arguments in any order as `Name1, Value1, ..., NameN, ValueN`.

Use `Name`, `Value` arguments to specify additional properties on page 1-753 of `idtf` models during model creation. For example,

`idtf(num,den,'InputName','Voltage')` creates an `idtf` model with the `InputName` property set to `Voltage`.

## Properties

`idtf` object properties include:

### Numerator

Values of transfer function numerator coefficients.

If you create an `idtf` model `sys` using the `idtf` command, `sys.Numerator` contains the initial values of numerator coefficients that you specify with the `num` input argument.

If you obtain an `idtf` model by identification using `tfest`, then `sys.Numerator` contains the estimated values of the numerator coefficients.

For an `idtf` model `sys`, the property `sys.Numerator` is an alias for the value of the property `sys.Structure.Numerator.Value`.

For SISO transfer functions, the values of the numerator coefficients are stored as a row vector in order of:

- Descending powers of  $s$  or  $p$  (for continuous-time transfer functions)
- Ascending powers of  $z^{-1}$  or  $q^{-1}$  (for discrete-time transfer functions)

Any coefficient whose initial value is not known is stored as `NaN`.

For MIMO transfer functions with `Ny` outputs and `Nu` inputs, `Numerator` is a `Ny`-by-`Nu` cell array of numerator coefficients for each input/output pair.

### Denominator

Values of transfer function denominator coefficients.

If you create an `idtf` model `sys` using the `idtf` command, `sys.Denominator` contains the initial values of denominator coefficients that you specify with the `den` input argument.

If you obtain an `idtf` model `sys` by identification using `tfest`, then `sys.Denominator` contains the estimated values of the denominator coefficients.

For an `idtf` model `sys`, the property `sys.Denominator` is an alias for the value of the property `sys.Structure.Denominator.Value`.

For SISO transfer functions, the values of the denominator coefficients are stored as a row vector in order of:

- Descending powers of  $s$  or  $p$  (for continuous-time transfer functions)
- Ascending powers of  $z^{-1}$  or  $q^{-1}$  (for discrete-time transfer functions)

The leading coefficient in `Denominator` is fixed to 1. Any coefficient whose initial value is not known is stored as `NaN`.

For MIMO transfer functions with  $N_y$  outputs and  $N_u$  inputs, `Denominator` is a  $N_y$ -by- $N_u$  cell array of denominator coefficients for each input/output pair.

## Variable

Transfer function display variable, specified as one of the following values:

- 's' — Default for continuous-time models
- 'p' — Equivalent to 's'
- 'z^-1' — Default for discrete-time models
- 'q^-1' — Equivalent to 'z^-1'

The value of `Variable` is reflected in the display, and also affects the interpretation of the `num` and `den` coefficient vectors for discrete-time models. For `Variable = 'z^-1'` or `'q^-1'`, the coefficient vectors are ordered as ascending powers of the variable.

## IODElay

Transport delays. `IODElay` is a numeric array specifying a separate transport delay for each input/output pair.

If you create an `idtf` model `sys` using the `idtf` command, `sys.IODElay` contains the initial values of the transport delay that you specify with a `Name, Value` argument pair.

If you obtain an `idtf` model `sys` by identification using `tfest`, then `sys.IODElay` contains the estimated values of the transport delay.

For an `idtf` model `sys`, the property `sys.IODElay` is an alias for the value of the property `sys.Structure.IODElay.Value`.

For continuous-time systems, transport delays are expressed in the time unit stored in the `TimeUnit` property. For discrete-time systems, transport delays are expressed as integers denoting delay of a multiple of the sample time `Ts`.

For a MIMO system with `Ny` outputs and `Nu` inputs, set `IODelay` as a `Ny`-by-`Nu` array. Each entry of this array is a numerical value representing the transport delay for the corresponding input/output pair. You can set `IODelay` to a scalar value to apply the same delay to all input/output pairs.

**Default:** 0 for all input/output pairs

### Structure

Information about the estimable parameters of the `idtf` model. `Structure.Numerator`, `Structure.Denominator`, and `Structure.IODelay` contain information about the numerator coefficients, denominator coefficients, and transport delay, respectively. Each contains the following fields:

- **Value** — Parameter values. For example, `sys.Structure.Numerator.Value` contains the initial or estimated values of the numerator coefficients.

`NaN` represents unknown parameter values. For denominators, the value of the leading coefficient, specified by `sys.Structure.Denominator.Value(1)` is fixed to 1.

For SISO models, `sys.Numerator`, `sys.Denominator`, and `sys.IODelay` are aliases for `sys.Structure.Numerator.Value`, `sys.Structure.Denominator.Value`, and `sys.Structure.IODelay.Value`, respectively.

For MIMO models, `sys.Numerator{i,j}` is an alias for `sys.Structure(i,j).Numerator.Value`, and `sys.Denominator{i,j}` is an alias for `sys.Structure(i,j).Denominator.Value`. Additionally, `sys.IODelay(i,j)` is an alias for `sys.Structure(i,j).IODelay.Value`

- **Minimum** — Minimum value that the parameter can assume during estimation. For example, `sys.Structure.IODelay.Minimum = 0.1` constrains the transport delay to values greater than or equal to 0.1.

`sys.Structure.IODelay.Minimum` must be greater than or equal to zero.

- **Maximum** — Maximum value that the parameter can assume during estimation.
- **Free** — Boolean specifying whether the parameter is a free estimation variable. If you want to fix the value of a parameter during estimation, set the corresponding `Free =`

`false`. For example, `sys.Structure.Denominator.Free = false` fixes all of the denominator coefficients in `sys` to the values specified in `sys.Structure.Denominator.Value`.

For denominators, the value of `Free` for the leading coefficient, specified by `sys.Structure.Denominator.Free(1)`, is always `false` (the leading denominator coefficient is always fixed to 1).

- **Scale** — Scale of the parameter's value. `Scale` is not used in estimation.
- **Info** — Structure array for storing parameter units and labels. The structure has `Label` and `Unit` fields.

Specify parameter units and labels as character vectors. For example, `'Time'`.

For a MIMO model with  $N_y$  outputs and  $N_u$  input, `Structure` is an  $N_y$ -by- $N_u$  array. The element `Structure(i, j)` contains information corresponding to the transfer function for the  $(i, j)$  input-output pair.

### NoiseVariance

The variance (covariance matrix) of the model innovations  $e$ .

An identified model includes a white, Gaussian noise component  $e(t)$ . `NoiseVariance` is the variance of this noise component. Typically, the model estimation function (such as `tfest`) determines this variance.

For SISO models, `NoiseVariance` is a scalar. For MIMO models, `NoiseVariance` is a  $N_y$ -by- $N_y$  matrix, where  $N_y$  is the number of outputs in the system.

### Report

Summary report that contains information about the estimation options and results when the transfer function model is obtained using estimation commands, such as `tfest` and `impulseest`. Use `Report` to query a model for how it was estimated, including its:

- Estimation method
- Estimation options
- Search termination conditions
- Estimation data fit and other quality metrics

The contents of `Report` are irrelevant if the model was created by construction.

```

m = idtf([1 4],[1 20 5]);
m.Report.OptionsUsed

ans =

     []

```

If you obtain the transfer function model using estimation commands, the fields of `Report` contain information on the estimation data, options, and results.

```

load iddata2 z2;
m = tfest(z2,3);
m.Report.OptionsUsed

    InitializeMethod: 'iv'
    InitializeOptions: [1x1 struct]
    InitialCondition: 'auto'
                   Focus: 'simulation'
    EstimateCovariance: 1
                   Display: 'off'
    InputOffset: []
    OutputOffset: []
    Regularization: [1x1 struct]
    SearchMethod: 'auto'
    SearchOptions: [1x1 idoptions.search.identsolver]
    OutputWeight: []
    Advanced: [1x1 struct]

```

`Report` is a read-only property.

For more information on this property and how to use it, see the [Output Arguments](#) section of the corresponding estimation command reference page and “[Estimation Report](#)”.

### InputDelay

Input delays. `InputDelay` is a numeric vector specifying a time delay for each input channel. For continuous-time systems, specify input delays in the time unit stored in the `TimeUnit` property. For discrete-time systems, specify input delays in integer multiples of the sample time  $T_s$ . For example, `InputDelay = 3` means a delay of three sample times.

For a system with  $N_u$  inputs, set `InputDelay` to an  $N_u$ -by-1 vector. Each entry of this vector is a numerical value representing the input delay for the corresponding input channel. You can also set `InputDelay` to a scalar value to apply the same delay to all channels.

Estimation treats `InputDelay` as a fixed constant of the model. Estimation uses the `IODelay` property for estimating time delays. To specify initial values and constraints for estimation of time delays, use `sys.Structure.IODelay`.

**Default:** 0 for all input channels

## OutputDelay

Output delays.

For identified systems, like `idtf`, `OutputDelay` is fixed to zero.

## Ts

Sample time. For continuous-time models,  $T_s = 0$ . For discrete-time models,  $T_s$  is a positive scalar representing the sampling period. This value is expressed in the unit specified by the `TimeUnit` property of the model. To denote a discrete-time model with unspecified sample time, set  $T_s = -1$ .

Changing this property does not discretize or resample the model. Use `c2d` and `d2c` to convert between continuous- and discrete-time representations. Use `d2d` to change the sample time of a discrete-time system.

**Default:** 0 (continuous time)

## TimeUnit

Units for the time variable, the sample time  $T_s$ , and any time delays in the model, specified as one of the following values:

- 'nanoseconds'
- 'microseconds'
- 'milliseconds'
- 'seconds'
- 'minutes'
- 'hours'
- 'days'
- 'weeks'
- 'months'

- 'years'

Changing this property has no effect on other properties, and therefore changes the overall system behavior. Use `chgTimeUnit` to convert between time units without modifying system behavior.

**Default:** 'seconds'

### **InputName**

Input channel names, specified as one of the following:

- Character vector — For single-input models, for example, 'controls'.
- Cell array of character vectors — For multi-input models.

Alternatively, use automatic vector expansion to assign input names for multi-input models. For example, if `sys` is a two-input model, enter:

```
sys.InputName = 'controls';
```

The input names automatically expand to {'controls(1)'; 'controls(2)'}

When you estimate a model using an `iddata` object, `data`, the software automatically sets `InputName` to `data.InputName`.

You can use the shorthand notation `u` to refer to the `InputName` property. For example, `sys.u` is equivalent to `sys.InputName`.

Input channel names have several uses, including:

- Identifying channels on model display and plots
- Extracting subsystems of MIMO systems
- Specifying connection points when interconnecting models

**Default:** '' for all input channels

### **InputUnit**

Input channel units, specified as one of the following:

- Character vector — For single-input models, for example, 'seconds'.

- Cell array of character vectors — For multi-input models.

Use `InputUnit` to keep track of input signal units. `InputUnit` has no effect on system behavior.

**Default:** '' for all input channels

### **InputGroup**

Input channel groups. The `InputGroup` property lets you assign the input channels of MIMO systems into groups and refer to each group by name. Specify input groups as a structure. In this structure, field names are the group names, and field values are the input channels belonging to each group. For example:

```
sys.InputGroup.controls = [1 2];  
sys.InputGroup.noise = [3 5];
```

creates input groups named `controls` and `noise` that include input channels 1, 2 and 3, 5, respectively. You can then extract the subsystem from the `controls` inputs to all outputs using:

```
sys(:, 'controls')
```

**Default:** Struct with no fields

### **OutputName**

Output channel names, specified as one of the following:

- Character vector — For single-output models. For example, 'measurements'.
- Cell array of character vectors — For multi-output models.

Alternatively, use automatic vector expansion to assign output names for multi-output models. For example, if `sys` is a two-output model, enter:

```
sys.OutputName = 'measurements';
```

The output names automatically expand to `{'measurements(1)'; 'measurements(2)'}`.

When you estimate a model using an `iddata` object, `data`, the software automatically sets `OutputName` to `data.OutputName`.

You can use the shorthand notation `y` to refer to the `OutputName` property. For example, `sys.y` is equivalent to `sys.OutputName`.

Output channel names have several uses, including:

- Identifying channels on model display and plots
- Extracting subsystems of MIMO systems
- Specifying connection points when interconnecting models

**Default:** `''` for all output channels

### OutputUnit

Output channel units, specified as one of the following:

- Character vector — For single-output models. For example, `'seconds'`.
- Cell array of character vectors — For multi-output models.

Use `OutputUnit` to keep track of output signal units. `OutputUnit` has no effect on system behavior.

**Default:** `''` for all output channels

### OutputGroup

Output channel groups. The `OutputGroup` property lets you assign the output channels of MIMO systems into groups and refer to each group by name. Specify output groups as a structure. In this structure, field names are the group names, and field values are the output channels belonging to each group. For example:

```
sys.OutputGroup.temperature = [1];  
sys.InputGroup.measurement = [3 5];
```

creates output groups named `temperature` and `measurement` that include output channels 1, and 3, 5, respectively. You can then extract the subsystem from all inputs to the measurement outputs using:

```
sys('measurement', :)
```

**Default:** Struct with no fields

## Name

System name, specified as a character vector. For example, 'system\_1'.

**Default:** ''

## Notes

Any text that you want to associate with the system, stored as a string or a cell array of character vectors. The property stores whichever data type you provide. For instance, if `sys1` and `sys2` are dynamic system models, you can set their `Notes` properties as follows:

```
sys1.Notes = "sys1 has a string.";
sys2.Notes = 'sys2 has a character vector.';
sys1.Notes
sys2.Notes
```

```
ans =
```

```
    "sys1 has a string."
```

```
ans =
```

```
    'sys2 has a character vector.'
```

**Default:** [0×1 string]

## UserData

Any type of data you want to associate with system, specified as any MATLAB data type.

**Default:** []

## SamplingGrid

Sampling grid for model arrays, specified as a data structure.

For arrays of identified linear (IDLTI) models that are derived by sampling one or more independent variables, this property tracks the variable values associated with each

model. This information appears when you display or plot the model array. Use this information to trace results back to the independent variables.

Set the field names of the data structure to the names of the sampling variables. Set the field values to the sampled variable values associated with each model in the array. All sampling variables should be numeric and scalar valued, and all arrays of sampled values should match the dimensions of the model array.

For example, if you collect data at various operating points of a system, you can identify a model for each operating point separately and then stack the results together into a single system array. You can tag the individual models in the array with information regarding the operating point:

```
nominal_engine_rpm = [1000 5000 10000];  
sys.SamplingGrid = struct('rpm', nominal_engine_rpm)
```

where `sys` is an array containing three identified models obtained at rpms 1000, 5000 and 10000, respectively.

For model arrays generated by linearizing a Simulink model at multiple parameter values or operating points, the software populates `SamplingGrid` automatically with the variable values that correspond to each entry in the array. For example, the Simulink Control Design commands `linearize` and `sLinearizer` populate `SamplingGrid` in this way.

**Default:** []

## See Also

`getcov` | `getpar` | `idfrd` | `idpoly` | `idproc` | `idss` | `oe` | `tfdata` | `tfest` | `translatecov`

## Topics

“Estimate Transfer Function Models With Prior Knowledge of Model Structure and Constraints”

“Dynamic System Models”

**Introduced in R2012a**

## ifft

Transform `iddata` objects from frequency to time domain

## Syntax

```
dat = ifft(Datf)
```

## Description

`ifft` transforms a frequency-domain `iddata` object to the time domain. It requires the frequencies on `Datf` to be equally spaced from frequency 0 to the Nyquist frequency. This means that if there are  $N$  frequencies in `Datf` and the sample time is  $T_s$ , then

`Datf.Frequency = [0:df:F]`, where  $F$  is  $\pi/T_s$  if  $N$  is odd and  $F = \pi/T_s * (1 - 1/N)$  if  $N$  is even.

## See Also

`fft` | `iddata`

**Introduced in R2007a**

# impulse

Impulse response plot of dynamic system; impulse response data

## Syntax

```
impulse(sys)
impulse(sys,Tfinal)
impulse(sys,t)
impulse(sys1,sys2,...,sysN)
impulse(sys1,sys2,...,sysN,Tfinal)
impulse(sys1,sys2,...,sysN,t)
[y,t] = impulse(sys)
[y,t] = impulse(sys,Tfinal)
y = impulse(sys,t)
[y,t,x] = impulse(sys)
[y,t,x,yzd] = impulse(sys)
```

## Description

`impulse` calculates the unit impulse response of a dynamic system model. For continuous-time dynamic systems, the impulse response is the response to a Dirac input  $\delta(t)$ . For discrete-time systems, the impulse response is the response to a unit area pulse of length  $T_s$  and height  $1/T_s$ , where  $T_s$  is the sample time of the system. (This pulse approaches  $\delta(t)$  as  $T_s$  approaches zero.) For state-space models, `impulse` assumes initial state values are zero.

`impulse(sys)` plots the impulse response of the dynamic system model `sys`. This model can be continuous or discrete, and SISO or MIMO. The impulse response of multi-input systems is the collection of impulse responses for each input channel. The duration of simulation is determined automatically to display the transient behavior of the response.

`impulse(sys,Tfinal)` simulates the impulse response from  $t = 0$  to the final time  $t = T_{\text{final}}$ . Express `Tfinal` in the system time units, specified in the `TimeUnit` property of `sys`. For discrete-time systems with unspecified sample time ( $T_s = -1$ ), `impulse` interprets `Tfinal` as the number of sampling periods to simulate.

`impulse(sys,t)` uses the user-supplied time vector `t` for simulation. Express `t` in the system time units, specified in the `TimeUnit` property of `sys`. For discrete-time models, `t` should be of the form `Ti:Ts:Tf`, where `Ts` is the sample time. For continuous-time models, `t` should be of the form `Ti:dt:Tf`, where `dt` becomes the sample time of a discrete approximation to the continuous system (see "Algorithms" on page 1-770). The `impulse` command always applies the impulse at `t=0`, regardless of `Ti`.

To plot the impulse responses of several models `sys1,..., sysN` on a single figure, use:

```
impulse(sys1,sys2,...,sysN)
```

```
impulse(sys1,sys2,...,sysN,Tfinal)
```

```
impulse(sys1,sys2,...,sysN,t)
```

As with `bode` or `plot`, you can specify a particular color, linestyle, and/or marker for each system, for example,

```
impulse(sys1,'y:',sys2,'g--')
```

See "Plotting and Comparing Multiple Systems" and the `bode` entry in this section for more details.

When invoked with output arguments:

```
[y,t] = impulse(sys)
```

```
[y,t] = impulse(sys,Tfinal)
```

```
y = impulse(sys,t)
```

`impulse` returns the output response `y` and the time vector `t` used for simulation (if not supplied as an argument to `impulse`). No plot is drawn on the screen. For single-input systems, `y` has as many rows as time samples (length of `t`), and as many columns as outputs. In the multi-input case, the impulse responses of each input channel are stacked up along the third dimension of `y`. The dimensions of `y` are then

For state-space models only:

```
[y,t,x] = impulse(sys)
```

(length of `t`) × (number of outputs) × (number of inputs)

and  $y(:, :, j)$  gives the response to an impulse disturbance entering the  $j$ th input channel. Similarly, the dimensions of  $x$  are

(length of  $t$ )  $\times$  (number of states)  $\times$  (number of inputs)

`[y,t,x,ysd] = impulse(sys)` returns the standard deviation YSD of the response  $Y$  of an identified system  $SYS$ . YSD is empty if  $SYS$  does not contain parameter covariance information.

## Examples

### Impulse Response Plot of Second-Order State-Space Model

Plot the impulse response of the second-order state-space model

$$\begin{bmatrix} \dot{x}_1 \\ \dot{x}_2 \end{bmatrix} = \begin{bmatrix} -0.5572 & -0.7814 \\ 0.7814 & 0 \end{bmatrix} \begin{bmatrix} x_1 \\ x_2 \end{bmatrix} + \begin{bmatrix} 1 & -1 \\ 0 & 2 \end{bmatrix} \begin{bmatrix} u_1 \\ u_2 \end{bmatrix}$$

$$y = [1.9691 \quad 6.4493] \begin{bmatrix} x_1 \\ x_2 \end{bmatrix}$$

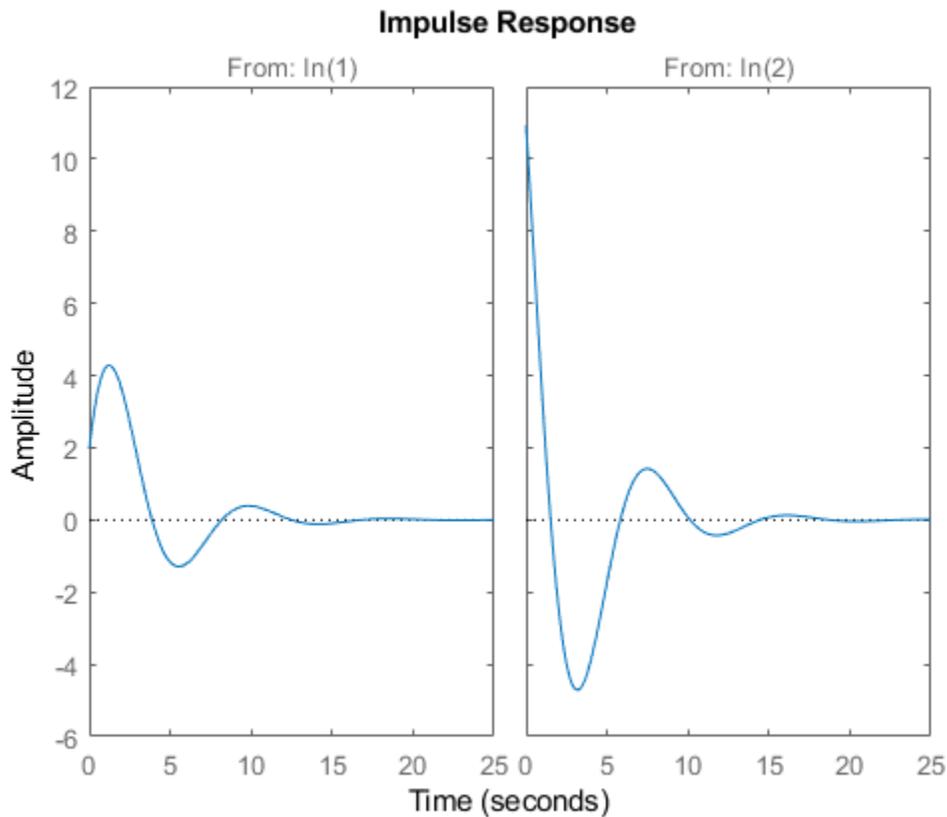
```
a = [-0.5572 -0.7814;0.7814 0];
```

```
b = [1 -1;0 2];
```

```
c = [1.9691 6.4493];
```

```
sys = ss(a,b,c,0);
```

```
impulse(sys)
```



The left plot shows the impulse response of the first input channel, and the right plot shows the impulse response of the second input channel.

You can store the impulse response data in MATLAB arrays by

```
[y,t] = impulse(sys);
```

Because this system has two inputs, y is a 3-D array with dimensions

```
size(y)
```

```
ans = 1x3
```

```
139    1    2
```

(the first dimension is the length of  $t$ ). The impulse response of the first input channel is then accessed by

```
ch1 = y(:, :, 1);
size(ch1)
```

```
ans = 1x2
    139     1
```

## Impulse Data from Identified System

Fetch the impulse response and the corresponding 1 std uncertainty of an identified linear system .

```
load(fullfile(matlabroot, 'toolbox', 'ident', 'iddemos', 'data', 'dcmotordata'));
z = iddata(y, u, 0.1, 'Name', 'DC-motor');
set(z, 'InputName', 'Voltage', 'InputUnit', 'V');
set(z, 'OutputName', {'Angular position', 'Angular velocity'});
set(z, 'OutputUnit', {'rad', 'rad/s'});
set(z, 'Tstart', 0, 'TimeUnit', 's');

model = tfest(z, 2);
[y, t, ~, ysd] = impulse(model, 2);

% Plot 3 std uncertainty
subplot(211)
plot(t, y(:, 1), t, y(:, 1) + 3 * ysd(:, 1), 'k:', t, y(:, 1) - 3 * ysd(:, 1), 'k:')
subplot(212)
plot(t, y(:, 2), t, y(:, 2) + 3 * ysd(:, 2), 'k:', t, y(:, 2) - 3 * ysd(:, 2), 'k:')
```

## Limitations

The impulse response of a continuous system with nonzero  $D$  matrix is infinite at  $t = 0$ . impulse ignores this discontinuity and returns the lower continuity value  $Cb$  at  $t = 0$ .

## Tips

You can change the properties of your plot, for example the units. For information on the ways to change properties of your plots, see “Ways to Customize Plots” (Control System Toolbox).

## Algorithms

Continuous-time models are first converted to state space. The impulse response of a single-input state-space model

$$\begin{aligned}\dot{x} &= Ax + bu \\ y &= Cx\end{aligned}$$

is equivalent to the following unforced response with initial state  $b$ .

$$\begin{aligned}\dot{x} &= Ax, \quad x(0) = b \\ y &= Cx\end{aligned}$$

To simulate this response, the system is discretized using zero-order hold on the inputs. The sample time is chosen automatically based on the system dynamics, except when a time vector  $t = 0:dt:Tf$  is supplied ( $dt$  is then used as sample time).

## See Also

**Linear System Analyzer** | `impulseest` | `lsim` | `step`

**Introduced before R2006a**

# impulseest

Nonparametric impulse response estimation

## Syntax

```
sys = impulseest(data)
sys = impulseest(data,N)
sys = impulseest(data,N,NK)
sys = impulseest( ____,options)
```

## Description

`sys = impulseest(data)` estimates an impulse response model, `sys`, using time- or frequency-domain data, `data`. The model order (number of nonzero impulse response coefficients) is determined automatically using persistence of excitation analysis on the input data.

`sys = impulseest(data,N)` estimates an  $N$ th order impulse response model, corresponding to the time range  $0 : T_s : (N - 1) * T_s$ , where  $T_s$  is the data sample time.

`sys = impulseest(data,N,NK)` specifies a transport delay of  $NK$  samples in the estimated impulse response.

`sys = impulseest( ____,options)` specifies estimation options using the options set `options`.

Use nonparametric impulse response to analyze data for feedback effects, delays and significant time constants.

## Input Arguments

### **data**

Estimation data with at least one input signal and nonzero sample time.

For time domain estimation, `data` is an `iddata` object containing the input and output signal values.

For frequency domain estimation, `data` can be one of the following:

- Frequency response data (`frd` or `idfrd`)
- `iddata` object with its properties specified as follows:
  - `InputData` — Fourier transform of the input signal
  - `OutputData` — Fourier transform of the output signal
  - `Domain` — 'Frequency'

### N

Order of the FIR model. Must be one of the following:

- A positive integer.

For data containing  $N_u$  inputs and  $N_y$  outputs, you can also specify `N` as an  $N_y$ -by- $N_u$  matrix of positive integers, such that  $N(i,j)$  represents the length of impulse response from input  $j$  to output  $i$ .

- `[]` — Determines the order automatically using persistence of excitation analysis on the input data.

### NK

Transport delay in the estimated impulse response, specified as a scalar integer. For data containing  $N_u$  inputs and  $N_y$  outputs, you can also specify a  $N_y$ -by- $N_u$  matrix.

- To generate the impulse response coefficients for negative time values, which is useful for feedback analysis, use a negative integer. If you specify a negative value, the value must be the same across all output channels.

You can also use `NK = 'negative'` to automatically pick negative lags for all input/output channels of the model.

- Specify `NK = 0` if the delay is unknown. The true delay is then be indicated by insignificant impulse response values in the beginning of the response.
- Specify `NK = 1` to create a system whose leading numerator coefficient is zero.

Positive values of  $NK$  greater than 1 are stored in the `IODelay` property of `sys` (`sys.IODelay = max(NK-1, 0)`). Negative values are stored in the `InputDelay` property.

The impulse response (input  $j$  to output  $i$ ) coefficients correspond to the time span  $NK(i,j)*Ts : Ts : (N(i,j)+NK(i,j)-1)*Ts$ .

**Default:** `zeros(Ny, Nu)`

### options

Estimation options that specify the following:

- Prefilter order
- Regularization algorithm
- Input and output data offsets

Use `impulseestOptions` to create the options set.

## Output Arguments

### sys

Estimated impulse response model, returned as an `idtf` model, which encapsulates an FIR model.

Information about the estimation results and options used is stored in the `Report` property of the model. `Report` has the following fields:

Report Field	Description
Status	Summary of the model status, which indicates whether the model was created by construction or obtained by estimation.
Method	Estimation command used.

Report Field	Description																		
Fit	<p>Quantitative assessment of the estimation, returned as a structure. See “Loss Function and Model Quality Metrics” for more information on these quality metrics. The structure has the following fields:</p> <table border="1" data-bbox="795 487 1334 1269"> <thead> <tr> <th data-bbox="795 487 869 562">Field</th> <th data-bbox="869 487 1334 562">Description</th> </tr> </thead> <tbody> <tr> <td data-bbox="795 562 869 736">Fit Percent</td> <td data-bbox="869 562 1334 736">Normalized root mean squared error (NRMSE) measure of how well the response of the model fits the estimation data, expressed as a percentage.</td> </tr> <tr> <td data-bbox="795 736 869 843">Loss Function</td> <td data-bbox="869 736 1334 843">Value of the loss function when the estimation completes.</td> </tr> <tr> <td data-bbox="795 843 869 951">MSE</td> <td data-bbox="869 843 1334 951">Mean squared error (MSE) measure of how well the response of the model fits the estimation data.</td> </tr> <tr> <td data-bbox="795 951 869 996">FPE</td> <td data-bbox="869 951 1334 996">Final prediction error for the model.</td> </tr> <tr> <td data-bbox="795 996 869 1072">AIC</td> <td data-bbox="869 996 1334 1072">Raw Akaike Information Criteria (AIC) measure of model quality.</td> </tr> <tr> <td data-bbox="795 1072 869 1149">AICc</td> <td data-bbox="869 1072 1334 1149">Small sample-size corrected AIC.</td> </tr> <tr> <td data-bbox="795 1149 869 1225">nAIC</td> <td data-bbox="869 1149 1334 1225">Normalized AIC.</td> </tr> <tr> <td data-bbox="795 1225 869 1269">BIC</td> <td data-bbox="869 1225 1334 1269">Bayesian Information Criteria (BIC).</td> </tr> </tbody> </table>	Field	Description	Fit Percent	Normalized root mean squared error (NRMSE) measure of how well the response of the model fits the estimation data, expressed as a percentage.	Loss Function	Value of the loss function when the estimation completes.	MSE	Mean squared error (MSE) measure of how well the response of the model fits the estimation data.	FPE	Final prediction error for the model.	AIC	Raw Akaike Information Criteria (AIC) measure of model quality.	AICc	Small sample-size corrected AIC.	nAIC	Normalized AIC.	BIC	Bayesian Information Criteria (BIC).
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Parameters	Estimated values of model parameters.																		
OptionsUsed	Option set used for estimation. If no custom options were configured, this is a set of default options. See <code>impulseeestOptions</code> for more information.																		

<b>Report Field</b>	<b>Description</b>
RandState	State of the random number stream at the start of estimation. Empty, [], if randomization was not used during estimation. For more information, see <code>rng</code> in the MATLAB documentation.

Report Field	Description																
DataUsed	Attributes of the data used for estimation, returned as a structure with the following fields:																
	<table border="1"> <thead> <tr> <th data-bbox="795 427 869 496">Field</th> <th data-bbox="869 427 1334 496">Description</th> </tr> </thead> <tbody> <tr> <td data-bbox="795 496 869 574">Name</td> <td data-bbox="869 496 1334 574">Name of the data set.</td> </tr> <tr> <td data-bbox="795 574 869 652">Type</td> <td data-bbox="869 574 1334 652">Data type.</td> </tr> <tr> <td data-bbox="795 652 869 730">Length</td> <td data-bbox="869 652 1334 730">Number of data samples.</td> </tr> <tr> <td data-bbox="795 730 869 774">Ts</td> <td data-bbox="869 730 1334 774">Sample time.</td> </tr> <tr> <td data-bbox="795 774 869 1216">InterSample</td> <td data-bbox="869 774 1334 1216">                     Input intersample behavior, returned as one of the following values:                     <ul style="list-style-type: none"> <li>• 'zoh' — Zero-order hold maintains a piecewise-constant input signal between samples.</li> <li>• 'foh' — First-order hold maintains a piecewise-linear input signal between samples.</li> <li>• 'bl' — Band-limited behavior specifies that the continuous-time input signal has zero power above the Nyquist frequency.</li> </ul> </td> </tr> <tr> <td data-bbox="795 1216 869 1355">InputOffset</td> <td data-bbox="869 1216 1334 1355">Offset removed from time-domain input data during estimation. For nonlinear models, it is [ ].</td> </tr> <tr> <td data-bbox="795 1355 869 1496">OutputOffset</td> <td data-bbox="869 1355 1334 1496">Offset removed from time-domain output data during estimation. For nonlinear models, it is [ ].</td> </tr> </tbody> </table>	Field	Description	Name	Name of the data set.	Type	Data type.	Length	Number of data samples.	Ts	Sample time.	InterSample	Input intersample behavior, returned as one of the following values: <ul style="list-style-type: none"> <li>• 'zoh' — Zero-order hold maintains a piecewise-constant input signal between samples.</li> <li>• 'foh' — First-order hold maintains a piecewise-linear input signal between samples.</li> <li>• 'bl' — Band-limited behavior specifies that the continuous-time input signal has zero power above the Nyquist frequency.</li> </ul>	InputOffset	Offset removed from time-domain input data during estimation. For nonlinear models, it is [ ].	OutputOffset	Offset removed from time-domain output data during estimation. For nonlinear models, it is [ ].
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For more information on using Report, see “Estimation Report”.

## Examples

### Identify Nonparametric Impulse Response Model from Data

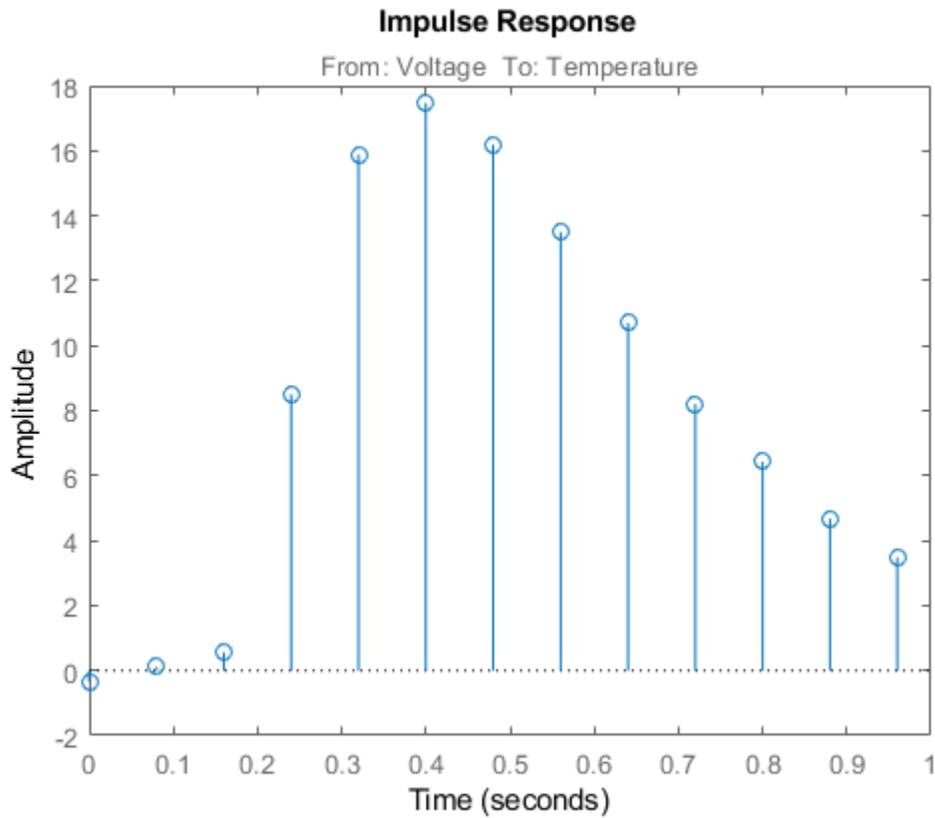
Compute a nonparametric impulse response model using data from a hair dryer. The input is the voltage applied to the heater and the output is the heater temperature. Use the first 500 samples for estimation.

```
load dry2
ze = dry2(1:500);
sys = impulseest(ze);
```

`ze` is an `iddata` object that contains time-domain data. `sys`, the identified nonparametric impulse response model, is an `idtf` model.

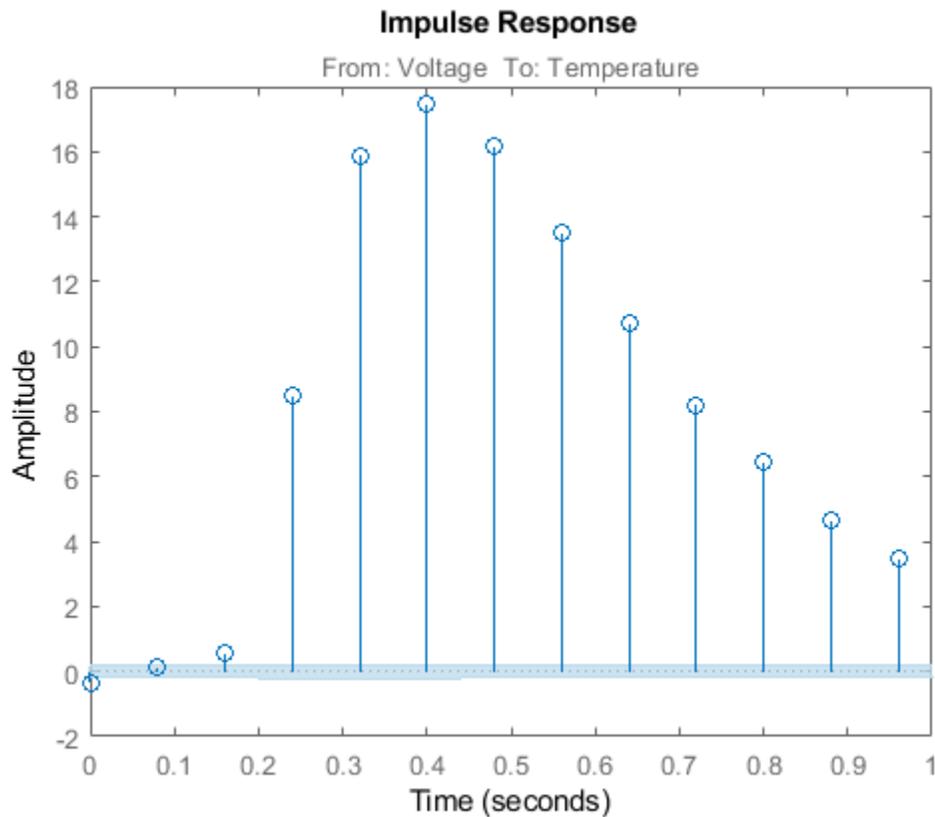
Analyze the impulse response of the identified model from time 0 to 1.

```
h = impulseplot(sys,1);
```



Right-click the plot and select **Characteristics > Confidence Region** to view the statistically zero-response region. Alternatively, you can use the `showConfidence` command.

```
showConfidence(h);
```



The first significantly nonzero response value occurs at 0.24 seconds, or, the third lag. This implies that the transport delay is 3 samples. To generate a model where the 3-sample delay is imposed, set the transport delay to 3:

```
sys = impulseest(ze,[],3)
```

### Specify Order of FIR Model

Load estimation data

```
load iddata3 z3;
```

Estimate a 35th order FIR model.

```
sys = impulseest(z3,35);
```

## Specify Transport Delay in FIR Model

Estimate an impulse response model with transport delay of 3 samples.

If you know about the presence of delay in the input/output data in advance, use the value as a transport delay for impulse response estimation.

Generate data with 3-sample input to output lag. Create a random input signal and use an `idpoly` model to simulate the output data.

```
u = rand(100,1);  
sys = idpoly([1 .1 .4],[0 0 0 4 -2],[1 1 .1]);  
opt = simOptions('AddNoise',true);  
y = sim(sys,u,opt);  
data = iddata(y,u,1);
```

Estimate a 20th order model with a 3-sample transport delay.

```
model = impulseest(data,20,3);
```

## Obtain Regularized Estimate of Impulse Response Model

Obtain regularized estimates of impulse response model using the regularizing kernel estimation option.

Estimate a model using regularization.

```
load iddata3 z3;  
sys1 = impulseest(z3);
```

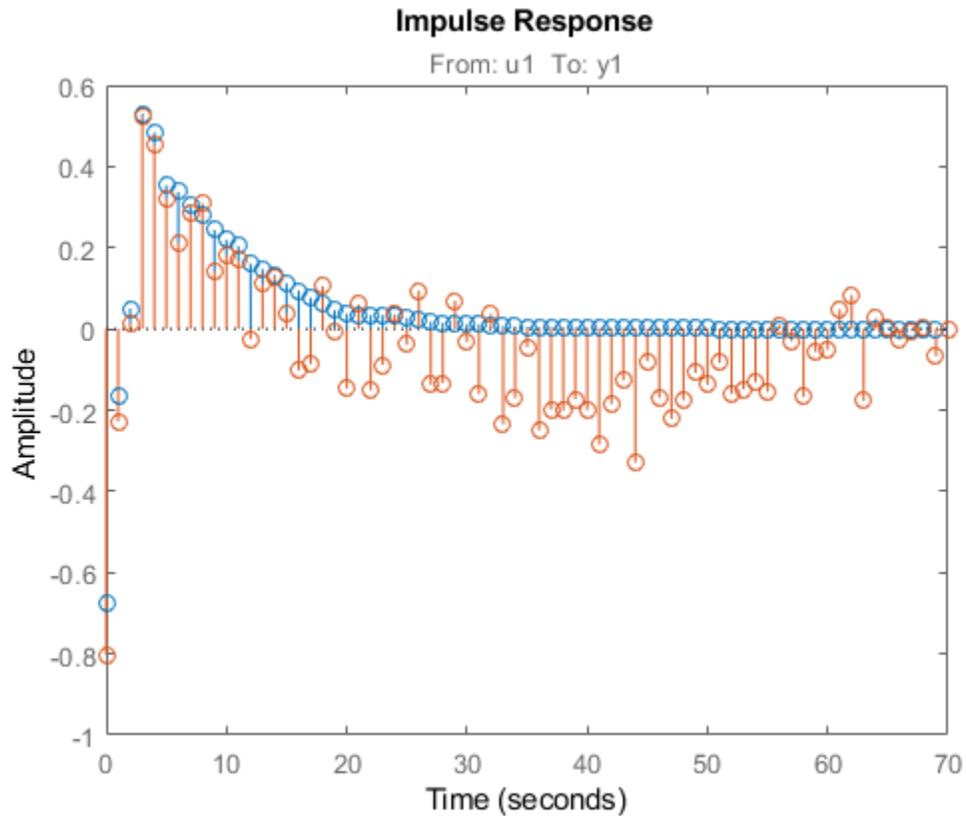
By default, tuned and correlated kernel ('TC') is used for regularization.

Estimate a model with no regularization.

```
opt = impulseestOptions('RegularizationKernel','none');  
sys2 = impulseest(z3,opt);
```

Compare the impulse response of both models.

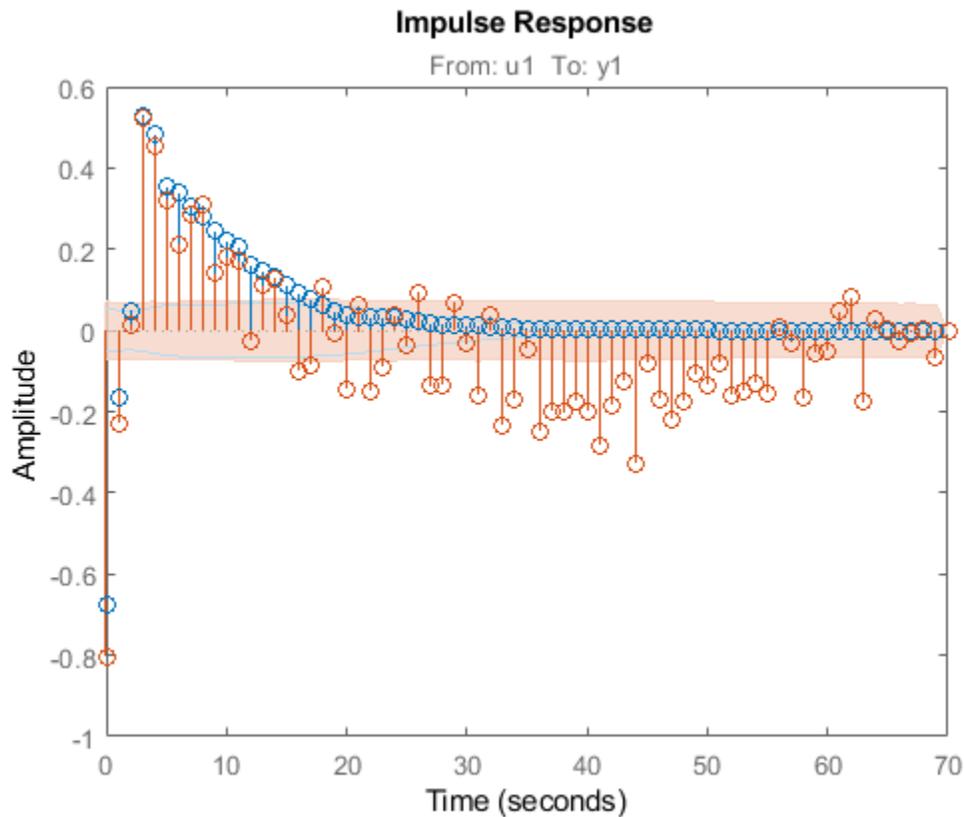
```
h = impulseplot(sys1,sys2,70);
```



As the plot shows, using regularization makes the response smoother.

Plot the confidence interval.

```
showConfidence(h);
```



The uncertainty in the computed response is reduced at larger lags for the model using regularization. Regularization decreases variance at the price of some bias. The tuning of the regularization is such that the bias is dominated by the variance error though.

### Use Regularized Impulse Response Model to Estimate State-Space Model

Load data.

```
load regularizationExampleData eData;
```

Create a transfer function model used for generating the estimation data (true system).

```
trueSys = idtf([0.02008 0.04017 0.02008],[1 -1.561 0.6414],1);
```

Obtain regularized impulse response (FIR) model.

```
opt = impulseestOptions('RegularizationKernel','DC');  
m0 = impulseest(eData,70,opt);
```

Convert the model into a state-space model and reduce the model order.

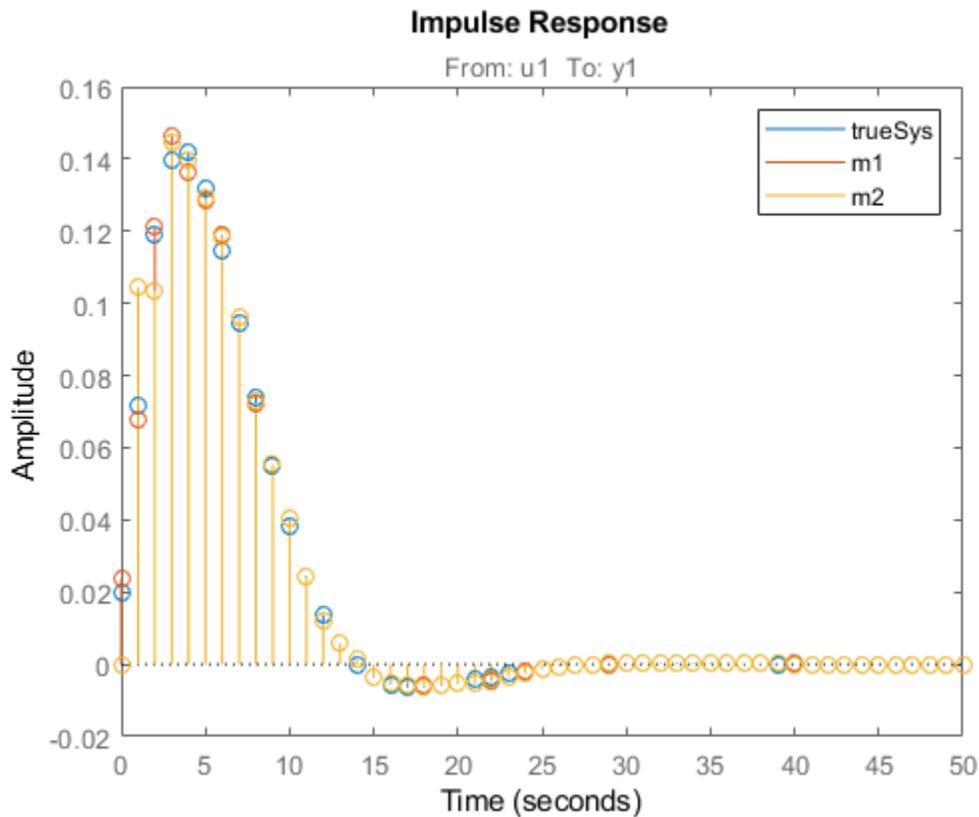
```
m1 = balred(idss(m0),15);
```

Obtain a second state-space model using regularized reduction of an ARX model.

```
m2 = ssregest(eData,15);
```

Compare the impulse responses of the true system and the estimated models.

```
impulse(trueSys,m1,m2,50);  
legend('trueSys','m1','m2');
```



**Test Measured Data for Feedback Effects**

Use the empirical impulse response of the measured data to verify whether there are feedback effects. Significant amplitude of the impulse response for negative time values indicates feedback effects in data.

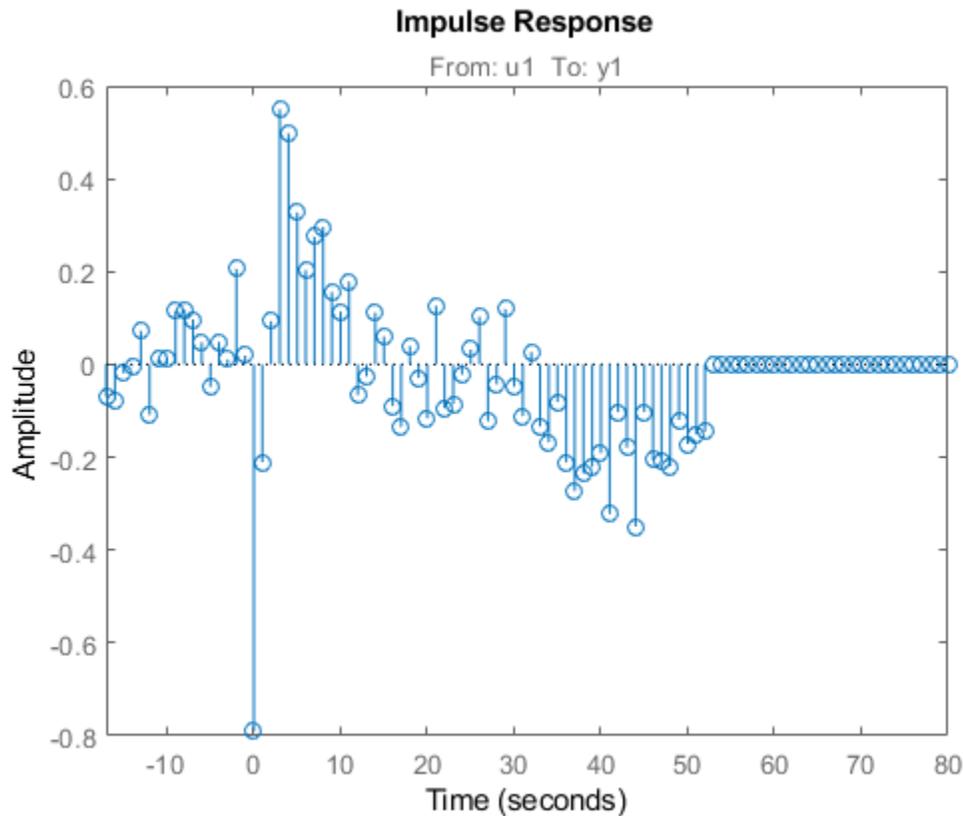
Compute the noncausal impulse response using a fourth-order prewhitening filter, automatically chosen order and negative lag using nonregularized estimation.

```
load iddata3 z3;  
opt = impulseeestOptions('pw',4,'RegularizationKernel','none');  
sys = impulseeest(z3,[],'negative',opt);
```

sys is a noncausal model containing response values for negative time.

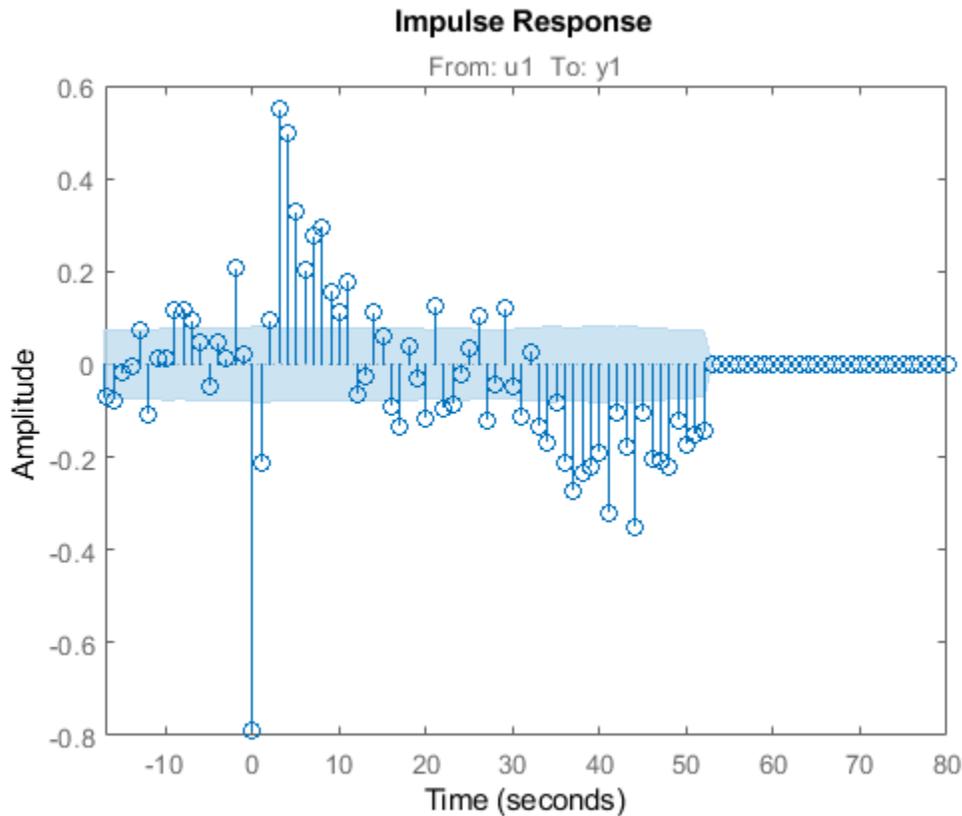
Analyze the impulse response of the identified model.

```
h = impulseplot(sys);
```



View the statistically zero-response region by right-clicking on the plot and selecting **Characteristics > Confidence Region**. Alternatively, you can use the showConfidence command.

```
showConfidence(h);
```



The large response value at  $t=0$  (zero lag) suggests that the data comes from a process containing feedthrough. That is, the input affects the output instantaneously. There could also be a direct feedback effect (proportional control without some delay that  $u(t)$  is determined partly by  $y(t)$ ).

Also, the response values are significant for some negative time lags, such as at -7 seconds and -9 seconds. Such significant negative values suggest the possibility of feedback in the data.

### Compute Impulse Response on Frequency Response Data

Compute an impulse response model for frequency response data.

```
load demofr;  
zfr = AMP.*exp(1i*PHA*pi/180);  
Ts = 0.1;  
data = idfrd(zfr,W,Ts);  
sys = impulseest(data);
```

### Compare Identified Nonparametric and Parametric Models

Identify parametric and nonparametric models for a data set, and compare their step response.

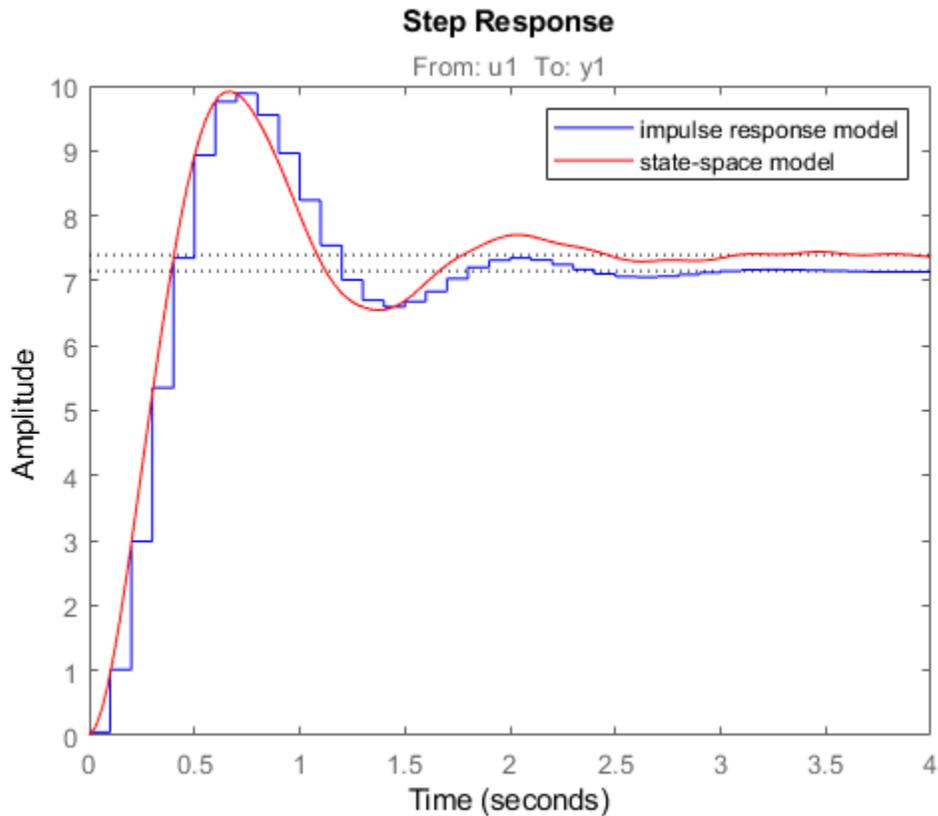
Identify the impulse response model (nonparametric) and state-space model (parametric), based on a data set.

```
load iddata1 z1;  
sys1 = impulseest(z1);  
sys2 = ssest(z1,4);
```

sys1 is a discrete-time identified transfer function model. sys2 is a continuous-time identified state-space model.

Compare the step response for sys1 and sys2.

```
step(sys1,'b',sys2,'r');  
legend('impulse response model','state-space model');
```



## Tips

- To view the impulse or step response of `sys`, use either `impzplot` or `stepplot`, respectively.
- A significant value of the impulse response of `sys` for negative time values indicates the presence of feedback in the data.
- To view the region of insignificant impulse response (statistically zero) in a plot, right-click on the plot and select **Characteristics > Confidence Region**. A patch depicting the zero-response region appears on the plot. The impulse response at any time value is significant only if it lies outside the zero response region. The level of significance

depends on the number of standard deviations specified in ShowConfidence or options in the property editor. A common choice is 3 standard deviations, which gives 99.7% significance.

## Algorithms

Correlation analysis refers to methods that estimate the impulse response of a linear model, without specific assumptions about model orders.

The impulse response,  $g$ , is the system's output when the input is an impulse signal. The output response to a general input,  $u(t)$ , is obtained as the convolution with the impulse response. In continuous time:

$$y(t) = \int_{-\infty}^t g(\tau) u(t - \tau) d\tau$$

In discrete-time:

$$y(t) = \sum_{k=1}^{\infty} g(k) u(t - k)$$

The values of  $g(k)$  are the discrete time impulse response coefficients.

You can estimate the values from observed input-output data in several different ways. `impulseeest` estimates the first  $n$  coefficients using the least-squares method to obtain a finite impulse response (FIR) model of order  $n$ .

Several important options are associated with the estimate:

- **Prewhitening** — The input can be pre-whitened by applying an input-whitening filter of order `PW` to the data. This minimizes the effect of the neglected tail ( $k > n$ ) of the impulse response.

- 1 A filter of order `PW` is applied such that it whitens the input signal  $u$ :

$$1/A = A(u) e, \text{ where } A \text{ is a polynomial and } e \text{ is white noise.}$$

- 2 The inputs and outputs are filtered using the filter:

$$u_f = Au, y_f = Ay$$

3 The filtered signals  $u_f$  and  $y_f$  are used for estimation.

You can specify prewhitening using the `PW` name-value pair argument of `impulseestOptions`.

- **Regularization** — The least-squares estimate can be regularized. This means that a prior estimate of the decay and mutual correlation among  $g(k)$  is formed and used to merge with the information about  $g$  from the observed data. This gives an estimate with less variance, at the price of some bias. You can choose one of the several kernels to encode the prior estimate.

This option is essential because, often, the model order  $n$  can be quite large. In cases where there is no regularization,  $n$  can be automatically decreased to secure a reasonable variance.

You can specify the regularizing kernel using the `RegularizationKernel` Name-Value pair argument of `impulseestOptions`.

- **Autoregressive Parameters** — The basic underlying FIR model can be complemented by `NA` autoregressive parameters, making it an ARX model.

$$y(t) = \sum_{k=1}^n g(k)u(t-k) - \sum_{k=1}^{NA} a_k y(t-k)$$

This gives both better results for small  $n$  and allows unbiased estimates when data are generated in closed loop. `impulseest` uses `NA = 5` for  $t > 0$  and `NA = 0` (no autoregressive component) for  $t < 0$ .

- **Noncausal effects** — Response for negative lags. It may happen that the data has been generated partly by output feedback:

$$u(t) = \sum_{k=0}^{\infty} h(k)y(t-k) + r(t)$$

where  $h(k)$  is the impulse response of the regulator and  $r$  is a setpoint or disturbance term. The existence and character of such feedback  $h$  can be estimated in the same way as  $g$ , simply by trading places between  $y$  and  $u$  in the estimation call. Using `impulseest` with an indication of negative delays,

`mi = impulseest(data, nk, nb)`,  $nk < 0$ , returns a model `mi` with an impulse response

$$[h(-nk), h(-nk-1), \dots, h(0), g(1), g(2), \dots, g(nb+nk)]$$

aligned so that it corresponds to lags  $[nk, nk + 1, \dots, 0, 1, 2, \dots, nb + nk]$ . This is achieved because the input delay (`InputDelay`) of model `mi` is `nk`.

For a multi-input multi-output system, the impulse response  $g(k)$  is an  $ny$ -by- $nu$  matrix, where  $ny$  is the number of outputs and  $nu$  is the number of inputs. The  $i$ - $j$  element of the matrix  $g(k)$  describes the behavior of the  $i$ th output after an impulse in the  $j$ th input.

## See Also

[cra](#) | [impulse](#) | [impulseestOptions](#) | [spa](#) | [step](#)

## Topics

“What Is Time-Domain Correlation Analysis?”

**Introduced in R2012a**

## impulseestOptions

Options set for `impulseest`

### Syntax

```
options = impulseestOptions  
options = impulseestOptions(Name,Value)
```

### Description

`options = impulseestOptions` creates a default options set for `impulseest`.

`options = impulseestOptions(Name,Value)` creates an options set with the options specified by one or more `Name,Value` pair arguments.

### Input Arguments

#### Name-Value Pair Arguments

Specify optional comma-separated pairs of `Name,Value` arguments. `Name` is the argument name and `Value` is the corresponding value. `Name` must appear inside quotes. You can specify several name and value pair arguments in any order as `Name1,Value1,...,NameN,ValueN`.

#### RegularizationKernel

Regularizing kernel, used for regularized estimates of impulse response for all input-output channels. Regularization reduces variance of estimated model coefficients and produces a smoother response by trading variance for bias. For more information, see [1].

Regularization is specified as one of the following values:

- 'TC' — Tuned and correlated kernel

- 'none' — No regularization is used
- 'CS' — Cubic spline kernel
- 'SE' — Squared exponential kernel
- 'SS' — Stable spline kernel
- 'HF' — High frequency stable spline kernel
- 'DI' — Diagonal kernel
- 'DC' — Diagonal and correlated kernel

**Default:** 'TC'

### **PW**

Order of the input prewhitening filter. Must be one of the following:

- 'auto' — Uses a filter of order 10 when `RegularizationKernel` is 'none'; otherwise, 0.
- Nonnegative integer

Use a nonzero value of prewhitening only for unregularized estimation (`RegularizationKernel` is 'none').

**Default:** 'auto'

### **InputOffset**

Input signal offset level of time-domain estimation data. Must be one of the following:

- An  $N_u$ -element column vector, where  $N_u$  is the number of inputs. For multi-experiment data, specify a  $N_u$ -by- $N_e$  matrix, where  $N_e$  is the number of experiments. The offset value `InputOffset(i, j)` is subtracted from the  $i^{\text{th}}$  input signal of the  $j^{\text{th}}$  experiment.
- [] — No offsets.

**Default:** []

### **OutputOffset**

Output signal offset level of time-domain estimation data. Must be one of the following:

- An  $N_y$ -element column vector, where  $N_y$  is the number of outputs. For multi-experiment data, specify a  $N_y$ -by- $N_e$  matrix, where  $N_e$  is the number of experiments. The offset value `OutputOffset(i, j)` is subtracted from the  $i^{\text{th}}$  output signal of the  $j^{\text{th}}$  experiment.
- `[]` — No offsets.

**Default:** `[]`

### Advanced

Structure, used during regularized estimation, with the following fields:

- `MaxSize` — Maximum allowable size of Jacobian matrices formed during estimation. Specify a large positive number.

**Default:** `250e3`

- `SearchMethod` — Search method for estimating regularization parameters, specified as one of the following values:
  - `'fmincon'`: Trust-region-reflective constrained minimizer. In general, `'fmincon'` is better than `'gn'` for handling bounds on regularization parameters that are imposed automatically during estimation.
  - `'gn'`: Quasi-Newton line search.

`SearchMethod` is used only when `RegularizationKernel` is not `'none'`.

**Default:** `'fmincon'`

- `AROrder` — Order of the AR-part in the model from input to output. Specify as a positive integer.

An `order>0` allows more accurate models of the impulse response in case of feedback and non-white output disturbances.

**Default:** `5`

- `FeedthroughInSys` — Specify whether the impulse response value at zero lag must be attributed to feedthrough in the system (`true`) or to feedback effects (`false`). Applies only when you compute the response values for negative lags.

**Default:** `false`

## Output Arguments

### options

Option set containing the specified options for `impulseest`.

## Examples

### Create Default Options Set for Impulse Response Estimation

Create a default options set for `impulseest`.

```
options = impulseestOptions;
```

### Specify Regularizing Kernel and Prewhitening Options for Impulse Response Estimation

Specify 'HF' regularizing kernel and order of prewhitening filter for `impulseest`.

```
options = impulseestOptions('RegularizationKernel','HF','PW',5);
```

Alternatively, use dot notation to specify these options.

```
options = impulseestOptions;  
options.RegularizationKernel = 'HF';  
options.PW = 5;
```

## Tips

- A linear model cannot describe arbitrary input-output offsets. Therefore, before using the data, you must either detrend it or remove the levels using `InputOffset` and `OutputOffset`. You can reintroduce the removed data during simulations by using the `InputOffset` and `OutputOffset` simulation options. For more information, see `simOptions`.

- Estimating the impulse response by specifying either `InputOffset`, `OutputOffset` or both is equivalent to detrending the data using `getTrend` and `detrend`. For example:

```
opt = impulseestOptions('InputOffset',in_off,'OutputOffset',out_off);  
impulseest(data,opt);
```

is the same as:

```
Tr = getTrend(data),  
Tr.InputOffset = in_off  
TR.OutputOffset = out_off  
dataT = detrend(data,Tr)  
impulseest(dataT)
```

## References

- [1] T. Chen, H. Ohlsson, and L. Ljung. “On the Estimation of Transfer Functions, Regularizations and Gaussian Processes - Revisited”, *Automatica*, Volume 48, August 2012.

## See Also

`impulseest`

**Introduced in R2012b**

# impulseplot

Plot impulse response and return plot handle

## Syntax

```
impulseplot(sys)
impulseplot(sys,Tfinal)
impulseplot(sys,t)
impulseplot(sys1,sys2,...,sysN)
impulseplot(sys1,sys2,...,sysN,Tfinal)
impulseplot(sys1,sys2,...,sysN,t)
impulseplot(AX,...)
impulseplot(..., plotoptions)
h = impulseplot(...)
```

## Description

`impulseplot` plots the impulse response of the dynamic system model `sys`. For multi-input models, independent impulse commands are applied to each input channel. The time range and number of points are chosen automatically. For continuous systems with direct feedthrough, the infinite pulse at  $t=0$  is disregarded. `impulseplot` can also return the plot handle, `h`. You can use this handle to customize the plot with the `getoptions` and `setoptions` commands. Type

```
help timeoptions
```

for a list of available plot options.

`impulseplot(sys)` plots the impulse response of the LTI model without returning the plot handle.

`impulseplot(sys,Tfinal)` simulates the impulse response from  $t = 0$  to the final time  $t = Tfinal$ . Express `Tfinal` in the system time units, specified in the `TimeUnit` property of `sys`. For discrete-time systems with unspecified sample time ( $Ts = -1$ ), `impulseplot` interprets `Tfinal` as the number of sampling intervals to simulate.

`impzplot(sys,t)` uses the user-supplied time vector `t` for simulation. Express `t` in the system time units, specified in the `TimeUnit` property of `sys`. For discrete-time models, `t` should be of the form `Ti:Ts:Tf`, where `Ts` is the sample time. For continuous-time models, `t` should be of the form `Ti:dt:Tf`, where `dt` becomes the sample time of a discrete approximation to the continuous system (see `impz`). The `impzplot` command always applies the impulse at `t=0`, regardless of `Ti`.

To plot the impulse response of multiple LTI models `sys1,sys2,...` on a single plot, use:

```
impzplot(sys1,sys2,...,sysN)
```

```
impzplot(sys1,sys2,...,sysN,Tfinal)
```

```
impzplot(sys1,sys2,...,sysN,t)
```

You can also specify a color, line style, and marker for each system, as in

```
impzplot(sys1,'r',sys2,'y--',sys3,'gx')
```

`impzplot(AX,...)` plots into the axes with handle `AX`.

`impzplot(..., plotoptions)` plots the impulse response with the options specified in `plotoptions`. Type

```
help timeoptions
```

for more detail.

`h = impzplot(...)` plots the impulse response and returns the plot handle `h`.

## Examples

### Example 1

Normalize the impulse response of a third-order system.

```
sys = rss(3);  
h = impzplot(sys);  
% Normalize responses  
setoptions(h,'Normalize','on');
```

## Example 2

Plot the impulse response and the corresponding 1 std "zero interval" of an identified linear system.

```
load(fullfile(matlabroot, 'toolbox', 'ident', 'iddemos', 'data', 'dcmotordata'));
z = iddata(y, u, 0.1, 'Name', 'DC-motor');
set(z, 'InputName', 'Voltage', 'InputUnit', 'V');
set(z, 'OutputName', {'Angular position', 'Angular velocity'});
set(z, 'OutputUnit', {'rad', 'rad/s'});
set(z, 'Tstart', 0, 'TimeUnit', 's');
model = n4sid(z,4,n4sidOptions('Focus', 'simulation'));
h = impulseplot(model,2);
showConfidence(h);
```

## Tips

You can change the properties of your plot, for example the units. For information on the ways to change properties of your plots, see “Ways to Customize Plots” (Control System Toolbox).

## See Also

`getoptions` | `impulse` | `setoptions` | `showConfidence`

**Introduced in R2012a**

## init

Set or randomize initial parameter values

## Syntax

```
m = init(m0)
m = init(m0,R,pars,sp)
```

## Description

`m = init(m0)` randomizes initial parameter estimates for model structures `m0` for any linear or nonlinear identified model. It does not support `idnlgrey` models. `m` is the same model structure as `m0`, but with a different nominal parameter vector. This vector is used as the initial estimate by `pem`.

`m = init(m0,R,pars,sp)` randomizes parameters around `pars` with variances given by the row vector `R`. Parameter number  $k$  is randomized as  $\text{pars}(k) + e \cdot \sqrt{R(k)}$ , where  $e$  is a normal random variable with zero mean and a variance of 1. The default value of `R` is all ones, and the default value of `pars` is the nominal parameter vector in `m0`.

Only models that give stable predictors are accepted. If `sp = 'b'`, only models that are both stable and have stable predictors are accepted.

`sp = 's'` requires stability only of the model, and `sp = 'p'` requires stability only of the predictor. `sp = 'p'` is the default.

Sufficiently free parameterizations can be stabilized by direct means without any random search. To just stabilize such an initial model, set `R = 0`. With `R > 0`, randomization is also done.

For model structures where a random search is necessary to find a stable model/predictor, a maximum of 100 trials is made by `init`. It can be difficult to find a stable predictor for high-order systems by trial and error.

## See Also

`idnlarx` | `idnlhw` | `rsmple` | `simsd`

**Introduced before R2006a**

## initialize

Initialize the state of the particle filter

### Syntax

```
initialize(pf,numParticles,mean,covariance)
initialize(pf,numParticles,stateBounds)
initialize( ____,Name,Value)
```

### Description

`initialize(pf,numParticles,mean,covariance)` initializes a particle filter object with a specified number of particles. The initial states of the particles in the state space are determined by sampling from the multivariate normal distribution with the specified mean and covariance. The number of state variables (`NumStateVariables`) is retrieved automatically based on the length of the mean vector.

`initialize(pf,numParticles,stateBounds)` determines the initial location of `numParticles` particles by sampling from the multivariate uniform distribution with the given `stateBounds`.

`initialize( ____,Name,Value)` initializes the particles with additional options specified by one or more `Name,Value` pair arguments.

### Examples

#### Initialize Particle Filter Object for Online State Estimation

To create a particle filter object for estimating the states of your system, create appropriate state transition function and measurement function for the system.

In this example, the functions `vdpParticleFilterStateFcn` and `vdpMeasurementLikelihoodFcn` describe a discrete-approximation to van der Pol oscillator with nonlinearity parameter, `mu`, equal to 1.

Create the particle filter object. Use function handles to provide the state transition and measurement likelihood functions to the object.

```
myPF = particleFilter(@vdpParticleFilterStateFcn,@vdpMeasurementLikelihoodFcn);
```

Initialize the particle filter at state [2; 0] with unit covariance, and use 1000 particles.

```
initialize(myPF, 1000, [2;0], eye(2));
```

```
myPF
```

```
myPF =
```

```
  particleFilter with properties:
```

```
      NumStateVariables: 2
      NumParticles: 1000
      StateTransitionFcn: @vdpParticleFilterStateFcn
      MeasurementLikelihoodFcn: @vdpMeasurementLikelihoodFcn
      IsStateVariableCircular: [0 0]
      ResamplingPolicy: [1x1 particleResamplingPolicy]
      ResamplingMethod: 'multinomial'
      StateEstimationMethod: 'mean'
      StateOrientation: 'column'
      Particles: [2x1000 double]
      Weights: [1x1000 double]
      State: 'Use the getStateEstimate function to see the value.'
      StateCovariance: 'Use the getStateEstimate function to see the value.'
```

To estimate the states and state estimation error covariance from the constructed object, use the `predict` and `correct` commands.

## Input Arguments

### **pf** — Particle filter

particleFilter object

Particle filter, specified as a object. See `particleFilter` for more information.

### **numParticles** — Number of particles used in the filter

scalar

Number of particles used in the filter, specified as a scalar.

Unless performance is an issue, do not use fewer than 1000 particles. A higher number of particles can improve the estimate but sacrifices performance speed, because the algorithm has to process more particles. Tuning the number of particles is the best way to improve the tracking of your particle filter.

### **mean — Mean of particle distribution**

vector

Mean of particle distribution, specified as a vector. The `NumStateVariables` property of `pf` is set based on the length of this vector.

### **covariance — Covariance of particle distribution**

*N*-by-*N* matrix

Covariance of particle distribution, specified as an *N*-by-*N* matrix, where *N* is the value of `NumStateVariables` property from `pf`.

### **stateBounds — Bounds of state variables**

*n*-by-2 matrix

Bounds of state variables, specified as an *n*-by-2 matrix. The `NumStateVariables` property of `pf` is set based on the value of *n*. Each row corresponds to the lower and upper limit of the corresponding state variable. The number of state variables (`NumStateVariables`) is retrieved automatically based on the number of rows of the `stateBounds` array.

## **Name-Value Pair Arguments**

Specify optional comma-separated pairs of `Name`, `Value` arguments. `Name` is the argument name and `Value` is the corresponding value. `Name` must appear inside single quotes ( ' '). You can specify several name and value pair arguments in any order as `Name1, Value1, ..., NameN, ValueN`.

Example: `... 'StateOrientation', 'row'`

### **CircularVariables — Circular variables**

logical vector

Circular variables, the comma-separated pair consisting of `CircularVariables` and specified as a logical vector. Each state variable that uses circular or angular coordinates is indicated with a 1. The length of the vector is equal to the `NumStateVariables` property of `particleFilter`.

## **StateOrientation — Orientation of states**

'column' (default) | 'row'

Orientation of states, specified as the comma-separated pair consisting of `StateOrientation` as one of these values: 'column' or 'row'. If it is 'column', `State` property and `getStateEstimate` method of the object `pf` returns the states as a column vector, and the `Particles` property has dimensions `NumStateVariables-by-NumParticles`. If it is 'row', the states have the row orientation and `Particles` has dimensions `NumParticles-by-NumStateVariables`.

## **See Also**

`clone` | `correct` | `extendedKalmanFilter` | `particleFilter` | `predict` | `unscentedKalmanFilter`

## **Topics**

“Nonlinear State Estimation Using Unscented Kalman Filter and Particle Filter”

“What Is Online Estimation?”

“Generate Code for Online State Estimation in MATLAB”

**Introduced in R2017b**

## interp

Interpolate FRD model

### Syntax

```
isys = interp(sys, freqs)
```

### Description

`isys = interp(sys, freqs)` interpolates the frequency response data contained in the FRD model `sys` at the frequencies `freqs`. `interp`, which is an overloaded version of the MATLAB function `interp`, uses linear interpolation and returns an FRD model `isys` containing the interpolated data at the new frequencies `freqs`. If `sys` is an IDFRD model, the noise spectrum, if non-empty, is also interpolated. The response and noise covariance data, if available, are also interpolated.

You should express the frequency values `freqs` in the same units as `sys`. frequency. The frequency values must lie between the smallest and largest frequency points in `sys` (extrapolation is not supported).

### See Also

`frd` | `freqresp` | `idfrd`

**Introduced in R2012a**

# iopzmap

Plot pole-zero map for I/O pairs of model

## Syntax

```
iopzmap(sys)  
iopzmap(sys1,sys2,...)
```

## Description

`iopzmap(sys)` computes and plots the poles and zeros of each input/output pair of the dynamic system model `sys`. The poles are plotted as x's and the zeros are plotted as o's.

`iopzmap(sys1,sys2,...)` shows the poles and zeros of multiple models `sys1,sys2,...` on a single plot. You can specify distinctive colors for each model, as in `iopzmap(sys1, 'r',sys2, 'y',sys3, 'g')`.

The functions `sgrid` or `zgrid` can be used to plot lines of constant damping ratio and natural frequency in the  $s$  or  $z$  plane.

For model arrays, `iopzmap` plots the poles and zeros of each model in the array on the same diagram.

## Examples

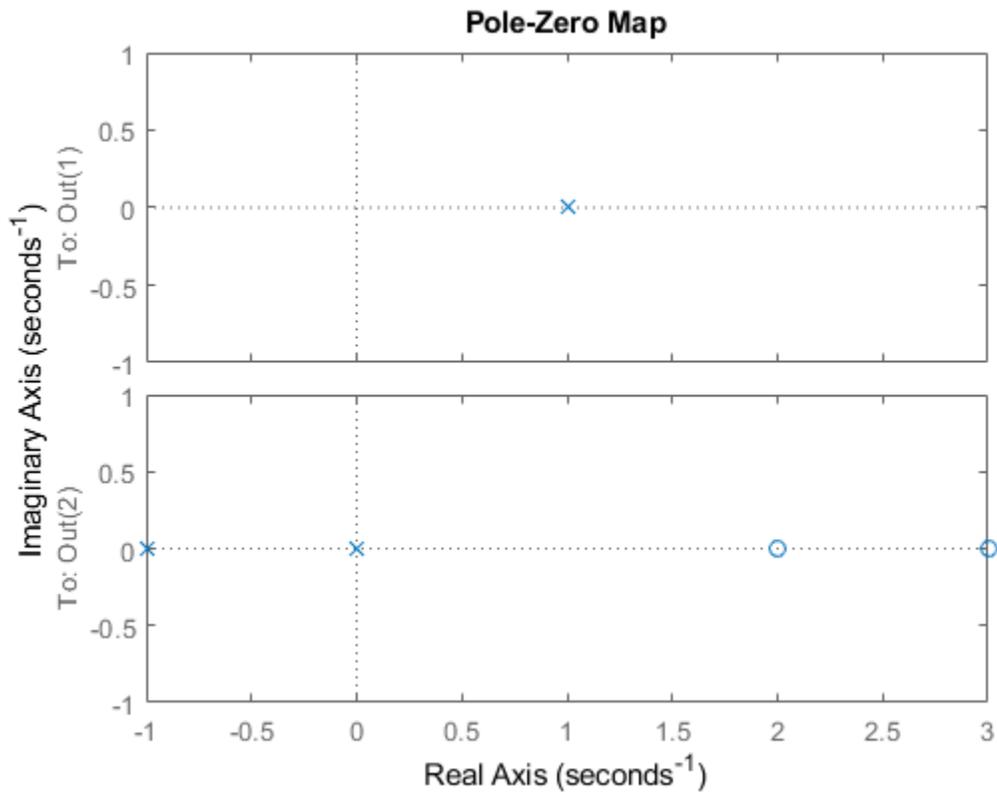
### Pole-Zero Map for MIMO System

Create a one-input, two-output dynamic system.

```
H = [tf(-5 ,[1 -1]); tf([1 -5 6],[1 1 0])];
```

Plot a pole-zero map.

```
iopzmap(H)
```

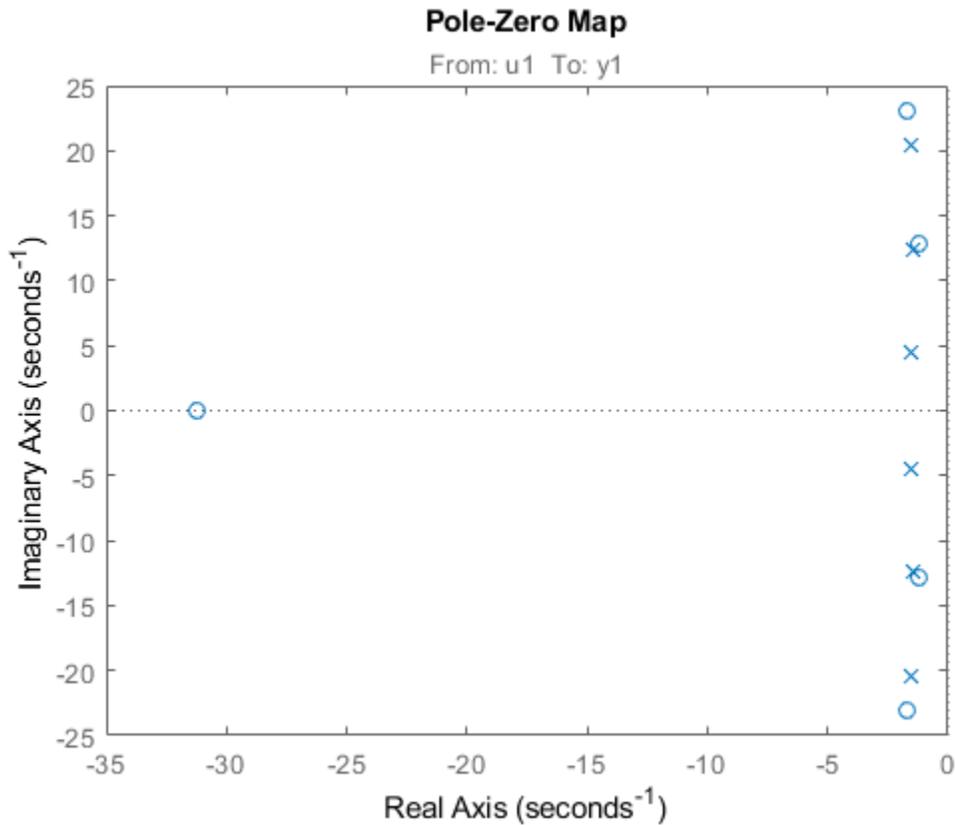


`iopzmap` generates a separate map for each I/O pair in the system.

### Pole-Zero Map of Identified Model

View the poles and zeros of an over-parameterized state-space model estimated from input-output data. (Requires System Identification Toolbox™).

```
load iddata1
sys = ssest(z1,6,ssestOptions('focus','simulation'));
iopzmap(sys)
```



The plot shows that there are two pole-zero pairs that almost overlap, which hints are their potential redundancy.

## Tips

For additional options for customizing the appearance of the pole-zero plot, use `iopzplot`.

## See Also

`iopzplot` | `pole` | `pzmap` | `sgrid` | `zero` | `zgrid`

**Introduced in R2012a**

# iopzplot

Plot pole-zero map for I/O pairs and return plot handle

## Syntax

```
h = iopzplot(sys)
iopzplot(sys1,sys2,...)
iopzplot(AX,...)
iopzplot(..., plotoptions)
```

## Description

`h = iopzplot(sys)` computes and plots the poles and zeros of each input/output pair of the dynamic system model `sys`. The poles are plotted as x's and the zeros are plotted as o's. It also returns the plot handle `h`. You can use this handle to customize the plot with the `getoptions` and `setoptions` commands. Type

```
help pzoptions
```

for a list of available plot options. For more information on the ways to change properties of your plots, see “Ways to Customize Plots” (Control System Toolbox).

`iopzplot(sys1,sys2,...)` shows the poles and zeros of multiple dynamic system models `sys1,sys2,...` on a single plot. You can specify distinctive colors for each model, as in

```
iopzplot(sys1,'r',sys2,'y',sys3,'g')
```

`iopzplot(AX,...)` plots into the axes with handle `AX`.

`iopzplot(..., plotoptions)` plots the poles and zeros with the options specified in `plotoptions`. Type

```
help pzoptions
```

for more detail.

The function `sgrid` or `zgrid` can be used to plot lines of constant damping ratio and natural frequency in the  $s$  or  $z$  plane.

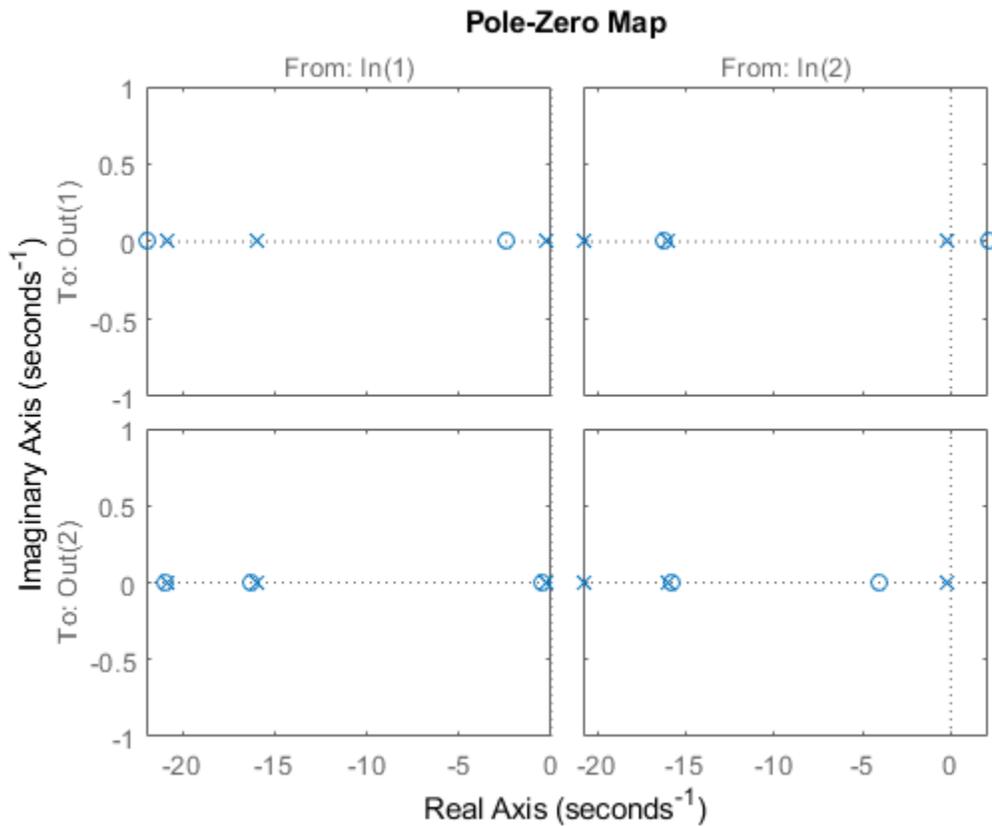
For arrays `sys` of LTI models, `iopzplot` plots the poles and zeros of each model in the array on the same diagram.

## Examples

### Change I/O Grouping on Pole/Zero Map

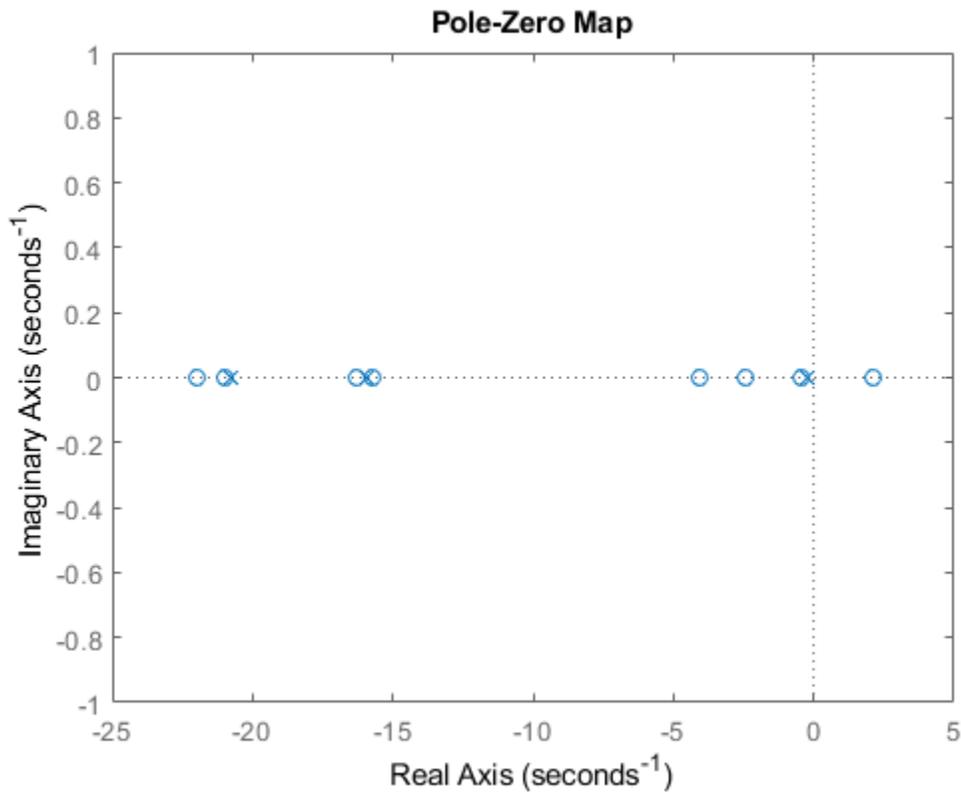
Create a pole/zero map of a two-input, two-output dynamic system.

```
sys = rss(3,2,2);  
h = iopzplot(sys);
```



By default, the plot displays the poles and zeros of each I/O pair on its own axis. Use the plot handle to view all I/Os on a single axis.

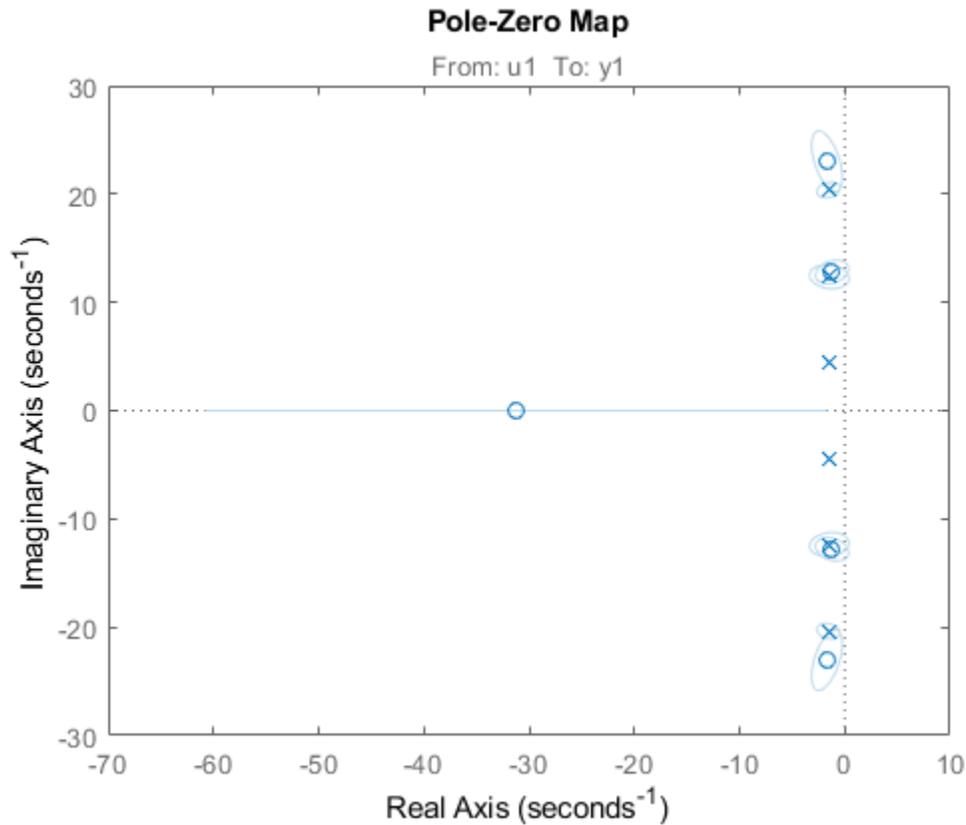
```
setoptions(h, 'IOGrouping', 'all')
```



### Use Pole-Zero Map to Examine Identified Model

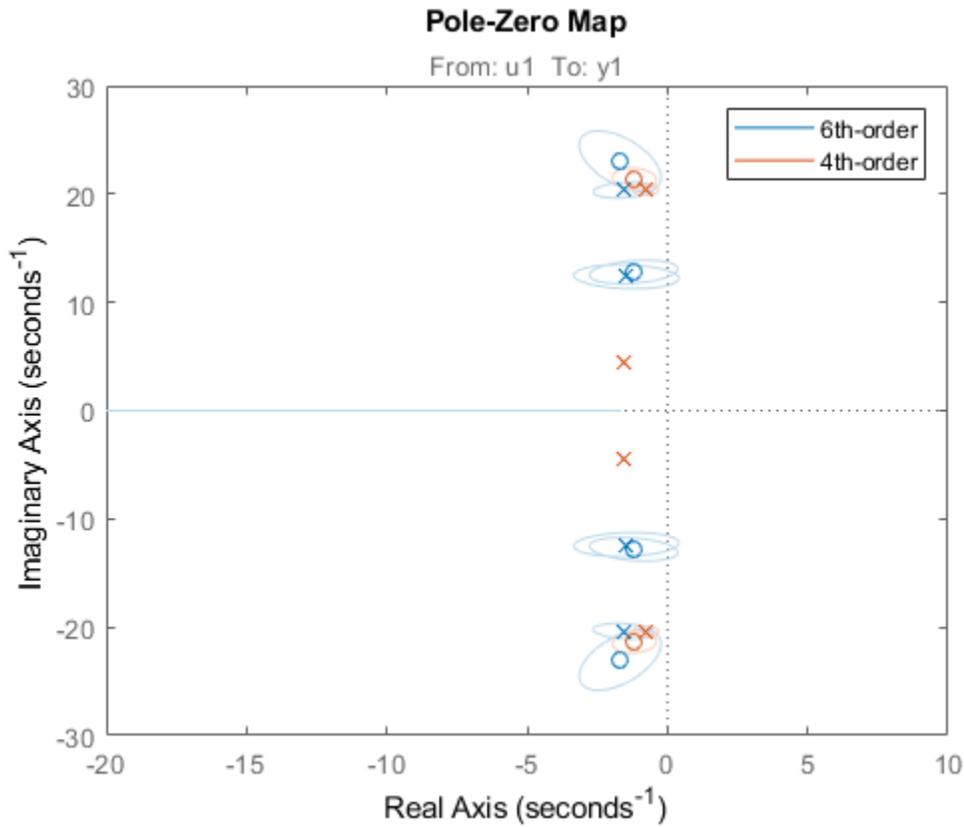
View the poles and zeros of a sixth-order state-space model estimated from input-output data. Use the plot handle to display the confidence intervals of the identified model's pole and zero locations.

```
load iddata1
sys = ssest(z1,6,ssestOptions('focus','simulation'));
h = iopzplot(sys);
showConfidence(h)
```



There is at least one pair of complex-conjugate poles whose locations overlap with those of a complex zero, within the 1- $\sigma$  confidence region. This suggests their redundancy. Hence, a lower (4th) order model might be more robust for the given data.

```
sys2 = ssest(z1,4,ssestOptions('focus','simulation'));
h = iopzplot(sys,sys2);
showConfidence(h)
legend('6th-order','4th-order')
axis([-20, 10 -30 30])
```



The fourth-order model sys2 shows less variability in the pole-zero locations.

## See Also

`getoptions` | `iopzmap` | `setoptions` | `showConfidence`

**Introduced in R2012a**

## isct

Determine if dynamic system model is in continuous time

### Syntax

```
bool = isct(sys)
```

### Description

`bool = isct(sys)` returns a logical value of 1 (`true`) if the dynamic system model `sys` is a continuous-time model. The function returns a logical value of 0 (`false`) otherwise.

### Input Arguments

#### **sys**

Dynamic system model or array of such models.

### Output Arguments

#### **bool**

Logical value indicating whether `sys` is a continuous-time model.

`bool = 1 (true)` if `sys` is a continuous-time model (`sys.Ts = 0`). If `sys` is a discrete-time model, `bool = 0 (false)`.

For a static gain, both `isct` and `isdtd` return `true` unless you explicitly set the sample time to a nonzero value. If you do so, `isdtd` returns `true` and `isct` returns `false`.

For arrays of models, `bool` is `true` if the models in the array are continuous.

## **See Also**

isdt | isstable

**Introduced in R2012a**

## isdt

Determine if dynamic system model is in discrete time

### Syntax

```
bool = isdt(sys)
```

### Description

`bool = isdt(sys)` returns a logical value of 1 (`true`) if the dynamic system model `sys` is a discrete-time model. The function returns a logical value of 0 (`false`) otherwise.

### Input Arguments

**sys**

Dynamic system model or array of such models.

### Output Arguments

**bool**

Logical value indicating whether `sys` is a discrete-time model.

`bool = 1 (true)` if `sys` is a discrete-time model (`sys.Ts ≠ 0`). If `sys` is a continuous-time model, `bool = 0 (false)`.

For a static gain, both `isct` and `isdt` return `true` unless you explicitly set the sample time to a nonzero value. If you do so, `isdt` returns `true` and `isct` returns `false`.

For arrays of models, `bool` is `true` if the models in the array are discrete.

## **See Also**

isct | isstable

**Introduced in R2012a**

## isempty

Determine whether dynamic system model is empty

### Syntax

```
isempty(sys)
```

### Description

`isempty(sys)` returns a logical value of 1 (`true`) if the dynamic system model `sys` has no input or no output, and a logical value of 0 (`false`) otherwise. Where `sys` is a `frd` model, `isempty(sys)` returns 1 when the frequency vector is empty. Where `sys` is a model array, `isempty(sys)` returns 1 when the array has empty dimensions or when the LTI models in the array are empty.

### Examples

#### Determine Whether Dynamic Model Is Empty

Create a continuous-time state-space model with 1 input and no outputs. In this example, specify the A and B matrices as 1 and 2, respectively.

```
sys1 = ss(1,2,[],[]);
```

Determine whether `sys1` is empty.

```
isempty(sys1)
```

```
ans = logical  
     1
```

The `isempty` command returns 1 because the system does not have any outputs.

Similarly, `isempty` returns 1 for an empty transfer function.

```
isempty(tf)
ans = logical
     1
```

Now create a state-space model with 1 input and 1 output. In this example, specify the A, B, C, and D matrices as 1, 2, 3, and 4, respectively.

```
sys2 = ss(1,2,3,4);
```

Determine whether sys2 is empty.

```
isempty(sys2)
ans = logical
     0
```

The command returns 0 because the system has inputs and outputs.

## See Also

[issiso](#) | [size](#)

**Introduced before R2006a**

# isLocked

Locked status of online parameter estimation System object

## Syntax

```
L = isLocked(obj)
```

## Description

`L = isLocked(obj)` returns the locked status of online parameter estimation System object, `obj`.

## Examples

### Check Locked Status of Online Estimation System Object

Create a System object™ for online estimation of an ARMAX model with default properties.

```
obj = recursiveARMAX;
```

Check the locked status of the object.

```
L = isLocked(obj)
```

```
L = logical  
    0
```

Estimate model parameters online using `step` and input-output data.

```
[A,B,C,EstimatedOutput] = step(obj,1,1);
```

Check the locked status of the object again.

```
L = isLocked(obj)
```

L = *logical*  
1

step puts the object in a locked state.

## Input Arguments

### **obj** – System object for online parameter estimation

recursiveAR object | recursiveARMA object | recursiveARX object | recursiveARMAX object | recursiveOE object | recursiveBJ object | recursiveLS object

System object for online parameter estimation, created using one of the following commands:

- recursiveAR
- recursiveARMA
- recursiveARX
- recursiveARMAX
- recursiveOE
- recursiveBJ
- recursiveLS

## Output Arguments

### **L** – Locked status of online estimation System object

*logical*

Locked status of online estimation System object, returned as a logical value. L is true if obj is locked.

## See Also

clone | recursiveAR | recursiveARMA | recursiveARMAX | recursiveARX | recursiveBJ | recursiveLS | recursiveOE | release | reset | step

## **Topics**

“What Is Online Estimation?”

**Introduced in R2015b**

## isnlarx

Detect nonlinearity in estimation data

### Syntax

```
isnlarx(Data,Orders)  
isnlarx(Data,Orders,Ky)  
isnlarx( ____,Name,Value)
```

```
NLHyp = isnlarx( ____)  
[NLHyp,NLValue,NLRegs,NoiseSigma,DetectRatio] = isnlarx( ____)
```

### Description

`isnlarx(Data,Orders)` detects nonlinearity in `Data` by testing whether a nonlinear ARX model with the indicated `Orders` produces a better estimate of `Data` than a linear ARX model. The nonlinear model uses a default `treepartition` nonlinearity estimator.

The result of the test is printed to the Command Window and indicates whether a nonlinearity is detected. Use the printed detection ratio to assess the reliability of the nonlinearity detection test:

- Larger values (>2) indicate that a significant nonlinearity was detected.
- Smaller values (<0.5) indicate that any error unexplained by the linear model is mostly noise. That is, no significant nonlinearity was detected.
- Values close to 1 indicate that the nonlinearity detection test is not reliable and that a weak nonlinearity may be present.

`isnlarx(Data,Orders,Ky)` restricts the nonlinearity test to output channel `Ky` for multi-output data.

`isnlarx( ____,Name,Value)` specifies additional nonlinear ARX model options using one or more `Name,Value` pair arguments.

`NLHyp = isnlarx( ____)` returns the result of the nonlinearity test and suppresses the command window output.

[NLHyp,NLValue,NLRegs,NoiseSigma,DetectRatio] = isnlrx( \_\_\_ )  
additionally returns the test quantities behind the evaluation.

## Examples

### Detect Nonlinearity in Estimation Data

Load the signal transmission data set.

```
load(fullfile(matlabroot,'toolbox','ident','iddemos','data','frictiondata'))
```

Construct an iddata object from the estimation data.

```
z = iddata(f1,v,1);
```

Specify the model orders and delays.

```
orders = [1 1 0];
```

Run the test to detect nonlinearity.

```
isnlrx(z,orders);
```

```
Nonlinearity is detected in data set z.  
Detection ratio: 525.05  
Estimated discrepancy of the linear model found: 0.0064966  
Estimated noise standard deviation: 0.00080938
```

The large detection ratio indicates that the test was robust and a significant nonlinearity was detected. Additionally, the estimated discrepancy of the linear model that was found, that is the data explained by the nonlinearity, is significantly greater than the noise error, which can indicate a significant nonlinearity.

### Detect Nonlinearity in Estimation Data Output Channel

Load the CSTR data set.

```
load(fullfile(matlabroot,'toolbox','ident','iddemos','data','cstrdata'))
```

Construct an `iddata` object from the estimation data using a sample time of 0.1 seconds.

```
z = iddata(y1,u1,0.1);
```

Specify the model orders and delays.

```
orders = [3*ones(2,2),ones(2,3),2*ones(2,3)];
```

Run the test to detect nonlinearity on the second output channel.

```
isnlarx(z,orders,2);
```

```
ISNLARX results for dataset z
```

```
Nonlinearity is not detected in channel (2).
```

```
However, the test may be on the edge of detecting the nonlinearity.
```

```
Detection ratio: 0.36446
```

```
Searching for best nonlinear regressors may provide more reliable results.
```

```
-
```

A detection ratio less than 1 indicates that no nonlinearity was detected. However, since this value is near 0.5, there may be a weak nonlinearity that was not detected by the test.

### Search for Best Regressors When Detecting Nonlinearity

Load the signal transmission data set.

```
load(fullfile(matlabroot,'toolbox','ident','iddemos','data','signaltransmissiondata'))
```

Construct an `iddata` object from the estimation data using a sample time of 0.1 seconds.

```
z = iddata(vout,vin,0.1);
```

Specify the model orders and delays.

```
orders = [3 0 2];
```

Display the model regressors for an `idnlarx` model with the given orders.

```
getreg(idnlarx(orders));
```

```
Regressors:
```

```
  y1(t-1)
```

```
y1(t-2)
y1(t-3)
```

Detect nonlinearities in the data, and search for the best nonlinear regressor combination.

```
isnlarx(z,orders,'NonlinearRegressors','search');
```

```
Nonlinearity is detected in data set z.
Detection ratio: 1.4691
Estimated discrepancy of the linear model found: 0.74371
Estimated noise standard deviation: 0.74935
Corresponding NonlinearRegressors parameter: [1          2          3]
```

The regressor search found that using the first two regressors produces the best nonlinear estimation of the given data.

A detection ratio greater than 1 but less than 2 means that a nonlinearity was detected, but the test was not robust. This result may indicate that the detected nonlinearity is not significant. Additionally, the data explained by the nonlinearity is smaller than the noise error, which can be an indication of a weak nonlinearity.

### Return Nonlinearity Detection Result

Load the estimation data.

```
load(fullfile(matlabroot,'toolbox','ident','iddemos','data','cstrdata'))
```

Construct an `iddata` object using the estimation data.

```
z = iddata(y1,u1,0.1);
```

Specify the model orders and delays.

```
orders = [3*ones(2,2),ones(2,3),2*ones(2,3)];
```

Detect nonlinearities in the data, and determine the test quantities behind the evaluation.

```
NLHyp = isnlarx(z,orders);
```

## Return Nonlinearity Detection Test Quantities

Load the estimation data.

```
load(fullfile(matlabroot, 'toolbox', 'ident', 'iddemos', 'data', 'narendralidata'))
```

Construct an `iddata` object using the estimation data.

```
z = iddata(u,y1,1);
```

Specify the model orders and delays.

```
orders = [1 1 2];
```

Detect nonlinearities in the data, and determine the test quantities behind the evaluation.

```
[NLHyp,NLValue,NLRegs,NoiseSigma,DetectRatio] = isnlarx(z,orders);
```

## Input Arguments

### Data — Time-domain estimation data

`iddata` object

Time-domain estimation data, specified as an `iddata` object. Data can have one or more output channels and zero or more input channels. Data must be uniformly sampled and cannot contain missing (NaN) samples.

### Orders — Model orders and delays

1-by-3 vector of positive integers | 1-by-3 vector of matrices

Model orders and delays for defining the regressor configuration, specified as a 1-by-3 vector, `[na nb nk]`.

For a model with  $n_y$  output channels and  $n_u$  input channels:

- `na` is an  $n_y$ -by- $n_y$  matrix, where `na(i,j)` specifies the number of regressors from the  $j$ th output used to predict the  $i$ th output.
- `nb` is an  $n_y$ -by- $n_u$  matrix, where `nb(i,j)` specifies the number of regressors from the  $j$ th input used to predict the  $i$ th output.
- `nk` is an  $n_y$ -by- $n_u$  matrix, where `nk(i,j)` specifies the lag in the  $j$ th input used to predict the  $i$ th output.

```
na = [1 2; 2 3]
nb = [1 2 3; 2 3 1];
nk = [2 0 3; 1 0 5];
```

The estimation data for this system has three inputs ( $u_1$ ,  $u_2$ ,  $u_3$ ) and two outputs ( $y_1$ ,  $y_2$ ). Consider the regressors used to predict output,  $y_2(t)$ :

- Since  $na(2, :)$  is  $[2 \ 3]$ , the contributing regressors from the outputs are:
  - $y_1(t-1)$  and  $y_1(t-2)$
  - $y_2(t-1)$ ,  $y_2(t-2)$ , and  $y_2(t-3)$
- Since  $nb(2, :)$  is  $[2 \ 3 \ 1]$  and  $nk(2, :)$  is  $[1 \ 0 \ 5]$ , the contributing regressors from the inputs are:
  - $u_1(t-1)$  and  $u_1(t-2)$
  - $u_2(t)$ ,  $u_2(t-1)$ , and  $u_2(t-2)$
  - $u_3(t-5)$

---

**Note** The minimum lag for regressors based on output variables is always 1, while the minimum lag for regressors based on input variables is dictated by  $nk$ . Use `getreg` to view the complete set of regressors used by the nonlinear ARX model.

---

### **Ky** — Output channel number in estimation data

positive integer in the range  $[0, n_y]$

Output channel number in estimation data, specified as a positive integer in the range  $[1, n_y]$ , where  $n_y$  is the number of output channels.

## **Name-Value Pair Arguments**

Specify optional comma-separated pairs of `Name`, `Value` arguments. `Name` is the argument name and `Value` is the corresponding value. `Name` must appear inside quotes. You can specify several name and value pair arguments in any order as `Name1, Value1, ..., NameN, ValueN`.

Example: `'NonlinearRegressors', 'output'` specifies that only the regressors containing output variables are used as inputs to the nonlinear block of the model.

### **TimeVariable** — Independent variable name

't' (default) | character vector

Independent variable name, specified as the comma-separated pair consisting of 'TimeVariable' and a character vector. For example, 't'.

### **CustomRegressors — Regressors constructed from combinations of inputs and outputs**

{ } (default) | cell array of character vectors | array of customreg objects

Regressors constructed from combinations of inputs and outputs, specified as the comma-separated pair consisting of 'CustomRegressors' and one of the following for single-output systems:

- Cell array of character vectors. For example:
  - {'y1(t-3)^3', 'y2(t-1)\*u1(t-3)', 'sin(u3(t-2))' }

Each character vector must represent a valid formula for a regressor contributing towards the prediction of the model output. The formula must be written using the input and output names and the time variable name as variables.

- Array of custom regressor objects, created using `customreg` or `polyreg`.

For a model with  $n_y$  outputs, specify an  $n_y$ -by-1 cell array of `customreg` object arrays or character arrays.

These regressors are in addition to the standard regressors based on `Orders`.

Example: 'CustomRegressors', {'y1(t-3)^3', 'y2(t-1)\*u1(t-3)' }

Example: 'CustomRegressors', {'sin(u3(t-2))' }

### **NonlinearRegressors — Subset of regressors that enter as inputs to the nonlinear block of the model**

'all' (default) | 'output' | 'input' | 'standard' | 'custom' | 'search' | vector of positive integers | [] | cell array

Subset of regressors that enter as inputs to the nonlinear block of the model, specified as the comma-separated pair consisting of 'NonlinearRegressors' and one of the following values:

- 'all' — All regressors
- 'output' — Regressors containing output variables
- 'input' — Regressors containing input variables

- 'standard' — Standard regressors
- 'custom' — Custom regressors
- 'search' — The estimation algorithm performs a search for the best regressor subset. This is useful when you want to reduce a large number of regressors entering the nonlinear function block of the nonlinearity estimator. This option must be applied to all output models simultaneously.
- [] — No regressors. This creates a linear-in-regressor model.
- Vector of regressor indices. To determine the number and order of regressors, use `getreg`.

For a model with multiple outputs, specify a cell array of  $n_y$  elements, where  $n_y$  is the number of output channels. For each output, specify one of the preceding options. Alternatively, to apply the same regressor subset to all model outputs, specify [] or any of the character vector options alone, for example 'standard'.

Example: 'NonlinearRegressors', 'search' performs a best regressor search for the only output of a single output model, or all of the outputs of a multiple output model.

Example: 'NonlinearReg', 'input' applies only input regressors to the inputs of the nonlinear function.

Example: 'NonlinearRegressors', {'input', 'output'} applies input regressors to the first output, and output regressors to the second output of a model with two outputs.

## Output Arguments

### NLHyp — Result of the nonlinearity test

0 | 1 | logical vector

Result of the nonlinearity test, returned as a logical vector with length equal to the number of output channels. The elements of NLHyp are 1 if nonlinearities were detected for the corresponding output. A value of 0 indicates that nonlinearities were not detected.

### NLValue — Estimated standard deviation of the data explained by the nonlinearity

vector of nonnegative scalars

Estimated standard deviation of the data explained by the nonlinearity, returned as a vector of nonnegative scalars with length equal to the number of output channels. The

elements of `NLValue` are 0 if nonlinearities are not detected for the corresponding output.

### **NLRegs — Regressors that should enter nonlinearly in the model**

vector of indices | [] | cell array

Regressors that should enter nonlinearly in the model, returned as a vector of indices for single output models. For multi-output models, `NLRegs` is returned as a cell array, with elements corresponding to each output channel. `NLRegs` is empty, [], if nonlinearities are not detected.

See the 'NonlinearRegressors' Name, Value argument for more information.

### **NoiseSigma — Estimated standard deviation of the unexplained error**

vector of nonnegative scalars

Estimated standard deviation of the unexplained error, returned as a vector of nonnegative scalars with length equal to the number of output channels. The elements of `NoiseSigma` are 0 if nonlinearities are not detected for the corresponding output.

### **DetectRatio — Ratio of the test statistic and the detection threshold**

vector

Ratio of the test statistic and the detection threshold, returned as a vector with length equal to the number of output channels. Use the elements of `DetectRatio` to assess the reliability of the nonlinearity detection test for the corresponding output:

- Larger values (>2) indicate that a significant nonlinearity was detected.
- Smaller values (<0.5) indicate that any error unexplained by the linear model is mostly noise. That is, no significant nonlinearity was detected.
- Values close to 1 indicate that the nonlinearity detection test is not reliable and that a weak nonlinearity may be present.

## **Algorithms**

`isnlarx` estimates a nonlinear ARX model using the given data and a `treepartition` nonlinearity estimator.

The estimation data can be described as  $Y(t) = L(t) + F_n(t) + E(t)$ , where:

- $L(t)$  is the portion of the data explained by the linear function of the nonlinear ARX model.
- $F_n(t)$  is the portion of the data explained by the nonlinear function of the nonlinear ARX model. The output argument `NLValue` is an estimate of the standard deviation of  $F_n(t)$ . If the nonlinear function explains a significant portion of the data beyond the data explained by the linear function, a nonlinearity is detected.
- $E(t)$  is the remaining error that is unexplained by the nonlinear ARX model and is typically white noise. The output argument `NoiseSigma` is an estimate of the standard deviation of  $E(t)$ .

## See Also

`getreg` | `idnlarx` | `nlarx` | `treepartition`

## Topics

“Structure of Nonlinear ARX Models”

**Introduced in R2007a**

## isproper

Determine if dynamic system model is proper

### Syntax

```
B = isproper(sys)
B = isproper(sys, 'elem')
[B,sysr] = isproper(sys)
```

### Description

`B = isproper(sys)` returns a logical value of 1 (`true`) if the dynamic system model `sys` is proper and a logical value of 0 (`false`) otherwise.

A proper model has relative degree  $\leq 0$  and is causal. SISO transfer functions and zero-pole-gain models are proper if the degree of their numerator is less than or equal to the degree of their denominator (in other words, if they have at least as many poles as zeroes). MIMO transfer functions are proper if all their SISO entries are proper. Regular state-space models (state-space models having no `E` matrix) are always proper. A descriptor state-space model that has an invertible `E` matrix is always proper. A descriptor state-space model having a singular (non-invertible) `E` matrix is proper if the model has at least as many poles as zeroes.

If `sys` is a model array, then `B` is 1 if all models in the array are proper.

`B = isproper(sys, 'elem')` checks each model in a model array `sys` and returns a logical array of the same size as `sys`. The logical array indicates which models in `sys` are proper.

`[B,sysr] = isproper(sys)` also returns an equivalent model `sysr` with fewer states (reduced order) and a non-singular `E` matrix, if `sys` is a proper descriptor state-space model with a non-invertible `E` matrix. If `sys` is not proper, `sysr = sys`.

### Examples

## Examine Whether Models are Proper

Create a SISO continuous-time transfer function,  $H_1 = s$

```
H1 = tf([1 0],1);
```

Check whether H1 is proper.

```
B1 = isproper(H1)
```

```
B1 = logical
      0
```

SISO transfer functions are proper if the degree of their numerator is less than or equal to the degree of their denominator. That is, if the transfer function has at least as many poles as zeroes. Since H1 has one zero and no poles, the `isproper` command returns 0.

Now create a transfer function with one pole and one zero,  $H_2 = s/(s + 1)$

```
H2 = tf([1 0],[1 1]);
```

Check whether H2 is proper.

```
B2 = isproper(H2)
```

```
B2 = logical
      1
```

Since H2 has equal number of poles and zeros, `isproper` returns 1.

## Compute Equivalent Lower-Order Model

Combining state-space models sometimes yields results that include more states than necessary. Use `isproper` to compute an equivalent lower-order model.

```
H1 = ss(tf([1 1],[1 2 5]));
H2 = ss(tf([1 7],[1]));
H = H1*H2;
size(H)
```

State-space model with 1 outputs, 1 inputs, and 4 states.

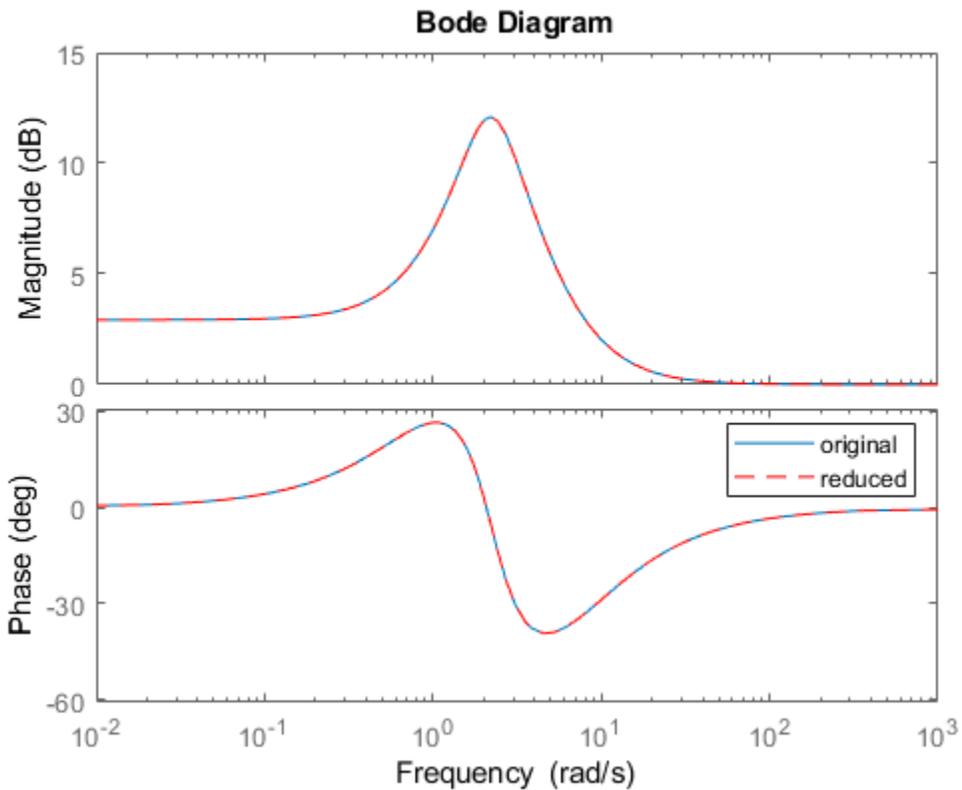
H is proper and reducible. `isproper` returns the reduced model.

```
[isprop,Hr] = isproper(H);  
size(Hr)
```

State-space model with 1 outputs, 1 inputs, and 2 states.

H and Hr are equivalent, as a Bode plot demonstrates.

```
bodeplot(H,Hr,'r--')  
legend('original','reduced')
```



## **See Also**

dss | ss

**Introduced before R2006a**

## **isreal**

Determine whether model parameters or data values are real

### **Syntax**

```
isreal(Data)  
isreal(Model)
```

### **Description**

`isreal(Data)` returns 1 if all signals of the data set are real. `Data` is an `iddata` object.

`isreal(Model)` returns 1 if all parameters of the model are real. `Model` is any linear identified model.

### **See Also**

`realdata`

**Introduced before R2006a**

## issiso

Determine if dynamic system model is single-input/single-output (SISO)

### Syntax

```
issiso(sys)
```

### Description

`issiso(sys)` returns a logical value of 1 (`true`) if the dynamic system model `sys` is SISO and a logical value of 0 (`false`) otherwise.

### See Also

`isempty` | `size`

**Introduced in R2012a**

## isstable

Determine whether system is stable

### Syntax

```
B = isstable(sys)
B = isstable(sys, 'elem')
```

### Description

`B = isstable(sys)` returns a logical value of 1 (`true`) if the dynamic system model `sys` has stable dynamics, and a logical value of 0 (`false`) otherwise. If `sys` is a model array, then `B = 1` only if all models in `sys` are stable.

`B = isstable(sys, 'elem')` returns a logical array of the same dimensions as the model array `sys`. The logical array indicates which models in `sys` are stable.

`isstable` is only supported for analytical models with a finite number of poles.

### Examples

#### Determine Stability of Models in Model Array

Create an array of SISO transfer function models with poles varying from -2 to 2. To do so, first initialize an array of dimension `[1, length(a)]` with zero-valued SISO transfer functions.

```
a = [-2:2];
sys = tf(zeros(1,1,1,length(a)));
```

Populate this array with transfer functions of the form  $1/(s-a)$ .

```
for j = 1:length(a)
    sys(1,1,1,j) = tf(1,[1 -a(j)]);
```

```
end  
sys.SamplingGrid = struct('a',a);
```

Examine the stability of the model array.

```
B_all = isstable(sys)  
  
B_all = logical  
      0
```

By default, `isstable` returns a single Boolean value that is 1 (true) only if all models in the array are stable. `sys` contains some models with nonnegative poles, which are not stable. Therefore, `isstable` returns 0 (false) for the entire array.

Examine stability of each model in the array, element by element.

```
B_elem = isstable(sys,'elem')  
  
B_elem = 1x5 logical array  
  
      1      1      0      0      0
```

The 'elem' flag causes `isstable` to return an array of Boolean values, which indicate the stability of the corresponding entry in the model array. For example, `B_elem(2) = 1`, which indicates that `sys(1,1,1,2)` is stable. This result is expected, because `sys(1,1,1,2)` has `a = -1`.

## See Also

`pole`

**Introduced in R2012a**

## ivar

AR model estimation using instrumental variable method

### Syntax

```
sys = ivar(data,na)
sys = ivar(data,na,nc)
sys = ivar(data,na,nc,max_size)
```

### Description

`sys = ivar(data,na)` estimates an AR polynomial model, `sys`, using the instrumental variable method and the time series data `data`. `na` specifies the order of the  $A$  polynomial.

An AR model is represented by the equation:

$$A(q)y(t) = e(t)$$

In the above model,  $e(t)$  is an arbitrary process, assumed to be a moving average process of order `nc`, possibly time varying. `nc` is assumed to be equal to `na`. Instruments are chosen as appropriately filtered outputs, delayed `nc` steps.

`sys = ivar(data,na,nc)` specifies the value of the moving average process order, `nc`, separately.

`sys = ivar(data,na,nc,max_size)` specifies the maximum size of matrices formed during estimation.

### Input Arguments

#### **data**

Estimation time series data.

data must be an `iddata` object with scalar output data only.

**na**

Order of the  $A$  polynomial

**nc**

Order of the moving average process representing  $e(t)$ .

**max\_size**

Maximum matrix size.

`max_size` specifies the maximum size of any matrix formed by the algorithm for estimation.

Specify `max_size` as a reasonably large positive integer.

**Default:** 250000

## Output Arguments

**sys**

Identified polynomial model.

`sys` is an AR `idpoly` model which encapsulates the identified polynomial model.

## Examples

Compare spectra for sinusoids in noise, estimated by the IV method and by the forward-backward least squares method.

```
y = iddata(sin([1:500]'*1.2) + sin([1:500]'*1.5) + ...  
           0.2*randn(500,1),[]);  
miv = ivar(y,4);  
mls = ar(y,4);  
spectrum(miv,mls)
```

## References

- [1] Stoica, P., et al. *Optimal Instrumental Variable Estimates of the AR-parameters of an ARMA Process*, IEEE Trans. Autom. Control, Volume AC-30, 1985, pp. 1066-1074.

## See Also

ar | arx | etfe | idpoly | polyest | spa | spectrum | step

**Introduced before R2006a**

## ivstruc

Compute loss functions for sets of ARX model structures using instrumental variable method

### Syntax

```
v = ivstruc(ze,zv,NN)
v = ivstruc(ze,zv,NN,p,maxsize)
```

### Description

`v = ivstruc(ze,zv,NN)` computes the loss functions for sets of single-output ARX model structures. `NN` is a matrix that defines a number of different structures of the ARX type. Each row of `NN` is of the form

```
nn = [na nb nk]
```

with the same interpretation as described for `arx`. See `struc` for easy generation of typical `NN` matrices.

`ze` and `zv` are `iddata` objects containing input-output data. Only time-domain data is supported. Models for each model structure defined in `NN` are estimated using the instrumental variable (IV) method on data set `ze`. The estimated models are simulated using the inputs from data set `zv`. The normalized quadratic fit between the simulated output and the measured output in `zv` is formed and returned in `v`. The rows below the first row in `v` are the transpose of `NN`, and the last row contains the logarithms of the condition numbers of the IV matrix

$$\sum \zeta(t)\varphi^T(t)$$

A large condition number indicates that the structure is of unnecessarily high order (see Ljung, L. *System Identification: Theory for the User*, Upper Saddle River, NJ, Prentice-Hall PTR, 1999, p. 498).

The information in `v` is best analyzed using `selstruc`.

The routine is for single-output systems only.

`v = ivstruc(ze,zv,NN,p,maxsize)` specifies the computation of condition numbers and the size of largest matrix formed during computations. If `p` is equal to zero, the computation of condition numbers is suppressed. `maxsize` affects the speed/memory trade-off.

---

**Note** The IV method used does not guarantee that the models obtained are stable. The output-error fit calculated in `v` can then be misleading.

---

## Examples

### Generate Model-Order Combinations and Estimate ARX Model Using IV Method

Create estimation and validation data sets

```
load iddata1;  
ze = z1(1:150);  
zv = z1(151:300);
```

Generate model-order combinations for estimation, specifying ranges for model orders and delays.

```
NN = struc(1:3,1:2,2:4);
```

Estimate ARX models using the instrumental variable method, and compute the loss function for each model order combination.

```
V = ivstruc(ze,zv,NN);
```

Select the model order with the best fit to the validation data.

```
order = selstruc(V,0);
```

Estimate an ARX model of selected order.

```
M = iv4(ze,order);
```

## Suppress Condition Number Computation When Determining ARX Loss Functions

Create estimation and validation data sets.

```
load iddata1;  
ze = z1(1:150);  
zv = z1(151:300);
```

Generate model-order combinations for estimation, specifying ranges for model orders and a delay of 2 for all model configurations.

```
NN = struc(2:3,1:2,2);
```

Compute the loss function for each model order combination. Suppress the computation of condition numbers.

```
V = ivstruc(ze,zv,NN,0);
```

## Algorithms

A maximum-order ARX model is computed using the least squares method. Instruments are generated by filtering the input(s) through this model. The models are subsequently obtained by operating on submatrices in the corresponding large IV matrix.

## References

- [1] Ljung, L. *System Identification: Theory for the User*, Upper Saddle River, NJ, Prentice-Hall PTR, 1999.

## See Also

arxstruc | iv4 | selstruc | struc

Introduced before R2006a

## **ivx**

ARX model estimation using instrumental variable method with arbitrary instruments

### **Syntax**

```
sys = ivx(data,[na nb nk],x)
sys = ivx(data,[na nb nk],x,max_size)
```

### **Description**

`sys = ivx(data,[na nb nk],x)` estimates an ARX polynomial model, `sys`, using the instrumental variable method with arbitrary instruments. The model is estimated for the time series data `data`. `[na nb nk]` specifies the ARX structure orders of the  $A$  and  $B$  polynomials and the input to output delay, expressed in the number of samples.

An ARX model is represented as:

$$A(q)y(t) = B(q)u(t - nk) + v(t)$$

`sys = ivx(data,[na nb nk],x,max_size)` specifies the maximum size of matrices formed during estimation.

### **Input Arguments**

#### **data**

Estimation data. The data can be:

- Time- or frequency-domain input-output data
- Time-series data
- Frequency-response data

`data` must be an `iddata`, `idfrd`, or `frd` object.

When using frequency-domain data, the number of outputs must be 1.

**[na nb nk]**

ARX model orders.

For more details on the ARX model structure, see `arx`.

**x**

Instrument variable matrix.

`x` is a matrix containing the arbitrary instruments for use in the instrumental variable method.

`x` must be of the same size as the output data, `data.y`. For multi-experiment data, specify `x` as a cell array with one entry for each experiment.

The instruments used are analogous to the regression vector, with `y` replaced by `x`.

**max\_size**

Maximum matrix size.

`max_size` specifies the maximum size of any matrix formed by the algorithm for estimation.

Specify `max_size` as a reasonably large positive integer.

**Default:** 250000

## Output Arguments

**sys**

ARX model that fits the estimation data, returned as a discrete-time `idpoly` object. This model is created using the specified model orders, delays, and estimation options. `ivx` does not return any estimated covariance information for `sys`.

Information about the estimation results and options used is stored in the `Report` property of the model. `Report` has the following fields:

<b>Report Field</b>	<b>Description</b>																		
Status	Summary of the model status, which indicates whether the model was created by construction or obtained by estimation.																		
Method	Estimation command used.																		
InitialCondition	<p>Handling of initial conditions during model estimation, returned as one of the following values:</p> <ul style="list-style-type: none"> <li>'zero' — The initial conditions were set to zero.</li> <li>'estimate' — The initial conditions were treated as independent estimation parameters.</li> <li>'backcast' — The initial conditions were estimated using the best least squares fit.</li> </ul> <p>This field is especially useful to view how the initial conditions were handled when the InitialCondition option in the estimation option set is 'auto'.</p>																		
Fit	<p>Quantitative assessment of the estimation, returned as a structure. See “Loss Function and Model Quality Metrics” for more information on these quality metrics. The structure has the following fields:</p> <table border="1"> <thead> <tr> <th><b>Field</b></th> <th><b>Description</b></th> </tr> </thead> <tbody> <tr> <td>FitPercent</td> <td>Normalized root mean squared error (NRMSE) measure of how well the response of the model fits the estimation data, expressed as a percentage.</td> </tr> <tr> <td>LossFcn</td> <td>Value of the loss function when the estimation completes.</td> </tr> <tr> <td>MSE</td> <td>Mean squared error (MSE) measure of how well the response of the model fits the estimation data.</td> </tr> <tr> <td>FPE</td> <td>Final prediction error for the model.</td> </tr> <tr> <td>AIC</td> <td>Raw Akaike Information Criteria (AIC) measure of model quality.</td> </tr> <tr> <td>AICc</td> <td>Small sample-size corrected AIC.</td> </tr> <tr> <td>nAIC</td> <td>Normalized AIC.</td> </tr> <tr> <td>BIC</td> <td>Bayesian Information Criteria (BIC).</td> </tr> </tbody> </table>	<b>Field</b>	<b>Description</b>	FitPercent	Normalized root mean squared error (NRMSE) measure of how well the response of the model fits the estimation data, expressed as a percentage.	LossFcn	Value of the loss function when the estimation completes.	MSE	Mean squared error (MSE) measure of how well the response of the model fits the estimation data.	FPE	Final prediction error for the model.	AIC	Raw Akaike Information Criteria (AIC) measure of model quality.	AICc	Small sample-size corrected AIC.	nAIC	Normalized AIC.	BIC	Bayesian Information Criteria (BIC).
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BIC	Bayesian Information Criteria (BIC).																		
Parameters	Estimated values of model parameters.																		

Report Field	Description																
OptionsUsed	Option set used for estimation. If no custom options were configured, this is a set of default options. See <code>arxOptions</code> for more information.																
RandState	State of the random number stream at the start of estimation. Empty, <code>[]</code> , if randomization was not used during estimation. For more information, see <code>rng</code> in the MATLAB documentation.																
DataUsed	Attributes of the data used for estimation, returned as a structure with the following fields: <table border="1" data-bbox="387 604 1332 1289"> <thead> <tr> <th>Field</th> <th>Description</th> </tr> </thead> <tbody> <tr> <td>Name</td> <td>Name of the data set.</td> </tr> <tr> <td>Type</td> <td>Data type.</td> </tr> <tr> <td>Length</td> <td>Number of data samples.</td> </tr> <tr> <td>Ts</td> <td>Sample time.</td> </tr> <tr> <td>InterSample</td> <td>Input intersample behavior, returned as one of the following values: <ul style="list-style-type: none"> <li>'zoh' — Zero-order hold maintains a piecewise-constant input signal between samples.</li> <li>'foh' — First-order hold maintains a piecewise-linear input signal between samples.</li> <li>'bl' — Band-limited behavior specifies that the continuous-time input signal has zero power above the Nyquist frequency.</li> </ul> </td> </tr> <tr> <td>InputOffset</td> <td>Offset removed from time-domain input data during estimation. For nonlinear models, it is <code>[]</code>.</td> </tr> <tr> <td>OutputOffset</td> <td>Offset removed from time-domain output data during estimation. For nonlinear models, it is <code>[]</code>.</td> </tr> </tbody> </table>	Field	Description	Name	Name of the data set.	Type	Data type.	Length	Number of data samples.	Ts	Sample time.	InterSample	Input intersample behavior, returned as one of the following values: <ul style="list-style-type: none"> <li>'zoh' — Zero-order hold maintains a piecewise-constant input signal between samples.</li> <li>'foh' — First-order hold maintains a piecewise-linear input signal between samples.</li> <li>'bl' — Band-limited behavior specifies that the continuous-time input signal has zero power above the Nyquist frequency.</li> </ul>	InputOffset	Offset removed from time-domain input data during estimation. For nonlinear models, it is <code>[]</code> .	OutputOffset	Offset removed from time-domain output data during estimation. For nonlinear models, it is <code>[]</code> .
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OutputOffset	Offset removed from time-domain output data during estimation. For nonlinear models, it is <code>[]</code> .																

For more information on using `Report`, see “Estimation Report”.

## Tips

- Use `iv4` first for IV estimation to identify ARX polynomial models where the instruments `x` are chosen automatically. Use `ivx` for nonstandard situations. For

example, when there is feedback present in the data, or, when other instruments need to be tried. You can also use `iv` to automatically generate instruments from certain custom defined filters.

### References

- [1] Ljung, L. *System Identification: Theory for the User*, page 222, Upper Saddle River, NJ, Prentice-Hall PTR, 1999.

### See Also

`arx` | `arxstruc` | `idpoly` | `iv4` | `ivar` | `polyest`

**Introduced before R2006a**

## iv4

ARX model estimation using four-stage instrumental variable method

### Syntax

```
sys = iv4(data,[na nb nk])
sys = iv4(data,'na',na,'nb',nb,'nk',nk)
sys = iv4( ____,Name,Value)
sys = iv4( ____,opt)
```

### Description

`sys = iv4(data,[na nb nk])` estimates an ARX polynomial model, `sys`, using the four-stage instrumental variable method, for the data object `data`. `[na nb nk]` specifies the ARX structure orders of the  $A$  and  $B$  polynomials and the input to output delay. The estimation algorithm is insensitive to the color of the noise term.

`sys` is an ARX model:

$$A(q)y(t) = B(q)u(t - nk) + v(t)$$

`sys = iv4(data,'na',na,'nb',nb,'nk',nk)` alternatively specify the ARX model orders separately.

`sys = iv4( ____,Name,Value)` estimates an ARX polynomial with additional options specified by one or more `Name,Value` pair arguments.

`sys = iv4( ____,opt)` uses the option set, `opt`, to configure the estimation behavior.

### Input Arguments

#### **data**

Estimation data. The data can be:

- Time- or frequency-domain input-output data
- Time-series data
- Frequency-response data

`data` must be an `iddata`, `idfrd`, or `frd` object.

`data` must be discrete-time ( $T_s > 0$ ) for frequency domain.

### **[na nb nk]**

ARX polynomial orders.

For multi-output model, `[na nb nk]` contains one row for every output. In particular, specify `na` as an  $N_y$ -by- $N_y$  matrix, where each entry is the polynomial order relating the corresponding output pair. Here,  $N_y$  is the number of outputs. Specify `nb` and `nk` as  $N_y$ -by- $N_u$  matrices, where  $N_u$  is the number of inputs. For more details on the ARX model structure, see `arx`.

### **opt**

Estimation options.

`opt` is an options set that configures the estimation options. These options include:

- estimation focus
- handling of initial conditions
- handling of data offsets

Use `iv4options` to create the options set.

## **Name-Value Pair Arguments**

Specify optional comma-separated pairs of `Name`, `Value` arguments. `Name` is the argument name and `Value` is the corresponding value. `Name` must appear inside quotes. You can specify several name and value pair arguments in any order as `Name1, Value1, . . . , NameN, ValueN`.

### **InputDelay**

Input delay for each input channel, specified as a scalar value or numeric vector. For continuous-time systems, specify input delays in the time unit stored in the `TimeUnit`

property. For discrete-time systems, specify input delays in integer multiples of the sample time  $T_s$ . For example, `InputDelay = 3` means a delay of three sample times.

For a system with  $N_u$  inputs, set `InputDelay` to an  $N_u$ -by-1 vector. Each entry of this vector is a numerical value that represents the input delay for the corresponding input channel.

You can also set `InputDelay` to a scalar value to apply the same delay to all channels.

**Default:** 0

### **IODelay**

Transport delays. `IODelay` is a numeric array specifying a separate transport delay for each input/output pair.

For continuous-time systems, specify transport delays in the time unit stored in the `TimeUnit` property. For discrete-time systems, specify transport delays in integer multiples of the sample time,  $T_s$ .

For a MIMO system with  $N_y$  outputs and  $N_u$  inputs, set `IODelay` to a  $N_y$ -by- $N_u$  array. Each entry of this array is a numerical value that represents the transport delay for the corresponding input/output pair. You can also set `IODelay` to a scalar value to apply the same delay to all input/output pairs.

**Default:** 0 for all input/output pairs

### **IntegrateNoise**

Specify integrators in the noise channels.

Adding an integrator creates an ARIX model represented by:

$$A(q)y(t) = B(q)u(t - nk) + \frac{1}{1 - q^{-1}} e(t)$$

where,  $\frac{1}{1 - q^{-1}}$  is the integrator in the noise channel,  $e(t)$ .

`IntegrateNoise` is a logical vector of length  $N_y$ , where  $N_y$  is the number of outputs.

**Default:** `false(Ny, 1)`, where `Ny` is the number of outputs

## Output Arguments

### `sys`

ARX model that fits the estimation data, returned as a discrete-time `idpoly` object. This model is created using the specified model orders, delays, and estimation options.

Information about the estimation results and options used is stored in the `Report` property of the model. `Report` has the following fields:

Report Field	Description
Status	Summary of the model status, which indicates whether the model was created by construction or obtained by estimation.
Method	Estimation command used.
InitialCondition	Handling of initial conditions during model estimation, returned as one of the following values: <ul style="list-style-type: none"><li>'zero' — The initial conditions were set to zero.</li><li>'estimate' — The initial conditions were treated as independent estimation parameters.</li></ul> <p>This field is especially useful to view how the initial conditions were handled when the <code>InitialCondition</code> option in the estimation option set is 'auto'.</p>

Report Field	Description																		
Fit	<p>Quantitative assessment of the estimation, returned as a structure. See “Loss Function and Model Quality Metrics” for more information on these quality metrics. The structure has the following fields:</p> <table border="1" data-bbox="387 456 1331 944"> <thead> <tr> <th data-bbox="387 456 516 499">Field</th> <th data-bbox="516 456 1331 499">Description</th> </tr> </thead> <tbody> <tr> <td data-bbox="387 499 516 604">FitPercent</td> <td data-bbox="516 499 1331 604">Normalized root mean squared error (NRMSE) measure of how well the response of the model fits the estimation data, expressed as a percentage.</td> </tr> <tr> <td data-bbox="387 604 516 647">LossFcn</td> <td data-bbox="516 604 1331 647">Value of the loss function when the estimation completes.</td> </tr> <tr> <td data-bbox="387 647 516 725">MSE</td> <td data-bbox="516 647 1331 725">Mean squared error (MSE) measure of how well the response of the model fits the estimation data.</td> </tr> <tr> <td data-bbox="387 725 516 769">FPE</td> <td data-bbox="516 725 1331 769">Final prediction error for the model.</td> </tr> <tr> <td data-bbox="387 769 516 812">AIC</td> <td data-bbox="516 769 1331 812">Raw Akaike Information Criteria (AIC) measure of model quality.</td> </tr> <tr> <td data-bbox="387 812 516 855">AICc</td> <td data-bbox="516 812 1331 855">Small sample-size corrected AIC.</td> </tr> <tr> <td data-bbox="387 855 516 899">nAIC</td> <td data-bbox="516 855 1331 899">Normalized AIC.</td> </tr> <tr> <td data-bbox="387 899 516 944">BIC</td> <td data-bbox="516 899 1331 944">Bayesian Information Criteria (BIC).</td> </tr> </tbody> </table>	Field	Description	FitPercent	Normalized root mean squared error (NRMSE) measure of how well the response of the model fits the estimation data, expressed as a percentage.	LossFcn	Value of the loss function when the estimation completes.	MSE	Mean squared error (MSE) measure of how well the response of the model fits the estimation data.	FPE	Final prediction error for the model.	AIC	Raw Akaike Information Criteria (AIC) measure of model quality.	AICc	Small sample-size corrected AIC.	nAIC	Normalized AIC.	BIC	Bayesian Information Criteria (BIC).
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nAIC	Normalized AIC.																		
BIC	Bayesian Information Criteria (BIC).																		
Parameters	Estimated values of model parameters.																		
OptionsUsed	Option set used for estimation. If no custom options were configured, this is a set of default options. See <code>iv4options</code> for more information.																		
RandState	State of the random number stream at the start of estimation. Empty, [], if randomization was not used during estimation. For more information, see <code>rng</code> in the MATLAB documentation.																		

Report Field	Description																
DataUsed	Attributes of the data used for estimation, returned as a structure with the following fields:																
	<table border="1"> <thead> <tr> <th>Field</th> <th>Description</th> </tr> </thead> <tbody> <tr> <td>Name</td> <td>Name of the data set.</td> </tr> <tr> <td>Type</td> <td>Data type.</td> </tr> <tr> <td>Length</td> <td>Number of data samples.</td> </tr> <tr> <td>Ts</td> <td>Sample time.</td> </tr> <tr> <td>InterSample</td> <td>Input intersample behavior, returned as one of the following values: <ul style="list-style-type: none"> <li>'zoh' — Zero-order hold maintains a piecewise-constant input signal between samples.</li> <li>'foh' — First-order hold maintains a piecewise-linear input signal between samples.</li> <li>'bl' — Band-limited behavior specifies that the continuous-time input signal has zero power above the Nyquist frequency.</li> </ul> </td> </tr> <tr> <td>InputOffset</td> <td>Offset removed from time-domain input data during estimation. For nonlinear models, it is [ ].</td> </tr> <tr> <td>OutputOffset</td> <td>Offset removed from time-domain output data during estimation. For nonlinear models, it is [ ].</td> </tr> </tbody> </table>	Field	Description	Name	Name of the data set.	Type	Data type.	Length	Number of data samples.	Ts	Sample time.	InterSample	Input intersample behavior, returned as one of the following values: <ul style="list-style-type: none"> <li>'zoh' — Zero-order hold maintains a piecewise-constant input signal between samples.</li> <li>'foh' — First-order hold maintains a piecewise-linear input signal between samples.</li> <li>'bl' — Band-limited behavior specifies that the continuous-time input signal has zero power above the Nyquist frequency.</li> </ul>	InputOffset	Offset removed from time-domain input data during estimation. For nonlinear models, it is [ ].	OutputOffset	Offset removed from time-domain output data during estimation. For nonlinear models, it is [ ].
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For more information on using Report, see “Estimation Report”.

## Examples

### Estimate ARX Model Using Four-Stage Instrumental Variable Method

Load estimation data.

```
load iddata7;
```

This data has two inputs, u1 and u2, and one output, y1.

Specify the ARX model orders, using the same orders for both inputs.

```
na = 2;  
nb = [2 2];
```

Specify a delay of 2 samples for input u2 and no delay for input u1.

```
nk = [0 2];
```

Estimate an ARX model using the four-stage instrumental variable method.

```
m = iv4(z7,[na nb nk]);
```

## Algorithms

Estimation is performed in 4 stages. The first stage uses the `arx` function. The resulting model generates the instruments for a second-stage IV estimate. The residuals obtained from this model are modeled as a high-order AR model. At the fourth stage, the input-output data is filtered through this AR model and then subjected to the IV function with the same instrument filters as in the second stage.

For the multiple-output case, optimal instruments are obtained only if the noise sources at the different outputs have the same color. The estimates obtained with the routine are reasonably accurate, however, even in other cases.

## References

- [1] Ljung, L. *System Identification: Theory for the User*, equations (15.21) through (15.26), Upper Saddle River, NJ, Prentice-Hall PTR, 1999.

## See Also

`armax` | `arx` | `bj` | `idpoly` | `iv4options` | `ivx` | `n4sid` | `oe` | `polyest`

**Introduced before R2006a**

## iv4Options

Option set for iv4

### Syntax

```
opt = iv4options  
opt = iv4options(Name,Value)
```

### Description

`opt = iv4options` creates the default options set for iv4.

`opt = iv4options(Name,Value)` creates an option set with the options specified by one or more `Name,Value` pair arguments.

### Input Arguments

#### Name-Value Pair Arguments

Specify optional comma-separated pairs of `Name,Value` arguments. `Name` is the argument name and `Value` is the corresponding value. `Name` must appear inside quotes. You can specify several name and value pair arguments in any order as `Name1,Value1,...,NameN,ValueN`.

#### InitialCondition — Handling of initial conditions

```
'auto' (default) | 'zero' | 'estimate'
```

Handling of initial conditions during estimation, specified as one of the following values:

- 'zero' — The initial condition is set to zero.
- 'estimate' — The initial condition is treated as an independent estimation parameter.
- 'auto' — The software chooses the initial condition handling method based on the estimation data.

**Focus — Error to be minimized**

'prediction' (default) | 'simulation'

Error to be minimized in the loss function during estimation, specified as the comma-separated pair consisting of 'Focus' and one of the following values:

- 'prediction' — The one-step ahead prediction error between measured and predicted outputs is minimized during estimation. As a result, the estimation focuses on producing a good predictor model.
- 'simulation' — The simulation error between measured and simulated outputs is minimized during estimation. As a result, the estimation focuses on making a good fit for simulation of model response with the current inputs.

The Focus option can be interpreted as a weighting filter in the loss function. For more information, see “Loss Function and Model Quality Metrics”.

**WeightingFilter — Weighting prefilter**

[] (default) | vector | matrix | cell array | linear system

Weighting prefilter applied to the loss function to be minimized during estimation. To understand the effect of WeightingFilter on the loss function, see “Loss Function and Model Quality Metrics”.

Specify WeightingFilter as one of the following values:

- [] — No weighting prefilter is used.
- Passbands — Specify a row vector or matrix containing frequency values that define desired passbands. You select a frequency band where the fit between estimated model and estimation data is optimized. For example, [w<sub>l</sub>, w<sub>h</sub>] where w<sub>l</sub> and w<sub>h</sub> represent lower and upper limits of a passband. For a matrix with several rows defining frequency passbands, [w<sub>1l</sub>, w<sub>1h</sub>; w<sub>2l</sub>, w<sub>2h</sub>; w<sub>3l</sub>, w<sub>3h</sub>; . . . ], the estimation algorithm uses the union of the frequency ranges to define the estimation passband.

Passbands are expressed in rad/TimeUnit for time-domain data and in FrequencyUnit for frequency-domain data, where TimeUnit and FrequencyUnit are the time and frequency units of the estimation data.

- SISO filter — Specify a single-input-single-output (SISO) linear filter in one of the following ways:
  - A SISO LTI model

- `{A,B,C,D}` format, which specifies the state-space matrices of a filter with the same sample time as estimation data.
- `{numerator,denominator}` format, which specifies the numerator and denominator of the filter as a transfer function with same sample time as estimation data.

This option calculates the weighting function as a product of the filter and the input spectrum to estimate the transfer function.

- **Weighting vector** — Applicable for frequency-domain data only. Specify a column vector of weights. This vector must have the same length as the frequency vector of the data set, `Data.Frequency`. Each input and output response in the data is multiplied by the corresponding weight at that frequency.

### **EnforceStability** — Control whether to enforce stability of model

`false` (default) | `true`

Control whether to enforce stability of estimated model, specified as the comma-separated pair consisting of 'EnforceStability' and either `true` or `false`.

Data Types: `logical`

### **EstimateCovariance** — Control whether to generate parameter covariance data

`true` (default) | `false`

Controls whether parameter covariance data is generated, specified as `true` or `false`.

If `EstimateCovariance` is `true`, then use `getcov` to fetch the covariance matrix from the estimated model.

### **Display** — Specify whether to display the estimation progress

`'off'` (default) | `'on'`

Specify whether to display the estimation progress, specified as one of the following values:

- `'on'` — Information on model structure and estimation results are displayed in a progress-viewer window.
- `'off'` — No progress or results information is displayed.

### **InputOffset** — Removal of offset from time-domain input data during estimation

`[]` (default) | vector of positive integers | matrix

Removal of offset from time-domain input data during estimation, specified as the comma-separated pair consisting of 'InputOffset' and one of the following:

- A column vector of positive integers of length  $Nu$ , where  $Nu$  is the number of inputs.
- `[]` — Indicates no offset.
- $Nu$ -by- $Ne$  matrix — For multi-experiment data, specify `InputOffset` as an  $Nu$ -by- $Ne$  matrix.  $Nu$  is the number of inputs, and  $Ne$  is the number of experiments.

Each entry specified by `InputOffset` is subtracted from the corresponding input data.

### **OutputOffset — Removal of offset from time-domain output data during estimation**

`[]` (default) | vector | matrix

Removal of offset from time-domain output data during estimation, specified as the comma-separated pair consisting of 'OutputOffset' and one of the following:

- A column vector of length  $Ny$ , where  $Ny$  is the number of outputs.
- `[]` — Indicates no offset.
- $Ny$ -by- $Ne$  matrix — For multi-experiment data, specify `OutputOffset` as a  $Ny$ -by- $Ne$  matrix.  $Ny$  is the number of outputs, and  $Ne$  is the number of experiments.

Each entry specified by `OutputOffset` is subtracted from the corresponding output data.

### **Advanced — Additional advanced options**

structure

Additional advanced options, specified as a structure with the following fields:

- `MaxSize` — Specifies the maximum number of elements in a segment when input-output data is split into segments.

`MaxSize` must be a positive integer.

**Default:** 250000

- `StabilityThreshold` — Specifies thresholds for stability tests.

`StabilityThreshold` is a structure with the following fields:

- `s` — Specifies the location of the right-most pole to test the stability of continuous-time models. A model is considered stable when its right-most pole is to the left of `s`.

**Default:** 0

- `z` — Specifies the maximum distance of all poles from the origin to test stability of discrete-time models. A model is considered stable if all poles are within the distance `z` from the origin.

**Default:** `1+sqrt(eps)`

## Output Arguments

**opt** — Options set for `iv4`

`iv4options` option set

Option set for `iv4`, returned as an `iv4options` option set.

## Examples

### Create Default Options Set for ARX Model Estimation Using 4-Stage Instrument Variable Method

```
opt = iv4options;
```

### Specify Options for ARX Model Estimation Using 4-Stage Instrument Variable Method

Create an options set for `iv4` using the `'backcast'` algorithm to initialize the state. Set `Display` to `'on'`.

```
opt = iv4options('InitialCondition','backcast','Display','on');
```

Alternatively, use dot notation to set the values of `opt`.

```
opt = iv4options;  
opt.InitialCondition = 'backcast';  
opt.Display = 'on';
```

## See Also

iv4

## Topics

“Loss Function and Model Quality Metrics”

**Introduced in R2012a**

## linapp

Linear approximation of nonlinear ARX and Hammerstein-Wiener models for given input

### Syntax

```
lm = linapp(nlmodel,u)
lm = linapp(nlmodel,umin,umax,nsample)
```

### Description

`lm = linapp(nlmodel,u)` computes a linear approximation of a nonlinear ARX or Hammerstein-Wiener model by simulating the model output for the input signal `u`, and estimating a linear model `lm` from `u` and the simulated output signal. `lm` is an `idpoly` model.

`lm = linapp(nlmodel,umin,umax,nsample)` computes a linear approximation of a nonlinear ARX or Hammerstein-Wiener model by first generating the input signal as a uniformly distributed white noise from the magnitude range `umin` and `umax` and (optionally) the number of samples.

### Input Arguments

`nlmodel`

Name of the `idnlarx` or `idnlhw` model object you want to linearize.

`u`

Input signal as an `iddata` object or a real matrix.

Dimensions of `u` must match the number of inputs in `nlmodel`.

`[umin,umax]`

Minimum and maximum input values for generating white-noise input with a magnitude in this rectangular range. The sample length of this signal is `nsample`.

nsample

Optional argument when you specify [umin, umax]. Specifies the length of the white-noise input.

**Default:** 1024.

## See Also

idnlarx | idnlarx/findop | idnlarx/linearize | idnlhw | idnlhw/findop | idnlhw/linearize

## Topics

“Linear Approximation of Nonlinear Black-Box Models”

**Introduced in R2007a**

## **linear**

Class representing linear nonlinearity estimator for nonlinear ARX models

### **Syntax**

```
lin=linear  
lin=linear('Parameters',Par)
```

### **Description**

`linear` is an object that stores the linear nonlinearity estimator for estimating nonlinear ARX models.

`lin=linear` instantiates the `linear` object.

`lin=linear('Parameters',Par)` instantiates the `linear` object and specifies optional values in the `Par` structure. For more information about this structure, see “linear Properties” on page 1-870.

### **linear Properties**

You can include property-value pairs in the constructor to specify the object.

After creating the object, you can use `get` or dot notation to access the object property values. For example:

```
% List Parameters values  
get(lin)  
% Get value of Parameters property  
lin.Parameters
```

Property Name	Description
Parameters	Structure containing the following fields: <ul style="list-style-type: none"> <li>LinearCoef: m-by-1 vector <math>L</math>.</li> <li>OutputOffset: Scalar <math>d</math>.</li> </ul>

## Examples

Estimate a nonlinear ARX model using the `linear` estimator with custom regressors for the following system:

$$y(t) = a_1y(t-1) + a_2y(t-2) + a_3u(t-1) + a_4y(t-1)u(t-2) + a_5|u(t)|u(t-3) + a_6,$$

where  $u$  is the input and  $y$  is the output.

```
% Create regressors y(t-1), y(t-2) and u(t-1).
orders = [2 1 1];
% Create an idnlarx model using linear estimator with custom regressors.
model = idnlarx(orders, linear, 'InputName', 'u', 'OutputName', 'y', ...
    'CustomRegressors', {'y(t-1)*u(t-2)', 'abs(u(t))*u(t-3)'});
% Estimate the model parameters a1, a2, ... a6.
EstimatedModel = nlarx(data, model)
```

---

**Note** The nonlinearity in the model is described by custom regressors only.

---

## Tutorials

“Estimate Nonlinear ARX Models at the Command Line”

## Tips

- `linear` is a linear (affine) function  $y = F(x)$ , defined as follows:

$$F(x) = xL + d$$

$y$  is scalar, and  $x$  is a 1-by-m vector.

- Use `evaluate(lin,x)` to compute the value of the function defined by the `linear` object `lin` at `x`.
- When creating a nonlinear ARX model using the constructor (`idnlarx`) or estimator (`nlarx`), you can specify a linear nonlinearity estimator using `[]`, instead of entering `linear` explicitly. For example:

```
m=idnlarx(orders,[]);
```

## Algorithms

When the `Focus` option in `nlarxOptions` is `'prediction'`, `linear` uses a fast, noniterative initialization and iterative search technique for estimating parameters. In most cases, iterative search requires only a few iterations.

When the `idnlarx` property `Focus='Simulation'`, `linear` uses an iterative technique for estimating parameters.

## See Also

`customreg` | `nlarx`

**Introduced in R2007a**

# idnlarx/linearize

Linearize nonlinear ARX model

## Syntax

```
SYS = linearize(NLSYS,U0,X0)
```

## Description

`SYS = linearize(NLSYS,U0,X0)` linearizes a nonlinear ARX model about the specified operating point `U0` and `X0`. The linearization is based on tangent linearization. For more information about the definition of states for `idnlarx` models, see “Definition of `idnlarx` States” on page 1-631.

## Input Arguments

- `NLSYS`: `idnlarx` model.
- `U0`: Matrix containing the constant input values for the model.
- `X0`: Model state values. The states of a nonlinear ARX model are defined by the time-delayed samples of input and output variables. For more information about the states of nonlinear ARX models, see the `getDelayInfo` reference page.

---

**Note** To estimate `U0` and `X0` from operating point specifications, use the `findop` command.

---

## Output Arguments

- `SYS` is an `idss` model.

When the Control System Toolbox product is installed, `SYS` is an LTI object.

## Examples

### Linearize Nonlinear ARX Model at Simulation Snapshot

Linearize a nonlinear ARX model around an operating point corresponding to a simulation snapshot at a specific time.

Load sample data.

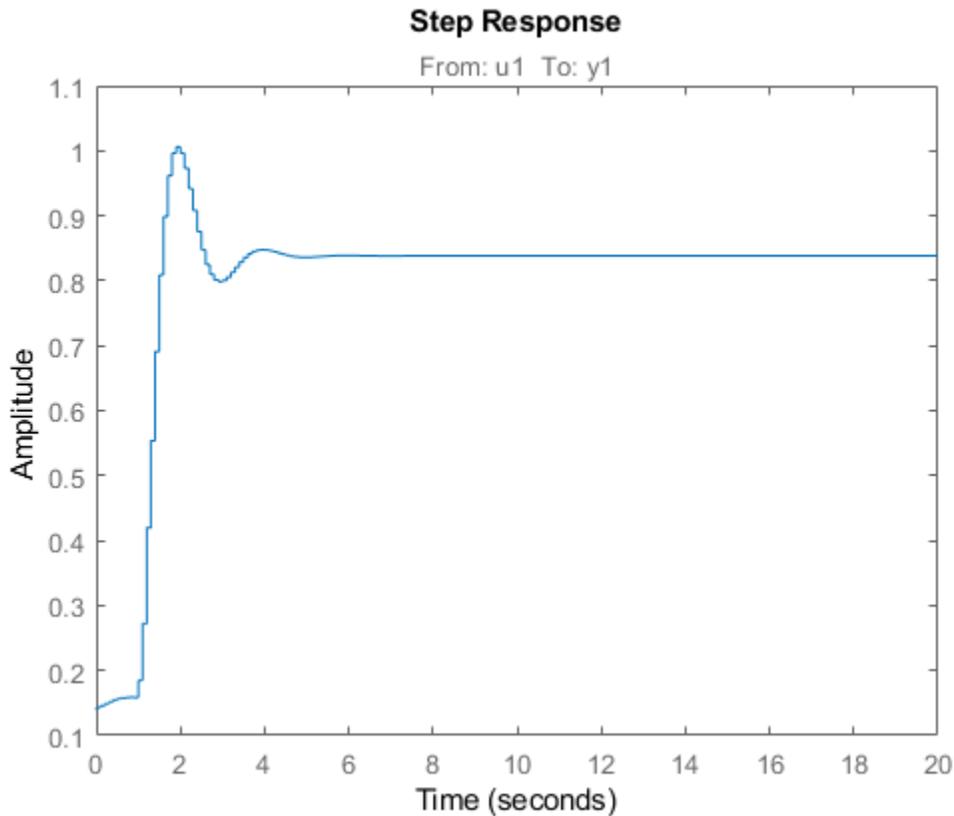
```
load iddata2
```

Estimate nonlinear ARX model from sample data.

```
nlsys = nlarx(z2,[4 3 10], 'tree', 'custom', ...  
    {'sin(y1(t-2)*u1(t))+y1(t-2)*u1(t)+u1(t).*u1(t-13)', ...  
    'y1(t-5)*y1(t-5)*y1(t-1)'}, 'nlr', [1:5, 7 9]);
```

Plot the response of the model for a step input.

```
step(nlsys, 20)
```



The step response is a steady-state value of 0.8383 at  $T = 20$  seconds.

Compute the operating point corresponding to  $T = 20$ .

```
stepinput = iddata([], [zeros(10,1); ones(200,1)], nlsys.Ts);
[x,u] = findop(nlsys, 'snapshot', 20, stepinput);
```

Linearize the model about the operating point corresponding to the model snapshot at  $T = 20$ .

```
sys = linearize(nlsys,u,x);
```

Validate the linear model.

Apply a small perturbation `delta_u` to the steady-state input of the nonlinear model `nlsys`. If the linear approximation is accurate, the following should match:

- The response of the nonlinear model `y_nl` to an input that is the sum of the equilibrium level and the perturbation `delta_u`.
- The sum of the response of the linear model to a perturbation input `delta_u` and the output equilibrium level.

Generate a 200-sample perturbation step signal with amplitude 0.1.

```
delta_u = [zeros(10,1); 0.1*ones(190,1)];
```

For a nonlinear system with a steady-state input of 1 and a steady-state output of 0.8383, compute the steady-state response `y_nl` to the perturbed input `u_nl`. Use equilibrium state values `x` computed previously as initial conditions.

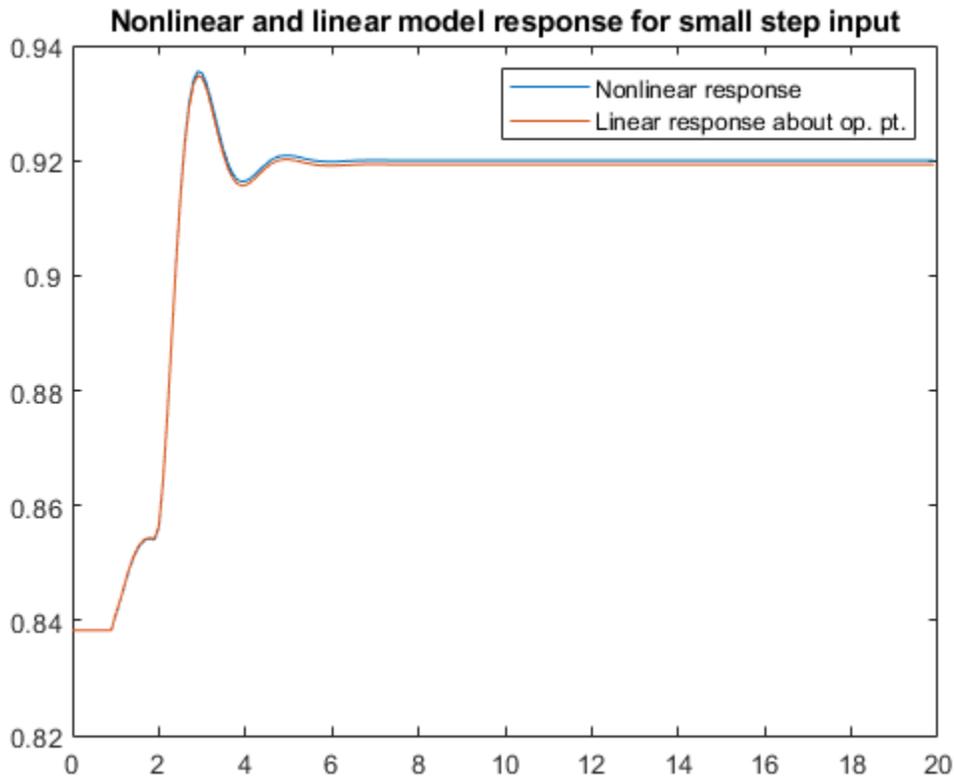
```
u_nl = 1 + delta_u;  
y_nl = sim(nlsys,u_nl,x);
```

Compute response of linear model to perturbation input and add it to the output equilibrium level.

```
y_lin = 0.8383 + lsim(sys,delta_u);
```

Compare the response of nonlinear and linear models.

```
time = [0:0.1:19.9]';  
plot(time,y_nl,time,y_lin)  
legend('Nonlinear response','Linear response about op. pt.')  
title(['Nonlinear and linear model response for small step input'])
```



## Algorithms

The following equations govern the dynamics of an `idnlarx` model:

$$X(t+1) = AX(t) + B\tilde{u}(t)$$

$$y(t) = f(X, u)$$

where  $X(t)$  is a state vector,  $u(t)$  is the input, and  $y(t)$  is the output.  $A$  and  $B$  are constant matrices.  $\tilde{u}(t)$  is  $[y(t), u(t)]^T$ .

The output at the operating point is given by

$$y^* = f(X^*, u^*)$$

where  $X^*$  and  $u^*$  are the state vector and input at the operating point.

The linear approximation of the model response is as follows:

$$\begin{aligned}\Delta X(t+1) &= (A + B_1 f_X) \Delta X(t) + (B_1 f_u + B_2) \Delta u(t) \\ \Delta y(t) &= f_X \Delta X(t) + f_u \Delta u(t)\end{aligned}$$

where

- $\Delta X(t) = X(t) - X^*(t)$
- $\Delta u(t) = u(t) - u^*(t)$
- $\Delta y(t) = y(t) - y^*(t)$
- $B\tilde{U} = [B_1, B_2] \begin{bmatrix} Y \\ U \end{bmatrix} = B_1 Y + B_2 U$
- $f_X = \left. \frac{\partial}{\partial X} f(X, U) \right|_{X^*, U^*}$
- $f_U = \left. \frac{\partial}{\partial U} f(X, U) \right|_{X^*, U^*}$

---

**Note** For linear approximations over larger input ranges, use `linapp`.

---

## See Also

`getDelayInfo` | `idnlarx` | `idnlarx/findop` | `linapp`

## Topics

“Linear Approximation of Nonlinear Black-Box Models”

**Introduced in R2014b**

## idnlhw/linearize

Linearize Hammerstein-Wiener model

### Syntax

```
SYS = linearize(NLSYS,U0)  
SYS = linearize(NLSYS,U0,X0)
```

### Description

`SYS = linearize(NLSYS,U0)` linearizes a Hammerstein-Wiener model around the equilibrium operating point. When using this syntax, equilibrium state values for the linearization are calculated automatically using `U0`.

`SYS = linearize(NLSYS,U0,X0)` linearizes the `idnlhw` model `NLSYS` around the operating point specified by the input `U0` and state values `X0`. In this usage, `X0` need not contain equilibrium state values. For more information about the definition of states for `idnlhw` models, see “Definition of `idnlhw` States” on page 1-672.

The output is a linear model that is the best linear approximation for inputs that vary in a small neighborhood of a constant input  $u(t) = U$ . The linearization is based on tangent linearization.

### Input Arguments

- `NLSYS`: `idnlhw` model.
- `U0`: Matrix containing the constant input values for the model.
- `X0`: Operating point state values for the model.

---

**Note** To estimate `U0` and `X0` from operating point specifications, use the `findop` command.

---

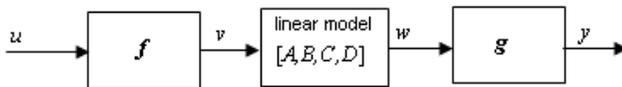
## Output Arguments

- SYS is an idss model.

When the Control System Toolbox product is installed, SYS is an LTI object.

## Algorithms

The `idnllhw` model structure represents a nonlinear system using a linear system connected in series with one or two static nonlinear systems. For example, you can use a static nonlinearity to simulate saturation or dead-zone behavior. The following figure shows the nonlinear system as a linear system that is modified by static input and output nonlinearities, where function  $f$  represents the input nonlinearity,  $g$  represents the output nonlinearity, and  $[A,B,C,D]$  represents a state-space parameterization of the linear model.



The following equations govern the dynamics of an `idnllhw` model:

$$v(t) = f(u(t))$$

$$X(t+1) = AX(t) + Bv(t)$$

$$w(t) = CX(t) + Dv(t)$$

$$y(t) = g(w(t))$$

where

- $u$  is the input signal
- $v$  and  $w$  are intermediate signals (outputs of the input nonlinearity and linear model respectively)
- $y$  is the model output

The linear approximation of the Hammerstein-Wiener model around an operating point  $(X^*, u^*)$  is as follows:

$$\begin{aligned}\Delta X(t+1) &= A\Delta X(t) + Bf_u \Delta u(t) \\ \Delta y(t) &\approx g_w C\Delta X(t) + g_w Df_u \Delta u(t)\end{aligned}$$

where

- $\Delta X(t) = X(t) - X^*(t)$
- $\Delta u(t) = u(t) - u^*(t)$
- $\Delta y(t) = y(t) - y^*(t)$
- $f_u = \left. \frac{\partial}{\partial u} f(u) \right|_{u=u^*}$
- $g_w = \left. \frac{\partial}{\partial w} g(w) \right|_{w=w^*}$

where  $y^*$  is the output of the model corresponding to input  $u^*$  and state vector  $X^*$ ,  $v^* = f(u^*)$ , and  $w^*$  is the response of the linear model for input  $v^*$  and state  $X^*$ .

---

**Note** For linear approximations over larger input ranges, use `linapp`. For more information, see the `linapp` reference page.

---

## See Also

`idnlhw` | `idnlhw/findop` | `linapp`

## Topics

“Linear Approximation of Nonlinear Black-Box Models”

**Introduced in R2014b**

# lsim

Simulate time response of dynamic system to arbitrary inputs

## Syntax

```
lsim(sys,u,t)
lsim(sys,u,t,x0)
lsim(sys,u,t,x0,method)
lsim(sys1,...,sysn,u,t)
lsim(sys1,LineStyle1,...,sysN,LineStyleN,u,t)
y = lsim(____)
[y,t,x] = lsim(____)
lsim(sys)
```

## Description

`lsim` simulates the (time) response of continuous or discrete linear systems to arbitrary inputs. When invoked without left-hand arguments, `lsim` plots the response on the screen.

`lsim(sys,u,t)` produces a plot of the time response of the dynamic system model `sys` to the input history, `t,u`. The vector `t` specifies the time samples for the simulation (in system time units, specified in the `TimeUnit` property of `sys`), and consists of regularly spaced time samples:

```
t = 0:dt:Tfinal
```

The input `u` is an array having as many rows as time samples (`length(t)`) and as many columns as system inputs. For instance, if `sys` is a SISO system, then `u` is a t-by-1 vector. If `sys` has three inputs, then `u` is a t-by-3 array. Each row `u(i,:)` specifies the input value(s) at the time sample `t(i)`. The signal `u` also appears on the plot.

The model `sys` can be continuous or discrete, SISO or MIMO. In discrete time, `u` must be sampled at the same rate as the system. In this case, the input `t` is redundant and can be omitted or set to an empty matrix. In continuous time, the time sampling `dt = t(2) - t(1)` is used to discretize the continuous model. If `dt` is too large (undersampling), `lsim`

issues a warning suggesting that you use a more appropriate sample time, but will use the specified sample time. See “Algorithms” on page 1-888 for a discussion of sample times.

`lsim(sys,u,t,x0)` further specifies an initial condition `x0` for the system states. This syntax applies only when `sys` is a state-space model. `x0` is a vector whose entries are the initial values of the corresponding states of `sys`.

`lsim(sys,u,t,x0,method)` explicitly specifies how the input values should be interpolated between samples, when `sys` is a continuous-time system. Specify `method` as one of the following values:

- 'zoh' — Use zero-order hold
- 'foh' — Use linear interpolation (first-order hold)

If you do not specify a method, `lsim` selects the interpolation method automatically based on the smoothness of the signal `u`.

`lsim(sys1,...,sysn,u,t)` simulates the responses of several dynamic system models to the same input history `t,u` and plots these responses on a single figure. You can also use the `x0` and `method` input arguments when computing the responses of multiple models.

`lsim(sys1,LineStyle1,...,sysN,LineStyleN,u,t)` specifies the line style, marker, and color of each of the system responses in the plot. You can also use the `x0` and `method` input arguments with this syntax. Each `LineStyle` argument is specified as a vector of one, two, or three characters. The characters can appear in any order. For example, the following code plots the response of `sys1` as a yellow dotted line and the response of `sys2` as a green dashed line:

```
lsim(sys1,'y:',sys2,'g--',u,t,x0)
```

For more information about configuring this argument, see the `LineStyle` input argument of the `plot` function.

`y = lsim(____)` returns the system response `y`, sampled at the same times as the input (`t`). The output `y` is an array having as many rows as time samples (`length(t)`) and as many columns as system outputs. No plot is drawn on the screen. You can use this syntax with any of the input arguments described in previous syntaxes except the `LineStyle` arguments.

`[y,t,x] = lsim(____)` also returns the time vector `t` used for simulation and the state trajectories `x` (for state-space models only). The output `x` has as many rows as time

samples (`length(t)`) and as many columns as system states. You can use this syntax with any of the input arguments described in previous syntaxes except the `LineSpec` arguments.

`lsim(sys)` opens the Linear Simulation Tool GUI. For more information about working with this GUI, see [Working with the Linear Simulation Tool \(Control System Toolbox\)](#).

## Examples

### Simulate Response to Square Wave

Simulate and plot the response of the following system to a square wave with period of four seconds:

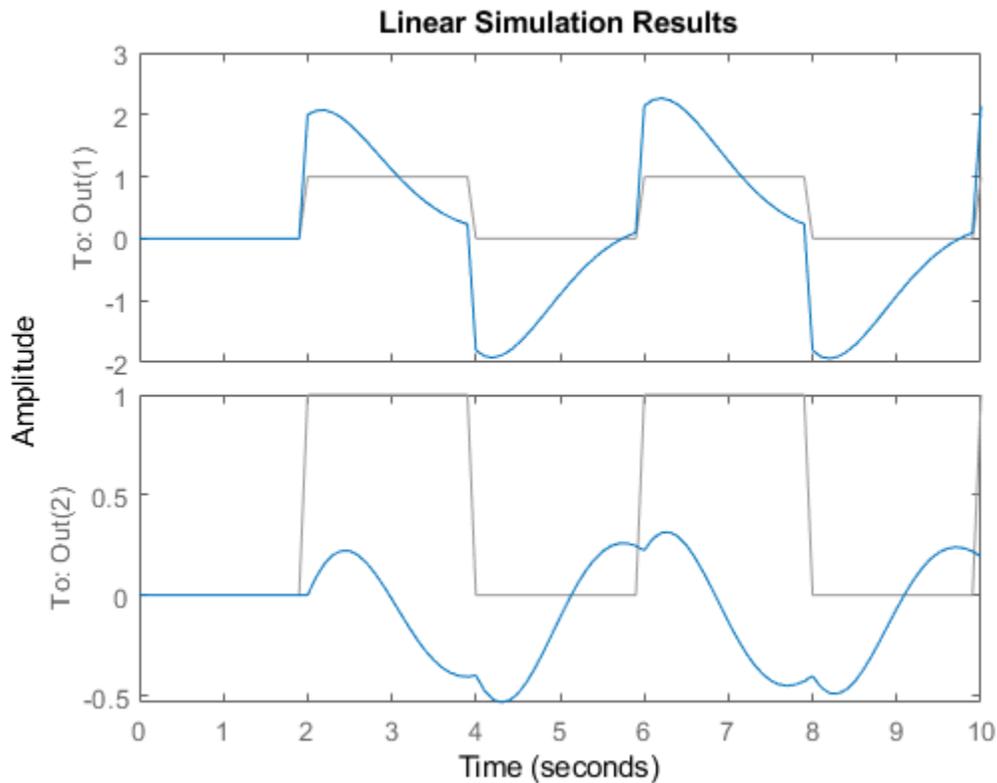
$$H(s) = \begin{bmatrix} \frac{2s^2 + 5s + 1}{s^2 + 2s + 3} \\ \frac{s - 1}{s^2 + s + 5} \end{bmatrix}.$$

Create the transfer function, and generate the square wave with `gensig`. Sample every 0.1 second during 10 seconds.

```
H = [tf([2 5 1],[1 2 3]);tf([1 -1],[1 1 5])];  
[u,t] = gensig('square',4,10,0.1);
```

Then simulate with `lsim`.

```
lsim(H,u,t)
```



The plot displays both the applied signal and the response.

### Simulate Response of Identified Model

Load estimation data to estimate a model.

```
load(fullfile(matlabroot, 'toolbox', 'ident', 'iddemos', 'data', 'dcmotordata'));  
z = iddata(y,u,0.1, 'Name', 'DC-motor');
```

`z` is an `iddata` object that stores the 1-input 2-output estimation data with sample time 0.1 seconds.

Estimate a state-space model of order 4 using estimation data `z`.

```
[sys,x0] = n4sid(z,4);
```

`sys` is the estimated model and `x0` are the estimated initial states.

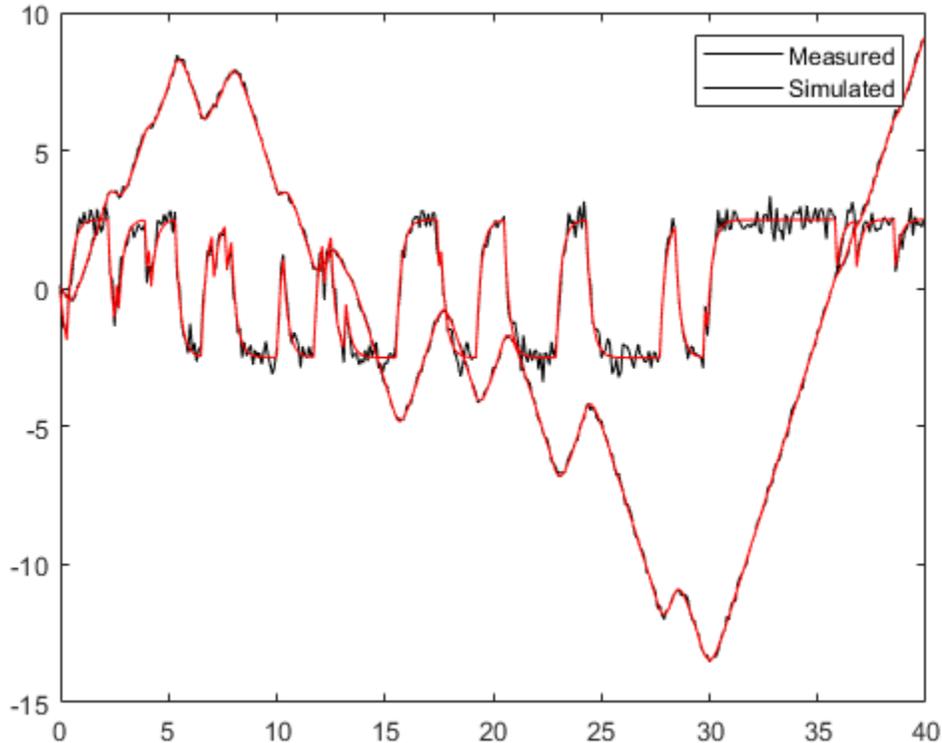
Simulate the response of `sys` using the same input data as the one used for estimation and the initial states returned by the estimation command.

```
[y,t,x] = lsim(sys,z.InputData,[],x0);
```

Here, `y` is the system response, `t` is the time vector used for simulation, and `x` is the state trajectory.

Compare the simulated response `y` to measured response `z.OutputData` for both outputs.

```
plot(t,z.OutputData,'k',t,y,'r')  
legend('Measured','Simulated')
```



## Algorithms

Discrete-time systems are simulated with `ltitr` (state space) or `filter` (transfer function and zero-pole-gain).

Continuous-time systems are discretized with `c2d` using either the `'zoh'` or `'foh'` method (`'foh'` is used for smooth input signals and `'zoh'` for discontinuous signals such as pulses or square waves). The sample time is set to the spacing `dt` between the user-supplied time samples `t`.

The choice of sample time can drastically affect simulation results. To illustrate why, consider the second-order model

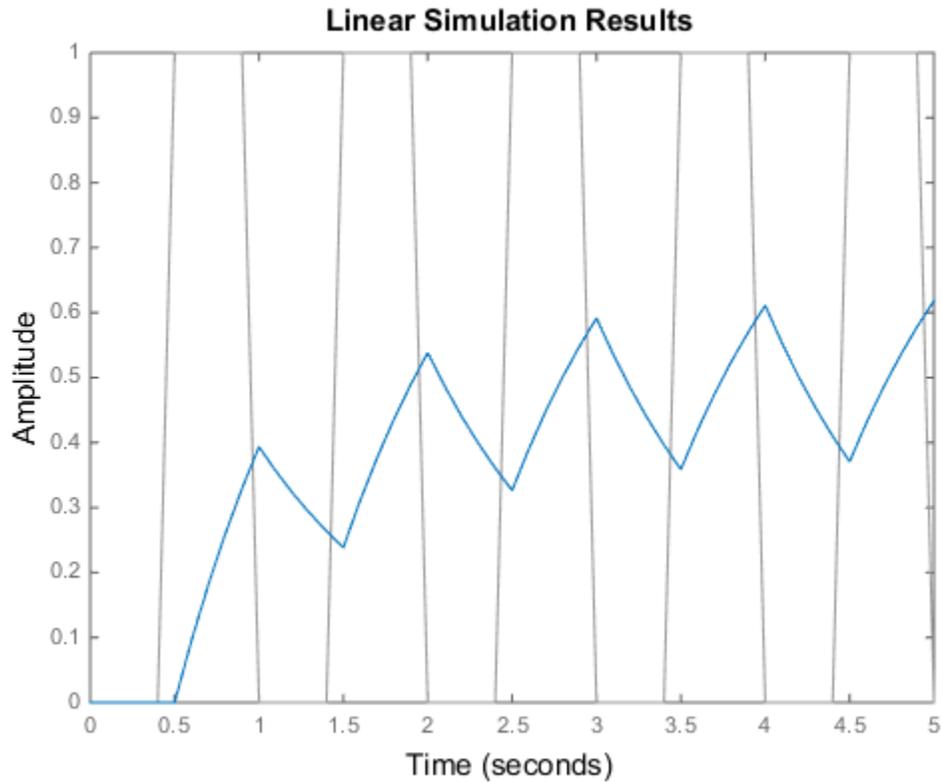
$$H(s) = \frac{\omega^2}{s^2 + 2s + \omega^2}, \quad \omega = 62.83$$

To simulate its response to a square wave with period 1 second, you can proceed as follows:

```
w2 = 62.83^2;
h = tf(w2,[1 2 w2]);
t = 0:0.1:5;           % vector of time samples
u = (rem(t,1) >= 0.5); % square wave values
lsim(h,u,t)
```

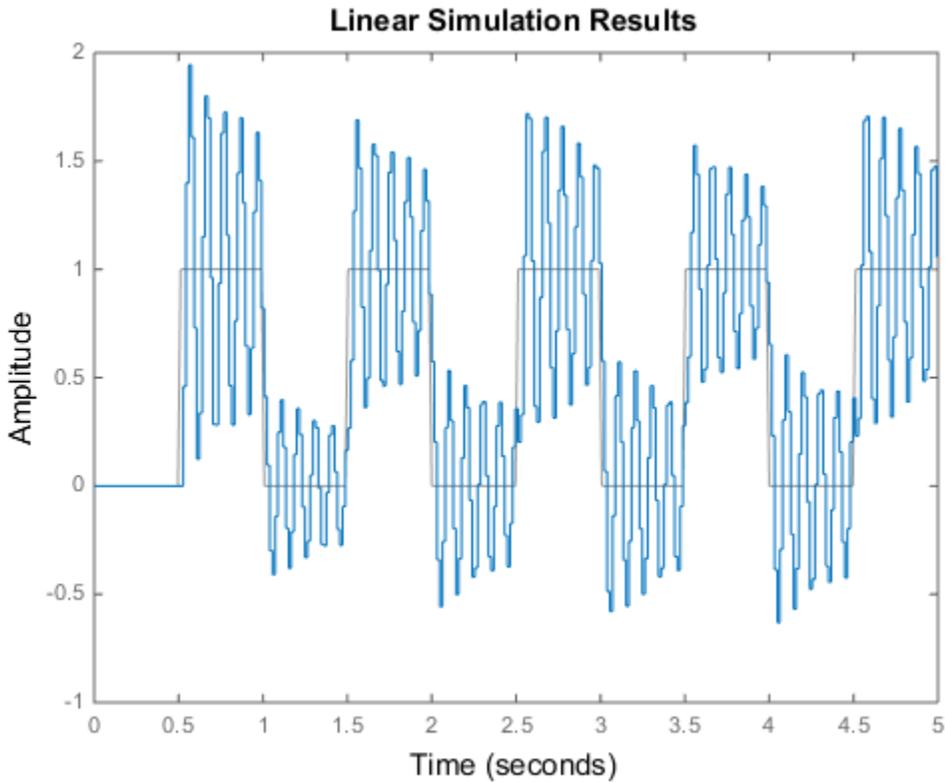
`lsim` evaluates the specified sample time, and issues a warning:

```
Warning: Input signal is undersampled. Sample every 0.016 sec or
faster.
```



To improve on this response, discretize  $H(s)$  using the recommended sample time:

```
dt = 0.016;  
ts = 0:dt:5;  
us = (rem(ts,1) >= 0.5);  
hd = c2d(h,dt);  
lsim(hd,us,ts)
```



This response exhibits strong oscillatory behavior that is hidden in the undersampled version.

## See Also

[Linear System Analyzer](#) | [gensig](#) | [impulse](#) | [initial](#) | [lsiminfo](#) | [sim](#) | [step](#)

**Introduced in R2012a**

## lsiminfo

Compute linear response characteristics

### Syntax

```
S = lsiminfo(y,t,yfinal)
S = lsiminfo(y,t)
S = lsiminfo(...,'SettlingTimeThreshold',ST)
```

### Description

`S = lsiminfo(y,t,yfinal)` takes the response data  $(t,y)$  and a steady-state value `yfinal` and returns a structure `S` containing the following performance indicators:

- `SettlingTime` — Settling time
- `Min` — Minimum value of `Y`
- `MinTime` — Time at which the minimum value is reached
- `Max` — Maximum value of `Y`
- `MaxTime` — Time at which the maximum value is reached

For SISO responses, `t` and `y` are vectors with the same length `NS`. For responses with `NY` outputs, you can specify `y` as an `NS`-by-`NY` array and `yfinal` as a `NY`-by-1 array. `lsiminfo` then returns an `NY`-by-1 structure array `S` of performance metrics for each output channel.

`S = lsiminfo(y,t)` uses the last sample value of `y` as steady-state value `yfinal`. `s = lsiminfo(y)` assumes `t = 1:NS`.

`S = lsiminfo(...,'SettlingTimeThreshold',ST)` lets you specify the threshold `ST` used in the settling time calculation. The response has settled when the error  $|y(t) - y_{\text{final}}|$  becomes smaller than a fraction `ST` of its peak value. The default value is `ST=0.02` (2%).

## Examples

### Compute Response Characteristics of a Transfer Function

Create the following continuous-time transfer function:

$$H(s) = \frac{s - 1}{s^3 + 2s^2 + 3s + 4}$$

```
sys = tf([1 -1],[1 2 3 4]);
```

Calculate the impulse response.

```
[y,t] = impulse(sys);
```

`impulse` returns the output response `y` and the time vector `t` used for simulation.

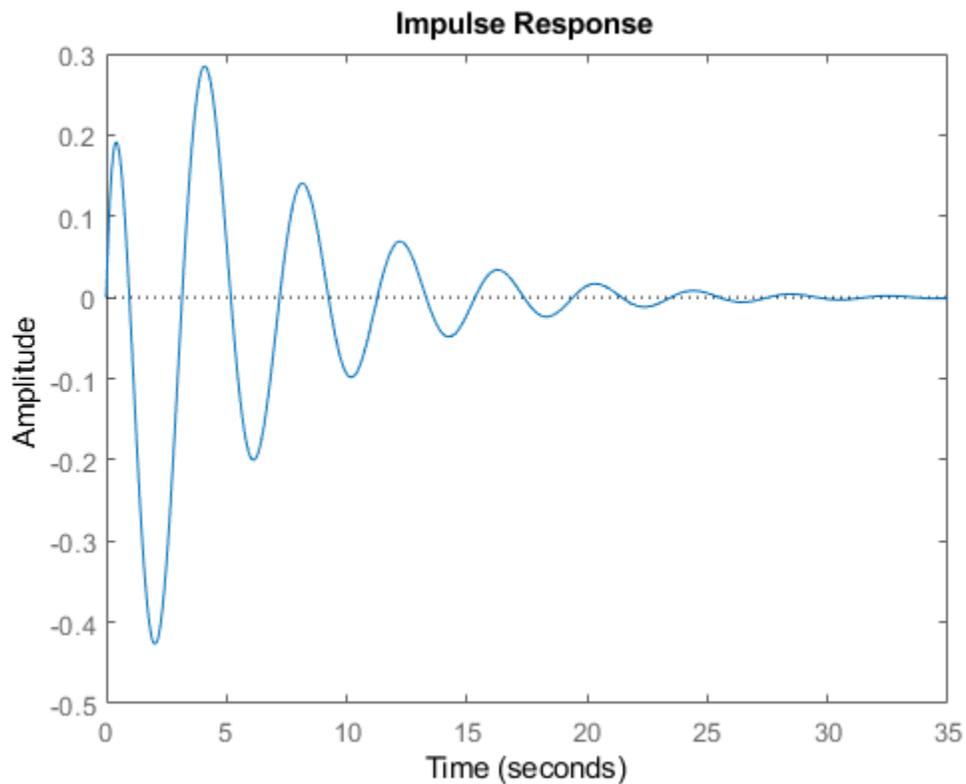
Compute the response characteristics using a final response value of 0.

```
s = lsiminfo(y,t,0)
```

```
s = struct with fields:
  SettlingTime: 22.8700
             Min: -0.4268
             MinTime: 2.0088
             Max: 0.2847
             MaxTime: 4.0733
```

You can plot the impulse response and verify these response characteristics. For example, the time at which the minimum response value (`MinTime`) is reached is approximately 2 seconds.

```
impulse(sys)
```



## See Also

`impulse` | `initial` | `lsim` | `stepinfo`

**Introduced in R2012a**

# lsimplot

Simulate response of dynamic system to arbitrary inputs and return plot handle

## Syntax

```
h = lsimplot(sys)
lsimplot(sys1,sys2,...)
lsimplot(sys,u,t)
lsimplot(sys,u,t,x0)
lsimplot(sys1,sys2,...,u,t,x0)
lsimplot(AX,...)
lsimplot(..., plotoptions)
lsimplot(sys,u,t,x0,'zoh')
lsimplot(sys,u,t,x0,'foh')
```

## Description

`h = lsimplot(sys)` opens the Linear Simulation Tool for the dynamic system model `sys`, which enables interactive specification of driving input(s), the time vector, and initial state. It also returns the plot handle `h`. You can use this handle to customize the plot with the `getoptions` and `setoptions` commands. Type

```
help timeoptions
```

for a list of available plot options.

`lsimplot(sys1,sys2,...)` opens the Linear Simulation Tool for multiple models `sys1,sys2,...`. Driving inputs are common to all specified systems but initial conditions can be specified separately for each.

`lsimplot(sys,u,t)` plots the time response of the model `sys` to the input signal described by `u` and `t`. The time vector `t` consists of regularly spaced time samples (in system time units, specified in the `TimeUnit` property of `sys`). For MIMO systems, `u` is a matrix with as many columns as inputs and whose `i`th row specifies the input value at time `t(i)`. For SISO systems `u` can be specified either as a row or column vector. For example,

```
t = 0:0.01:5;
u = sin(t);
lsimplot(sys,u,t)
```

simulates the response of a single-input model `sys` to the input  $u(t)=\sin(t)$  during 5 seconds.

For discrete-time models, `u` should be sampled at the same rate as `sys` (`t` is then redundant and can be omitted or set to the empty matrix).

For continuous-time models, choose the sampling period  $t(2) - t(1)$  small enough to accurately describe the input `u`. `lsim` issues a warning when `u` is undersampled, and hidden oscillations can occur.

`lsimplot(sys,u,t,x0)` specifies the initial state vector `x0` at time `t(1)` (for state-space models only). `x0` is set to zero when omitted.

`lsimplot(sys1,sys2,...,u,t,x0)` simulates the responses of multiple LTI models `sys1,sys2,...` on a single plot. The initial condition `x0` is optional. You can also specify a color, line style, and marker for each system, as in

```
lsimplot(sys1,'r',sys2,'y--',sys3,'gx',u,t)
```

`lsimplot(AX,...)` plots into the axes with handle `AX`.

`lsimplot(..., plotoptions)` plots the initial condition response with the options specified in `plotoptions`. Type

```
help timeoptions
```

for more detail.

For continuous-time models, `lsimplot(sys,u,t,x0,'zoh')` or `lsimplot(sys,u,t,x0,'foh')` explicitly specifies how the input values should be interpolated between samples (zero-order hold or linear interpolation). By default, `lsimplot` selects the interpolation method automatically based on the smoothness of the signal `u`.

## See Also

`getoptions` | `lsim` | `setoptions`

**Introduced in R2012a**

## **mag2db**

Convert magnitude to decibels (dB)

### **Syntax**

`ydb = mag2db(y)`

### **Description**

`ydb = mag2db(y)` returns the corresponding decibel (dB) value *ydb* for a given magnitude *y*. The relationship between magnitude and decibels is  $ydb = 20 \log_{10}(y)$ .

### **See Also**

`db2mag`

**Introduced in R2008a**

## merge (iddata)

Merge data sets into iddata object

### Syntax

```
dat = merge(dat1,dat2,...,datN)
```

### Description

`dat` collects the data sets in `dat1`, ..., `datN` into one `iddata` object, with several *experiments*. The number of experiments in `dat` will be the sum of the number of experiments in `datk`. For the merging to be allowed, a number of conditions must be satisfied:

- All of `datk` must have the same number of input channels, and the `InputNames` must be the same.
- All of `datk` must have the same number of output channels, and the `OutputNames` must be the same. If some input or output channel is lacking in one experiment, it can be replaced by a vector of NaNs to conform with these rules.
- If the `ExperimentNames` of `datk` have been specified as something other than the default 'Exp1', 'Exp2', etc., they must all be unique. If default names overlap, they are modified so that `dat` will have a list of unique `ExperimentNames`.

The sampling intervals, the number of observations, and the input properties (`Period`, `InterSample`) might be different in the different experiments.

You can retrieve the individual experiments by using the command `getexp`. You can also retrieve them by subreferencing with a fourth index.

```
dat1 = dat(:, :, :, ExperimentNumber)
```

or

```
dat1 = dat(:, :, :, ExperimentName)
```

Storing multiple experiments as one `iddata` object can be very useful for handling experimental data that has been collected on different occasions, or when a data set has

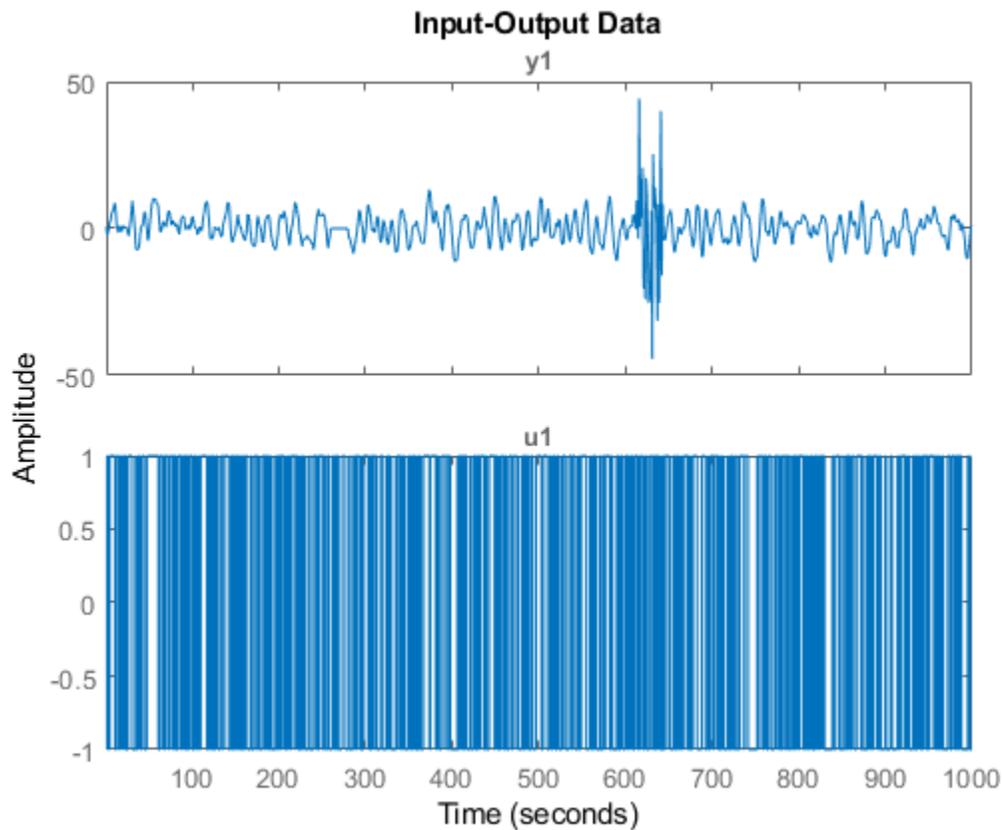
been split up to remove “bad” portions of the data. All the toolbox routines accept multiple-experiment data.

## Examples

### Merge Multiple Data Sets

Remove bad portions of data to estimate models without the bad data destroying the estimate.

```
load iddemo8;  
plot(dat);
```



Bad portions of data are detected around sample 250 to 280 and between samples 600 to 650. Cut out these bad portions to form a multiple-experiment data set and merge the data.

```
dat = merge(dat(1:250),dat(281:600),dat(651:1000));
```

You can use the first two experiments to estimate a model and the third experiment to validate the model.

```
dat_est = getexp(dat,[1,2]);
m = ssest(dat_est,2);
dat_val = getexp(dat,3);
```

## **See Also**

getexp | iddata | merge

## **Topics**

“Dealing with Multi-Experiment Data and Merging Models”

“Create Multiexperiment Data at the Command Line”

**Introduced before R2006a**

## merge

Merge estimated models

### Syntax

```
m = merge(m1,m2,...,mN)
[m,tv] = merge(m1,m2)
```

### Description

`m = merge(m1,m2,...,mN)` merges estimated models. The models `m1,m2,...,mN` must all be of the same structure, just differing in parameter values and covariance matrices. Then `m` is the merged model, where the parameter vector is a statistically weighted mean (using the covariance matrices to determine the weights) of the parameters of `mk`.

`[m,tv] = merge(m1,m2)` returns a test variable `tv`. When two models are merged,

```
[m, tv] = merge(m1,m2)
```

`tv` is  $\chi^2$  distributed with `n` degrees of freedom, if the parameters of `m1` and `m2` have the same means. Here `n` is the length of the parameter vector. A large value of `tv` thus indicates that it might be questionable to merge the models.

For `idfrd` models, `merge` is a statistical average of two responses in the individual models, weighted using inverse variances. You can only merge two `idfrd` models with responses at the same frequencies and nonzero covariances.

Merging models is an alternative to merging data sets and estimating a model for the merged data.

```
load iddata1 z1;
load iddata2 z2;
m1 = arx(z1,[2 3 4]);
m2 = arx(z2,[2 3 4]);
ma = merge(m1,m2);
```

and

```
mb = arx(merge(z1,z2),[2 3 4]);
```

result in models `ma` and `mb` that are related and should be close. The difference is that merging the data sets assumes that the signal-to-noise ratios are about the same in the two experiments. Merging the models allows one model to be much more uncertain, for example, due to more disturbances in that experiment. If the conditions are about the same, we recommend that you merge data rather than models, since this is more efficient and typically involves better conditioned calculations.

### **See Also**

`append`

**Introduced in R2007a**

# midprefs

Specify location for file containing System Identification app startup information

## Syntax

```
midprefs  
midprefs(path)
```

## Description

The **System Identification** app allows a large number of variables for customized choices. These include the window layout, the default choices of plot options, and names and directories of the four most recent sessions with the **System Identification** app. This information is stored in the file `idprefs.mat`, which should be placed on the user's `MATLABPATH`. The default, automatic location for this file is in the same folder as the user's `startup.m` file.

`midprefs` is used to select or change the folder where you store `idprefs.mat`. Either type `midprefs` and follow the instructions, or give the folder name as the argument. Include all folder delimiters, as in the PC case:

```
midprefs('c:\matlab\toolbox\local\')
```

or in the UNIX<sup>®</sup> case"

```
midprefs('/home/ljung/matlab/')
```

## See Also

**Apps**  
**System Identification**

**Introduced before R2006a**

## misdata

Reconstruct missing input and output data

### Syntax

```
Datae = misdata(Data)
Datae = misdata(Data,Model)
Datae = misdata(Data,MaxIterations,Tol)
```

### Description

`Datae = misdata(Data)` reconstructs missing input and output data. `Data` is time-domain input-output data in the `iddata` object format. Missing data samples (both in inputs and in outputs) are entered as NaNs. `Datae` is an `iddata` object where the missing data has been replaced by reasonable estimates.

`Datae = misdata(Data,Model)` specifies a model used for the reconstruction of missing data. `Model` is any linear identified model (`idtf`, `idproc`, `idgrey`, `idpoly`, `idss`). If no suitable model is known, it is estimated in an iterative fashion using default order state-space models.

`Datae = misdata(Data,MaxIterations,Tol)` specifies maximum number of iterations and tolerance. `MaxIterations` is the maximum number of iterations carried out (the default is 10). The iterations are terminated when the difference between two consecutive data estimates differs by less than `Tol`%. The default value of `Tol` is 1.

### Examples

#### Reconstruct Missing Data Using Specified Model

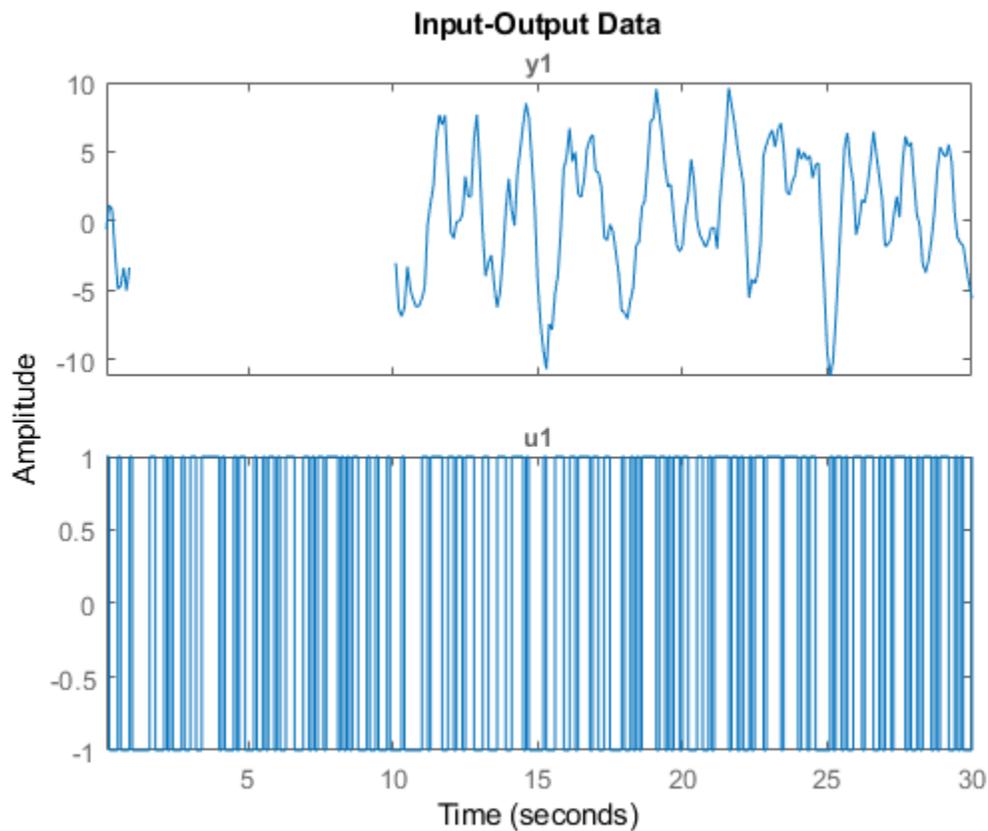
Load data with missing data points.

```
load(fullfile(matlabroot,'examples','ident','missing_data.mat'))
```

`missing_data` is an `iddata` object containing input-output data.

Plot the data.

```
plot(missing_data)
```



The output data contains missing data between indices 10 and 100.

To reconstruct missing data using a specified model, estimate the model using measured data that has no missing samples. In this example, estimate a transfer function model with 2 poles.

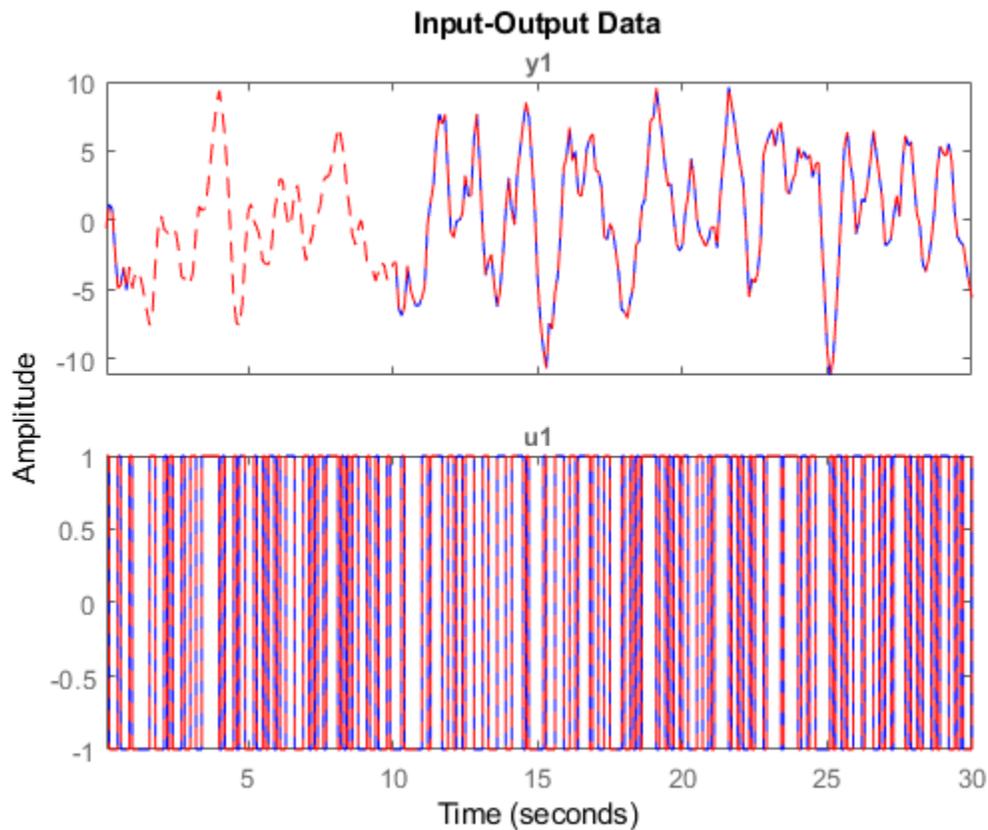
```
data2 = missing_data(101:end);  
model = tfest(data2,2);
```

Reconstruct the missing data.

```
datae = misdata(missing_data,model);
```

Plot the original and reconstructed data.

```
plot(missing_data,'b',datae,'--r')
```



If you do not specify a model for reconstructing the data, the software alternates between estimating missing data and estimating models, based on the current data reconstruction.

## Algorithms

For a given model, the missing data is estimated as parameters so as to minimize the output prediction errors obtained from the reconstructed data. See Section 14.2 in Ljung (1999). Treating missing outputs as parameters is not the best approach from a statistical point of view, but is a good approximation in many cases.

When no model is given, the algorithm alternates between estimating missing data and estimating models, based on the current reconstruction.

## See Also

`advice` | `arx` | `pexcit` | `tfest`

**Introduced before R2006a**

## n4sid

Estimate state-space model using subspace method

### Syntax

```
sys = n4sid(data,nx)
sys = n4sid(data,nx,Name,Value)
sys = n4sid( ____,opt)
[sys,x0] = n4sid( ____,opt)
```

### Description

`sys = n4sid(data,nx)` estimates an  $n_x$  order state-space model, `sys`, using measured input-output data, `data`.

`sys` is an `idss` model representing the system:

$$\begin{aligned}\dot{x}(t) &= Ax(t) + Bu(t) + Ke(t) \\ y(t) &= Cx(t) + Du(t) + e(t)\end{aligned}$$

$A, B, C$ , and  $D$  are state-space matrices.  $K$  is the disturbance matrix.  $u(t)$  is the input,  $y(t)$  is the output,  $x(t)$  is the vector of  $n_x$  states and  $e(t)$  is the disturbance.

All the entries of the  $A$ ,  $B$ ,  $C$ , and  $K$  matrices are free estimation parameters by default.  $D$  is fixed to zero by default, meaning that there is no feedthrough, except for static systems ( $n_x=0$ ).

`sys = n4sid(data,nx,Name,Value)` specifies additional attributes of the state-space structure using one or more `Name, Value` pair arguments. Use the `Form`, `Feedthrough`, and `DisturbanceModel` name-value pair arguments to modify the default behavior of the  $A$ ,  $B$ ,  $C$ ,  $D$ , and  $K$  matrices.

`sys = n4sid( ____,opt)` specifies estimation options, `opt`, that configure the initial states, estimation objective, and subspace algorithm related choices to be used for estimation.

`[sys,x0] = n4sid( ___ )` also returns the estimated initial state.

## Input Arguments

### **data**

Estimation data.

For time domain estimation, `data` is an `iddata` object containing the input and output signal values.

For frequency domain estimation, `data` can be one of the following:

- Recorded frequency response data (`frd` or `idfrd`)
- `iddata` object with its properties specified as follows:
  - `InputData` — Fourier transform of the input signal
  - `OutputData` — Fourier transform of the output signal
  - `Domain` — 'Frequency'

For multiexperiment data, the sample times and intersample behavior of all the experiments must match.

You can only estimate continuous-time models using continuous-time frequency domain data. You can estimate both continuous-time and discrete-time models (of sample time matching that of `data`) using time-domain data and discrete-time frequency domain data.

### **nx**

Order of estimated model.

Specify `nx` as a positive integer. `nx` may be a scalar or a vector. If `nx` is a vector, then `n4sid` creates a plot which you can use to choose a suitable model order. The plot shows the Hankel singular values for models of different orders. States with relatively small Hankel singular values can be safely discarded. A default choice is suggested in the plot.

You can also specify `nx` as `'best'`, in which case the optimal order is automatically chosen from `nx = 1, ..., 10`.

## **opt**

Estimation options.

`opt` is an options set, created using `n4sidOptions`, which specifies options including:

- Estimation objective
- Handling of initial conditions
- Subspace algorithm related choices

## **Name-Value Pair Arguments**

Specify optional comma-separated pairs of `Name`, `Value` arguments. `Name` is the argument name and `Value` is the corresponding value. `Name` must appear inside quotes. You can specify several name and value pair arguments in any order as `Name1, Value1, ..., NameN, ValueN`.

### **Ts — Sample time**

`sample time of data (data.Ts) (default) | positive scalar | 0`

Sample time, specified as a positive scalar. For continuous-time models, use `Ts = 0`. For discrete-time models, specify `Ts` as a positive scalar whose value is equal to that of the data sample time.

### **Form — Type of canonical form**

`'free' (default) | 'modal' | 'companion' | 'canonical'`

Type of canonical form of `sys`, specified as one of the following values:

- `'modal'` — Obtain `sys` in modal form on page 1-919.
- `'companion'` — Obtain `sys` in companion form on page 1-920.
- `'free'` — All entries of the  $A$ ,  $B$ , and  $C$  matrices are estimated.
- `'canonical'` — Obtain `sys` in observability canonical form [1].

Use the `Form`, `Feedthrough`, and `DisturbanceModel` name-value pair arguments to modify the default behavior of the  $A$ ,  $B$ ,  $C$ ,  $D$ , and  $K$  matrices.

### **Feedthrough — Direct feedthrough from input to output**

`false(1, Nu) (default) | vector | scalar`

Logical specifying direct feedthrough from input to output, specified as a logical vector of length  $Nu$ , where  $Nu$  is the number of inputs.

If `Feedthrough` is specified as a logical scalar, this value is applied to all the inputs. If the model has no states, then `Feedthrough` is `true(1, Nu)`.

### **DisturbanceModel — Specify whether to estimate the $K$ matrix**

'estimate' (For time-domain data) (default) | 'none'

Specify estimation of the noise component ( $K$  matrix), specified as one of the following values:

- 'none' — Noise component is not estimated. The value of the  $K$  matrix, is fixed to zero value.
- 'estimate' — The  $K$  matrix is treated as a free parameter.

`DisturbanceModel` must be 'none' when using frequency domain data.

### **InputDelay — Input delays**

0 (default) | scalar | vector

Input delay for each input channel, specified as a numeric vector. For continuous-time systems, specify input delays in the time unit stored in the `TimeUnit` property. For discrete-time systems, specify input delays in integer multiples of the sample time  $T_s$ . For example, `InputDelay = 3` means a delay of three sampling periods.

For a system with  $Nu$  inputs, set `InputDelay` to an  $Nu$ -by-1 vector. Each entry of this vector is a numerical value that represents the input delay for the corresponding input channel.

You can also set `InputDelay` to a scalar value to apply the same delay to all channels.

## **Output Arguments**

### **sys**

Identified state-space model, returned as a `idss` model. This model is created using the specified model orders, delays, and estimation options.

Information about the estimation results and options used is stored in the `Report` property of the model. `Report` has the following fields:

<b>Report Field</b>	<b>Description</b>
Status	Summary of the model status, which indicates whether the model was created by construction or obtained by estimation.
Method	Estimation command used.
InitialState	<p>How initial states were handled during estimation, returned as one of the following values:</p> <ul style="list-style-type: none"> <li>• 'zero' — The initial state is set to zero.</li> <li>• 'estimate' — The initial state is treated as an independent estimation parameter.</li> </ul> <p>This field is especially useful when the <code>InitialState</code> option in the estimation option set is 'auto'.</p>
N4Weight	<p>Weighting scheme used for singular-value decomposition by the N4SID algorithm, returned as one of the following values:</p> <ul style="list-style-type: none"> <li>• 'MOESP' — Uses the MOESP algorithm.</li> <li>• 'CVA' — Uses the Canonical Variable Algorithm.</li> <li>• 'SSARX' — A subspace identification method that uses an ARX estimation based algorithm to compute the weighting.</li> </ul> <p>This option is especially useful when the <code>N4Weight</code> option in the estimation option set is 'auto'.</p>
N4Horizon	Forward and backward prediction horizons used by the N4SID algorithm, returned as a row vector with three elements — $[r \ sy \ su]$ , where $r$ is the maximum forward prediction horizon. $sy$ is the number of past outputs, and $su$ is the number of past inputs that are used for the predictions.

Report Field	Description																		
Fit	<p>Quantitative assessment of the estimation, returned as a structure. See “Loss Function and Model Quality Metrics” for more information on these quality metrics. The structure has the following fields:</p> <table border="1" data-bbox="387 456 1328 951"> <thead> <tr> <th data-bbox="387 456 516 499">Field</th> <th data-bbox="516 456 1328 499">Description</th> </tr> </thead> <tbody> <tr> <td data-bbox="387 499 516 609">FitPercent</td> <td data-bbox="516 499 1328 609">Normalized root mean squared error (NRMSE) measure of how well the response of the model fits the estimation data, expressed as a percentage.</td> </tr> <tr> <td data-bbox="387 609 516 652">LossFcn</td> <td data-bbox="516 609 1328 652">Value of the loss function when the estimation completes.</td> </tr> <tr> <td data-bbox="387 652 516 730">MSE</td> <td data-bbox="516 652 1328 730">Mean squared error (MSE) measure of how well the response of the model fits the estimation data.</td> </tr> <tr> <td data-bbox="387 730 516 774">FPE</td> <td data-bbox="516 730 1328 774">Final prediction error for the model.</td> </tr> <tr> <td data-bbox="387 774 516 817">AIC</td> <td data-bbox="516 774 1328 817">Raw Akaike Information Criteria (AIC) measure of model quality.</td> </tr> <tr> <td data-bbox="387 817 516 861">AICc</td> <td data-bbox="516 817 1328 861">Small sample-size corrected AIC.</td> </tr> <tr> <td data-bbox="387 861 516 904">nAIC</td> <td data-bbox="516 861 1328 904">Normalized AIC.</td> </tr> <tr> <td data-bbox="387 904 516 951">BIC</td> <td data-bbox="516 904 1328 951">Bayesian Information Criteria (BIC).</td> </tr> </tbody> </table>	Field	Description	FitPercent	Normalized root mean squared error (NRMSE) measure of how well the response of the model fits the estimation data, expressed as a percentage.	LossFcn	Value of the loss function when the estimation completes.	MSE	Mean squared error (MSE) measure of how well the response of the model fits the estimation data.	FPE	Final prediction error for the model.	AIC	Raw Akaike Information Criteria (AIC) measure of model quality.	AICc	Small sample-size corrected AIC.	nAIC	Normalized AIC.	BIC	Bayesian Information Criteria (BIC).
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nAIC	Normalized AIC.																		
BIC	Bayesian Information Criteria (BIC).																		
Parameters	Estimated values of model parameters.																		
OptionsUsed	Option set used for estimation. If no custom options were configured, this is a set of default options. See <code>n4sidOptions</code> for more information.																		
RandState	State of the random number stream at the start of estimation. Empty, [], if randomization was not used during estimation. For more information, see <code>rng</code> in the MATLAB documentation.																		

Report Field	Description	
DataUsed	Attributes of the data used for estimation, returned as a structure with the following fields:	
	Field	Description
	Name	Name of the data set.
	Type	Data type.
	Length	Number of data samples.
	Ts	Sample time.
	InterSample	Input intersample behavior, returned as one of the following values: <ul style="list-style-type: none"> <li>• 'zoh' — Zero-order hold maintains a piecewise-constant input signal between samples.</li> <li>• 'foh' — First-order hold maintains a piecewise-linear input signal between samples.</li> <li>• 'bl' — Band-limited behavior specifies that the continuous-time input signal has zero power above the Nyquist frequency.</li> </ul>
	InputOffset	Offset removed from time-domain input data during estimation. For nonlinear models, it is [ ].
OutputOffset	Offset removed from time-domain output data during estimation. For nonlinear models, it is [ ].	

For more information on using Report, see “Estimation Report”.

**x0**

Initial states computed during the estimator of sys.

If data contains multiple experiments, then x0 is an array with each column corresponding to an experiment.

## Examples

## Estimate State-Space Model and Specify Estimation Options

Load estimation data.

```
load iddata2 z2
```

Specify the estimation options.

```
opt = n4sidOptions('Focus','simulation','Display','on');
```

Estimate the model.

```
nx = 3;
sys = n4sid(z2,nx,opt);
```

sys is a third-order, state-space model.

## Estimate State-Space Model from Closed-Loop Data

Estimate a state-space model from closed-loop data using the subspace algorithm SSARX. This algorithm is better at capturing feedback effects than other weighting algorithms.

Generate closed-loop estimation data for a second-order system corrupted by white noise.

```
N = 1000;
K = 0.5;
rng('default');
w = randn(N,1);
z = zeros(N,1);
u = zeros(N,1);
y = zeros(N,1);
e = randn(N,1);
v = filter([1 0.5],[1 1.5 0.7],e);
for k = 3:N
    u(k-1) = -K*y(k-2) + w(k);
    u(k-1) = -K*y(k-1) + w(k);
    z(k) = 1.5*z(k-1) - 0.7*z(k-2) + u(k-1) + 0.5*u(k-2);
    y(k) = z(k) + 0.8*v(k);
end
dat = iddata(y, u, 1);
```

Specify the weighting scheme used by the N4SID algorithm. In one options set, specify the algorithm as CVA and in the other, specify as SSARX.

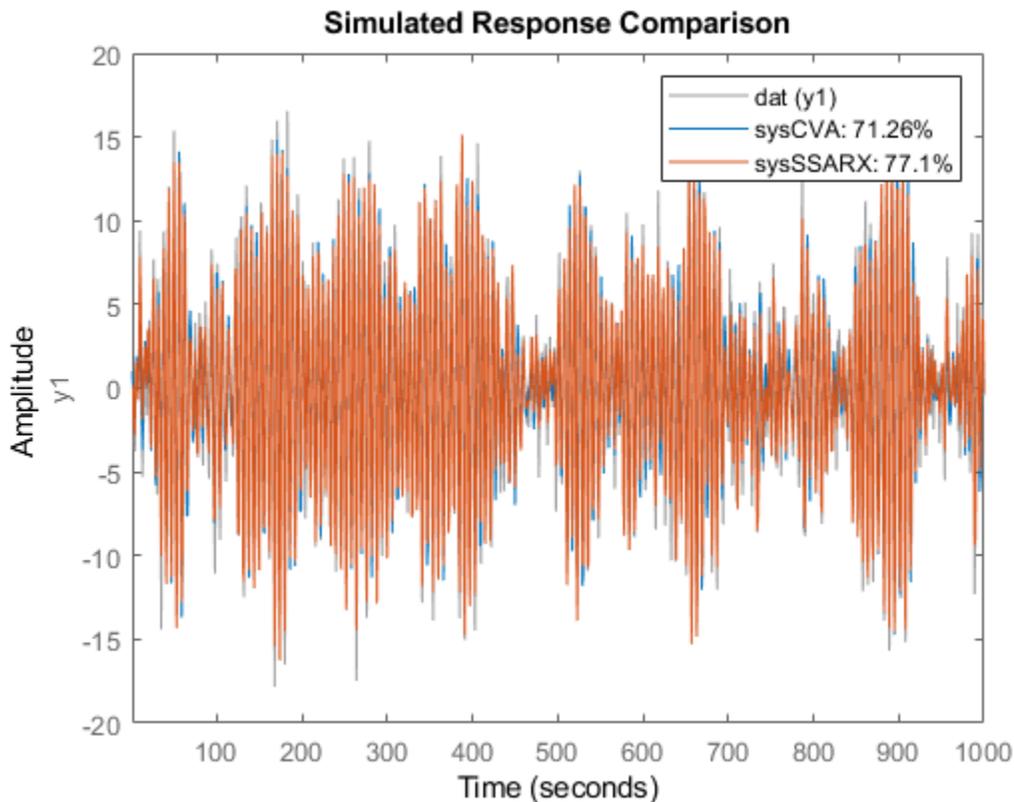
```
optCVA = n4sidOptions('N4weight','CVA');  
optSSARX = n4sidOptions('N4weight','SSARX');
```

Estimate state-space models using the options sets.

```
sysCVA = n4sid(dat,2,optCVA);  
sysSSARX = n4sid(dat,2,optSSARX);
```

Compare the fit of the two models with the estimation data.

```
compare(dat,sysCVA,sysSSARX);
```



From the plot, you see that the model estimated using the SSARX algorithm produces a better fit than the CVA algorithm.

## Estimate a Canonical-Form, Continuous-Time Model

Estimate a continuous-time, canonical-form model.

Load estimation data.

```
load iddata1 z1
```

Specify the estimation options.

```
opt = n4sidOptions('Focus','simulation','Display','on');
```

Estimate the model.

```
nx = 2;
sys = n4sid(z1,nx,'Ts',0,'Form','canonical',opt);
```

sys is a second-order, continuous-time, state-space model in the canonical form.

## Definitions

### Modal Form

In modal form,  $A$  is a block-diagonal matrix. The block size is typically 1-by-1 for real eigenvalues and 2-by-2 for complex eigenvalues. However, if there are repeated eigenvalues or clusters of nearby eigenvalues, the block size can be larger.

For example, for a system with eigenvalues  $(\lambda_1, \sigma \pm j\omega, \lambda_2)$ , the modal  $A$  matrix is of the form

$$\begin{bmatrix} \lambda_1 & 0 & 0 & 0 \\ 0 & \sigma & \omega & 0 \\ 0 & -\omega & \sigma & 0 \\ 0 & 0 & 0 & \lambda_2 \end{bmatrix}$$

## Companion Form

In the companion realization, the characteristic polynomial of the system appears explicitly in the right-most column of the  $A$  matrix.

For a system with characteristic polynomial

$$p(s) = s^n + \alpha_1 s^{n-1} + \dots + \alpha_{n-1} s + \alpha_n$$

the corresponding companion  $A$  matrix is

$$A = \begin{bmatrix} 0 & 0 & \dots & \dots & 0 & -\alpha_n \\ 1 & 0 & 0 & \dots & 0 & -\alpha_{n-1} \\ 0 & 1 & 0 & \dots & \vdots & \vdots \\ \vdots & 0 & \dots & \dots & \vdots & \vdots \\ 0 & \dots & \dots & 1 & 0 & -\alpha_2 \\ 0 & \dots & \dots & 0 & 1 & -\alpha_1 \end{bmatrix}$$

The companion transformation requires that the system be controllable from the first input. The companion form is poorly conditioned for most state-space computations; avoid using it when possible.

## References

- [1] Ljung, L. *System Identification: Theory for the User*, Appendix 4A, Second Edition, pp. 132-134. Upper Saddle River, NJ: Prentice Hall PTR, 1999.
- [2] van Overschee, P., and B. De Moor. *Subspace Identification of Linear Systems: Theory, Implementation, Applications*. Springer Publishing: 1996.
- [3] Verhaegen, M. "Identification of the deterministic part of MIMO state space models." *Automatica*, 1994, Vol. 30, pp. 61-74.
- [4] Larimore, W.E. "Canonical variate analysis in identification, filtering and adaptive control." *Proceedings of the 29th IEEE Conference on Decision and Control*, 1990, pp. 596-604.

## See Also

canon | iddata | idfrd | idgrey | idss | n4sidOptions | pem | polyest | procest |  
ssest | tfest

**Introduced before R2006a**

## n4sidOptions

Option set for n4sid

### Syntax

```
opt = n4sidOptions  
opt = n4sidOptions(Name,Value)
```

### Description

`opt = n4sidOptions` creates the default options set for `n4sid`.

`opt = n4sidOptions(Name,Value)` creates an option set with the options specified by one or more `Name,Value` pair arguments.

### Input Arguments

#### Name-Value Pair Arguments

Specify optional comma-separated pairs of `Name,Value` arguments. `Name` is the argument name and `Value` is the corresponding value. `Name` must appear inside quotes. You can specify several name and value pair arguments in any order as `Name1,Value1,...,NameN,ValueN`.

#### InitialState — Handling of initial states

```
'estimate' (default) | 'zero'
```

Handling of initial states during estimation, specified as one of the following values:

- 'zero' — The initial state is set to zero.
- 'estimate' — The initial state is treated as an independent estimation parameter.

### **N4Weight — Weighting scheme used for singular-value decomposition by the N4SID algorithm**

'auto' (default) | 'MOESP' | 'CVA' | 'SSARX'

Weighting scheme used for singular-value decomposition by the N4SID algorithm, specified as one of the following values:

- 'MOESP' — Uses the MOESP algorithm by Verhaegen [2].
- 'CVA' — Uses the Canonical Variable Algorithm by Larimore [1].

Estimation using frequency-domain data always uses 'CVA'.

- 'SSARX' — A subspace identification method that uses an ARX estimation based algorithm to compute the weighting.

Specifying this option allows unbiased estimates when using data that is collected in closed-loop operation. For more information about the algorithm, see [4].

- 'auto' — The estimating function chooses between the MOESP, CVA and SSARX algorithms.

### **N4Horizon — Forward- and backward-prediction horizons used by the N4SID algorithm**

'auto' (default) | vector [r sy su] | k-by-3 matrix

Forward- and backward-prediction horizons used by the N4SID algorithm, specified as one of the following values:

- A row vector with three elements — [r sy su], where r is the maximum forward prediction horizon, using up to r step-ahead predictors. sy is the number of past outputs, and su is the number of past inputs that are used for the predictions. See pages 209 and 210 in [3] for more information. These numbers can have a substantial influence on the quality of the resulting model, and there are no simple rules for choosing them. Making 'N4Horizon' a k-by-3 matrix means that each row of 'N4Horizon' is tried, and the value that gives the best (prediction) fit to data is selected. k is the number of guesses of [r sy su] combinations. If you specify N4Horizon as a single column, r = sy = su is used.
- 'auto' — The software uses an Akaike Information Criterion (AIC) for the selection of sy and su.

### **Focus — Error to be minimized**

'prediction' (default) | 'simulation'

Error to be minimized in the loss function during estimation, specified as the comma-separated pair consisting of 'Focus' and one of the following values:

- 'prediction' — The one-step ahead prediction error between measured and predicted outputs is minimized during estimation. As a result, the estimation focuses on producing a good predictor model.
- 'simulation' — The simulation error between measured and simulated outputs is minimized during estimation. As a result, the estimation focuses on making a good fit for simulation of model response with the current inputs.

The Focus option can be interpreted as a weighting filter in the loss function. For more information, see “Loss Function and Model Quality Metrics”.

### **WeightingFilter — Weighting prefilter**

[ ] (default) | vector | matrix | cell array | linear system

Weighting prefilter applied to the loss function to be minimized during estimation. To understand the effect of `WeightingFilter` on the loss function, see “Loss Function and Model Quality Metrics”.

Specify `WeightingFilter` as one of the following values:

- [ ] — No weighting prefilter is used.
- Passbands — Specify a row vector or matrix containing frequency values that define desired passbands. You select a frequency band where the fit between estimated model and estimation data is optimized. For example, `[wl, wh]` where `wl` and `wh` represent lower and upper limits of a passband. For a matrix with several rows defining frequency passbands, `[w1l, w1h; w2l, w2h; w3l, w3h; ...]`, the estimation algorithm uses the union of the frequency ranges to define the estimation passband.

Passbands are expressed in `rad/TimeUnit` for time-domain data and in `FrequencyUnit` for frequency-domain data, where `TimeUnit` and `FrequencyUnit` are the time and frequency units of the estimation data.

- SISO filter — Specify a single-input-single-output (SISO) linear filter in one of the following ways:
  - A SISO LTI model
  - `{A, B, C, D}` format, which specifies the state-space matrices of a filter with the same sample time as estimation data.

- `{numerator,denominator}` format, which specifies the numerator and denominator of the filter as a transfer function with same sample time as estimation data.

This option calculates the weighting function as a product of the filter and the input spectrum to estimate the transfer function.

- **Weighting vector** — Applicable for frequency-domain data only. Specify a column vector of weights. This vector must have the same length as the frequency vector of the data set, `Data.Frequency`. Each input and output response in the data is multiplied by the corresponding weight at that frequency.

### **EnforceStability** — Control whether to enforce stability of model

`false` (default) | `true`

Control whether to enforce stability of estimated model, specified as the comma-separated pair consisting of 'EnforceStability' and either `true` or `false`.

Data Types: `logical`

### **EstimateCovariance** — Control whether to generate parameter covariance data

`true` (default) | `false`

Controls whether parameter covariance data is generated, specified as `true` or `false`.

If `EstimateCovariance` is `true`, then use `getcov` to fetch the covariance matrix from the estimated model.

### **Display** — Specify whether to display the estimation progress

`'off'` (default) | `'on'`

Specify whether to display the estimation progress, specified as one of the following values:

- `'on'` — Information on model structure and estimation results are displayed in a progress-viewer window.
- `'off'` — No progress or results information is displayed.

### **InputOffset** — Removal of offset from time-domain input data during estimation

`[]` (default) | vector of positive integers | matrix

Removal of offset from time-domain input data during estimation, specified as the comma-separated pair consisting of 'InputOffset' and one of the following:

- A column vector of positive integers of length  $Nu$ , where  $Nu$  is the number of inputs.
- [] — Indicates no offset.
- $Nu$ -by- $Ne$  matrix — For multi-experiment data, specify InputOffset as an  $Nu$ -by- $Ne$  matrix.  $Nu$  is the number of inputs, and  $Ne$  is the number of experiments.

Each entry specified by InputOffset is subtracted from the corresponding input data.

### **OutputOffset — Removal of offset from time-domain output data during estimation**

[] (default) | vector | matrix

Removal of offset from time-domain output data during estimation, specified as the comma-separated pair consisting of 'OutputOffset' and one of the following:

- A column vector of length  $Ny$ , where  $Ny$  is the number of outputs.
- [] — Indicates no offset.
- $Ny$ -by- $Ne$  matrix — For multi-experiment data, specify OutputOffset as a  $Ny$ -by- $Ne$  matrix.  $Ny$  is the number of outputs, and  $Ne$  is the number of experiments.

Each entry specified by OutputOffset is subtracted from the corresponding output data.

### **OutputWeight — Weighting of prediction errors in multi-output estimations**

[] (default) | 'noise' | positive semidefinite symmetric matrix

Weighting of prediction errors in multi-output estimations, specified as one of the following values:

- 'noise' — Minimize  $\det(E^*E/N)$ , where  $E$  represents the prediction error and  $N$  is the number of data samples. This choice is optimal in a statistical sense and leads to the maximum likelihood estimates in case no data is available about the variance of the noise. This option uses the inverse of the estimated noise variance as the weighting function.
- Positive semidefinite symmetric matrix ( $W$ ) — Minimize the trace of the weighted prediction error matrix  $\text{trace}(E^*EW/N)$  where:
  - $E$  is the matrix of prediction errors, with one column for each output.  $W$  is the positive semidefinite symmetric matrix of size equal to the number of outputs. Use

W to specify the relative importance of outputs in multiple-output models, or the reliability of corresponding data.

- N is the number of data samples.
- [] — The software chooses between the 'noise' or using the identity matrix for W.

This option is relevant only for multi-output models.

### **Advanced — Additional advanced options**

structure

Additional advanced options, specified as a structure with the field `MaxSize`. `MaxSize` specifies the maximum number of elements in a segment when input-output data is split into segments.

`MaxSize` must be a positive integer.

**Default:** 250000

## **Output Arguments**

### **opt — Option set for n4sid**

n4sidOptions option set

Option set for `n4sid`, returned as an `n4sidOptions` option set.

## **Examples**

### **Create Default Options Set for State-Space Estimation Using Subspace Method**

```
opt = n4sidOptions;
```

### **Specify Options for State-Space Estimation Using Subspace Method**

Create an options set for `n4sid` using the 'zero' option to initialize the state. Set the `Display` to 'on'.

```
opt = n4sidOptions('InitialState', 'zero', 'Display', 'on');
```

Alternatively, use dot notation to set the values of `opt`.

```
opt = n4sidOptions;  
opt.InitialState = 'zero';  
opt.Display = 'on';
```

## References

- [1] Larimore, W.E. "Canonical variate analysis in identification, filtering and adaptive control." *Proceedings of the 29th IEEE Conference on Decision and Control*, pp. 596-604, 1990.
- [2] Verhaegen, M. "Identification of the deterministic part of MIMO state space models." *Automatica*, Vol. 30, 1994, pp. 61-74.
- [3] Ljung, L. *System Identification: Theory for the User*. Upper Saddle River, NJ: Prentice-Hall PTR, 1999.
- [4] Jansson, M. "Subspace identification and ARX modeling." *13th IFAC Symposium on System Identification*, Rotterdam, The Netherlands, 2003.

## See Also

`idfilt` | `idpar` | `n4sid`

## Topics

"Loss Function and Model Quality Metrics"

**Introduced in R2012a**

# ndims

Query number of dimensions of dynamic system model or model array

## Syntax

```
n = ndims(sys)
```

## Description

`n = ndims(sys)` is the number of dimensions of a dynamic system model or a model array `sys`. A single model has two dimensions (one for outputs, and one for inputs). A model array has  $2 + p$  dimensions, where  $p \geq 2$  is the number of array dimensions. For example, a 2-by-3-by-4 array of models has  $2 + 3 = 5$  dimensions.

```
ndims(sys) = length(size(sys))
```

## Examples

### Determine Dimensions of Model Array

Create a 3-by-1 array of random state-space models, each with 4 states, 1 input, and 1 output.

```
sys = rss(4,1,1,3);
```

Compute the number of dimensions of the model array.

```
ndims(sys)
```

```
ans = 4
```

The number of dimensions is  $2+p$ , where  $p$  is the number of array dimensions. In this example,  $p$  is 2 because `sys` is 3-by-1.

## **See Also**

size

**Introduced in R2012a**

# neuralnet

Class representing neural network nonlinearity estimator for nonlinear ARX models

## Syntax

```
net_estimator = neuralnet(Network)
```

## Description

`neuralnet` is the class that encapsulates the neural network nonlinearity estimator. A `neuralnet` object lets you use networks, created using Deep Learning Toolbox software, in nonlinear ARX models.

The neural network nonlinearity estimator defines a nonlinear function  $y = F(x)$ , where  $F$  is a multilayer feed-forward (static) neural network, as defined in the Deep Learning Toolbox software.  $y$  is a scalar and  $x$  is an  $m$ -dimensional row vector.

You create multi-layer feed-forward neural networks using Deep Learning Toolbox commands such as `feedforwardnet`, `cascadeforwardnet` and `linearlayer`. When you create the network:

- Designate the input and output sizes to be unknown by leaving them at the default value of zero (recommended method). When estimating a nonlinear ARX model using the `nlarx` command, the software automatically determines the input-output sizes of the network.
- Initialize the sizes manually by setting input and output ranges to  $m$ -by-2 and 1-by-2 matrices, respectively, where  $m$  is the number of nonlinear ARX model regressors and the range values are minimum and maximum values of regressors and output data, respectively.

See “Examples” on page 1-0 for more information.

Use `evaluate(net_estimator, x)` to compute the value of the function defined by the `neuralnet` object `net_estimator` at input value  $x$ . When used for nonlinear ARX model

estimation,  $x$  represents the model regressors for the output for which the `neuralnet` object is assigned as the nonlinearity estimator.

You cannot use `neuralnet` when the `Focus` option in `nlarxOptions` is `'simulation'` because this nonlinearity estimator is considered to be nondifferentiable for estimation. Minimization of simulation error requires differentiable nonlinear functions.

## Construction

`net_estimator = neuralnet(Network)` creates a neural network nonlinearity estimator based on the feed-forward (static) network object `Network` created using Deep Learning Toolbox commands `feedforwardnet`, `cascadeforwardnet`, and `linearlayer`. `Network` must represent a static mapping between the inputs and output without I/O delays or feedback. The number of outputs of the network, if assigned, must be one. For a multiple-output nonlinear ARX models, create a separate `neuralnet` object for each output—that is, each estimator must represent a single-output network object.

## Properties

Network	Neural network object, typically created using the Deep Learning Toolbox commands <code>feedforwardnet</code> , <code>cascadeforwardnet</code> , and <code>linearlayer</code> .
---------	---

After creating the object, you can use `get` or dot notation to access the object property values. For example:

```
% List Network property value
get(n)
n.Network
```

You can also use the `set` function to set the value of particular properties. For example:

```
set(d, 'Network', net_obj)
```

The first argument to `set` must be the name of a MATLAB variable.

## Examples

## Create a Nonlinearity Estimator Using Feed-Forward Neural Network

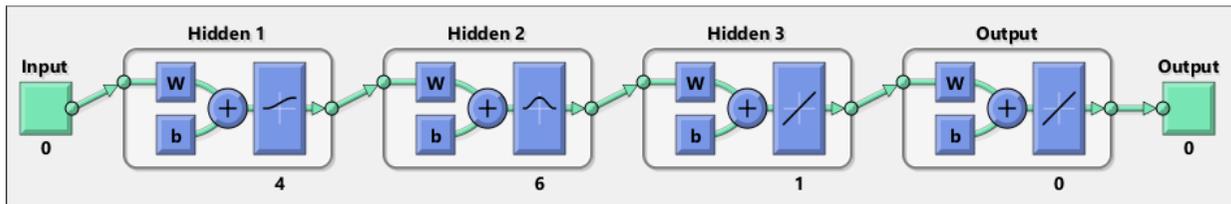
Create a neural network nonlinearity estimator using a feed-forward neural network with three hidden layers; transfer functions of types `logsig`, `radbas`, and `purelin`; and unknown input and output sizes.

Create a neural network.

```
net = feedforwardnet([4 6 1]);
net.layers{1}.transferFcn = 'logsig';
net.layers{2}.transferFcn = 'radbas';
net.layers{3}.transferFcn = 'purelin';
```

View the network diagram.

```
view(net)
```



Create a neural network estimator.

```
net_estimator = neuralnet(net);
```

## Estimate Nonlinear ARX Model Using a Neural Network Nonlinearity Estimator

Create a single-layer, cascade-forward network with unknown input and output sizes and use this network for nonlinear ARX model estimation.

Create a cascade-forward neural network with 20 neurons and unknown input-output sizes.

```
net = cascadeforwardnet(20);
```

Create a neural network nonlinearity estimator.

```
net_estimator = neuralnet(net);
```

Load estimation data.

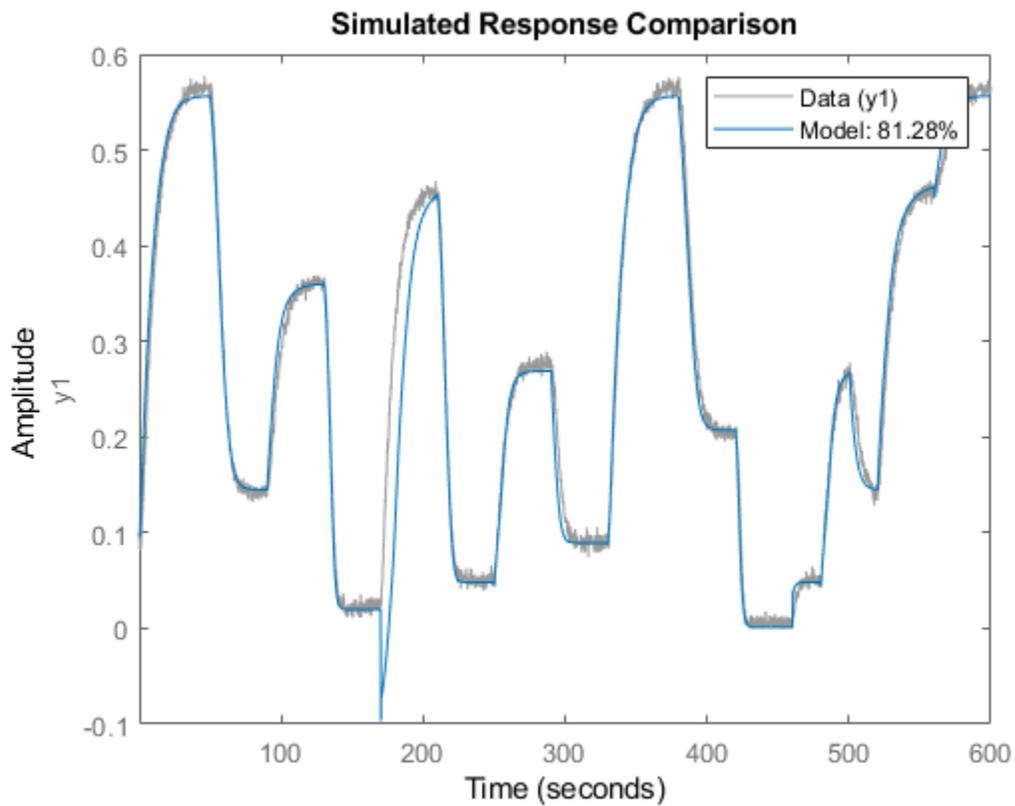
```
load twotankdata  
Data = iddata(y,u,0.2);
```

Estimate nonlinear ARX model.

```
Model = nlarx(Data,[2 2 1],net_estimator);
```

Compare model response to measured output signal.

```
compare(Data,Model)
```



## Initialize Input-Output Sizes of Neural Network Nonlinearity Estimator

Initialize the input-output sizes of a two-layer feed-forward neural network based on estimation data, and use this network for nonlinear ARX estimation.

Load estimation data.

```
load iddata7 z7
z7 = z7(1:200);
```

Create a template nonlinear ARX model with no nonlinearity.

```
model = idnlarx([4 4 4 1 1],[]);
```

This model has six regressors and is used to define the regressors. The range of regressor values for input-output data in `z7` is then used to set the input ranges in the neural network object, as shown in the next steps.

Obtain the model regressor values.

```
R = getreg(model, 'all', z7);
```

Create a two-layer, feed-forward neural network and initialize the network input and output dimensions to 2 and 1, respectively. Use 5 neurons for first layer and 7 for second layer.

```
net = feedforwardnet([5 7]);
```

Determine input range.

```
InputRange = [min(R);max(R)].';
```

Initialize input dimensions of estimator.

```
net.inputs{1}.range = InputRange;
```

Determine output range.

```
OutputRange = [min(z7.OutputData),max(z7.OutputData)];
```

Initialize output dimensions of estimator.

```
net.outputs{net.outputConnect}.range = OutputRange;
```

Create a neural network nonlinearity estimator.

```
net_estimator = neuralnet(net);
```

Specify the nonlinearity estimator in the model.

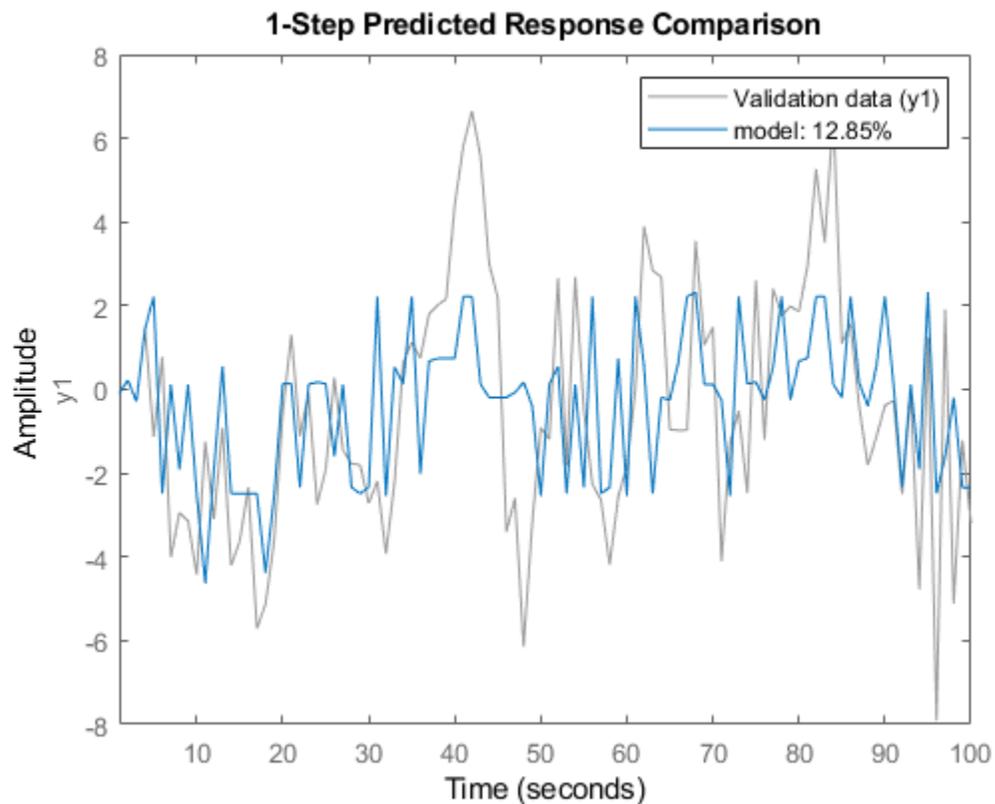
```
model.Nonlinearity = net_estimator;
```

Estimate the parameters of the network to minimize the prediction error between data and model. Estimate model.

```
model = nlarx(z7,model);
```

Compare model's predicted response to measured output signal.

```
compare(z7(1:100),model,1)
```



## Algorithms

The `nlarx` command uses the `train` method of the `network` object, defined in the Deep Learning Toolbox software, to compute the network parameter values.

## See Also

`cascadeforwardnet` | `customnet` | `feedforwardnet` | `linearlayer` | `nlarx` | `sigmoidnet` | `treepartition` | `wavenet`

## Topics

“Identifying Nonlinear ARX Models”

**Introduced in R2007a**

## nkshift

Shift data sequences

### Syntax

```
Datas = nkshift(Data,nk)
```

### Description

Data contains input-output data in the `iddata` format.

`nk` is a row vector with the same length as the number of input channels in `Data`.

`Datas` is an `iddata` object where the input channels in `Data` have been shifted according to `nk`. A positive value of `nk(ku)` means that input channel number `ku` is delayed `nk(ku)` samples.

`nkshift` supports both frequency- and time-domain data. For frequency-domain data it multiplies with  $e^{ink\omega T}$  to obtain the same effect as shifting in the time domain. For continuous-time frequency-domain data ( $T_s = 0$ ), `nk` should be interpreted as the shift in seconds.

`nkshift` lives in symbiosis with the `InputDelay` property of linear identified models:

```
m1 = ssest(dat,4,'InputDelay',nk)
```

is related to

```
m2 = ssest(nkshift(dat,nk),4);
```

such that `m1` and `m2` are the same models, but `m1` stores the delay information and uses this information when computing the frequency response, for example. When using `m2`, the delay value must be accounted for separately when computing time and frequency responses.

## **See Also**

absorbDelay | delayest | idpoly | idss

**Introduced before R2006a**

## **nlarx**

Estimate parameters of nonlinear ARX model

### **Syntax**

```
sys = nlarx(Data,Orders)
sys = nlarx(Data,Orders,Nonlinearity)
sys = nlarx(Data,Orders,Nonlinearity,Name,Value)

sys = nlarx(Data,LinModel)
sys = nlarx(Data,LinModel,Nonlinearity)
sys = nlarx(Data,LinModel,Nonlinearity,Name,Value)

sys = nlarx(Data,sys0)

sys = nlarx( ____,Options)
```

### **Description**

`sys = nlarx(Data,Orders)` estimates a nonlinear ARX model to fit the given estimation data using the specified orders and a default wavelet network nonlinearity estimator.

`sys = nlarx(Data,Orders,Nonlinearity)` specifies the nonlinearity to use for model estimation.

`sys = nlarx(Data,Orders,Nonlinearity,Name,Value)` specifies additional attributes of the estimated model using one or more `Name,Value` pair arguments. These attributes include the nonlinear and custom regressor structure, and the data properties on page 1-621 of the `idnlarx` model.

`sys = nlarx(Data,LinModel)` uses a linear ARX model, `LinModel`, to specify the model orders and the initial values of the linear coefficients of the model.

`sys = nlarx(Data,LinModel,Nonlinearity)` specifies the nonlinearity to use for model estimation.

`sys = nlarx(Data, LinModel, Nonlinearity, Name, Value)` specifies additional attributes of the `idnlarx` model structure using one or more `Name, Value` pair arguments.

`sys = nlarx(Data, sys0)` refines the parameters of the nonlinear ARX model, `sys0`.

Use this syntax to:

- Update the parameters of a previously estimated model to improve the fit to the estimation data. In this case, the estimation algorithm uses the parameters of `sys0` as initial guesses.
- Estimate the parameters of a model previously created using the `idnlarx` constructor. Prior to estimation, you can configure the model properties using dot notation.

`sys = nlarx( ___, Options)` specifies additional configuration options for the model estimation. Use `Options` with any of the previous syntaxes.

## Examples

### Estimate Nonlinear ARX Model with Default Settings

Load the estimation data.

```
load twotankdata;
```

Create an `iddata` object from the estimation data with a sample time of 0.2 min.

```
Ts = 0.2;  
z = iddata(y,u,Ts);
```

Estimate the nonlinear ARX model.

```
sys = nlarx(z,[4 4 1]);
```

### Estimate Nonlinear ARX Model from Time Series Data

Create time and data arrays.

```
dt = 0.01;  
t = 0:dt:10;  
y = 10*sin(2*pi*t)+rand(size(t));
```

Create an `iddata` object with no input signal specified.

```
z = iddata(y', [], dt);
```

Estimate the nonlinear ARX model.

```
sys = nlarx(z, 2);
```

### **Estimate Nonlinear ARX Model with Specific Nonlinearity**

Load the estimation data.

```
load twotankdata;
```

Create an `iddata` object from the estimation data.

```
z = iddata(y, u, 0.2);
```

Create a wavelet network nonlinearity estimator with 5 units.

```
NL = wavenet('NumberOfUnits', 5);
```

Estimate the nonlinear ARX model.

```
sys = nlarx(z, [4 4 1], NL);
```

### **Estimate Nonlinear ARX Model Using Custom Network Nonlinearity**

Generating a custom network nonlinearity requires the definition of a user-defined unit function.

Define the unit function and save it as `gaussunit.m`.

```
% Copyright 2015 The MathWorks, Inc.
```

```
function [f, g, a] = gaussunit(x)
f = exp(-x.*x);
if nargin>1
    g = -2*x.*f;
    a = 0.2;
end
```

Create a custom network nonlinearity using the `gaussunit` function.

```
H = @gaussunit;
CNet = customnet(H);
```

Load the estimation data.

```
load iddata1;
```

Estimate a nonlinear ARX model using the custom network.

```
sys = nlarx(z1,[1 2 1],CNet);
```

### Estimate MIMO Nonlinear ARX Model

Load the estimation data.

```
load motorizedcamera;
```

Create an `iddata` object.

```
z = iddata(y,u,0.02,'Name','Motorized Camera','TimeUnit','s');
```

`z` is an `iddata` object with 6 inputs and 2 outputs.

Specify the model orders.

```
Orders = [ones(2,2),2*ones(2,6),ones(2,6)];
```

Specify different nonlinearity estimators for each output channel.

```
NL = [wavenet('NumberOfUnits',2),linear];
```

Estimate the nonlinear ARX model.

```
sys = nlarx(z,Orders,NL);
```

### Estimate MIMO Nonlinear ARX Model with Same Nonlinearity for All Outputs

Load the estimation data and create an `iddata` object.

```
load motorizedcamera;  
z = iddata(y,u,0.02);
```

Specify the model orders.

```
Orders = [ones(2,2),2*ones(2,6),ones(2,6)];
```

Estimate a nonlinear ARX model using a `sigmoidnet` nonlinearity with 4 units for all output channels.

```
m = nlarx(z,Orders,sigmoidnet('numberOfUnits',4));
```

### Estimate Nonlinear ARX Model with Custom Regressors

Load the estimation data.

```
load iddata1;
```

Create a cell array with two custom regressors.

```
C = {'y1(t-1)^2', 'y1(t-2)*u1(t-3)'};
```

Estimate a nonlinear ARX model with custom regressors and no standard regressors.

```
sys = nlarx(z1,[0 0 0], 'linear', 'CustomRegressors',C);
```

### Estimate Nonlinear ARX Model with Custom Regressor Objects

Load the estimation data.

```
load iddata1;
```

Define a custom regressor object for  $y1(t-1)^2$ .

```
C1 = customreg(@(x)x^2,{'y1'},[1]);
```

Define a custom regressor object for  $y1(t-2)*u1(t-3)$ .

```
C2 = customreg(@(x,y)x*y,{'y1','u1'},[2 3]);
```

Create a custom regressor object array.

```
C = [C1,C2];
```

Estimate a nonlinear ARX model with custom regressors.

```
sys = nlarx(z1,[0 0 0], 'linear', 'CustomRegressors',C);
```

List the model regressors.

```
getreg(sys);
```

Regressors:

```
  y1(t-1)^2
  y1(t-2)*u1(t-3)
```

### Estimate Nonlinear ARX Model Searching for Optimum Nonlinear Regressors

Load the estimation data.

```
load iddata1;
```

Estimate a Nonlinear ARX model using the 'search' option.

```
sys = nlarx(z1,[4 4 1], 'sigmoidnet', 'NonlinearRegressors', 'search');
```

List the model nonlinear regressor indices.

```
sys.NonlinearRegressors
```

```
ans = 1×4
```

```
     3     5     6     7
```

List all of the model regressors.

```
getreg(sys)
```

```
Regressors:
```

```
  y1(t-1)
```

```
  y1(t-2)
```

```
  y1(t-3)
```

```
  y1(t-4)
```

```
  u1(t-1)
```

```
  u1(t-2)
```

```
  u1(t-3)
```

```
  u1(t-4)
```

The optimum set of nonlinear regressors for this model includes  $y1(t-3)$ ,  $u1(t-1)$ ,  $u1(t-2)$ , and  $u1(t-3)$ .

### Estimate Nonlinear ARX Model with No Linear Term in Nonlinearity Estimator

Load the estimation data.

```
load iddata1;
```

Create a sigmoid network nonlinearity estimator with no linear term.

```
SNL = sigmoidnet('LinearTerm','off');
```

Estimate the nonlinear ARX model.

```
sys = nlarx(z1,[2 2 1],SNL);
```

### Specify Nonlinear ARX Orders and Linear Parameters Using Linear ARX Model

Load the estimation data.

```
load throttledata;
```

Detrend the data.

```
Tr = getTrend(ThrottleData);
```

```
Tr.OutputOffset = 15;
```

```
DetrendedData = detrend(ThrottleData,Tr);
```

Estimate the linear ARX model.

```
LinearModel = arx(DetrendedData,[2 1 1]);
```

Estimate the nonlinear ARX model using the linear model. The model orders, delays, and linear parameters of `NonlinearModel` are derived from `LinearModel`.

```
NonlinearModel = nlarx(ThrottleData,LinearModel);
```

### **Estimate Nonlinear ARX Model Using Constructed `idnlarx` Object**

Load the estimation data.

```
load iddata1;
```

Create an `idnlarx` model.

```
sys = idnlarx([2 2 1]);
```

Configure the model using dot notation to set the following parameters:

- Use a sigmoid network nonlinearity
- Search for an optimum nonlinear regressor subset

```
sys.Nonlinearity = 'sigmoidnet';  
sys.NonlinearRegressors = 'search';
```

Estimate a nonlinear ARX model with the structure and properties specified in the `idnlarx` object.

```
sys = nlarx(z1,sys);
```

### **Estimate Nonlinear ARX Model and Avoid Local Minima**

If an estimation stops at a local minimum, you can perturb the model using `init` and reestimate the model.

Load the estimation data.

```
load iddata1;
```

Estimate the initial nonlinear model using specific nonlinear regressors.

```
sys1 = nlarx(z1,[4 2 1], 'sigmoidnet', 'NonlinearRegressors', [1:3]);
```

Randomly perturb the model parameters to avoid local minima.

```
sys2 = init(sys1);
```

Estimate the new nonlinear model with the perturbed values.

```
sys2 = nlarx(z1,sys2);
```

### **Estimate Nonlinear ARX Model Using Specific Options**

Load the estimation data.

```
load twotankdata;
```

Create an `iddata` object from the estimation data.

```
z = iddata(y,u,0.2);
```

Create an `nlarxOptions` option set specifying a simulation error minimization objective and a maximum of 50 estimation iterations.

```
opt = nlarxOptions;  
opt.Focus = 'simulation';  
opt.SearchOptions.MaxIterations = 50;
```

Estimate the nonlinear ARX model.

```
sys = nlarx(z,[4 4 1], 'sigmoidnet', opt);
```

### **Estimate Regularized Nonlinear ARX Model with Large Number of Units**

Load the regularization example data.

```
load regularizationExampleData.mat nldata;
```

Create a `sigmoidnet` nonlinearity with 30 units, and specify the model orders.

```
NL = sigmoidnet('NumberOfUnits',30);  
Orders = [1 2 1];
```

Create an estimation option set and set the estimation search method to `lm`.

```
opt = nlarxOptions('SearchMethod','lm');
```

Estimate an unregularized model.

```
sys = nlarx(nldata,Orders,NL,opt);
```

Configure the regularization Lambda parameter.

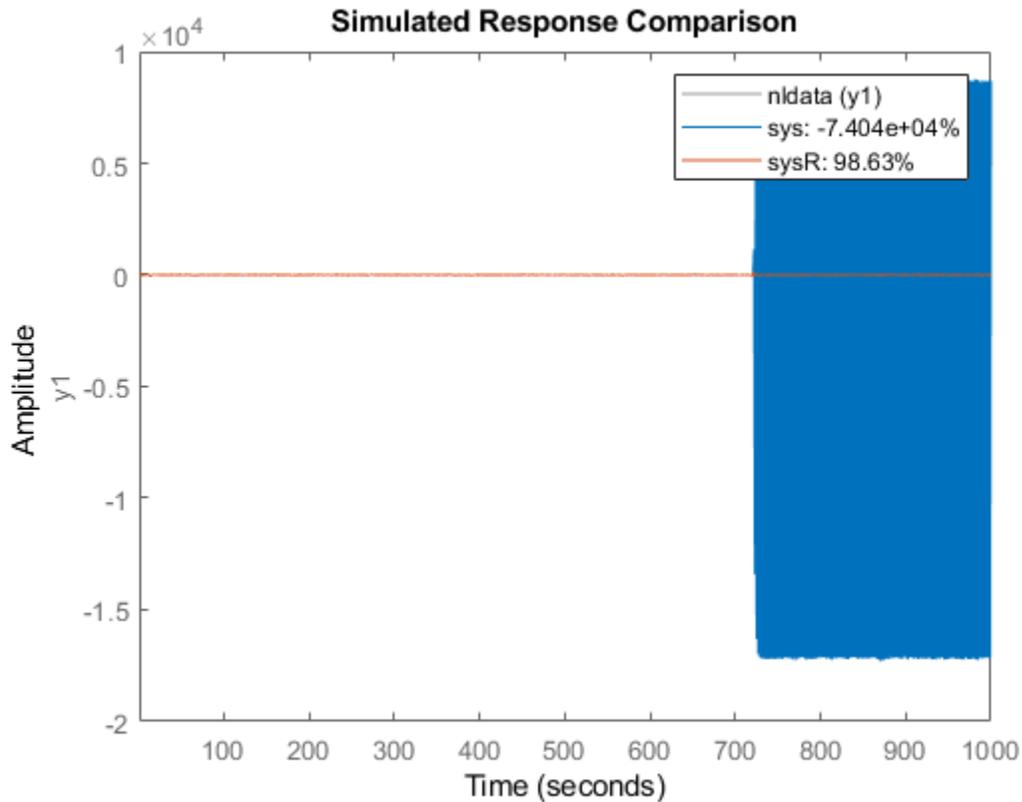
```
opt.Regularization.Lambda = 1e-8;
```

Estimate a regularized model.

```
sysR = nlarx(nldata,Orders,NL,opt);
```

Compare the two models.

```
compare(nldata,sys,sysR)
```



The large negative fit result for the unregularized model indicates a poor fit to the data. Estimating a regularized model produces a significantly better result.

## Input Arguments

### Data — Time-domain estimation data

`iddata` object

Time-domain estimation data, specified as an `iddata` object. Data can have one or more output channels and zero or more input channels. Data must be uniformly sampled and cannot contain missing (NaN) samples.

## Orders — Model orders and delays

1-by-3 vector of positive integers | 1-by-3 vector of matrices

Model orders and delays for defining the regressor configuration, specified as a 1-by-3 vector, `[na nb nk]`.

For a model with  $n_y$  output channels and  $n_u$  input channels:

- `na` is an  $n_y$ -by- $n_y$  matrix, where `na(i, j)` specifies the number of regressors from the  $j$ th output used to predict the  $i$ th output.
- `nb` is an  $n_y$ -by- $n_u$  matrix, where `nb(i, j)` specifies the number of regressors from the  $j$ th input used to predict the  $i$ th output.
- `nk` is an  $n_y$ -by- $n_u$  matrix, where `nk(i, j)` specifies the lag in the  $j$ th input used to predict the  $i$ th output.

```
na = [1 2; 2 3]
nb = [1 2 3; 2 3 1];
nk = [2 0 3; 1 0 5];
```

The estimation data for this system has three inputs (`u1`, `u2`, `u3`) and two outputs (`y1`, `y2`). Consider the regressors used to predict output, `y2(t)`:

- Since `na(2, :)` is `[2 3]`, the contributing regressors from the outputs are:
  - `y1(t-1)` and `y1(t-2)`
  - `y2(t-1)`, `y2(t-2)`, and `y2(t-3)`
- Since `nb(2, :)` is `[2 3 1]` and `nk(2, :)` is `[1 0 5]`, the contributing regressors from the inputs are:
  - `u1(t-1)` and `u1(t-2)`
  - `u2(t)`, `u2(t-1)`, and `u2(t-2)`
  - `u3(t-5)`

---

**Note** The minimum lag for regressors based on output variables is always 1, while the minimum lag for regressors based on input variables is dictated by `nk`. Use `getreg` to view the complete set of regressors used by the nonlinear ARX model.

---

**Nonlinearity — Nonlinearity estimator**

'wavenet' (default) | 'sigmoidnet' | 'treepartition' | 'linear' | nonlinearity estimator object | array of nonlinearity estimator objects

Nonlinearity estimator, specified as one of the following:

'wavenet' or wavenet object	Wavelet network
'sigmoidnet' or sigmoidnet object	Sigmoid network
'treepartition' or treepartition object	Binary-tree
'linear' or [] or linear object	Linear function
neuralnet object	Neural network — Requires Deep Learning Toolbox.
customnet object	Custom network — Similar to <code>sigmoidnet</code> , but with a user-defined replacement for the sigmoid function.

For more information, see “Available Nonlinearity Estimators for Nonlinear ARX Models”.

Specifying a character vector, for example 'sigmoidnet', creates a nonlinearity estimator object with default settings. Alternatively, you can specify nonlinearity estimator settings in two ways:

- Use the associated nonlinearity estimator function with Name-Value pair arguments.

```
NL = sigmoidnet('NumberOfUnits',10);
```

- Create and modify a default nonlinearity estimator object.

```
NL = sigmoidnet;  
NL.NumberOfUnits = 10;
```

For `ny` output channels, you can specify nonlinear estimators individually for each channel by setting `Nonlinearity` to an `ny`-by-1 array of nonlinearity estimator objects. To specify the same nonlinearity for all outputs, specify `Nonlinearity` as a character vector or a single nonlinearity estimator object.

Example: 'sigmoidnet' specifies a sigmoid network nonlinearity with a default configuration.

Example: `treepartition('NumberOfUnits',5)` specifies a binary-tree nonlinearity with 5 terms in the binary tree expansion.

Example: `[wavenet('NumberOfUnits',10);sigmoidnet]` specifies different nonlinearity estimators for two output channels.

### **LinModel — Discrete time input-output polynomial model of ARX structure**

`idpoly` model

Discrete time input-output polynomial model of ARX structure, specified as an `idpoly` model. Create this object using the `idpoly` constructor or estimate it using the `arx` command.

### **sys0 — Nonlinear ARX model**

`idnlarx` model

Nonlinear ARX model, specified as an `idnlarx` model. `sys0` can be:

- A model previously estimated using `nlarx`. The estimation algorithm uses the parameters of `sys0` as initial guesses. In this case, use `init` to slightly perturb the model properties to avoid being trapped in local minima.

```
sys = init(sys);
sys = nlarx(data,sys);
```

- A model previously created using `idnlarx` and with properties set using dot notation. Use this method to avoid complicated Name-Value pair syntax when configuring multiple model properties. For example, use

```
sys1 = idnlarx([4 3 1]);
sys1.Nonlinearity = 'treepartition';
sys1.CustomRegressors = {'sin(u1(t-1))'};
sys1.NonlinearRegressors = 'search';
sys2 = nlarx(data,sys1);
```

in place of the equivalent

```
sys2 = nlarx(data,[4,3,1],'treepartition','CustomRegressors',...
    {'sin(u1(t-1))'},'NonlinearRegressors','search');
```

### **Options — Estimation options**

`nlarxOptions` option set

Estimation options for nonlinear ARX model identification, specified as an `nlarxOptions` option set.

## Name-Value Pair Arguments

Specify optional comma-separated pairs of `Name`, `Value` arguments. `Name` is the argument name and `Value` is the corresponding value. `Name` must appear inside quotes. You can specify several name and value pair arguments in any order as `Name1,Value1,...,NameN,ValueN`.

Example: `'NonlinearRegressors','output'` specifies that only the regressors containing output variables are used as inputs to the nonlinear block of the model.

### **TimeVariable** — Independent variable name

`'t'` (default) | character vector

Independent variable name, specified as the comma-separated pair consisting of `'TimeVariable'` and a character vector. For example, `'t'`.

### **CustomRegressors** — Regressors constructed from combinations of inputs and outputs

`{}` (default) | cell array of character vectors | array of `customreg` objects

Regressors constructed from combinations of inputs and outputs, specified as the comma-separated pair consisting of `'CustomRegressors'` and one of the following for single-output systems:

- Cell array of character vectors. For example:
  - `{'y1(t-3)^3','y2(t-1)*u1(t-3)','sin(u3(t-2))'}`

Each character vector must represent a valid formula for a regressor contributing towards the prediction of the model output. The formula must be written using the input and output names and the time variable name as variables.

- Array of custom regressor objects, created using `customreg` or `polyreg`.

For a model with  $n_y$  outputs, specify an  $n_y$ -by-1 cell array of `customreg` object arrays or character arrays.

These regressors are in addition to the standard regressors based on `Orders`.

Example: `'CustomRegressors',{ 'y1(t-3)^3', 'y2(t-1)*u1(t-3)' }`

Example: `'CustomRegressors',{ 'sin(u3(t-2))' }`

**NonlinearRegressors — Subset of regressors that enter as inputs to the nonlinear block of the model**

'all' (default) | 'output' | 'input' | 'standard' | 'custom' | 'search' | vector of positive integers | [] | cell array

Subset of regressors that enter as inputs to the nonlinear block of the model, specified as the comma-separated pair consisting of 'NonlinearRegressors' and one of the following values:

- 'all' — All regressors
- 'output' — Regressors containing output variables
- 'input' — Regressors containing input variables
- 'standard' — Standard regressors
- 'custom' — Custom regressors
- 'search' — The estimation algorithm performs a search for the best regressor subset. This is useful when you want to reduce a large number of regressors entering the nonlinear function block of the nonlinearity estimator. This option must be applied to all output models simultaneously.
- [] — No regressors. This creates a linear-in-regressor model.
- Vector of regressor indices. To determine the number and order of regressors, use `getreg`.

For a model with multiple outputs, specify a cell array of  $n_y$  elements, where  $n_y$  is the number of output channels. For each output, specify one of the preceding options. Alternatively, to apply the same regressor subset to all model outputs, specify [] or any of the character vector options alone, for example 'standard'.

Example: 'NonlinearRegressors', 'search' performs a best regressor search for the only output of a single output model, or all of the outputs of a multiple output model.

Example: 'NonlinearReg', 'input' applies only input regressors to the inputs of the nonlinear function.

Example: 'NonlinearRegressors', {'input', 'output'} applies input regressors to the first output, and output regressors to the second output of a model with two outputs.

## Output Arguments

### **sys** — Nonlinear ARX model

`idnlarx` object

Nonlinear ARX model that fits the given estimation data, returned as an `idnlarx` object. This model is created using the specified model orders, nonlinearity estimator, and estimation options.

Information about the estimation results and options used is stored in the `Report` property of the model. The contents of `Report` depend upon the choice of nonlinearity and estimation focus you specified for `nlarx`. `Report` has the following fields:

<b>Report Field</b>	<b>Description</b>	
Status	Summary of the model status, which indicates whether the model was created by construction or obtained by estimation.	
Method	Estimation command used.	
Fit	Quantitative assessment of the estimation, returned as a structure. See “Loss Function and Model Quality Metrics” for more information on these quality metrics. The structure has the following fields:	
	<b>Field</b>	<b>Description</b>
	FitPercent	Normalized root mean squared error (NRMSE) measure of how well the response of the model fits the estimation data, expressed as a percentage.
	LossFcn	Value of the loss function when the estimation completes.
	MSE	Mean squared error (MSE) measure of how well the response of the model fits the estimation data.
	FPE	Final prediction error for the model.
	AIC	Raw Akaike Information Criteria (AIC) measure of model quality.
	AICc	Small sample-size corrected AIC.
	nAIC	Normalized AIC.
	BIC	Bayesian Information Criteria (BIC).

Report Field	Description																
Parameters	Estimated values of model parameters.																
OptionsUsed	Option set used for estimation. If no custom options were configured, this is a set of default options. See <code>nlrxOptions</code> for more information.																
RandState	State of the random number stream at the start of estimation. Empty, <code>[]</code> , if randomization was not used during estimation. For more information, see <code>rng</code> in the MATLAB documentation.																
DataUsed	Attributes of the data used for estimation, returned as a structure with the following fields: <table border="1" data-bbox="387 683 1334 1367"> <thead> <tr> <th>Field</th> <th>Description</th> </tr> </thead> <tbody> <tr> <td>Name</td> <td>Name of the data set.</td> </tr> <tr> <td>Type</td> <td>Data type.</td> </tr> <tr> <td>Length</td> <td>Number of data samples.</td> </tr> <tr> <td>Ts</td> <td>Sample time.</td> </tr> <tr> <td>InterSample</td> <td>Input intersample behavior, returned as one of the following values: <ul style="list-style-type: none"> <li>'zoh' — Zero-order hold maintains a piecewise-constant input signal between samples.</li> <li>'foh' — First-order hold maintains a piecewise-linear input signal between samples.</li> <li>'bl' — Band-limited behavior specifies that the continuous-time input signal has zero power above the Nyquist frequency.</li> </ul> </td> </tr> <tr> <td>InputOffset</td> <td>Offset removed from time-domain input data during estimation. For nonlinear models, it is <code>[]</code>.</td> </tr> <tr> <td>OutputOffset</td> <td>Offset removed from time-domain output data during estimation. For nonlinear models, it is <code>[]</code>.</td> </tr> </tbody> </table>	Field	Description	Name	Name of the data set.	Type	Data type.	Length	Number of data samples.	Ts	Sample time.	InterSample	Input intersample behavior, returned as one of the following values: <ul style="list-style-type: none"> <li>'zoh' — Zero-order hold maintains a piecewise-constant input signal between samples.</li> <li>'foh' — First-order hold maintains a piecewise-linear input signal between samples.</li> <li>'bl' — Band-limited behavior specifies that the continuous-time input signal has zero power above the Nyquist frequency.</li> </ul>	InputOffset	Offset removed from time-domain input data during estimation. For nonlinear models, it is <code>[]</code> .	OutputOffset	Offset removed from time-domain output data during estimation. For nonlinear models, it is <code>[]</code> .
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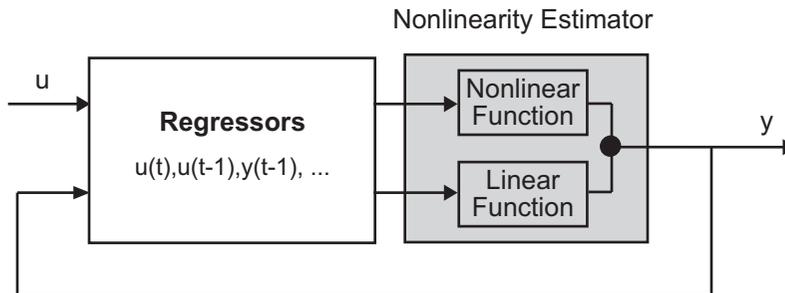
Report Field	Description																
Termination	Termination conditions for the iterative search used for prediction error minimization. Structure with the following fields:																
	<table border="1"> <thead> <tr> <th data-bbox="387 425 516 465">Field</th> <th data-bbox="516 425 1335 465">Description</th> </tr> </thead> <tbody> <tr> <td data-bbox="387 465 516 512">WhyStop</td> <td data-bbox="516 465 1335 512">Reason for terminating the numerical search.</td> </tr> <tr> <td data-bbox="387 512 516 586">Iterations</td> <td data-bbox="516 512 1335 586">Number of search iterations performed by the estimation algorithm.</td> </tr> <tr> <td data-bbox="387 586 516 696">FirstOrderOptimality</td> <td data-bbox="516 586 1335 696"><math>\infty</math>-norm of the gradient search vector when the search algorithm terminates.</td> </tr> <tr> <td data-bbox="387 696 516 770">FcnCount</td> <td data-bbox="516 696 1335 770">Number of times the objective function was called.</td> </tr> <tr> <td data-bbox="387 770 516 845">UpdateNorm</td> <td data-bbox="516 770 1335 845">Norm of the gradient search vector in the last iteration. Omitted when the search method is 'lsqnonlin' or 'fmincon'.</td> </tr> <tr> <td data-bbox="387 845 516 954">LastImprovement</td> <td data-bbox="516 845 1335 954">Criterion improvement in the last iteration, expressed as a percentage. Omitted when the search method is 'lsqnonlin' or 'fmincon'.</td> </tr> <tr> <td data-bbox="387 954 516 1029">Algorithm</td> <td data-bbox="516 954 1335 1029">Algorithm used by 'lsqnonlin' or 'fmincon' search method. Omitted when other search methods are used.</td> </tr> </tbody> </table>	Field	Description	WhyStop	Reason for terminating the numerical search.	Iterations	Number of search iterations performed by the estimation algorithm.	FirstOrderOptimality	$\infty$ -norm of the gradient search vector when the search algorithm terminates.	FcnCount	Number of times the objective function was called.	UpdateNorm	Norm of the gradient search vector in the last iteration. Omitted when the search method is 'lsqnonlin' or 'fmincon'.	LastImprovement	Criterion improvement in the last iteration, expressed as a percentage. Omitted when the search method is 'lsqnonlin' or 'fmincon'.	Algorithm	Algorithm used by 'lsqnonlin' or 'fmincon' search method. Omitted when other search methods are used.
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For estimation methods that do not require numerical search optimization, the Termination field is omitted.																	

For more information on using Report, see “Estimation Report”.

## Algorithms

### Nonlinear ARX Model Structure

A nonlinear ARX model consists of model regressors and a nonlinearity estimator. The nonlinearity estimator comprises both linear and nonlinear functions that act on the model regressors to give the model output. This block diagram represents the structure of a nonlinear ARX model in a simulation scenario.



The software computes the nonlinear ARX model output  $y$  in two stages:

- 1 It computes regressor values from the current and past input values and past output data.

In the simplest case, regressors are delayed inputs and outputs, such as  $u(t-1)$  and  $y(t-3)$ . These kind of regressors are called *standard regressors*. You specify the standard regressors using the model orders and delay. For more information, see “Nonlinear ARX Model Orders and Delay”. You can also specify *custom* regressors, which are nonlinear functions of delayed inputs and outputs. For example,  $u(t-1)*y(t-3)$ . To create a set of polynomial type regressors, use `polyreg`.

By default, all regressors are inputs to both the linear and the nonlinear function blocks of the nonlinearity estimator. You can choose a subset of regressors as inputs to the nonlinear function block.

- 2 It maps the regressors to the model output using the nonlinearity estimator block. The nonlinearity estimator block can include linear and nonlinear blocks in parallel. For example:

$$F(x) = L^T(x-r) + d + g(Q(x-r))$$

Here,  $x$  is a vector of the regressors, and  $r$  is the mean of the regressors  $x$ .  $L^T(x-r) + d$  is the output of the linear function block and is affine when  $d \neq 0$ .  $d$  is a scalar offset.

$g(Q(x-r))$  represents the output of the nonlinear function block.  $Q$  is a projection matrix that makes the calculations well conditioned. The exact form of  $F(x)$  depends on your choice of the nonlinearity estimator. You can select from available nonlinearity estimators, such as tree-partition networks, wavelet networks, and

multilayer neural networks. You can also exclude either the linear or the nonlinear function block from the nonlinearity estimator.

When estimating a nonlinear ARX model, the software computes the model parameter values, such as  $L$ ,  $r$ ,  $d$ ,  $Q$ , and other parameters specifying  $g$ .

Resulting nonlinear ARX models are `idnlarx` objects that store all model data, including model regressors and parameters of the nonlinearity estimator. For more information about these objects, see “Nonlinear Model Structures”.

## Extended Capabilities

### Automatic Parallel Support

Accelerate code by automatically running computation in parallel using Parallel Computing Toolbox™.

Parallel computing support is available for estimation using the `lsqnonlin` search method (requires Optimization Toolbox). To enable parallel computing, use `nlarxOptions`, set `SearchMethod` to `'lsqnonlin'`, and set `SearchOptions.Advanced.UseParallel` to `true`.

For example:

```
opt = nlarxOptions;  
opt.SearchMethod = 'lsqnonlin';  
opt.SearchOptions.Advanced.UseParallel = true;
```

### See Also

`aic` | `fpe` | `godnessofFit` | `idnlarx` | `isnlarx` | `nlarxOptions`

### Topics

“Estimate Nonlinear ARX Models at the Command Line”

“Estimate Nonlinear ARX Models Initialized Using Linear ARX Models”

“Identifying Nonlinear ARX Models”

“Validate Nonlinear ARX Models”

“Using Nonlinear ARX Models”

"Loss Function and Model Quality Metrics"  
"Regularized Estimates of Model Parameters"  
"Estimation Report"

**Introduced in R2007a**

## **nlarxOptions**

Option set for `nlarx`

### **Syntax**

```
opt = nlarxOptions  
opt = nlarxOptions(Name,Value)
```

### **Description**

`opt = nlarxOptions` creates the default option set for `nlarx`. Use dot notation to modify this option set for your specific application. Any options that you do not modify retain their default values.

`opt = nlarxOptions(Name,Value)` creates an option set with options specified by one or more `Name,Value` pair arguments.

### **Examples**

#### **Create Default Option Set for Nonlinear ARX Estimation**

```
opt = nlarxOptions;
```

#### **Create and Modify Default Nonlinear ARX Option Set**

Create a default option set for `nlarx`, and use dot notation to modify specific options.

```
opt = nlarxOptions;
```

Turn on the estimation progress display.

```
opt.Display = 'on';
```

Minimize the norm of the simulation error.

```
opt.Focus = 'simulation';
```

Use a subspace Gauss-Newton least squares search with a maximum of 25 iterations.

```
opt.SearchMethod = 'gn';  
opt.SearchOptions.MaxIterations = 25;
```

## Specify Options for Nonlinear ARX Estimation

Create an option set for `nlarx` specifying the following options:

- Turn off iterative estimation for the default wavelet network estimation.
- Turn on the estimation progress-viewer display.

```
opt = nlarxOptions('IterativeWavenet','off','Display','on');
```

# Input Arguments

## Name-Value Pair Arguments

Specify optional comma-separated pairs of `Name`, `Value` arguments. `Name` is the argument name and `Value` is the corresponding value. `Name` must appear inside quotes. You can specify several name and value pair arguments in any order as `Name1,Value1,...,NameN,ValueN`.

Example: `'Focus','simulation','SearchMethod','grad'` specifies that the norm of the simulation error is minimized using a steepest descent least squares search.

### Focus — Minimization objective

```
'prediction' (default) | 'simulation'
```

Minimization objective, specified as the comma-separated pair consisting of `'Focus'` and one of the following:

- `'prediction'` — Minimize the norm of the prediction error, which is defined as the difference between the measured output and the one-step ahead predicted response of the model.
- `'simulation'` — Minimize the norm of the simulation error, which is defined as the difference between the measured output and simulated response of the model.

### **Display — Estimation progress display setting**

`'off'` (default) | `'on'`

Estimation progress display setting, specified as the comma-separated pair consisting of `'Display'` and one of the following:

- `'off'` — No progress or results information is displayed.
- `'on'` — Information on model structure and estimation results are displayed in a progress-viewer window.

### **OutputWeight — Weighting of prediction error in multi-output estimations**

`'noise'` (default) | positive semidefinite matrix

Weighting of prediction error in multi-output model estimations, specified as the comma-separated pair consisting of `'OutputWeight'` and one of the following:

- `'noise'` — Optimal weighting is automatically computed as the inverse of the estimated noise variance. This weighting minimizes  $\det(E'E)$ , where  $E$  is the matrix of prediction errors. This option is not available when using `'lsqnonlin'` as a `'SearchMethod'`.
- A positive semidefinite matrix,  $W$ , of size equal to the number of outputs. This weighting minimizes  $\text{trace}(E'E*W/N)$ , where  $E$  is the matrix of prediction errors and  $N$  is the number of data samples.

### **IterativeWavenet — Iterative wavenet estimation setting**

`'auto'` (default) | `'on'` | `'off'`

Iterative wavenet estimation setting, specified as the comma-separated pair consisting of `'IterativeWavenet'` and one of the following:

- `'auto'` — First estimation is noniterative and subsequent estimations are iterative.
- `'on'` — Perform iterative estimation only.
- `'off'` — Perform noniterative estimation only.

This option applies only when using a wavenet nonlinearity estimator.

## Regularization — Options for regularized estimation of model parameters structure

Options for regularized estimation of model parameters, specified as the comma-separated pair consisting of 'Regularization' and a structure with fields:

Field Name	Description	Default
Lambda	Bias versus variance trade-off constant, specified as a nonnegative scalar.	0 — Indicates no regularization.
R	Weighting matrix, specified as a vector of nonnegative scalars or a square positive semidefinite matrix. The length must be equal to the number of free parameters in the model, np. Use the nparams command to determine the number of model parameters.	1 — Indicates a value of eye(np).
Nominal	The nominal value towards which the free parameters are pulled during estimation, specified as one of the following: <ul style="list-style-type: none"> <li>'zero' — Pull parameters towards zero.</li> <li>'model' — Pull parameters towards preexisting values in the initial model. Use this option only when you have a well-initialized idnlarx model with finite parameter values.</li> </ul>	'zero'

To specify field values in Regularization, create a default nlarxOptions set and modify the fields using dot notation. Any fields that you do not modify retain their default values.

```
opt = nlarxOptions;
opt.Regularization.Lambda = 1.2;
opt.Regularization.R = 0.5*eye(np);
```

Regularization is a technique for specifying model flexibility constraints, which reduce uncertainty in the estimated parameter values. For more information, see “Regularized Estimates of Model Parameters”.

## SearchMethod — Numerical search method used for iterative parameter estimation

'auto' (default) | 'gn' | 'gna' | 'lm' | 'grad' | 'lsqnonlin' | 'fmincon'

Numerical search method used for iterative parameter estimation, specified as the comma-separated pair consisting of 'SearchMethod' and one of the following:

- 'auto' — A combination of the line search algorithms, 'gn', 'lm', 'gna', and 'grad' methods is tried in sequence at each iteration. The first descent direction leading to a reduction in estimation cost is used.
- 'gn' — Subspace Gauss-Newton least squares search. Singular values of the Jacobian matrix less than  $\text{GnPinvConstant} \cdot \text{eps} \cdot \max(\text{size}(J)) \cdot \text{norm}(J)$  are discarded when computing the search direction.  $J$  is the Jacobian matrix. The Hessian matrix is approximated as  $J^T J$ . If there is no improvement in this direction, the function tries the gradient direction.
- 'gna' — Adaptive subspace Gauss-Newton search. Eigenvalues less than  $\text{gamma} \cdot \max(\text{sv})$  of the Hessian are ignored, where  $\text{sv}$  contains the singular values of the Hessian. The Gauss-Newton direction is computed in the remaining subspace.  $\text{gamma}$  has the initial value `InitialGnaTolerance` (see `Advanced` in 'SearchOptions' for more information). This value is increased by the factor `LMStep` each time the search fails to find a lower value of the criterion in fewer than five bisections. This value is decreased by the factor  $2 \cdot \text{LMStep}$  each time a search is successful without any bisections.
- 'lm' — Levenberg-Marquardt least squares search, where the next parameter value is  $-\text{pinv}(H+d \cdot I) \cdot \text{grad}$  from the previous one.  $H$  is the Hessian,  $I$  is the identity matrix, and  $\text{grad}$  is the gradient.  $d$  is a number that is increased until a lower value of the criterion is found.
- 'grad' — Steepest descent least squares search.
- 'lsqnonlin' — Trust-region-reflective algorithm of `lsqnonlin`. Requires Optimization Toolbox software.
- 'fmincon' — Constrained nonlinear solvers. You can use the sequential quadratic programming (SQP) and trust-region-reflective algorithms of the `fmincon` solver. If you have Optimization Toolbox software, you can also use the interior-point and active-set algorithms of the `fmincon` solver. Specify the algorithm in the `SearchOptions.Algorithm` option. The `fmincon` algorithms may result in improved estimation results in the following scenarios:
  - Constrained minimization problems when there are bounds imposed on the model parameters.
  - Model structures where the loss function is a nonlinear or non smooth function of the parameters.

- Multi-output model estimation. A determinant loss function is minimized by default for multi-output model estimation. `fmincon` algorithms are able to minimize such loss functions directly. The other search methods such as `'lm'` and `'gn'` minimize the determinant loss function by alternately estimating the noise variance and reducing the loss value for a given noise variance value. Hence, the `fmincon` algorithms can offer better efficiency and accuracy for multi-output model estimations.

### **SearchOptions — Options set for the search algorithm**

search option set

Options set for the search algorithm, specified as the comma-separated pair consisting of `'SearchOptions'` and a search option set with fields that depend on the value of `SearchMethod`:

**SearchOptions Structure When SearchMethod is Specified as 'gn', 'gna', 'lm', 'grad', or 'auto'**

Field Name	Description	Default
Tolerance	<p>Minimum percentage difference between the current value of the loss function and its expected improvement after the next iteration, specified as a positive scalar. When the percentage of expected improvement is less than <code>Tolerance</code>, the iterations stop. The estimate of the expected loss-function improvement at the next iteration is based on the Gauss-Newton vector computed for the current parameter value.</p>	1e-5
MaxIterations	<p>Maximum number of iterations during loss-function minimization, specified as a positive integer. The iterations stop when <code>MaxIterations</code> is reached or another stopping criterion is satisfied, such as <code>Tolerance</code>.</p> <p>Setting <code>MaxIterations = 0</code> returns the result of the start-up procedure.</p> <p>Use <code>sys.Report.Termination.Iterations</code> to get the actual number of iterations during an estimation, where <code>sys</code> is an <code>idtf</code> model.</p>	20

Field Name	Description	Default
Advanced	Advanced search settings, specified as a structure with the following fields:	
Field Name	Description	Default
GnPinvConstant	Jacobian matrix singular value threshold, specified as a positive scalar. Singular values of the Jacobian matrix that are smaller than $\text{GnPinvConstant} * \max(\text{size}(J) * \text{norm}(J) * \text{eps})$ are discarded when computing the search direction. Applicable when SearchMethod is 'gn'.	10000
InitialGnaTolerance	Initial value of $\gamma$ , specified as a positive scalar. Applicable when SearchMethod is 'gna'.	0.0001
LMStartValue	Starting value of search-direction length $d$ in the Levenberg-Marquardt method, specified as a positive scalar. Applicable when SearchMethod is 'lm'.	0.001
LMStep	Size of the Levenberg-Marquardt step, specified as a positive integer. The next value of the search-direction length $d$ in the Levenberg-Marquardt method is LMStep times the previous one. Applicable when SearchMethod is 'lm'.	2
MaxBisections	Maximum number of bisections used for line search along the search direction, specified as a positive integer.	25
MaxFunctionEvaluations	Maximum number of calls to the model file, specified as a positive integer. Iterations stop if the number of calls to the model file exceeds this value.	Inf
MinParameterChange	Smallest parameter update allowed per iteration, specified as a nonnegative scalar.	0
RelativeImprovement	Relative improvement threshold, specified as a nonnegative scalar. Iterations stop if the relative	0

<b>Field Name</b>	<b>Description</b>		<b>Default</b>
	<b>Field Name</b>	<b>Description</b>	<b>Default</b>
		improvement of the criterion function is less than this value.	
	StepReduction	<p>Step reduction factor, specified as a positive scalar that is greater than 1. The suggested parameter update is reduced by the factor <code>StepReduction</code> after each try. This reduction continues until <code>MaxBisections</code> tries are completed or a lower value of the criterion function is obtained.</p> <p><code>StepReduction</code> is not applicable for SearchMethod 'lm' (Levenberg-Marquardt method).</p>	2

**SearchOptions Structure When SearchMethod is Specified as 'lsqnonlin'**

<b>Field Name</b>	<b>Description</b>	<b>Default</b>
FunctionTolerance	Termination tolerance on the loss function that the software minimizes to determine the estimated parameter values, specified as a positive scalar.  The value of FunctionTolerance is the same as that of <code>opt.SearchOptions.Advanced.TolFun</code> .	1e-5
StepTolerance	Termination tolerance on the estimated parameter values, specified as a positive scalar.  The value of StepTolerance is the same as that of <code>opt.SearchOptions.Advanced.TolX</code> .	1e-6
MaxIterations	Maximum number of iterations during loss-function minimization, specified as a positive integer. The iterations stop when MaxIterations is reached or another stopping criterion is satisfied, such as FunctionTolerance.  The value of MaxIterations is the same as that of <code>opt.SearchOptions.Advanced.MaxIter</code> .	20
Advanced	Advanced search settings, specified as an option set for <code>lsqnonlin</code> .  For more information, see the Optimization Options table in "Optimization Options" (Optimization Toolbox).	Use <code>optimset('lsqnonlin')</code> to create a default option set.

**SearchOptions Structure When SearchMethod is Specified as 'fmincon'**

Field Name	Description	Default
Algorithm	<p>fmincon optimization algorithm, specified as one of the following:</p> <ul style="list-style-type: none"> <li>• 'sqp' — Sequential quadratic programming algorithm. The algorithm satisfies bounds at all iterations, and it can recover from NaN or Inf results. It is not a large-scale algorithm. For more information, see “Large-Scale vs. Medium-Scale Algorithms” (Optimization Toolbox).</li> <li>• 'trust-region-reflective' — Subspace trust-region method based on the interior-reflective Newton method. It is a large-scale algorithm.</li> <li>• 'interior-point' — Large-scale algorithm that requires Optimization Toolbox software. The algorithm satisfies bounds at all iterations, and it can recover from NaN or Inf results.</li> <li>• 'active-set' — Requires Optimization Toolbox software. The</li> </ul>	'sqp'

Field Name	Description	Default
	<p>algorithm can take large steps, which adds speed. It is not a large-scale algorithm.</p> <p>For more information about the algorithms, see “Constrained Nonlinear Optimization Algorithms” (Optimization Toolbox) and “Choosing the Algorithm” (Optimization Toolbox).</p>	
FunctionTolerance	Termination tolerance on the loss function that the software minimizes to determine the estimated parameter values, specified as a positive scalar.	1e-6
StepTolerance	Termination tolerance on the estimated parameter values, specified as a positive scalar.	1e-6
MaxIterations	Maximum number of iterations during loss function minimization, specified as a positive integer. The iterations stop when MaxIterations is reached or another stopping criterion is satisfied, such as FunctionTolerance.	100

To specify field values in `SearchOptions`, create a default `nlrxOptions` set and modify the fields using dot notation. Any fields that you do not modify retain their default values.

```
opt = nlarxOptions;
opt.SearchOptions.MaxIter = 15;
opt.SearchOptions.Advanced.RelImprovement = 0.5;
```

**Advanced — Additional advanced options**

structure

Additional advanced options, specified as the comma-separated pair consisting of 'Advanced' and a structure with fields:

Field Name	Description	Default
ErrorThreshold	Threshold for when to adjust the weight of large errors from quadratic to linear, specified as a nonnegative scalar. Errors larger than ErrorThreshold times the estimated standard deviation have a linear weight in the loss function. The standard deviation is estimated robustly as the median of the absolute deviations from the median of the prediction errors, divided by 0.7. If your estimation data contains outliers, try setting ErrorThreshold to 1.6.	0 — Leads to a purely quadratic loss function.
MaxSize	Maximum number of elements in a segment when input-output data is split into segments, specified as a positive integer.	250000

To specify field values in Advanced, create a default nlarxOptions set and modify the fields using dot notation. Any fields that you do not modify retain their default values.

```
opt = nlarxOptions;
opt.Advanced.ErrorThreshold = 1.2;
```

## Output Arguments

**opt — Option set for nlarx command**

nlarxOptions option set

Option set for nlarx command, returned as an nlarxOptions option set.

## **See Also**

nlarx

**Introduced in R2015a**

## nlgreyest

Estimate nonlinear grey-box model parameters

### Syntax

```
sys= nlgreyest(data,init_sys)  
sys= nlgreyest(data,init_sys,options)
```

### Description

`sys= nlgreyest(data,init_sys)` estimates the parameters of a nonlinear grey-box model, `init_sys`, using time-domain data, `data`.

`sys= nlgreyest(data,init_sys,options)` specifies additional model estimation options.

### Examples

#### Selectively Estimate Parameters of Nonlinear Grey-Box Model

Load data.

```
load(fullfile(matlabroot,'toolbox','ident','iddemos','data','twotankdata'));  
z = iddata(y,u,0.2,'Name','Two tanks');
```

The data contains 3000 input-output data samples of a two tank system. The input is the voltage applied to a pump, and the output is the liquid level of the lower tank.

Specify file describing the model structure for a two-tank system. The file specifies the state derivatives and model outputs as a function of time, states, inputs, and model parameters.

```
FileName = 'twotanks_c';
```

Specify model orders [ny nu nx].

```
Order = [1 1 2];
```

Specify initial parameters (Np = 6).

```
Parameters = {0.5;0.0035;0.019; ...
              9.81;0.25;0.016};
```

Specify initial initial states.

```
InitialStates = [0;0.1];
```

Specify as continuous system.

```
Ts = 0;
```

Create idnlgrey model object.

```
nlgr = idnlgrey(FileName,Order,Parameters,InitialStates,Ts, ...
               'Name','Two tanks');
```

Set some parameters as constant.

```
nlgr.Parameters(1).Fixed = true;
nlgr.Parameters(4).Fixed = true;
nlgr.Parameters(5).Fixed = true;
```

Estimate the model parameters.

```
nlgr = nlgreyest(z,nlgr);
```

### Estimate a Nonlinear Grey-Box Model Using Specific Options

Create estimation option set for nlgreyest to view estimation progress, and to set the maximum iteration steps to 50.

```
opt = nlgreyestOptions;
opt.Display = 'on';
opt.SearchOptions.MaxIterations = 50;
```

Load data.

```
load(fullfile(matlabroot,'toolbox','ident','iddemos','data','dcmotordata'));
z = iddata(y,u,0.1,'Name','DC-motor');
```

The data is from a linear DC motor with one input (voltage), and two outputs (angular position and angular velocity). The structure of the model is specified by `dcmotor_m.m` file.

Create a nonlinear grey-box model.

```
file_name = 'dcmotor_m';
Order = [2 1 2];
Parameters = [1;0.28];
InitialStates = [0;0];

init_sys = idnlgrey(file_name,Order,Parameters,InitialStates,0, ...
    'Name','DC-motor');
```

Estimate the model parameters using the estimation options.

```
sys = nlgreyest(z,init_sys,opt);
```

## Input Arguments

### **data** — Time domain data

`iddata` object

Time-domain estimation data, specified as an `iddata` object. `data` has the same input and output dimensions as `init_sys`.

If you specify the `InterSample` property of `data` as `'bl'` (band-limited) and the model is continuous-time, the software treats data as first-order-hold (foh) interpolated for estimation.

### **init\_sys** — Constructed nonlinear grey-box model

`idnlgrey` object

Constructed nonlinear grey-box model that configures the initial parameterization of `sys`, specified as an `idnlgrey` object. `init_sys` has the same input and output dimensions as `data`. Create `init_sys` using `idnlgrey`.

### **options** — Estimation options

`nlgreyestOptions` option set

Estimation options for nonlinear grey-box model identification, specified as an `nlgreyestOptions` option set.

## Output Arguments

### **sys** — Estimated nonlinear grey-box model

`idnlgrey` object

Nonlinear grey-box model with the same structure as `init_sys`, returned as an `idnlgrey` object. The parameters of `sys` are estimated such that the response of `sys` matches the output signal in the estimation data.

Information about the estimation results and options used is stored in the `Report` property of the model. `Report` has the following fields:

<b>Report Field</b>	<b>Description</b>
Status	Summary of the model status, which indicates whether the model was created by construction or obtained by estimation.
Method	Name of the simulation solver and the search method used during estimation.

Report Field	Description																		
Fit	<p>Quantitative assessment of the estimation, returned as a structure. See “Loss Function and Model Quality Metrics” for more information on these quality metrics. The structure has the following fields:</p> <table border="1" data-bbox="387 456 1332 951"> <thead> <tr> <th data-bbox="387 456 516 499">Field</th> <th data-bbox="516 456 1332 499">Description</th> </tr> </thead> <tbody> <tr> <td data-bbox="387 499 516 609">FitPercent</td> <td data-bbox="516 499 1332 609">Normalized root mean squared error (NRMSE) measure of how well the response of the model fits the estimation data, expressed as a percentage.</td> </tr> <tr> <td data-bbox="387 609 516 652">LossFcn</td> <td data-bbox="516 609 1332 652">Value of the loss function when the estimation completes.</td> </tr> <tr> <td data-bbox="387 652 516 730">MSE</td> <td data-bbox="516 652 1332 730">Mean squared error (MSE) measure of how well the response of the model fits the estimation data.</td> </tr> <tr> <td data-bbox="387 730 516 774">FPE</td> <td data-bbox="516 730 1332 774">Final prediction error for the model.</td> </tr> <tr> <td data-bbox="387 774 516 817">AIC</td> <td data-bbox="516 774 1332 817">Raw Akaike Information Criteria (AIC) measure of model quality.</td> </tr> <tr> <td data-bbox="387 817 516 861">AICc</td> <td data-bbox="516 817 1332 861">Small sample-size corrected AIC.</td> </tr> <tr> <td data-bbox="387 861 516 904">nAIC</td> <td data-bbox="516 861 1332 904">Normalized AIC.</td> </tr> <tr> <td data-bbox="387 904 516 951">BIC</td> <td data-bbox="516 904 1332 951">Bayesian Information Criteria (BIC).</td> </tr> </tbody> </table>	Field	Description	FitPercent	Normalized root mean squared error (NRMSE) measure of how well the response of the model fits the estimation data, expressed as a percentage.	LossFcn	Value of the loss function when the estimation completes.	MSE	Mean squared error (MSE) measure of how well the response of the model fits the estimation data.	FPE	Final prediction error for the model.	AIC	Raw Akaike Information Criteria (AIC) measure of model quality.	AICc	Small sample-size corrected AIC.	nAIC	Normalized AIC.	BIC	Bayesian Information Criteria (BIC).
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Parameters	<p>Estimated values of the model parameters. Structure with the following fields:</p> <table border="1" data-bbox="387 1052 1332 1459"> <thead> <tr> <th data-bbox="387 1052 639 1095">Field</th> <th data-bbox="639 1052 1332 1095">Description</th> </tr> </thead> <tbody> <tr> <td data-bbox="387 1095 639 1173">InitialValues</td> <td data-bbox="639 1095 1332 1173">Structure with values of parameters and initial states before estimation.</td> </tr> <tr> <td data-bbox="387 1173 639 1216">ParVector</td> <td data-bbox="639 1173 1332 1216">Value of parameters after estimation.</td> </tr> <tr> <td data-bbox="387 1216 639 1295">Free</td> <td data-bbox="639 1216 1332 1295">Logical vector specifying the fixed or free status of parameters during estimation</td> </tr> <tr> <td data-bbox="387 1295 639 1373">FreeParCovariance</td> <td data-bbox="639 1295 1332 1373">Covariance of the free parameters.</td> </tr> <tr> <td data-bbox="387 1373 639 1416">X0</td> <td data-bbox="639 1373 1332 1416">Value of initial states after estimation.</td> </tr> <tr> <td data-bbox="387 1416 639 1459">X0Covariance</td> <td data-bbox="639 1416 1332 1459">Covariance of the initial states.</td> </tr> </tbody> </table>	Field	Description	InitialValues	Structure with values of parameters and initial states before estimation.	ParVector	Value of parameters after estimation.	Free	Logical vector specifying the fixed or free status of parameters during estimation	FreeParCovariance	Covariance of the free parameters.	X0	Value of initial states after estimation.	X0Covariance	Covariance of the initial states.				
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OptionsUsed	Option set used for estimation. If no custom options were configured, this is a set of default options. See <code>nlgreyestOptions</code> for more information.																		

Report Field	Description																
RandState	State of the random number stream at the start of estimation. Empty, [], if randomization was not used during estimation. For more information, see <code>rng</code> in the MATLAB documentation.																
DataUsed	Attributes of the data used for estimation — Structure with the following fields:																
	<table border="1"> <thead> <tr> <th data-bbox="385 534 516 574">Field</th> <th data-bbox="516 534 1337 574">Description</th> </tr> </thead> <tbody> <tr> <td data-bbox="385 574 516 618">Name</td> <td data-bbox="516 574 1337 618">Name of the data set.</td> </tr> <tr> <td data-bbox="385 618 516 696">Type</td> <td data-bbox="516 618 1337 696">Data type — For <code>idnlgrey</code> models, this is set to 'Time domain data'.</td> </tr> <tr> <td data-bbox="385 696 516 739">Length</td> <td data-bbox="516 696 1337 739">Number of data samples.</td> </tr> <tr> <td data-bbox="385 739 516 782">Ts</td> <td data-bbox="516 739 1337 782">Sample time. This is equivalent to <code>data.Ts</code>.</td> </tr> <tr> <td data-bbox="385 782 516 1156">InterSample</td> <td data-bbox="516 782 1337 1156">           Input intersample behavior. One of the following values:           <ul style="list-style-type: none"> <li>• 'zoh' — Zero-order hold maintains a piecewise-constant input signal between samples.</li> <li>• 'foh' — First-order hold maintains a piecewise-linear input signal between samples.</li> <li>• 'bl' — Band-limited behavior specifies that the continuous-time input signal has zero power above the Nyquist frequency.</li> </ul>           The value of <code>Intersample</code> has no effect on estimation results for discrete-time models.         </td> </tr> <tr> <td data-bbox="385 1156 516 1234">InputOffset</td> <td data-bbox="516 1156 1337 1234">Empty, [], for nonlinear estimation methods.</td> </tr> <tr> <td data-bbox="385 1234 516 1310">OutputOffset</td> <td data-bbox="516 1234 1337 1310">Empty, [], for nonlinear estimation methods.</td> </tr> </tbody> </table>	Field	Description	Name	Name of the data set.	Type	Data type — For <code>idnlgrey</code> models, this is set to 'Time domain data'.	Length	Number of data samples.	Ts	Sample time. This is equivalent to <code>data.Ts</code> .	InterSample	Input intersample behavior. One of the following values: <ul style="list-style-type: none"> <li>• 'zoh' — Zero-order hold maintains a piecewise-constant input signal between samples.</li> <li>• 'foh' — First-order hold maintains a piecewise-linear input signal between samples.</li> <li>• 'bl' — Band-limited behavior specifies that the continuous-time input signal has zero power above the Nyquist frequency.</li> </ul> The value of <code>Intersample</code> has no effect on estimation results for discrete-time models.	InputOffset	Empty, [], for nonlinear estimation methods.	OutputOffset	Empty, [], for nonlinear estimation methods.
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InputOffset	Empty, [], for nonlinear estimation methods.																
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Report Field	Description																
Termination	Termination conditions for the iterative search used for prediction error minimization. Structure with the following fields:																
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For estimation methods that do not require numerical search optimization, the Termination field is omitted.																	

For more information, see “Estimation Report”.

## Extended Capabilities

### Automatic Parallel Support

Accelerate code by automatically running computation in parallel using Parallel Computing Toolbox™.

Parallel computing support is available for estimation using the `lsqnonlin` search method (requires Optimization Toolbox). To enable parallel computing, use `nlgreyestOptions`, set `SearchMethod` to `'lsqnonlin'`, and set `SearchOptions.Advanced.UseParallel` to `true`.

For example:

```
opt = nlgreyestOptions;  
opt.SearchMethod = 'lsqnonlin';  
opt.SearchOptions.Advanced.UseParallel = true;
```

## See Also

[aic](#) | [fpe](#) | [goodnessofFit](#) | [idnlgrey](#) | [nlgreyestOptions](#) | [pem](#)

## Topics

[“Creating IDNLGREY Model Files”](#)

[“Represent Nonlinear Dynamics Using MATLAB File for Grey-Box Estimation”](#)

[“Estimate Nonlinear Grey-Box Models”](#)

[“Loss Function and Model Quality Metrics”](#)

[“Regularized Estimates of Model Parameters”](#)

[“Estimation Report”](#)

**Introduced in R2015a**

## **nlgreyestOptions**

Option set for `nlgreyest`

### **Syntax**

```
opt = nlgreyestOptions  
opt = nlgreyestOptions(Name,Value)
```

### **Description**

`opt = nlgreyestOptions` creates the default option set for `nlgreyest`. Use dot notation to customize the option set, if needed.

`opt = nlgreyestOptions(Name,Value)` creates an option set with options specified by one or more `Name, Value` pair arguments. The options that you do not specify retain their default value.

### **Examples**

#### **Create Default Option Set for Nonlinear Grey-Box Model Estimation**

```
opt = nlgreyestOptions;
```

#### **Estimate a Nonlinear Grey-Box Model Using Specific Options**

Create estimation option set for `nlgreyest` to view estimation progress, and to set the maximum iteration steps to 50.

```
opt = nlgreyestOptions;  
opt.Display = 'on';  
opt.SearchOptions.MaxIterations = 50;
```

Load data.

```
load(fullfile(matlabroot, 'toolbox', 'ident', 'iddemos', 'data', 'dcmotordata'));
z = iddata(y,u,0.1, 'Name', 'DC-motor');
```

The data is from a linear DC motor with one input (voltage), and two outputs (angular position and angular velocity). The structure of the model is specified by `dcmotor_m.m` file.

Create a nonlinear grey-box model.

```
file_name = 'dcmotor_m';
Order = [2 1 2];
Parameters = [1;0.28];
InitialStates = [0;0];

init_sys = idnlgrey(file_name,Order,Parameters,InitialStates,0, ...
    'Name', 'DC-motor');
```

Estimate the model parameters using the estimation options.

```
sys = nlgreyest(z,init_sys,opt);
```

## Specify Options for Nonlinear Grey-Box Model Estimation

Create an option set for `nlgreyest` where:

- Parameter covariance data is not generated.
- Subspace Gauss-Newton least squares method is used for estimation.

```
opt = nlgreyestOptions('EstimateCovariance',false, 'SearchMethod', 'gn');
```

## Input Arguments

### Name-Value Pair Arguments

Specify optional comma-separated pairs of `Name`, `Value` arguments. `Name` is the argument name and `Value` is the corresponding value. `Name` must appear inside quotes.

You can specify several name and value pair arguments in any order as `Name1, Value1, ..., NameN, ValueN`.

Example: `nlgreyestOptions('Display','on')`

**GradientOptions — Options for computing Jacobians and gradients**  
structure

Options for computing Jacobians and gradients, specified as the comma-separated pair consisting of 'GradientOptions' and a structure with fields:

Field Name	Description	Default
MaxDifference	Largest allowed parameter perturbation when computing numerical derivatives. Specified as a positive real value > 'MinDifference'.	Inf
MinDifference	Smallest allowed parameter perturbation when computing numerical derivatives. Specified as a positive real value < 'MaxDifference'.	0.01*sqrt(eps)
DifferenceScheme	Method for computing numerical derivatives with respect to the components of the parameters and/or the initial state(s) to form the Jacobian. Specified as one of the following: <ul style="list-style-type: none"> <li>'Auto' - Automatically chooses from the following methods.</li> <li>'Central approximation'</li> <li>'Forward approximation'</li> <li>'Backward approximation'</li> </ul>	'Auto'

Field Name	Description	Default
Type	<p>Method used when computing derivatives (Jacobian) of the parameters or the initial states to be estimated. Specified as one of the following:</p> <ul style="list-style-type: none"> <li>• 'Auto' — Automatically chooses from the following methods.</li> <li>• 'Basic' — Individually computes all numerical derivatives required to form each column of the Jacobian.</li> <li>• 'Refined' — Simultaneously computes all numerical derivatives required to form each column of the Jacobian.</li> </ul>	'Auto'

To specify field values in `GradientOptions`, create a default `nlgreyestOptions` set and modify the fields using dot notation. Any fields that you do not modify retain their default values.

```
opt = nlgreyestOptions;
opt.GradientOptions.Type = 'Basic';
```

### **EstimateCovariance — Parameter covariance data generation setting**

1 or true (default) | 0 or false

Controls whether parameter covariance data is generated, specified as true (1) or false (0).

### **Display — Estimation progress display setting**

'off' (default) | 'on'

Estimation progress display setting, specified as the comma-separated pair consisting of 'Display' and one of the following:

- 'off' — No progress or results information is displayed.
- 'on' — Information on model structure and estimation results are displayed in a progress-viewer window.

### **Regularization — Options for regularized estimation of model parameters**

structure

Options for regularized estimation of model parameters, specified as the comma-separated pair consisting of 'Regularization' and a structure with fields:

Field Name	Description	Default
Lambda	Bias versus variance trade-off constant, specified as a nonnegative scalar.	0 — Indicates no regularization.
R	Weighting matrix, specified as a vector of nonnegative scalars or a square positive semi-definite matrix. The length must be equal to the number of free parameters in the model, np. Use the nparams command to determine the number of model parameters.	1 — Indicates a value of eye(np).
Nominal	The nominal value towards which the free parameters are pulled during estimation specified as one of the following: <ul style="list-style-type: none"> <li>'zero' — Pull parameters towards zero.</li> <li>'model' — Pull parameters towards pre-existing values in the initial model.</li> </ul>	'zero'

To specify field values in Regularization, create a default nlgreyestOptions set and modify the fields using dot notation. Any fields that you do not modify retain their default values.

```
opt = nlgreyestOptions;
opt.Regularization.Lambda = 1.2;
opt.Regularization.R = 0.5*eye(np);
```

Regularization is a technique for specifying model flexibility constraints, which reduce uncertainty in the estimated parameter values. For more information, see “Regularized Estimates of Model Parameters”.

**SearchMethod — Numerical search method used for iterative parameter estimation**

```
'auto' (default) | 'gn' | 'gna' | 'lm' | 'grad' | 'lsqnonlin'
```

Numerical search method used for iterative parameter estimation, specified as the comma-separated pair consisting of 'SearchMethod' and one of the following:

- 'auto' — If Optimization Toolbox is available, 'lsqnonlin' is used. Otherwise, a combination of the line search algorithms, 'gn', 'lm', 'gna', and 'grad' methods

is tried in sequence at each iteration. The first descent direction leading to a reduction in estimation cost is used.

- 'gn' — Subspace Gauss-Newton least squares search. Singular values of the Jacobian matrix less than  $\text{GnPinvConstant} \cdot \text{eps} \cdot \max(\text{size}(J)) \cdot \text{norm}(J)$  are discarded when computing the search direction.  $J$  is the Jacobian matrix. The Hessian matrix is approximated by  $J^T J$ . If there is no improvement in this direction, the function tries the gradient direction.
- 'gna' — Adaptive subspace Gauss-Newton search. Eigenvalues less than  $\text{gamma} \cdot \max(\text{sv})$  of the Hessian are ignored, where  $\text{sv}$  are the singular values of the Hessian. The Gauss-Newton direction is computed in the remaining subspace.  $\text{gamma}$  has the initial value `InitialGnaTolerance` (see Advanced in 'SearchOptions' for more information). This value is increased by the factor `LMStep` each time the search fails to find a lower value of the criterion in fewer than five bisections. This value is decreased by the factor  $2 \cdot \text{LMStep}$  each time a search is successful without any bisections.
- 'lm' — Levenberg-Marquardt least squares search, where the next parameter value is  $-\text{pinv}(H+d \cdot I) \cdot \text{grad}$  from the previous one.  $H$  is the Hessian,  $I$  is the identity matrix, and  $\text{grad}$  is the gradient.  $d$  is a number that is increased until a lower value of the criterion is found.
- 'grad' — Steepest descent least squares search.
- 'lsqnonlin' — Trust-region-reflective algorithm of `lsqnonlin`. Requires Optimization Toolbox software.
- 'fmincon' — Constrained nonlinear solvers. You can use the sequential quadratic programming (SQP) and trust-region-reflective algorithms of the `fmincon` solver. If you have Optimization Toolbox software, you can also use the interior-point and active-set algorithms of the `fmincon` solver. Specify the algorithm in the `SearchOptions.Algorithm` option. The `fmincon` algorithms may result in improved estimation results in the following scenarios:
  - Constrained minimization problems when there are bounds imposed on the model parameters.
  - Model structures where the loss function is a nonlinear or non smooth function of the parameters.
  - Multi-output model estimation. A determinant loss function is minimized by default for MIMO model estimation. `fmincon` algorithms are able to minimize such loss functions directly. The other available search methods such as 'lm' and 'gn' minimize the determinant loss function by alternately estimating the noise variance and reducing the loss value for a given noise variance value. Hence, the `fmincon`

algorithms can offer better efficiency and accuracy for multi-output model estimations.

### **SearchOptions — Option set for the search algorithm**

search option set

Option set for the search algorithm, specified as the comma-separated pair consisting of 'SearchOptions' and a search option set with fields that depend on the value of SearchMethod.

### SearchOptions Structure When SearchMethod Is Specified as 'lsqnonlin' or 'auto', When Optimization Toolbox Is Available

Field Name	Description	Default
FunctionTolerance	Termination tolerance on the loss function that the software minimizes to determine the estimated parameter values, specified as a positive scalar.  The value of FunctionTolerance is the same as that of <code>opt.SearchOptions.Advanced.TolFun</code> .	1e-5
StepTolerance	Termination tolerance on the estimated parameter values, specified as a positive scalar.  The value of StepTolerance is the same as that of <code>opt.SearchOptions.Advanced.TolX</code> .	1e-6
MaxIterations	Maximum number of iterations during loss-function minimization, specified as a positive integer. The iterations stop when MaxIterations is reached or another stopping criterion is satisfied, such as FunctionTolerance.  The value of MaxIterations is the same as that of <code>opt.SearchOptions.Advanced.MaxIter</code> .	20
Advanced	Advanced search settings, specified as an option set for <code>lsqnonlin</code> .  For more information, see the Optimization Options table in "Optimization Options" (Optimization Toolbox).	Use <code>optimset('lsqnonlin')</code> to create a default option set.

**SearchOptions Structure When SearchMethod Is Specified as 'gn', 'gna', 'lm', 'grad', or 'auto', When Optimization Toolbox Is Not Available**

Field Name	Description	Default
Tolerance	Minimum percentage difference between the current value of the loss function and its expected improvement after the next iteration, specified as a positive scalar. When the percentage of expected improvement is less than <code>Tolerance</code> , the iterations stop. The estimate of the expected loss-function improvement at the next iteration is based on the Gauss-Newton vector computed for the current parameter value.	1e-5
MaxIterations	<p>Maximum number of iterations during loss-function minimization, specified as a positive integer. The iterations stop when <code>MaxIterations</code> is reached or another stopping criterion is satisfied, such as <code>Tolerance</code>.</p> <p>Setting <code>MaxIterations = 0</code> returns the result of the start-up procedure.</p> <p>Use <code>sys.Report.Termination.Iterations</code> to get the actual number of iterations during an estimation, where <code>sys</code> is an <code>idtf</code> model.</p>	20

Field Name	Description	Default	
Advanced	Advanced search settings, specified as a structure with the following fields:		
	Field Name	Description	Default
	GnPinvConstant	Jacobian matrix singular value threshold, specified as a positive scalar. Singular values of the Jacobian matrix that are smaller than $GnPinvConstant * \max(\text{size}(J) * \text{norm}(J) * \text{eps})$ are discarded when computing the search direction. Applicable when SearchMethod is 'gn'.	10000
	InitialGnaTolerance	Initial value of <i>gamma</i> , specified as a positive scalar. Applicable when SearchMethod is 'gna'.	0.0001
	LMStartValue	Starting value of search-direction length <i>d</i> in the Levenberg-Marquardt method, specified as a positive scalar. Applicable when SearchMethod is 'lm'.	0.001
	LMStep	Size of the Levenberg-Marquardt step, specified as a positive integer. The next value of the search-direction length <i>d</i> in the Levenberg-Marquardt method is LMStep times the previous one. Applicable when SearchMethod is 'lm'.	2
	MaxBisections	Maximum number of bisections used for line search along the search direction, specified as a positive integer.	25
	MaxFunctionEvaluations	Maximum number of calls to the model file, specified as a positive integer. Iterations stop if the number of calls to the model file exceeds this value.	Inf
MinParameterChange	Smallest parameter update allowed per iteration, specified as a nonnegative scalar.	0	

<b>Field Name</b>	<b>Description</b>		<b>Default</b>
	<b>Field Name</b>	<b>Description</b>	<b>Default</b>
	RelativeImprovement	Relative improvement threshold, specified as a nonnegative scalar. Iterations stop if the relative improvement of the criterion function is less than this value.	0
	StepReduction	Step reduction factor, specified as a positive scalar that is greater than 1. The suggested parameter update is reduced by the factor <code>StepReduction</code> after each try. This reduction continues until either <code>MaxBisections</code> tries are completed or a lower value of the criterion function is obtained.  StepReduction is not applicable for SearchMethod 'lm' (Levenberg-Marquardt method).	2

**SearchOptions Structure When SearchMethod is Specified as 'fmincon'**

Field Name	Description	Default
Algorithm	<p>fmincon optimization algorithm, specified as one of the following:</p> <ul style="list-style-type: none"> <li>• 'sqp' — Sequential quadratic programming algorithm. The algorithm satisfies bounds at all iterations, and it can recover from NaN or Inf results. It is not a large-scale algorithm. For more information, see “Large-Scale vs. Medium-Scale Algorithms” (Optimization Toolbox).</li> <li>• 'trust-region-reflective' — Subspace trust-region method based on the interior-reflective Newton method. It is a large-scale algorithm.</li> <li>• 'interior-point' — Large-scale algorithm that requires Optimization Toolbox software. The algorithm satisfies bounds at all iterations, and it can recover from NaN or Inf results.</li> <li>• 'active-set' — Requires Optimization Toolbox software. The</li> </ul>	'sqp'

Field Name	Description	Default
	<p>algorithm can take large steps, which adds speed. It is not a large-scale algorithm.</p> <p>For more information about the algorithms, see “Constrained Nonlinear Optimization Algorithms” (Optimization Toolbox) and “Choosing the Algorithm” (Optimization Toolbox).</p>	
FunctionTolerance	Termination tolerance on the loss function that the software minimizes to determine the estimated parameter values, specified as a positive scalar.	1e-6
StepTolerance	Termination tolerance on the estimated parameter values, specified as a positive scalar.	1e-6
MaxIterations	Maximum number of iterations during loss function minimization, specified as a positive integer. The iterations stop when MaxIterations is reached or another stopping criterion is satisfied, such as FunctionTolerance.	100

To specify field values in `SearchOptions`, create a default `nlgreyestOptions` set and modify the fields using dot notation. Any fields that you do not modify retain their default values.

```
opt = nlgreyestOptions('SearchMethod','gna');
opt.SearchOptions.MaxIterations = 50;
opt.SearchOptions.Advanced.RelImprovement = 0.5;
```

### **OutputWeight — Weighting of prediction error in multi-output estimations**

`[]` (default) | 'noise' | matrix

Weighting of prediction error in multi-output model estimations, specified as the comma-separated pair consisting of 'OutputWeight' and one of the following:

- `[]` — No weighting is used. Specifying as `[]` is the same as `eye(Ny)`, where `Ny` is the number of outputs.
- 'noise' — Optimal weighting is automatically computed as the inverse of the estimated noise variance. This weighting minimizes  $\det(E' * E / N)$ , where `E` is the matrix of prediction errors and `N` is the number of data samples. This option is not available when using 'lsqnonlin' as a 'SearchMethod'.
- A positive semidefinite matrix, `W`, of size equal to the number of outputs. This weighting minimizes  $\text{trace}(E' * E * W / N)$ , where `E` is the matrix of prediction errors and `N` is the number of data samples.

### **Advanced — Additional advanced options**

structure

Additional advanced options, specified as the comma-separated pair consisting of 'Advanced' and a structure with field:

Field Name	Description	Default
ErrorThreshold	Threshold for when to adjust the weight of large errors from quadratic to linear, specified as a nonnegative scalar. Errors larger than ErrorThreshold times the estimated standard deviation have a linear weight in the loss function. The standard deviation is estimated robustly as the median of the absolute deviations from the median of the prediction errors divided by 0.7. If your estimation data contains outliers, try setting ErrorThreshold to 1.6.	0 — Leads to a purely quadratic loss function.

To specify field values in `Advanced`, create a default `nlgreyestOptions` set and modify the fields using dot notation. Any fields that you do not modify retain their default values.

```
opt = nlgreyestOptions;  
opt.Advanced.ErrorThreshold = 1.2;
```

## Output Arguments

**opt** — Option set for `nlgreyest`

`nlgreyestOptions` option set

Option set for `nlgreyest`, returned as an `nlgreyestOptions` option set.

## See Also

`nlgreyest`

**Introduced in R2015a**

# nlhw

Estimate Hammerstein-Wiener model

## Syntax

```
sys = nlhw(Data,Orders)
sys = nlhw(Data,Orders,InputNL,OutputNL)

sys = nlhw(Data,LinModel)
sys = nlhw(Data,LinModel,InputNL,OutputNL)

sys = nlhw(Data,sys0)

sys = nlhw( ____,Options)
```

## Description

`sys = nlhw(Data,Orders)` creates and estimates a Hammerstein-Wiener model using the estimation data, model orders and delays, and default piecewise linear functions as input and output nonlinearity estimators.

`sys = nlhw(Data,Orders,InputNL,OutputNL)` specifies `InputNL` and `OutputNL` as the input and output nonlinearity estimators, respectively.

`sys = nlhw(Data,LinModel)` uses a linear model to specify the model orders and delays, and default piecewise linear functions for the input and output nonlinearity estimators.

`sys = nlhw(Data,LinModel,InputNL,OutputNL)` specifies `InputNL` and `OutputNL` as the input and output nonlinearity estimators, respectively.

`sys = nlhw(Data,sys0)` refines or estimates the parameters of a Hammerstein-Wiener model, `sys0`, using the estimation data.

Use this syntax to:

- Update the parameters of a previously estimated model to improve the fit to the estimation data. In this case, the estimation algorithm uses the parameters of `sys0` as initial guesses.
- Estimate the parameters of a model previously created using the `idnlhw` constructor. Prior to estimation, you can configure the model properties using dot notation.

`sys = nlhw( ____, Options)` specifies additional model estimation options. Use `Options` with any of the previous syntaxes.

## Examples

### Estimate a Hammerstein-Wiener Model

```
load iddata3
m1 = nlhw(z3,[4 2 1]);
```

### Estimate a Hammerstein Model with Saturation

Load data.

```
load twotankdata;
z = iddata(y,u,0.2,'Name','Two tank system');
z1 = z(1:1000);
```

Create a saturation object with lower limit of 0 and upper limit of 5.

```
InputNL = saturation('LinearInterval',[0 5]);
```

Estimate model with no output nonlinearity.

```
m = nlhw(z1,[2 3 0],InputNL,[]);
```

### Estimate Hammerstein-Wiener Model with a Custom Network Nonlinearity

Generating a custom network nonlinearity requires the definition of a user-defined unit function.

Define the unit function and save it as `gaussunit.m`.

```
% Copyright 2015 The MathWorks, Inc.
```

```
function [f, g, a] = gaussunit(x)
f = exp(-x.*x);
if nargin>1
    g = -2*x.*f;
    a = 0.2;
end
```

Create a custom network nonlinearity using the `gaussunit` function.

```
H = @gaussunit;
CNet = customnet(H);
```

Load the estimation data.

```
load twotankdata;
z = iddata(y,u,0.2,'Name','Two tank system');
z1 = z(1:1000);
```

Estimate a Hammerstein-Wiener model using the custom network.

```
m = nlhw(z1,[5 1 3],CNet,[]);
```

### Estimate Default Hammerstein-Wiener Model Using an Input-Output Polynomial Model of OE Structure

Estimate linear OE model.

```
load throttledata.mat
Tr = getTrend(ThrottleData);
Tr.OutputOffset = 15;
DetrendedData = detrend(ThrottleData, Tr);
opt = oeOptions('Focus','simulation');
LinearModel = oe(DetrendedData,[1 2 1],opt);
```

Estimate Hammerstein-Wiener model using OE model as its linear component and saturation as its output nonlinearity.

```
sys = nlhw(ThrottleData,LinearModel,[],'saturation');
```

## Estimate a Hammerstein-Wiener Model Using `idnlhw` to first Define the Model Properties

Load the estimation data.

```
load iddata1
```

Construct a Hammerstein-Wiener model using `idnlhw` to define the model properties B and F.

```
sys0 = idnlhw([2,2,0],[],'wavenet');  
sys0.B{1} = [0.8,1];  
sys0.F{1} = [1,-1.2,0.5];
```

Estimate the model.

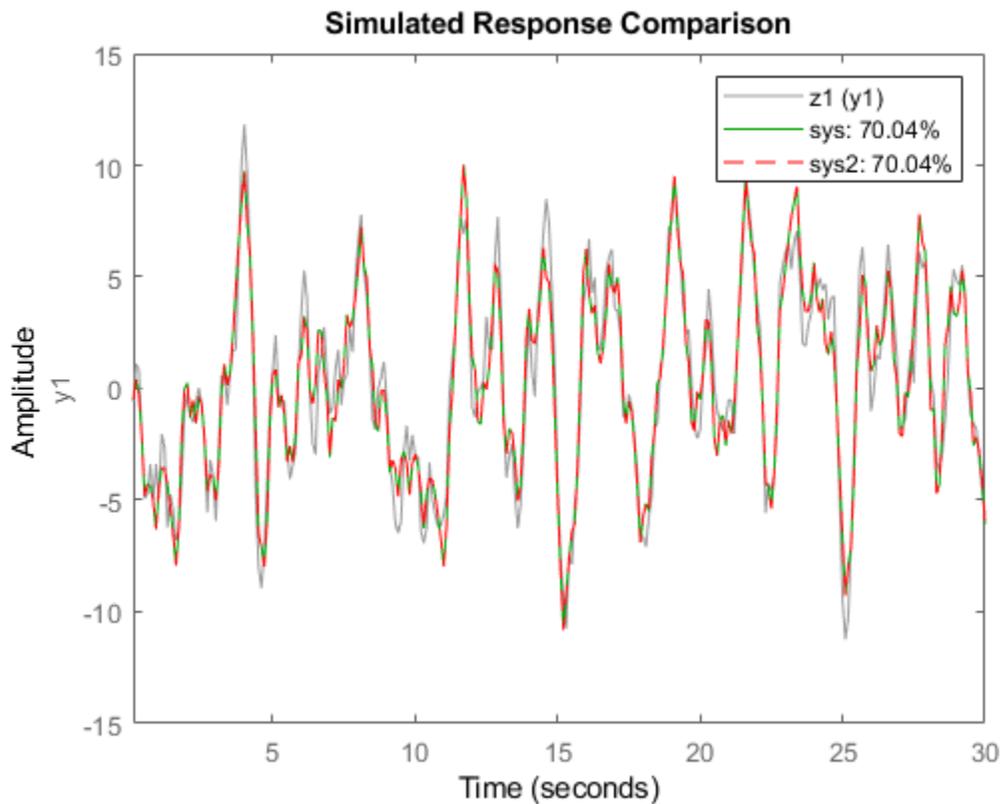
```
sys = nlhw(z1,sys0);
```

Estimate a Hammerstein-Wiener model using `nlhw` to define the model properties B and F.

```
sys2 = nlhw(z1,[2,2,0],[],'wavenet','B',{[0.8,1]},'F',{[1,-1.2,0.5]});
```

Compare the two estimated models to see that they are equivalent.

```
compare(z1,sys,'g',sys2,'r--');
```



### Refine a Hammerstein-Wiener Model Using Successive Calls of nlhw

Estimate a Hammerstein-Wiener Model.

```
load iddata3  
sys = nlhw(z3,[4 2 1], 'sigmoidnet', 'wavenet');
```

Refine the model, sys.

```
sys = nlhw(z3,sys);
```

## Estimate Hammerstein-Wiener Model Using an Estimation Option Set

Create estimation option set for `nlhw` to view estimation progress and to set the maximum iteration steps to 50.

```
opt = nlhwOptions;  
opt.Display = 'on';  
opt.SearchOptions.MaxIterations = 50;
```

Load data and estimate the model.

```
load iddata3  
sys = nlhw(z3,[4 2 1], 'sigmoidnet', 'deadzone', opt);
```

## Input Arguments

### Data — Time domain data

`iddata` object

Time-domain estimation data, specified as an `iddata`.

### Orders — Order and delays of the linear subsystem transfer function

`[nb nf nk]` vector of positive integers | `[nb nf nk]` vector of matrices

Order and delays of the linear subsystem transfer function, specified as a `[nb nf nk]` vector.

Dimensions of `Orders`:

- For a SISO transfer function, `Orders` is a vector of positive integers.

`nb` is the number of zeros plus 1, `nf` is the number of poles, and `nk` is the input delay.

- For a MIMO transfer function with  $n_u$  inputs and  $n_y$  outputs, `Orders` is a vector of matrices.

`nb`, `nf`, and `nk` are  $n_y$ -by- $n_u$  matrices whose  $i$ - $j$ th entry specifies the orders and delay of the transfer function from the  $j$ th input to the  $i$ th output.

## InputNL — Input static nonlinearity

'pwnlinear' (default) | 'sigmoidnet' | 'wavenet' | 'saturation' | 'deadzone' | 'poly1d' | 'unitgain' | nonlinearity estimator object | array of nonlinearity estimators

Input static nonlinearity estimator, specified as one of the following.

'pwnlinear' or pwnlinear object (default)	Piecewise linear function
'sigmoidnet' or sigmoidnet object	Sigmoid network
'wavenet' or wavenet object	Wavelet network
'saturation' or saturation object	Saturation
'deadzone' or deadzone object	Dead zone
'poly1d' or poly1d object	One-dimensional polynomial
'unitgain' or [] or unitgain object	Unit gain
customnet object	Custom network — Similar to <b>sigmoidnet</b> , but with a user-defined replacement for the sigmoid function.

Specifying a character vector, for example 'sigmoidnet', creates a nonlinearity estimator object with default settings. Use object representation instead to configure the properties of a nonlinearity estimator.

```
InputNL = wavenet;
InputNL.NumberOfUnits = 10;
```

Alternatively, use the associated input nonlinearity estimator function with Name-Value pair arguments.

```
InputNL = wavenet('NumberOfUnits',10);
```

For  $n_u$  input channels, you can specify nonlinear estimators individually for each input channel by setting InputNL to an  $n_u$ -by-1 array of nonlinearity estimators.

```
InputNL = [sigmoidnet('NumberOfUnits',5); deadzone([-1,2])]
```

To specify the same nonlinearity for all inputs, specify a single input nonlinearity estimator.

**OutputNL — Output static nonlinearity**

'pwlinear' (default) | 'sigmoidnet' | 'wavenet' | 'saturation' | 'deadzone' | 'poly1d' | 'unitgain' | nonlinearity estimator object | array of nonlinearity estimators

Output static nonlinearity estimator, specified as one of the following:

'pwlinear' or pwlinear object (default)	Piecewise linear function
'sigmoidnet' or sigmoidnet object	Sigmoid network
'wavenet' or wavenet object	Wavelet network
'saturation' or saturation object	Saturation
'deadzone' or deadzone object	Dead zone
'poly1d' or poly1d object	One-dimensional polynomial
'unitgain' or [] or unitgain object	Unit gain
customnet object	Custom network — Similar to <code>sigmoidnet</code> , but with a user-defined replacement for the sigmoid function.

Specifying a character vector creates a nonlinearity estimator object with default settings. Use object representation instead to configure the properties of a nonlinearity estimator.

```
OutputNL = sigmoidnet;  
OutputNL.NumberOfUnits = 10;
```

Alternatively, use the associated input nonlinearity estimator function with Name-Value pair arguments.

```
OutputNL = sigmoidnet('NumberOfUnits',10);
```

For  $n_y$  output channels, you can specify nonlinear estimators individually for each output channel by setting `OutputNL` to an  $n_y$ -by-1 array of nonlinearity estimators. To specify the same nonlinearity for all outputs, specify a single output nonlinearity estimator.

**LinModel — Discrete time linear model**

idpoly | idss with  $K = 0$  | idtf

Discrete-time linear model used to specify the linear subsystem, specified as one of the following:

- Input-output polynomial model of Output-Error (OE) structure (`idpoly`)
- State-space model with no disturbance component (`idss` with  $K = 0$ )
- Transfer function model (`idtf`)

Typically, you estimate the model using `oe`, `n4sid`, or `tfest`.

### **sys0 — Hammerstein-Wiener model**

`idnlhw` object

Hammerstein-Wiener model, specified as an `idnlhw` object. `sys0` can be:

- A model previously created using `idnlhw` to specify model properties.
- A model previously estimated using `nlhw`, that you want to update using a new estimation data set.

You can also refine `sys0` using the original estimation data set. If the previous estimation stopped when the numerical search was stuck at a local minima of the cost function, use `init` to first randomize the parameters of `sys0`. See `sys0.Report.Termination` for search stopping conditions. Using `init` does not guarantee a better solution on further refinement.

### **Options — Estimation options**

`nlhwOptions` option set

Estimation options for Hammerstein-Wiener model identification, specified as an `nlhwOptions` option set.

## **Output Arguments**

### **sys — Estimated Hammerstein-Wiener model**

`idnlhw` object

Estimated Hammerstein-Wiener model, returned as an `idnlhw` object. The model is estimated using the specified model orders and delays, input and output nonlinearity estimators, and estimation options.

Information about the estimation results and options used is stored in the `Report` property of the model. `Report` has the following fields:

<b>Report Field</b>	<b>Description</b>																		
Status	Summary of the model status, which indicates whether the model was created by construction or obtained by estimation.																		
Method	Estimation command used.																		
Fit	Quantitative assessment of the estimation, returned as a structure. See “Loss Function and Model Quality Metrics” for more information on these quality metrics. The structure has the following fields: <table border="1" data-bbox="387 574 1334 1072"> <thead> <tr> <th><b>Field</b></th> <th><b>Description</b></th> </tr> </thead> <tbody> <tr> <td>FitPercent</td> <td>Normalized root mean squared error (NRMSE) measure of how well the response of the model fits the estimation data, expressed as a percentage.</td> </tr> <tr> <td>LossFcn</td> <td>Value of the loss function when the estimation completes.</td> </tr> <tr> <td>MSE</td> <td>Mean squared error (MSE) measure of how well the response of the model fits the estimation data.</td> </tr> <tr> <td>FPE</td> <td>Final prediction error for the model.</td> </tr> <tr> <td>AIC</td> <td>Raw Akaike Information Criteria (AIC) measure of model quality.</td> </tr> <tr> <td>AICc</td> <td>Small sample-size corrected AIC.</td> </tr> <tr> <td>nAIC</td> <td>Normalized AIC.</td> </tr> <tr> <td>BIC</td> <td>Bayesian Information Criteria (BIC).</td> </tr> </tbody> </table>	<b>Field</b>	<b>Description</b>	FitPercent	Normalized root mean squared error (NRMSE) measure of how well the response of the model fits the estimation data, expressed as a percentage.	LossFcn	Value of the loss function when the estimation completes.	MSE	Mean squared error (MSE) measure of how well the response of the model fits the estimation data.	FPE	Final prediction error for the model.	AIC	Raw Akaike Information Criteria (AIC) measure of model quality.	AICc	Small sample-size corrected AIC.	nAIC	Normalized AIC.	BIC	Bayesian Information Criteria (BIC).
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BIC	Bayesian Information Criteria (BIC).																		
Parameters	Estimated values of model parameters.																		
OptionsUsed	Option set used for estimation. If no custom options were configured, this is a set of default options. See <code>nlhwOptions</code> for more information.																		
RandState	State of the random number stream at the start of estimation. Empty, [], if randomization was not used during estimation. For more information, see <code>rng</code> in the MATLAB documentation.																		

Report Field	Description																
DataUsed	Attributes of the data used for estimation, returned as a structure with the following fields:																
	<table border="1"> <thead> <tr> <th data-bbox="385 425 516 465">Field</th> <th data-bbox="516 425 1335 465">Description</th> </tr> </thead> <tbody> <tr> <td data-bbox="385 465 516 512">Name</td> <td data-bbox="516 465 1335 512">Name of the data set.</td> </tr> <tr> <td data-bbox="385 512 516 558">Type</td> <td data-bbox="516 512 1335 558">Data type.</td> </tr> <tr> <td data-bbox="385 558 516 605">Length</td> <td data-bbox="516 558 1335 605">Number of data samples.</td> </tr> <tr> <td data-bbox="385 605 516 652">Ts</td> <td data-bbox="516 605 1335 652">Sample time.</td> </tr> <tr> <td data-bbox="385 652 516 956">InterSample</td> <td data-bbox="516 652 1335 956">           Input intersample behavior, returned as one of the following values:           <ul style="list-style-type: none"> <li>• 'zoh' — Zero-order hold maintains a piecewise-constant input signal between samples.</li> <li>• 'foh' — First-order hold maintains a piecewise-linear input signal between samples.</li> <li>• 'bl' — Band-limited behavior specifies that the continuous-time input signal has zero power above the Nyquist frequency.</li> </ul> </td> </tr> <tr> <td data-bbox="385 956 516 1038">InputOffset</td> <td data-bbox="516 956 1335 1038">Offset removed from time-domain input data during estimation. For nonlinear models, it is [ ].</td> </tr> <tr> <td data-bbox="385 1038 516 1111">OutputOffset</td> <td data-bbox="516 1038 1335 1111">Offset removed from time-domain output data during estimation. For nonlinear models, it is [ ].</td> </tr> </tbody> </table>	Field	Description	Name	Name of the data set.	Type	Data type.	Length	Number of data samples.	Ts	Sample time.	InterSample	Input intersample behavior, returned as one of the following values: <ul style="list-style-type: none"> <li>• 'zoh' — Zero-order hold maintains a piecewise-constant input signal between samples.</li> <li>• 'foh' — First-order hold maintains a piecewise-linear input signal between samples.</li> <li>• 'bl' — Band-limited behavior specifies that the continuous-time input signal has zero power above the Nyquist frequency.</li> </ul>	InputOffset	Offset removed from time-domain input data during estimation. For nonlinear models, it is [ ].	OutputOffset	Offset removed from time-domain output data during estimation. For nonlinear models, it is [ ].
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OutputOffset	Offset removed from time-domain output data during estimation. For nonlinear models, it is [ ].																

Report Field	Description																
Termination	Termination conditions for the iterative search used for prediction error minimization. Structure with the following fields:																
	<table border="1"> <thead> <tr> <th data-bbox="387 427 516 465">Field</th> <th data-bbox="516 427 1335 465">Description</th> </tr> </thead> <tbody> <tr> <td data-bbox="387 465 516 512">WhyStop</td> <td data-bbox="516 465 1335 512">Reason for terminating the numerical search.</td> </tr> <tr> <td data-bbox="387 512 516 586">Iterations</td> <td data-bbox="516 512 1335 586">Number of search iterations performed by the estimation algorithm.</td> </tr> <tr> <td data-bbox="387 586 516 690">FirstOrderOptimality</td> <td data-bbox="516 586 1335 690"><math>\infty</math>-norm of the gradient search vector when the search algorithm terminates.</td> </tr> <tr> <td data-bbox="387 690 516 765">FcnCount</td> <td data-bbox="516 690 1335 765">Number of times the objective function was called.</td> </tr> <tr> <td data-bbox="387 765 516 840">UpdateNorm</td> <td data-bbox="516 765 1335 840">Norm of the gradient search vector in the last iteration. Omitted when the search method is 'lsqnonlin' or 'fmincon'.</td> </tr> <tr> <td data-bbox="387 840 516 944">LastImprovement</td> <td data-bbox="516 840 1335 944">Criterion improvement in the last iteration, expressed as a percentage. Omitted when the search method is 'lsqnonlin' or 'fmincon'.</td> </tr> <tr> <td data-bbox="387 944 516 1031">Algorithm</td> <td data-bbox="516 944 1335 1031">Algorithm used by 'lsqnonlin' or 'fmincon' search method. Omitted when other search methods are used.</td> </tr> </tbody> </table>	Field	Description	WhyStop	Reason for terminating the numerical search.	Iterations	Number of search iterations performed by the estimation algorithm.	FirstOrderOptimality	$\infty$ -norm of the gradient search vector when the search algorithm terminates.	FcnCount	Number of times the objective function was called.	UpdateNorm	Norm of the gradient search vector in the last iteration. Omitted when the search method is 'lsqnonlin' or 'fmincon'.	LastImprovement	Criterion improvement in the last iteration, expressed as a percentage. Omitted when the search method is 'lsqnonlin' or 'fmincon'.	Algorithm	Algorithm used by 'lsqnonlin' or 'fmincon' search method. Omitted when other search methods are used.
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For estimation methods that do not require numerical search optimization, the Termination field is omitted.																	

For more information, see “Estimation Report”.

## Extended Capabilities

### Automatic Parallel Support

Accelerate code by automatically running computation in parallel using Parallel Computing Toolbox™.

Parallel computing support is available for estimation using the `lsqnonlin` search method (requires Optimization Toolbox). To enable parallel computing, use `nlhwOptions`, set `SearchMethod` to `'lsqnonlin'`, and set `SearchOptions.Advanced.UseParallel` to `true`.

For example:

```
opt = nlhwOptions;  
opt.SearchMethod = 'lsqnonlin';  
opt.SearchOptions.Advanced.UseParallel = true;
```

## See Also

`aic` | `customnet` | `deadzone` | `fpe` | `goodnessofFit` | `idnlhw` | `idnlhw/findop` | `init` | `linapp` | `linearize` | `n4sid` | `nlhwOptions` | `oe` | `pem` | `poly1d` | `pwlinear` | `saturation` | `sigmoidnet` | `tfest` | `unitgain` | `wavenet`

## Topics

“Estimate Multiple Hammerstein-Wiener Models”

“Estimate Hammerstein-Wiener Models Initialized Using Linear OE Models”

“Identifying Hammerstein-Wiener Models”

“Initialize Hammerstein-Wiener Estimation Using Linear Model”

“Loss Function and Model Quality Metrics”

“Regularized Estimates of Model Parameters”

“Estimation Report”

## Introduced in R2007a

## nlhwOptions

Option set for nlhw

### Syntax

```
opt = nlhwOptions  
opt = nlhwOptions(Name,Value)
```

### Description

`opt = nlhwOptions` creates the default option set for `nlhw`. Use dot notation to customize the option set, if needed.

`opt = nlhwOptions(Name,Value)` creates an option set with options specified by one or more `Name,Value` pair arguments. The options that you do not specify retain their default value.

### Examples

#### Estimate Hammerstein-Wiener Model Using an Estimation Option Set

Create estimation option set for `nlhw` to view estimation progress and to set the maximum iteration steps to 50.

```
opt = nlhwOptions;  
opt.Display = 'on';  
opt.SearchOptions.MaxIterations = 50;
```

Load data and estimate the model.

```
load iddata3  
sys = nlhw(z3,[4 2 1], 'sigmoidnet', 'deadzone', opt);
```

## Specify an Option Set for Hammerstein-Wiener Model Estimation

Create an options set for `nlhw` where:

- Initial conditions are estimated from the estimation data.
- Subspace Gauss-Newton least squares method is used for estimation.

```
opt = nlhwOptions('InitialCondition','estimate','SearchMethod','gn');
```

## Input Arguments

### Name-Value Pair Arguments

Specify optional comma-separated pairs of `Name`, `Value` arguments. `Name` is the argument name and `Value` is the corresponding value. `Name` must appear inside quotes. You can specify several name and value pair arguments in any order as `Name1,Value1,...,NameN,ValueN`.

Example: `nlhwOptions('InitialCondition','estimate')`

### InitialCondition — Handling of initial conditions

'zero' (default) | 'estimate'

Handling of initial conditions during estimation using `nlhw`, specified as the comma-separated pair consisting of `InitialCondition` and one of the following:

- 'zero' — The initial conditions are set to zero.
- 'estimate' — The initial conditions are treated as independent estimation parameters.

### Display — Estimation progress display setting

'off' (default) | 'on'

Estimation progress display setting, specified as the comma-separated pair consisting of `Display` and one of the following:

- 'off' — No progress or results information is displayed.
- 'on' — Information on model structure and estimation results are displayed in a progress-viewer window.

**OutputWeight — Weighting of prediction error in multi-output estimations**

'noise' (default) | positive semidefinite matrix

Weighting of prediction error in multi-output model estimations, specified as the comma-separated pair consisting of 'OutputWeight' and one of the following:

- 'noise' — Optimal weighting is automatically computed as the inverse of the estimated noise variance. This weighting minimizes  $\det(E' * E)$ , where E is the matrix of prediction errors. This option is not available when using 'lsqnonlin' as a 'SearchMethod'.
- A positive semidefinite matrix, W, of size equal to the number of outputs. This weighting minimizes  $\text{trace}(E' * E * W / N)$ , where E is the matrix of prediction errors and N is the number of data samples.

**Regularization — Options for regularized estimation of model parameters**

structure

Options for regularized estimation of model parameters, specified as the comma-separated pair consisting of 'Regularization' and a structure with fields:

Field Name	Description	Default
Lambda	Bias versus variance trade-off constant, specified as a nonnegative scalar.	0 — Indicates no regularization.
R	Weighting matrix, specified as a vector of nonnegative scalars or a square positive semi-definite matrix. The length must be equal to the number of free parameters in the model, np. Use the nparams command to determine the number of model parameters.	1 — Indicates a value of eye(np).
Nominal	The nominal value towards which the free parameters are pulled during estimation, specified as one of the following: <ul style="list-style-type: none"> <li>• 'zero' — Pull parameters towards zero.</li> <li>• 'model' — Pull parameters towards pre-existing values in the initial model. Use this option only when you have a well-initialized idnlhw model with finite parameter values.</li> </ul>	'zero'

To specify field values in `Regularization`, create a default `nlhwOptions` set and modify the fields using dot notation. Any fields that you do not modify retain their default values.

```
opt = nlhwOptions;
opt.Regularization.Lambda = 1.2;
opt.Regularization.R = 0.5*eye(np);
```

Regularization is a technique for specifying model flexibility constraints, which reduce uncertainty in the estimated parameter values. For more information, see “Regularized Estimates of Model Parameters”.

### **SearchMethod — Numerical search method used for iterative parameter estimation**

'auto' (default) | 'gn' | 'gna' | 'lm' | 'grad' | 'lsqnonlin' | 'fmincon'

Numerical search method used for iterative parameter estimation, specified as the comma-separated pair consisting of 'SearchMethod' and one of the following:

- 'auto' — A combination of the line search algorithms, 'gn', 'lm', 'gna', and 'grad' methods is tried in sequence at each iteration. The first descent direction leading to a reduction in estimation cost is used.
- 'gn' — Subspace Gauss-Newton least squares search. Singular values of the Jacobian matrix less than  $GnPinvConstant * \epsilon * \max(\text{size}(J)) * \text{norm}(J)$  are discarded when computing the search direction.  $J$  is the Jacobian matrix. The Hessian matrix is approximated as  $J^T J$ . If there is no improvement in this direction, the function tries the gradient direction.
- 'gna' — Adaptive subspace Gauss-Newton search. Eigenvalues less than  $\gamma * \max(sv)$  of the Hessian are ignored, where  $sv$  contains the singular values of the Hessian. The Gauss-Newton direction is computed in the remaining subspace.  $\gamma$  has the initial value `InitialGnaTolerance` (see `Advanced in 'SearchOptions'` for more information). This value is increased by the factor `LMStep` each time the search fails to find a lower value of the criterion in fewer than five bisections. This value is decreased by the factor  $2 * LMStep$  each time a search is successful without any bisections.
- 'lm' — Levenberg-Marquardt least squares search, where the next parameter value is  $-\text{pinv}(H+d*I) * \text{grad}$  from the previous one.  $H$  is the Hessian,  $I$  is the identity matrix, and  $\text{grad}$  is the gradient.  $d$  is a number that is increased until a lower value of the criterion is found.
- 'grad' — Steepest descent least squares search.

- 'lsqnonlin' — Trust-region-reflective algorithm of lsqnonlin. Requires Optimization Toolbox software.
- 'fmincon' — Constrained nonlinear solvers. You can use the sequential quadratic programming (SQP) and trust-region-reflective algorithms of the fmincon solver. If you have Optimization Toolbox software, you can also use the interior-point and active-set algorithms of the fmincon solver. Specify the algorithm in the SearchOptions.Algorithm option. The fmincon algorithms may result in improved estimation results in the following scenarios:
  - Constrained minimization problems when there are bounds imposed on the model parameters.
  - Model structures where the loss function is a nonlinear or non smooth function of the parameters.
  - Multi-output model estimation. A determinant loss function is minimized by default for multi-output model estimation. fmincon algorithms are able to minimize such loss functions directly. The other search methods such as 'lm' and 'gn' minimize the determinant loss function by alternately estimating the noise variance and reducing the loss value for a given noise variance value. Hence, the fmincon algorithms can offer better efficiency and accuracy for multi-output model estimations.

### **SearchOptions — Option set for the search algorithm**

search option set

Option set for the search algorithm, specified as the comma-separated pair consisting of 'SearchOptions' and a search option set with fields that depend on the value of SearchMethod.

**SearchOptions Structure When SearchMethod is Specified as 'gn', 'gna', 'lm', 'grad', or 'auto'**

Field Name	Description	Default
Tolerance	Minimum percentage difference between the current value of the loss function and its expected improvement after the next iteration, specified as a positive scalar. When the percentage of expected improvement is less than <code>Tolerance</code> , the iterations stop. The estimate of the expected loss-function improvement at the next iteration is based on the Gauss-Newton vector computed for the current parameter value.	1e-5
MaxIterations	<p>Maximum number of iterations during loss-function minimization, specified as a positive integer. The iterations stop when <code>MaxIterations</code> is reached or another stopping criterion is satisfied, such as <code>Tolerance</code>.</p> <p>Setting <code>MaxIterations = 0</code> returns the result of the start-up procedure.</p> <p>Use <code>sys.Report.Termination.Iterations</code> to get the actual number of iterations during an estimation, where <code>sys</code> is an <code>idtf</code> model.</p>	20

Field Name	Description	Default
Advanced	Advanced search settings, specified as a structure with the following fields:	
<b>Field Name</b>	<b>Description</b>	<b>Default</b>
GnPinvConstant	Jacobian matrix singular value threshold, specified as a positive scalar. Singular values of the Jacobian matrix that are smaller than $\text{GnPinvConstant} * \max(\text{size}(J) * \text{norm}(J) * \text{eps})$ are discarded when computing the search direction. Applicable when SearchMethod is 'gn'.	10000
InitialGnaTolerance	Initial value of $\gamma$ , specified as a positive scalar. Applicable when SearchMethod is 'gna'.	0.0001
LMStartValue	Starting value of search-direction length $d$ in the Levenberg-Marquardt method, specified as a positive scalar. Applicable when SearchMethod is 'lm'.	0.001
LMStep	Size of the Levenberg-Marquardt step, specified as a positive integer. The next value of the search-direction length $d$ in the Levenberg-Marquardt method is LMStep times the previous one. Applicable when SearchMethod is 'lm'.	2
MaxBisections	Maximum number of bisections used for line search along the search direction, specified as a positive integer.	25
MaxFunctionEvaluations	Maximum number of calls to the model file, specified as a positive integer. Iterations stop if the number of calls to the model file exceeds this value.	Inf
MinParameterChange	Smallest parameter update allowed per iteration, specified as a nonnegative scalar.	0
RelativeImprovement	Relative improvement threshold, specified as a nonnegative scalar. Iterations stop if the relative	0

Field Name	Description		Default
	<b>Field Name</b>	<b>Description</b>	<b>Default</b>
		improvement of the criterion function is less than this value.	
	StepReduction	<p>Step reduction factor, specified as a positive scalar that is greater than 1. The suggested parameter update is reduced by the factor <code>StepReduction</code> after each try. This reduction continues until <code>MaxBisections</code> tries are completed or a lower value of the criterion function is obtained.</p> <p><code>StepReduction</code> is not applicable for <code>SearchMethod 'lm'</code> (Levenberg-Marquardt method).</p>	2

**SearchOptions Structure When SearchMethod is Specified as 'lsqnonlin'**

<b>Field Name</b>	<b>Description</b>	<b>Default</b>
FunctionTolerance	Termination tolerance on the loss function that the software minimizes to determine the estimated parameter values, specified as a positive scalar.  The value of FunctionTolerance is the same as that of <code>opt.SearchOptions.Advanced.TolFun</code> .	1e-5
StepTolerance	Termination tolerance on the estimated parameter values, specified as a positive scalar.  The value of StepTolerance is the same as that of <code>opt.SearchOptions.Advanced.TolX</code> .	1e-6
MaxIterations	Maximum number of iterations during loss-function minimization, specified as a positive integer. The iterations stop when MaxIterations is reached or another stopping criterion is satisfied, such as FunctionTolerance.  The value of MaxIterations is the same as that of <code>opt.SearchOptions.Advanced.MaxIter</code> .	20
Advanced	Advanced search settings, specified as an option set for <code>lsqnonlin</code> .  For more information, see the Optimization Options table in "Optimization Options" (Optimization Toolbox).	Use <code>optimset('lsqnonlin')</code> to create a default option set.

### SearchOptions Structure When SearchMethod is Specified as 'fmincon'

Field Name	Description	Default
Algorithm	<p>fmincon optimization algorithm, specified as one of the following:</p> <ul style="list-style-type: none"> <li>'sqp' — Sequential quadratic programming algorithm. The algorithm satisfies bounds at all iterations, and it can recover from NaN or Inf results. It is not a large-scale algorithm. For more information, see “Large-Scale vs. Medium-Scale Algorithms” (Optimization Toolbox).</li> <li>'trust-region-reflective' — Subspace trust-region method based on the interior-reflective Newton method. It is a large-scale algorithm.</li> <li>'interior-point' — Large-scale algorithm that requires Optimization Toolbox software. The algorithm satisfies bounds at all iterations, and it can recover from NaN or Inf results.</li> <li>'active-set' — Requires Optimization Toolbox software. The</li> </ul>	'sqp'

Field Name	Description	Default
	<p>algorithm can take large steps, which adds speed. It is not a large-scale algorithm.</p> <p>For more information about the algorithms, see “Constrained Nonlinear Optimization Algorithms” (Optimization Toolbox) and “Choosing the Algorithm” (Optimization Toolbox).</p>	
FunctionTolerance	Termination tolerance on the loss function that the software minimizes to determine the estimated parameter values, specified as a positive scalar.	1e-6
StepTolerance	Termination tolerance on the estimated parameter values, specified as a positive scalar.	1e-6
MaxIterations	Maximum number of iterations during loss function minimization, specified as a positive integer. The iterations stop when MaxIterations is reached or another stopping criterion is satisfied, such as FunctionTolerance.	100

To specify field values in `SearchOptions`, create a default `nllwOptions` set and modify the fields that you do not modify retain their default values.

```
opt = nllwOptions;
opt.SearchOptions.MaxIterations = 50;
opt.SearchOptions.Advanced.RelImprovement = 0.5;
```

## Advanced — Additional advanced options

structure

Additional advanced options, specified as the comma-separated pair consisting of 'Advanced' and a structure with fields:

Field Name	Description	Default
ErrorThreshold	Threshold for when to adjust the weight of large errors from quadratic to linear, specified as a nonnegative scalar. Errors larger than ErrorThreshold times the estimated standard deviation have a linear weight in the loss function. The standard deviation is estimated robustly as the median of the absolute deviations from the median of the prediction errors, divided by 0.7. If your estimation data contains outliers, try setting ErrorThreshold to 1.6.	0 — Leads to a purely quadratic loss function.
MaxSize	Maximum number of elements in a segment when input-output data is split into segments, specified as a positive integer.	250000

To specify field values in Advanced, create a default nlhwOptions set and modify the fields using dot notation. Any fields that you do not modify retain their default values.

```
opt = nlhwOptions;
opt.Advanced.ErrorThreshold = 1.2;
```

## Output Arguments

**opt** — Option set for nlhw

nlhwOptions option set

Option set for nlhw, returned as an nlhwOptions option set.

## See Also

nlhw

**Introduced in R2015a**

# noise2meas

Noise component of model

## Syntax

```
noise_model = noise2meas(sys)
noise_model = noise2meas(sys,noise)
```

## Description

`noise_model = noise2meas(sys)` returns the noise component, `noise_model`, of a linear identified model, `sys`. Use `noise2meas` to convert a time-series model (no inputs) to an input/output model. The converted model can be used for linear analysis, including viewing pole/zero maps, and plotting the step response.

`noise_model = noise2meas(sys,noise)` specifies the noise variance normalization method.

## Input Arguments

### **sys**

Identified linear model.

### **noise**

Noise variance normalization method, specified as one of the following values:

- 'innovations' — Noise sources are not normalized and remain as the innovations process.
- 'normalize' — Noise sources are normalized to be independent and of unit variance.

**Default:** 'innovations'

## Output Arguments

### `noise_model`

Noise component of `sys`.

`sys` represents the system

$$y(t) = Gu(t) + He(t)$$

$G$  is the transfer function between the measured input,  $u(t)$ , and the output,  $y(t)$ .  $H$  is the noise model and describes the effect of the disturbance,  $e(t)$ , on the model's response.

An equivalent state-space representation of `sys` is

$$\dot{x}(t) = Ax(t) + Bu(t) + Ke(t)$$

$$y(t) = Cx(t) + Du(t) + e(t)$$

$$e(t) = Lv(t)$$

$v(t)$  is white noise with independent channels and unit variances. The white-noise signal  $e(t)$  represents the model's innovations and has variance  $LL^T$ . The noise-variance data is stored using the `NoiseVariance` property of `sys`.

- If `noise` is 'innovations', then `noise2meas` returns  $H$  and `noise_model` represents the system

$$y(t) = He(t)$$

An equivalent state-space representation of `noise_model` is

$$\dot{x}(t) = Ax(t) + Ke(t)$$

$$y(t) = Cx(t) + e(t)$$

`noise2meas` returns the noise channels of `sys` as the input channels of `noise_model`. The input channels are named using the format 'e@yk', where `yk` corresponds to the `OutputName` property of an output. The measured input channels of `sys` are discarded and the noise variance is set to zero.

- If `noise` is 'normalize', then `noise2meas` first normalizes

$$e(t) = Lv(t)$$

`noise_model` represents the system

$$y(t) = HLv(t)$$

or, equivalently, in state-space representation

$$\dot{x}(t) = Ax(t) + KLv(t)$$

$$y(t) = Cx(t) + Lv(t)$$

The input channels are named using the format 'v@yk', where yk corresponds to the `OutputName` property of an output.

The model type of `noise_model` depends on the model type of `sys`.

- `noise_model` is an `idtf` model if `sys` is an `idproc` model.
- `noise_model` is an `idss` model if `sys` is an `idgrey` model.
- `noise_model` is the same type of model as `sys` for all other model types.

To obtain the model coefficients of `noise_model` in state-space form, use `ssdata`. Similarly, to obtain the model coefficients in transfer-function form, use `tfddata`.

## Examples

### Convert Noise Component of Linear Identified Model into Input/Output Model

Convert a time-series model to an input/output model that may be used by linear analysis tools.

Identify a time-series model.

```
load iddata9 z9
sys = ar(z9,4,'ls');
```

`sys` is an `idpoly` model with no inputs.

Convert `sys` to a measured model.

```
noise_model = noise2meas(sys);
```

`noise_model` is an `idpoly` model with one input.

You can use `noise_model` for linear analysis functions such as `step`, `iopzmap`, etc.

## Normalizing Noise Variance

Convert an identified linear model to an input/output model, and normalize its noise variance.

Identify a linear model using data.

```
load twotankdata;  
z = iddata(y,u,0.2);  
sys = ssest(z,4);
```

`sys` is an `idss` model, with a noise variance of  $6.6211e-06$ . The value of  $L$  is `sqrt(sys.NoiseVariance)`, which is 0.0026.

View the disturbance matrix.

```
sys.K
```

```
ans = 4×1  
  
    0.2719  
    1.6570  
   -0.6318  
   -0.2877
```

Obtain a model that absorbs the noise variance of `sys`.

```
noise_model_normalize = noise2meas(sys,'normalize');
```

`noise_model_normalize` is an `idpoly` model.

View the  $B$  matrix for `noise_model_normalize`

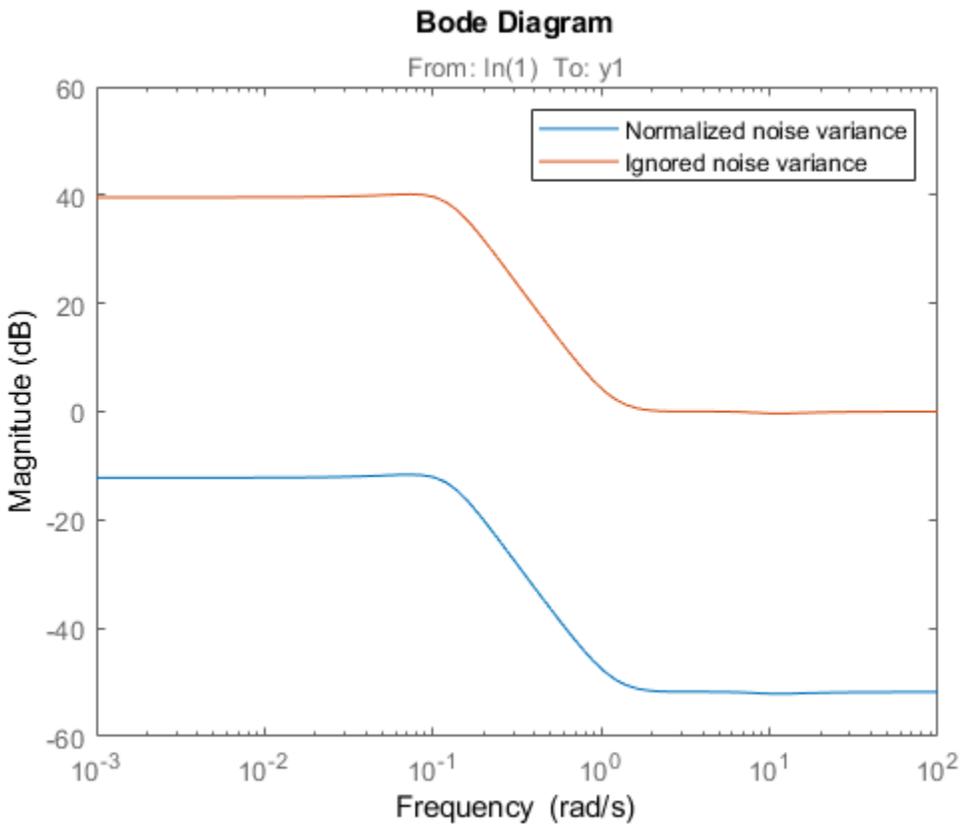
```
noise_model_normalize.B
```

```
ans = 4×1  
  
    0.0007  
    0.0043  
   -0.0016  
   -0.0007
```

As expected, `noise_model_normalize.B` is equal to  $L \cdot \text{sys} \cdot K$ .

Compare the bode response with a model that ignores the noise variance of `sys`.

```
noise_model_innovation = noise2meas(sys,'innovations');  
bodemag(noise_model_normalize,noise_model_innovation);  
legend('Normalized noise variance','Ignored noise variance');
```



The difference between the bode magnitudes of the `noise_model_innovation` and `noise_model_normalized` is approximately 51 dB. As expected, the magnitude difference is approximately equal to  $20 \cdot \log_{10}(L)$ .

## See Also

`idssdata` | `noisecnv` | `spectrum` | `tfdata` | `zpkdata`

**Introduced in R2012a**

## noisecnv

Transform identified linear model with noise channels to model with measured channels only

### Syntax

```
mod1 = noisecnv(mod)
mod2 = noisecnv(mod, 'normalize')
```

### Description

`mod1 = noisecnv(mod)` and `mod2 = noisecnv(mod, 'normalize')` transform an identified linear model with noise channels to a model with measured channels only.

`mod` is any linear identified model, `idproc`, `idtf`, `idgrey`, `idpoly`, or `idss`.

The noise input channels in `mod` are converted as follows: Consider a model with both measured input channels  $u$  ( $nu$  channels) and noise channels  $e$  ( $ny$  channels) with covariance matrix  $\Lambda$ :

$$y = Gu + He$$

$$\text{cov}(e) = \Lambda = LL'$$

where  $L$  is a lower triangular matrix. Note that `mod.NoiseVariance =  $\Lambda$` . The model can also be described with unit variance, using a normalized noise source  $v$ :

$$y = Gu + HLv$$

$$\text{cov}(v) = I$$

- `mod1 = noisecnv(mod)` converts the model to a representation of the system  $[G H]$  with  $nu+ny$  inputs and  $ny$  outputs. All inputs are treated as measured, and `mod1` does not have any noise model. The former noise input channels have names `e@yname`, where `yname` is the name of the corresponding output.
- `mod2 = noisecnv(mod, 'norm')` converts the model to a representation of the system  $[G HL]$  with  $nu+ny$  inputs and  $ny$  outputs. All inputs are treated as measured,

and `mod2` does not have any noise model. The former noise input channels have names `v@yname`, where `yname` is the name of the corresponding output. Note that the noise variance matrix factor  $L$  typically is uncertain (has a nonzero covariance). This is taken into account in the uncertainty description of `mod2`.

- If `mod` is a time series, that is,  $nu = 0$ , `mod1` is a model that describes the transfer function  $H$  with measured input channels. Analogously, `mod2` describes the transfer function  $HL$ .

Note the difference with subreferencing:

- `mod(:, [])` gives a description of the noise model characteristics as a time-series model, that is, it describes  $H$  and also the covariance of  $e$ . In contrast, `noisecnv(m(:, []))` or `noise2meas(m)` describe just the transfer function  $H$ . To obtain a description of the normalized transfer function  $HL$ , use `noisecnv(m(:, []), 'normalize')` or `noise2meas('normalize')`.

Converting the noise channels to measured inputs is useful to study the properties of the individual transfer functions from noise to output. It is also useful for transforming identified linear models to representations that do not handle disturbance descriptions explicitly.

## Examples

Identify a model with a measured component ( $G$ ) and a non-trivial noise component ( $H$ ). Compare the amplitude of the measured component's frequency response to the noise component's spectrum amplitude. You must convert the noise component into a measured one by using `noisecnv` if you want to compare its behavior against a truly measured component.

```
load iddata2 z2
sys1 = armax(z2,[2 2 2 1]); % model with noise component
sys2 = tfest(z2,3); % model with a trivial noise component

sys1 = noisecnv(sys1);
sys2 = noisecnv(sys2);
bodemag(sys1,sys2)
```

## See Also

`idssdata` | `noise2meas` | `tfdata` | `zpkdata`

## **Topics**

“Treating Noise Channels as Measured Inputs”

**Introduced before R2006a**

## norm

Norm of linear model

### Syntax

```
n = norm(sys)
n = norm(sys,2)

n = norm(sys,Inf)
[n, fpeak] = norm(sys,Inf)
[n, fpeak] = norm(sys,Inf, tol)
```

### Description

`n = norm(sys)` or `n = norm(sys,2)` returns the root-mean-squares of the impulse response of the linear dynamic system model `sys`. This value is equivalent to the  $H_2$  norm on page 1-1036 of `sys`.

`n = norm(sys,Inf)` returns the  $L_\infty$  norm (Control System Toolbox) of `sys`, which is the peak gain of the frequency response of `sys` across frequencies. For MIMO systems, this quantity is the peak gain over all frequencies and all input directions, which corresponds to the peak value of the largest singular value of `sys`. For stable systems, the  $L_\infty$  norm is equivalent to the  $H_\infty$  norm. For more information, see `hinfnorm`.

`[n, fpeak] = norm(sys,Inf)` also returns the frequency `fpeak` at which the gain reaches its peak value.

`[n, fpeak] = norm(sys,Inf, tol)` sets the relative accuracy of the  $L_\infty$  norm to `tol`.

This command requires a Control System Toolbox license.

### Examples

## Compute Norm of Discrete-Time Linear System

Compute the  $H_2$  and  $L_\infty$  norms of the following discrete-time transfer function, with sample time 0.1 second.

$$\text{sys}(z) = \frac{z^3 - 2.841z^2 + 2.875z - 1.004}{z^3 - 2.417z^2 + 2.003z - 0.5488}$$

Compute the  $H_2$  norm of the transfer function. The  $H_2$  norm is the root-mean-square of the impulse response of `sys`.

```
sys = tf([1 -2.841 2.875 -1.004],[1 -2.417 2.003 -0.5488],0.1);
n2 = norm(sys)
```

```
n2 = 1.2438
```

Compute the  $L_\infty$  norm of the transfer function.

```
[ninf, fpeak] = norm(sys, Inf)
```

```
ninf = 2.5715
```

```
fpeak = 3.0051
```

Because `sys` is a stable system, `ninf` is the peak gain of the frequency response of `sys`, and `fpeak` is the frequency at which the peak gain occurs. Confirm these values using `getPeakGain`.

```
[gpeak, fpeak] = getPeakGain(sys)
```

```
gpeak = 2.5715
```

```
fpeak = 3.0051
```

## Input Arguments

### **sys** — Dynamic system

dynamic system model | model array

Input dynamic system, specified as any SISO or MIMO linear dynamic system model or model array. `sys` can be continuous-time or discrete-time.

### **tol — Relative accuracy**

0.01 (default) | positive real scalar

Relative accuracy of the  $H_\infty$  norm, specified as a positive real scalar value.

## Output Arguments

### **n — $H_2$ or $L_\infty$ norm**

scalar | array

$H_2$  norm or  $L_\infty$  norm of `sys`, returned as a scalar or an array.

- If `sys` is a single model, then `n` is a scalar value.
- If `sys` is a model array, then `n` is an array of the same size as `sys`, where `n(k) = norm(sys(:, :, k))`.

### **fpeak — Frequency of peak gain**

nonnegative real scalar | array of nonnegative real values

Frequency at which the gain achieves the peak value `gpeak`, returned as a nonnegative real scalar value or an array of nonnegative real values. The frequency is expressed in units of rad/TimeUnit, relative to the `TimeUnit` property of `sys`.

- If `sys` is a single model, then `fpeak` is a scalar.
- If `sys` is a model array, then `fpeak` is an array of the same size as `sys`, where `fpeak(k)` is the peak gain frequency of `sys(:, :, k)`.

## Definitions

### **$H_2$ norm**

The  $H_2$  norm of a stable system  $H$  is the root-mean-square of the impulse response of the system. The  $H_2$  norm measures the steady-state covariance (or power) of the output response  $y = Hw$  to unit white noise inputs  $w$ :

$$\|H\|_2^2 = \lim_{t \rightarrow \infty} E\{y(t)^T y(t)\}, \quad E(u(t)u(\tau)^T) = \delta(t - \tau) I.$$

The  $H_2$  norm of a continuous-time system with transfer function  $H(s)$  is given by:

$$\|H\|_2 = \sqrt{\frac{1}{2\pi} \int_{-\infty}^{\infty} \text{Trace} [H(j\omega)^H H(j\omega)] d\omega}.$$

For a discrete-time system with transfer function  $H(z)$ , the  $H_2$  norm is given by:

$$\|H\|_2 = \sqrt{\frac{1}{2\pi} \int_{-\pi}^{\pi} \text{Trace} [H(e^{j\omega})^H H(e^{j\omega})] d\omega}.$$

The  $H_2$  norm is infinite in the following cases:

- **sys** is unstable.
- **sys** is continuous and has a nonzero feedthrough (that is, nonzero gain at the frequency  $\omega = \infty$ ).

Using `norm(sys)` produces the same result as `sqrt(trace(covar(sys,1)))`.

## L-infinity norm

The  $L_\infty$  norm of a SISO linear system is the peak gain of the frequency response. For a MIMO system, the  $L_\infty$  norm is the peak gain across all input/output channels.

For a continuous-time system  $H(s)$ , this definition means:

$$\|H(s)\|_{L_\infty} = \max_{\omega \in R} |H(j\omega)| \quad (\text{SISO})$$

$$\|H(s)\|_{L_\infty} = \max_{\omega \in R} \sigma_{\max}(H(j\omega)) \quad (\text{MIMO})$$

where  $\sigma_{\max}(\cdot)$  denotes the largest singular value of a matrix.

For a discrete-time system  $H(z)$ , the definition means:

$$\|H(z)\|_{L_\infty} = \max_{\theta \in [0, 2\pi]} |H(e^{j\theta})| \quad (\text{SISO})$$

$$\|H(z)\|_{L_\infty} = \max_{\theta \in [0, 2\pi]} \sigma_{\max}(H(e^{j\theta})) \quad (\text{MIMO})$$

For stable systems, the  $L_\infty$  norm is equivalent to the  $H_\infty$  norm. For more information, see `hinfnorm`. For a system with unstable poles, the  $H_\infty$  norm is infinite. For all systems, `norm` returns the  $L_\infty$  norm, which is the peak gain without regard to system stability.

## Algorithms

After converting `sys` to a state space model, `norm` uses the same algorithm as `covar` for the  $H_2$  norm. For the  $L_\infty$  norm, `norm` uses the algorithm of [1]. `norm` computes the peak gain using the SLICOT library. For more information about the SLICOT library, see <http://slicot.org>.

## References

- [1] Bruisma, N.A. and M. Steinbuch, "A Fast Algorithm to Compute the  $H_\infty$ -Norm of a Transfer Function Matrix," *System Control Letters*, 14 (1990), pp. 287-293.

## See Also

`freqresp` | `getPeakGain` | `hinfnorm` | `sigma`

**Introduced before R2006a**

# nparams

Number of model parameters

## Syntax

```
np = nparams(sys)
np = nparams(sys, 'free')
```

## Description

`np = nparams(sys)` returns the number of parameters in the identified model `sys`.

`np = nparams(sys, 'free')` returns the number free estimation parameters in the identified model `sys`.

---

**Note** Not all model coefficients are parameters, such as the leading entry of the denominator polynomials in `idpoly` and `idtf` models.

---

## Input Arguments

**sys**

Identified linear model.

## Output Arguments

**np**

Number of parameters of `sys`.

For the syntax `np = nparams(sys, 'free')`, `np` is the number of free estimation parameters of `sys`.

`idgrey` models can contain non-scalar parameters. `nparams` accounts for each individual entry of the non-scalar parameters in the total parameter count.

### Examples

Obtain the number of parameters of a transfer function model.

```
sys = idtf(1,[1 2]);  
np = nparams(sys);
```

Obtain the number of free estimation parameters of a transfer function model.

```
sys0 = idtf([1 0],[1 2 0]);  
sys0.Structure.Denominator.Free(3) = false;  
np = nparams(sys,'free');
```

### See Also

`idfrd` | `idgrey` | `idpoly` | `idproc` | `idss` | `idtf` | `size`

**Introduced in R2012a**

# nuderst

Set step size for numerical differentiation

## Syntax

```
nds = nuderst(pars)
```

## Description

Many estimation functions use numerical differentiation with respect to the model parameters to compute their values.

The step size used in these numerical derivatives is determined by the `nuderst` command. The output argument `nds` is a row vector whose  $k$ th entry gives the increment to be used when differentiating with respect to the  $k$ th element of the parameter vector `pars`.

The default version of `nuderst` uses a very simple method. The step size is the maximum of  $10^{-4}$  times the absolute value of the current parameter and  $10^{-7}$ . You can adjust this to the actual value of the corresponding parameter by editing `nuderst`. Note that the nominal value, for example 0, of a parameter might not reflect its normal size.

**Introduced before R2006a**

## nyquist

Nyquist plot of frequency response

### Syntax

```
nyquist(sys)
nyquist(sys,w)
nyquist(sys1,sys2,...,sysN)
nyquist(sys1,sys2,...,sysN,w)
nyquist(sys1,'PlotStyle1',...,sysN,'PlotStyleN')
[re,im,w] = nyquist(sys)
[re,im] = nyquist(sys,w)
[re,im,w,sdre,sdim] = nyquist(sys)
```

### Description

`nyquist` creates a Nyquist plot of the frequency response of a dynamic system model. When invoked without left-hand arguments, `nyquist` produces a Nyquist plot on the screen. Nyquist plots are used to analyze system properties including gain margin, phase margin, and stability.

`nyquist(sys)` creates a Nyquist plot of a dynamic system `sys`. This model can be continuous or discrete, and SISO or MIMO. In the MIMO case, `nyquist` produces an array of Nyquist plots, each plot showing the response of one particular I/O channel. The frequency points are chosen automatically based on the system poles and zeros.

`nyquist(sys,w)` explicitly specifies the frequency range or frequency points to be used for the plot. To focus on a particular frequency interval, set `w = {wmin,wmax}`. To use particular frequency points, set `w` to the vector of desired frequencies. Use `logspace` to generate logarithmically spaced frequency vectors. Frequencies must be in `rad/TimeUnit`, where `TimeUnit` is the time units of the input dynamic system, specified in the `TimeUnit` property of `sys`.

`nyquist(sys1,sys2,...,sysN)` or `nyquist(sys1,sys2,...,sysN,w)` superimposes the Nyquist plots of several LTI models on a single figure. All systems must

have the same number of inputs and outputs, but may otherwise be a mix of continuous- and discrete-time systems. You can also specify a distinctive color, linestyle, and/or marker for each system plot with the syntax `nyquist(sys1, 'PlotStyle1', ..., sysN, 'PlotStyleN')`.

`[re,im,w] = nyquist(sys)` and `[re,im] = nyquist(sys,w)` return the real and imaginary parts of the frequency response at the frequencies `w` (in rad/TimeUnit). `re` and `im` are 3-D arrays (see "Arguments" below for details).

`[re,im,w,sdre,sdim] = nyquist(sys)` also returns the standard deviations of `re` and `im` for the identified system `sys`.

## Arguments

The output arguments `re` and `im` are 3-D arrays with dimensions

$$(\text{number of outputs}) \times (\text{number of inputs}) \times (\text{length of } w)$$

For SISO systems, the scalars `re(1,1,k)` and `im(1,1,k)` are the real and imaginary parts of the response at the frequency  $\omega_k = w(k)$ .

$$\text{re}(1,1,k) = \text{Re}(h(j\omega_k))$$

$$\text{im}(1,1,k) = \text{Im}(h(j\omega_k))$$

For MIMO systems with transfer function  $H(s)$ , `re(:, :, k)` and `im(:, :, k)` give the real and imaginary parts of  $H(j\omega_k)$  (both arrays with as many rows as outputs and as many columns as inputs). Thus,

$$\text{re}(i,j,k) = \text{Re}(h_{ij}(j\omega_k))$$

$$\text{im}(i,j,k) = \text{Im}(h_{ij}(j\omega_k))$$

where  $h_{ij}$  is the transfer function from input  $j$  to output  $i$ .

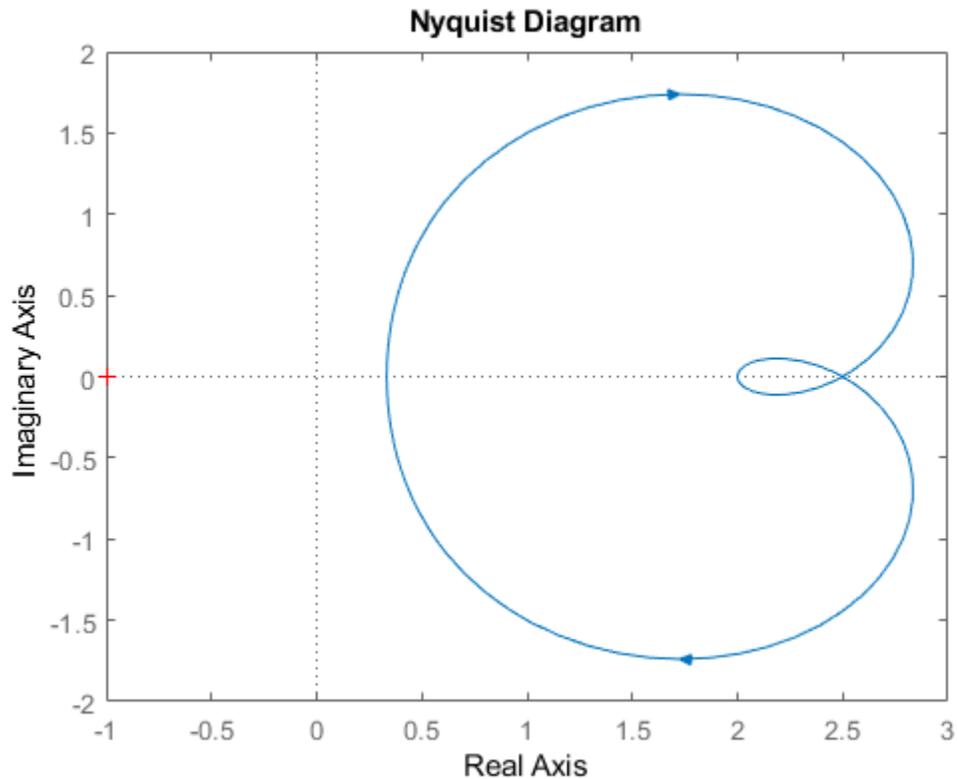
## Examples

### Nyquist Plot of Dynamic System

Plot the Nyquist response of the system

$$H(s) = \frac{2s^2 + 5s + 1}{s^2 + 2s + 3}$$

```
H = tf([2 5 1],[1 2 3]);  
nyquist(H)
```



The nyquist function has support for M-circles, which are the contours of the constant closed-loop magnitude. M-circles are defined as the locus of complex numbers where

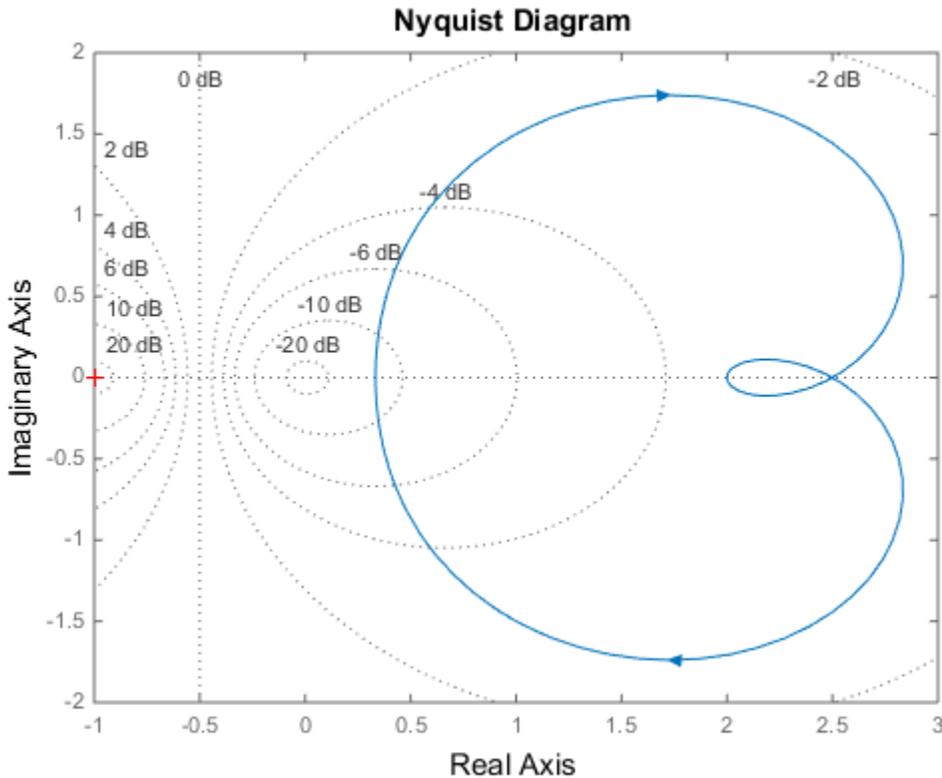
$$T(j\omega) = \left| \frac{G(j\omega)}{1 + G(j\omega)} \right|$$

is a constant value. In this equation,  $\omega$  is the frequency in radians/TimeUnit, where TimeUnit is the system time units, and  $G$  is the collection of complex numbers that satisfy the constant magnitude requirement.

To activate the grid, select **Grid** from the right-click menu or type

grid

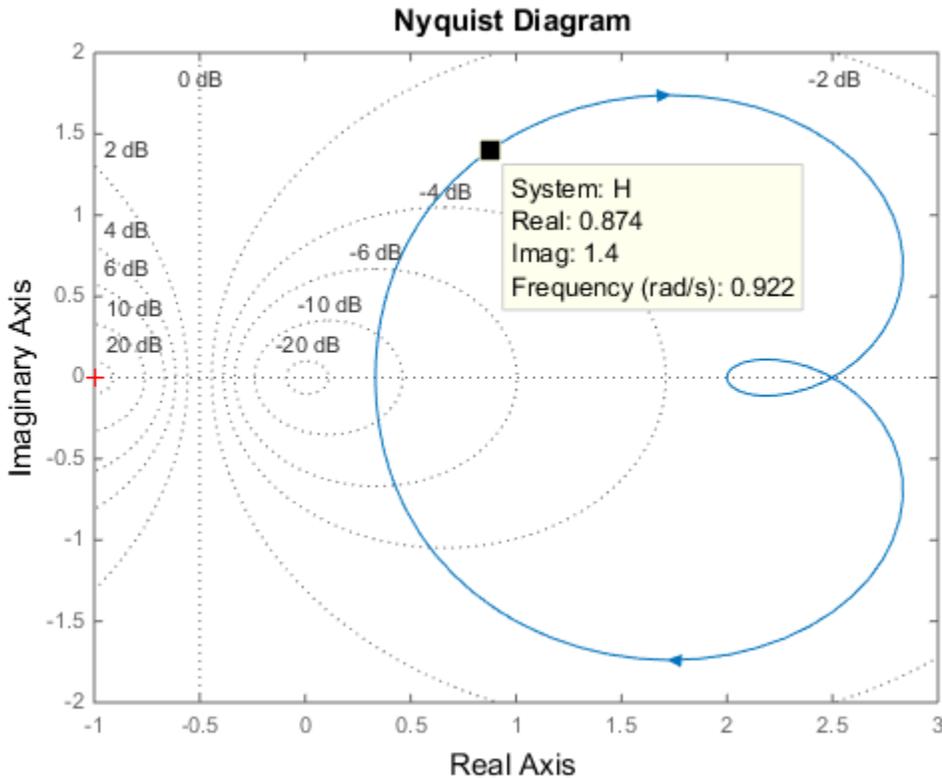
at the MATLAB prompt. This figure shows the M circles for transfer function  $H$ .



You have two zoom options available from the right-click menu that apply specifically to Nyquist plots:

- **Tight** —Clips unbounded branches of the Nyquist plot, but still includes the critical point (-1, 0)
- **On (-1,0)** — Zooms around the critical point (-1,0)

Also, click anywhere on the curve to activate data markers that display the real and imaginary values at a given frequency. This figure shows the nyquist plot with a data marker.



## Create Nyquist Plot of Identified Model With Response Uncertainty

Compute the standard deviations of the real and imaginary parts of the frequency response of an identified model. Use this data to create a  $3\sigma$  plot of the response uncertainty.

Load estimation data `z2`.

```
load iddata2 z2;
```

Identify a transfer function model using the data.

```
sys_p = tfest(z2,2);
```

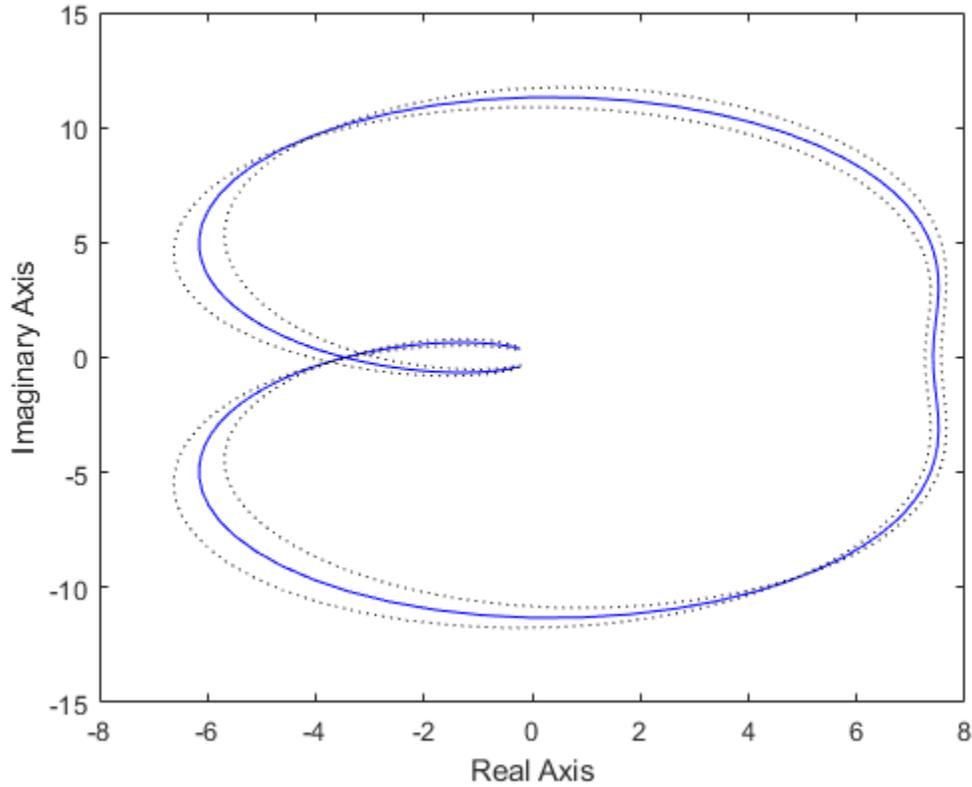
Obtain the standard deviations for the real and imaginary parts of the frequency response for a set of 512 frequencies,  $w$ .

```
w = linspace(-10*pi,10*pi,512);  
[re,im,wout,sdre,sdim] = nyquist(sys_p,w);
```

Here  $re$  and  $im$  are the real and imaginary parts of the frequency response, and  $sdre$  and  $sdim$  are their standard deviations, respectively. The frequencies in  $wout$  are the same as the frequencies you specified in  $w$ .

Create a Nyquist plot showing the response and its  $3\sigma$  uncertainty.

```
re = squeeze(re);  
im = squeeze(im);  
sdre = squeeze(sdre);  
sdim = squeeze(sdim);  
plot(re,im,'b',re+3*sdre,im+3*sdim,'k:',re-3*sdre,im-3*sdim,'k:')  
xlabel('Real Axis');  
ylabel('Imaginary Axis');
```



## Tips

You can change the properties of your plot, for example the units. For information on the ways to change properties of your plots, see “Ways to Customize Plots” (Control System Toolbox).

## Algorithms

See `bode`.

## **See Also**

**Linear System Analyzer** | `bode` | `evalfr` | `freqresp` | `nichols` | `sigma`

**Introduced before R2006a**

# nyquistoptions

List of Nyquist plot options

## Syntax

P = nyquistoptions  
 P = nyquistoptions('cstprefs')

## Description

P = nyquistoptions returns the default options for Nyquist plots. You can use these options to customize the Nyquist plot appearance using the command line.

P = nyquistoptions('cstprefs') initializes the plot options with the options you selected in the Control System and System Identification Toolbox Preferences Editor. For more information about the editor, see “Toolbox Preferences Editor”.

The following table summarizes the Nyquist plot options.

Option	Description
Title, XLabel, YLabel	Label text and style
TickLabel	Tick label style
Grid	Show or hide the grid Specified as one of the following values: 'off'   'on' <b>Default:</b> 'off'
GridColor	Color of the grid lines Specified as one of the following: Vector of RGB values in the range [0,1]   character vector of color name   'none'. For example, for yellow color, specify as one of the following: [1 1 0], 'yellow', or 'y'. <b>Default:</b> [0.15,0.15,0.15]
XlimMode, YlimMode	Limit modes

Option	Description
Xlim, Ylim	Axes limits
IOGrouping	Grouping of input-output pairs Specified as one of the following values: 'none'  'inputs' 'outputs' 'all' <b>Default:</b> 'none'
InputLabel, OutputLabel	Input and output label styles
InputVisible, OutputVisible	Visibility of input and output channels

Option	Description
FreqUnits	<p>Frequency units, specified as one of the following values:</p> <ul style="list-style-type: none"> <li>• 'Hz'</li> <li>• 'rad/second'</li> <li>• 'rpm'</li> <li>• 'kHz'</li> <li>• 'MHz'</li> <li>• 'GHz'</li> <li>• 'rad/nanosecond'</li> <li>• 'rad/microsecond'</li> <li>• 'rad/millisecond'</li> <li>• 'rad/minute'</li> <li>• 'rad/hour'</li> <li>• 'rad/day'</li> <li>• 'rad/week'</li> <li>• 'rad/month'</li> <li>• 'rad/year'</li> <li>• 'cycles/nanosecond'</li> <li>• 'cycles/microsecond'</li> <li>• 'cycles/millisecond'</li> <li>• 'cycles/hour'</li> <li>• 'cycles/day'</li> <li>• 'cycles/week'</li> <li>• 'cycles/month'</li> <li>• 'cycles/year'</li> </ul> <p><b>Default:</b> 'rad/s'</p> <p>You can also specify 'auto' which uses frequency units rad/TimeUnit relative to system time units specified in the TimeUnit property. For</p>

Option	Description
	multiple systems with different time units, the units of the first system are used.
MagUnits	Magnitude units Specified as one of the following values: 'dB'   'abs' <b>Default:</b> 'dB'
PhaseUnits	Phase units Specified as one of the following values: 'deg'   'rad' <b>Default:</b> 'deg'
ShowFullContour	Show response for negative frequencies Specified as one of the following values: 'on'   'off' <b>Default:</b> 'on'
ConfidenceRegionNumberSD	Number of standard deviations to use to plotting the response confidence region (identified models only). <b>Default:</b> 1.
ConfidenceRegionDisplaySpacing	The frequency spacing of confidence ellipses. For identified models only. <b>Default:</b> 5, which means the confidence ellipses are shown at every 5th frequency sample.

## Examples

This example shows how to create a Nyquist plot displaying the full contour (the response for both positive and negative frequencies).

```
P = nyquistoptions;  
P.ShowFullContour = 'on';  
h = nyquistplot(tf(1,[1,.2,1]),P);
```

## See Also

getoptions | nyquist | nyquistplot | setoptions | setoptions | showConfidence

**Introduced in R2012a**

# nyquistplot

Nyquist plot with additional plot customization options

## Syntax

```
h = nyquistplot(sys)
nyquistplot(sys, {wmin, wmax})
nyquistplot(sys, w)
nyquistplot(sys1, sys2, ..., w)
nyquistplot(AX, ...)
nyquistplot(..., plotoptions)
```

## Description

`h = nyquistplot(sys)` draws the Nyquist plot of the dynamic system model `sys`. It also returns the plot handle `h`. You can use this handle to customize the plot with the `getoptions` and `setoptions` commands. Type

```
help nyquistoptions
```

for a list of available plot options.

The frequency range and number of points are chosen automatically. See `bode` for details on the notion of frequency in discrete time.

`nyquistplot(sys, {wmin, wmax})` draws the Nyquist plot for frequencies between `wmin` and `wmax` (in `rad/TimeUnit`, where `TimeUnit` is the time units of the input dynamic system, specified in the `TimeUnit` property of `sys`).

`nyquistplot(sys, w)` uses the user-supplied vector `w` of frequencies (in `rad/TimeUnit`, where `TimeUnit` is the time units of the input dynamic system, specified in the `TimeUnit` property of `sys`) at which the Nyquist response is to be evaluated. See `logspace` to generate logarithmically spaced frequency vectors.

`nyquistplot(sys1, sys2, ..., w)` draws the Nyquist plots of multiple models `sys1, sys2, ...` on a single plot. The frequency vector `w` is optional. You can also specify a color, line style, and marker for each system, as in

```
nyquistplot(sys1,'r',sys2,'y--',sys3,'gx')
```

nyquistplot(Ax,...) plots into the axes with handle AX.

nyquistplot(..., plotoptions) plots the Nyquist response with the options specified in plotoptions. Type

```
help nyquistoptions
```

for more details.

## Examples

### Example 1

#### Customize Nyquist Plot Frequency Units

Plot the Nyquist frequency response and change the units to rad/s.

```
sys = rss(5);  
h = nyquistplot(sys);  
% Change units to radians per second.  
setoptions(h,'FreqUnits','rad/s');
```

### Example 2

Compare the frequency responses of identified state-space models of order 2 and 6 along with their 1-std confidence regions rendered at every 50th frequency sample.

```
load iddata1  
sys1 = n4sid(z1, 2) % discrete-time IDSS model of order 2  
sys2 = n4sid(z1, 6) % discrete-time IDSS model of order 6
```

Both models produce about 76% fit to data. However, sys2 shows higher uncertainty in its frequency response, especially close to Nyquist frequency as shown by the plot:

```
w = linspace(10,10*pi,256);  
h = nyquistplot(sys1,sys2,w);  
setoptions(h,'ConfidenceRegionDisplaySpacing',50,'ShowFullContour','off');
```

Right-click to turn on the confidence region characteristic by using the **Characteristics-> Confidence Region**.

## Tips

You can change the properties of your plot, for example the units. For information on the ways to change properties of your plots, see “Ways to Customize Plots” (Control System Toolbox).

## See Also

`getoptions` | `nyquist` | `setoptions` | `showConfidence`

**Introduced in R2012a**

## oe

Estimate Output-Error polynomial model using time or frequency domain data

### Syntax

```
sys = oe(data, [nb nf nk])  
sys = oe(data, [nb nf nk], Name, Value)  
sys = oe(data, init_sys)  
sys = oe(data, ___, opt)
```

### Description

`sys = oe(data, [nb nf nk])` estimates an Output-Error model, `sys`, represented by:

$$y(t) = \frac{B(q)}{F(q)}u(t - nk) + e(t)$$

$y(t)$  is the output,  $u(t)$  is the input, and  $e(t)$  is the error.

`sys` is estimated for the time- or frequency-domain, measured input-output data, `data`. The orders, `[nb nf nk]`, parameterize the estimated polynomial.

`sys = oe(data, [nb nf nk], Name, Value)` specifies model structure attributes using additional options specified by one or more `Name, Value` pair arguments.

`sys = oe(data, init_sys)` uses the linear system `init_sys` to configure the initial parameterization of `sys`.

`sys = oe(data, ___, opt)` estimates a polynomial model using the option set, `opt`, to specify estimation behavior.

## Input Arguments

### data

Estimation data.

For time domain estimation, `data` is an `iddata` object containing the input and output signal values.

For frequency domain estimation, `data` can be one of the following:

- Recorded frequency response data (`frd` or `idfrd`)
- `iddata` object with its properties specified as follows:
  - `InputData` — Fourier transform of the input signal
  - `OutputData` — Fourier transform of the output signal
  - `Domain` — 'Frequency'

For multi-experiment data, the sample times and inter-sample behavior of all the experiments must match.

### [nb nf nk]

Output error model orders.

For a system represented by:

$$y(t) = \frac{B(q)}{F(q)} u(t - nk) + e(t)$$

where  $y(t)$  is the output,  $u(t)$  is the input and  $e(t)$  is the error.

- `nb` — Order of the  $B$  polynomial + 1. `nb` is an  $N_y$ -by- $N_u$  matrix.  $N_y$  is the number of outputs and  $N_u$  is the number of inputs.
- `nf` — Order of the  $F$  polynomial. `nf` is an  $N_y$ -by- $N_u$  matrix.  $N_y$  is the number of outputs and  $N_u$  is the number of inputs.
- `nk` — Input delay, expressed as the number of samples. `nk` is an  $N_y$ -by- $N_u$  matrix.  $N_y$  is the number of outputs and  $N_u$  is the number of inputs. The delay appears as leading zeros of the  $B$  polynomial.

For estimation using continuous-time data, only specify [nb nf] and omit nk.

## **init\_sys**

Linear system that configures the initial parameterization of `sys`.

You obtain `init_sys` by either performing an estimation using measured data or by direct construction.

If `init_sys` is an `idpoly` model of Output-Error structure, `oe` uses the parameter values of `init_sys` as the initial guess for estimating `sys`.

Use the `Structure` property of `init_sys` to configure initial guesses and constraints for  $B(q)$  and  $F(q)$ . For example:

- To specify an initial guess for the  $F(q)$  term of `init_sys`, set `init_sys.Structure.F.Value` as the initial guess.
- To specify constraints for the  $B(q)$  term of `init_sys`:
  - Set `init_sys.Structure.B.Minimum` to the minimum  $B(q)$  coefficient values
  - Set `init_sys.Structure.B.Maximum` to the maximum  $B(q)$  coefficient values
  - Set `init_sys.Structure.B.Free` to indicate which  $B(q)$  coefficients are free for estimation

If `init_sys` is not a polynomial model of Output-Error structure, the software first converts `init_sys` to an Output-Error structure model. `oe` uses the parameters of the resulting model as the initial guess for estimating `sys`.

If `opt` is not specified, and `init_sys` was created by estimation, then the estimation options from `init_sys.Report.OptionsUsed` are used.

## **opt**

Estimation options.

`opt` is an option set, created using `oeOptions`, that specifies estimation options including:

- Estimation objective
- Handling of initial conditions

- Numerical search method and the associated options

## Name-Value Pair Arguments

Specify optional comma-separated pairs of `Name`, `Value` arguments. `Name` is the argument name and `Value` is the corresponding value. `Name` must appear inside quotes. You can specify several name and value pair arguments in any order as `Name1, Value1, . . . , NameN, ValueN`.

### InputDelay

Input delay for each input channel, specified as a scalar value or numeric vector. For continuous-time systems, specify input delays in the time unit stored in the `TimeUnit` property. For discrete-time systems, specify input delays in integer multiples of the sample time `Ts`. For example, `InputDelay = 3` means a delay of three sample times.

For a system with `Nu` inputs, set `InputDelay` to an `Nu`-by-1 vector. Each entry of this vector is a numerical value that represents the input delay for the corresponding input channel.

You can also set `InputDelay` to a scalar value to apply the same delay to all channels.

**Default:** 0

### IODElay

Transport delays. `IODElay` is a numeric array specifying a separate transport delay for each input/output pair.

For continuous-time systems, specify transport delays in the time unit stored in the `TimeUnit` property. For discrete-time systems, specify transport delays as integers denoting delay of a multiple of the sample time `Ts`. You can specify `IODElay` as an alternative to the `nk` value. Doing so simplifies the model structure by reducing the number of leading zeros the  $B$  polynomial. In particular, you can represent  $\max(nk - 1, 0)$  leading zeros as input/output delays using `IODElay` instead.

For a MIMO system with `Ny` outputs and `Nu` inputs, set `IODElay` to a `Ny`-by-`Nu` array. Each entry of this array is a numerical value that represents the transport delay for the corresponding input/output pair. You can also set `IODElay` to a scalar value to apply the same delay to all input/output pairs.

**Default:** 0 for all input/output pairs

## Output Arguments

### sys

Output-Error polynomial model that fits the estimation data, returned as a `idpoly` model object. This model is created using the specified model orders, delays, and estimation options.

Information about the estimation results and options used is stored in the `Report` property of the model. `Report` has the following fields:

Report Field	Description
Status	Summary of the model status, which indicates whether the model was created by construction or obtained by estimation.
Method	Estimation command used.
InitialCondition	Handling of initial conditions during model estimation, returned as one of the following values: <ul style="list-style-type: none"><li>'zero' — The initial conditions were set to zero.</li><li>'estimate' — The initial conditions were treated as independent estimation parameters.</li><li>'backcast' — The initial conditions were estimated using the best least squares fit.</li></ul> <p>This field is especially useful to view how the initial conditions were handled when the <code>InitialCondition</code> option in the estimation option set is 'auto'.</p>

Report Field	Description																		
Fit	<p>Quantitative assessment of the estimation, returned as a structure. See “Loss Function and Model Quality Metrics” for more information on these quality metrics. The structure has the following fields:</p> <table border="1" data-bbox="387 456 1331 944"> <thead> <tr> <th data-bbox="387 456 516 499">Field</th> <th data-bbox="516 456 1331 499">Description</th> </tr> </thead> <tbody> <tr> <td data-bbox="387 499 516 604">FitPercent</td> <td data-bbox="516 499 1331 604">Normalized root mean squared error (NRMSE) measure of how well the response of the model fits the estimation data, expressed as a percentage.</td> </tr> <tr> <td data-bbox="387 604 516 647">LossFcn</td> <td data-bbox="516 604 1331 647">Value of the loss function when the estimation completes.</td> </tr> <tr> <td data-bbox="387 647 516 725">MSE</td> <td data-bbox="516 647 1331 725">Mean squared error (MSE) measure of how well the response of the model fits the estimation data.</td> </tr> <tr> <td data-bbox="387 725 516 769">FPE</td> <td data-bbox="516 725 1331 769">Final prediction error for the model.</td> </tr> <tr> <td data-bbox="387 769 516 812">AIC</td> <td data-bbox="516 769 1331 812">Raw Akaike Information Criteria (AIC) measure of model quality.</td> </tr> <tr> <td data-bbox="387 812 516 855">AICc</td> <td data-bbox="516 812 1331 855">Small sample-size corrected AIC.</td> </tr> <tr> <td data-bbox="387 855 516 899">nAIC</td> <td data-bbox="516 855 1331 899">Normalized AIC.</td> </tr> <tr> <td data-bbox="387 899 516 944">BIC</td> <td data-bbox="516 899 1331 944">Bayesian Information Criteria (BIC).</td> </tr> </tbody> </table>	Field	Description	FitPercent	Normalized root mean squared error (NRMSE) measure of how well the response of the model fits the estimation data, expressed as a percentage.	LossFcn	Value of the loss function when the estimation completes.	MSE	Mean squared error (MSE) measure of how well the response of the model fits the estimation data.	FPE	Final prediction error for the model.	AIC	Raw Akaike Information Criteria (AIC) measure of model quality.	AICc	Small sample-size corrected AIC.	nAIC	Normalized AIC.	BIC	Bayesian Information Criteria (BIC).
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BIC	Bayesian Information Criteria (BIC).																		
Parameters	Estimated values of model parameters.																		
OptionsUsed	Option set used for estimation. If no custom options were configured, this is a set of default options. See <code>oeOptions</code> for more information.																		
RandState	State of the random number stream at the start of estimation. Empty, <code>[]</code> , if randomization was not used during estimation. For more information, see <code>rng</code> in the MATLAB documentation.																		

Report Field	Description																
DataUsed	Attributes of the data used for estimation, returned as a structure with the following fields:																
	<table border="1"> <thead> <tr> <th data-bbox="385 425 516 465">Field</th> <th data-bbox="516 425 1335 465">Description</th> </tr> </thead> <tbody> <tr> <td data-bbox="385 465 516 512">Name</td> <td data-bbox="516 465 1335 512">Name of the data set.</td> </tr> <tr> <td data-bbox="385 512 516 558">Type</td> <td data-bbox="516 512 1335 558">Data type.</td> </tr> <tr> <td data-bbox="385 558 516 605">Length</td> <td data-bbox="516 558 1335 605">Number of data samples.</td> </tr> <tr> <td data-bbox="385 605 516 652">Ts</td> <td data-bbox="516 605 1335 652">Sample time.</td> </tr> <tr> <td data-bbox="385 652 516 956">InterSample</td> <td data-bbox="516 652 1335 956">                     Input intersample behavior, returned as one of the following values:                     <ul style="list-style-type: none"> <li>• 'zoh' — Zero-order hold maintains a piecewise-constant input signal between samples.</li> <li>• 'foh' — First-order hold maintains a piecewise-linear input signal between samples.</li> <li>• 'bl' — Band-limited behavior specifies that the continuous-time input signal has zero power above the Nyquist frequency.</li> </ul> </td> </tr> <tr> <td data-bbox="385 956 516 1038">InputOffset</td> <td data-bbox="516 956 1335 1038">Offset removed from time-domain input data during estimation. For nonlinear models, it is [ ].</td> </tr> <tr> <td data-bbox="385 1038 516 1112">OutputOffset</td> <td data-bbox="516 1038 1335 1112">Offset removed from time-domain output data during estimation. For nonlinear models, it is [ ].</td> </tr> </tbody> </table>	Field	Description	Name	Name of the data set.	Type	Data type.	Length	Number of data samples.	Ts	Sample time.	InterSample	Input intersample behavior, returned as one of the following values: <ul style="list-style-type: none"> <li>• 'zoh' — Zero-order hold maintains a piecewise-constant input signal between samples.</li> <li>• 'foh' — First-order hold maintains a piecewise-linear input signal between samples.</li> <li>• 'bl' — Band-limited behavior specifies that the continuous-time input signal has zero power above the Nyquist frequency.</li> </ul>	InputOffset	Offset removed from time-domain input data during estimation. For nonlinear models, it is [ ].	OutputOffset	Offset removed from time-domain output data during estimation. For nonlinear models, it is [ ].
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OutputOffset	Offset removed from time-domain output data during estimation. For nonlinear models, it is [ ].																

Report Field	Description	
Termination	Termination conditions for the iterative search used for prediction error minimization. Structure with the following fields:	
	Field	Description
	WhyStop	Reason for terminating the numerical search.
	Iterations	Number of search iterations performed by the estimation algorithm.
	FirstOrderOptimality	$\infty$ -norm of the gradient search vector when the search algorithm terminates.
	FcnCount	Number of times the objective function was called.
	UpdateNorm	Norm of the gradient search vector in the last iteration. Omitted when the search method is 'lsqnonlin' or 'fmincon'.
	LastImprovement	Criterion improvement in the last iteration, expressed as a percentage. Omitted when the search method is 'lsqnonlin' or 'fmincon'.
	Algorithm	Algorithm used by 'lsqnonlin' or 'fmincon' search method. Omitted when other search methods are used.
For estimation methods that do not require numerical search optimization, the Termination field is omitted.		

For more information on using Report, see “Estimation Report”.

## Examples

### Estimate Continuous-Time Model Using Frequency Response

Obtain the estimation data.

```
filename = fullfile(matlabroot, 'examples', 'ident', 'oe_data1.mat');
load(filename);
```

`data`, an `idfrd` object, contains the continuous-time frequency response for the following model:

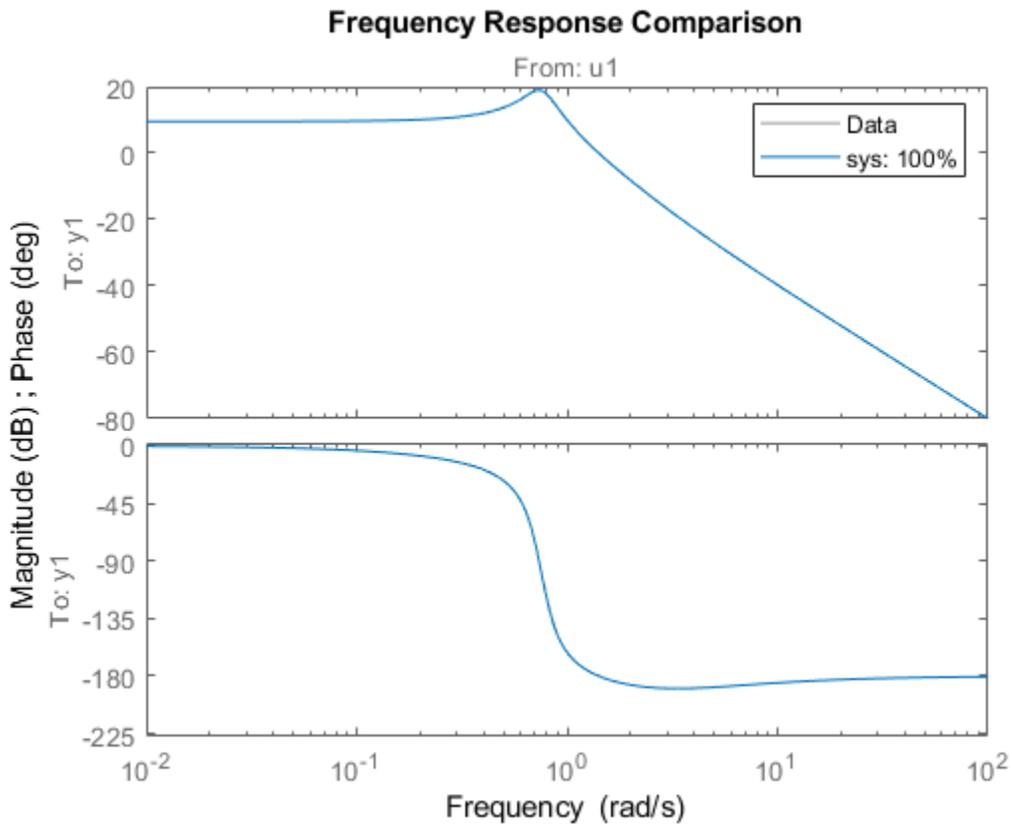
$$G(s) = \frac{s + 3}{s^3 + 2s^2 + s + 1}$$

Estimate the model.

```
nb = 2;  
nk = 3;  
sys = oe(data, [nb nk]);
```

Evaluate the goodness of the fit.

```
compare(data, sys);
```



### Estimate Output-Error Model Using Regularization

Estimate a high-order OE model from data collected by simulating a high-order system. Determine the regularization constants by trial and error and use the values for model estimation.

Load data.

```
load regularizationExampleData.mat m0simdata
```

Estimate an unregularized OE model of order 30.

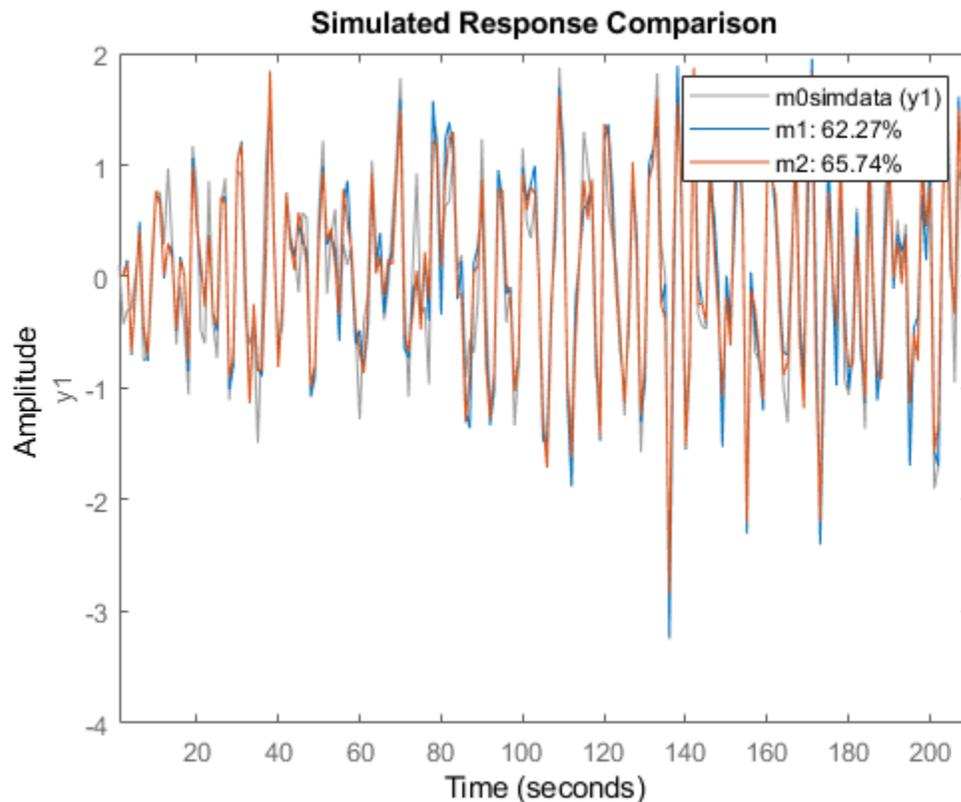
```
m1 = oe(m0simdata,[30 30 1]);
```

Obtain a regularized OE model by determining Lambda value using trial and error.

```
opt = oeOptions;  
opt.Regularization.Lambda = 1;  
m2 = oe(m0simdata,[30 30 1],opt);
```

Compare the model outputs with the estimation data.

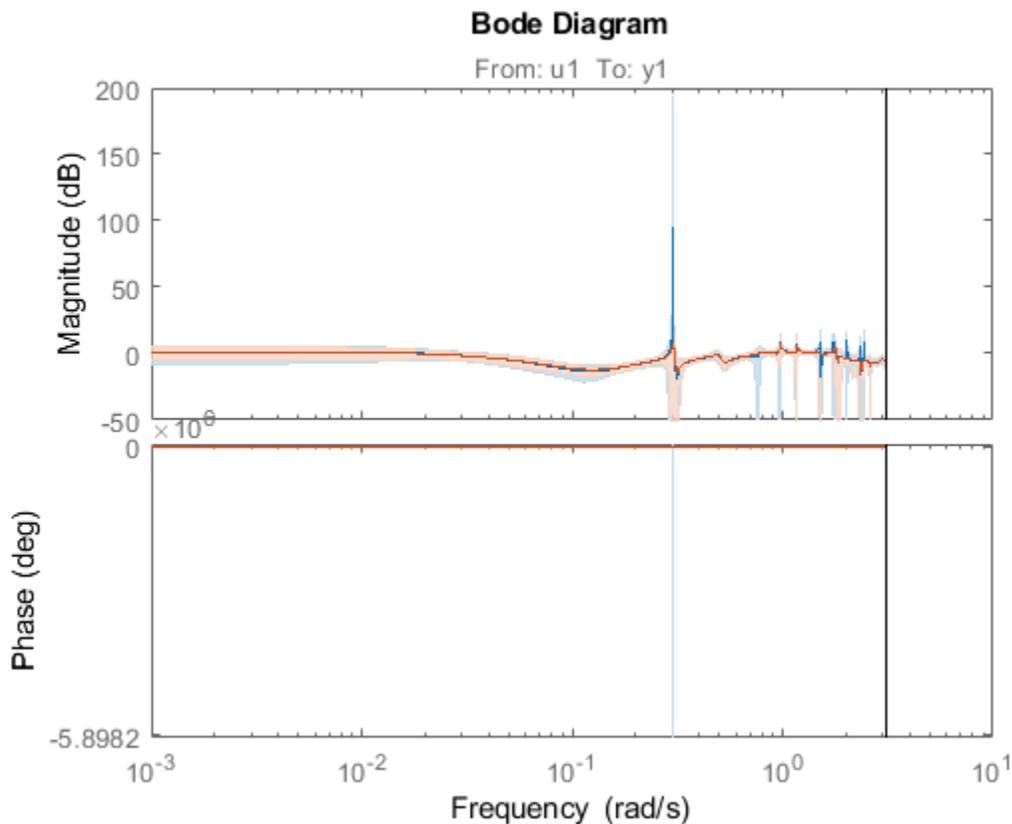
```
opt = compareOptions('InitialCondition','z');  
compare(m0simdata,m1,m2,opt);
```



The regularized model, m2, produces a better fit than the unregularized model, m1.

Compare the variance in the model responses.

```
h = bodeplot(m1,m2);  
opt = getoptions(h);  
opt.PhaseMatching = 'on';  
opt.ConfidenceRegionNumberSD = 3;  
opt.PhaseMatching = 'on';  
setoptions(h,opt);  
showConfidence(h);
```



The variance of the regularized model m2 is reduced compared to the unregularized model m1.

## Estimate Model Using Band-Limited Discrete-Time Frequency-Domain Data

Obtain the estimation data.

```
filename = fullfile(matlabroot, 'examples', 'ident', 'oe_data2.mat');  
load(filename, 'data', 'Ts');
```

`data`, an `iddata` object, contains the discrete-time frequency response for the following model:

$$G(s) = \frac{1000}{s + 500}$$

The sample time for `data`, `Ts`, is 0.001 seconds.

Treat `data` as continuous-time data. When you plot `data`, the input/output signals are band-limited, which allows you to treat `data` as continuous-time data. You can now obtain a continuous-time model.

```
data.Ts = 0;
```

Specify the estimation options.

```
opt = oeOptions('WeightingFilter',[0 0.5*pi/Ts]);
```

This prefilter choice directs the software to ignore the response values for frequencies higher than  $0.5\pi/Ts$  rad/s.

Estimate the model.

```
nb = 1;  
nf = 3;  
sys = oe(data,[nb nf],opt);
```

## Definitions

### Output-Error (OE) Model

The general Output-Error model structure is:

$$y(t) = \frac{B(q)}{F(q)} u(t - n_k) + e(t)$$

The orders of the Output-Error model are:

$$nb: B(q) = b_1 + b_2 q^{-1} + \dots + b_{nb} q^{-nb+1}$$

$$nf: F(q) = 1 + f_1 q^{-1} + \dots + f_{nf} q^{-nf}$$

## Continuous-Time, Output-Error Model

If `data` is continuous-time frequency-domain data, `oe` estimates a continuous-time model with transfer function:

$$G(s) = \frac{B(s)}{F(s)} = \frac{b_{nb} s^{(nb-1)} + b_{nb-1} s^{(nb-2)} + \dots + b_1}{s^{nf} + f_{nf} s^{(nf-1)} + \dots + f_1}$$

The orders of the numerator and denominator are `nb` and `nf`, similar to the discrete-time case. However, the delay `nk` has no meaning and you should omit it when specifying model orders for estimation. Use `model = oe(data, [nb nf])`. Use the `IODelay` model property to specify any input-output delays. For example, use `model = oe(data, [nb nf], 'IODelay', iod)` instead.

## Tips

- To estimate a continuous-time model when `data` represents continuous-time frequency response data, omit `nk`.

For example, use `sys = oe(data, [nb nf])`.

## Algorithms

The estimation algorithm minimizes prediction errors.

## Alternatives

Output-Error models are a special configuration of polynomial models, having only two active polynomials -  $B$  and  $F$ . For such models, it may be more convenient to use a transfer function (`idtf`) model and its estimation command, `tfest`.

Also, `tfest` is the recommended command for estimating continuous-time models.

## Extended Capabilities

### Automatic Parallel Support

Accelerate code by automatically running computation in parallel using Parallel Computing Toolbox™.

Parallel computing support is available for estimation using the `lsqnonlin` search method (requires Optimization Toolbox). To enable parallel computing, use `oeOptions`, set `SearchMethod` to `'lsqnonlin'`, and set `SearchOptions.Advanced.UseParallel` to `true`.

For example:

```
opt = oeOptions;  
opt.SearchMethod = 'lsqnonlin';  
opt.SearchOptions.Advanced.UseParallel = true;
```

### See Also

[armax](#) | [arx](#) | [bj](#) | [compare](#) | [iddata](#) | [idfrd](#) | [idpoly](#) | [iv4](#) | [n4sid](#) | [oeOptions](#) | [polyest](#) | [sim](#) | [tfest](#)

### Topics

[“What Are Polynomial Models?”](#)

[“Regularized Estimates of Model Parameters”](#)

**Introduced before R2006a**

# oeOptions

Option set for oe

## Syntax

```
opt = oeOptions  
opt = oeOptions(Name,Value)
```

## Description

`opt = oeOptions` creates the default options set for `oe`.

`opt = oeOptions(Name,Value)` creates an option set with the options specified by one or more `Name,Value` pair arguments.

## Input Arguments

### Name-Value Pair Arguments

Specify optional comma-separated pairs of `Name,Value` arguments. `Name` is the argument name and `Value` is the corresponding value. `Name` must appear inside quotes. You can specify several name and value pair arguments in any order as `Name1,Value1,...,NameN,ValueN`.

### InitialCondition — Handling of initial conditions

```
'auto' (default) | 'zero' | 'estimate' | 'backcast'
```

Handling of initial conditions during estimation, specified as one of the following values:

- `'zero'` — The initial conditions are set to zero.
- `'estimate'` — The initial conditions are treated as independent estimation parameters.
- `'backcast'` — The initial conditions are estimated using the best least squares fit.

- 'auto' — The software chooses the method to handle initial conditions based on the estimation data.

## **WeightingFilter — Weighting prefilter**

[ ] (default) | vector | matrix | cell array | linear system

Weighting prefilter applied to the loss function to be minimized during estimation. To understand the effect of `WeightingFilter` on the loss function, see “Loss Function and Model Quality Metrics”.

Specify `WeightingFilter` as one of the following values:

- [ ] — No weighting prefilter is used.
- Passbands — Specify a row vector or matrix containing frequency values that define desired passbands. You select a frequency band where the fit between estimated model and estimation data is optimized. For example, `[wl,wh]` where `wl` and `wh` represent lower and upper limits of a passband. For a matrix with several rows defining frequency passbands, `[w1l,w1h;w2l,w2h;w3l,w3h;...]`, the estimation algorithm uses the union of the frequency ranges to define the estimation passband.

Passbands are expressed in `rad/TimeUnit` for time-domain data and in `FrequencyUnit` for frequency-domain data, where `TimeUnit` and `FrequencyUnit` are the time and frequency units of the estimation data.

- SISO filter — Specify a single-input-single-output (SISO) linear filter in one of the following ways:
  - A SISO LTI model
  - `{A,B,C,D}` format, which specifies the state-space matrices of a filter with the same sample time as estimation data.
  - `{numerator,denominator}` format, which specifies the numerator and denominator of the filter as a transfer function with same sample time as estimation data.

This option calculates the weighting function as a product of the filter and the input spectrum to estimate the transfer function.

- Weighting vector — Applicable for frequency-domain data only. Specify a column vector of weights. This vector must have the same length as the frequency vector of the data set, `Data.Frequency`. Each input and output response in the data is multiplied by the corresponding weight at that frequency.

**EnforceStability — Control whether to enforce stability of model**`false` (default) | `true`

Control whether to enforce stability of estimated model, specified as the comma-separated pair consisting of 'EnforceStability' and either `true` or `false`.

Use this option when estimating models using frequency-domain data. Models estimated using time-domain data are always stable.

Data Types: `logical`

**EstimateCovariance — Control whether to generate parameter covariance data**`true` (default) | `false`

Controls whether parameter covariance data is generated, specified as `true` or `false`.

If `EstimateCovariance` is `true`, then use `getcov` to fetch the covariance matrix from the estimated model.

**Display — Specify whether to display the estimation progress**`'off'` (default) | `'on'`

Specify whether to display the estimation progress, specified as one of the following values:

- `'on'` — Information on model structure and estimation results are displayed in a progress-viewer window.
- `'off'` — No progress or results information is displayed.

**InputOffset — Removal of offset from time-domain input data during estimation**`[]` (default) | vector of positive integers | matrix

Removal of offset from time-domain input data during estimation, specified as the comma-separated pair consisting of 'InputOffset' and one of the following:

- A column vector of positive integers of length  $Nu$ , where  $Nu$  is the number of inputs.
- `[]` — Indicates no offset.
- $Nu$ -by- $Ne$  matrix — For multi-experiment data, specify `InputOffset` as an  $Nu$ -by- $Ne$  matrix.  $Nu$  is the number of inputs, and  $Ne$  is the number of experiments.

Each entry specified by `InputOffset` is subtracted from the corresponding input data.

## **OutputOffset — Removal of offset from time-domain output data during estimation**

[ ] (default) | vector | matrix

Removal of offset from time-domain output data during estimation, specified as the comma-separated pair consisting of 'OutputOffset' and one of the following:

- A column vector of length  $N_y$ , where  $N_y$  is the number of outputs.
- [ ] — Indicates no offset.
- $N_y$ -by- $N_e$  matrix — For multi-experiment data, specify OutputOffset as a  $N_y$ -by- $N_e$  matrix.  $N_y$  is the number of outputs, and  $N_e$  is the number of experiments.

Each entry specified by OutputOffset is subtracted from the corresponding output data.

## **Regularization — Options for regularized estimation of model parameters**

structure

Options for regularized estimation of model parameters. For more information on regularization, see “Regularized Estimates of Model Parameters”.

Regularization is a structure with the following fields:

- Lambda — Constant that determines the bias versus variance tradeoff.

Specify a positive scalar to add the regularization term to the estimation cost.

The default value of zero implies no regularization.

**Default:** 0

- R — Weighting matrix.

Specify a vector of nonnegative numbers or a square positive semi-definite matrix. The length must be equal to the number of free parameters of the model.

For black-box models, using the default value is recommended. For structured and grey-box models, you can also specify a vector of  $np$  positive numbers such that each entry denotes the confidence in the value of the associated parameter.

The default value of 1 implies a value of  $\text{eye}(np_{\text{free}})$ , where  $np_{\text{free}}$  is the number of free parameters.

**Default: 1**

- **Nominal** — The nominal value towards which the free parameters are pulled during estimation.

The default value of zero implies that the parameter values are pulled towards zero. If you are refining a model, you can set the value to 'model' to pull the parameters towards the parameter values of the initial model. The initial parameter values must be finite for this setting to work.

**Default: 0****SearchMethod — Numerical search method used for iterative parameter estimation**

'auto' (default) | 'gn' | 'gna' | 'lm' | 'grad' | 'lsqnonlin' | 'fmincon'

Numerical search method used for iterative parameter estimation, specified as the comma-separated pair consisting of 'SearchMethod' and one of the following:

- 'auto' — A combination of the line search algorithms, 'gn', 'lm', 'gna', and 'grad' methods is tried in sequence at each iteration. The first descent direction leading to a reduction in estimation cost is used.
- 'gn' — Subspace Gauss-Newton least squares search. Singular values of the Jacobian matrix less than  $GnPinvConstant * \epsilon * \max(\text{size}(J)) * \text{norm}(J)$  are discarded when computing the search direction.  $J$  is the Jacobian matrix. The Hessian matrix is approximated as  $J^T J$ . If there is no improvement in this direction, the function tries the gradient direction.
- 'gna' — Adaptive subspace Gauss-Newton search. Eigenvalues less than  $\text{gamma} * \max(sv)$  of the Hessian are ignored, where  $sv$  contains the singular values of the Hessian. The Gauss-Newton direction is computed in the remaining subspace.  $\text{gamma}$  has the initial value `InitialGnaTolerance` (see Advanced in 'SearchOptions' for more information). This value is increased by the factor `LMStep` each time the search fails to find a lower value of the criterion in fewer than five bisections. This value is decreased by the factor  $2 * \text{LMStep}$  each time a search is successful without any bisections.
- 'lm' — Levenberg-Marquardt least squares search, where the next parameter value is  $-\text{pinv}(H+d*I) * \text{grad}$  from the previous one.  $H$  is the Hessian,  $I$  is the identity matrix, and  $\text{grad}$  is the gradient.  $d$  is a number that is increased until a lower value of the criterion is found.
- 'grad' — Steepest descent least squares search.

- 'lsqnonlin' — Trust-region-reflective algorithm of lsqnonlin. Requires Optimization Toolbox software.
- 'fmincon' — Constrained nonlinear solvers. You can use the sequential quadratic programming (SQP) and trust-region-reflective algorithms of the fmincon solver. If you have Optimization Toolbox software, you can also use the interior-point and active-set algorithms of the fmincon solver. Specify the algorithm in the SearchOptions.Algorithm option. The fmincon algorithms may result in improved estimation results in the following scenarios:
  - Constrained minimization problems when there are bounds imposed on the model parameters.
  - Model structures where the loss function is a nonlinear or non smooth function of the parameters.
  - Multi-output model estimation. A determinant loss function is minimized by default for multi-output model estimation. fmincon algorithms are able to minimize such loss functions directly. The other search methods such as 'lm' and 'gn' minimize the determinant loss function by alternately estimating the noise variance and reducing the loss value for a given noise variance value. Hence, the fmincon algorithms can offer better efficiency and accuracy for multi-output model estimations.

### **SearchOptions — Option set for the search algorithm**

search option set

Option set for the search algorithm, specified as the comma-separated pair consisting of 'SearchOptions' and a search option set with fields that depend on the value of SearchMethod.

**SearchOptions Structure When SearchMethod is Specified as 'gn', 'gna', 'lm', 'grad', or 'auto'**

Field Name	Description	Default
Tolerance	Minimum percentage difference between the current value of the loss function and its expected improvement after the next iteration, specified as a positive scalar. When the percentage of expected improvement is less than <code>Tolerance</code> , the iterations stop. The estimate of the expected loss-function improvement at the next iteration is based on the Gauss-Newton vector computed for the current parameter value.	0.01
MaxIterations	<p>Maximum number of iterations during loss-function minimization, specified as a positive integer. The iterations stop when <code>MaxIterations</code> is reached or another stopping criterion is satisfied, such as <code>Tolerance</code>.</p> <p>Setting <code>MaxIterations = 0</code> returns the result of the start-up procedure.</p> <p>Use <code>sys.Report.Termination.Iterations</code> to get the actual number of iterations during an estimation, where <code>sys</code> is an <code>idtf</code> model.</p>	20

Field Name	Description	Default
Advanced	Advanced search settings, specified as a structure with the following fields:	
<b>Field Name</b>	<b>Description</b>	<b>Default</b>
GnPinvConstant	Jacobian matrix singular value threshold, specified as a positive scalar. Singular values of the Jacobian matrix that are smaller than $\text{GnPinvConstant} \times \max(\text{size}(J) \times \text{norm}(J) \times \text{eps})$ are discarded when computing the search direction. Applicable when SearchMethod is 'gn'.	10000
InitialGnaTolerance	Initial value of $\gamma$ , specified as a positive scalar. Applicable when SearchMethod is 'gna'.	0.0001
LMStartValue	Starting value of search-direction length $d$ in the Levenberg-Marquardt method, specified as a positive scalar. Applicable when SearchMethod is 'lm'.	0.001
LMStep	Size of the Levenberg-Marquardt step, specified as a positive integer. The next value of the search-direction length $d$ in the Levenberg-Marquardt method is LMStep times the previous one. Applicable when SearchMethod is 'lm'.	2
MaxBisections	Maximum number of bisections used for line search along the search direction, specified as a positive integer.	25
MaxFunctionEvaluations	Maximum number of calls to the model file, specified as a positive integer. Iterations stop if the number of calls to the model file exceeds this value.	Inf
MinParameterChange	Smallest parameter update allowed per iteration, specified as a nonnegative scalar.	0
RelativeImprovement	Relative improvement threshold, specified as a nonnegative scalar. Iterations stop if the relative	0

Field Name	Description		Default
	Field Name	Description	Default
		improvement of the criterion function is less than this value.	
	StepReduction	<p>Step reduction factor, specified as a positive scalar that is greater than 1. The suggested parameter update is reduced by the factor <code>StepReduction</code> after each try. This reduction continues until <code>MaxBisections</code> tries are completed or a lower value of the criterion function is obtained.</p> <p><code>StepReduction</code> is not applicable for SearchMethod 'lm' (Levenberg-Marquardt method).</p>	2

**SearchOptions Structure When SearchMethod is Specified as 'lsqnonlin'**

<b>Field Name</b>	<b>Description</b>	<b>Default</b>
FunctionTolerance	Termination tolerance on the loss function that the software minimizes to determine the estimated parameter values, specified as a positive scalar.  The value of FunctionTolerance is the same as that of <code>opt.SearchOptions.Advanced.TolFun</code> .	1e-5
StepTolerance	Termination tolerance on the estimated parameter values, specified as a positive scalar.  The value of StepTolerance is the same as that of <code>opt.SearchOptions.Advanced.TolX</code> .	1e-6
MaxIterations	Maximum number of iterations during loss-function minimization, specified as a positive integer. The iterations stop when MaxIterations is reached or another stopping criterion is satisfied, such as FunctionTolerance.  The value of MaxIterations is the same as that of <code>opt.SearchOptions.Advanced.MaxIter</code> .	20
Advanced	Advanced search settings, specified as an option set for <code>lsqnonlin</code> .  For more information, see the Optimization Options table in "Optimization Options" (Optimization Toolbox).	Use <code>optimset('lsqnonlin')</code> to create a default option set.

### SearchOptions Structure When SearchMethod is Specified as 'fmincon'

Field Name	Description	Default
Algorithm	<p>fmincon optimization algorithm, specified as one of the following:</p> <ul style="list-style-type: none"> <li>'sqp' — Sequential quadratic programming algorithm. The algorithm satisfies bounds at all iterations, and it can recover from NaN or Inf results. It is not a large-scale algorithm. For more information, see “Large-Scale vs. Medium-Scale Algorithms” (Optimization Toolbox).</li> <li>'trust-region-reflective' — Subspace trust-region method based on the interior-reflective Newton method. It is a large-scale algorithm.</li> <li>'interior-point' — Large-scale algorithm that requires Optimization Toolbox software. The algorithm satisfies bounds at all iterations, and it can recover from NaN or Inf results.</li> <li>'active-set' — Requires Optimization Toolbox software. The</li> </ul>	'sqp'

Field Name	Description	Default
	<p>algorithm can take large steps, which adds speed. It is not a large-scale algorithm.</p> <p>For more information about the algorithms, see “Constrained Nonlinear Optimization Algorithms” (Optimization Toolbox) and “Choosing the Algorithm” (Optimization Toolbox).</p>	
FunctionTolerance	Termination tolerance on the loss function that the software minimizes to determine the estimated parameter values, specified as a positive scalar.	1e-6
StepTolerance	Termination tolerance on the estimated parameter values, specified as a positive scalar.	1e-6
MaxIterations	Maximum number of iterations during loss function minimization, specified as a positive integer. The iterations stop when MaxIterations is reached or another stopping criterion is satisfied, such as FunctionTolerance.	100

**Advanced – Additional advanced options**  
structure

Additional advanced options, specified as a structure with the following fields:

- `ErrorThreshold` — Specifies when to adjust the weight of large errors from quadratic to linear.

Errors larger than `ErrorThreshold` times the estimated standard deviation have a linear weight in the loss function. The standard deviation is estimated robustly as the median of the absolute deviations from the median of the prediction errors, divided by 0.7. For more information on robust norm choices, see section 15.2 of [2].

`ErrorThreshold = 0` disables robustification and leads to a purely quadratic loss function. When estimating with frequency-domain data, the software sets `ErrorThreshold` to zero. For time-domain data that contains outliers, try setting `ErrorThreshold` to 1.6.

**Default:** 0

- `MaxSize` — Specifies the maximum number of elements in a segment when input-output data is split into segments.

`MaxSize` must be a positive integer.

**Default:** 250000

- `StabilityThreshold` — Specifies thresholds for stability tests.

`StabilityThreshold` is a structure with the following fields:

- `s` — Specifies the location of the right-most pole to test the stability of continuous-time models. A model is considered stable when its right-most pole is to the left of `s`.

**Default:** 0

- `z` — Specifies the maximum distance of all poles from the origin to test stability of discrete-time models. A model is considered stable if all poles are within the distance `z` from the origin.

**Default:** `1+sqrt(eps)`

- `AutoInitThreshold` — Specifies when to automatically estimate the initial condition.

The initial condition is estimated when

$$\frac{\|y_{p,z} - y_{meas}\|}{\|y_{p,e} - y_{meas}\|} > \text{AutoInitThreshold}$$

- $y_{meas}$  is the measured output.
- $y_{p,z}$  is the predicted output of a model estimated using zero initial conditions.
- $y_{p,e}$  is the predicted output of a model estimated using estimated initial conditions.

Applicable when `InitialCondition` is 'auto'.

**Default:** 1.05

## Output Arguments

### **opt** — Options set for oe

`oeOptions` option set

Option set for oe, returned as an `oeOptions` option set.

## Examples

### **Create Default Options Set for Output-Error Estimation**

```
opt = oeOptions;
```

### **Specify Options for Output-Error Estimation**

Create an options set for oe using the 'backcast' algorithm to initialize the condition and set the `Display` to 'on'.

```
opt = oeOptions('InitialCondition', 'backcast', 'Display', 'on');
```

Alternatively, use dot notation to set the values of `opt`.

```
opt = oeOptions;  
opt.InitialCondition = 'backcast';  
opt.Display = 'on';
```

## References

- [1] Wills, Adrian, B. Ninness, and S. Gibson. "On Gradient-Based Search for Multivariable System Estimates". *Proceedings of the 16th IFAC World Congress, Prague, Czech Republic, July 3-8, 2005*. Oxford, UK: Elsevier Ltd., 2005.
- [2] Ljung, L. *System Identification: Theory for the User*. Upper Saddle River, NJ: Prentice-Hall PTR, 1999.

## See Also

idfilt | oe

## Topics

"Loss Function and Model Quality Metrics"

**Introduced in R2012a**

## idnlarx/operspec

Construct operating point specification object for `idnlarx` model

### Syntax

```
spec = operspec(nlsys)
```

### Description

`spec = operspec(nlsys)` creates a default operating point specification object for the `idnlarx` model `nlsys`. This object is used with `findop` and specifies constraints on the model input and output signal values. Modify the default specifications using dot notation.

### Input Arguments

**nlsys** — Nonlinear ARX model

`idnlarx` object

Nonlinear ARX model, specified as an `idnlarx` object.

### Output Arguments

**spec** — Operating point specification

operating point specification object

Operating point specification, used to determine an operating point of the `idnlarx` model using `findop`, returned as an object containing the following:

- **Input** — Structure with fields:

Field	Description	Default for Each Input
Value	Initial guesses or fixed levels for the values of the model inputs, specified as a vector with length equal to the number of input signals.	0
Min	Minimum value constraints on the model inputs, specified as a vector with length equal to the number of input signals.	-Inf
Max	Maximum value constraints on the model inputs, specified as a vector with length equal to the number of input signals.	Inf
Known	Known value indicator, specified as a logical vector with length equal to the number of input signals and with the following values: <ul style="list-style-type: none"> <li>• true — findop will set the corresponding input signal to Value.</li> <li>• false — findop will estimate the corresponding input signal using Value as an initial guess.</li> </ul>	true

- Output — Structure with fields:

Field	Description	Default for Each Output
Value	Initial guesses for the values of the model outputs, specified as a vector with length equal to the number of output signals.	0
Min	Minimum value constraints on the model outputs, specified as a vector with length equal to the number of output signals.	-Inf
Max	Maximum value constraints on the model outputs, specified as a vector with length equal to the number of output signals.	Inf

## See Also

idnlarx/findop

**Introduced in R2008a**

# idnlhw/operspec

Construct operating point specification object for `idnlhw` model

## Syntax

```
spec = operspec(nlsys)
```

## Description

`spec = operspec(nlsys)` creates a default operating point specification object for the `idnlhw` model `nlsys`. This object is used with `findop` and specifies constraints on the model input and output signal values. Modify the default specifications using dot notation.

## Input Arguments

**nlsys** — Nonlinear Hammerstein-Wiener model

`idnlhw` object

Nonlinear Hammerstein-Wiener model, specified as an `idnlhw` object.

## Output Arguments

**spec** — Operating point specification

operating point specification object

Operating point specification, used to determine an operating point of the `idnlhw` model using `findop`, returned as an object containing the following:

- **Input** — Structure with fields:

Field	Description	Default for Each Input
Value	Initial guesses or fixed levels for the values of the model inputs, specified as a vector with length equal to the number of input signals.	0
Min	Minimum value constraints on the model inputs, specified as a vector with length equal to the number of input signals.	-Inf
Max	Maximum value constraints on the model inputs, specified as a vector with length equal to the number of input signals.	Inf
Known	Known value indicator, specified as a logical vector with length equal to the number of input signals and with the following values: <ul style="list-style-type: none"> <li>• true — findop will set the corresponding input signal to Value.</li> <li>• false — findop will estimate the corresponding input signal using Value as an initial guess.</li> </ul>	true

- Output — Structure with fields:

Field	Description	Default for Each Input
Value	Target values the model outputs, specified as a vector with length equal to the number of output signals.	0
Min	Minimum value constraints on the model outputs, specified as a vector with length equal to the number of output signals.	-Inf
Max	Maximum value constraints on the model outputs, specified as a vector with length equal to the number of output signals.	Inf

Field	Description	Default for Each Input
Known	<p>Known value indicator, specified as a logical vector with length equal to the number of output signals and with the following values:</p> <ul style="list-style-type: none"> <li>• <code>true</code> — <code>findop</code> will use <code>Value</code> as an estimation target for the corresponding output.</li> <li>• <code>false</code> — <code>findop</code> will keep the corresponding output within the constraints specified by <code>Min</code> and <code>Max</code>.</li> </ul>	false

---

### Note

- 1 If `Input.Known` is `true` for all model inputs, then the initial state values are determined using the input specifications only. In this case, `findop` ignores the specifications in the `Output` structure.
  - 2 Otherwise, `findop` uses the output specifications to meet the objectives indicated by `Output.Known`.
- 

## See Also

idnlhw/findop

**Introduced in R2008a**

## order

Query model order

## Syntax

```
NS = order(sys)
```

## Description

`NS = order(sys)` returns the model order `NS`. The order of a dynamic system model is the number of poles (for proper transfer functions) or the number of states (for state-space models). For improper transfer functions, the order is defined as the minimum number of states needed to build an equivalent state-space model (ignoring pole/zero cancellations).

`order(sys)` is an overloaded method that accepts SS, TF, and ZPK models. For LTI arrays, `NS` is an array of the same size listing the orders of each model in `sys`.

## Caveat

`order` does not attempt to find minimal realizations of MIMO systems. For example, consider this 2-by-2 MIMO system:

```
s=tf('s');  
h = [1, 1/(s*(s+1)); 1/(s+2), 1/(s*(s+1)*(s+2))];  
order(h)  
ans =
```

```
6
```

Although `h` has a 3rd order realization, `order` returns 6. Use

```
order(ss(h, 'min'))
```

to find the minimal realization order.

## **See Also**

balred | pole

**Introduced in R2012a**

## particleFilter

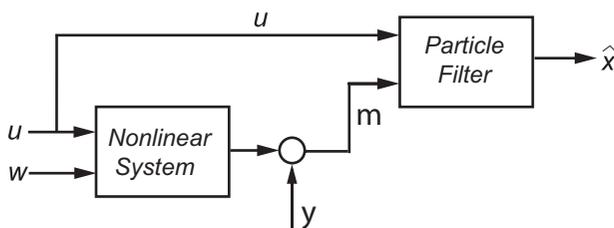
Particle filter object for online state estimation

### Description

A particle filter is a recursive, Bayesian state estimator that uses discrete particles to approximate the posterior distribution of an estimated state. It is useful for online state estimation when measurements and a system model, that relates model states to the measurements, are available. The particle filter algorithm computes the state estimates recursively and involves initialization, prediction, and correction steps.

`particleFilter` creates an object for online state estimation of a discrete-time nonlinear system using the discrete-time particle filter algorithm.

Consider a plant with states  $x$ , input  $u$ , output  $m$ , process noise  $w$ , and measurement  $y$ . Assume that you can represent the plant as a nonlinear system.



The algorithm computes the state estimates  $\hat{x}$  of the nonlinear system using the state transition and measurement likelihood functions you specify.

The software supports arbitrary nonlinear state transition and measurement models, with arbitrary process and measurement noise distributions.

To perform online state estimation, create the nonlinear state transition function and measurement likelihood function. Then construct the `particleFilter` object using these nonlinear functions. After you create the object:

- 1 Initialize the particles using the `initialize` command.

- 2 Predict state estimates at the next step using the `predict` command.
- 3 Correct the state estimates using the `correct` command.

The prediction step uses the latest state to predict the next state based on the state transition model you provide. The correction step uses the current sensor measurement to correct the state estimate. The algorithm optionally redistributes, or resamples, the particles in the state space to match the posterior distribution of the estimated state. Each particle represents a discrete state hypothesis of these state variables. The set of all particles is used to help determine the state estimate.

## Creation

## Syntax

```
pf = particleFilter(StateTransitionFcn,MeasurementLikelihoodFcn)
```

## Object Description

```
pf = particleFilter(StateTransitionFcn,MeasurementLikelihoodFcn)
```

creates a particle filter object for online state estimation of a discrete-time nonlinear system. `StateTransitionFcn` is a function that calculates the particles (state hypotheses) at the next time step, given the state vector at a time step. `MeasurementLikelihoodFcn` is a function that calculates the likelihood of each particle based on sensor measurements.

After creating the object, use the `initialize` command to initialize the particles with a known mean and covariance or uniformly distributed particles within defined bounds. Then, use the `correct` and `predict` commands to update particles (and hence the state estimate) using sensor measurements.

## Input Arguments

### **StateTransitionFcn** — State transition function

function handle

State transition function, specified as a function handle, determines the transition of particles (state hypotheses) between time steps. Also a property of the `particleFilter` object. For more information, see “Properties” on page 1-1098.

## **MeasurementLikelihoodFcn — Measurement likelihood function**

function handle

Measurement likelihood function, specified as a function handle, is used to calculate the likelihood of particles (state hypotheses) from sensor measurements. Also a property of the `particleFilter` object. For more information, see “Properties” on page 1-1098.

## **Properties**

### **NumStateVariables — Number of state variables**

[ ] (default) | scalar

Number of state variables, specified as a scalar. This property is read-only and is set using `initialize`. The number of states is implicit based on the specified matrices for the initial mean of particles, or the state bounds.

### **NumParticles — Number of particles used in the filter**

[ ] (default) | scalar

Number of particles used in the filter, specified as a scalar. Each particle represents a state hypothesis. You specify this property only by using `initialize`.

### **StateTransitionFcn — State transition function**

function handle

State transition function, specified as a function handle, determines the transition of particles (state hypotheses) between time steps. This function calculates the particles at the next time step, including the process noise, given particles at a time step.

In contrast, the state transition function for the `extendedKalmanFilter` and `unscentedKalmanFilter` generates a single state estimate at a given time step.

You write and save the state transition function for your nonlinear system, and specify it as a function handle when constructing the `particleFilter` object. For example, if `vdpParticleFilterStateFcn.m` is the state transition function, specify `StateTransitionFcn` as `@vdpParticleFilterStateFcn`. You can also specify `StateTransitionFcn` as a function handle to an anonymous function.

The function signature is as follows:

```
function predictedParticles = myStateTransitionFcn(previousParticles,varargin)
```

The `StateTransitionFcn` function accepts at least one input argument. The first argument is the set of particles `previousParticles` that represents the state hypotheses at the previous time step. The optional use of `varargin` in the function enables you to input any extra parameters that are relevant for predicting the next state, using `predict`, as follows:

```
predict(pf, arg1, arg2)
```

If `StateOrientation` is 'column', then `previousParticles` is a `NumStateVariables-by-NumParticles` array. If `StateOrientation` is 'row', then `previousParticles` is a `NumParticles-by-NumStateVariables` array.

`StateTransitionFcn` must return exactly one output, `predictedParticles`, which is the set of predicted particle locations for the current time step (array with same dimensions as `previousParticles`).

`StateTransitionFcn` must include the random process noise (from any distribution suitable for your application) in the `predictedParticles`.

To see an example of a state transition function with the `StateOrientation` property set to 'column', type `edit vdpParticleFilterStateFcn` at the command line.

### **MeasurementLikelihoodFcn — Measurement likelihood function**

`function handle`

Measurement likelihood function, specified as a function handle, is used to calculate the likelihood of particles (state hypotheses) using the sensor measurements. For each state hypothesis (particle), the function first calculates an N-element measurement hypothesis vector. Then the likelihood of each measurement hypothesis is calculated based on the sensor measurement and the measurement noise probability distribution.

In contrast, the measurement function for `extendedKalmanFilter` and `unscentedKalmanFilter` takes a single state hypothesis and returns a single measurement estimate.

You write and save the measurement likelihood function based on your measurement model, and use it to construct the object. For example, if `vdpMeasurementLikelihoodFcn.m` is the measurement likelihood function, specify

MeasurementLikelihoodFcn as @vdpMeasurementLikelihoodFcn. You can also specify MeasurementLikelihoodFcn as a function handle to an anonymous function.

The function signature is as follows:

```
function likelihood = myMeasurementLikelihoodFcn(predictedParticles,measurement,varargin)
```

The MeasurementLikelihoodFcn function accepts at least two input arguments. The first argument is the set of particles predictedParticles that represents the predicted state hypothesis. If StateOrientation is 'column', then predictedParticles is a NumStateVariables-by-NumParticles array. If StateOrientation is 'row', then predictedParticles is a NumParticles-by-NumStateVariables array. The second argument, measurement, is the N-element sensor measurement at the current time step. You can provide additional input arguments using varargin.

The MeasurementLikelihoodFcn must return exactly one output, likelihood, a vector with NumParticles length, which is the likelihood of the given measurement for each particle (state hypothesis).

To see an example of a measurement likelihood function, type edit vdpMeasurementLikelihoodFcn at the command line.

## **IsStateVariableCircular — Whether the state variables have a circular distribution**

[ ] (default) | logical array

Whether the state variables have a circular distribution, specified as a logical array.

This is a read-only property and is set using initialize.

Circular (or angular) distributions use a probability density function with a range of  $[-\pi, \pi]$ . IsStateVariableCircular is a row-vector with NumStateVariables elements. Each vector element indicates whether the associated state variable is circular.

## **ResamplingPolicy — Policy settings that determine when to trigger resampling**

particleResamplingPolicy object

Policy settings that determine when to trigger resampling, specified as a particleResamplingPolicy object.

The resampling of particles is a vital step in estimating states using a particle filter. It enables you to select particles based on the current state, instead of using the particle distribution given at initialization. By continuously resampling the particles around the

current estimate, you can get more accurate tracking and improve long-term performance.

You can trigger resampling either at fixed intervals or dynamically, based on the number of effective particles. The minimum effective particle ratio is a measure of how well the current set of particles approximates the posterior distribution. The number of effective particles is calculated by:

$$N_{eff} = \frac{1}{\sum_{i=1}^N (w^i)^2}$$

In this equation,  $N$  is the number of particles, and  $w$  is the normalized weight of each particle. The effective particle ratio is then  $N_{eff} / NumParticles$ . Therefore, the effective particle ratio is a function of the weights of all the particles. After the weights of the particles reach a low enough value, they are not contributing to the state estimation. This low value triggers resampling, so the particles are closer to the current state estimation and have higher weights.

The following properties of the `particleResamplingPolicy` object can be modified to control when resampling is triggered:

<b>Description</b>
<p><b>r</b> <b>p</b> <b>p</b> <b>e</b> <b>r</b> <b>t</b> <b>y</b></p> <p>It is a method to determine when resampling occurs, based on the value chosen. The 'interval' value triggers resampling at regular time steps of the particle filter operation. The 'ratio' value triggers resampling based on the ratio of effective total particles.</p> <p><b>r</b> <b>e</b> <b>t</b> <b>M</b> <b>d</b> <b>e</b> <b>h</b> <b>a</b> <b>d</b> <b>b</b> <b>t</b> <b>)</b> <b>,</b> <b>i</b> <b>n</b> <b>t</b> <b>e</b> <b>r</b> <b>v</b> <b>a</b> <b>l</b> <b>,</b></p>

Description
<p data-bbox="238 522 1338 656">Fixed interval between resampling, specified as a scalar. This interval determines during which correction steps the resampling is executed. For example, a value of 2 means the resampling is executed every second correction step. A value of <code>inf</code> means that resampling is never executed.</p> <p data-bbox="238 656 1338 713">This property only applies with the <code>TriggerMethod</code> is set to <code>'interval'</code>.</p>

Description
<p><b>M</b> <b>p</b> <b>p</b> <b>e</b> <b>r</b> <b>t</b> <b>y</b></p> <p><b>M</b> is the minimum desired ratio of the effective number of particles to the total number of particles <code>NumParticles</code>. The effective number of particles is a measure of how well the current set of particles approximates the posterior distribution. A lower effective particle ratio implies that a lower number of particles are contributing to the estimation and resampling is required.</p> <p><b>e</b> <b>f</b> <b>N</b>umParticles falls below the <code>MinEffectiveParticleRatio</code>, a resampling step is triggered.</p> <p><b>v</b> <b>e</b> <b>p</b> <b>a</b> <b>r</b> <b>t</b> <b>i</b> <b>c</b> <b>l</b> <b>e</b> <b>R</b> <b>a</b> <b>t</b> <b>i</b> <b>o</b></p>

**ResamplingMethod – Method used for particle resampling**

'multinomial' (default) | 'residual' | 'stratified' | 'systematic'

Method used for particle resampling, specified as one of the following:

- 'multinomial'

- 'residual'
- 'stratified'
- 'systematic'

### **StateEstimationMethod — Method used for extracting a state estimate from particles**

'mean' (default) | 'maxweight'

Method used for extracting a state estimate from particles, specified as one of the following:

- 'mean' - The object outputs the weighted mean of the particles, depending on the properties `Weights` and `Particles`, as the state estimate.
- 'maxweight' - The object outputs the particle with the highest weight as the state estimate.

### **Particles — Array of particle values**

[ ] (default) | array

Array of particle values, specified as an array based on the `StateOrientation` property:

- If `StateOrientation` is 'row' then `Particles` is an `NumParticles`-by-`NumStateVariables` array.
- If `StateOrientation` is 'column' then `Particles` is an `NumStateVariables`-by-`NumParticles` array.

Each row or column corresponds to a state hypothesis (a single particle).

### **Weights — Particle weights**

[ ] (default) | vector

Particle weights, defined as a vector based on the value of the `StateOrientation` property:

- If `StateOrientation` is 'row' then `Weights` is a `NumParticles`-by-1 vector, where each weight is associated with the particle in the same row in the `Particles` property.
- If `StateOrientation` is 'column' then `Weights` is a 1-by-`NumParticles` vector, where each weight is associated with the particle in the same column in the `Particles` property.

## State — Current state estimate

[ ] (default) | vector

Current state estimate, defined as a vector based on the value of the `StateOrientation` property:

- If `StateOrientation` is 'row' then `State` is a 1-by-`NumStateVariables` vector
- If `StateOrientation` is 'column' then `State` is a `NumStateVariables`-by-1 vector

`State` is a read-only property, and is derived from `Particles` based on the `StateEstimationMethod` property. Refer to “`StateEstimationMethod`” on page 1-0 for details on how the value of `State` is determined.

`State` along with `StateCovariance` can also be determined using `getStateEstimate`.

## StateCovariance — Current estimate of state estimation error covariance

`NumStateVariables`-by-`NumStateVariables` array (default) | [ ] | array

Current estimate of state estimation error covariance, defined as an `NumStateVariables`-by-`NumStateVariables` array. `StateCovariance` is a read-only property and is calculated based on the `StateEstimationMethod`. If you specify a state estimation method that does not support covariance, then the function returns `StateCovariance` as [ ].

`StateCovariance` and `State` can be determined together using `getStateEstimate`.

## Object Functions

<code>initialize</code>	Initialize the state of the particle filter
<code>predict</code>	Predict state and state estimation error covariance at next time step using extended or unscented Kalman filter, or particle filter
<code>correct</code>	Correct state and state estimation error covariance using extended or unscented Kalman filter, or particle filter and measurements
<code>getStateEstimate</code>	Extract best state estimate and covariance from particles
<code>clone</code>	Copy online state estimation object

## Examples

## Create Particle Filter Object for Online State Estimation

To create a particle filter object for estimating the states of your system, create appropriate state transition function and measurement likelihood function for the system.

In this example, the function `vdpParticleFilterStateFcn` describes a discrete-time approximation to van der Pol oscillator with nonlinearity parameter,  $\mu$ , equal to 1. In addition, it models Gaussian process noise. `vdpMeasurementLikelihood` function calculates the likelihood of particles from the noisy measurements of the first state, assuming a Gaussian measurement noise distribution.

Create the particle filter object. Use function handles to provide the state transition and measurement likelihood functions to the object.

```
myPF = particleFilter(@vdpParticleFilterStateFcn,@vdpMeasurementLikelihoodFcn);
```

To initialize and estimate the states and state estimation error covariance from the constructed object, use the `initialize`, `predict`, and `correct` commands.

## Estimate States Online using Particle Filter

Load the van der Pol ODE data, and specify the sample time.

`vdpODEdata.mat` contains a simulation of the van der Pol ODE with nonlinearity parameter  $\mu=1$ , using `ode45`, with initial conditions  $[2;0]$ . The true state was extracted with sample time  $dt = 0.05$ .

```
addpath(fullfile(matlabroot,'examples','ident','main')) % add example data
```

```
load ('vdpODEdata.mat','xTrue','dt')
tSpan = 0:dt:5;
```

Get the measurements. For this example, a sensor measures the first state with a Gaussian noise with standard deviation  $0.04$ .

```
sqrtr = 0.04;
yMeas = xTrue(:,1) + sqrtr*randn(numel(tSpan),1);
```

Create a particle filter, and set the state transition and measurement likelihood functions.

```
myPF = particleFilter(@vdpParticleFilterStateFcn,@vdpMeasurementLikelihoodFcn);
```

Initialize the particle filter at state  $[2; 0]$  with unit covariance, and use 1000 particles.

```
initialize(myPF,1000,[2;0],eye(2));
```

Pick the mean state estimation and systematic resampling methods.

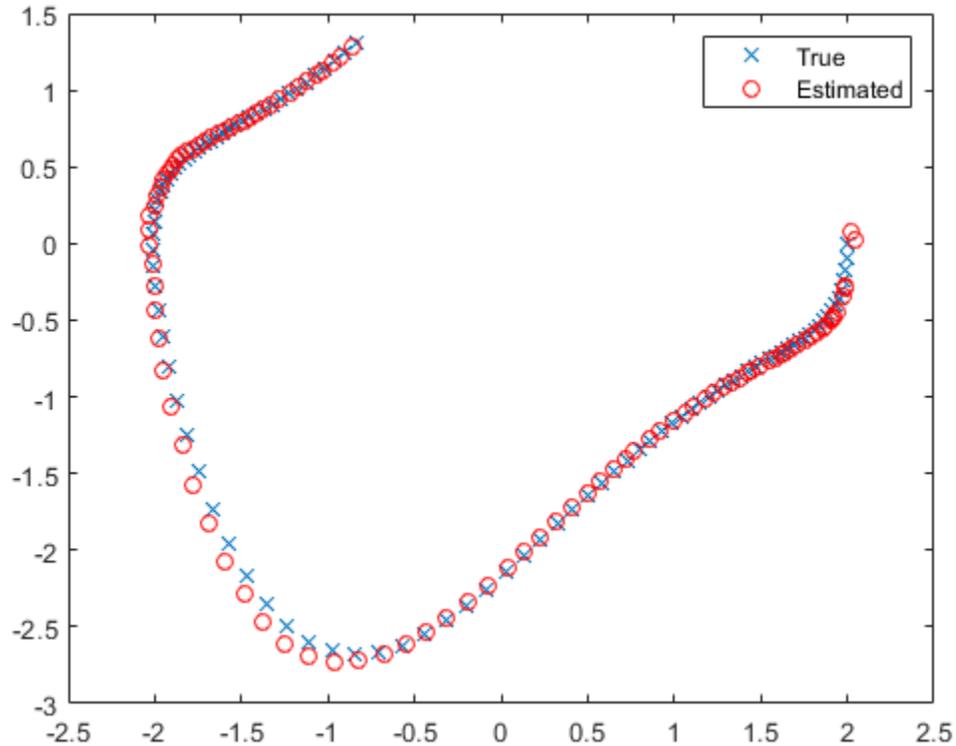
```
myPF.StateEstimationMethod = 'mean';  
myPF.ResamplingMethod = 'systematic';
```

Estimate the states using the `correct` and `predict` commands, and store the estimated states.

```
xEst = zeros(size(xTrue));  
for k=1:size(xTrue,1)  
    xEst(k,:) = correct(myPF,yMeas(k));  
    predict(myPF);  
end
```

Plot the results, and compare the estimated and true states.

```
figure(1)  
plot(xTrue(:,1),xTrue(:,2),'x',xEst(:,1),xEst(:,2),'ro')  
legend('True','Estimated')
```



```
rmpath(fullfile(matlabroot, 'examples', 'ident', 'main')) % remove example data
```

## Extended Capabilities

### C/C++ Code Generation

Generate C and C++ code using MATLAB® Coder™.

Usage notes and limitations:

For more information, see “Generate Code for Online State Estimation in MATLAB”.

Supports MATLAB Function block: No

## See Also

### Functions

`clone` | `correct` | `extendedKalmanFilter` | `initialize` | `predict` | `unscentedKalmanFilter`

### Topics

“Nonlinear State Estimation Using Unscented Kalman Filter and Particle Filter”

“Generate Code for Online State Estimation in MATLAB”

“What Is Online Estimation?”

“Validate Online State Estimation at the Command Line”

“Troubleshoot Online State Estimation”

**Introduced in R2017b**

## pe

Prediction error for identified model

## Syntax

```
err = pe(sys,data,K)
err = pe(sys,data,K,opt)
[err,x0e,sys_pred] = pe( ___ )
pe(sys,data,K, ___ )
pe(sys,Linespec,data,K, ___ )
pe(sys1,...,sysN,data,K, ___ )
pe(sys1,Linespec1,...,sysN,LinespecN,data,K, ___ )
```

## Description

`err = pe(sys,data,K)` returns the K-step prediction error for the output of the identified model `sys`. The prediction error is determined by subtracting the K-step ahead predicted response from the measured output. The prediction error is calculated for the time span covered by `data`. For more information on the computation of predicted response, see `predict`.

`err = pe(sys,data,K,opt)` returns the prediction error using the option set, `opt`, to specify prediction error calculation behavior.

`[err,x0e,sys_pred] = pe( ___ )` also returns the estimated initial state, `x0e`, and a predictor system, `sys_pred`.

`pe(sys,data,K, ___ )` plots the prediction error. Use with any of the previous input argument combinations. To change display options in the plot, right-click the plot to access the context menu. For more details about the menu, see “Tips” on page 1-1116.

`pe(sys,Linespec,data,K, ___ )` uses `Linespec` to specify the line type, marker symbol, and color.

`pe(sys1,...,sysN,data,K, ___ )` plots the prediction errors for multiple identified models. `pe` automatically chooses colors and line styles.

`pe(sys1,Linespec1,...,sysN,LinespecN,data,K, ___)` uses the line type, marker symbol, and color specified for each model.

## Input Arguments

### **sys**

Identified model.

### **data**

Measured input-output history.

If `sys` is a time-series model, which has no input signals, then specify `data` as an `iddata` object with no inputs. In this case, you can also specify `data` as a matrix of the past time-series values.

### **K**

Prediction horizon.

Specify `K` as a positive integer that is a multiple of the data sample time. Use `K = Inf` to compute the pure simulation error.

**Default:** 1

### **opt**

Prediction options.

`opt` is an option set, created using `peOptions`, that configures the computation of the predicted response. Options that you can specify include:

- Handling of initial conditions
- Data offsets

### **Linespec**

Line style, marker, and color

Line style, marker, and color, specified as a character vector. For example, `'b'` or `'b+:'`.

For more information about configuring `Linespec`, see `plot`.

## Output Arguments

### **err**

Prediction error.

`err` is returned as an `iddata` object or matrix, depending on how you specify `data`. For example, if `data` is an `iddata` object, then so is `err`.

Outputs up to the time `t-K` and inputs up to the time instant `t` are used to calculate the prediction error at the time instant `t`.

When `K = Inf`, the predicted output is a pure simulation of the system.

For multi-experiment data, `err` contains the prediction error data for each experiment. The time span of the prediction error matches that of the observed data.

### **x0e**

Estimated initial states.

`x0e` is returned only for state-space systems.

### **sys\_pred**

Predictor system.

`sys_pred` is a dynamic system. When you simulate `sys_pred`, using `[data.OutputData data.InputData]` as the input, the output, `yp`, is such that `err.OutputData = data.OutputData - yp`. For state-space models, the software uses `x0e` as the initial condition when simulating `sys_pred`.

For discrete-time data, `sys_pred` is always a discrete-time model.

For multi-experiment data, `sys_pred` is an array of models, with one entry for each experiment.

## Examples

### Compute Prediction Error for an ARIX Model

Compute the prediction error for an ARIX model.

Use the error data to compute the variance of the noise source  $e(t)$ .

Obtain noisy data.

```
noise = [(1:150)';(151:-1:2)'];
```

```
load iddata1 z1;  
z1.y = z1.y+noise;
```

`noise` is a triangular wave that is added to the output signal of `z1`, an `iddata` object.

Estimate an ARIX model for the noisy data.

```
sys = arx(z1,[2 2 1], 'IntegrateNoise', true);
```

Compute the prediction error of the estimated model.

```
K = 1;  
err = pe(z1,sys,K);
```

`pe` computes the one-step prediction error for the output of the identified model, `sys`.

Compute the variance of the noise source,  $e(t)$ .

```
noise_var = err.y'*err.y/(299-nparams(sys)-order(sys));
```

Compare the computed value with model's noise variance.

```
sys.NoiseVariance
```

The output of `sys.NoiseVariance` matches the computed variance.

### Plot Prediction Error for Multiple Models

Load the estimation data.

```
load iddata1;  
data = z1;
```

Estimate an ARX model of order [2 2 1].

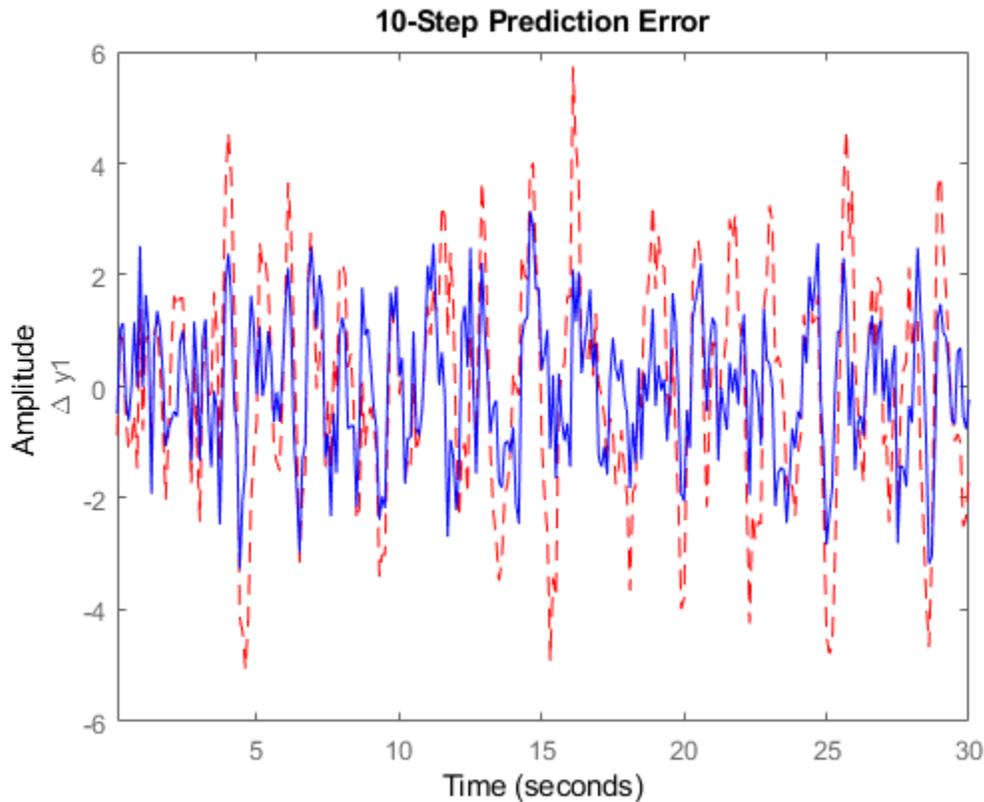
```
sys1 = arx(data,[2 2 1]);
```

Estimate a transfer function with 2 poles.

```
sys2 = tfest(data,2);
```

Plot the prediction error for the estimated models. Specify prediction horizon as 10, and specify the line styles for plotting the prediction error of each system.

```
pe(sys1, 'r--', sys2, 'b', data, 10);
```



To change the display options, right-click the plot to access the context menu. For example, to view the estimation data, select **Show Validation Data** from the context menu. To view the predicted outputs, select **Predicted Response Plot**.

## Tips

- Right-clicking the plot of the prediction error opens the context menu, where you can access the following options:
  - **Systems** — Select systems to view prediction error. By default, the prediction error of all systems is plotted.

- **Data Experiment** — For multi-experiment data only. Toggle between data from different experiments.
- **Characteristics** — View the following data characteristics:
  - **Peak Value** — View the absolute peak value of the data. Applicable for time-domain data only.
  - **Peak Response** — View peak response of the data. Applicable for frequency-response data only.
  - **Mean Value** — View mean value of the data. Applicable for time-domain data only.
- **Show** — For frequency-domain and frequency-response data only.
  - **Magnitude** — View magnitude of frequency response of the system.
  - **Phase** — View phase of frequency response of the system.
- **Show Validation Data** — Plot data used to compute the prediction error.
- **I/O Grouping** — For datasets containing more than one input or output channel. Select grouping of input and output channels on the plot.
  - **None** — Plot input-output channels in their own separate axes.
  - **All** — Group all input channels together and all output channels together.
- **I/O Selector** — For datasets containing more than one input or output channel. Select a subset of the input and output channels to plot. By default, all output channels are plotted.
- **Grid** — Add grids to the plot.
- **Normalize** — Normalize the y-scale of all data in the plot.
- **Full View** — Return to full view. By default, the plot is scaled to full view.
- **Prediction Horizon** — Set the prediction horizon, or choose simulation.
- **Initial Condition** — Specify handling of initial conditions. Not applicable for frequency-response data.

Specify as one of the following:

- **Estimate** — Treat the initial conditions as estimation parameters.
- **Zero** — Set all initial conditions to zero.
- **Absorb delays and estimate** — Absorb nonzero delays into the model coefficients and treat the initial conditions as estimation parameters. Use this option for discrete-time models only.

- **Predicted Response Plot** — Plot the predicted model response.
- **Prediction Error Plot** — Plot the error between the model response and prediction data. By default, the error plot is shown.
- **Properties** — Open the Property Editor dialog box to customize plot attributes.

## See Also

ar | arx | compare | iddata | n4sid | peOptions | predict | resid | sim

**Introduced before R2006a**

# peOptions

Option set for pe

## Syntax

```
opt = peOptions  
opt = peOptions(Name,Value)
```

## Description

`opt = peOptions` creates the default options set for `pe`.

`opt = peOptions(Name,Value)` creates an option set with the options specified by one or more `Name,Value` pair arguments.

## Input Arguments

### Name-Value Pair Arguments

Specify optional comma-separated pairs of `Name,Value` arguments. `Name` is the argument name and `Value` is the corresponding value. `Name` must appear inside quotes. You can specify several name and value pair arguments in any order as `Name1,Value1,...,NameN,ValueN`.

### InitialCondition

Handling of initial conditions.

Specify `InitialCondition` as one of the following:

- 'z' — Zero initial conditions.
- 'e' — Estimate initial conditions such that the prediction error for observed output is minimized.

For nonlinear grey-box models, only those initial states *i* that are designated as free in the model (`sys.InitialStates(i).Fixed = false`) are estimated. To estimate all the states of the model, first specify all the *N<sub>x</sub>* states of the `idnlgrey` model `sys` as free.

```
for i = 1:Nx
    sys.InitialStates(i).Fixed = false;
end
```

Similarly, to fix all the initial states to values specified in `sys.InitialStates`, first specify all the states as fixed in the `sys.InitialStates` property of the nonlinear grey-box model.

- 'd' — Similar to 'e', but absorbs nonzero delays into the model coefficients. The delays are first converted to explicit model states, and the initial values of those states are also estimated and returned.

Use this option for linear models only.

- Vector or Matrix — Initial guess for state values, specified as a numerical column vector of length equal to the number of states. For multi-experiment data, specify a matrix with *N<sub>e</sub>* columns, where *N<sub>e</sub>* is the number of experiments. Otherwise, use a column vector to specify the same initial conditions for all experiments. Use this option for state-space (`idss` and `idgrey`) and nonlinear models (`idnlarx`, `idnlhw`, and `idnlgrey`) only.
- Structure with the following fields, which contain the historical input and output values for a time interval immediately before the start time of the data used by `pe`:

Field	Description
Input	Input history, specified as a matrix with <i>N<sub>u</sub></i> columns, where <i>N<sub>u</sub></i> is the number of input channels. For time series models, use [ ]. The number of rows must be greater than or equal to the model order.
Output	Output history, specified as a matrix with <i>N<sub>y</sub></i> columns, where <i>N<sub>y</sub></i> is the number of output channels. The number of rows must be greater than or equal to the model order.

For multi-experiment data, configure the initial conditions separately for each experiment by specifying `InitialCondition` as a structure array with *N<sub>e</sub>* elements. To specify the same initial conditions for all experiments, use a single structure.

The software uses `data2state` to map the historical data to states. If your model is not `idss`, `idgrey`, `idnlgrey`, or `idnlarx`, the software first converts the model to its

state-space representation and then maps the data to states. If conversion of your model to `idss` is not possible, the estimated states are returned empty.

- `x0obj` — Specification object created using `idpar`. Use this object for discrete-time state-space (`idss` and `idgrey`) and nonlinear grey-box (`idnlgrey`) models only. Use `x0obj` to impose constraints on the initial states by fixing their value or specifying minimum or maximum bounds.

**Default:** 'e'

### **InputOffset**

Removes offset from time domain input data during prediction-error calculation.

Specify as a column vector of length  $N_u$ , where  $N_u$  is the number of inputs.

For multi-experiment data, specify `InputOffset` as an  $N_u$ -by- $N_e$  matrix.  $N_u$  is the number of inputs, and  $N_e$  is the number of experiments.

Each entry specified by `InputOffset` is subtracted from the corresponding input data.

Specify input offset for only time domain data.

**Default:** []

### **OutputOffset**

Removes offset from time domain output data during prediction-error calculation.

Specify as a column vector of length  $N_y$ , where  $N_y$  is the number of outputs.

In case of multi-experiment data, specify `OutputOffset` as a  $N_y$ -by- $N_e$  matrix.  $N_y$  is the number of outputs, and  $N_e$  is the number of experiments.

Each entry specified by `OutputOffset` is subtracted from the corresponding output data.

Specify output offset for only time domain data.

**Default:** []

### **OutputWeight**

Weight of output for initial condition estimation.

OutputWeight takes one of the following:

- [] — No weighting is used. This value is the same as using `eye(Ny)` for the output weight, where `Ny` is the number of outputs.
- 'noise' — Inverse of the noise variance stored with the model.
- matrix — A positive, semidefinite matrix of dimension `Ny-by-Ny`, where `Ny` is the number of outputs.

**Default:** []

## Output Arguments

**opt**

Option set containing the specified options for `pe`.

## Examples

### Create Default Options Set for Prediction-Error Calculation

```
opt = peOptions;
```

### Specify Options for Prediction-Error Calculation

Create an options set for `pe` using zero initial conditions, and set the input offset to 5.

```
opt = peOptions('InitialCondition','z','InputOffset',5);
```

Alternatively, use dot notation to set the values of `opt`.

```
opt = peOptions;  
opt.InitialCondition = 'z';  
opt.InputOffset = 5;
```

## **See Also**

idpar | pe

**Introduced in R2012a**

## pem

Prediction error estimate for linear and nonlinear model

### Syntax

```
sys = pem(data,init_sys)
sys = pem(data,init_sys,opt)
```

### Description

`sys = pem(data,init_sys)` updates the parameters of an initial model to fit the estimation data. The function uses prediction-error minimization algorithm to update the parameters of the initial model. Use this command to refine the parameters of a previously estimated model.

`sys = pem(data,init_sys,opt)` specifies estimation options using an option set.

### Examples

#### Refine Estimated State-Space Model

Estimate a discrete-time state-space model using the subspace method. Then, refine it by minimizing the prediction error.

Estimate a discrete-time state-space model using `n4sid`, which applies the subspace method.

```
load iddata7 z7;
z7a = z7(1:300);
opt = n4sidOptions('Focus','simulation');
init_sys = n4sid(z7a,4,opt);
```

`init_sys` provides a 73.85% fit to the estimation data.

```
init_sys.Report.Fit.FitPercent
```

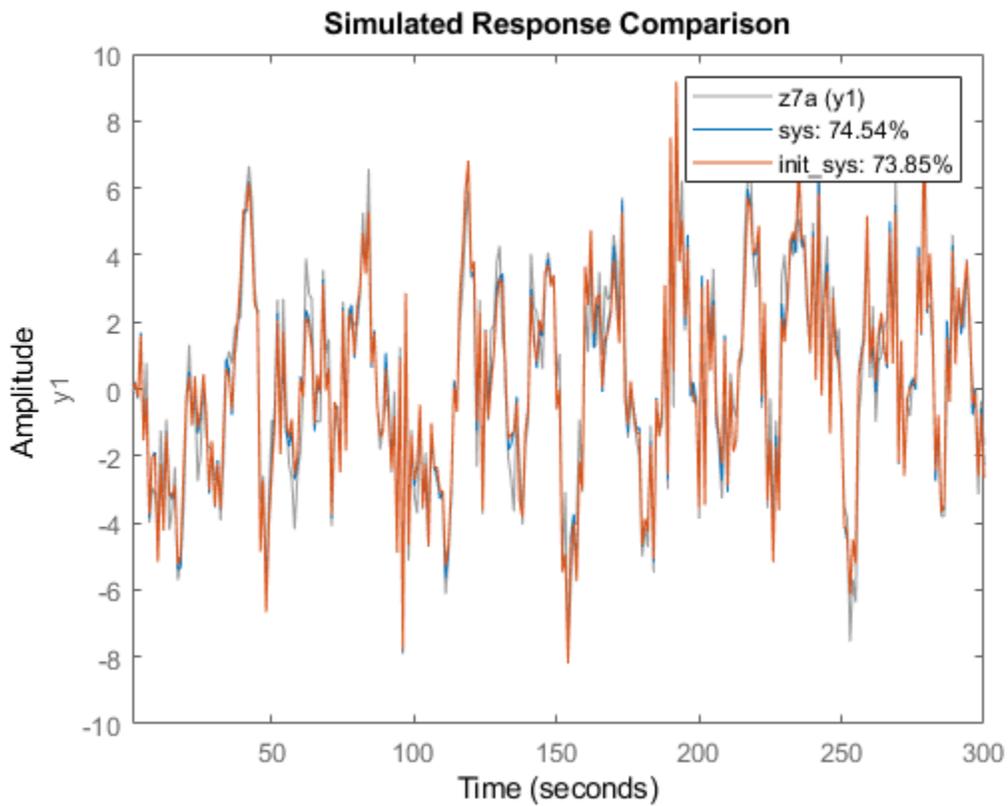
```
ans = 73.8490
```

Use `pem` to improve the closeness of the fit.

```
sys = pem(z7a,init_sys);
```

Analyze the results.

```
compare(z7a,sys,init_sys);
```



`sys` provides a 74.54% fit to the estimation data.

## Estimate Nonlinear Grey-Box Model

Estimate the parameters of a nonlinear grey-box model to fit DC motor data.

Load the experimental data, and specify the signal attributes such as start time and units.

```
load(fullfile(matlabroot, 'toolbox', 'ident', 'iddemos', 'data', 'dcmotordata'));  
data = iddata(y, u, 0.1);  
data.Tstart = 0;  
data.TimeUnit = 's';
```

Configure the nonlinear grey-box model (`idnlgrey`) model.

For this example, use `dcmotor_m.m` file. To view this file, type `edit dcmotor_m.m` at the MATLAB® command prompt.

```
file_name = 'dcmotor_m';  
order = [2 1 2];  
parameters = [1;0.28];  
initial_states = [0;0];  
Ts = 0;  
init_sys = idnlgrey(file_name,order,parameters,initial_states,Ts);  
init_sys.TimeUnit = 's';
```

```
setinit(init_sys, 'Fixed', {false false});
```

`init_sys` is a nonlinear grey-box model with its structure described by `dcmotor_m.m`. The model has one input, two outputs and two states, as specified by `order`.

`setinit(init_sys, 'Fixed', {false false})` specifies that the initial states of `init_sys` are free estimation parameters.

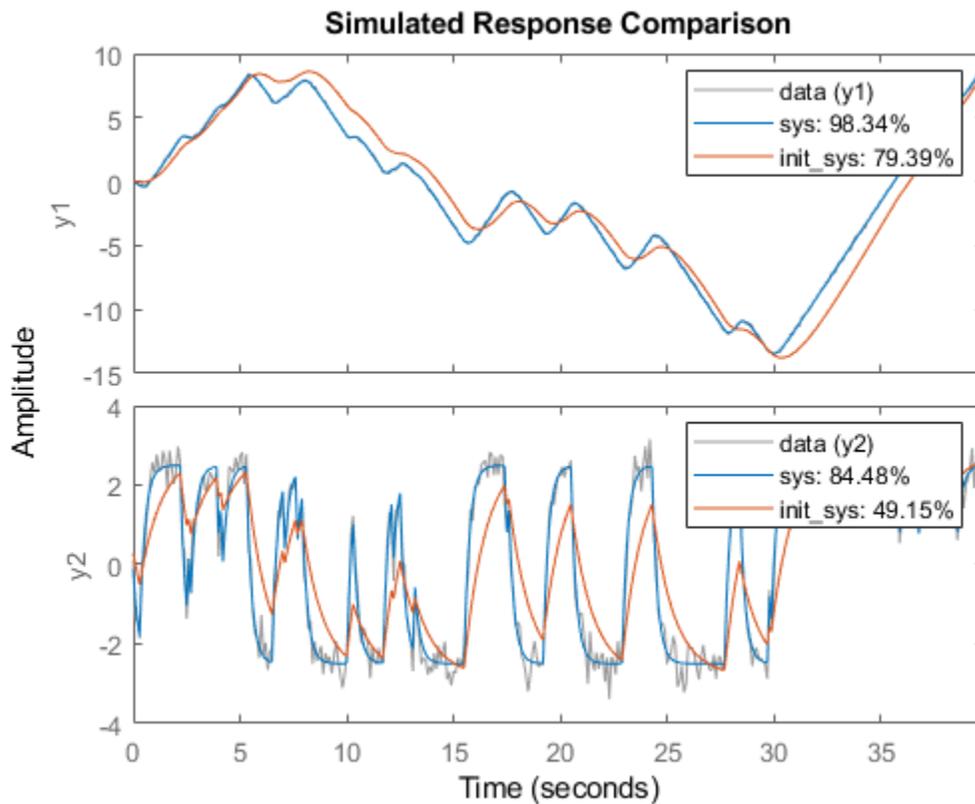
Estimate the model parameters and initial states.

```
sys = pem(data,init_sys);
```

`sys` is an `idnlgrey` model, which encapsulates the estimated parameters and their covariance.

Analyze the estimation result.

```
compare(data,sys,init_sys);
```



sys provides a 98.34% fit to the estimation data.

### Configure Estimation Using Process Model

Create a process model structure and update its parameter values to minimize prediction error.

Initialize the coefficients of a process model.

```
init_sys = idproc('P2UDZ');  
init_sys.Kp = 10;
```

```
init_sys.Tw = 0.4;  
init_sys.Zeta = 0.5;  
init_sys.Td = 0.1;  
init_sys.Tz = 0.01;
```

The Kp, Tw, Zeta, Td, and Tz coefficients of `init_sys` are configured with their initial guesses.

Use `init_sys` to configure the estimation of a prediction error minimizing model using measured data. Because `init_sys` is an `idproc` model, use `procestOptions` to create the option set.

```
load iddata1 z1;  
opt = procestOptions('Display','on','SearchMethod','lm');  
sys = pem(z1,init_sys,opt);
```

Examine the model fit.

```
sys.Report.Fit.FitPercent
```

```
ans = 70.6330
```

`sys` provides a 70.63% fit to the measured data.

## Input Arguments

### **data** — Estimation data

`iddata` | `idfrd`

Estimation data that contains measured input-output data, specified as an `iddata` or `idfrd` object. You can use frequency-domain data only when `init_sys` is a linear model.

The input-output dimensions of `data` and `init_sys` must match.

### **init\_sys** — Identified model that configures the initial parameterization of `sys`

linear model | nonlinear model

Identified model that configures the initial parameterization of `sys`, specified as a linear, or nonlinear model. You can obtain `init_sys` by performing an estimation using measured data or by direct construction.

`init_sys` must have finite parameter values. You can configure initial guesses, specify minimum/maximum bounds, and fix or free for estimating any parameter of `init_sys`:

- For linear models, use the `Structure` property. For more information, see “Imposing Constraints on Model Parameter Values”.
- For nonlinear grey-box models, use the `InitialStates` and `Parameters` properties. Parameter constraints cannot be specified for nonlinear ARX and Hammerstein-Wiener models.

### **opt** — Estimation options

option set

Estimation options that configure the algorithm settings, handling of estimation focus, initial conditions, and data offsets, specified as an option set. The command used to create the option set depends on the initial model type:

<b>Model Type</b>	<b>Use</b>
<code>idss</code>	<code>ssestOptions</code>
<code>idtf</code>	<code>tfestOptions</code>
<code>idproc</code>	<code>procestOptions</code>
<code>idpoly</code>	<code>polyestOptions</code>
<code>idgrey</code>	<code>greyestOptions</code>
<code>idnlarx</code>	<code>nlarxOptions</code>
<code>idnlhw</code>	<code>nlhwOptions</code>
<code>idnlgrey</code>	<code>nlgreyestOptions</code>

## **Output Arguments**

### **sys** — Identified model

linear model | nonlinear model

Identified model, returned as the same model type as `init_sys`. The model is obtained by estimating the free parameters of `init_sys` using the prediction error minimization algorithm.

## Algorithms

PEM uses numerical optimization to minimize the *cost function*, a weighted norm of the prediction error, defined as follows for scalar outputs:

$$V_N(G, H) = \sum_{t=1}^N e^2(t)$$

where  $e(t)$  is the difference between the measured output and the predicted output of the model. For a linear model, the error is defined as:

$$e(t) = H^{-1}(q)[y(t) - G(q)u(t)]$$

where  $e(t)$  is a vector and the cost function  $V_N(G, H)$  is a scalar value. The subscript  $N$  indicates that the cost function is a function of the number of data samples and becomes more accurate for larger values of  $N$ . For multiple-output models, the previous equation is more complex. For more information, see chapter 7 in *System Identification: Theory for the User*, Second Edition, by Lennart Ljung, Prentice Hall PTR, 1999.

## Alternative Functionality

You can achieve the same results as `pem` by using dedicated estimation commands for the various model structures. For example, use `ssest(data, init_sys)` for estimating state-space models.

## See Also

`armax` | `bj` | `greyest` | `n4sid` | `nlarx` | `nlgreyest` | `nlhw` | `oe` | `polyest` | `procest` | `ssest` | `tfest`

## Topics

“Refine Linear Parametric Models”

**Introduced before R2006a**

# pexcit

Level of excitation of input signals

## Syntax

```
Ped = pexcit(Data)  
[Ped,Maxnr] = pexcit(Data,Maxnr,Threshold)
```

## Description

`Ped = pexcit(Data)` tests the degree of persistence of excitation for the input. `Data` is an `iddata` object with time- or frequency-domain signals. `Ped` is the degree or order of excitation of the inputs in `Data` and is a row vector of integers with as many components as there are inputs in `Data`. The intuitive interpretation of the degree of excitation in an input is the order of a model that the input is capable of estimating in an unambiguous way.

`[Ped,Maxnr] = pexcit(Data,Maxnr,Threshold)` specifies the maximum order tested and threshold level used to measure which singular values are significant. Default value of `Maxnr` is  $\min(N/3, 50)$ , where `N` is the number of input data. Default value of `Threshold` is  $1e-9$ .

## References

Section 13.2 in Ljung (1999).

## See Also

[advice](#) | [feedback](#) | [iddata](#) | [idnlarx](#)

**Introduced before R2006a**

## iddata/plot

Plot input-output data

### Syntax

```
plot(data)
plot(data,LineStyle)
plot(data1,...,dataN)
plot(data1,LineStyle1...,dataN,LineStyleN)
```

```
plot(axes_handle, ___)
```

```
plot(___,plotoptions)
```

```
h = plot(___)
```

### Description

`plot(data)` plots the input and output channels of an `iddata` object. The outputs are plotted on the top axes and the inputs on the bottom axes.

- For time-domain data, the input and output signals are plotted as a function of time. Depending on the `InterSample` property of the `iddata` object, the input signals are plotted as linearly interpolated curves or as staircase plots. For example, if `data.InterSample = 'zoh'`, the input is piecewise constant between sampling points, and it is then plotted accordingly.
- For frequency-domain data, the magnitude and phase of each input and output signal is plotted over the available frequency span.

To plot a subselection of the data, use subreferencing:

- `plot(data(201:300))` plots the samples 201 to 300 in the dataset `data`.
- `plot(data(201:300,'Altitude',{'Angle_of_attack','Speed'}))` plots the chosen samples of output named `Altitude` and inputs named `Angle_of_attack` and `Speed`.

- `plot(data(:, [3 4], [3:7]))` plots all samples of output channel numbers 3 and 4 and input numbers 3 through 7.

`plot(data, LineSpec)` specifies the color, line style and marker symbol for the dataset.

`plot(data1, ..., dataN)` plots multiple datasets. The number of plot axes are determined by the number of unique input and output names among all the datasets.

`plot(data1, LineSpec1, ..., dataN, LineSpecN)` specifies the line style, marker type, and color for each dataset. You can mix `data, LineSpec` pairs with `data`. For example, `plot(data1, data2, LineSpec2, data3)`.

`plot(axes_handle, ___)` plots into the axes with handle `axes_handle` instead of into the current axes (`gca`). This syntax can include any of the input argument combinations in the previous syntaxes.

`plot(___, plotoptions)` specifies the plot options. This syntax can include any of the input argument combinations in the previous syntaxes.

`h = plot(___)` returns the handle to the plot. You can use this handle to customize the plot with `getoptions` and `setoptions`.

## Input Arguments

### **data** — Input-output data

iddata object

Input-output data, specified as an `iddata` object. The data can be time-domain or frequency-domain. It can be a single- or multi-channel data, or single- or multi-experiment data.

### **LineSpec** — Line style, marker symbol, and color

character vector

Line style, marker symbol, and color, specified as a character vector. `LineSpec` takes values such as `'b'`, `'b+:'`. For more information, see the `plot` reference page in the MATLAB documentation.

Data Types: `char`

### **axes\_handle** — Axes handle

handle

Axes handle, which is the reference to an axes object. Use the `gca` function to get the handle to the current axes, for example, `axes_handle= gca`.

## **plotoptions — Plot options**

structure

Plot options, specified as an option set created using `iddataPlotOptions`.

## **Output Arguments**

### **h — Lineseries handle**

scalar | vector

Lineseries handle, returned as a scalar or vector. These are unique identifiers, which you can use to query and modify properties of a specific plot.

## **Examples**

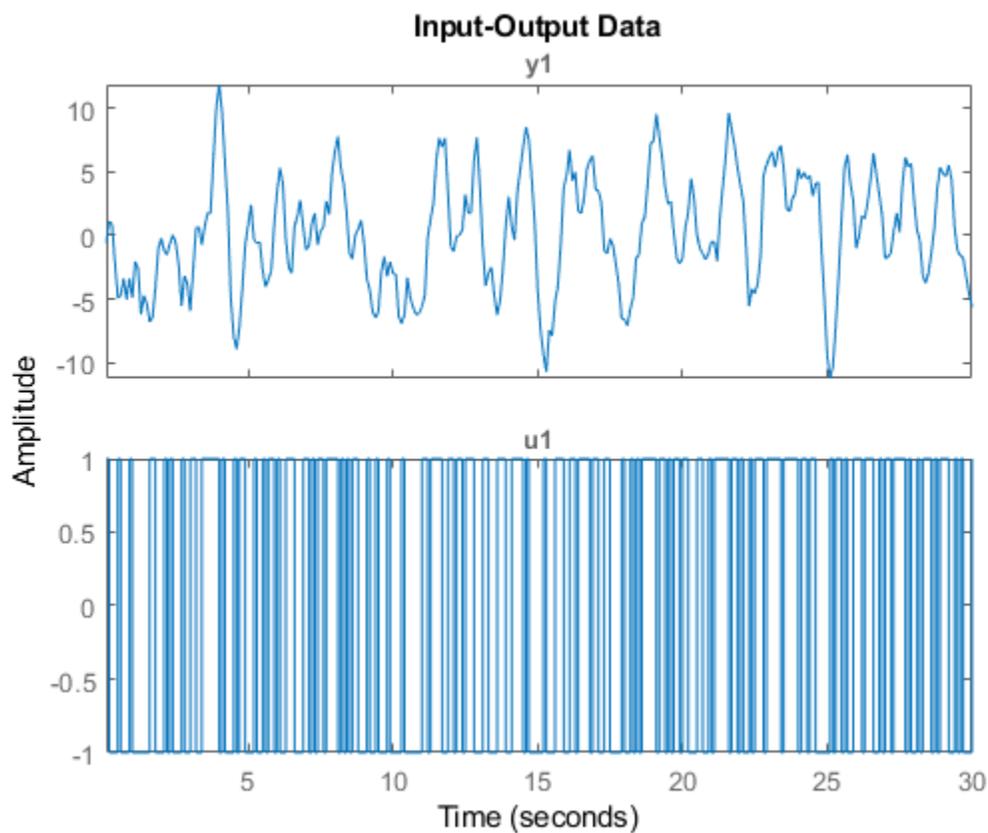
### **Plot Time-Domain Input-Output Data**

Load the data.

```
load iddata1 z1;
```

Plot the data.

```
plot(z1)
```



The output is plotted on the top axes and the input on the bottom axes.

You can right-click the plot to explore its characteristics such as peak and mean values.

### Plot Frequency-Domain Input-Output Data

Load the data.

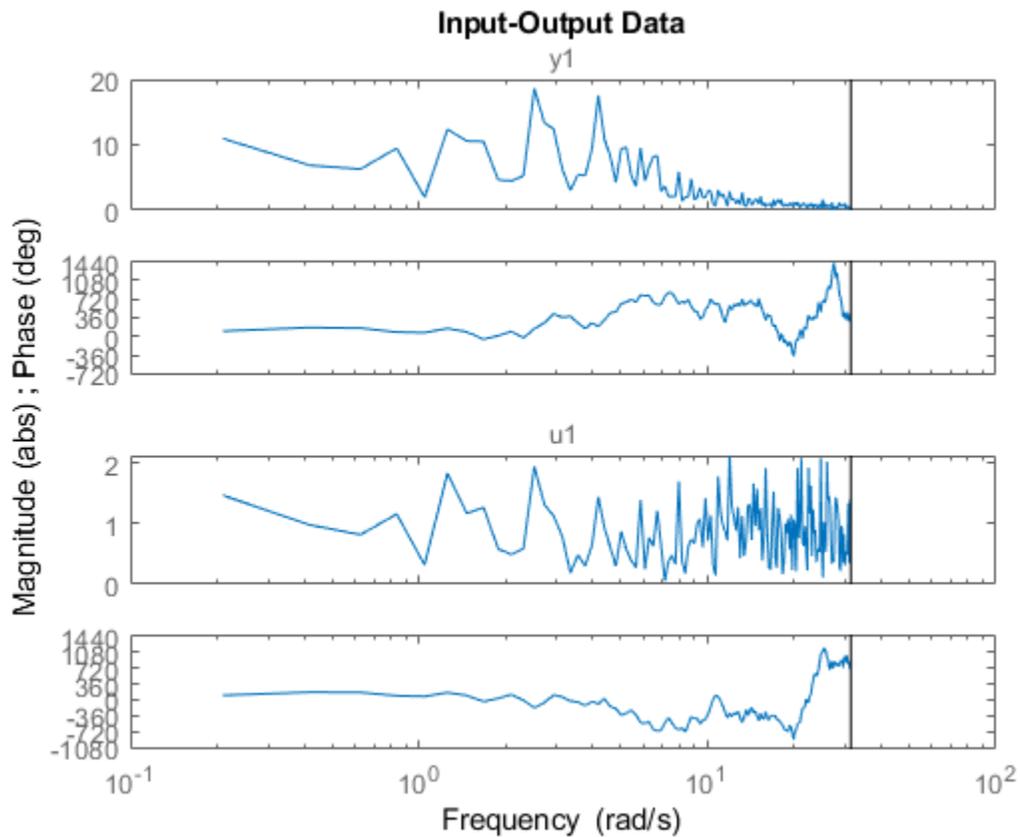
```
load iddata1 z1
```

Convert the data to frequency domain.

```
zf = fft(z1);
```

Plot the data.

```
plot(zf);
```



### Plot Input Data, Output Data and Input-Output Data

Generate input data.

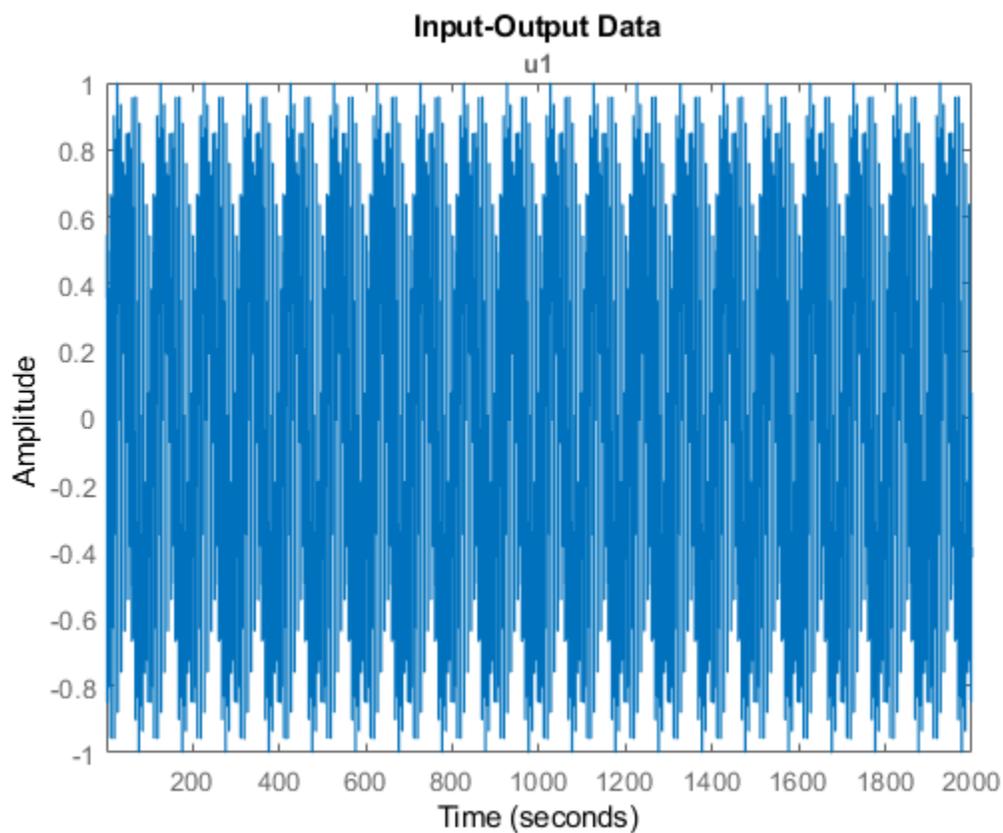
```
u = idinput([100 1 20], 'sine', [], [], [5 10 1]);  
u = iddata([], u, 1, 'per', 100);
```

Generate output data.

```
sys = idtf(1,[1 2 1]);  
y = sim(sys,u);
```

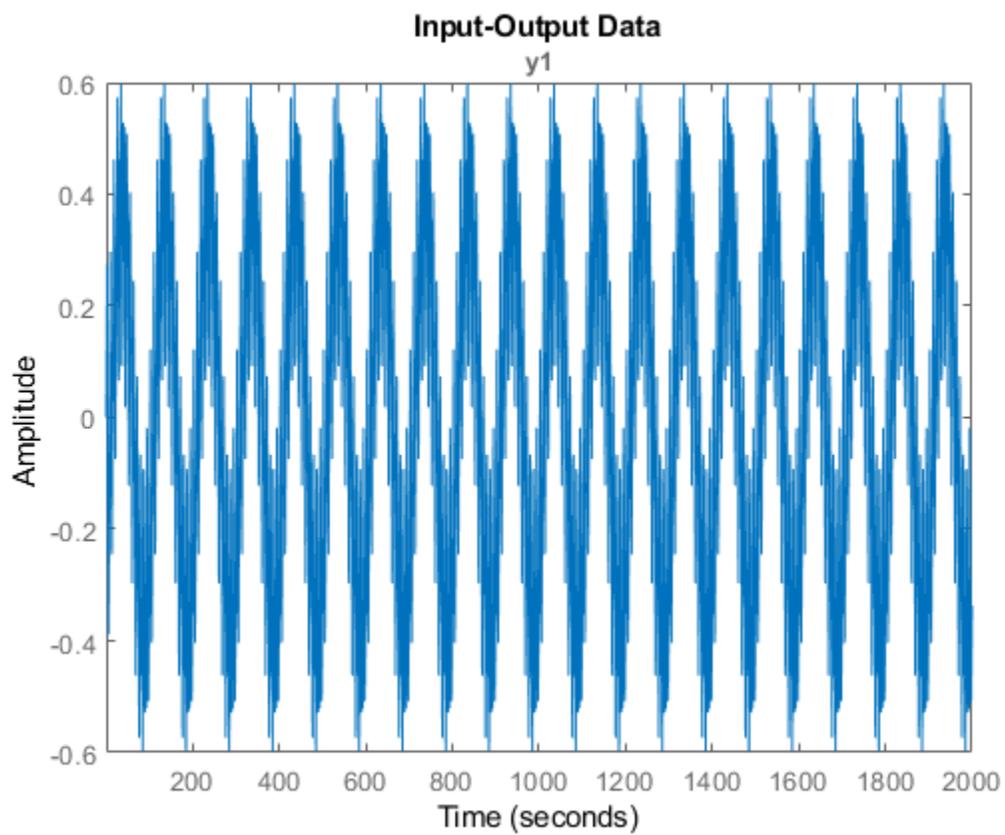
Plot only the input.

```
plot(u)
```



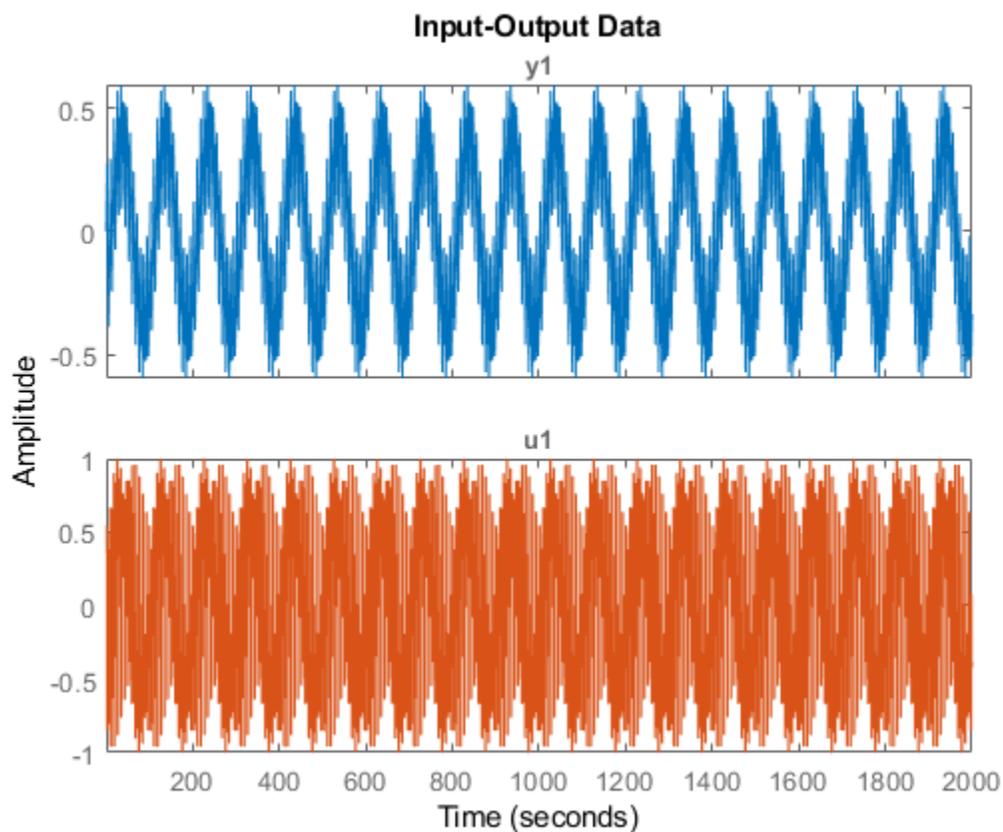
Plot only the output.

```
plot(y)
```



Plot the input and output together.

```
plot(y,u)
```



Alternatively, you can use `plot(iddata(y,u))` .

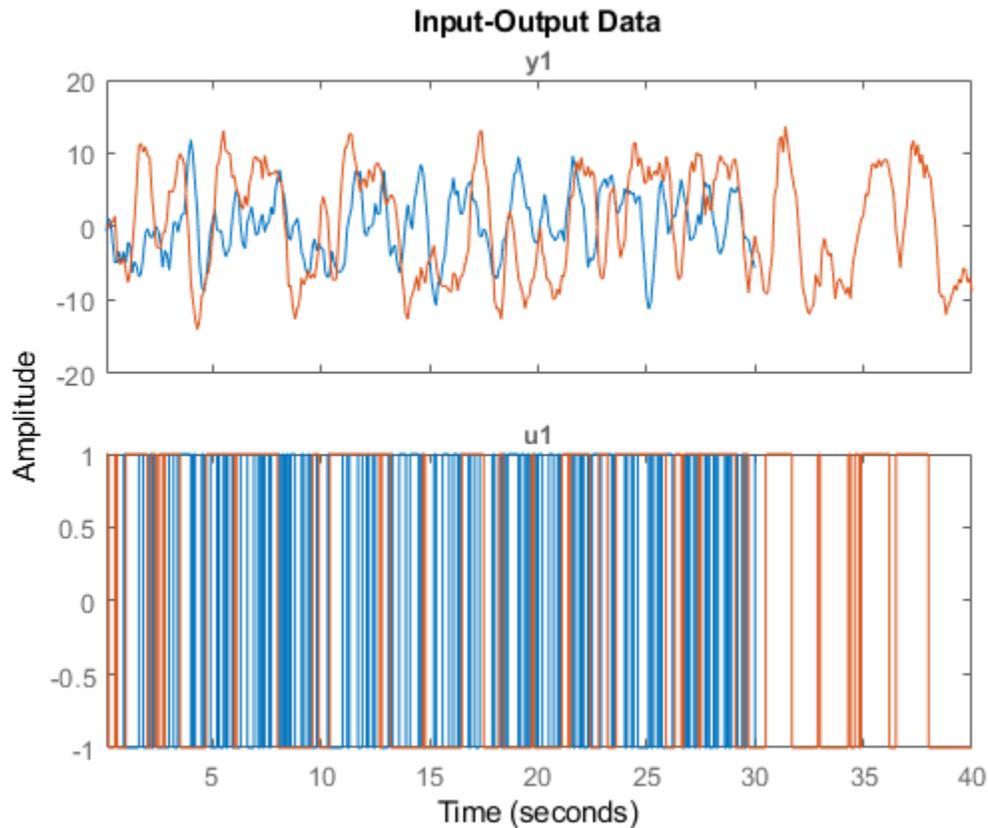
### Plot Multi-Experiment Data

Create a multi-experiment data set.

```
load iddata1 z1
load iddata2 z2
zm = merge(z1,z2);
```

Plot the data.

```
plot(zm);
```



For multi-experiment data, each experiment is treated as a separate data set. You can right-click the plots to view their characteristics.

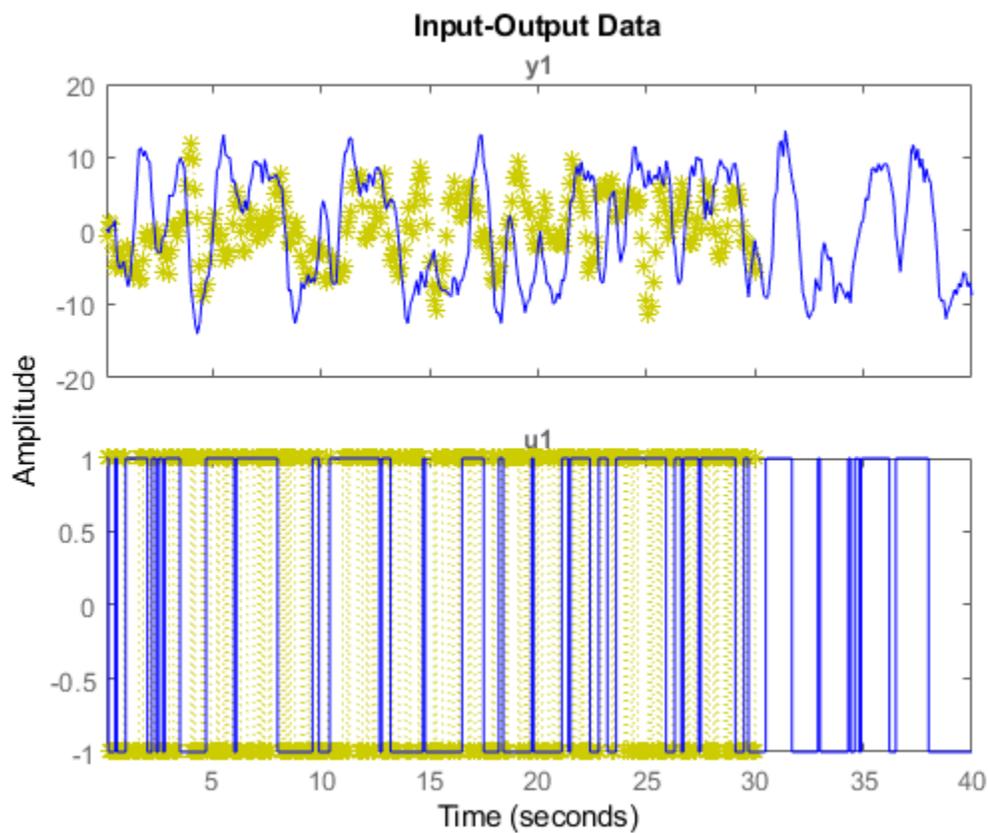
## Specify Line Style, Marker Symbol and Color

Load multiple datasets.

```
load iddata1 z1;  
load iddata2 z2;
```

Specify the linestyle properties.

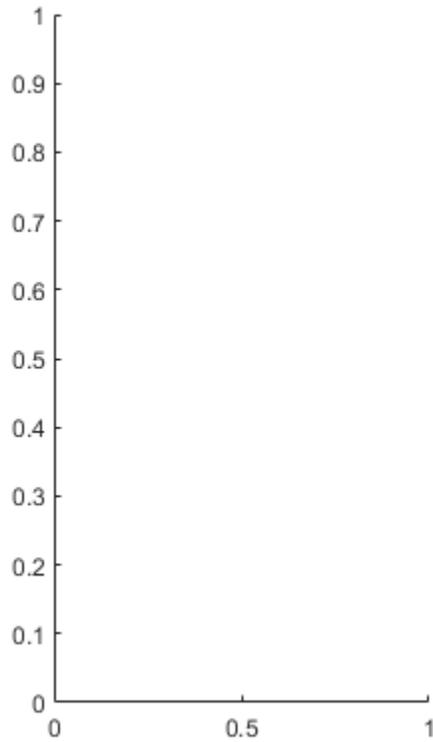
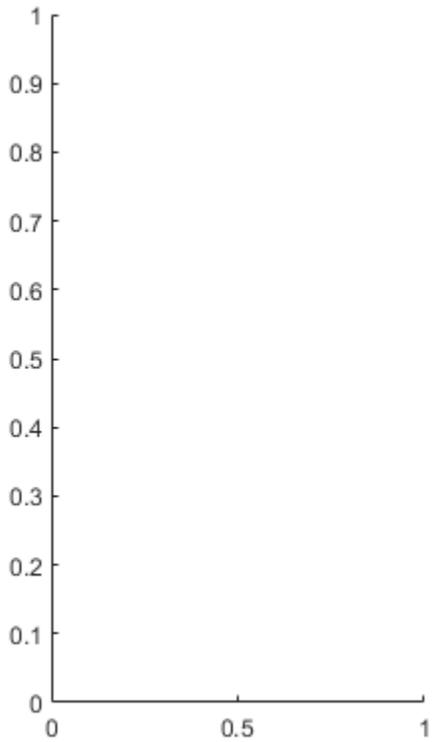
```
plot(z1,'y:*',z2,'b')
```



### Specify Axes Handle

Create a figure with two subplots and return the handle to each subplot axes, `s(1)` and `s(2)`.

```
figure; % new figure
s(1) = subplot(1,2,1); % left subplot
s(2) = subplot(1,2,2); % right subplot
```

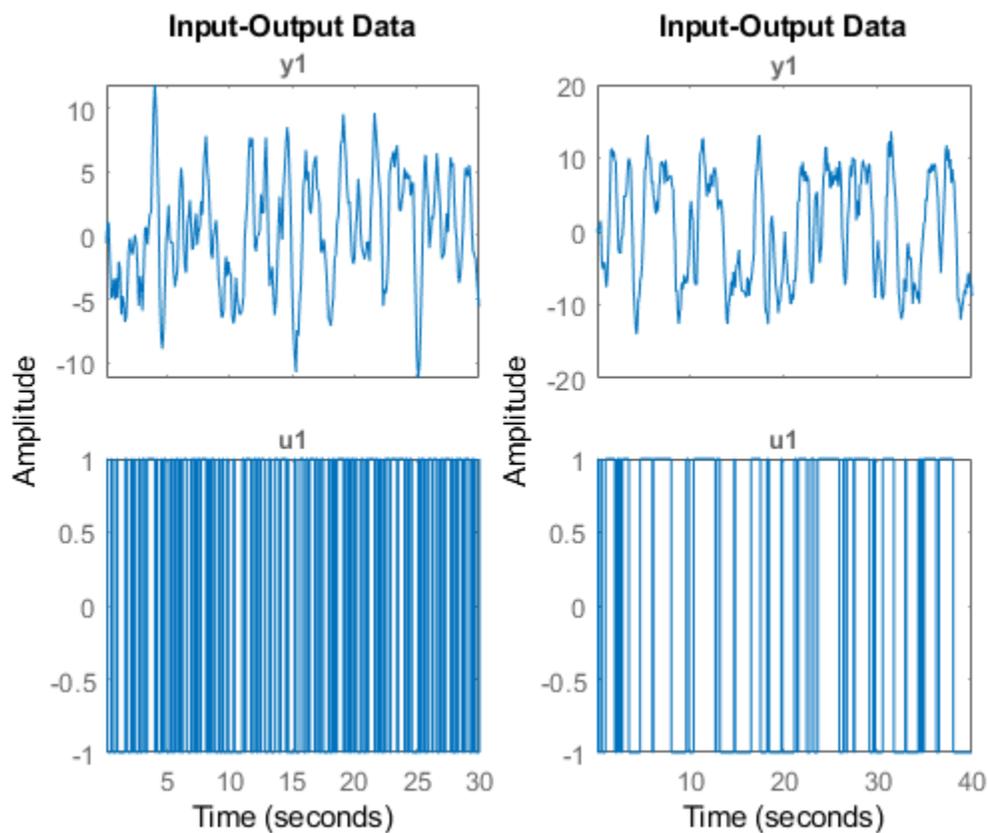


Load the data sets.

```
load iddata1;  
load iddata2;
```

Create a data plot in each axes referring to the axes handles.

```
plot(s(1),z1)  
plot(s(2),z2)
```



### Specify Plot Options

Configure a time plot.

```
opt = iddataPlotOptions('time');
```

View the plot in minutes time units.

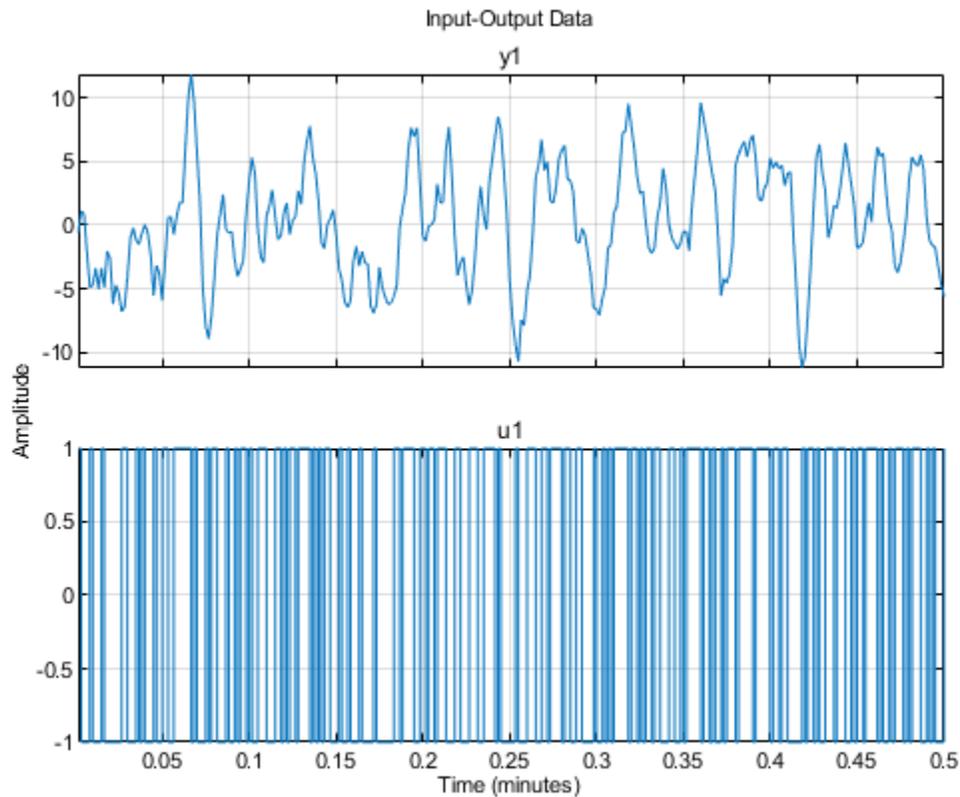
```
opt.TimeUnits = 'minutes';
```

Turn grid on.

```
opt.Grid = 'on';
```

Create plot with the options specified by opt.

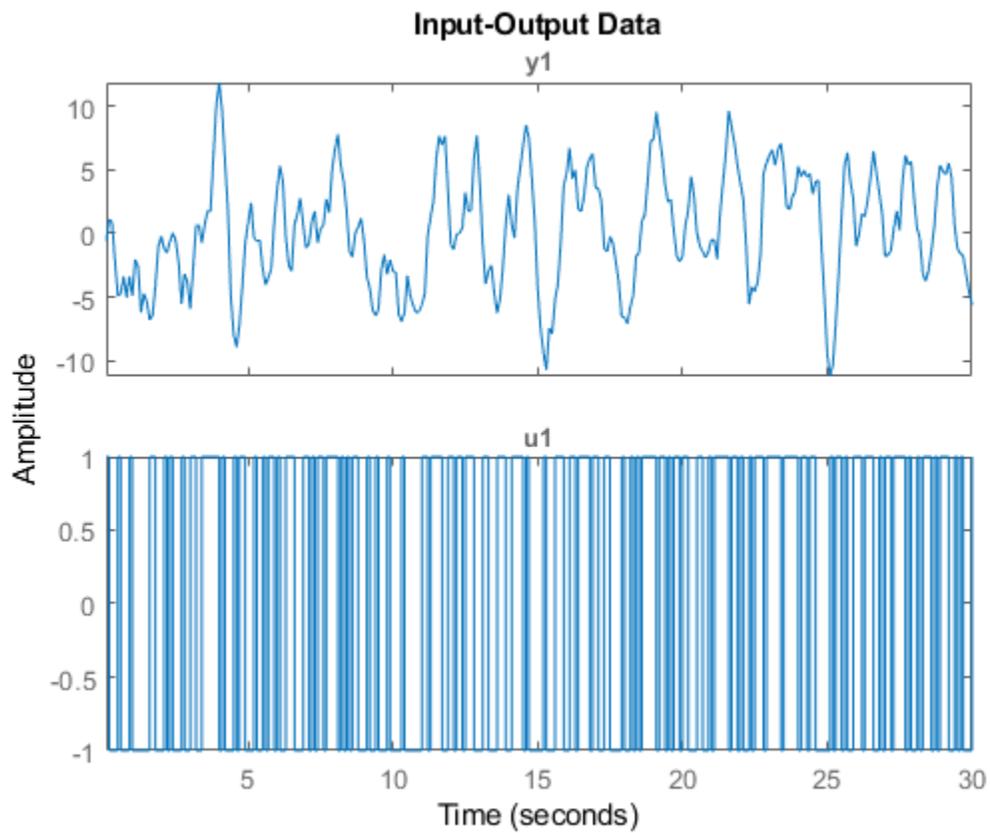
```
load iddata1 z1  
plot(z1, opt);
```



## Change Plot Properties Using Handles

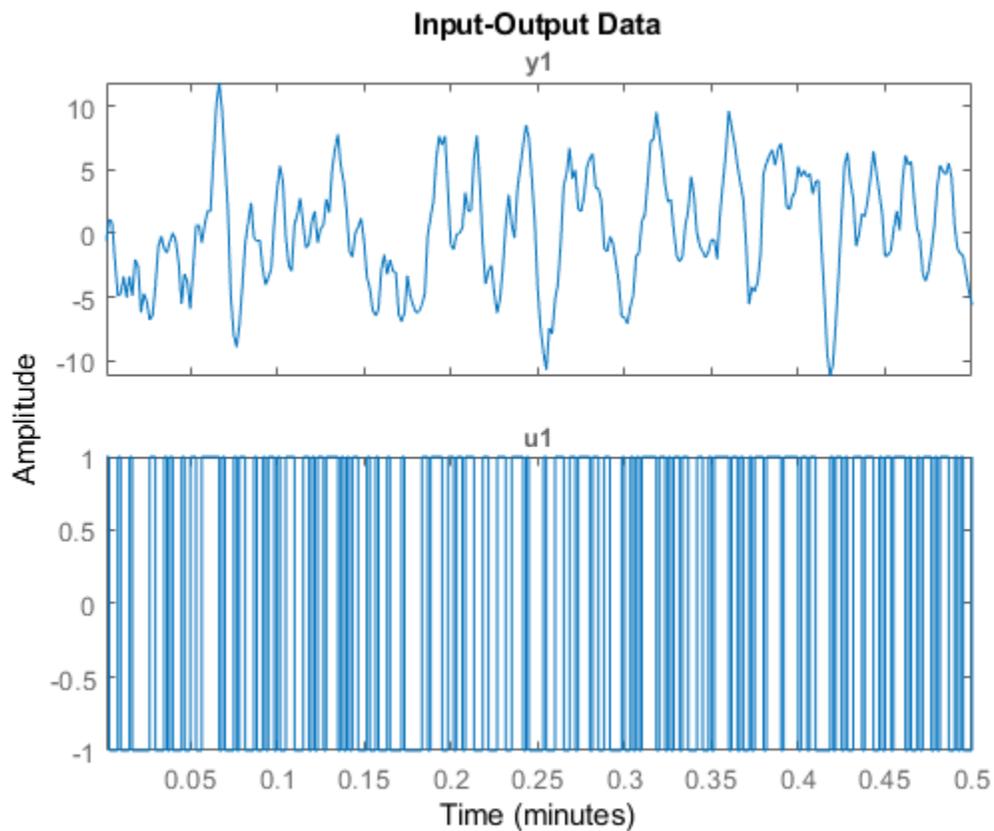
Create a data plot and return the handle.

```
load iddata1;  
h = plot(z1);
```



Set the time unit on the plot.

```
setoptions(h, 'TimeUnits', 'minutes');
```



## Change Orientation of Input-Output Data Axes

Generate data with two inputs and one output.

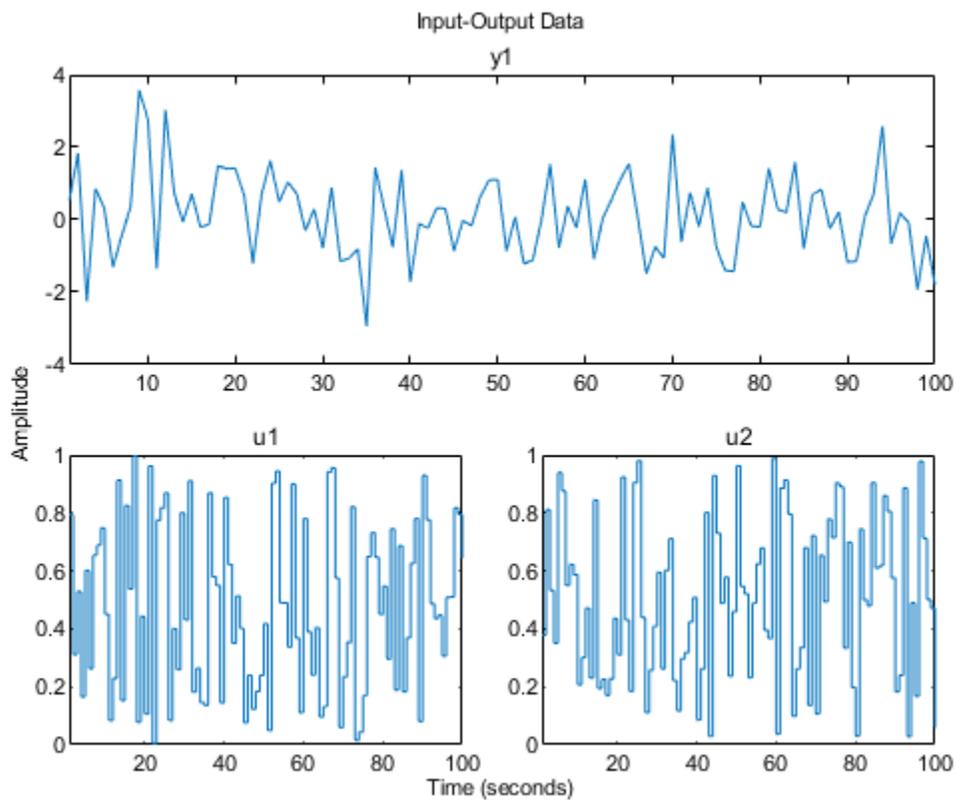
```
z = iddata(randn(100,1),rand(100,2));
```

Configure a time plot.

```
opt = iddataPlotOptions('time');
```

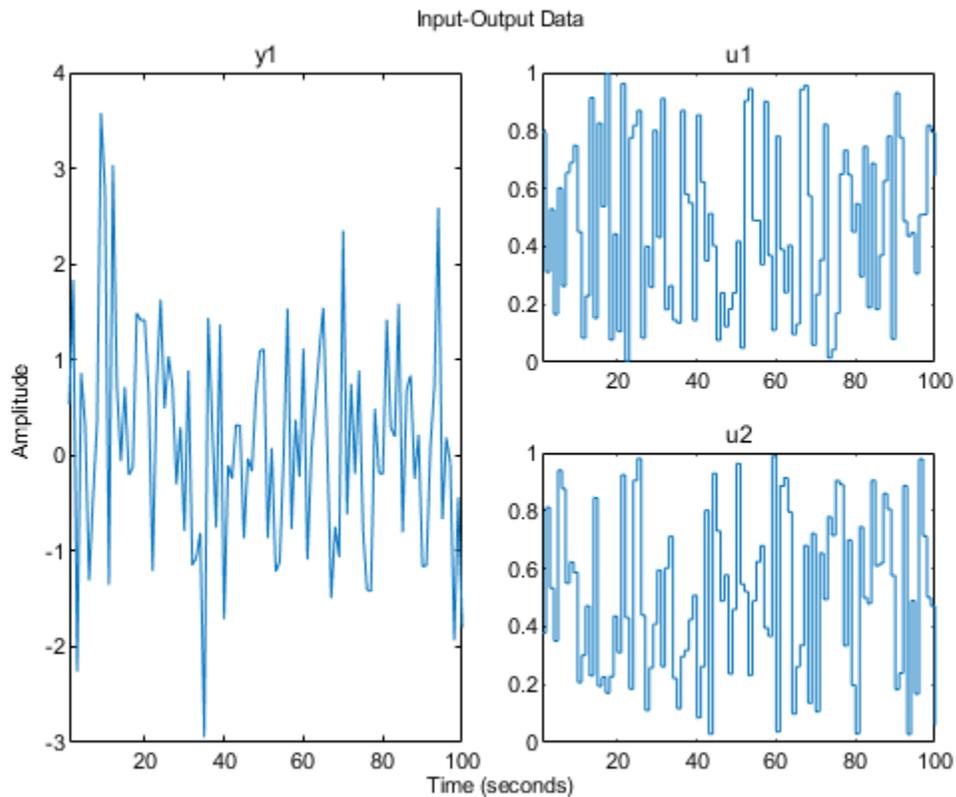
Plot the data.

```
h = plot(z,opt);
```



Change the orientation of the plots such that all inputs are plotted in one column, and all outputs are in a second column.

```
opt.Orientation = 'two-column';  
h = plot(z,opt);
```



Alternatively, use `setoptions`.

```
setoptions(h, 'Orientation', 'two-column')
```

You can also change the orientation by right-clicking the plot and choosing **Orientation** in the context menu.

## Tips

- Right-clicking the plot opens the context menu where you can access the following options and plot controls:

- **Datasets** — View the datasets used in the plot.
- **Characteristics** — View data characteristics.
  - **Peak Value** — Peak value of the data. Useful for transient data.
  - **Mean Value** — Mean value of the data. Useful for steady-state data.
- **Orientation** — Choose orientation of the input and output plots.
  - **Output row and input row** — (For datasets containing more than one input or output channel) Plot all outputs in one row and all inputs in a second row.
  - **Output column and input column** — (For datasets containing more than one input or output channel) Plot all outputs in one column and all inputs in a second column.
  - **Single row** — Plot all inputs and outputs in one row.
  - **Single column** — Plot all inputs and outputs in one column.
- **I/O Grouping** — (For datasets containing more than one input or output channel) Group input and output channels on the plot.
- **I/O Selector** — (For datasets containing more than one input or output channel) Select a subset of the input and output channels to plot. By default, all input and output channels are plotted.
- **Grid** — Add grids to your plot.
- **Normalize** — Normalize the y-scale of all data in the plot.
- **Properties** — Open the Property Editor dialog box, where you can customize plot attributes.

## See Also

iddata | iddataPlotOptions | identpref

**Introduced in R2014a**

## idnlarx/plot

Plot nonlinearity of nonlinear ARX model

### Syntax

```
plot(model)
plot(model,color)
plot(model1,...,modelN)
plot(model1,color1,...,modelN,colorN)
plot( ____, 'NumberofSamples',N)
```

### Description

`plot(model)` plots the nonlinearity of a nonlinear ARX model on a nonlinear ARX plot on page 1-1156. The plot shows the nonlinearity for all outputs of the model as a function of its input regressors.

`plot(model,color)` specifies the color to use.

`plot(model1,...,modelN)` generates the plot for multiple models.

`plot(model1,color1,...,modelN,colorN)` specifies the color for each model. You do not need to specify the color for all models.

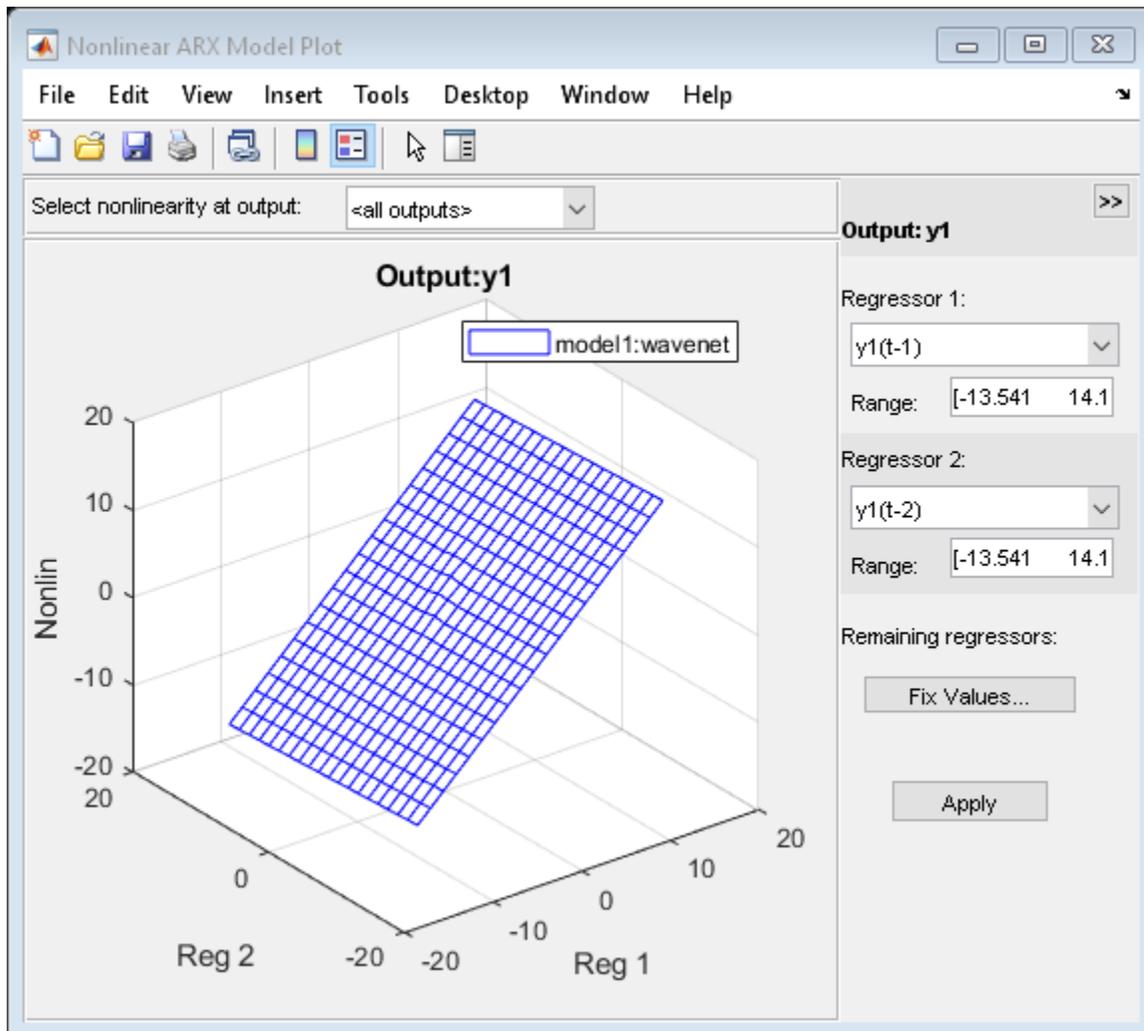
`plot( ____, 'NumberofSamples',N)` specifies the number of samples to use to grid the regressor space on each axis. This syntax can include any of the input argument combinations in the previous syntaxes.

### Examples

#### Plot Nonlinearity of a Nonlinear ARX Model

Estimate a nonlinear ARX model and plot its nonlinearity.

```
load iddata1
model1 = nlarx(z1,[4 2 1], 'wave', 'nlr', [1:3]);
plot(model1)
```

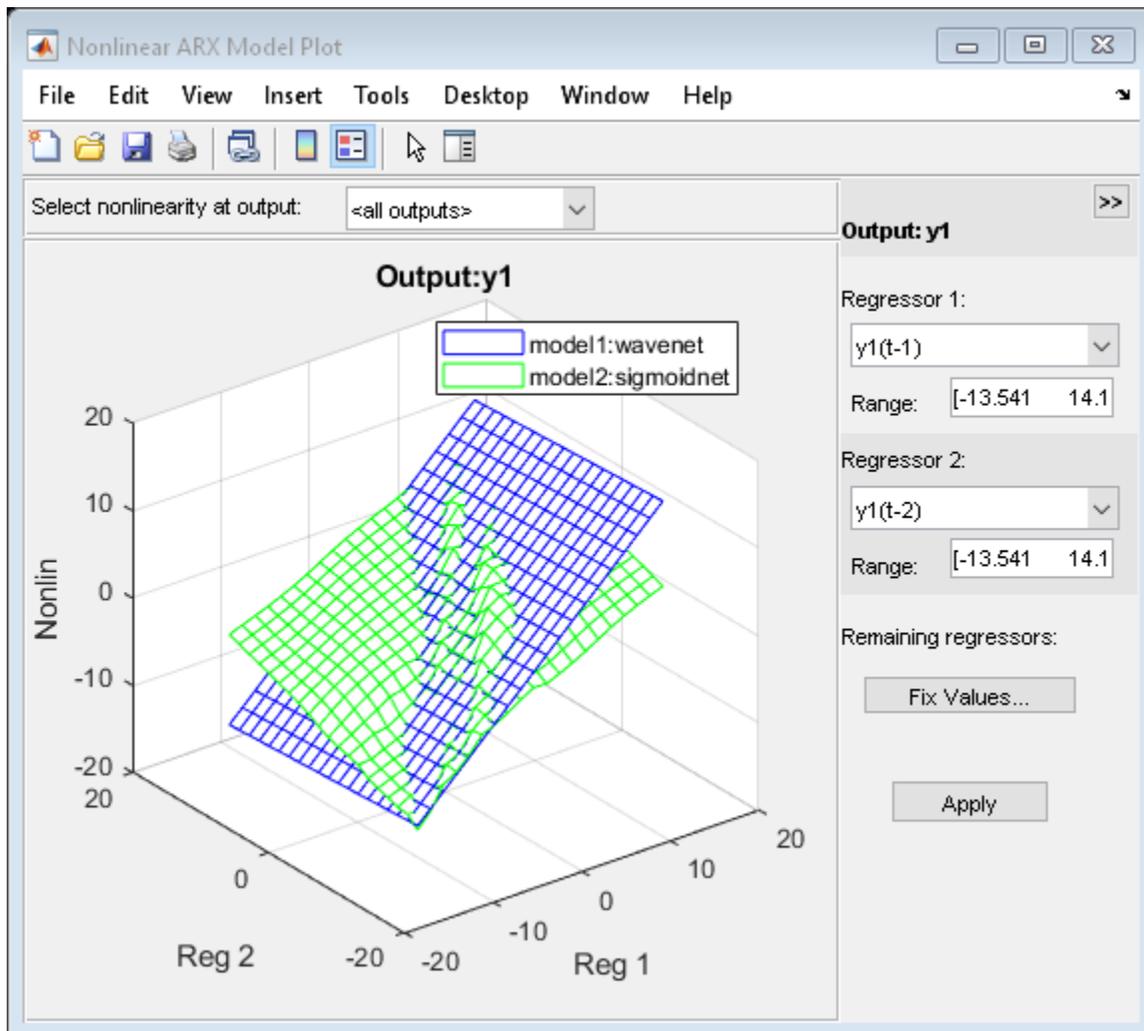


In the plot window, you can choose:

- The regressors to use on the plot axes, and specify the center points for the other regressors in the configuration panel. For multi-output models, each output is plotted separately.
- The output to view from the drop-down list located at the top of the plot.

### Specify Line Style for Multiple Models

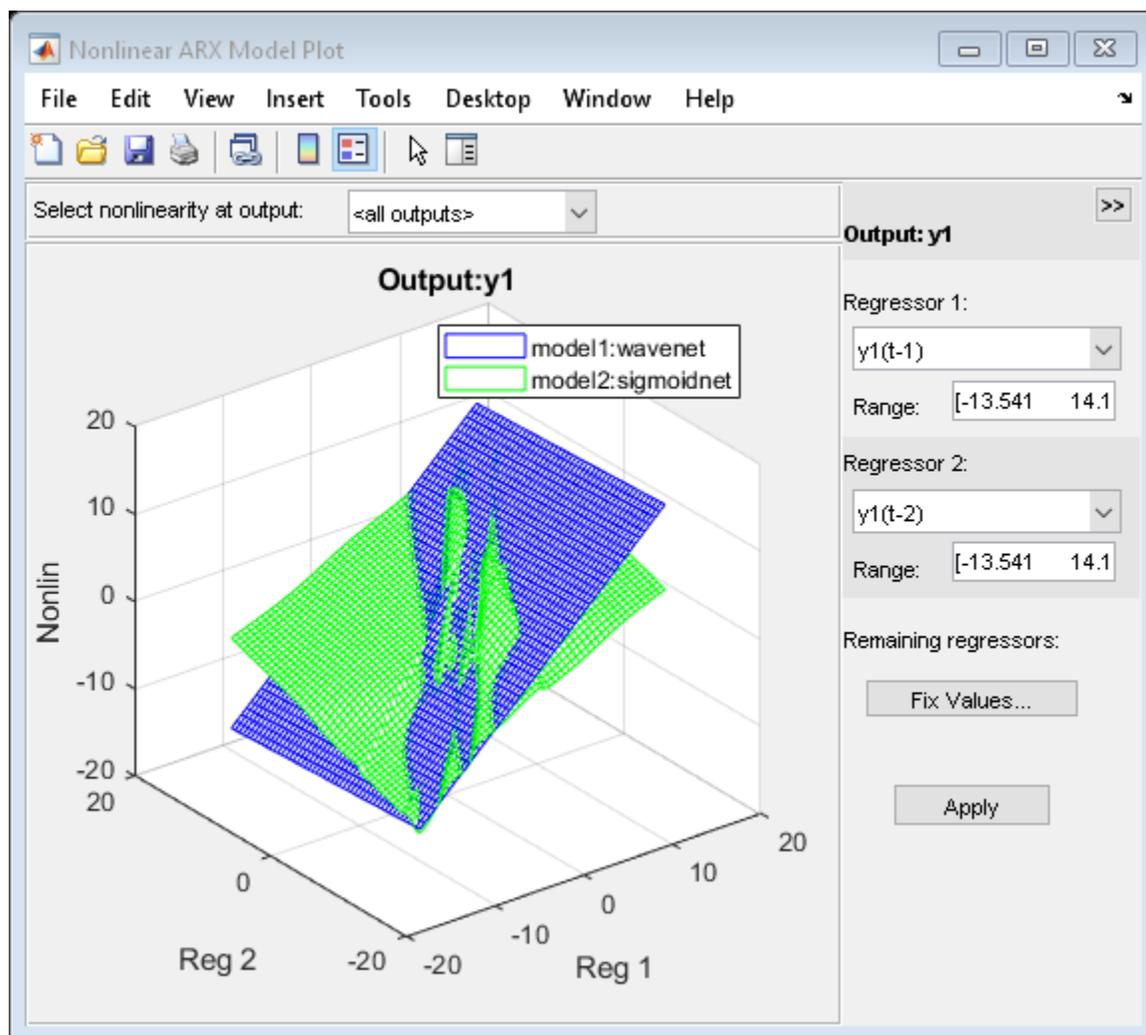
```
load iddata1
model1 = nlarx(z1,[4 2 1], 'wave', 'nlr', [1:3]);
model2 = nlarx(z1,[4 2 1], 'sigmoidnet', 'nlr', [1:3]);
plot(model1, 'b', model2, 'g')
```



### Specify Number of Samples

```
load iddata1
model1 = nlarx(z1,[4 2 1], 'wave', 'nlr', [1:3]);
```

```
model2 = nlarx(z1,[4 2 1], 'sigmoidnet', 'nlr',[1:3]);  
plot(model1,'b', model2, 'g', 'NumberofSamples',50)
```



## Input Arguments

### **model** — Estimated nonlinear ARX model

`idnlarx model`

Estimated nonlinear ARX model, specified as an `idnlarx` model object. Use `nlarx` to estimate the model.

### **color** — Color to use

character vector of color name | vector of doubles

Color to use to plot the regressors, specified as one of the following:

- Character vector of color name, specified as one of the following:
  - 'b'
  - 'y'
  - 'm'
  - 'c'
  - 'r'
  - 'g'
  - 'w'
- 3-element double vector of RGB values

By default, the colors are automatically chosen.

Data Types: `double` | `char`

### **N** — Number of points

20 (default) | positive integer

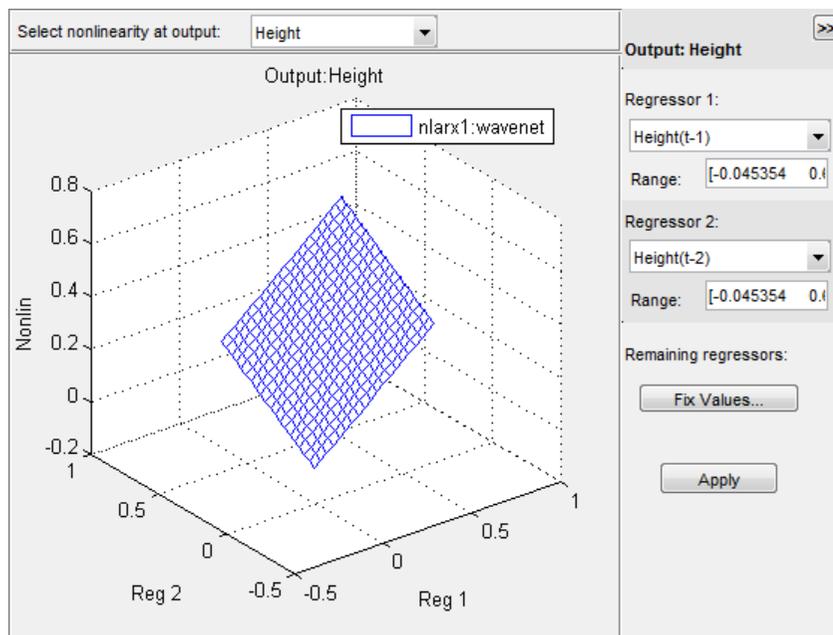
Number of points used on the regressor axis to display the regressor samples, specified as a positive integer.

Data Types: `double`

## Definitions

### What is a Nonlinear ARX Plot?

A nonlinear ARX plot displays the evaluated model nonlinearity for a chosen model output as a function of one or two model regressors. For a model  $M$ , the model nonlinearity ( $M.Nonlinearity$ ) is a nonlinearity estimator function, such as `wavenet`, `sigmoidnet`, or `treepartition`, that uses model regressors as its inputs.



To understand what is plotted, suppose that  $\{r_1, r_2, \dots, r_N\}$  are the  $N$  regressors used by a nonlinear ARX model  $M$  with nonlinearity  $n_l$  corresponding to a model output. You can use `getreg(M)` to view these regressors. The expression `Nonlin = evaluate(nl, [v1, v2, ..., vN])` returns the model output for given values of these regressors, that is,  $r_1 = v_1, r_2 = v_2, \dots, r_N = v_N$ . For plotting the nonlinearities, you select one or two of the  $N$  regressors, for example,  $r_{sub} = \{r_1, r_4\}$ . The software varies the values of these regressors in a specified range, while fixing the value of the remaining regressors, and generates the plot of `Nonlin` vs.  $r_{sub}$ . By default, the software sets the values of the remaining fixed regressors to their estimated means, but you can change these values.

The regressor means are stored in the `Nonlinearity.Parameters.RegressorMean` property of the model.

Examining a nonlinear ARX plot can help you gain insight into which regressors have the strongest effect on the model output. Understanding the relative importance of the regressors on the output can help you decide which regressors to include in the nonlinear function for that output. If the shape of the plot looks like a plane for all the chosen regressor values, then the model is probably linear in those regressors. In this case, you can remove the corresponding regressors from nonlinear block, and repeat the estimation.

Furthermore, you can create several nonlinear models for the same data using different nonlinearity estimators, such a `wavenet` network and `treepartition`, and then compare the nonlinear surfaces of these models. Agreement between plots for various models increases the confidence that these nonlinear models capture the true dynamics of the system.

To learn more about configuring the plot, see “Configuring a Nonlinear ARX Plot”.

## See Also

`evaluate` | `getreg` | `idnlarx` | `nlarx`

## Topics

“Structure of Nonlinear ARX Models”

“Validate Nonlinear ARX Models”

**Introduced in R2014a**

## idnlhw/plot

Plot input and output nonlinearity, and linear responses of Hammerstein-Wiener model

### Syntax

```
plot(model)
plot(model,LineStyle)
plot(model1,...,modelN)
plot(model1,LineStyle1...,modelN,LineStyleN)

plot( ____,Name,Value)
```

### Description

`plot(model)` plots the input and output nonlinearity, and linear responses of a Hammerstein-Wiener model on a Hammerstein-Wiener plot on page 1-1169. The plot shows the responses of the input and output nonlinearity, and linear blocks that represent the model.

`plot(model,LineStyle)` specifies the line style.

`plot(model1,...,modelN)` generates the plot for multiple models.

`plot(model1,LineStyle1...,modelN,LineStyleN)` specifies the line style for each model. You do not need to specify the line style for all models.

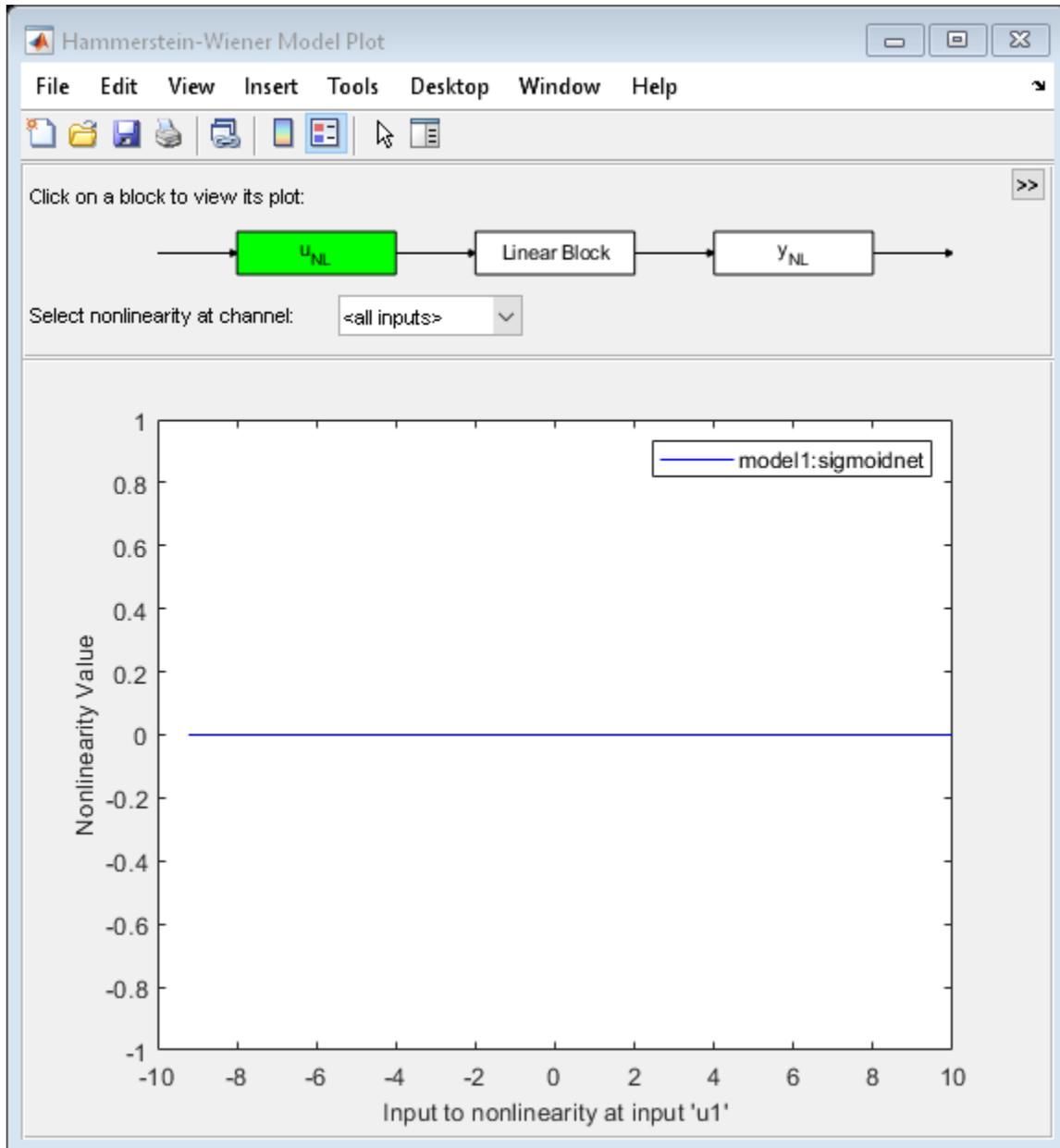
`plot( ____,Name,Value)` specifies plot properties using additional options specified by one or more `Name,Value` pair arguments. This syntax can include any of the input argument combinations in the previous syntaxes.

### Examples

## Plot Input and Output Nonlinearity and Linear Response of a Hammerstein-Wiener Model

Estimate a Hammerstein-Wiener Model and plot responses of its input and output nonlinearity and linear blocks.

```
load iddata3
model1 = nlhw(z3,[4 2 1], 'sigmoidnet', 'deadzone');
plot(model1)
```

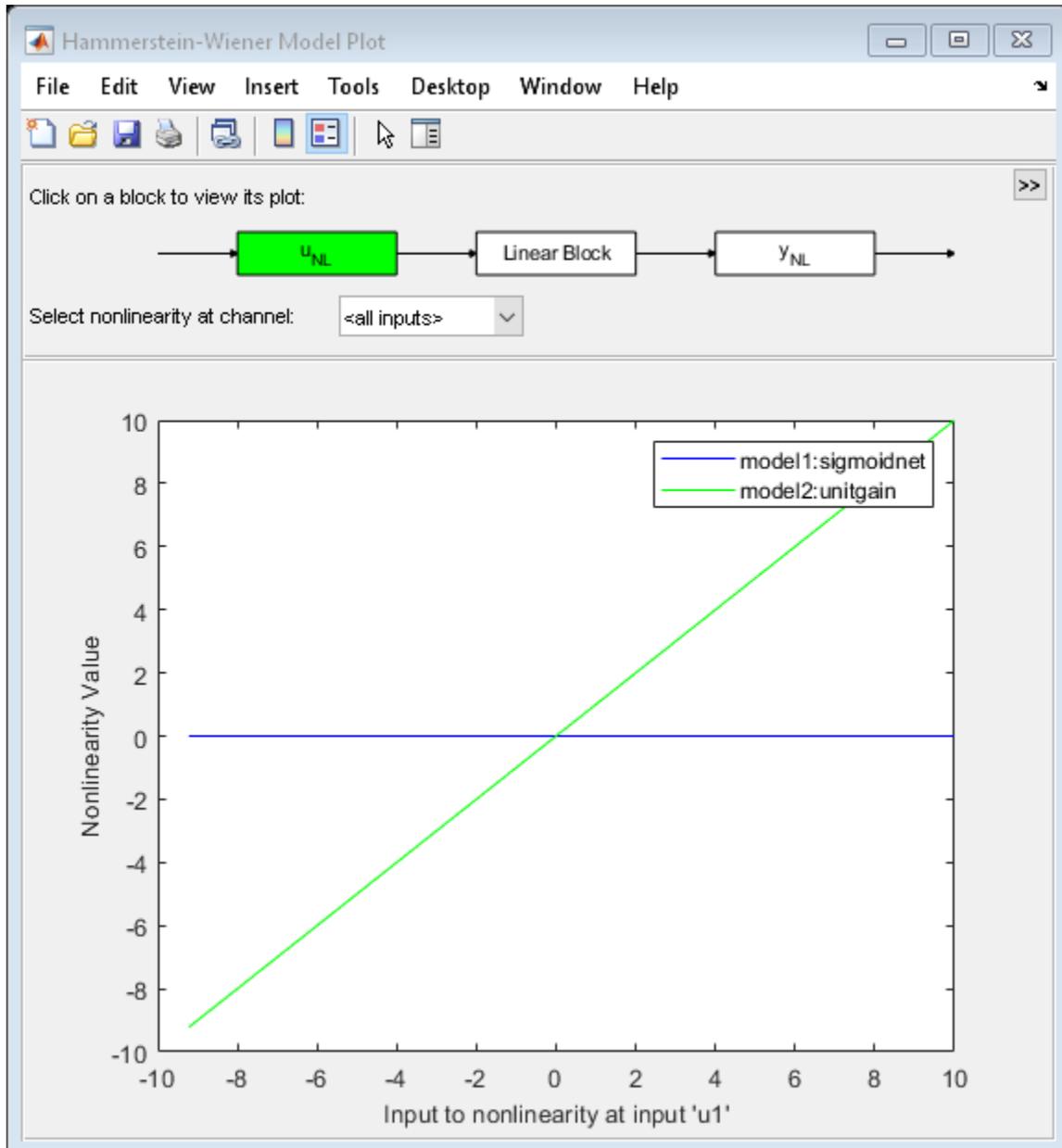


Explore the various plots in the plot window by clicking one of the three blocks that represent the model:

- uNL - Input nonlinearity, representing the static nonlinearity at the input ( `model.InputNonlinearity` ) to the LinearBlock.
- Linear Block - Step, impulse, Bode and pole-zero plots of the embedded linear model ( `model.LinearModel` ). By default, a step plot is displayed.
- yNL - Output nonlinearity, representing the static nonlinearity at the output ( `model.OutputNonlinearity` ) of the Linear Block.

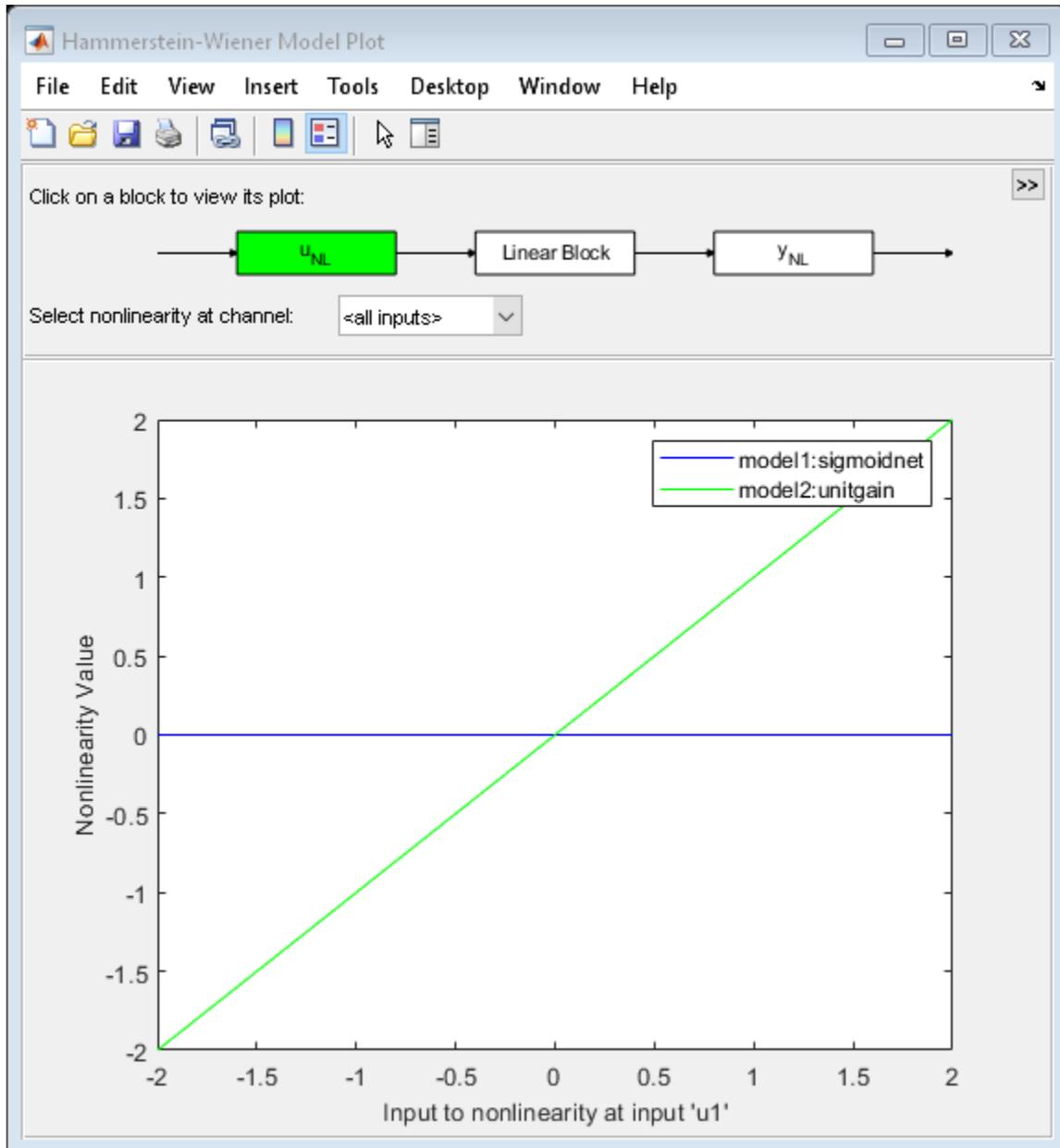
### Specify Line Style for Multiple Hammerstein-Weiner Models

```
load iddata3
model1 = nlhw(z3,[4 2 1], 'sigmoidnet', 'deadzone');
model2 = nlhw(z3, [4 2 1],[], 'sigmoidnet');
plot(model1, 'b-', model2, 'g')
```



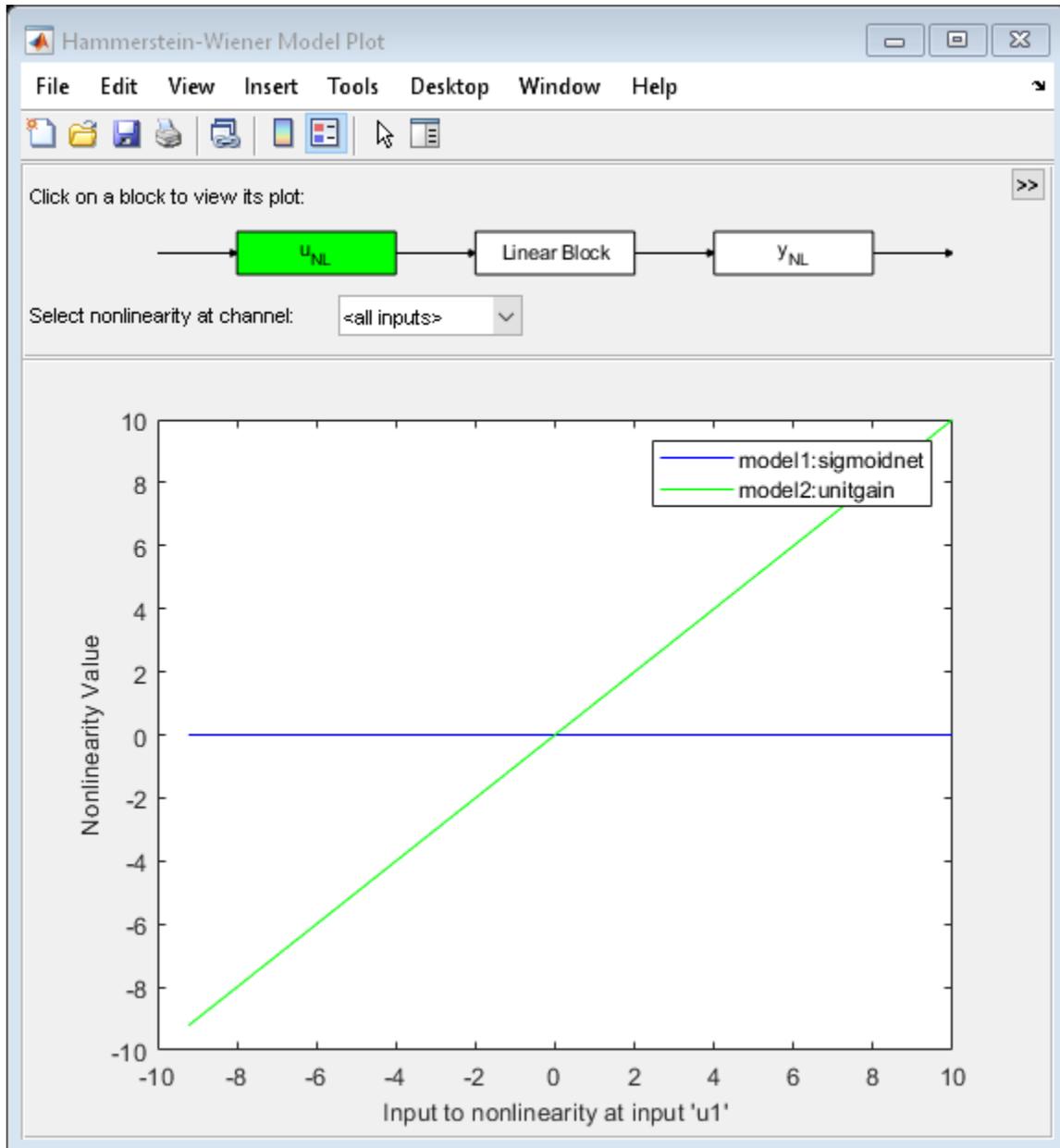
**Specify Number of Samples, Time Samples, and Range of Input Nonlinearity**

```
load iddata3
model1 = nlhw(z3,[4 2 1], 'sigmoidnet', 'deadzone');
model2 = nlhw(z3, [4 2 1],[], 'sigmoidnet');
plot(model1, 'b-', model2, 'g', 'NumberOfSamples', 50, 'time', 10, 'InputRange', [-2 2]);
```



### Specify Time Samples, Frequency, and Range of Output Nonlinearity

```
load iddata3
model1 = nlhw(z3,[4 2 1], 'sigmoidnet', 'deadzone');
model2 = nlhw(z3, [4 2 1], [], 'sigmoidnet');
plot(model1,model2, 'time', 1:500, 'freq', {0.01,100}, 'OutputRange', [0 1000]);
```



## Input Arguments

### **model** — Estimated Hammerstein-Wiener model

`idnlhw` model

Estimated Hammerstein-Wiener model, specified as an `idnlhw` model object. Use `n_lhw` to estimate the model.

### **LineStyle** — Line style, marker symbol, and color

character vector

Line style, marker symbol, and color, specified as a character vector. `LineStyle` takes values such as `'b'`, `'b+:'`. For more information, see the `plot` reference page in the MATLAB documentation.

Data Types: `char`

## Name-Value Pair Arguments

Specify optional comma-separated pairs of `Name`, `Value` arguments. `Name` is the argument name and `Value` is the corresponding value. `Name` must appear inside quotes. You can specify several name and value pair arguments in any order as `Name1, Value1, ..., NameN, ValueN`.

Example: `plot(model, 'NumberOfSamples', 10)` specifies to use 10 data points for the input regressors.

### **NumberOfSamples** — Number of data points to use for input regressors

100 (default) | positive integer

Number of data points to use for the input regressors when evaluating the nonlinearities at individual input or output channels, specified as a positive integer. This property does not affect the plots of the linear block.

Data Types: `double`

### **InputRange** — Minimum and maximum regressor values for evaluating input nonlinearities

range of regressor values used during each model's estimation. (default) | positive integer | vector

Minimum and maximum regressor values to use when evaluating the nonlinearities at each input channel, specified as positive integers or [min max] vector, where minimum value is less than the maximum value.

You can use 'uRange' as a shortcut name for this property.

Data Types: double

### **OutputRange — Minimum and maximum regressor values for evaluating output nonlinearities**

range of regressor values used during each model's estimation (default) | positive integer | vector

Minimum and maximum regressor values to use when evaluating the nonlinearities at each output channel, specified as positive integers or [min max] vector, where minimum value is less than the maximum value.

You can use 'yRange' as a shortcut name for this property.

Data Types: double

### **Time — Time samples to compute transient responses of the linear block**

each model's dynamics determine the time samples used (default) | positive scalar | vector

The time samples at which the transient responses (step and impulse) of the linear block of the `idnlhw` model must be computed, specified as one of the following values:

- Positive scalar — Denotes end time for transient responses of all models. For example, 10.
- Vector of time instants — A double vector of equi-sampled values denotes the time samples at which the transient response must be computed. For example, [0:0.1:10].

This property takes the same values as the `step` command on the model.

### **Frequency — Frequencies at which to compute the Bode response**

automatically chosen inside the Nyquist frequency range (default) | [min max] range of positive scalars | vector of positive integers

Frequencies at which to compute the Bode response, specified as one of the following values:

- `[Wmin Wmax]` range — Frequency interval between `Wmin` and `Wmax` (in units `rad/(model.TimeUnit)`) covered using logarithmically placed points.
- Vector of non-negative frequency values — Allows computation of bode response at those frequencies.

By default, the response is computed at some automatically chosen frequencies inside the Nyquist frequency range. Frequencies above Nyquist frequency ( $\pi/\text{model.Ts}$ ) are ignored.

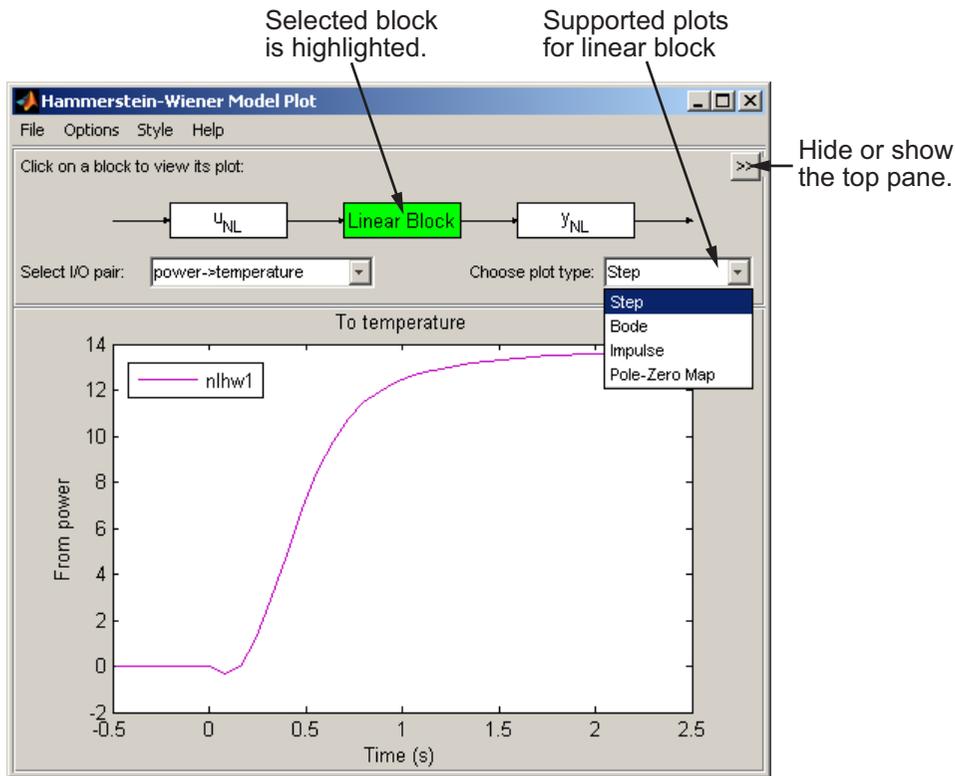
This property takes the same values as the `bode` command on the model.

## Definitions

### What is a Hammerstein-Wiener Plot?

A Hammerstein-Wiener plot displays the static input and output nonlinearities and linear responses of a Hammerstein-Wiener model.

Examining a Hammerstein-Wiener plot can help you determine whether you have selected a complicated nonlinearity for modeling your system. For example, suppose you use a piecewise-linear input nonlinearity to estimate your model, but the plot indicates saturation behavior. You can estimate a new model using the simpler saturation nonlinearity instead. For multivariable systems, you can use the Hammerstein-Wiener plot to determine whether to exclude nonlinearities for specific channels. If the nonlinearity for a specific input or output channel does not exhibit strong nonlinear behavior, you can estimate a new model after setting the nonlinearity at that channel to unit gain.



You can generate these plots in the **System Identification** app and at the command line. In the plot window, you can view the nonlinearities and linear responses by clicking one of the three blocks that represent the model:

- $u_{NL}$  (*input nonlinearity*)— Click this block to view the static nonlinearity at the input to the **Linear Block**. The plot displays `evaluate(M.InputNonlinearity,u)` where **M** is the Hammerstein-Wiener model, and **u** is the input to the input nonlinearity block. For information about the blocks, see “Structure of Hammerstein-Wiener Models”.
- **Linear Block** — Click this block to view the Step, impulse, Bode, and pole-zero response plots of the embedded linear model (**M.LinearModel**). By default, a step plot of the linear model is displayed.
- $y_{NL}$  (*output nonlinearity*) — Click this block to view the static nonlinearity at the output of the **Linear Block**. The plot displays `evaluate(M.OutputNonlinearity,x)`, where **x** is the output of the linear block.

To learn more about how to configure the linear and nonlinear blocks plots, see “Configuring a Hammerstein-Wiener Plot”.

## See Also

`idnlhw` | `nlhw`

## Topics

“Structure of Hammerstein-Wiener Models”

“Validating Hammerstein-Wiener Models”

**Introduced in R2014a**

## pole

Poles of dynamic system

### Syntax

```
P = pole(sys)
P = pole(sys,J1,...,JN)
```

### Description

`P = pole(sys)` returns the poles of the SISO or MIMO dynamic system model `sys`. The output is expressed as the reciprocal of the time units specified in `sys.TimeUnit`. The poles of a dynamic system determine the stability and response of the system.

An open-loop linear time-invariant system is stable if:

- In continuous-time, all the poles of the transfer function have negative real parts. When the poles are visualized on the complex s-plane, then they must all lie in the left-half plane (LHP) to ensure stability.
- In discrete-time, all the poles must have a magnitude strictly smaller than one, that is they must all lie inside the unit circle.

`P = pole(sys,J1,...,JN)` returns the poles `P` of the entries in model array `sys` with subscripts `(J1,...,JN)`.

### Examples

#### Poles of Discrete-Time Transfer Function

Compute the poles of the following discrete-time transfer function:

$$\text{sys}(z) = \frac{0.0478z - 0.0464}{z^2 - 1.81z + 0.9048}$$

```
sys = tf([0.04798 0.0464],[1 -1.81 0.9048],0.1);
P = pole(sys)
```

*P = 2×1 complex*

```
0.9050 + 0.2929i
0.9050 - 0.2929i
```

For stable discrete systems, all their poles must have a magnitude strictly smaller than one, that is they must all lie inside the unit circle. The poles in this example are a pair of complex conjugates, and lie inside the unit circle. Hence, the system `sys` is stable.

### Poles of Transfer Function

Calculate the poles of following transfer function:

$$sys(s) = \frac{4.2s^2 + 0.25s - 0.004}{s^2 + 9.6s + 17}$$

```
sys = tf([4.2,0.25,-0.004],[1,9.6,17]);
P = pole(sys)
```

*P = 2×1*

```
-7.2576
-2.3424
```

For stable continuous systems, all their poles must have negative real parts. `sys` is stable since the poles are negative, that is, they lie in the left half of the complex plane.

### Poles of Models in an Array

For this example, load `invertedPendulumArray.mat`, which contains a 3-by-3 array of inverted pendulum models. The mass of the pendulum varies as you move from model to model along a single column of `sys`, and the length of the pendulum varies as you move

along a single row. The mass values used are 100g, 200g and 300g, and the pendulum lengths used are 3m, 2m and 1m respectively.

	<i>Column 1</i>	<i>Column 2</i>	<i>Column 3</i>
<i>Row 1</i>	100g, 3m	100g, 2m	100g, 1m
<i>Row 2</i>	200g, 3m	200g, 2m	200g, 1m
<i>Row 3</i>	300g, 3m	300g, 2m	300g, 1m

```
load('invertedPendulumArray.mat', 'sys');  
size(sys)
```

3x3 array of transfer functions.  
Each model has 1 outputs and 1 inputs.

Find poles of the model array.

```
P = pole(sys);  
P(:, :, 2, 1)
```

```
ans = 3x1  
  
    2.1071  
   -2.1642  
   -0.1426
```

`P(:, :, 2, 1)` corresponds to the poles of the model with 200g pendulum weight and 3m length.

## Input Arguments

### **sys** — Dynamic system

dynamic system model | model array

Dynamic system, specified as a SISO or MIMO dynamic system model, or an array of SISO or MIMO dynamic system models. Dynamic systems that you can use include continuous-time or discrete-time numeric LTI models such as `tf`, `zpk`, or `ss` models.

If `sys` is a generalized state-space model `genss` or an uncertain state-space model `uss`, `pole` returns the poles of the current or nominal value of `sys`. If `sys` is an array of

models, `pole` returns the poles of the model corresponding to its subscript `J1, . . . , JN` in `sys`. For more information on model arrays, see “Model Arrays” (Control System Toolbox).

### **J1, . . . , JN — Indices of models in array whose poles you want to extract**

positive integer

Indices of models in array whose poles you want to extract, specified as a positive integer. You can provide as many indices as there are array dimensions in `sys`. For example, if `sys` is a 4-by-5 array of dynamic system models, the following command extracts the poles for entry (2,3) in the array.

```
P = pole(sys,2,3);
```

## **Output Arguments**

### **P — Poles of the dynamic system**

column vector | array

Poles of the dynamic system, returned as a scalar or an array. If `sys` is:

- A single model, then `P` is a column vector of poles of the dynamic system model `sys`
- A model array, then `P` is an array of poles of each model in `sys`

`P` is expressed as the reciprocal of the time units specified in `sys.TimeUnit`. For example, `pole` is expressed in 1/minute if `sys.TimeUnit = 'minutes'`.

Depending on the type of system model, poles are computed in the following way:

- For state-space models, the poles are the eigenvalues of the  $A$  matrix, or the generalized eigenvalues of  $A - \lambda E$  in the descriptor case.
- For SISO transfer functions or zero-pole-gain models, the poles are the denominator roots. For more information, see `roots`.
- For MIMO transfer functions (or zero-pole-gain models), the poles are returned as the union of the poles for each SISO entry. If some I/O pairs have a common denominator, the roots of such I/O pair denominator are counted only once.

## Limitations

- Multiple poles are numerically sensitive and cannot be computed with high accuracy. A pole  $\lambda$  with multiplicity  $m$  typically results in a cluster of computed poles distributed on a circle with center  $\lambda$  and radius of order

$$\rho \approx \varepsilon^{1/m},$$

where  $\varepsilon$  is the relative machine precision (`eps`).

For more information on multiple poles, see “Sensitivity of Multiple Roots” (Control System Toolbox).

- If `sys` has internal delays, poles are obtained by first setting all internal delays to zero so that the system has a finite number of poles, thereby creating a zero-order Padé approximation. For some systems, setting delays to zero creates singular algebraic loops, which result in either improper or ill-defined, zero-delay approximations. For these systems, `pole` returns an error.

To assess the stability of models with internal delays, use `step` or `impulse`.

## See Also

`damp` | `impulse` | `pzmap` | `pzplot` | `step` | `zero`

## Topics

“Pole and Zero Locations” (Control System Toolbox)

“Sensitivity of Multiple Roots” (Control System Toolbox)

**Introduced in R2012a**

# polydata

Access polynomial coefficients and uncertainties of identified model

## Syntax

```
[A,B,C,D,F] = polydata(sys)
[A,B,C,D,F,dA,dB,dC,dD,dF] = polydata(sys)
[___] = polydata(sys,J1,...,JN)
[___] = polydata(___, 'cell')
```

## Description

`[A,B,C,D,F] = polydata(sys)` returns the coefficients of the polynomials A, B, C, D, and F that describe the identified model `sys`. The polynomials describe the `idpoly` representation of `sys` as follows.

- For discrete-time `sys`:

$$A(q^{-1})y(t) = \frac{B(q^{-1})}{F(q^{-1})}u(t-nk) + \frac{C(q^{-1})}{D(q^{-1})}e(t).$$

$u(t)$  are the inputs to `sys`.  $y(t)$  are the outputs.  $e(t)$  is a white noise disturbance.

- For continuous-time `sys`:

$$A(s)Y(s) = \frac{B(s)}{F(s)}U(s)e^{-\tau s} + \frac{C(s)}{D(s)}E(s).$$

$U(s)$  are the Laplace transformed inputs to `sys`.  $Y(s)$  are the Laplace transformed outputs.  $E(s)$  is the Laplace transform of a white noise disturbance.

If `sys` is an identified model that is not an `idpoly` model, `polydata` converts `sys` to `idpoly` form to extract the polynomial coefficients.

`[A,B,C,D,F,dA,dB,dC,dD,dF] = polydata(sys)` also returns the uncertainties `dA`, `dB`, `dC`, `dD`, and `dF` of each of the corresponding polynomial coefficients of `sys`.

`[___] = polydata(sys,J1,...,JN)` returns the polynomial coefficients for the `J1, ..., JN` entry in the array `sys` of identified models.

`[___] = polydata(___, 'cell')` returns all polynomials as cell arrays of double vectors, regardless of the input and output dimensions of `sys`.

## Input Arguments

### **sys**

Identified model or array of identified models. `sys` can be continuous-time or discrete-time. `sys` can be SISO or MIMO.

### **J1, ..., JN**

Indices selecting a particular model from an N-dimensional array `sys` of identified models.

## Output Arguments

### **A,B,C,D,F**

Polynomial coefficients of the `idpoly` representation of `sys`.

- If `sys` is a SISO model, each of `A`, `B`, `C`, `D`, and `F` is a row vector. The length of each row vector is the order of the corresponding polynomial.
  - For discrete-time `sys`, the coefficients are ordered in ascending powers of  $q^{-1}$ . For example, `B = [1 -4 9]` means that  $B(q^{-1}) = 1 - 4q^{-1} + 9q^{-2}$ .
  - For continuous-time `sys`, the coefficients are ordered in descending powers of  $s$ . For example, `B = [1 -4 9]` means that  $B(s) = s^2 - 4s + 9$ .
- If `sys` is a MIMO model, each of `A`, `B`, `C`, `D`, and `F` is a cell array. The dimensions of the cell arrays are determined by the input and output dimensions of `sys` as follows:
  - `A` —  $N_y$ -by- $N_y$  cell array

- B, F —  $N_y$ -by- $N_u$  cell array
- C, D —  $N_y$ -by-1 cell array

$N_y$  is the number of outputs of `sys`, and  $N_u$  is the number of inputs.

Each entry in a cell array is a row vector that contains the coefficients of the corresponding polynomial. The polynomial coefficients are ordered the same way as the SISO case.

### **dA, dB, dC, dD, dF**

Uncertainties in the estimated polynomial coefficients of `sys`.

dA, dB, dC, dD, and dF are row vectors or cell arrays whose dimensions exactly match the corresponding A, B, C, D, and F outputs.

Each entry in dA, dB, dC, dD, and dF gives the standard deviation of the corresponding estimated coefficient. For example, `dA{1,1}(2)` gives the standard deviation of the estimated coefficient returned at `A{1,1}(2)`.

## **Examples**

### **Extract Polynomial Coefficients and Uncertainties from Identified Model**

Load system data and estimate a 2-input, 2-output model.

```
load iddata1 z1
load iddata2 z2
data = [z1 z2(1:300)];

nk = [1 1; 1 0];
na = [2 2; 1 3];
nb = [2 3; 1 4];
nc = [2;3];
nd = [1;2];
nf = [2 2;2 1];

sys = polyest(data,[na nb nc nd nf nk]);
```

The data loaded into `z1` and `z2` is discrete-time `iddata` with a sample time of 0.1 s. Therefore, `sys` is a two-input, two-output discrete-time `idpoly` model of the form:

$$A(q^{-1})y(t) = \frac{B(q^{-1})}{F(q^{-1})}u(t - nk) + \frac{C(q^{-1})}{D(q^{-1})}e(t)$$

The inputs to `polyest` set the order of each polynomial in `sys`.

Access the estimated polynomial coefficients of `sys` and the uncertainties in those coefficients.

```
[A,B,C,D,F,dA,dB,dC,dD,dF] = polydata(sys);
```

The outputs `A`, `B`, `C`, `D`, and `F` are cell arrays of coefficient vectors. The dimensions of the cell arrays are determined by the input and output dimensions of `sys`. For example, `A` is a 2-by-2 cell array because `sys` has two inputs and two outputs. Each entry in `A` is a row vector containing identified polynomial coefficients. For example, examine the second diagonal entry in `A`.

```
A{2,2}
```

```
ans = 1×4
```

```
1.0000    -0.8825    -0.2030     0.4364
```

For discrete-time `sys`, the coefficients are arranged in order of increasing powers of  $q^{-1}$ . Therefore, `A{2,2}` corresponds to the polynomial

$$1 - 0.8682q^{-1} - 0.2244q^{-2} + 0.4467q^{-3}.$$

The dimensions of `dA` match those of `A`. Each entry in `dA` gives the standard deviation of the corresponding estimated polynomial coefficient in `A`. For example, examine the uncertainties of the second diagonal entry in `A`.

```
dA{2,2}
```

```
ans = 1×4
```

```
0    0.2849    0.4269    0.2056
```

The lead coefficient of `A{2,2}` is fixed at 1, and therefore has no uncertainty. The remaining entries in `dA{2,2}` are the uncertainties in the  $q^{-1}$ ,  $q^{-2}$ , and  $q^{-3}$  coefficients, respectively.

## **See Also**

`iddata` | `idpoly` | `idssdata` | `polyest` | `tfdata` | `zpkdata`

**Introduced before R2006a**

## polyest

Estimate polynomial model using time- or frequency-domain data

### Syntax

```
sys = polyest(data,[na nb nc nd nf nk])  
sys = polyest(data,[na nb nc nd nf nk],Name,Value)  
sys = polyest(data,init_sys)  
sys = polyest( ___, opt)
```

### Description

`sys = polyest(data,[na nb nc nd nf nk])` estimates a polynomial model, `sys`, using the time- or frequency-domain data, `data`.

`sys` is of the form

$$A(q)y(t) = \frac{B(q)}{F(q)}u(t - nk) + \frac{C(q)}{D(q)}e(t)$$

$A(q)$ ,  $B(q)$ ,  $F(q)$ ,  $C(q)$  and  $D(q)$  are polynomial matrices.  $u(t)$  is the input, and  $nk$  is the input delay.  $y(t)$  is the output and  $e(t)$  is the disturbance signal. `na`, `nb`, `nc`, `nd` and `nf` are the orders of the  $A(q)$ ,  $B(q)$ ,  $C(q)$ ,  $D(q)$  and  $F(q)$  polynomials, respectively.

`sys = polyest(data,[na nb nc nd nf nk],Name,Value)` estimates a polynomial model with additional attributes of the estimated model structure specified by one or more `Name, Value` pair arguments.

`sys = polyest(data,init_sys)` estimates a polynomial model using the linear system `init_sys` to configure the initial parameterization.

`sys = polyest( ___, opt)` estimates a polynomial model using the option set, `opt`, to specify estimation behavior.

## Input Arguments

### **data**

Estimation data.

For time-domain estimation, `data` is an `iddata` object containing the input and output signal values.

You can estimate only discrete-time models using time-domain data. For estimating continuous-time models using time-domain data, see `tfest`.

For frequency-domain estimation, `data` can be one of the following:

- Recorded frequency response data (`frd` or `idfrd`)
- `iddata` object with its properties specified as follows:
  - `InputData` — Fourier transform of the input signal
  - `OutputData` — Fourier transform of the output signal
  - `Domain` — 'Frequency'

It may be more convenient to use `oe` or `tfest` to estimate a model for frequency-domain data.

### **na**

Order of the polynomial  $A(q)$ .

`na` is an  $N_y$ -by- $N_y$  matrix of nonnegative integers.  $N_y$  is the number of outputs, and  $N_u$  is the number of inputs.

`na` must be zero if you are estimating a model using frequency-domain data.

### **nb**

Order of the polynomial  $B(q) + 1$ .

`nb` is an  $N_y$ -by- $N_u$  matrix of nonnegative integers.  $N_y$  is the number of outputs, and  $N_u$  is the number of inputs.

## **nc**

Order of the polynomial  $C(q)$ .

$nc$  is a column vector of nonnegative integers of length  $N_y$ .  $N_y$  is the number of outputs.

$nc$  must be zero if you are estimating a model using frequency-domain data.

## **nd**

Order of the polynomial  $D(q)$ .

$nd$  is a column vector of nonnegative integers of length  $N_y$ .  $N_y$  is the number of outputs.

$nd$  must be zero if you are estimating a model using frequency-domain data.

## **nf**

Order of the polynomial  $F(q)$ .

$nf$  is an  $N_y$ -by- $N_u$  matrix of nonnegative integers.  $N_y$  is the number of outputs, and  $N_u$  is the number of inputs.

## **nk**

Input delay in number of samples, expressed as fixed leading zeros of the  $B$  polynomial.

$nk$  is an  $N_y$ -by- $N_u$  matrix of nonnegative integers.

$nk$  must be zero when estimating a continuous-time model.

## **opt**

Estimation options.

$opt$  is an options set, created using `polyestOptions`, that specifies estimation options including:

- Estimation objective
- Handling of initial conditions
- Numerical search method to be used in estimation

## **init\_sys**

Linear system that configures the initial parameterization of `sys`.

You obtain `init_sys` by either performing an estimation using measured data or by direct construction.

If `init_sys` is an `idpoly` model, `polyest` uses the parameters and constraints defined in `init_sys` as the initial guess for estimating `sys`.

Use the `Structure` property of `init_sys` to configure initial guesses and constraints for  $A(q)$ ,  $B(q)$ ,  $F(q)$ ,  $C(q)$ , and  $D(q)$ . For example:

- To specify an initial guess for the  $A(q)$  term of `init_sys`, set `init_sys.Structure.A.Value` as the initial guess.
- To specify constraints for the  $B(q)$  term of `init_sys`:
  - Set `init_sys.Structure.B.Minimum` to the minimum  $B(q)$  coefficient values.
  - Set `init_sys.Structure.B.Maximum` to the maximum  $B(q)$  coefficient values.
  - Set `init_sys.Structure.B.Free` to indicate which  $B(q)$  coefficients are free for estimation.

If `init_sys` is not an `idpoly` model, the software first converts `init_sys` to a polynomial model. `polyest` uses the parameters of the resulting model as the initial guess for estimation.

If `opt` is not specified, and `init_sys` is created by estimation, then the estimation options from `init_sys.Report.OptionsUsed` are used.

## **Name-Value Pair Arguments**

Specify optional comma-separated pairs of `Name`, `Value` arguments. `Name` is the argument name and `Value` is the corresponding value. `Name` must appear inside quotes. You can specify several name and value pair arguments in any order as `Name1, Value1, ..., NameN, ValueN`.

### **IODElay**

Transport delays. `IODElay` is a numeric array specifying a separate transport delay for each input/output pair.

For continuous-time systems, specify transport delays in the time unit stored in the `TimeUnit` property. For discrete-time systems, specify transport delays in integer multiples of the sample time, `Ts`.

For a MIMO system with `Ny` outputs and `Nu` inputs, set `IODelay` to a `Ny`-by-`Nu` array. Each entry of this array is a numerical value that represents the transport delay for the corresponding input/output pair. You can also set `IODelay` to a scalar value to apply the same delay to all input/output pairs.

**Default:** 0 for all input/output pairs

### **InputDelay**

Input delay for each input channel, specified as a scalar value or numeric vector. For continuous-time systems, specify input delays in the time unit stored in the `TimeUnit` property. For discrete-time systems, specify input delays in integer multiples of the sample time `Ts`. For example, `InputDelay = 3` means a delay of three sample times.

For a system with `Nu` inputs, set `InputDelay` to an `Nu`-by-1 vector. Each entry of this vector is a numerical value that represents the input delay for the corresponding input channel.

You can also set `InputDelay` to a scalar value to apply the same delay to all channels.

**Default:** 0

### **IntegrateNoise**

Logical vector specifying integrators in the noise channel.

`IntegrateNoise` is a logical vector of length `Ny`, where `Ny` is the number of outputs.

Setting `IntegrateNoise` to `true` for a particular output results in the model:

$$A(q)y(t) = \frac{B(q)}{F(q)}u(t - nk) + \frac{C(q)}{D(q)}\frac{e(t)}{1 - q^{-1}}$$

Where,  $\frac{1}{1 - q^{-1}}$  is the integrator in the noise channel,  $e(t)$ .

Use `IntegrateNoise` to create an ARIMAX model.

For example,

```
load iddata1 z1;
z1 = iddata(cumsum(z1.y),cumsum(z1.u),z1.Ts,'InterSample','foh');
sys = polyest(z1, [2 2 2 0 0 1],'IntegrateNoise',true);
```

## Output Arguments

### sys

Polynomial model, returned as an `idpoly` model. This model is created using the specified model orders, delays, and estimation options.

If `data.Ts` is zero, `sys` is a continuous-time model representing:

$$Y(s) = \frac{B(s)}{F(s)}U(s) + E(s)$$

$Y(s)$ ,  $U(s)$  and  $E(s)$  are the Laplace transforms of the time-domain signals  $y(t)$ ,  $u(t)$  and  $e(t)$ , respectively.

Information about the estimation results and options used is stored in the `Report` property of the model. `Report` have the following fields:

Report Field	Description
Status	Summary of the model status, which indicates whether the model was created by construction or obtained by estimation.
Method	Estimation command used.

Report Field	Description																		
InitialCondition	<p>Handling of initial conditions during model estimation, returned as one of the following values:</p> <ul style="list-style-type: none"> <li>• 'zero' — The initial conditions were set to zero.</li> <li>• 'estimate' — The initial conditions were treated as independent estimation parameters.</li> <li>• 'backcast' — The initial conditions were estimated using the best least squares fit.</li> </ul> <p>This field is especially useful to view how the initial conditions were handled when the InitialCondition option in the estimation option set is 'auto'.</p>																		
Fit	<p>Quantitative assessment of the estimation, returned as a structure. See “Loss Function and Model Quality Metrics” for more information on these quality metrics. The structure has the following fields:</p> <table border="1" data-bbox="387 829 1332 1322"> <thead> <tr> <th data-bbox="387 829 520 873">Field</th> <th data-bbox="520 829 1332 873">Description</th> </tr> </thead> <tbody> <tr> <td data-bbox="387 873 520 982">FitPercent</td> <td data-bbox="520 873 1332 982">Normalized root mean squared error (NRMSE) measure of how well the response of the model fits the estimation data, expressed as a percentage.</td> </tr> <tr> <td data-bbox="387 982 520 1025">LossFcn</td> <td data-bbox="520 982 1332 1025">Value of the loss function when the estimation completes.</td> </tr> <tr> <td data-bbox="387 1025 520 1104">MSE</td> <td data-bbox="520 1025 1332 1104">Mean squared error (MSE) measure of how well the response of the model fits the estimation data.</td> </tr> <tr> <td data-bbox="387 1104 520 1147">FPE</td> <td data-bbox="520 1104 1332 1147">Final prediction error for the model.</td> </tr> <tr> <td data-bbox="387 1147 520 1190">AIC</td> <td data-bbox="520 1147 1332 1190">Raw Akaike Information Criteria (AIC) measure of model quality.</td> </tr> <tr> <td data-bbox="387 1190 520 1234">AICc</td> <td data-bbox="520 1190 1332 1234">Small sample-size corrected AIC.</td> </tr> <tr> <td data-bbox="387 1234 520 1277">nAIC</td> <td data-bbox="520 1234 1332 1277">Normalized AIC.</td> </tr> <tr> <td data-bbox="387 1277 520 1322">BIC</td> <td data-bbox="520 1277 1332 1322">Bayesian Information Criteria (BIC).</td> </tr> </tbody> </table>	Field	Description	FitPercent	Normalized root mean squared error (NRMSE) measure of how well the response of the model fits the estimation data, expressed as a percentage.	LossFcn	Value of the loss function when the estimation completes.	MSE	Mean squared error (MSE) measure of how well the response of the model fits the estimation data.	FPE	Final prediction error for the model.	AIC	Raw Akaike Information Criteria (AIC) measure of model quality.	AICc	Small sample-size corrected AIC.	nAIC	Normalized AIC.	BIC	Bayesian Information Criteria (BIC).
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BIC	Bayesian Information Criteria (BIC).																		
Parameters	Estimated values of model parameters.																		
OptionsUsed	Option set used for estimation. If no custom options were configured, this is a set of default options. See polyestOptions for more information.																		

Report Field	Description																
RandState	State of the random number stream at the start of estimation. Empty, [], if randomization was not used during estimation. For more information, see <code>rng</code> in the MATLAB documentation.																
DataUsed	Attributes of the data used for estimation, returned as a structure with the following fields: <table border="1" data-bbox="385 529 1332 1218"> <thead> <tr> <th>Field</th> <th>Description</th> </tr> </thead> <tbody> <tr> <td>Name</td> <td>Name of the data set.</td> </tr> <tr> <td>Type</td> <td>Data type.</td> </tr> <tr> <td>Length</td> <td>Number of data samples.</td> </tr> <tr> <td>Ts</td> <td>Sample time.</td> </tr> <tr> <td>InterSample</td> <td>Input intersample behavior, returned as one of the following values: <ul style="list-style-type: none"> <li>'zoh' — Zero-order hold maintains a piecewise-constant input signal between samples.</li> <li>'foh' — First-order hold maintains a piecewise-linear input signal between samples.</li> <li>'bl' — Band-limited behavior specifies that the continuous-time input signal has zero power above the Nyquist frequency.</li> </ul> </td> </tr> <tr> <td>InputOffset</td> <td>Offset removed from time-domain input data during estimation. For nonlinear models, it is [].</td> </tr> <tr> <td>OutputOffset</td> <td>Offset removed from time-domain output data during estimation. For nonlinear models, it is [].</td> </tr> </tbody> </table>	Field	Description	Name	Name of the data set.	Type	Data type.	Length	Number of data samples.	Ts	Sample time.	InterSample	Input intersample behavior, returned as one of the following values: <ul style="list-style-type: none"> <li>'zoh' — Zero-order hold maintains a piecewise-constant input signal between samples.</li> <li>'foh' — First-order hold maintains a piecewise-linear input signal between samples.</li> <li>'bl' — Band-limited behavior specifies that the continuous-time input signal has zero power above the Nyquist frequency.</li> </ul>	InputOffset	Offset removed from time-domain input data during estimation. For nonlinear models, it is [].	OutputOffset	Offset removed from time-domain output data during estimation. For nonlinear models, it is [].
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Report Field	Description	
Termination	Termination conditions for the iterative search used for prediction error minimization. Structure with the following fields:	
	Field	Description
	WhyStop	Reason for terminating the numerical search.
	Iterations	Number of search iterations performed by the estimation algorithm.
	FirstOrderOptimality	$\infty$ -norm of the gradient search vector when the search algorithm terminates.
	FcnCount	Number of times the objective function was called.
	UpdateNorm	Norm of the gradient search vector in the last iteration. Omitted when the search method is 'lsqnonlin' or 'fmincon'.
	LastImprovement	Criterion improvement in the last iteration, expressed as a percentage. Omitted when the search method is 'lsqnonlin' or 'fmincon'.
	Algorithm	Algorithm used by 'lsqnonlin' or 'fmincon' search method. Omitted when other search methods are used.
For estimation methods that do not require numerical search optimization, the Termination field is omitted.		

For more information on using Report, see “Estimation Report”.

## Examples

### Estimate Polynomial Model with Redundant Parameterization

Estimate a model with redundant parameterization. That is, a model with all polynomials ( $A$ ,  $B$ ,  $C$ ,  $D$ , and  $F$ ) active.

Load estimation data.

```
load iddata2 z2;
```

Specify the model orders and delays.

```
na = 2;
nb = 2;
nc = 3;
nd = 3;
nf = 2;
nk = 1;
```

Estimate the model.

```
sys = polyest(z2,[na nb nc nd nf nk]);
```

### Estimate Polynomial Model Using Regularization

Estimate a regularized polynomial model by converting a regularized ARX model.

Load estimation data.

```
load regularizationExampleData.mat m0simdata;
```

Estimate an unregularized polynomial model of order 20.

```
m1 = polyest(m0simdata(1:150),[0 20 20 20 20 1]);
```

Estimate a regularized polynomial model of the same order. Determine the Lambda value by trial and error.

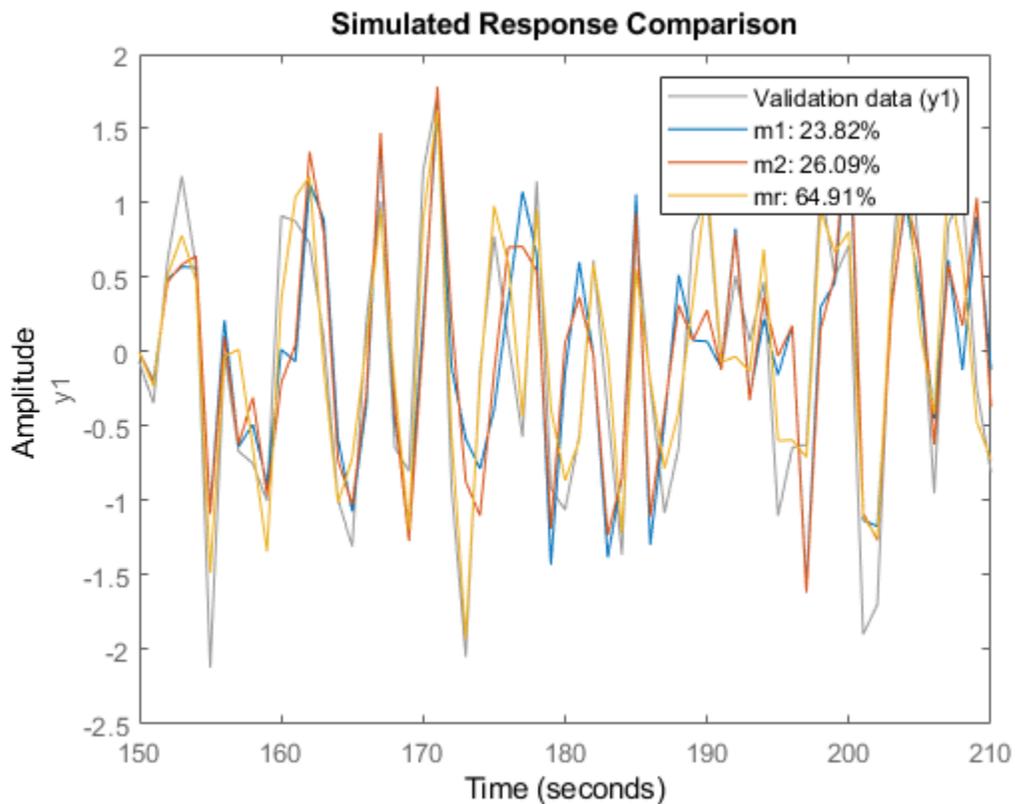
```
opt = polyestOptions;
opt.Regularization.Lambda = 1;
m2 = polyest(m0simdata(1:150),[0 20 20 20 20 1],opt);
```

Obtain a lower-order polynomial model by converting a regularized ARX model and reducing its order. Use `arxregul` to determine the regularization parameters.

```
[L,R] = arxRegul(m0simdata(1:150),[30 30 1]);
opt1 = arxOptions;
opt1.Regularization.Lambda = L;
opt1.Regularization.R = R;
m0 = arx(m0simdata(1:150),[30 30 1],opt1);
mr = idpoly(balred(idss(m0),7));
```

Compare the model outputs against the data.

```
opt2 = compareOptions('InitialCondition','z');  
compare(m0simdata(150:end),m1,m2,mr,opt2);
```



### Estimate ARIMAX model

Load input/output data and create cumulative sum input and output signals for estimation.

```
load iddata1 z1  
data = iddata(cumsum(z1.y),cumsum(z1.u),z1.Ts,'InterSample','foh');
```

Specify the model polynomial orders. Set the orders of the inactive polynomials,  $D$  and  $F$ , to 0.

```
na = 2;
nb = 2;
nc = 2;
nd = 0;
nf = 0;
nk = 1;
```

Identify an ARIMAX model by setting the 'IntegrateNoise' option to true.

```
sys = polyest(data,[na nb nc nd nf nk],'IntegrateNoise',true);
```

### Estimate Multi-Output ARMAX Model

Estimate a multi-output ARMAX model for a multi-input, multi-output data set.

Load estimation data.

```
load iddata1 z1
load iddata2 z2
data = [z1 z2(1:300)];
```

`data` is a data set with 2 inputs and 2 outputs. The first input affects only the first output. Similarly, the second input affects only the second output.

Specify the model orders and delays. The  $F$  and  $D$  polynomials are inactive.

```
na = [2 2; 2 2];
nb = [2 2; 3 4];
nk = [1 1; 0 0];
nc = [2;2];
nd = [0;0];
nf = [0 0; 0 0];
```

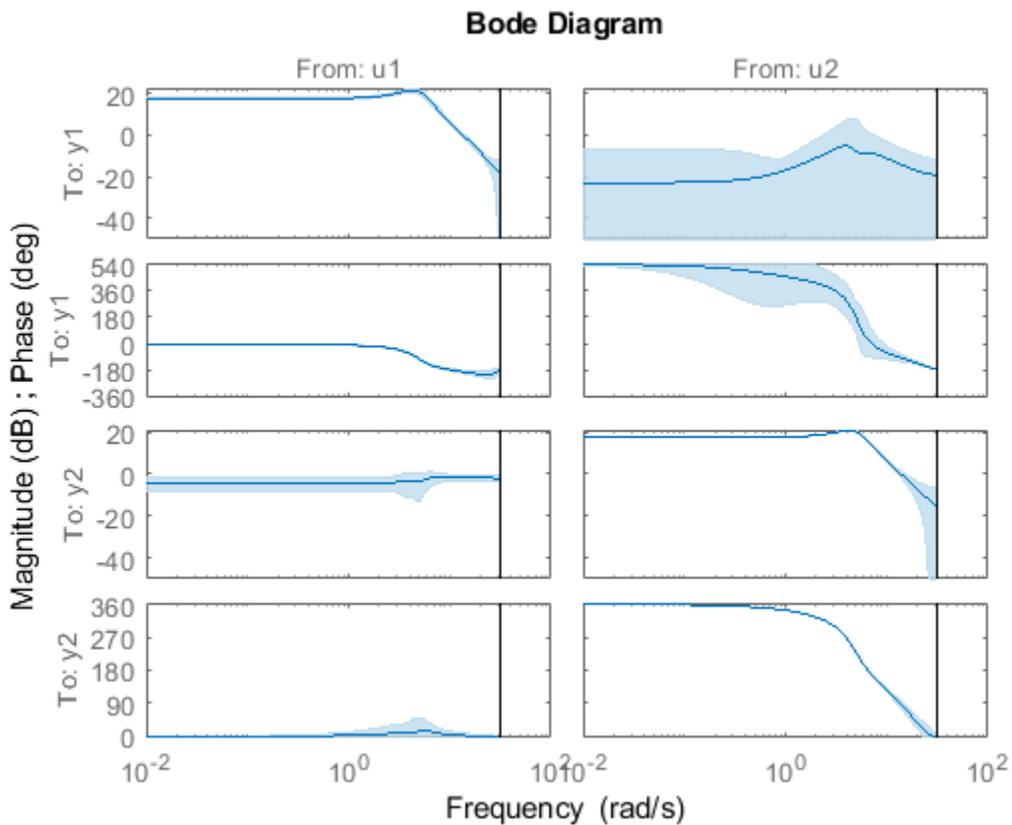
Estimate the model.

```
sys = polyest(data,[na nb nc nd nf nk]);
```

In the estimated ARMAX model, the cross terms, which model the effect of the first input on the second output and vice versa, are negligible. If you assigned higher orders to those dynamics, their estimation would show a high level of uncertainty.

Analyze the results.

```
h = bodeplot(sys);
showConfidence(h,3)
```



The responses from the cross terms show larger uncertainty.

## Tips

- In most situations, all the polynomials of an identified polynomial model are not simultaneously active. Set one or more of the orders `na`, `nc`, `nd` and `nf` to zero to simplify the model structure.

For example, you can estimate an Output-Error (OE) model by specifying `na`, `nc` and `nd` as zero.

Alternatively, you can use a dedicated estimating function for the simplified model structure. Linear polynomial estimation functions include `oe`, `bj`, `arx` and `armax`.

## Alternatives

- To estimate a polynomial model using time-series data, use `ar`.
- Use `polyest` to estimate a polynomial of arbitrary structure. If the structure of the estimated polynomial model is known, that is, you know which polynomials will be active, then use the appropriate dedicated estimating function. For examples, for an ARX model, use `arx`. Other polynomial model estimating functions include, `oe`, `armax`, and `bj`.
- To estimate a continuous-time transfer function, use `tfest`. You can also use `oe`, but only with continuous-time frequency-domain data.

## Extended Capabilities

### Automatic Parallel Support

Accelerate code by automatically running computation in parallel using Parallel Computing Toolbox™.

Parallel computing support is available for estimation using the `lsqnonlin` search method (requires Optimization Toolbox). To enable parallel computing, use `polyestOptions`, set `SearchMethod` to `'lsqnonlin'`, and set `SearchOptions.Advanced.UseParallel` to `true`.

For example:

```
opt = polyestOptions;  
opt.SearchMethod = 'lsqnonlin';  
opt.SearchOptions.Advanced.UseParallel = true;
```

## See Also

ar | armax | arx | bj | forecast | iddata | idpoly | oe | pem | polyestOptions |  
procest | ssest | tfest

## Topics

“Regularized Estimates of Model Parameters”

**Introduced in R2012a**

# polyestOptions

Option set for polyest

## Syntax

```
opt = polyestOptions  
opt = polyestOptions(Name,Value)
```

## Description

`opt = polyestOptions` creates the default option set for `polyest`.

`opt = polyestOptions(Name,Value)` creates an option set with the options specified by one or more `Name,Value` pair arguments.

## Input Arguments

### Name-Value Pair Arguments

Specify optional comma-separated pairs of `Name,Value` arguments. `Name` is the argument name and `Value` is the corresponding value. `Name` must appear inside quotes. You can specify several name and value pair arguments in any order as `Name1,Value1,...,NameN,ValueN`.

### InitialCondition — Handling of initial conditions

'auto' (default) | 'zero' | 'estimate' | 'backcast'

Handling of initial conditions during estimation, specified as one of the following values:

- 'zero' — The initial condition is set to zero.
- 'estimate' — The initial state is treated as an independent estimation parameter.
- 'backcast' — The initial state is estimated using the best least squares fit.

- 'auto' — The software chooses the method to handle initial states based on the estimation data.

### **Focus — Error to be minimized**

'prediction' (default) | 'simulation'

Error to be minimized in the loss function during estimation, specified as the comma-separated pair consisting of 'Focus' and one of the following values:

- 'prediction' — The one-step ahead prediction error between measured and predicted outputs is minimized during estimation. As a result, the estimation focuses on producing a good predictor model.
- 'simulation' — The simulation error between measured and simulated outputs is minimized during estimation. As a result, the estimation focuses on making a good fit for simulation of model response with the current inputs.

The Focus option can be interpreted as a weighting filter in the loss function. For more information, see “Loss Function and Model Quality Metrics”.

### **WeightingFilter — Weighting prefilter**

[] (default) | vector | matrix | cell array | linear system

Weighting prefilter applied to the loss function to be minimized during estimation. To understand the effect of WeightingFilter on the loss function, see “Loss Function and Model Quality Metrics”.

Specify WeightingFilter as one of the following values:

- [] — No weighting prefilter is used.
- Passbands — Specify a row vector or matrix containing frequency values that define desired passbands. You select a frequency band where the fit between estimated model and estimation data is optimized. For example, [w<sub>l</sub>, w<sub>h</sub>] where w<sub>l</sub> and w<sub>h</sub> represent lower and upper limits of a passband. For a matrix with several rows defining frequency passbands, [w<sub>1l</sub>, w<sub>1h</sub>; w<sub>2l</sub>, w<sub>2h</sub>; w<sub>3l</sub>, w<sub>3h</sub>; . . . ], the estimation algorithm uses the union of the frequency ranges to define the estimation passband.

Passbands are expressed in rad/TimeUnit for time-domain data and in FrequencyUnit for frequency-domain data, where TimeUnit and FrequencyUnit are the time and frequency units of the estimation data.

- SISO filter — Specify a single-input-single-output (SISO) linear filter in one of the following ways:

- A SISO LTI model
- `{A, B, C, D}` format, which specifies the state-space matrices of a filter with the same sample time as estimation data.
- `{numerator, denominator}` format, which specifies the numerator and denominator of the filter as a transfer function with same sample time as estimation data.

This option calculates the weighting function as a product of the filter and the input spectrum to estimate the transfer function.

- **Weighting vector** — Applicable for frequency-domain data only. Specify a column vector of weights. This vector must have the same length as the frequency vector of the data set, `Data.Frequency`. Each input and output response in the data is multiplied by the corresponding weight at that frequency.

### **EnforceStability** — Control whether to enforce stability of model

`false` (default) | `true`

Control whether to enforce stability of estimated model, specified as the comma-separated pair consisting of 'EnforceStability' and either `true` or `false`.

This option is not available for multi-output models with a non-diagonal *A* polynomial array.

Data Types: `logical`

### **EstimateCovariance** — Control whether to generate parameter covariance data

`true` (default) | `false`

Controls whether parameter covariance data is generated, specified as `true` or `false`.

If `EstimateCovariance` is `true`, then use `getcov` to fetch the covariance matrix from the estimated model.

### **Display** — Specify whether to display the estimation progress

'off' (default) | 'on'

Specify whether to display the estimation progress, specified as one of the following values:

- 'on' — Information on model structure and estimation results are displayed in a progress-viewer window.

- 'off' — No progress or results information is displayed.

## **InputOffset — Removal of offset from time-domain input data during estimation**

[ ] (default) | vector of positive integers | matrix

Removal of offset from time-domain input data during estimation, specified as the comma-separated pair consisting of 'InputOffset' and one of the following:

- A column vector of positive integers of length  $Nu$ , where  $Nu$  is the number of inputs.
- [ ] — Indicates no offset.
- $Nu$ -by- $Ne$  matrix — For multi-experiment data, specify InputOffset as an  $Nu$ -by- $Ne$  matrix.  $Nu$  is the number of inputs, and  $Ne$  is the number of experiments.

Each entry specified by InputOffset is subtracted from the corresponding input data.

## **OutputOffset — Removal of offset from time-domain output data during estimation**

[ ] (default) | vector | matrix

Removal of offset from time-domain output data during estimation, specified as the comma-separated pair consisting of 'OutputOffset' and one of the following:

- A column vector of length  $Ny$ , where  $Ny$  is the number of outputs.
- [ ] — Indicates no offset.
- $Ny$ -by- $Ne$  matrix — For multi-experiment data, specify OutputOffset as a  $Ny$ -by- $Ne$  matrix.  $Ny$  is the number of outputs, and  $Ne$  is the number of experiments.

Each entry specified by OutputOffset is subtracted from the corresponding output data.

## **Regularization — Options for regularized estimation of model parameters** structure

Options for regularized estimation of model parameters. For more information on regularization, see “Regularized Estimates of Model Parameters”.

Regularization is a structure with the following fields:

- Lambda — Constant that determines the bias versus variance tradeoff.

Specify a positive scalar to add the regularization term to the estimation cost.

The default value of zero implies no regularization.

**Default: 0**

- `R` — Weighting matrix.

Specify a vector of nonnegative numbers or a square positive semi-definite matrix. The length must be equal to the number of free parameters of the model.

For black-box models, using the default value is recommended. For structured and grey-box models, you can also specify a vector of `np` positive numbers such that each entry denotes the confidence in the value of the associated parameter.

The default value of 1 implies a value of `eye(npfree)`, where `npfree` is the number of free parameters.

**Default: 1**

- `Nominal` — The nominal value towards which the free parameters are pulled during estimation.

The default value of zero implies that the parameter values are pulled towards zero. If you are refining a model, you can set the value to `'model'` to pull the parameters towards the parameter values of the initial model. The initial parameter values must be finite for this setting to work.

**Default: 0**

**SearchMethod — Numerical search method used for iterative parameter estimation**

`'auto'` (default) | `'gn'` | `'gna'` | `'lm'` | `'grad'` | `'lsqnonlin'` | `'fmincon'`

Numerical search method used for iterative parameter estimation, specified as the comma-separated pair consisting of `'SearchMethod'` and one of the following:

- `'auto'` — A combination of the line search algorithms, `'gn'`, `'lm'`, `'gna'`, and `'grad'` methods is tried in sequence at each iteration. The first descent direction leading to a reduction in estimation cost is used.
- `'gn'` — Subspace Gauss-Newton least squares search. Singular values of the Jacobian matrix less than `GnPinvConstant*eps*max(size(J))*norm(J)` are discarded when computing the search direction.  $J$  is the Jacobian matrix. The Hessian matrix is approximated as  $J^T J$ . If there is no improvement in this direction, the function tries the gradient direction.

- 'gna' — Adaptive subspace Gauss-Newton search. Eigenvalues less than  $\gamma \cdot \max(sv)$  of the Hessian are ignored, where *sv* contains the singular values of the Hessian. The Gauss-Newton direction is computed in the remaining subspace. *gamma* has the initial value `InitialGnaTolerance` (see `Advanced in 'SearchOptions'` for more information). This value is increased by the factor `LMStep` each time the search fails to find a lower value of the criterion in fewer than five bisections. This value is decreased by the factor  $2 \cdot \text{LMStep}$  each time a search is successful without any bisections.
- 'lm' — Levenberg-Marquardt least squares search, where the next parameter value is  $-\text{pinv}(H+d \cdot I) \cdot \text{grad}$  from the previous one. *H* is the Hessian, *I* is the identity matrix, and *grad* is the gradient. *d* is a number that is increased until a lower value of the criterion is found.
- 'grad' — Steepest descent least squares search.
- 'lsqnonlin' — Trust-region-reflective algorithm of `lsqnonlin`. Requires Optimization Toolbox software.
- 'fmincon' — Constrained nonlinear solvers. You can use the sequential quadratic programming (SQP) and trust-region-reflective algorithms of the `fmincon` solver. If you have Optimization Toolbox software, you can also use the interior-point and active-set algorithms of the `fmincon` solver. Specify the algorithm in the `SearchOptions.Algorithm` option. The `fmincon` algorithms may result in improved estimation results in the following scenarios:
  - Constrained minimization problems when there are bounds imposed on the model parameters.
  - Model structures where the loss function is a nonlinear or non smooth function of the parameters.
  - Multi-output model estimation. A determinant loss function is minimized by default for multi-output model estimation. `fmincon` algorithms are able to minimize such loss functions directly. The other search methods such as 'lm' and 'gn' minimize the determinant loss function by alternately estimating the noise variance and reducing the loss value for a given noise variance value. Hence, the `fmincon` algorithms can offer better efficiency and accuracy for multi-output model estimations.

## **SearchOptions — Option set for the search algorithm**

search option set

Option set for the search algorithm, specified as the comma-separated pair consisting of 'SearchOptions' and a search option set with fields that depend on the value of SearchMethod.

**SearchOptions Structure When SearchMethod is Specified as 'gn', 'gna', 'lm', 'grad', or 'auto'**

Field Name	Description	Default
Tolerance	Minimum percentage difference between the current value of the loss function and its expected improvement after the next iteration, specified as a positive scalar. When the percentage of expected improvement is less than <code>Tolerance</code> , the iterations stop. The estimate of the expected loss-function improvement at the next iteration is based on the Gauss-Newton vector computed for the current parameter value.	0.01
MaxIterations	<p>Maximum number of iterations during loss-function minimization, specified as a positive integer. The iterations stop when <code>MaxIterations</code> is reached or another stopping criterion is satisfied, such as <code>Tolerance</code>.</p> <p>Setting <code>MaxIterations = 0</code> returns the result of the start-up procedure.</p> <p>Use <code>sys.Report.Termination.Iterations</code> to get the actual number of iterations during an estimation, where <code>sys</code> is an <code>idtf</code> model.</p>	20

Field Name	Description	Default
Advanced	Advanced search settings, specified as a structure with the following fields:	
Field Name	Description	Default
GnPinvConstant	Jacobian matrix singular value threshold, specified as a positive scalar. Singular values of the Jacobian matrix that are smaller than $\text{GnPinvConstant} * \max(\text{size}(J) * \text{norm}(J) * \text{eps})$ are discarded when computing the search direction. Applicable when SearchMethod is 'gn'.	10000
InitialGnaTolerance	Initial value of $\gamma$ , specified as a positive scalar. Applicable when SearchMethod is 'gna'.	0.0001
LMStartValue	Starting value of search-direction length $d$ in the Levenberg-Marquardt method, specified as a positive scalar. Applicable when SearchMethod is 'lm'.	0.001
LMStep	Size of the Levenberg-Marquardt step, specified as a positive integer. The next value of the search-direction length $d$ in the Levenberg-Marquardt method is LMStep times the previous one. Applicable when SearchMethod is 'lm'.	2
MaxBisections	Maximum number of bisections used for line search along the search direction, specified as a positive integer.	25
MaxFunctionEvaluations	Maximum number of calls to the model file, specified as a positive integer. Iterations stop if the number of calls to the model file exceeds this value.	Inf
MinParameterChange	Smallest parameter update allowed per iteration, specified as a nonnegative scalar.	0
RelativeImprovement	Relative improvement threshold, specified as a nonnegative scalar. Iterations stop if the relative	0

<b>Field Name</b>	<b>Description</b>		<b>Default</b>
	<b>Field Name</b>	<b>Description</b>	<b>Default</b>
		improvement of the criterion function is less than this value.	
	StepReduction	<p>Step reduction factor, specified as a positive scalar that is greater than 1. The suggested parameter update is reduced by the factor <code>StepReduction</code> after each try. This reduction continues until <code>MaxBisections</code> tries are completed or a lower value of the criterion function is obtained.</p> <p><code>StepReduction</code> is not applicable for SearchMethod 'lm' (Levenberg-Marquardt method).</p>	2

**SearchOptions Structure When SearchMethod is Specified as 'lsqnonlin'**

<b>Field Name</b>	<b>Description</b>	<b>Default</b>
FunctionTolerance	Termination tolerance on the loss function that the software minimizes to determine the estimated parameter values, specified as a positive scalar.  The value of FunctionTolerance is the same as that of <code>opt.SearchOptions.Advanced.TolFun</code> .	1e-5
StepTolerance	Termination tolerance on the estimated parameter values, specified as a positive scalar.  The value of StepTolerance is the same as that of <code>opt.SearchOptions.Advanced.TolX</code> .	1e-6
MaxIterations	Maximum number of iterations during loss-function minimization, specified as a positive integer. The iterations stop when MaxIterations is reached or another stopping criterion is satisfied, such as FunctionTolerance.  The value of MaxIterations is the same as that of <code>opt.SearchOptions.Advanced.MaxIter</code> .	20
Advanced	Advanced search settings, specified as an option set for <code>lsqnonlin</code> .  For more information, see the Optimization Options table in "Optimization Options" (Optimization Toolbox).	Use <code>optimset('lsqnonlin')</code> to create a default option set.

**SearchOptions Structure When SearchMethod is Specified as 'fmincon'**

Field Name	Description	Default
Algorithm	<p>fmincon optimization algorithm, specified as one of the following:</p> <ul style="list-style-type: none"> <li>• 'sqp' — Sequential quadratic programming algorithm. The algorithm satisfies bounds at all iterations, and it can recover from NaN or Inf results. It is not a large-scale algorithm. For more information, see “Large-Scale vs. Medium-Scale Algorithms” (Optimization Toolbox).</li> <li>• 'trust-region-reflective' — Subspace trust-region method based on the interior-reflective Newton method. It is a large-scale algorithm.</li> <li>• 'interior-point' — Large-scale algorithm that requires Optimization Toolbox software. The algorithm satisfies bounds at all iterations, and it can recover from NaN or Inf results.</li> <li>• 'active-set' — Requires Optimization Toolbox software. The</li> </ul>	'sqp'

Field Name	Description	Default
	<p>algorithm can take large steps, which adds speed. It is not a large-scale algorithm.</p> <p>For more information about the algorithms, see “Constrained Nonlinear Optimization Algorithms” (Optimization Toolbox) and “Choosing the Algorithm” (Optimization Toolbox).</p>	
FunctionTolerance	Termination tolerance on the loss function that the software minimizes to determine the estimated parameter values, specified as a positive scalar.	1e-6
StepTolerance	Termination tolerance on the estimated parameter values, specified as a positive scalar.	1e-6
MaxIterations	Maximum number of iterations during loss function minimization, specified as a positive integer. The iterations stop when MaxIterations is reached or another stopping criterion is satisfied, such as FunctionTolerance.	100

### Advanced – Additional advanced options

structure

Additional advanced options, specified as a structure with the following fields:

- `ErrorThreshold` — Specifies when to adjust the weight of large errors from quadratic to linear.

Errors larger than `ErrorThreshold` times the estimated standard deviation have a linear weight in the loss function. The standard deviation is estimated robustly as the median of the absolute deviations from the median of the prediction errors, divided by 0.7. For more information on robust norm choices, see section 15.2 of [2].

`ErrorThreshold = 0` disables robustification and leads to a purely quadratic loss function. When estimating with frequency-domain data, the software sets `ErrorThreshold` to zero. For time-domain data that contains outliers, try setting `ErrorThreshold` to 1.6.

**Default:** 0

- `MaxSize` — Specifies the maximum number of elements in a segment when input-output data is split into segments.

`MaxSize` must be a positive integer.

**Default:** 250000

- `StabilityThreshold` — Specifies thresholds for stability tests.

`StabilityThreshold` is a structure with the following fields:

- `s` — Specifies the location of the right-most pole to test the stability of continuous-time models. A model is considered stable when its right-most pole is to the left of `s`.

**Default:** 0

- `z` — Specifies the maximum distance of all poles from the origin to test stability of discrete-time models. A model is considered stable if all poles are within the distance `z` from the origin.

**Default:** `1+sqrt(eps)`

- `AutoInitThreshold` — Specifies when to automatically estimate the initial condition.

The initial condition is estimated when

$$\frac{\|y_{p,z} - y_{meas}\|}{\|y_{p,e} - y_{meas}\|} > \text{AutoInitThreshold}$$

- $y_{meas}$  is the measured output.
- $y_{p,z}$  is the predicted output of a model estimated using zero initial states.
- $y_{p,e}$  is the predicted output of a model estimated using estimated initial states.

Applicable when `InitialCondition` is 'auto'.

**Default:** 1.05

## Output Arguments

**opt** — Options set for `polyest`

`polyestOptions` option set

Option set for `polyest`, returned as an `polyestOptions` option set.

## Examples

### Create Default Option Set for Polynomial Estimation

```
opt = polyestOptions;
```

### Specify Options for Polynomial Estimation

Create an option set for `polyest` where you enforce model stability and set the `Display` to 'on'.

```
opt = polyestOptions('EnforceStability',true,'Display','on');
```

Alternatively, use dot notation to set the values of `opt`.

```
opt = polyestOptions;  
opt.EnforceStability = true;  
opt.Display = 'on';
```

## References

- [1] Wills, Adrian, B. Ninness, and S. Gibson. "On Gradient-Based Search for Multivariable System Estimates". *Proceedings of the 16th IFAC World Congress, Prague, Czech Republic, July 3-8, 2005*. Oxford, UK: Elsevier Ltd., 2005.
- [2] Ljung, L. *System Identification: Theory for the User*. Upper Saddle River, NJ: Prentice-Hall PTR, 1999.

## See Also

polyest

## Topics

"Loss Function and Model Quality Metrics"

**Introduced in R2012a**

## polyreg

Powers and products of standard regressors

### Syntax

```
R = polyreg(model)
R = polyreg(model, 'MaxPower', n)
R = polyreg(model, 'MaxPower', n, 'CrossTerm', CrossTermVal)
```

### Description

`R = polyreg(model)` creates an array *R* of polynomial regressors up to the power 2. If a model order has input *u* and output *y*, *na*=*nb*=2, and delay *nk*=1, polynomial regressors are  $y(t-1)^2$ ,  $u(t-1)^2$ ,  $y(t-2)^2$ ,  $u(t-2)^2$ . *model* is an `idnlarx` object. You must add these regressors to the *model* by assigning the `CustomRegressors` *model* property or by using `addreg`.

`R = polyreg(model, 'MaxPower', n)` creates an array *R* of polynomial regressors up to the power *n*. Excludes terms of power 1 and cross terms, such as  $y(t-1)*u(t-1)$ .

`R = polyreg(model, 'MaxPower', n, 'CrossTerm', CrossTermVal)` creates an array *R* of polynomial regressors up to the power *n* and includes cross terms (products of standards regressors) when *CrossTermVal* is 'on'. By default, *CrossTermVal* is 'off'.

## Examples

### Create Polynomial Regressors Up To Power 2

Estimate a nonlinear ARX model with *na* = 2, *nb* = 2, and *nk* = 1, and nonlinearity estimator `wavenet`.

```
load iddata1
m = nlarx(z1,[2 2 1]);
```

Create polynomial regressors.

```
R = polyreg(m);
```

Estimate the model.

```
m = nlarx(z1,[2 2 1], 'wavenet', 'CustomReg', R);
```

View all model regressors (standard and custom).

```
getreg(m)
```

Regressors:

```
y1(t-1)
y1(t-2)
u1(t-1)
u1(t-2)
y1(t-1).^2
y1(t-2).^2
u1(t-1).^2
u1(t-2).^2
```

## Create Polynomial Regressors Up To Power 3

Estimate a nonlinear ARX model with  $na = 2$ ,  $nb = 1$ , and  $nk = 1$ , and nonlinearity estimator `wavenet`.

```
load iddata1
m = nlarx(z1,[2 1 1]);
```

Create polynomial regressors.

```
R = polyreg(m, 'MaxPower', 3, 'CrossTerm', 'on')
```

16x1 array of Custom Regressors with fields: Function, Arguments, Delays, Vectorized

If the model `m` has three standard regressors `a`, `b` and `c`, then `R` includes the terms  $a^2$ ,  $b^2$ ,  $c^2$ ,  $ab$ ,  $ac$ ,  $bc$ ,  $a^2b$ ,  $a^2c$ ,  $ab^2$ ,  $abc$ ,  $ac^2$ ,  $b^2c$ ,  $bc^2$ ,  $a^3$ ,  $b^3$ , and  $c^3$ .

Estimate the model.

```
m = nlarx(z1,[2 1 1], 'wavenet', 'CustomReg',R);
```

## See Also

[addreg](#) | [customreg](#) | [getreg](#) | [idnlarx](#) | [nlarx](#)

## Topics

“Identifying Nonlinear ARX Models”

**Introduced in R2007a**

## poly1d

Class representing single-variable polynomial nonlinear estimator for Hammerstein-Wiener models

### Syntax

```
t=poly1d('Degree',n)
t=poly1d('Coefficients',C)
t=poly1d(n)
```

### Description

`poly1d` is an object that stores the single-variable polynomial nonlinear estimator for Hammerstein-Wiener models.

You can use the constructor to create the nonlinearity object, as follows:

`t=poly1d('Degree',n)` creates a polynomial nonlinearity estimator object of  $n$ th degree.

`t=poly1d('Coefficients',C)` creates a polynomial nonlinearity estimator object with coefficients  $C$ .

`t=poly1d(n)` a polynomial nonlinearity estimator object of  $n$ th degree.

Use `evaluate(p,x)` to compute the value of the function defined by the `poly1d` object  $p$  at  $x$ .

### poly1d Properties

After creating the object, you can use `get` or dot notation to access the object property values. For example:

```
% List all property values
get(p)
```

```
% Get value of Coefficients property
p.Coefficients
```

Property Name	Description
Degree	Positive integer specifies the degree of the polynomial Default=1.  For example: <code>poly1d('Degree',3)</code>
Coefficients	1-by-(n+1) matrix containing the polynomial coefficients.

## Examples

Use `poly1s` to specify the single-variable polynomial nonlinearity estimator in Hammerstein-Wiener models. For example:

```
m=nlhw(Data,Orders,poly1d('deg',3),[]);
```

where 'deg' is an abbreviation for the property 'Degree'.

## Tips

Use `poly1d` to define a nonlinear function  $y = F(x)$ , where  $F$  is a single-variable polynomial function of  $x$ :

$$F(x) = c(1)x^n + c(2)x^{(n-1)} + \dots + c(n)x + c(n+1)$$

## See Also

`nlhw`

**Introduced in R2007b**

## predict

Predict K-step ahead model output

This `predict` command computes the K-step ahead output of an identified model using measured input-output data. To identify the model, you first collect all the input-output data and then estimate the model parameters offline. To perform online state estimation of a nonlinear system using real-time data, use the `predict` command for extended and unscented Kalman filters instead.

## Syntax

```
yp = predict(sys,data,K)
yp = predict(sys,data,K,opt)
[yp,x0,sys_pred] = predict( ___ )

predict(sys,data,K, ___ )
predict(sys,Linespec,data,K, ___ )
predict(sys1,...,sysN,data,K, ___ )
predict(sys1,Linespec1,...,sysN,LinespecN,data,K, ___ )
```

## Description

`yp = predict(sys,data,K)` predicts the output of an identified model `sys`, `K` steps ahead using the measured input-output data.

`predict` command predicts the output response over the time span of measured data. In contrast, `forecast` performs prediction into the future in a time range beyond the last instant of measured data. Use `predict` to validate `sys` over the time span of measured data.

`yp = predict(sys,data,K,opt)` uses the option set `opt` to specify additional prediction options such as handling of initial conditions and data offsets.

`[yp,x0,sys_pred] = predict( ___ )` also returns the estimated values for initial states `x0` and a predictor model `sys_pred`. Use with any of the previous input argument combinations.

`predict(sys,data,K, ___)` plots the predicted output. Use with any of the previous input argument combinations. To change display options in the plot, right-click the plot to access the context menu. For more details about the menu, see “Tips” on page 1-1231.

You can also plot the predicted model response using the `compare` command. The `compare` command compares the prediction results with observed data and displays a quantitative goodness of fit.

`predict(sys,Linespec,data,K, ___)` uses `Linespec` to specify the line type, marker symbol, and color.

`predict(sys1,...,sysN,data,K, ___)` plots the predicted outputs for multiple identified models. `predict` automatically chooses colors and line styles.

`predict(sys1,Linespec1,...,sysN,LinespecN,data,K, ___)` uses the line type, marker symbol, and color specified for each model.

## Examples

### Predict Time Series Model Response

Simulate time-series data.

```
init_sys = idpoly([1 -0.99],[],[1 -1 0.2]);  
opt = simOptions('AddNoise',true);  
u = iddata([],zeros(400,0),1);  
data = sim(init_sys,u,opt);
```

`data` is an `iddata` object containing the simulated response data of a time series model.

Estimate an ARMAX model by using `data` as estimation data.

```
na = 1;  
nb = 2;  
sys = armax(data(1:200),[na nb]);
```

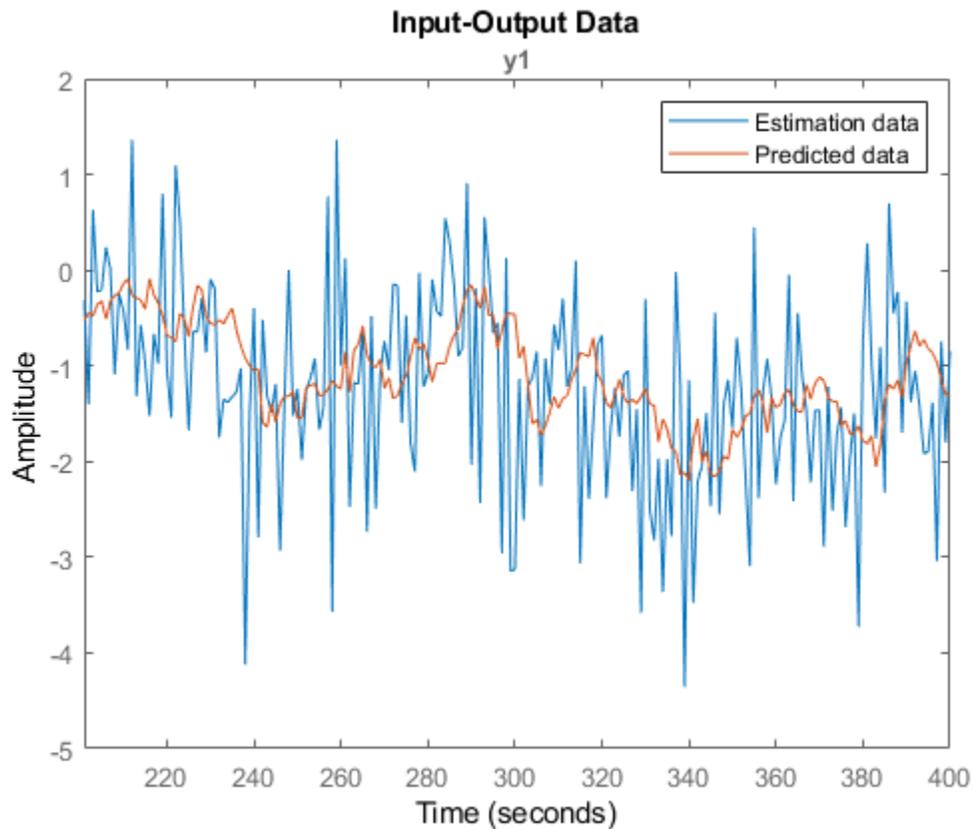
Predict the output of the model using a prediction horizon of 4.

```
K = 4;  
yp = predict(sys,data,K);
```

`yp` is an `iddata` object. The predicted output is returned in the `OutputData` property of the object.

Compare the predicted and estimated data outputs.

```
plot(data(201:400),yp(201:400));  
legend('Estimation data','Predicted data');
```



Alternatively, to plot the predicted response and estimation data, use `compare(sys,data,K)`.

## Plot Predicted Output for Multiple Models

Load the estimation data.

```
load iddata1;  
data = z1;
```

Estimate an ARX model of order [2 2 1].

```
sys1 = arx(data,[2 2 1]);
```

Estimate a transfer function with 2 poles.

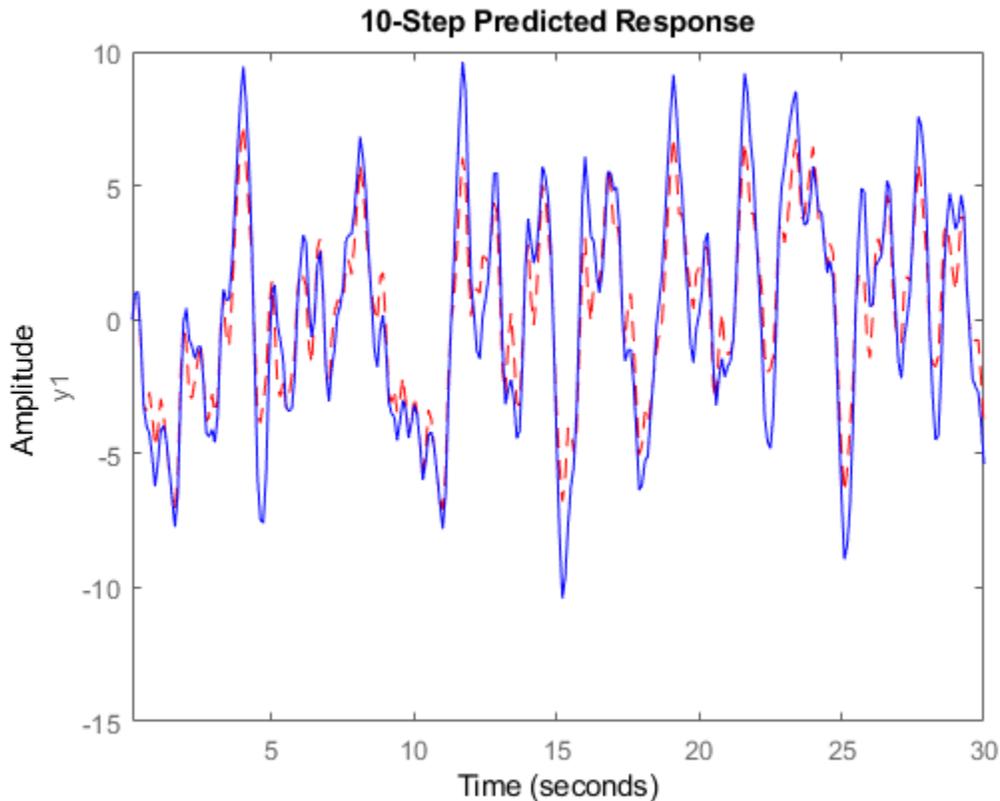
```
sys2 = tfest(data,2);
```

Create a `predict` option set to specify zero initial conditions for prediction.

```
opt = predictOptions('InitialCondition','z');
```

Plot the predicted outputs for the estimated models. Use the specified prediction option set, `opt`, and specify prediction horizon as 10. Specify line styles for plotting the predicted output of each system.

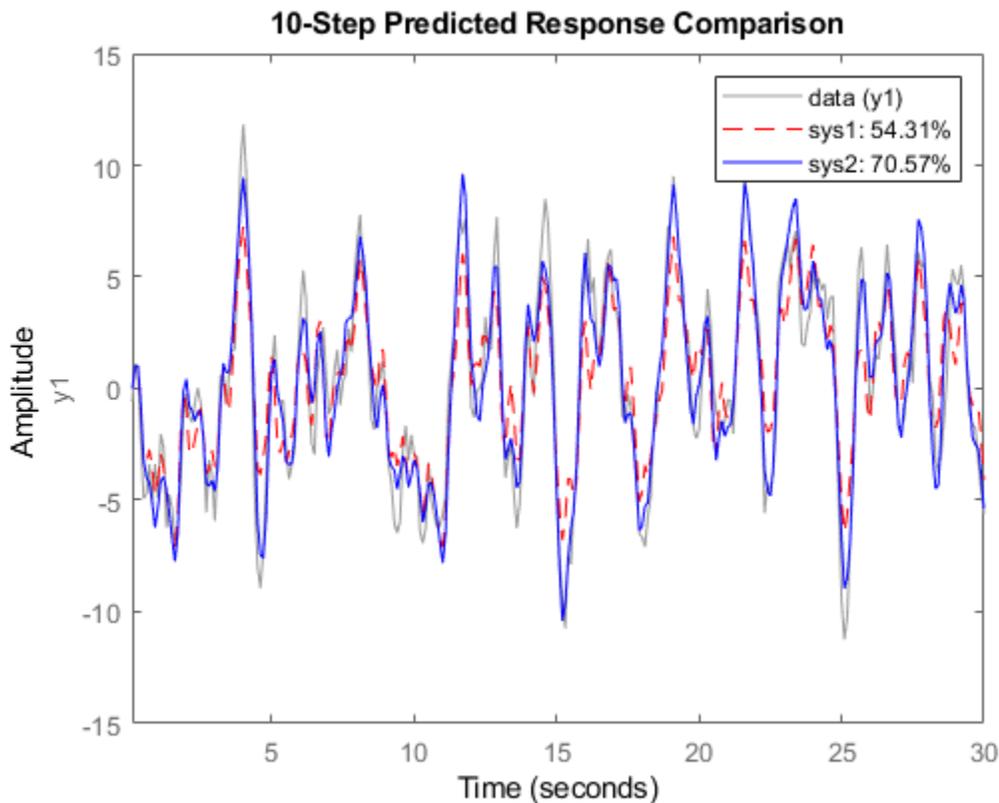
```
predict(sys1,'r--',sys2,'b',data,10,opt);
```



To change the display options, right-click the plot to access the context menu. For example, to view the estimation data, select **Show Validation Data** from the context menu. To view the prediction error, select **Prediction Error Plot**.

You can also plot the predicted response using the `compare` command. To do so, first create an option set for `compare` to specify the use of zero initial conditions.

```
opt = compareOptions('InitialCondition', 'z');  
compare(data, sys1, 'r--', sys2, 'b', 10, opt);
```



### Reproduce Prediction Results by Simulation

Use estimation data to estimate a model, and then compute the predicted model output and predictor model using the `predict` command. Simulate the predictor model to reproduce the predicted output.

Load estimation data.

```
load iddata3 z3  
data = z3;
```

Estimate a polynomial model from the data.

```
sys = polyest(z3,[2 2 2 0 0 1]);
```

Predict the system response using prediction horizon 4.

```
K = 4;  
[yp,x0,sysp] = predict(sys,data,K);
```

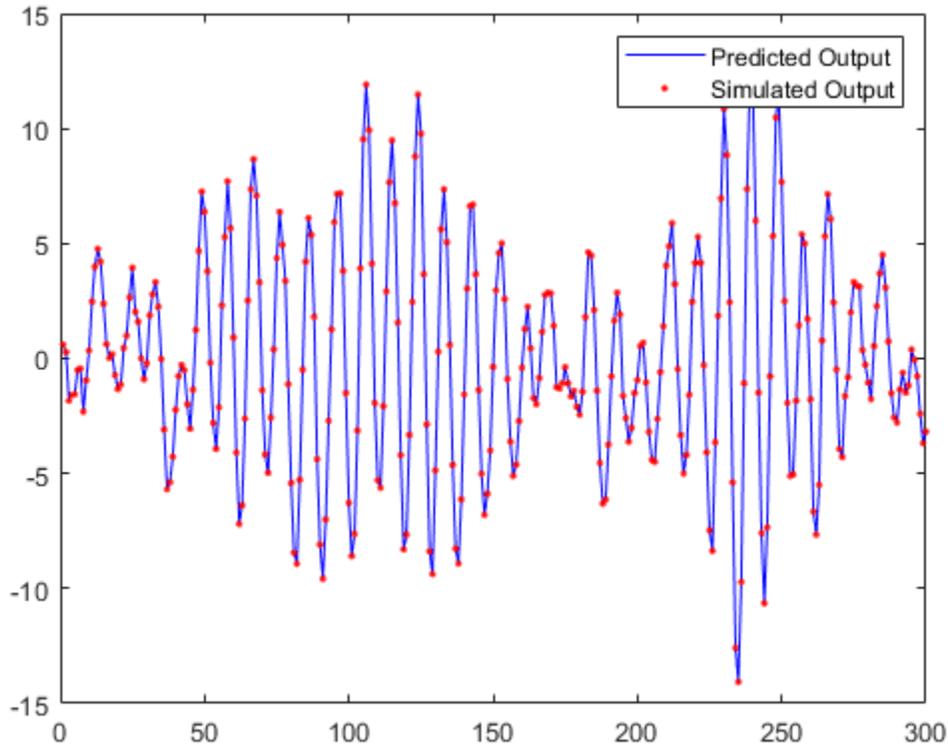
yp is the predicted model response, x0 is the estimated value for initial states, and sysp is the predictor model.

Simulate the predictor model with inputs [data.OutputData,data.InputData] and initial conditions x0.

```
opt = simOptions;  
opt.InitialCondition = x0;  
ys = sim(sysp,[data.OutputData,data.InputData],opt);
```

Plot the predicted and simulated outputs.

```
t = yp.SamplingInstants;  
plot(t,yp.OutputData,'b',t,ys,'.r');  
legend('Predicted Output','Simulated Output')
```



### Understand Use of Historical Data for Model Prediction

Perform model prediction using historical data to specify initial conditions. You first predict using the `predict` command and specify the historical data using the `predictOptions` option set. You then reproduce the predicted response by manually mapping the historical data to initial states.

Load a two-input, one-output dataset.

```
load iddata7 z7
```

Identify a fifth-order state-space model using the data.

```
sys = n4sid(z7,5);
```

Split the dataset into two parts.

```
zA = z7(1:15);  
zB = z7(16:end);
```

Suppose that you want to compute the 10-step-ahead prediction of the response of the identified system for data zB. For initial conditions, use the signal values in zA as the historical record. That is, the input and output values for the time immediately preceding data in zB.

```
I0 = struct('Input',zA.InputData,'Output',zA.OutputData);  
opt = predictOptions('InitialCondition',I0);
```

Generate the 10-step-ahead prediction for data zB using the specified initial conditions and `predict`.

```
[yp,x0,Predictor] = predict(sys,zB,10,opt);
```

`yp` is the predicted model response, `x0` are the initial states corresponding to the predictor model `Predictor`. You can simulate `Predictor` using `x0` as initial conditions to reproduce `yp.OutputData`.

Now reproduce the output by manually mapping the historical data to initial states. To do so, minimize 1-step prediction errors over the time span of zA.

```
x0est = data2state(sys,zA);
```

`x0est` contains the values of the five states of `sys` at the time instant immediately after the most recent data sample in zA.

The `Predictor` has more states than the original system due to the 10-step prediction horizon. Specify the additional states induced by the horizon to zero initial values, and then append `x0est`.

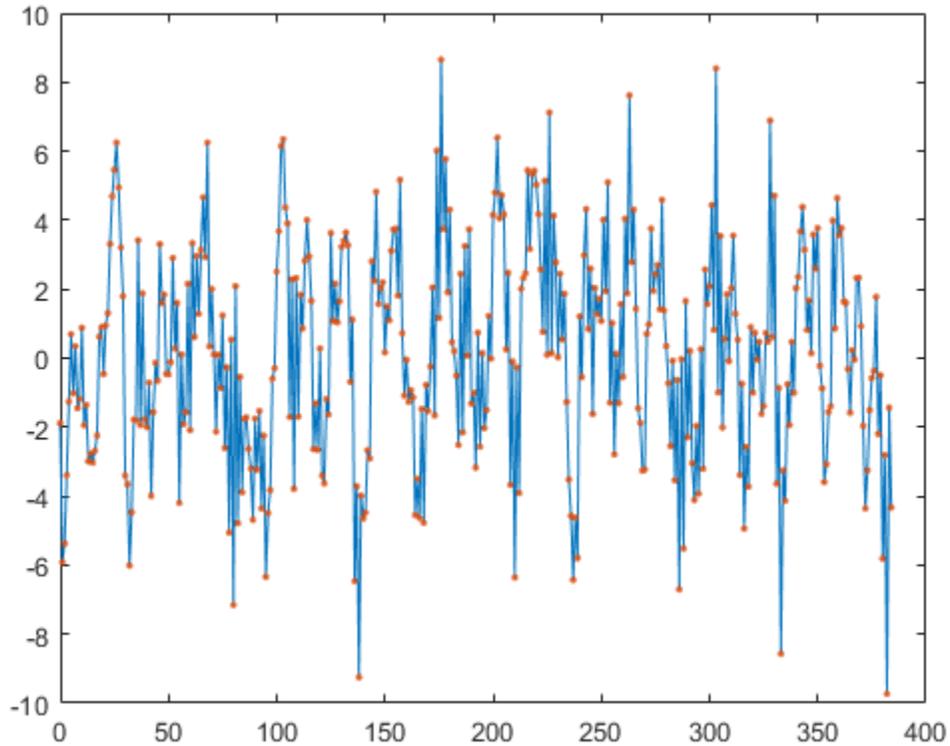
```
x0Predictor = zeros(order(Predictor),1);  
x0Predictor(end-4:end) = x0est;
```

Simulate the predictor using `[zB.OutputData,zB.InputData]` as the input signal and `x0Predictor` as initial conditions.

```
uData = [zB.OutputData,zB.InputData]; % signals required for prediction  
[ysim,t,xsim] = lsim(Predictor,uData,[],x0Predictor);
```

Plot the predicted output of the `predict` command `yp.OutputData` and the manually computed results `ysim`.

```
plot(t,yp.OutputData,t,ysim, 'r.')
```



`ysim` is the same as `yp.OutputData`.

## Input Arguments

**sys** — Identified model

linear model | nonlinear model

Identified model whose output is to be predicted, specified as one of the following:

- Linear model — `idpoly`, `idproc`, `idss`, `idtf`, or `idgrey`
- Nonlinear model — `idnlgrey`, `idnlhw`, or `idnlarx`

When `sys` is an `idnlhw` or `idnlgrey` model, the predicted output `yp` is the same as the simulated response computed using `data.InputData` as input.

If a model is unavailable, estimate `sys` from `data` using commands such as `ar`, `armax`, `tfest`, `nlarx`, and `ssest`.

### **data — Measured input-output data**

`iddata` object | matrix of doubles

Measured input-output data, specified as one of the following:

- `iddata` object — Use observed input and output signals to create an `iddata` object. For time-series data (no inputs), specify as an `iddata` object with no inputs `iddata(output, [])`.
- Matrix of doubles — For models with  $N_u$  inputs and  $N_y$  outputs, specify `data` as an  $N$ -by- $(N_y+N_u)$  matrix. Where,  $N$  is the number of observations.

For time series data, specify as an  $N$ -by- $N_y$  matrix.

### **K — Prediction horizon**

1 (default) | positive integer | `Inf`

Prediction horizon, specified as one of the following:

- Positive integer — Output `yp` is calculated  $K$  steps into the future, where  $K$  represents a multiple of `data` sample time.

The output at time instant  $t$  is calculated using previously measured outputs up to time  $t-K$  and inputs up to the time instant  $t$ .

- `Inf` — No previous outputs are used in the computation, and `predict` returns the same result as simulation using the `sim` command.

For Output-Error models, there is no difference between the  $K$  step-ahead predictions and the simulated output. This is because Output-Error models only use past inputs to predict future outputs.

---

**Note** For careful model validation, a one-step-ahead prediction ( $K = 1$ ) is usually not a good test for validating the model `sys` over the time span of measured data. Even the trivial one step-ahead predictor,  $\hat{y}(t) = y(t-1)$ , can give good predictions. So a poor model may look fine for one-step-ahead prediction of data that has a small sample time. Prediction with  $K = \text{Inf}$ , which is the same as performing simulation with `sim` command, can lead to diverging outputs because low-frequency disturbances in the data are emphasized, especially for models with integration. Use a  $K$  value between 1 and `Inf` to capture the mid-frequency behavior of the measured data.

---

### **opt** — Prediction options

`predictOptions` option set

Prediction options, specified as a `predictOptions` option set. Use the option set to specify prediction options such as handling of initial conditions and data offsets.

### **Linespec** — Line style, marker, and color

character vector

Line style, marker, and color, specified as a character vector. For example, 'b' or 'b+:'.

For more information about configuring `Linespec`, see the `Linespec` argument of `plot`.

## **Output Arguments**

### **yp** — Predicted output response

`iddata` object | matrix of doubles

Predicted output response, returned as one of the following:

- `iddata` object — When data is an `iddata` object. The `OutputData` property of `yp` stores the values of the predicted output. The time variable takes values in the range represented by `data.SamplingInstants`.
- Matrix of doubles — When data is a matrix of doubles.

The output at time instant `t` is calculated using previously measured outputs up to time `t-K` and inputs up to the time instant `t`. In other words, the predicted response at time point `r` of measured data is stored in the `r+K-1` sample of `yp`. Note that at time `r`, the future inputs `u(r+1)`, `u(r+2)`, ..., `u(r+K)` required for prediction are assumed to be

known. For multi-experiment data, `yp` contains a predicted data set for each experiment. The time span of the predicted outputs matches that of the observed data.

When `sys` is specified using an `idnlhw` or `idnlgrey` model, `yp` is the same as the simulated response computed using `data.InputData` as input.

## **x0 — Estimated initial states**

column vector | cell array

Estimated initial states corresponding to the predictor model `sys_pred`, returned as a column vector of size equal to number of states. If `data` is multi-experiment, `x0` is a cell array of size  $N_e$ , where  $N_e$  is the number of experiments. You can simulate `sys_pred` using `x0` as initial conditions, to reproduce the results of prediction. For an example, see “Reproduce Prediction Results by Simulation” on page 1-1223.

If you specify `opt.InitialCondition` as 'z', the initial conditions of data are not estimated and `x0` equals 0.

If `sys` is an `idnlarx` model, `x0` is returned empty.

## **sys\_pred — Predictor model**

dynamic system model | array of models

Predictor model, returned as a dynamic system model. For multi-experiment data, `sys_pred` is an array of models, with one entry for each experiment. You can use the predictor model `sys_pred` and estimated initial states `x0` to reproduce the results of prediction:

- If `sys` is a linear model, the predictor model is returned as either a model of the same type as `sys` or as a state-space version of the model (`idss`). To reproduce the results of prediction, simulate `sys_pred` using [`data.OutputData` `data.InputData`] as input and `x0` as initial conditions. The simulation output is the same as the predicted output `yp.OutputData`. For an example, see “Reproduce Prediction Results by Simulation” on page 1-1223.
- When `sys` is a nonlinear grey-box model (`idnlgrey`) or Hammerstein-Wiener model (`idnlhw`), the noise-component of the model is trivial, and so the predictor model is the same as the model. `sys_pred` is returned empty. To reproduce the results of prediction, simulate `sys` using initial conditions `x0`. For a definition of the states of `idnlhw` models, see “Definition of `idnlhw` States” on page 1-672.
- If `sys` is a nonlinear ARX model (`idnlarx`), `sys_pred` and `x0` are returned empty. You cannot reproduce the prediction results by simulation.

For discrete-time data that is time-domain or frequency-domain data with sample time  $T_s$  greater than zero, `sys_pred` is a discrete-time model, even if `sys` is a continuous-time model.

## Tips

- Right-clicking the plot of the predicted output opens the context menu, where you can access the following options:
  - **Systems** — Select systems to view predicted response. By default, the response of all systems is plotted.
  - **Data Experiment** — For multi-experiment data only. Toggle between data from different experiments.
  - **Characteristics** — View the following data characteristics:
    - **Peak Value** — View the absolute peak value of the data. Applicable for time-domain data only.
    - **Peak Response** — View peak response of the data. Applicable for frequency-response data only.
    - **Mean Value** — View mean value of the data. Applicable for time-domain data only.
  - **Show** — For frequency-domain and frequency-response data only.
    - **Magnitude** — View magnitude of frequency response of the system.
    - **Phase** — View phase of frequency response of the system.
  - **Show Validation Data** — Plot data used to predict the model response.
  - **I/O Grouping** — For datasets containing more than one input or output channel. Select grouping of input and output channels on the plot.
    - **None** — Plot input-output channels in their own separate axes.
    - **All** — Group all input channels together and all output channels together.
  - **I/O Selector** — For datasets containing more than one input or output channel. Select a subset of the input and output channels to plot. By default, all output channels are plotted.
  - **Grid** — Add grids to the plot.

- **Normalize** — Normalize the y-scale of all data in the plot.
- **Full View** — Return to full view. By default, the plot is scaled to full view.
- **Prediction Horizon** — Set the prediction horizon, or choose simulation.
- **Initial Condition** — Specify handling of initial conditions. Not applicable for frequency-response data.

Specify as one of the following:

- **Estimate** — Treat the initial conditions as estimation parameters.
- **Zero** — Set all initial conditions to zero.
- **Absorb delays and estimate** — Absorb nonzero delays into the model coefficients and treat the initial conditions as estimation parameters. Use this option for discrete-time models only.
- **Predicted Response Plot** — Plot the predicted model response. By default, the response plot is shown.
- **Prediction Error Plot** — Plot the error between the model response and prediction data.
- **Properties** — Open the Property Editor dialog box to customize plot attributes.

## See Also

`compare` | `forecast` | `iddata` | `pe` | `predictOptions` | `sim` | `simsd`

## Topics

“Simulate and Predict Identified Model Output”

“Simulation and Prediction at the Command Line”

**Introduced before R2006a**

## predict

Predict state and state estimation error covariance at next time step using extended or unscented Kalman filter, or particle filter

The `predict` command predicts the state and state estimation error covariance of an `extendedKalmanFilter`, `unscentedKalmanFilter` or `particleFilter` object at the next time step. To implement the extended or unscented Kalman filter algorithms, use the `predict` and `correct` commands together. If the current output measurement exists, you can use `predict` and `correct`. If the measurement is missing, you can only use `predict`. For information about the order in which to use the commands, see “Using `predict` and `correct` Commands” on page 1-1240.

Use this `predict` command for online state estimation using real-time data. When data is not available in real time, to compute the K-step ahead output of an identified model, use `predict` for offline estimation.

## Syntax

```
[PredictedState,PredictedStateCovariance] = predict(obj)
[PredictedState,PredictedStateCovariance] = predict(obj,Us1,...Usn)
```

## Description

`[PredictedState,PredictedStateCovariance] = predict(obj)` predicts state estimate and state estimation error covariance of an extended or unscented Kalman filter, or particle filter object `obj` at the next time step.

You create `obj` using the `extendedKalmanFilter`, `unscentedKalmanFilter` or `particleFilter` commands. You specify the state transition function and measurement function of your nonlinear system in `obj`. You also specify whether the process and measurement noise terms are additive or nonadditive in these functions. The `State` property of the object stores the latest estimated state value. Assume that at time step `k`, `obj.State` is  $\hat{x}[k|k]$ . This value is the state estimate for time `k`, estimated using measured outputs until time `k`. When you use the `predict` command, the software returns  $\hat{x}[k+1|k]$  in the `PredictedState` output. Where  $\hat{x}[k+1|k]$  is the state estimate

for time  $k+1$ , estimated using measured output until time  $k$ . The command returns the state estimation error covariance of  $\hat{x}[k+1|k]$  in the `PredictedStateCovariance` output. The software also updates the `State` and `StateCovariance` properties of `obj` with these corrected values.

Use this syntax if the state transition function  $f$  that you specified in `obj.StateTransitionFcn` has one of the following forms:

- $x(k) = f(x(k-1))$  — for additive process noise.
- $x(k) = f(x(k-1), w(k-1))$  — for nonadditive process noise.

Where  $x$  and  $w$  are the state and process noise of the system. The only inputs to  $f$  are the states and process noise.

`[PredictedState, PredictedStateCovariance] = predict(obj, Us1, ... Usn)` specifies additional input arguments, if the state transition function of the system requires these inputs. You can specify multiple arguments.

Use this syntax if your state transition function  $f$  has one of the following forms:

- $x(k) = f(x(k-1), Us1, \dots Usn)$  — for additive process noise.
- $x(k) = f(x(k-1), w(k-1), Us1, \dots Usn)$  — for nonadditive process noise.

## Examples

### Estimate States Online Using Unscented Kalman Filter

Estimate the states of a van der Pol oscillator using an unscented Kalman filter algorithm and measured output data. The oscillator has two states and one output.

Create an unscented Kalman filter object for the oscillator. Use previously written and saved state transition and measurement functions, `vdpStateFcn.m` and `vdpMeasurementFcn.m`. These functions describe a discrete-approximation to a van der Pol oscillator with nonlinearity parameter,  $\mu$ , equal to 1. The functions assume additive process and measurement noise in the system. Specify the initial state values for the two states as `[1;0]`. This is the guess for the state value at initial time  $k$ , using knowledge of system outputs until time  $k-1$ ,  $\hat{x}[k|k-1]$ .

```
obj = unscentedKalmanFilter(@vdpStateFcn,@vdpMeasurementFcn,[1;0]);
```

Load the measured output data,  $y$ , from the oscillator. In this example, use simulated static data for illustration. The data is stored in the `vdp_data.mat` file.

```
load vdp_data.mat y
```

Specify the process noise and measurement noise covariances of the oscillator.

```
obj.ProcessNoise = 0.01;
obj.MeasurementNoise = 0.16;
```

Implement the unscented Kalman filter algorithm to estimate the states of the oscillator by using the `correct` and `predict` commands. You first correct  $\hat{x}[k|k-1]$  using measurements at time  $k$  to get  $\hat{x}[k|k]$ . Then, you predict the state value at next time step,  $\hat{x}[k+1|k]$ , using  $\hat{x}[k|k]$ , the state estimate at time step  $k$  that is estimated using measurements until time  $k$ .

To simulate real-time data measurements, use the measured data one time step at a time.

```
for k = 1:size(y)
    [CorrectedState,CorrectedStateCovariance] = correct(obj,y(k));
    [PredictedState,PredictedStateCovariance] = predict(obj);
end
```

When you use the `correct` command, `obj.State` and `obj.StateCovariance` are updated with the corrected state and state estimation error covariance values for time step  $k$ , `CorrectedState` and `CorrectedStateCovariance`. When you use the `predict` command, `obj.State` and `obj.StateCovariance` are updated with the predicted values for time step  $k+1$ , `PredictedState` and `PredictedStateCovariance`.

In this example, you used `correct` before `predict` because the initial state value was  $\hat{x}[k|k-1]$ , a guess for the state value at initial time  $k$  using system outputs until time  $k-1$ . If your initial state value is  $\hat{x}[k-1|k-1]$ , the value at previous time  $k-1$  using measurement until  $k-1$ , then use the `predict` command first. For more information about the order of using `predict` and `correct`, see “Using `predict` and `correct` Commands” on page 1-1240.

## Estimate States Online using Particle Filter

Load the van der Pol ODE data, and specify the sample time.

`vdpODEdata.mat` contains a simulation of the van der Pol ODE with nonlinearity parameter  $\mu=1$ , using `ode45`, with initial conditions  $[2;0]$ . The true state was extracted with sample time  $dt = 0.05$ .

```
addpath(fullfile(matlabroot, 'examples', 'ident', 'main')) % add example data
```

```
load ('vdpODEdata.mat', 'xTrue', 'dt')  
tSpan = 0:dt:5;
```

Get the measurements. For this example, a sensor measures the first state with a Gaussian noise with standard deviation  $0.04$ .

```
sqrtr = 0.04;  
yMeas = xTrue(:,1) + sqrtr*randn(numel(tSpan),1);
```

Create a particle filter, and set the state transition and measurement likelihood functions.

```
myPF = particleFilter(@vdpParticleFilterStateFcn,@vdpMeasurementLikelihoodFcn);
```

Initialize the particle filter at state  $[2; 0]$  with unit covariance, and use 1000 particles.

```
initialize(myPF,1000,[2;0],eye(2));
```

Pick the mean state estimation and systematic resampling methods.

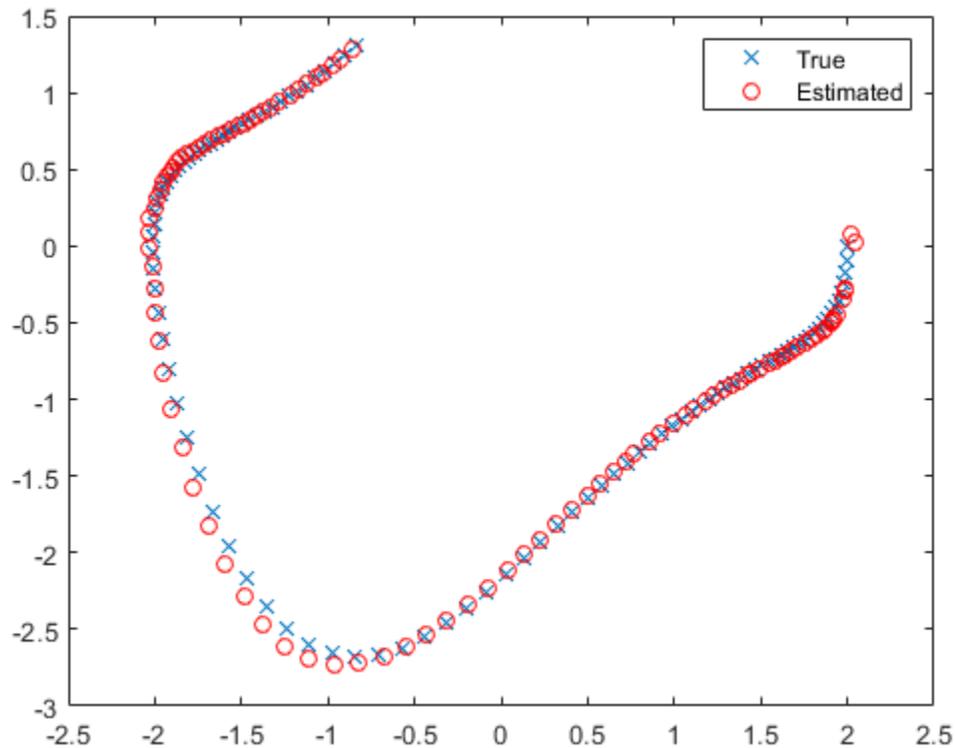
```
myPF.StateEstimationMethod = 'mean';  
myPF.ResamplingMethod = 'systematic';
```

Estimate the states using the `correct` and `predict` commands, and store the estimated states.

```
xEst = zeros(size(xTrue));  
for k=1:size(xTrue,1)  
    xEst(k,:) = correct(myPF,yMeas(k));  
    predict(myPF);  
end
```

Plot the results, and compare the estimated and true states.

```
figure(1)
plot(xTrue(:,1),xTrue(:,2),'x',xEst(:,1),xEst(:,2),'ro')
legend('True','Estimated')
```



```
rmpath(fullfile(matlabroot,'examples','ident','main')) % remove example data
```

### Specify State Transition and Measurement Functions with Additional Inputs

Consider a nonlinear system with input  $u$  whose state  $x$  and measurement  $y$  evolve according to the following state transition and measurement equations:

$$x[k] = \sqrt{x[k-1] + u[k-1]} + w[k-1]$$

$$y[k] = x[k] + 2 * u[k] + v[k]^2$$

The process noise  $w$  of the system is additive while the measurement noise  $v$  is nonadditive.

Create the state transition function and measurement function for the system. Specify the functions with an additional input  $u$ .

```
f = @(x,u)(sqrt(x+u));  
h = @(x,v,u)(x+2*u+v^2);
```

$f$  and  $h$  are function handles to the anonymous functions that store the state transition and measurement functions, respectively. In the measurement function, because the measurement noise is nonadditive,  $v$  is also specified as an input. Note that  $v$  is specified as an input before the additional input  $u$ .

Create an extended Kalman filter object for estimating the state of the nonlinear system using the specified functions. Specify the initial value of the state as 1, and the measurement noise as nonadditive.

```
obj = extendedKalmanFilter(f,h,1,'HasAdditiveMeasurementNoise',false);
```

Specify the measurement noise covariance.

```
obj.MeasurementNoise = 0.01;
```

You can now estimate the state of the system using the `predict` and `correct` commands. You pass the values of  $u$  to `predict` and `correct`, which in turn pass them to the state transition and measurement functions, respectively.

Correct the state estimate with measurement  $y[k]=0.8$  and input  $u[k]=0.2$  at time step  $k$ .

```
correct(obj,0.8,0.2)
```

Predict the state at next time step, given  $u[k]=0.2$ .

```
predict(obj,0.2)
```

## Input Arguments

### **obj** — Extended or unscented Kalman filter, or particle filter object

`extendedKalmanFilter` object | `unscentedKalmanFilter` object | `particleFilter` object

Extended or unscented Kalman filter, or particle filter object for online state estimation, created using one of the following commands:

- `extendedKalmanFilter` — Uses the extended Kalman filter algorithm.
- `unscentedKalmanFilter` — Uses the unscented Kalman filter algorithm.
- `particleFilter` — Uses the particle filter algorithm.

### **Us1, . . . Usn** — Additional input arguments to state transition function

input arguments of any type

Additional input arguments to state transition function, specified as input arguments of any type. The state transition function,  $f$ , is specified in the `StateTransitionFcn` property of the object. If the function requires input arguments in addition to the state and process noise values, you specify these inputs in the `predict` command syntax.

For example, suppose that your state transition function calculates the predicted state  $x$  at time step  $k$  using system inputs  $u(k-1)$  and time  $k-1$ , in addition to the state  $x(k-1)$ :

$$x(k) = f(x(k-1), u(k-1), k-1)$$

Then when you perform online state estimation at time step  $k$ , specify these additional inputs in the `predict` command syntax:

```
[PredictedState, PredictedStateCovariance] = predict(obj, u(k-1), k-1);
```

## Output Arguments

### **PredictedState** — Predicted state estimate

vector

Predicted state estimate, returned as a vector of size  $M$ , where  $M$  is the number of states of the system. If you specify the initial states of `obj` as a column vector then  $M$  is returned as a column vector, otherwise  $M$  is returned as a row vector.

For information about how to specify the initial states of the object, see the `extendedKalmanFilter`, `unscentedKalmanFilter` and `particleFilter` reference pages.

## **PredictedStateCovariance** — Predicted state estimation error covariance matrix

Predicted state estimation error covariance, returned as an  $M$ -by- $M$  matrix, where  $M$  is the number of states of the system.

## Definitions

### Using `predict` and `correct` Commands

After you have created an extended or unscented Kalman filter, or particle filter object, `obj`, to implement the estimation algorithms, use the `correct` and `predict` commands together.

At time step  $k$ , `correct` command returns the corrected value of states and state estimation error covariance using measured system outputs  $y[k]$  at the same time step. If your measurement function has additional input arguments  $U_m$ , you specify these as inputs to the `correct` command. The command passes these values to the measurement function.

```
[CorrectedState,CorrectedCovariance] = correct(obj,y,Um)
```

The `correct` command updates the `State` and `StateCovariance` properties of the object with the estimated values, `CorrectedState` and `CorrectedCovariance`.

The `predict` command returns the prediction of state and state estimation error covariance at the next time step. If your state transition function has additional input arguments  $U_s$ , you specify these as inputs to the `predict` command. The command passes these values to the state transition function.

```
[PredictedState,PredictedCovariance] = predict(obj,Us)
```

The `predict` command updates the `State` and `StateCovariance` properties of the object with the predicted values, `PredictedState` and `PredictedCovariance`.

If the current output measurement exists at a given time step, you can use `correct` and `predict`. If the measurement is missing, you can only use `predict`. For details about

how these commands implement the algorithms, see “Extended and Unscented Kalman Filter Algorithms for Online State Estimation”.

The order in which you implement the commands depends on the availability of measured data  $y$ ,  $U_s$ , and  $U_m$  for your system:

- **correct then predict** — Assume that at time step  $k$ , the value of `obj.State` is  $\hat{x}[k | k - 1]$ . This value is the state of the system at time  $k$ , estimated using measured outputs until time  $k - 1$ . You also have the measured output  $y[k]$  and inputs  $U_s[k]$  and  $U_m[k]$  at the same time step.

Then you first execute the `correct` command with measured system data  $y[k]$  and additional inputs  $U_m[k]$ . The command updates the value of `obj.State` to be  $\hat{x}[k | k]$ , the state estimate for time  $k$ , estimated using measured outputs up to time  $k$ . When you then execute the `predict` command with input  $U_s[k]$ , `obj.State` now stores

$\hat{x}[k + 1 | k]$ . The algorithm uses this state value as an input to the `correct` command in the next time step.

- **predict then correct** — Assume that at time step  $k$ , the value of `obj.State` is  $\hat{x}[k - 1 | k - 1]$ . You also have the measured output  $y[k]$  and input  $U_m[k]$  at the same time step but you have  $U_s[k - 1]$  from the previous time step.

Then you first execute the `predict` command with input  $U_s[k - 1]$ . The command updates the value of `obj.State` to  $\hat{x}[k | k - 1]$ . When you then execute the `correct` command with input arguments  $y[k]$  and  $U_m[k]$ , `obj.State` is updated with  $\hat{x}[k | k]$ . The algorithm uses this state value as an input to the `predict` command in the next time step.

Thus, while in both cases the state estimate for time  $k$ ,  $\hat{x}[k | k]$  is the same, if at time  $k$  you do not have access to the current state transition function inputs  $U_s[k]$ , and instead have  $U_s[k - 1]$ , then use `predict` first and then `correct`.

For an example of estimating states using the `predict` and `correct` commands, see “Estimate States Online Using Unscented Kalman Filter” on page 1-1234 or “Estimate States Online using Particle Filter” on page 1-1235.

## See Also

`clone` | `correct` | `extendedKalmanFilter` | `initialize` | `particleFilter` | `unscentedKalmanFilter`

## Topics

“Nonlinear State Estimation Using Unscented Kalman Filter and Particle Filter”

“Generate Code for Online State Estimation in MATLAB”

“What Is Online Estimation?”

“Extended and Unscented Kalman Filter Algorithms for Online State Estimation”

**Introduced in R2016b**

# predictOptions

Option set for predict

## Syntax

```
opt = predictOptions  
opt = predictOptions(Name,Value)
```

## Description

`opt = predictOptions` creates the default option set for `predict`. Use dot notation to modify this option set. Any options that you do not modify retain their default values.

`opt = predictOptions(Name,Value)` creates an option set with options specified by one or more `Name,Value` pair arguments.

## Examples

### Specify Output Offset for Predicting Model Response

Create a default option set for model prediction.

```
opt = predictOptions;
```

Specify the output offsets for a two-output model as 2 and 5, respectively.

```
opt.OutputOffset = [2;5];
```

The software subtracts the offset value `OutputOffset(i)` from the  $i$ th output signal before using the output to predict the model response. The software then adds back these offsets to the predicted response to give the final response.

## Specify Zero Initial Conditions for Model Prediction

Create an option set for `predict` using zero initial conditions.

```
opt = predictOptions('InitialCondition','z');
```

## Use Historical Data to Specify Initial Conditions for Model Prediction

Load a two-input, one-output dataset.

```
load iddata7 z7
```

Identify a fifth-order state-space model using the data.

```
sys = n4sid(z7,5);
```

Split the dataset into two parts.

```
zA = z7(1:15);  
zB = z7(16:end);
```

Suppose that you want to compute the 10-step-ahead prediction of the response of the identified system for data `zB`. For initial conditions, use the signal values in `zA` as the historical record. That is, the input and output values for the time immediately preceding data in `zB`.

```
I0 = struct('Input',zA.InputData,'Output',zA.OutputData);  
opt = predictOptions('InitialCondition',I0);
```

Generate the 10-step-ahead prediction for data `zB` using the specified initial conditions.

```
[yp,x0,Predictor] = predict(sys,zB,10,opt);
```

`yp` is the predicted model response, `x0` are the initial states corresponding to the predictor model `Predictor`. You can simulate `Predictor` using `x0` as initial conditions to reproduce `yp.OutputData`.

To understand how the past data is mapped to the initial states of the model, see “Understand Use of Historical Data for Model Prediction” on page 1-1225.

# Input Arguments

## Name-Value Pair Arguments

Specify optional comma-separated pairs of `Name`, `Value` arguments. `Name` is the argument name and `Value` is the corresponding value. `Name` must appear inside quotes. You can specify several name and value pair arguments in any order as `Name1, Value1, ..., NameN, ValueN`.

Example: `predictOptions('InitialCondition','z')` specifies zero initial conditions for the measured input-output data.

### InitialCondition — Handling of initial conditions

'e' (default) | 'z' | 'd' | column vector | matrix | structure | idpar object x00bj

Handling of initial conditions, specified as the comma-separated pair consisting of 'InitialCondition' and one of the following values:

- 'z' — Zero initial conditions.
- 'e' — Estimate initial conditions such that the prediction error for observed output is minimized.

For nonlinear grey-box models, only those initial states `i` that are designated as free in the model (`sys.InitialStates(i).Fixed = false`) are estimated. To estimate all the states of the model, first specify all the `Nx` states of the `idnlgrey` model `sys` as free.

```
for i = 1:Nx
    sys.InitialStates(i).Fixed = false;
end
```

Similarly, to fix all the initial states to values specified in `sys.InitialStates`, first specify all the states as fixed in the `sys.InitialStates` property of the nonlinear grey-box model.

- 'd' — Similar to 'e', but absorbs nonzero delays into the model coefficients. The delays are first converted to explicit model states, and the initial values of those states are also estimated and returned.

Use this option for linear models only.

- Vector or Matrix — Initial guess for state values, specified as a numerical column vector of length equal to the number of states. For multi-experiment data, specify a

matrix with  $N_e$  columns, where  $N_e$  is the number of experiments. Otherwise, use a column vector to specify the same initial conditions for all experiments. Use this option for state-space (`idss` and `idgrey`) and nonlinear models (`idnlarx`, `idnlhw`, and `idnlgrey`) only.

- Structure with the following fields, which contain the historical input and output values for a time interval immediately before the start time of the data used in the prediction:

Field	Description
Input	Input history, specified as a matrix with $N_u$ columns, where $N_u$ is the number of input channels. For time series models, use <code>[]</code> . The number of rows must be greater than or equal to the model order.
Output	Output history, specified as a matrix with $N_y$ columns, where $N_y$ is the number of output channels. The number of rows must be greater than or equal to the model order.

For an example, see “Use Historical Data to Specify Initial Conditions for Model Prediction” on page 1-1244.

For multi-experiment data, configure the initial conditions separately for each experiment by specifying `InitialCondition` as a structure array with  $N_e$  elements. To specify the same initial conditions for all experiments, use a single structure.

The software uses `data2state` to map the historical data to states. If your model is not `idss`, `idgrey`, `idnlgrey`, or `idnlarx`, the software first converts the model to its state-space representation and then maps the data to states. If conversion of your model to `idss` is not possible, the estimated states are returned empty.

- `x0obj` — Specification object created using `idpar`. Use this object for discrete-time state-space (`idss` and `idgrey`) and nonlinear grey-box (`idnlgrey`) models only. Use `x0obj` to impose constraints on the initial states by fixing their value or specifying minimum or maximum bounds.

### **InputOffset — Input signal offset**

`[]` (default) | column vector | matrix

Input signal offset for time-domain data, specified as the comma-separated pair consisting of `'InputOffset'` and one of the following values:

- `[]` — No input offsets.

- A column vector of length  $Nu$ , where  $Nu$  is the number of inputs. The software subtracts the offset value `InputOffset(i)` from the  $i$ th input signal before using the input to predict the model response.
- $Nu$ -by- $Ne$  matrix — For multi-experiment data, specify `InputOffset` as an  $Nu$ -by- $Ne$  matrix, where  $Ne$  is the number of experiments. The software subtracts the offset value `InputOffset(i, j)` from the  $i$ th input signal of the  $j$ th experiment before prediction.

If you specify a column vector of length  $Nu$ , then the offset value `InputOffset(i)` is subtracted from the  $i$ th input signal of all the experiments.

### **OutputOffset — Output signal offset**

`[]` (default) | column vector | matrix

Output signal offset for time-domain data, specified as the comma-separated pair consisting of 'OutputOffset' and one of the following values:

- `[]` — No output offsets.
- A column vector of length  $Ny$ , where  $Ny$  is the number of outputs. The software subtracts the offset value `OutputOffset(i)` from the  $i$ th output signal before using the output to predict the model response. After prediction, the software adds the offsets to the predicted response to give the final predicted response.
- $Ny$ -by- $Ne$  matrix — For multi-experiment data, specify `OutputOffset` as an  $Ny$ -by- $Ne$  matrix, where  $Ne$  is the number of experiments. The software subtracts the offset value `OutputOffset(i, j)` from the  $i$ th output signal of the  $j$ th experiment before prediction.

If you specify a column vector of length  $Ny$ , then the offset value `OutputOffset(i)` is subtracted from the  $i$ th output signal of all the experiments.

After prediction, the software adds the removed offsets to the predicted response to give the final predicted response.

### **OutputWeight — Weight of output for initial condition estimation**

`[]` (default) | 'noise' | matrix

Weight of output for initial condition estimation, specified as the comma-separated pair consisting of 'OutputWeight' and one of the following values:

- `[]` — No weighting is used by the software for initial condition estimation. This option is the same as using `eye(Ny)` for the output weight, where  $N_y$  is the number of outputs.
- `'noise'` — The software uses the inverse of the `NoiseVariance` property of the model as the weight.
- A positive, semidefinite matrix of dimension  $N_y$ -by- $N_y$ , where  $N_y$  is the number of outputs.

`OutputWeight` is relevant only for multi-output models.

## Output Arguments

**opt** — Option set for `predict`

`predictOptions` option set

Option set for `predict`, returned as a `predictOptions` option set.

## See Also

`absorbDelay` | `idpar` | `predict`

**Introduced in R2012a**

# present

Display model information, including estimated uncertainty

## Syntax

```
present(m)
```

## Description

`present(m)` displays the linear or nonlinear identified model `m` and the following information:

- Estimated one standard deviation of the parameters, which gives 68.27% confidence region
- Termination conditions for iterative estimation algorithms
- Status of the model — whether the model was constructed or estimated
- Fit to estimation data
- Akaike's Final Prediction Error (FPE) criterion
- Mean-square error (MSE)

## Examples

### Display Information About Identified Model

Estimate a transfer function model.

```
load iddata1 z1;  
np = 2;  
sys = tfest(z1,np);
```

Display model information.

```
present(sys)
```

sys =

```
From input "u1" to output "y1":  
    2.455 (+/- 1.101) s + 177 (+/- 10.73)  
-----  
s^2 + 3.163 (+/- 0.2522) s + 23.16 (+/- 1.115)
```

Continuous-time identified transfer function.

Parameterization:

Number of poles: 2    Number of zeros: 1

Number of free coefficients: 4

Use "tfdata", "getpvec", "getcov" for parameters and their uncertainties.

Status:

Termination condition: Near (local) minimum, (norm(g) < tol)..

Number of iterations: 1, Number of function evaluations: 3

Estimated using TFEST on time domain data "z1".

Fit to estimation data: 70.77%

FPE: 1.725, MSE: 1.658

More information in model's "Report" property.

## See Also

[frdata](#) | [getcov](#) | [getpvec](#) | [idssdata](#) | [polydata](#) | [ssdata](#) | [tfdata](#) | [zpkdata](#)

## Topics

["Estimation Report"](#)

["Loss Function and Model Quality Metrics"](#)

**Introduced before R2006a**

# procest

Estimate process model using time or frequency data

## Syntax

```
sys = procest(data,type)
sys = procest(data,type,'InputDelay',InputDelay)

sys = procest(data,init_sys)

sys = procest( ____,opt)

[sys,offset] = procest( ____ )
```

## Description

`sys = procest(data,type)` estimates a process model, `sys`, using time or frequency-domain data, `data`. `type` defines the structure of `sys`.

`sys = procest(data,type,'InputDelay',InputDelay)` specifies the input delay `InputDelay`.

`sys = procest(data,init_sys)` uses the process model `init_sys` to configure the initial parameterization.

`sys = procest( ____,opt)` specifies additional model estimation options. Use `opt` with any of the previous syntaxes.

`[sys,offset] = procest( ____ )` returns the estimated value of the offset in input signal. Input offset is automatically estimated when the model contains an integrator, or when you set the `InputOffset` estimation option to 'estimate' using `procestOptions`. Use `offset` with any of the previous syntaxes.

## Input Arguments

### **data — Estimation data**

`iddata` | `idfrd` | `frd`

Estimation data, specified as an `iddata` object containing the input and output signal values, for time-domain estimation. For frequency-domain estimation, `data` can be one of the following:

- Recorded frequency response data (`frd` or `idfrd`)
- `iddata` object with its properties specified as follows:
  - `InputData` — Fourier transform of the input signal
  - `OutputData` — Fourier transform of the output signal
  - `Domain` — 'Frequency'

`data` must have at least one input and one output.

Time-series models, which are models that contain no measured inputs, cannot be estimated using `procest`. Use `ar`, `arx`, or `armax` for time-series models instead.

### **type — Process model structure**

character vector | string | cell array of character vectors | string array

Process model structure, specified for SISO models as a character vector or string representing an acronym for the model structure, such as 'P1D' or 'P2DZ'. The acronym is made up of:

- **P** — All 'Type' acronyms start with this letter.
- **0, 1, 2, or 3** — Number of time constants (poles) to be modeled. Possible integrations (poles in the origin) are not included in this number.
- **I** — Integration is enforced (self-regulating process).
- **D** — Time delay (dead time).
- **Z** — Extra numerator term, a zero.
- **U** — Underdamped modes (complex-valued poles) permitted. If **U** is not included in `type`, all poles must be real. The number of poles must be 2 or 3.

For MIMO models, use an  $N_y$ -by- $N_u$  cell array of character vectors or string array, with one entry for each input-output pair. Here  $N_y$  is the number of inputs and  $N_u$  is the number of outputs.

For information regarding how `type` affects the structure of a process model, see `idproc`.

### **InputDelay — Input delays**

0 for all input channels (default) | numeric vector

Input delays, specified as a numeric vector specifying a time delay for each input channel. Specify input delays in the time unit stored in the `TimeUnit` property.

For a system with  $N_u$  inputs, set `InputDelay` to an  $N_u$ -by-1 vector. Each entry of this vector is a numerical value that represents the input delay for the corresponding input channel. You can also set `InputDelay` to a scalar value to apply the same delay to all channels.

### **init\_sys — System for configuring initial parametrization**

`idproc` object

System for configuring initial parametrization of `sys`, specified as an `idproc` object. You obtain `init_sys` by either performing an estimation using measured data or by direct construction using `idproc`. The software uses the parameters and constraints defined in `init_sys` as the initial guess for estimating `sys`.

Use the `Structure` property of `init_sys` to configure initial guesses and constraints for  $K_p$ ,  $T_{p1}$ ,  $T_{p2}$ ,  $T_{p3}$ ,  $T_w$ ,  $Zeta$ ,  $T_d$ , and  $T_z$ . For example:

- To specify an initial guess for the  $T_{p1}$  parameter of `init_sys`, set `init_sys.Structure.Tp1.Value` as the initial guess.
- To specify constraints for the  $T_{p2}$  parameter of `init_sys`:
  - Set `init_sys.Structure.Tp2.Minimum` to the minimum  $T_{p2}$  value.
  - Set `init_sys.Structure.Tp2.Maximum` to the maximum  $T_{p2}$  value.
  - Set `init_sys.Structure.Tp2.Free` to indicate if  $T_{p2}$  is a free parameter for estimation.

If `opt` is not specified, and `init_sys` was obtained by estimation, then the estimation options from `init_sys.Report.OptionsUsed` are used.

**opt — Estimation options**`procestOptions` option set

Estimation options, specified as an `procestOptions` option set. The estimation options include:

- Estimation objective
- Handling on initial conditions and disturbance component
- Numerical search method to be used in estimation

## Output Arguments

**sys — Identified process model**`idproc` model

Identified process model, returned as an `idproc` model of a structure defined by `type`.

Information about the estimation results and options used is stored in the model's `Report` property. `Report` has the following fields:

Report Field	Description
Status	Summary of the model status, which indicates whether the model was created by construction or obtained by estimation.
Method	Estimation command used.
InitialCondition	Handling of initial conditions during model estimation, returned as one of the following values: <ul style="list-style-type: none"><li>• 'zero' — The initial conditions were set to zero.</li><li>• 'estimate' — The initial conditions were treated as independent estimation parameters.</li><li>• 'backcast' — The initial conditions were estimated using the best least squares fit.</li></ul> <p>This field is especially useful to view how the initial conditions were handled when the <code>InitialCondition</code> option in the estimation option set is 'auto'.</p>

Report Field	Description	
Fit	Quantitative assessment of the estimation, returned as a structure. See “Loss Function and Model Quality Metrics” for more information on these quality metrics. The structure has the following fields:	
	<b>Field</b>	<b>Description</b>
	FitPercent	Normalized root mean squared error (NRMSE) measure of how well the response of the model fits the estimation data, expressed as a percentage.
	LossFunction	Value of the loss function when the estimation completes.
	MSE	Mean squared error (MSE) measure of how well the response of the model fits the estimation data.
	FPE	Final prediction error for the model.
	AIC	Raw Akaike Information Criteria (AIC) measure of model quality.
	AICc	Small sample-size corrected AIC.
	nAIC	Normalized AIC.
BIC	Bayesian Information Criteria (BIC).	
Parameters	Estimated values of model parameters.	
OptionsUsed	Option set used for estimation. If no custom options were configured, this is a set of default options. See <code>procestOptions</code> for more information.	
RandState	State of the random number stream at the start of estimation. Empty, [], if randomization was not used during estimation. For more information, see <code>rng</code> in the MATLAB documentation.	

Report Field	Description																
DataUsed	Attributes of the data used for estimation. Structure with the following fields:																
	<table border="1"> <thead> <tr> <th data-bbox="543 395 649 430">Field</th> <th data-bbox="649 395 1338 430">Description</th> </tr> </thead> <tbody> <tr> <td data-bbox="543 430 649 475">Name</td> <td data-bbox="649 430 1338 475">Name of the data set.</td> </tr> <tr> <td data-bbox="543 475 649 552">Type</td> <td data-bbox="649 475 1338 552">Data type. For <code>idnlarx</code> models, this is set to 'Time domain data'.</td> </tr> <tr> <td data-bbox="543 552 649 597">Length</td> <td data-bbox="649 552 1338 597">Number of data samples.</td> </tr> <tr> <td data-bbox="543 597 649 642">Ts</td> <td data-bbox="649 597 1338 642">Sample time. This is equivalent to <code>Data.Ts</code>.</td> </tr> <tr> <td data-bbox="543 642 649 1079">Intersample</td> <td data-bbox="649 642 1338 1079">                     Input intersample behavior. One of the following values:                     <ul style="list-style-type: none"> <li>• 'zoh' — Zero-order hold maintains a piecewise-constant input signal between samples.</li> <li>• 'foh' — First-order hold maintains a piecewise-linear input signal between samples.</li> <li>• 'bl' — Band-limited behavior specifies that the continuous-time input signal has zero power above the Nyquist frequency.</li> </ul>                     The value of <code>Intersample</code> has no effect on estimation results for discrete-time models.                 </td> </tr> <tr> <td data-bbox="543 1079 649 1156">InputOffset</td> <td data-bbox="649 1079 1338 1156">Empty, [], for nonlinear estimation methods.</td> </tr> <tr> <td data-bbox="543 1156 649 1234">OutputOffset</td> <td data-bbox="649 1156 1338 1234">Empty, [], for nonlinear estimation methods.</td> </tr> </tbody> </table>	Field	Description	Name	Name of the data set.	Type	Data type. For <code>idnlarx</code> models, this is set to 'Time domain data'.	Length	Number of data samples.	Ts	Sample time. This is equivalent to <code>Data.Ts</code> .	Intersample	Input intersample behavior. One of the following values: <ul style="list-style-type: none"> <li>• 'zoh' — Zero-order hold maintains a piecewise-constant input signal between samples.</li> <li>• 'foh' — First-order hold maintains a piecewise-linear input signal between samples.</li> <li>• 'bl' — Band-limited behavior specifies that the continuous-time input signal has zero power above the Nyquist frequency.</li> </ul> The value of <code>Intersample</code> has no effect on estimation results for discrete-time models.	InputOffset	Empty, [], for nonlinear estimation methods.	OutputOffset	Empty, [], for nonlinear estimation methods.
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	InputOffset	Empty, [], for nonlinear estimation methods.															
OutputOffset	Empty, [], for nonlinear estimation methods.																

Report Field	Description	
Termination	Termination conditions for the iterative search used for prediction error minimization. Structure with the following fields:	
	<b>Field</b>	<b>Description</b>
	WhySto p	Reason for terminating the numerical search.
	Iterat ions	Number of search iterations performed by the estimation algorithm.
	FirstO rderOp timali ty	$\infty$ -norm of the gradient search vector when the search algorithm terminates.
	FcnCou nt	Number of times the objective function was called.
	Update Norm	Norm of the gradient search vector in the last iteration. Omitted when the search method is 'lsqnonlin' or 'fmincon'.
	LastIm provem ent	Criterion improvement in the last iteration, expressed as a percentage. Omitted when the search method is 'lsqnonlin' or 'fmincon'.
	Algori thm	Algorithm used by 'lsqnonlin' or 'fmincon' search method. Omitted when other search methods are used.
For estimation methods that do not require numerical search optimization, the Termination field is omitted.		

For more information on using Report, see “Estimation Report”.

### **offset** — Estimated value of input offset

vector

Estimated value of input offset, returned as a vector. When data has multiple experiments, offset is a matrix where each column corresponds to an experiment.

## Examples

### Estimate a First Order Plus Dead Time Model

Obtain the measured input-output data.

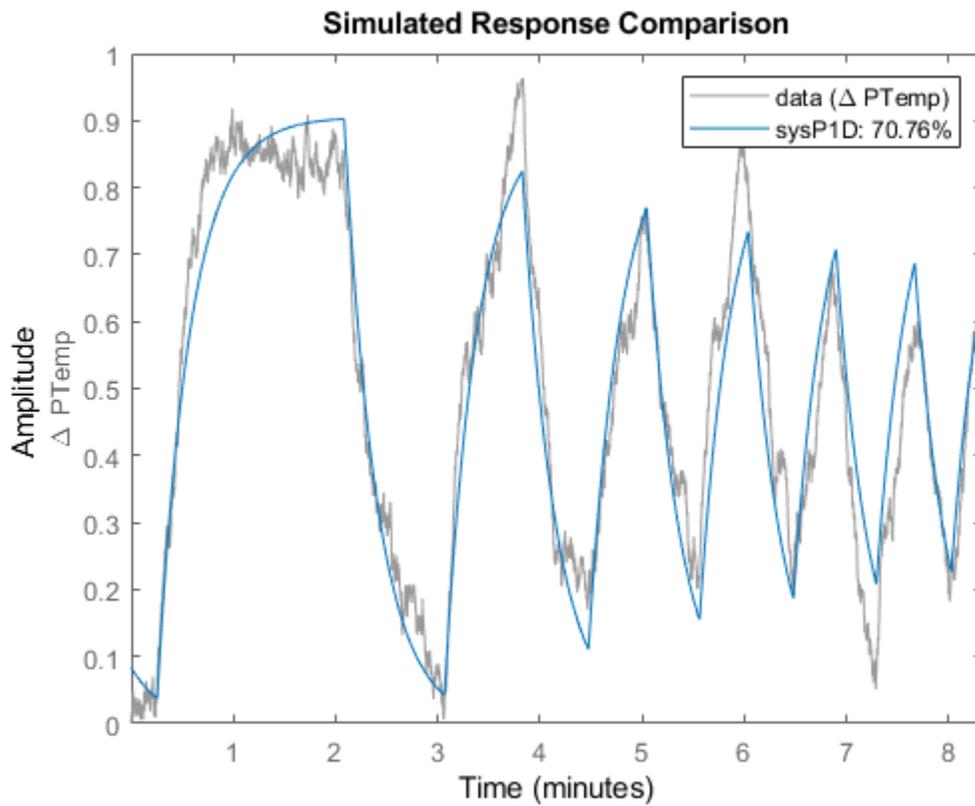
```
load iddemo_heatexchanger_data;  
data = iddata(pt,ct,Ts);  
data.InputName = '\Delta CTemp';  
data.InputUnit = 'C';  
data.OutputName = '\Delta PTemp';  
data.OutputUnit = 'C';  
data.TimeUnit = 'minutes';
```

Estimate a first-order plus dead time process model.

```
type = 'PID';  
sysPID = procest(data,type);
```

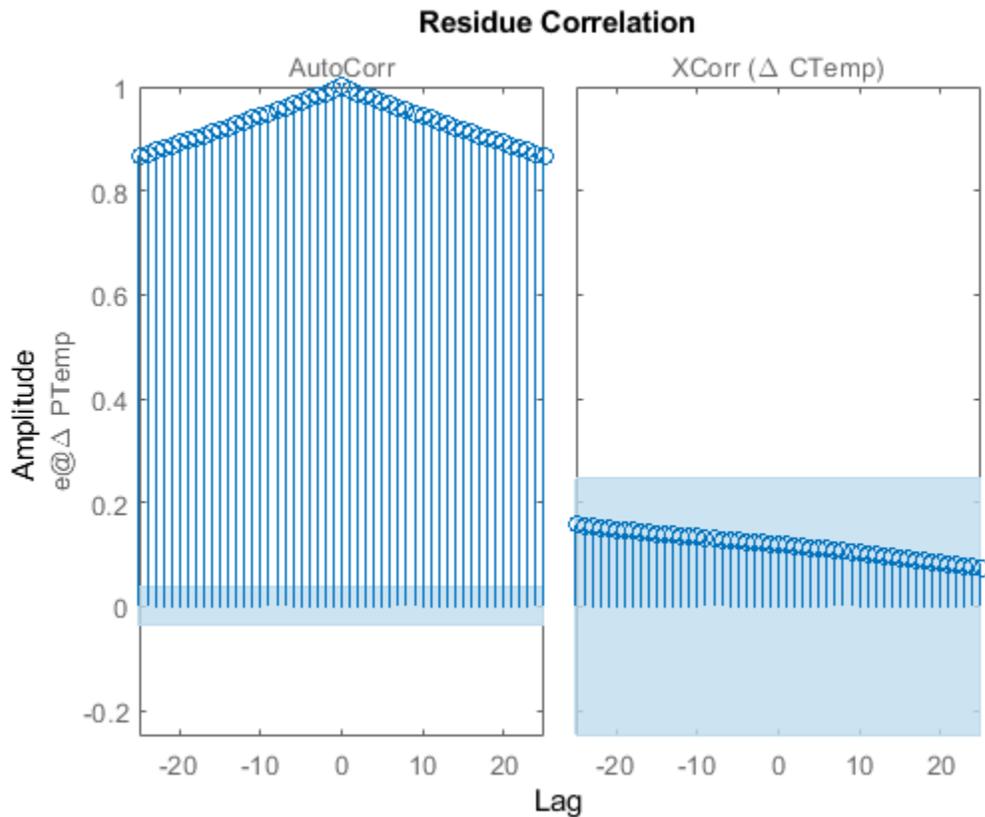
Compare the model with the data.

```
compare(data,sysPID)
```



Plot the model residuals.

```
figure  
resid(data,sysP1D);
```

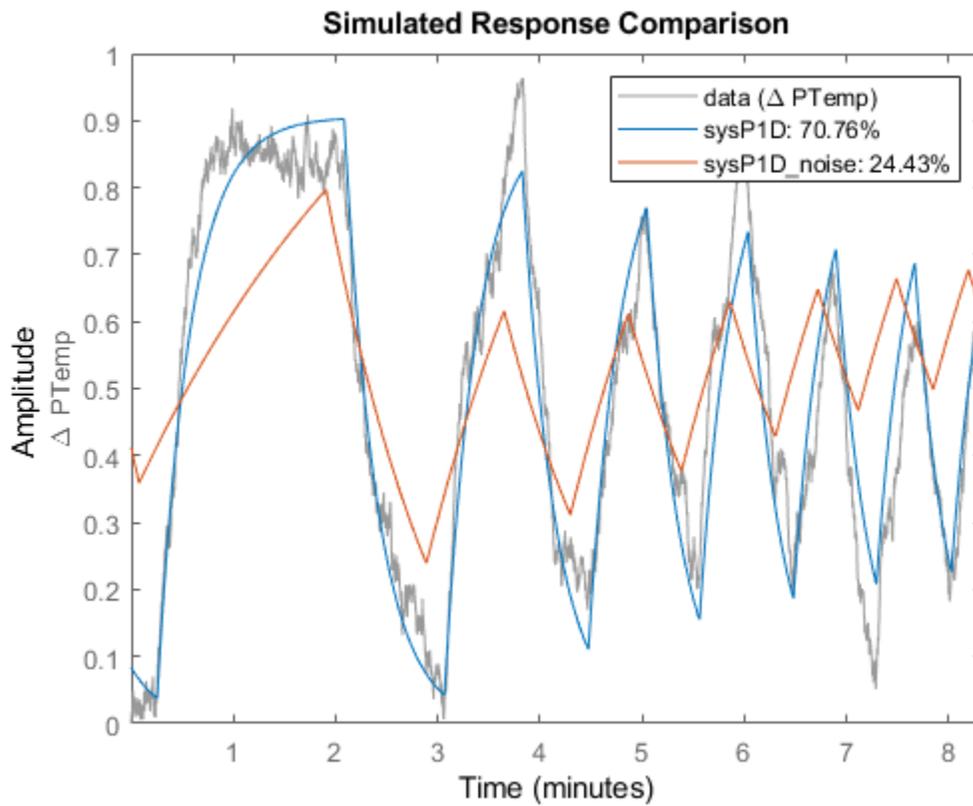


The figure shows that the residuals are correlated. To account for that, add a first order ARMA disturbance component to the process model.

```
opt = procestOptions('DisturbanceModel','ARMA1');
sysP1D_noise = procest(data,'p1d',opt);
```

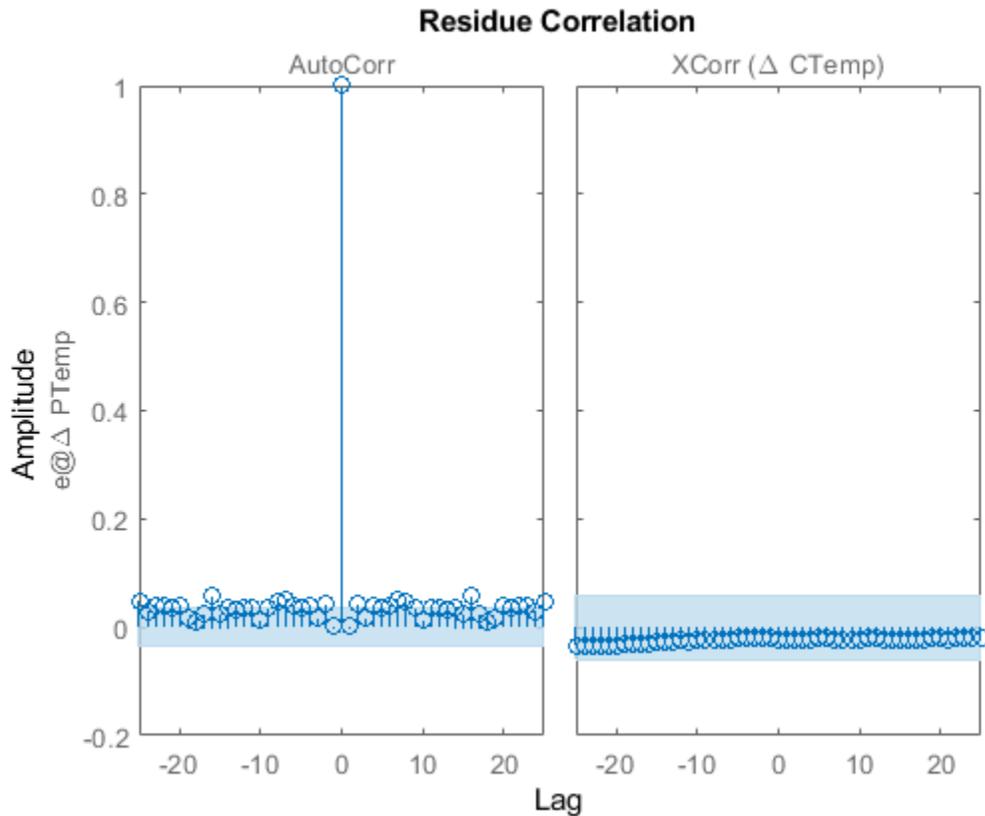
Compare the models.

```
compare(data,sysP1D,sysP1D_noise)
```



Plot the model residuals.

```
figure  
resid(data,sysP1D_noise);
```



The residues of `sysPID_noise` are uncorrelated.

### Estimate Over-parameterized Process Model Using Regularization

Use regularization to estimate parameters of an over-parameterized process model.

Assume that gain is known with a higher degree of confidence than other model parameters.

Load data.

```
load iddata1 z1;
```

Estimate an unregularized process model.

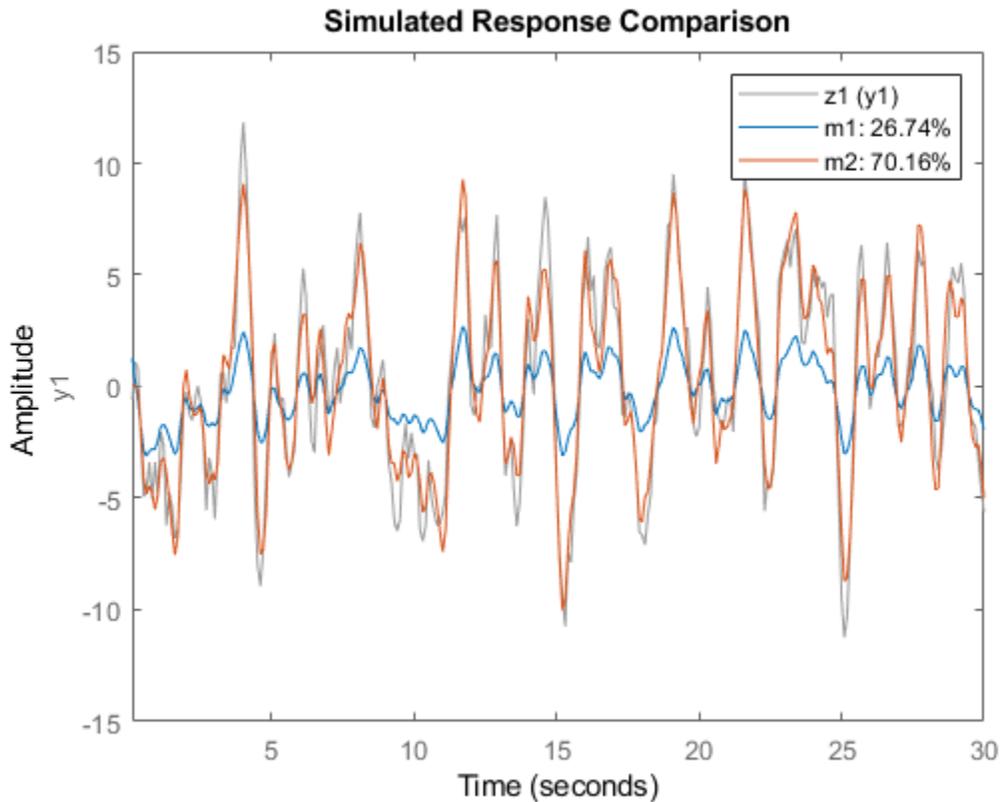
```
m = idproc('P3UZ','K',7.5,'Tw',0.25,'Zeta',0.3,'Tp3',20,'Tz',0.02);  
m1 = procest(z1,m);
```

Estimate a regularized process model.

```
opt = procestOptions;  
opt.Regularization.Nominal = 'model';  
opt.Regularization.R = [100;1;1;1;1];  
opt.Regularization.Lambda = 0.1;  
m2 = procest(z1,m,opt);
```

Compare the model outputs with data.

```
compare(z1,m1,m2);
```



Regularization helps steer the estimation process towards the correct parameter values.

### Specify Parameter Initial Values for Estimated Process Model

Estimate a process model after specifying initial guesses for parameter values and bounding them.

Obtain input/output data.

```
data = idfrd(idtf([10 2],[1 1.3 1.2], 'iod', 0.45), logspace(-2, 2, 256));
```

Specify the parameters of the estimation initialization model.

```
type = 'P2UZD';
init_sys = idproc(type);

init_sys.Structure.Kp.Value = 1;
init_sys.Structure.Tw.Value = 2;
init_sys.Structure.Zeta.Value = 0.1;
init_sys.Structure.Td.Value = 0;
init_sys.Structure.Tz.Value = 1;
init_sys.Structure.Kp.Minimum = 0.1;
init_sys.Structure.Kp.Maximum = 10;
init_sys.Structure.Td.Maximum = 1;
init_sys.Structure.Tz.Maximum = 10;
```

Specify the estimation options.

```
opt = procestOptions('Display','full','InitialCondition','Zero');
opt.SearchMethod = 'lm';
opt.SearchOptions.MaxIterations = 100;
```

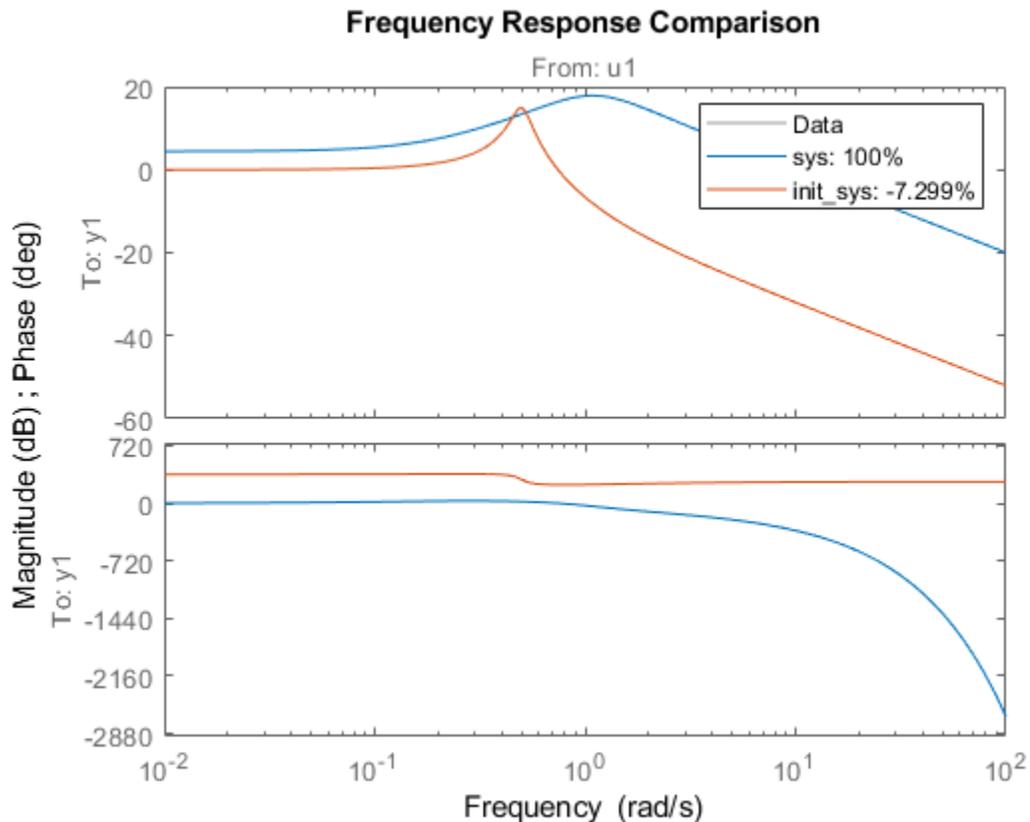
Estimate the process model.

```
sys = procest(data,init_sys,opt);
```

Since the 'Display' option is specified as 'full', the estimation progress is displayed in a separate **Plant Identification Progress** window.

Compare the data to the estimated model.

```
compare(data,sys,init_sys);
```



### Detect Overparameterization of Estimated Model

Obtain input/output data.

```
load iddata1 z1
load iddata2 z2
data = [z1 z2(1:300)];
```

`data` is a data set with 2 inputs and 2 outputs. The first input affects only the first output. Similarly, the second input affects only the second output.

In the estimated process model, the cross terms, modeling the effect of the first input on the second output and vice versa, should be negligible. If higher orders are assigned to those dynamics, their estimations show a high level of uncertainty.

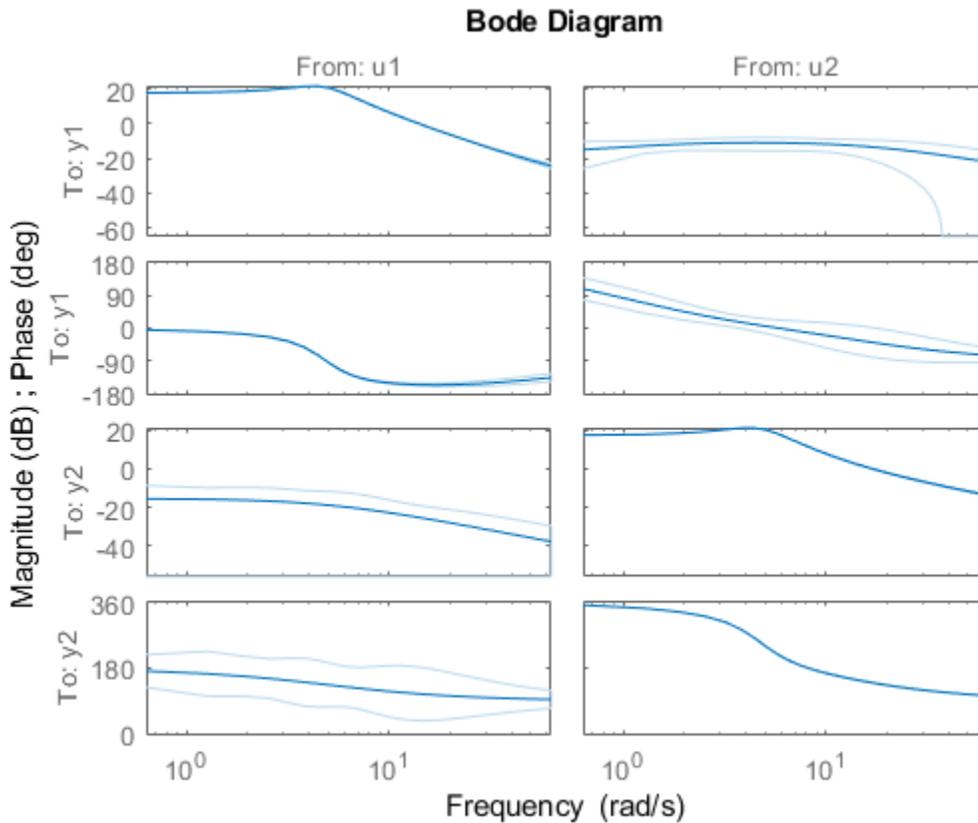
Estimate the process model.

```
type = 'P2UZ';  
sys = procest(data,type);
```

The `type` variable denotes a model with complex-conjugate pair of poles, a zero, and a delay.

To evaluate the uncertainties, plot the frequency response.

```
w = linspace(0,20*pi,100);  
h = bodeplot(sys,w);  
showConfidence(h);
```



### Return Input Offsets Estimated During Process Model Estimation

```
load iddata1
[sys,offset] = procest(z1,'PIDI');
offset

offset = 0.0412
```

## Extended Capabilities

### Automatic Parallel Support

Accelerate code by automatically running computation in parallel using Parallel Computing Toolbox™.

Parallel computing support is available for estimation using the `lsqnonlin` search method (requires Optimization Toolbox). To enable parallel computing, use `procestOptions`, set `SearchMethod` to `'lsqnonlin'`, and set `SearchOptions.Advanced.UseParallel` to `true`.

For example:

```
opt = procestOptions;  
opt.SearchMethod = 'lsqnonlin';  
opt.SearchOptions.Advanced.UseParallel = true;
```

### See Also

[ar](#) | [arx](#) | [bj](#) | [idproc](#) | [oe](#) | [polyest](#) | [procestOptions](#) | [srest](#) | [tfest](#)

### Topics

“What Is a Process Model?”

“Regularized Estimates of Model Parameters”

**Introduced in R2012a**

## procestOptions

Options set for procest

### Syntax

```
opt = procestOptions  
opt = procestOptions(Name,Value)
```

### Description

`opt = procestOptions` creates the default options set for `procest`.

`opt = procestOptions(Name,Value)` creates an option set with the options specified by one or more `Name,Value` pair arguments.

### Input Arguments

#### Name-Value Pair Arguments

Specify optional comma-separated pairs of `Name,Value` arguments. `Name` is the argument name and `Value` is the corresponding value. `Name` must appear inside quotes. You can specify several name and value pair arguments in any order as `Name1,Value1,...,NameN,ValueN`.

#### InitialCondition — Handling of initial conditions

'auto' (default) | 'zero' | 'estimate' | 'backcast'

Handling of initial conditions during estimation, specified as one of the following values:

- 'zero' — The initial condition is set to zero.
- 'estimate' — The initial condition is treated as an independent estimation parameter.
- 'backcast' — The initial condition is estimated using the best least squares fit.

- 'auto' — The software chooses the method to handle initial condition based on the estimation data.

### **DisturbanceModel — Handling of additive noise**

'estimate' (default) | 'none' | 'ARMA1' | 'ARMA2' | 'fixed'

Handling of additive noise ( $H$ ) during estimation for the model

$$y = G(s)u + H(s)e$$

$e$  is white noise,  $u$  is the input and  $y$  is the output.

$H(s)$  is stored in the NoiseTF property of the numerator and denominator of idproc models.

DisturbanceModel is specified as one of the following values:

- 'none' —  $H$  is fixed to one.
- 'estimate' —  $H$  is treated as an estimation parameter. The software uses the value of the NoiseTF property as the initial guess.
- 'ARMA1' — The software estimates  $H$  as a first-order ARMA model

$$\frac{1 + cs}{1 + ds}$$

- 'ARMA2' — The software estimates  $H$  as a second-order ARMA model

$$\frac{1 + c_1s + c_2s^2}{1 + d_1s + d_2s^2}$$

- 'fixed' — The software fixes the value of the NoiseTF property of the idproc model as the value of  $H$ .

---

**Note** A noise model cannot be estimated using frequency domain data.

---

### **Focus — Error to be minimized**

'prediction' (default) | 'simulation'

Error to be minimized in the loss function during estimation, specified as the comma-separated pair consisting of 'Focus' and one of the following values:

- 'prediction' — The one-step ahead prediction error between measured and predicted outputs is minimized during estimation. As a result, the estimation focuses on producing a good predictor model.
- 'simulation' — The simulation error between measured and simulated outputs is minimized during estimation. As a result, the estimation focuses on making a good fit for simulation of model response with the current inputs.

The Focus option can be interpreted as a weighting filter in the loss function. For more information, see “Loss Function and Model Quality Metrics”.

### **WeightingFilter — Weighting prefilter**

[ ] (default) | vector | matrix | cell array | linear system

Weighting prefilter applied to the loss function to be minimized during estimation. To understand the effect of `WeightingFilter` on the loss function, see “Loss Function and Model Quality Metrics”.

Specify `WeightingFilter` as one of the following values:

- [ ] — No weighting prefilter is used.
- Passbands — Specify a row vector or matrix containing frequency values that define desired passbands. You select a frequency band where the fit between estimated model and estimation data is optimized. For example, [w<sub>l</sub>, w<sub>h</sub>] where w<sub>l</sub> and w<sub>h</sub> represent lower and upper limits of a passband. For a matrix with several rows defining frequency passbands, [w<sub>1l</sub>, w<sub>1h</sub>; w<sub>2l</sub>, w<sub>2h</sub>; w<sub>3l</sub>, w<sub>3h</sub>; . . . ], the estimation algorithm uses the union of the frequency ranges to define the estimation passband.

Passbands are expressed in rad/TimeUnit for time-domain data and in FrequencyUnit for frequency-domain data, where TimeUnit and FrequencyUnit are the time and frequency units of the estimation data.

- SISO filter — Specify a single-input-single-output (SISO) linear filter in one of the following ways:
  - A SISO LTI model
  - {A, B, C, D} format, which specifies the state-space matrices of a filter with the same sample time as estimation data.

- `{numerator,denominator}` format, which specifies the numerator and denominator of the filter as a transfer function with same sample time as estimation data.

This option calculates the weighting function as a product of the filter and the input spectrum to estimate the transfer function.

- **Weighting vector** — Applicable for frequency-domain data only. Specify a column vector of weights. This vector must have the same length as the frequency vector of the data set, `Data.Frequency`. Each input and output response in the data is multiplied by the corresponding weight at that frequency.

### **EstimateCovariance — Control whether to generate parameter covariance data**

`true` (default) | `false`

Controls whether parameter covariance data is generated, specified as `true` or `false`.

If `EstimateCovariance` is `true`, then use `getcov` to fetch the covariance matrix from the estimated model.

### **Display — Specify whether to display the estimation progress**

`'off'` (default) | `'on'`

Specify whether to display the estimation progress, specified as one of the following values:

- `'on'` — Information on model structure and estimation results are displayed in a progress-viewer window.
- `'off'` — No progress or results information is displayed.

### **InputOffset — Removal of offset from time-domain input data**

`'auto'` (default) | `'estimate'` | `vector` | `matrix` | `object` | `[]`

Removal of offset from time-domain input data during estimation, specified as one of the following values:

- `'estimate'` — The software treats the input offsets as an estimation parameter.
- `'auto'` — The software chooses the method to handle input offsets based on the estimation data and the model structure. The estimation either assumes zero input offset or estimates the input offset.

For example, the software estimates the input offset for a model that contains an integrator.

- A column vector of length  $Nu$ , where  $Nu$  is the number of inputs.

Use `[]` to specify no offsets.

In case of multi-experiment data, specify `InputOffset` as a  $Nu$ -by- $Ne$  matrix.  $Nu$  is the number of inputs, and  $Ne$  is the number of experiments.

Each entry specified by `InputOffset` is subtracted from the corresponding input data.

- A parameter object, constructed using `param.Continuous`, that imposes constraints on how the software estimates the input offset.

For example, create a parameter object for a 2-input model estimation. Specify the first input offset as fixed to zero and the second input offset as an estimation parameter.

```
opt = procestOptions;  
u0 = param.Continuous('u0', [0;NaN]);  
u0.Free(1) = false;  
opt.Inputoffset = u0;
```

## **OutputOffset — Removal of offset from time-domain output data during estimation**

`[]` (default) | vector | matrix

Removal of offset from time-domain output data during estimation, specified as the comma-separated pair consisting of `'OutputOffset'` and one of the following:

- A column vector of length  $Ny$ , where  $Ny$  is the number of outputs.
- `[]` — Indicates no offset.
- $Ny$ -by- $Ne$  matrix — For multi-experiment data, specify `OutputOffset` as a  $Ny$ -by- $Ne$  matrix.  $Ny$  is the number of outputs, and  $Ne$  is the number of experiments.

Each entry specified by `OutputOffset` is subtracted from the corresponding output data.

## **OutputWeight — Weighting of prediction errors in multi-output estimations**

`[]` (default) | `'noise'` | positive semidefinite symmetric matrix

Weighting of prediction errors in multi-output estimations, specified as one of the following values:

- 'noise' — Minimize  $\det(E^*E / N)$ , where  $E$  represents the prediction error and  $N$  is the number of data samples. This choice is optimal in a statistical sense and leads to maximum likelihood estimates if nothing is known about the variance of the noise. It uses the inverse of the estimated noise variance as the weighting function.

---

**Note** OutputWeight must not be 'noise' if SearchMethod is 'lsqnonlin'.

---

- Positive semidefinite symmetric matrix ( $W$ ) — Minimize the trace of the weighted prediction error matrix  $\text{trace}(E^*E*W/N)$  where:
  - $E$  is the matrix of prediction errors, with one column for each output, and  $W$  is the positive semidefinite symmetric matrix of size equal to the number of outputs. Use  $W$  to specify the relative importance of outputs in multiple-output models, or the reliability of corresponding data.
  - $N$  is the number of data samples.
- [] — The software chooses between the 'noise' or using the identity matrix for  $W$ .

This option is relevant for only multi-output models.

### Regularization — Options for regularized estimation of model parameters

structure

Options for regularized estimation of model parameters. For more information on regularization, see “Regularized Estimates of Model Parameters”.

Regularization is a structure with the following fields:

- Lambda — Constant that determines the bias versus variance tradeoff.

Specify a positive scalar to add the regularization term to the estimation cost.

The default value of zero implies no regularization.

**Default:** 0

- R — Weighting matrix.

Specify a vector of nonnegative numbers or a square positive semi-definite matrix. The length must be equal to the number of free parameters of the model.

For black-box models, using the default value is recommended. For structured and grey-box models, you can also specify a vector of `np` positive numbers such that each entry denotes the confidence in the value of the associated parameter.

The default value of 1 implies a value of `eye(npfree)`, where `npfree` is the number of free parameters.

**Default:** 1

- **Nominal** — The nominal value towards which the free parameters are pulled during estimation.

The default value of zero implies that the parameter values are pulled towards zero. If you are refining a model, you can set the value to `'model'` to pull the parameters towards the parameter values of the initial model. The initial parameter values must be finite for this setting to work.

**Default:** 0

## **SearchMethod — Numerical search method used for iterative parameter estimation**

`'auto'` (default) | `'gn'` | `'gna'` | `'lm'` | `'grad'` | `'lsqnonlin'` | `'fmincon'`

Numerical search method used for iterative parameter estimation, specified as the comma-separated pair consisting of `'SearchMethod'` and one of the following:

- `'auto'` — A combination of the line search algorithms, `'gn'`, `'lm'`, `'gna'`, and `'grad'` methods is tried in sequence at each iteration. The first descent direction leading to a reduction in estimation cost is used.
- `'gn'` — Subspace Gauss-Newton least squares search. Singular values of the Jacobian matrix less than `GnPinvConstant*eps*max(size(J))*norm(J)` are discarded when computing the search direction.  $J$  is the Jacobian matrix. The Hessian matrix is approximated as  $J^T J$ . If there is no improvement in this direction, the function tries the gradient direction.
- `'gna'` — Adaptive subspace Gauss-Newton search. Eigenvalues less than `gamma*max(sv)` of the Hessian are ignored, where `sv` contains the singular values of the Hessian. The Gauss-Newton direction is computed in the remaining subspace. `gamma` has the initial value `InitialGnaTolerance` (see `Advanced in 'SearchOptions'` for more information). This value is increased by the factor `LMStep` each time the search fails to find a lower value of the criterion in fewer than five bisections. This value is decreased by the factor `2*LMStep` each time a search is successful without any bisections.

- 'lm' — Levenberg-Marquardt least squares search, where the next parameter value is  $-\text{pinv}(H+d*I)*\text{grad}$  from the previous one.  $H$  is the Hessian,  $I$  is the identity matrix, and  $\text{grad}$  is the gradient.  $d$  is a number that is increased until a lower value of the criterion is found.
- 'grad' — Steepest descent least squares search.
- 'lsqnonlin' — Trust-region-reflective algorithm of `lsqnonlin`. Requires Optimization Toolbox software.
- 'fmincon' — Constrained nonlinear solvers. You can use the sequential quadratic programming (SQP) and trust-region-reflective algorithms of the `fmincon` solver. If you have Optimization Toolbox software, you can also use the interior-point and active-set algorithms of the `fmincon` solver. Specify the algorithm in the `SearchOptions.Algorithm` option. The `fmincon` algorithms may result in improved estimation results in the following scenarios:
  - Constrained minimization problems when there are bounds imposed on the model parameters.
  - Model structures where the loss function is a nonlinear or non smooth function of the parameters.
  - Multi-output model estimation. A determinant loss function is minimized by default for multi-output model estimation. `fmincon` algorithms are able to minimize such loss functions directly. The other search methods such as 'lm' and 'gn' minimize the determinant loss function by alternately estimating the noise variance and reducing the loss value for a given noise variance value. Hence, the `fmincon` algorithms can offer better efficiency and accuracy for multi-output model estimations.

### **SearchOptions — Option set for the search algorithm**

search option set

Option set for the search algorithm, specified as the comma-separated pair consisting of 'SearchOptions' and a search option set with fields that depend on the value of SearchMethod.

**SearchOptions Structure When SearchMethod is Specified as 'gn', 'gna', 'lm', 'grad', or 'auto'**

Field Name	Description	Default
Tolerance	Minimum percentage difference between the current value of the loss function and its expected improvement after the next iteration, specified as a positive scalar. When the percentage of expected improvement is less than <code>Tolerance</code> , the iterations stop. The estimate of the expected loss-function improvement at the next iteration is based on the Gauss-Newton vector computed for the current parameter value.	0.01
MaxIterations	<p>Maximum number of iterations during loss-function minimization, specified as a positive integer. The iterations stop when <code>MaxIterations</code> is reached or another stopping criterion is satisfied, such as <code>Tolerance</code>.</p> <p>Setting <code>MaxIterations = 0</code> returns the result of the start-up procedure.</p> <p>Use <code>sys.Report.Termination.Iterations</code> to get the actual number of iterations during an estimation, where <code>sys</code> is an <code>idtf</code> model.</p>	20

Field Name	Description	Default
Advanced	Advanced search settings, specified as a structure with the following fields:	
Field Name	Description	Default
GnPinvConstant	Jacobian matrix singular value threshold, specified as a positive scalar. Singular values of the Jacobian matrix that are smaller than $\text{GnPinvConstant} * \max(\text{size}(J) * \text{norm}(J) * \text{eps})$ are discarded when computing the search direction. Applicable when SearchMethod is 'gn'.	10000
InitialGnaTolerance	Initial value of $\gamma$ , specified as a positive scalar. Applicable when SearchMethod is 'gna'.	0.0001
LMStartValue	Starting value of search-direction length $d$ in the Levenberg-Marquardt method, specified as a positive scalar. Applicable when SearchMethod is 'lm'.	0.001
LMStep	Size of the Levenberg-Marquardt step, specified as a positive integer. The next value of the search-direction length $d$ in the Levenberg-Marquardt method is LMStep times the previous one. Applicable when SearchMethod is 'lm'.	2
MaxBisections	Maximum number of bisections used for line search along the search direction, specified as a positive integer.	25
MaxFunctionEvaluations	Maximum number of calls to the model file, specified as a positive integer. Iterations stop if the number of calls to the model file exceeds this value.	Inf
MinParameterChange	Smallest parameter update allowed per iteration, specified as a nonnegative scalar.	0
RelativeImprovement	Relative improvement threshold, specified as a nonnegative scalar. Iterations stop if the relative	0

<b>Field Name</b>	<b>Description</b>		<b>Default</b>
	<b>Field Name</b>	<b>Description</b>	<b>Default</b>
		improvement of the criterion function is less than this value.	
	StepReduction	<p>Step reduction factor, specified as a positive scalar that is greater than 1. The suggested parameter update is reduced by the factor <code>StepReduction</code> after each try. This reduction continues until <code>MaxBisections</code> tries are completed or a lower value of the criterion function is obtained.</p> <p><code>StepReduction</code> is not applicable for SearchMethod 'lm' (Levenberg-Marquardt method).</p>	2

**SearchOptions Structure When SearchMethod is Specified as 'lsqnonlin'**

<b>Field Name</b>	<b>Description</b>	<b>Default</b>
FunctionTolerance	Termination tolerance on the loss function that the software minimizes to determine the estimated parameter values, specified as a positive scalar.  The value of FunctionTolerance is the same as that of <code>opt.SearchOptions.Advanced.TolFun</code> .	1e-5
StepTolerance	Termination tolerance on the estimated parameter values, specified as a positive scalar.  The value of StepTolerance is the same as that of <code>opt.SearchOptions.Advanced.TolX</code> .	1e-6
MaxIterations	Maximum number of iterations during loss-function minimization, specified as a positive integer. The iterations stop when MaxIterations is reached or another stopping criterion is satisfied, such as FunctionTolerance.  The value of MaxIterations is the same as that of <code>opt.SearchOptions.Advanced.MaxIter</code> .	20
Advanced	Advanced search settings, specified as an option set for <code>lsqnonlin</code> .  For more information, see the Optimization Options table in "Optimization Options" (Optimization Toolbox).	Use <code>optimset('lsqnonlin')</code> to create a default option set.

**SearchOptions Structure When SearchMethod is Specified as 'fmincon'**

Field Name	Description	Default
Algorithm	<p>fmincon optimization algorithm, specified as one of the following:</p> <ul style="list-style-type: none"> <li>• 'sqp' — Sequential quadratic programming algorithm. The algorithm satisfies bounds at all iterations, and it can recover from NaN or Inf results. It is not a large-scale algorithm. For more information, see “Large-Scale vs. Medium-Scale Algorithms” (Optimization Toolbox).</li> <li>• 'trust-region-reflective' — Subspace trust-region method based on the interior-reflective Newton method. It is a large-scale algorithm.</li> <li>• 'interior-point' — Large-scale algorithm that requires Optimization Toolbox software. The algorithm satisfies bounds at all iterations, and it can recover from NaN or Inf results.</li> <li>• 'active-set' — Requires Optimization Toolbox software. The</li> </ul>	'sqp'

Field Name	Description	Default
	<p>algorithm can take large steps, which adds speed. It is not a large-scale algorithm.</p> <p>For more information about the algorithms, see “Constrained Nonlinear Optimization Algorithms” (Optimization Toolbox) and “Choosing the Algorithm” (Optimization Toolbox).</p>	
FunctionTolerance	Termination tolerance on the loss function that the software minimizes to determine the estimated parameter values, specified as a positive scalar.	1e-6
StepTolerance	Termination tolerance on the estimated parameter values, specified as a positive scalar.	1e-6
MaxIterations	Maximum number of iterations during loss function minimization, specified as a positive integer. The iterations stop when MaxIterations is reached or another stopping criterion is satisfied, such as FunctionTolerance.	100

### Advanced – Additional advanced options structure

Advanced is a structure with the following fields:

- `ErrorThreshold` — Specifies when to adjust the weight of large errors from quadratic to linear.

Errors larger than `ErrorThreshold` times the estimated standard deviation have a linear weight in the loss function. The standard deviation is estimated robustly as the median of the absolute deviations from the median of the prediction errors, divided by 0.7. For more information on robust norm choices, see section 15.2 of [1].

`ErrorThreshold = 0` disables robustification and leads to a purely quadratic loss function. When estimating with frequency-domain data, the software sets `ErrorThreshold` to zero. For time-domain data that contains outliers, try setting `ErrorThreshold` to 1.6.

**Default:** 0

- `MaxSize` — Specifies the maximum number of elements in a segment when input-output data is split into segments.

`MaxSize` must be a positive integer.

**Default:** 250000

- `StabilityThreshold` — Specifies thresholds for stability tests.

`StabilityThreshold` is a structure with the following fields:

- `s` — Specifies the location of the right-most pole to test the stability of continuous-time models. A model is considered stable when its right-most pole is to the left of `s`.

**Default:** 0

- `z` — Specifies the maximum distance of all poles from the origin to test stability of discrete-time models. A model is considered stable if all poles are within the distance `z` from the origin.

**Default:** `1+sqrt(eps)`

- `AutoInitThreshold` — Specifies when to automatically estimate the initial condition.

The initial condition is estimated when

$$\frac{\|y_{p,z} - y_{meas}\|}{\|y_{p,e} - y_{meas}\|} > \text{AutoInitThreshold}$$

- $y_{meas}$  is the measured output.
- $y_{p,z}$  is the predicted output of a model estimated using zero initial states.
- $y_{p,e}$  is the predicted output of a model estimated using estimated initial states.

Applicable when `InitialCondition` is 'auto'.

**Default:** 1.05

## Output Arguments

**opt** — Option set for `procest`

`procestOptions` option set

Option set for `procest`, returned as a `procestOptions` option set.

## Examples

### Create Default Option Set for Process Model Estimation

```
opt = procestOptions;
```

### Specify Options for Process Model Estimation

Create an option set for `procest` setting `Focus` to 'simulation' and turning on the `Display`.

```
opt = procestOptions('Focus','simulation','Display','on');
```

Alternatively, use dot notation to set the values of `opt`.

```
opt = procestOptions;  
opt.Focus = 'simulation';  
opt.Display = 'on';
```

## References

- [1] Ljung, L. *System Identification: Theory for the User*. Upper Saddle River, NJ: Prentice-Hall PTR, 1999.
- [2] Wills, Adrian, B. Ninness, and S. Gibson. "On Gradient-Based Search for Multivariable System Estimates". *Proceedings of the 16th IFAC World Congress, Prague, Czech Republic, July 3-8, 2005*. Oxford, UK: Elsevier Ltd., 2005.

## See Also

idfilt | idproc | procest

## Topics

"Loss Function and Model Quality Metrics"

**Introduced in R2012a**

# pwlinear

Create a piecewise-linear nonlinearity estimator object

## Syntax

```
NL = pwlinear  
NL = pwlinear(Name, Value)
```

## Description

`NL = pwlinear` creates a default piecewise-linear nonlinearity estimator object with 10 break points for estimating Hammerstein-Wiener models. The value of the nonlinearity at the break points are set to []. The initial value of the nonlinearity is determined from the estimation data range during estimation using `n_lhw`. Use dot notation to customize the object properties, if needed.

`NL = pwlinear(Name, Value)` creates a piecewise-linear nonlinearity estimator object with properties specified by one or more `Name, Value` pair arguments. The properties that you do not specify retain their default value.

## Object Description

`pwlinear` is an object that stores the piecewise-linear nonlinearity estimator for estimating Hammerstein-Wiener models.

Use `pwlinear` to define a nonlinear function  $y = F(x, \theta)$ , where  $y$  and  $x$  are scalars, and  $\theta$  represents the parameters specifying the number of break points and the value of nonlinearity at the break points.

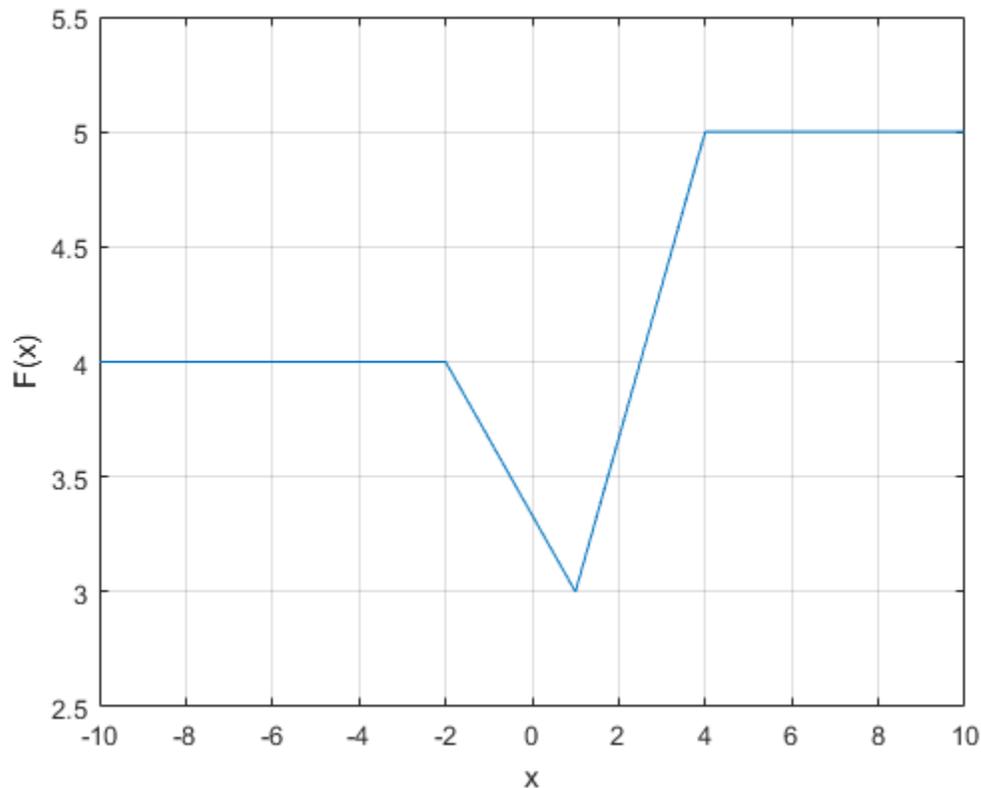
The nonlinearity function,  $F$ , is a piecewise-linear (affine) function of  $x$ . There are  $n$  breakpoints  $(x_k, y_k)$ ,  $k = 1, \dots, n$ , such that  $y_k = F(x_k)$ .  $F$  is linearly interpolated between the breakpoints.

$F$  is also linear to the left and right of the extreme breakpoints. The slope of these extensions is a function of  $x_i$  and  $y_i$  breakpoints. The breakpoints are ordered by

ascending x-values, which is important when you set a specific breakpoint to a different value.

There are minor difference between the breakpoint values you set and the values stored in the object because the toolbox has a different internal representation of breakpoints.

For example, in the following plot, the breakpoints are  $x_k = [-2,1,4]$  and the corresponding nonlinearity values are  $y_k = [4,3,5]$ .



The value  $F(x)$  is computed by `evaluate(NL, x)`, where `NL` is the `pwlinear` object. When using `evaluate`, the break points have to be initialized manually.

For `pwlinear` object properties, see “Properties” on page 1-1290.

## Examples

### Create a Default Piecewise-Linear Nonlinearity Estimator

```
NL = pwlinear;
```

Specify the number of break points.

```
NL.NumberOfUnits = 5;
```

### Estimate a Hammerstein Model with Piecewise-Linear Nonlinearity

Load estimation data.

```
load twotankdata;  
z = iddata(y,u,0.2,'Name','Two tank system');  
z1 = z(1:1000);
```

Create a pwlinear object, and specify the breakpoints.

```
InputNL = pwlinear('BreakPoints',[-2,1,4]);
```

Since BreakPoints is specified as a vector, the specified vector is interpreted as the x-values of the break points. The y-values of the break points are set to 0, and are determined during model estimation.

Estimate model with no output nonlinearity.

```
sys = nlhw(z1,[2 3 0],InputNL,[]);
```

## Input Arguments

### Name-Value Pair Arguments

Specify optional comma-separated pairs of Name, Value arguments. Name is the argument name and Value is the corresponding value. Name must appear inside quotes.

You can specify several name and value pair arguments in any order as `Name1, Value1, . . . , NameN, ValueN`.

Use `Name, Value` arguments to specify additional properties on page 1-662 of `pwnlinear` nonlinearity. For example, `NL= pwnlinear('NumberofUnits',5)` creates a piecewise-linear nonlinearity estimator object with 5 breakpoints.

## Properties

`pwnlinear` object properties include:

### **NumberofUnits**

Number of breakpoints, specified as an integer.

**Default:** 10

### **BreakPoints**

Break points,  $x_k$ , and the corresponding nonlinearity values at the breakpoints,  $y_k$ , specified as one of the following:

- 2-by- $n$  matrix — The  $x$  and  $y$  values for each of the  $n$  break points are specified as  $[x_1, x_2, \dots, x_n; y_1, y_2, \dots, y_n]$ .
- 1-by- $n$  vector — The specified vector is interpreted as the  $x$  values of the break points:  $x_1, x_2, \dots, x_n$ . All the  $y$  values of the break points are set to 0.

When the nonlinearity object is created, the breakpoints are ordered by ascending  $x$ -values. This is important to consider if you set a specific breakpoint to a different value after creating the object.

**Default:** []

## Output Arguments

### **NL — Piecewise-linear nonlinearity estimator object**

`pwnlinear` object

Piecewise-linear nonlinearity estimator object, returned as a `pwnlinear` object.

## **See Also**

n1hw

**Introduced in R2007a**

## pzmap

Pole-zero plot of dynamic system

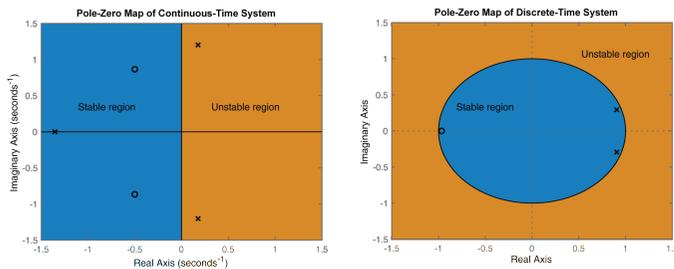
## Syntax

```
pzmap(sys)  
pzmap(sys1,sys2,...,sysN)
```

```
[p,z] = pzmap(sys)
```

## Description

`pzmap(sys)` creates a pole-zero plot of the continuous or discrete-time dynamic system model `sys`. `x` and `o` indicates the poles and zeros respectively, as shown in the following figure.



From the figure above, an open-loop linear time-invariant system is stable if:

- In continuous-time, all the poles on the complex s-plane must be in the left-half plane (blue region) to ensure stability. The system is marginally stable if distinct poles lie on the imaginary axis, that is, the real parts of the poles are zero.
- In discrete-time, all the poles in the complex z-plane must lie inside the unit circle (blue region). The system is marginally stable if it has one or more poles lying on the unit circle.

`pzmap(sys1, sys2, ..., sysN)` creates the pole-zero plot of multiple models on a single figure. The models can have different numbers of inputs and outputs and can be a mix of continuous and discrete systems. For SISO systems, `pzmap` plots the system poles and zeros. For MIMO systems, `pzmap` plots the system poles and transmission zeros.

`[p, z] = pzmap(sys)` returns the system poles and transmission zeros as column vectors `p` and `z`. The pole-zero plot is not displayed on the screen.

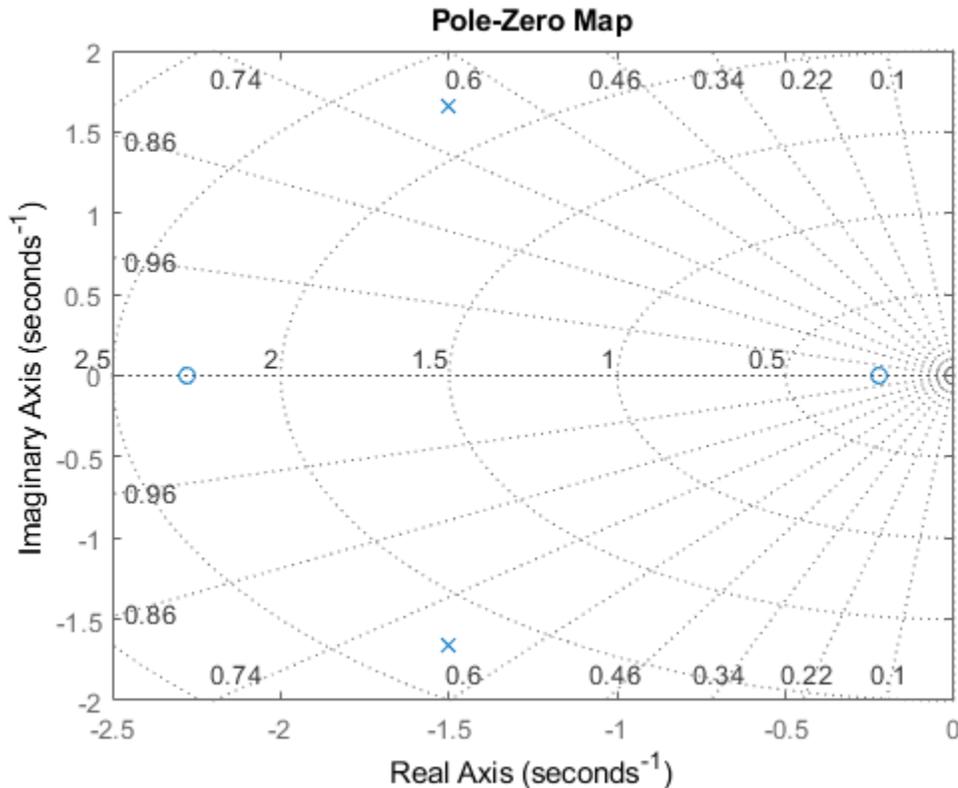
## Examples

### Pole-Zero Plot of Dynamic System

Plot the poles and zeros of the continuous-time system represented by the following transfer function:

$$H(s) = \frac{2s^2 + 5s + 1}{s^2 + 3s + 5}$$

```
H = tf([2 5 1],[1 3 5]);  
pzmap(H)  
grid on
```



Turning on the grid displays lines of constant damping ratio (zeta) and lines of constant natural frequency ( $\omega_n$ ). This system has two real zeros, marked by o on the plot. The system also has a pair of complex poles, marked by x.

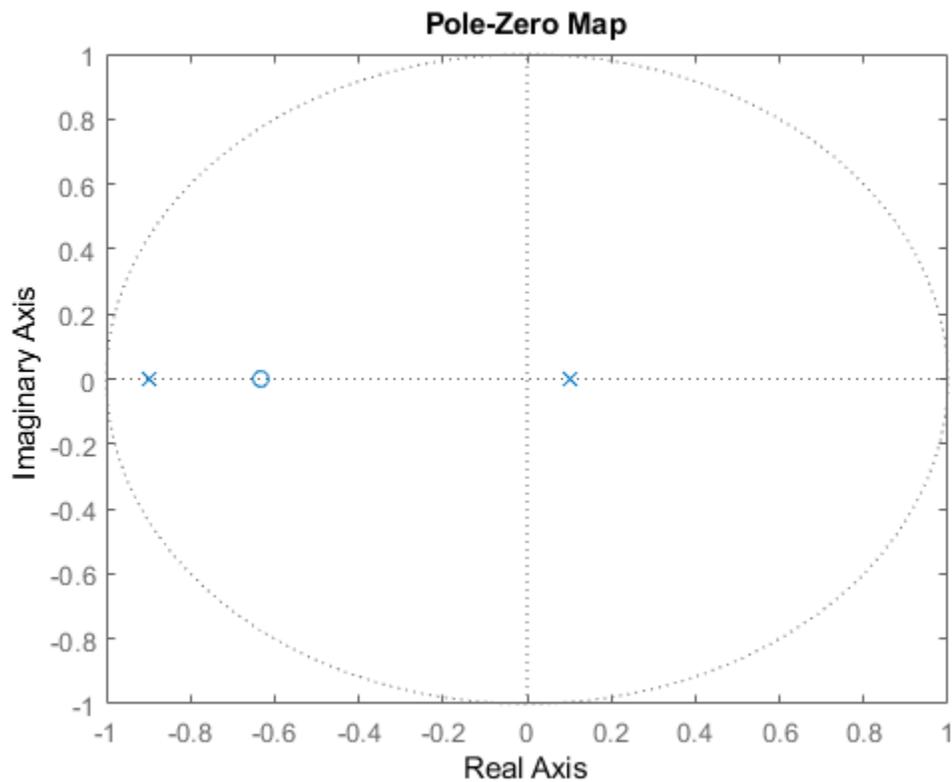
### Pole-Zero Plot of Identified System

Plot the pole-zero map of a discrete time identified state-space (idss) model. In practice you can obtain an idss model by estimation based on input-output measurements of a system. For this example, create one from state-space data.

```
A = [0.1 0; 0.2 -0.9];  
B = [.1 ; 0.1];  
C = [10 5];  
D = [0];  
sys = idss(A,B,C,D,'Ts',0.1);
```

Examine the pole-zero map.

```
pzmap(sys)
```



System poles are marked by x, and zeros are marked by o.

### Pole-Zero Map of Multiple Models

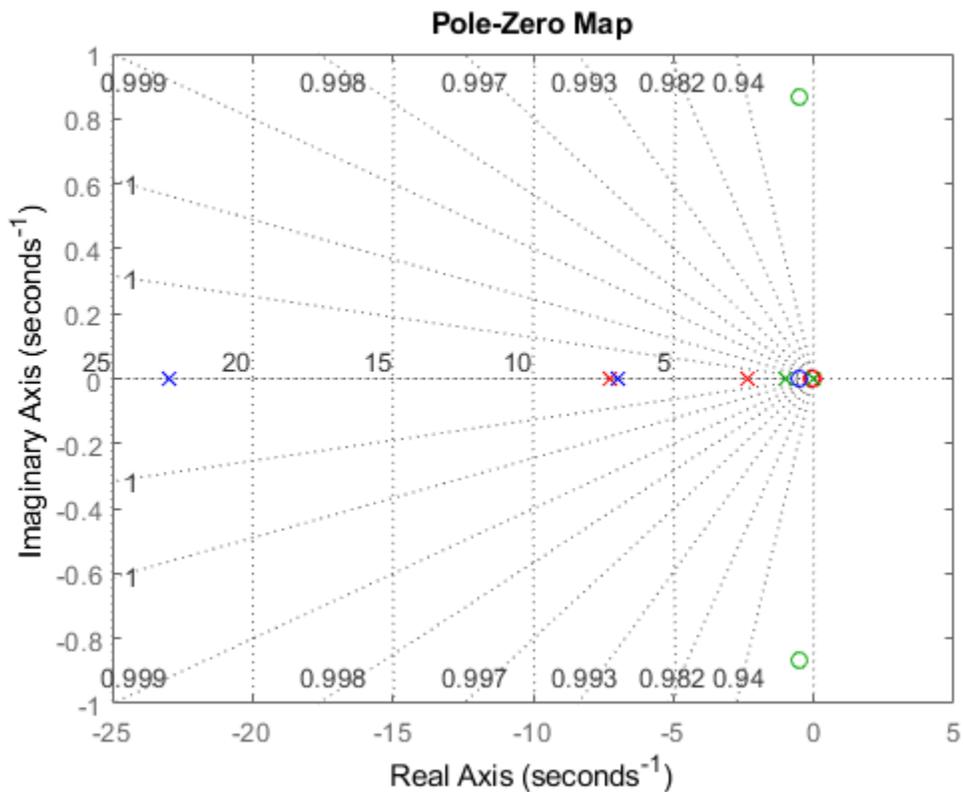
For this example, load a 3-by-1 array of transfer function models.

```
load('tfArray.mat', 'sys');
size(sys)
```

```
3x1 array of transfer functions.
Each model has 1 outputs and 1 inputs.
```

Plot the poles and zeros of each model in the array with distinct colors. For this example, use red for the first model, green for the second and blue for the third model in the array.

```
pzmap(sys(:, :, 1), 'r', sys(:, :, 2), 'g', sys(:, :, 3), 'b')
sgrid
```



sgrid plots lines of constant damping ratio and natural frequency in the s-plane of the pole-zero plot.

## Poles and Zeros of Transfer Function

Use pzmap to calculate the poles and zeros of the following transfer function:

$$\text{sys}(s) = \frac{4.2s^2 + 0.25s - 0.004}{s^2 + 9.6s + 17}$$

```
sys = tf([4.2,0.25,-0.004],[1,9.6,17]);
[p,z] = pzmap(sys)
```

```
p = 2×1
```

```
-7.2576
-2.3424
```

```
z = 2×1
```

```
-0.0726
0.0131
```

## Identify Near-Cancelling Pole-Zero Pairs

This example uses a model of a building with eight floors, each with three degrees of freedom: two displacements and one rotation. The I/O relationship for any one of these displacements is represented as a 48-state model, where each state represents a displacement or its rate of change (velocity).

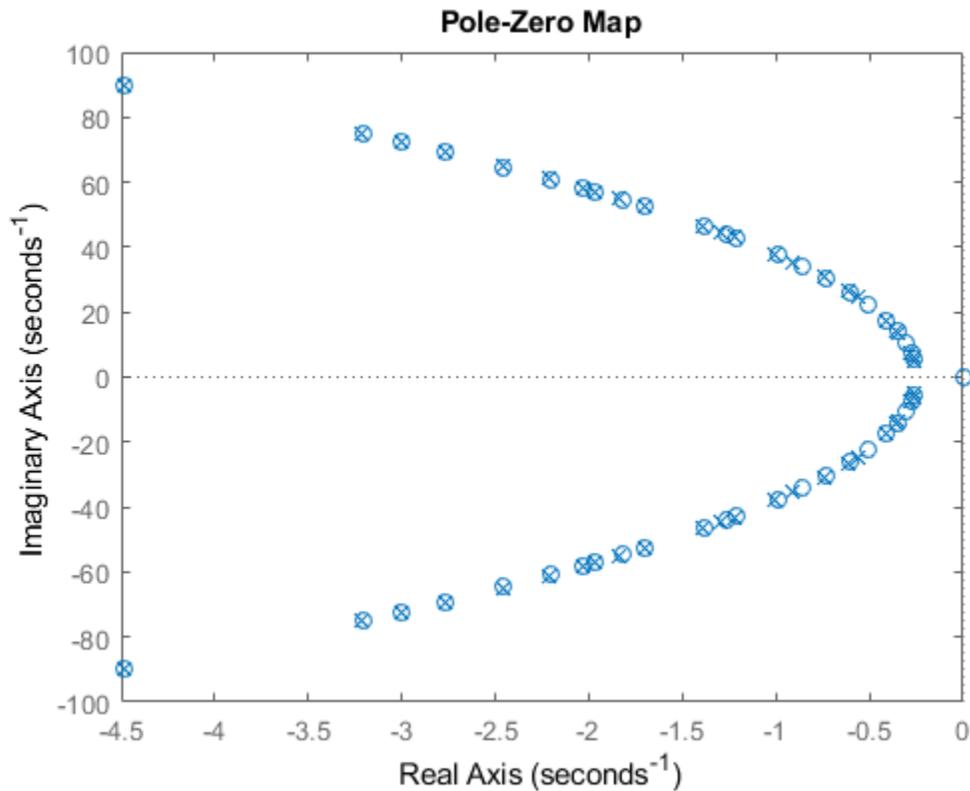
Load the building model.

```
load('building.mat');
size(G)
```

State-space model with 1 outputs, 1 inputs, and 48 states.

Plot the poles and zeros of the system.

pzmap(G)



From the plot, observe that there are numerous near-canceling pole-zero pairs that could be potentially eliminated to simplify the model, with no effect on the overall model response. pzmap is useful to visually identify such near-canceling pole-zero pairs to perform pole-zero simplification.

## Input Arguments

### **sys** — Dynamic system

dynamic system model | model array

Dynamic system, specified as a dynamic system model or model array. Dynamic systems that you can use include continuous-time or discrete-time numeric LTI models such as `tf`, `zpk`, or `ss` models.

If `sys` is an array of models, `pzmap` plots all the poles and zeros of every model in the array on the same plot.

## Output Arguments

### **p** — Poles of the system

column vector

Poles of the system, returned as a column vector, in order of its increasing natural frequency. `p` is the same as the output of `pole(sys)`, except for the order.

### **z** — Transmission zeros of the system

column vector

Transmission zeros of the system, returned as a column vector. `z` is the same as the output of `tzero(sys)`.

## Tips

- Use the functions `sgrid` or `zgrid` to plot lines of constant damping ratio and natural frequency in the  $s$ - or  $z$ -plane on the pole-zero plot.
- For MIMO models, `pzmap` displays all system poles and transmission zeros on a single plot. To map poles and zeros for individual I/O pairs, use `iopzmap`.
- For additional options to customize the appearance of the pole-zero plot, use `pzplot`.

## See Also

`damp` | `dsort` | `esort` | `iopzmap` | `pole` | `pzplot` | `rlocus` | `sgrid` | `zero` | `zgrid`

**Introduced before R2006a**

# pzoptions

Create list of pole/zero plot options

## Syntax

P = pzoptions

P = pzoption('cstprefs')

## Description

P = pzoptions returns a list of available options for pole/zero plots (pole/zero, input-output pole/zero and root locus) with default values set.. You can use these options to customize the pole/zero plot appearance from the command line.

P = pzoption('cstprefs') initializes the plot options with the options you selected in the Control System and System Identification Toolbox Preferences Editor. For more information about the editor, see “Toolbox Preferences Editor”.

This table summarizes the available pole/zero plot options.

Option	Description
Title, XLabel, YLabel	Label text and style
TickLabel	Tick label style
Grid	Show or hide the grid, specified as one of the following values: 'off'   'on' <b>Default:</b> 'off'
GridColor	Color of the grid lines, specified as one of the following: Vector of RGB values in the range [0,1]  character vector of color name   'none'. For example, for yellow color, specify as one of the following: [1 1 0], 'yellow', or 'y'. <b>Default:</b> [0.15,0.15,0.15]

<b>Option</b>	<b>Description</b>
XlimMode, YlimMode	Limit modes
Xlim, Ylim	Axes limits
IOGrouping	Grouping of input-output pairs, specified as one of the following values: 'none'   'inputs'   'outputs'   'all' <b>Default:</b> 'none'
InputLabel, OutputLabel	Input and output label styles
InputVisible, OutputVisible	Visibility of input and output channels

Option	Description
FreqUnits	<p>Frequency units, specified as one of the following values:</p> <ul style="list-style-type: none"><li>• 'Hz'</li><li>• 'rad/second'</li><li>• 'rpm'</li><li>• 'kHz'</li><li>• 'MHz'</li><li>• 'GHz'</li><li>• 'rad/nanosecond'</li><li>• 'rad/microsecond'</li><li>• 'rad/millisecond'</li><li>• 'rad/minute'</li><li>• 'rad/hour'</li><li>• 'rad/day'</li><li>• 'rad/week'</li><li>• 'rad/month'</li><li>• 'rad/year'</li><li>• 'cycles/nanosecond'</li><li>• 'cycles/microsecond'</li><li>• 'cycles/millisecond'</li><li>• 'cycles/hour'</li><li>• 'cycles/day'</li><li>• 'cycles/week'</li><li>• 'cycles/month'</li><li>• 'cycles/year'</li></ul> <p><b>Default:</b> 'rad/s'</p> <p>You can also specify 'auto' which uses frequency units rad/TimeUnit relative to system time units specified in the</p>

Option	Description
	TimeUnit property. For multiple systems with different time units, the units of the first system are used.
TimeUnits	<p>Time units, specified as one of the following values:</p> <ul style="list-style-type: none"> <li>• 'nanoseconds'</li> <li>• 'microseconds'</li> <li>• 'milliseconds'</li> <li>• 'seconds'</li> <li>• 'minutes'</li> <li>• 'hours'</li> <li>• 'days'</li> <li>• 'weeks'</li> <li>• 'months'</li> <li>• 'years'</li> </ul> <p><b>Default:</b> 'seconds'</p> <p>You can also specify 'auto' which uses time units specified in the TimeUnit property of the input system. For multiple systems with different time units, the units of the first system is used.</p>
ConfidenceRegionNumberSD	Number of standard deviations to use when displaying the confidence region characteristic for identified models (valid only iopzplot).

## Examples

## Display Grid in Root Locus Plot

Create a default option set for root locus plots.

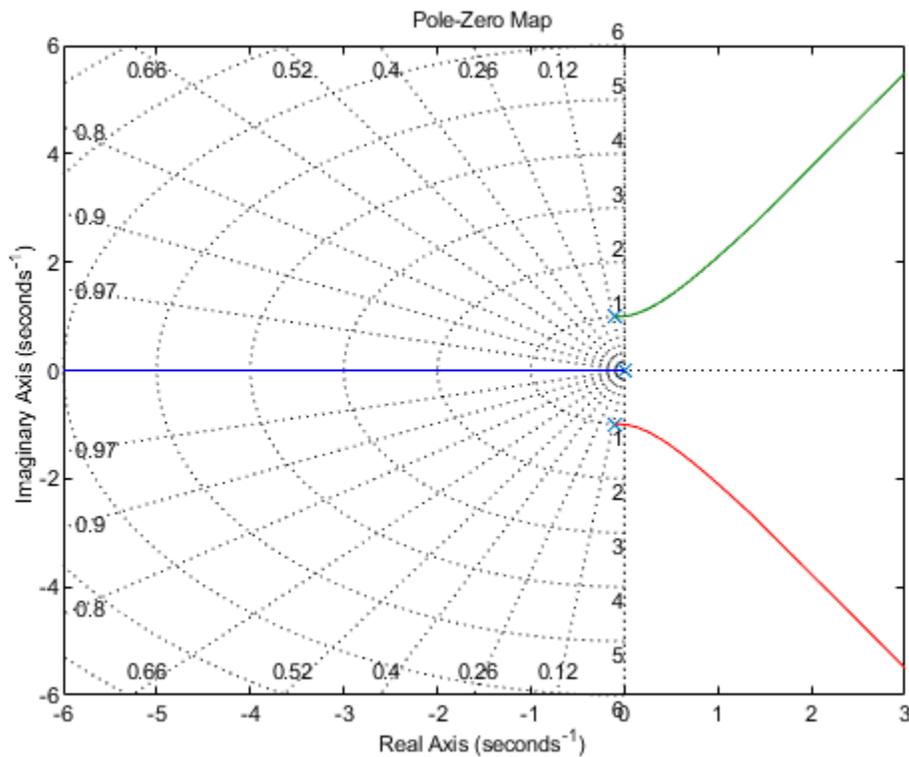
```
opt = pzoptions;
```

In the plot options, specify the grid as visible.

```
opt.Grid = 'on';
```

Use the specified options to plot the root locus of a transfer function.

```
h = rlocusplot(tf(1,[1,.2,1,0]),opt);
```



The root locus plot is generated with the grid enabled.

## **See Also**

`getoptions` | `iopzplot` | `pzplot` | `setoptions`

**Introduced in R2012a**

# pzplot

Pole-zero plot of dynamic system model with plot customization options

## Syntax

```
h = pzplot(sys)

pzplot(sys)
pzplot(sys1,sys2,...,sysN)
pzplot(ax,...)
pzplot(...,plotoptions)
```

## Description

`h = pzplot(sys)` plots the poles and transmission zeros of the dynamic system model `sys` and returns the plot handle `h` to the plot. Use this function to generate a pole-zero map with customizable plot options such as `FreqUnits`, `TimeUnits`, and `IOMGrouping`. For more information about using and interpreting pole-zero maps, see `pzmap`.

`pzplot(sys)` plots the poles and transmission zeros of the dynamic system model `sys`. `x` and `o` indicates poles and zeros respectively.

`pzplot(sys1,sys2,...,sysN)` displays the poles and transmission zeros of multiple models on a single plot. You can specify distinct colors for each model individually.

`pzplot(ax,...)` plots into the axes specified by `ax` instead of the current axis `gca`.

`pzplot(...,plotoptions)` plots the poles and transmission zeros with the options specified in `plotoptions`. For more information on the ways to change properties of your plots, see “Ways to Customize Plots” (Control System Toolbox).

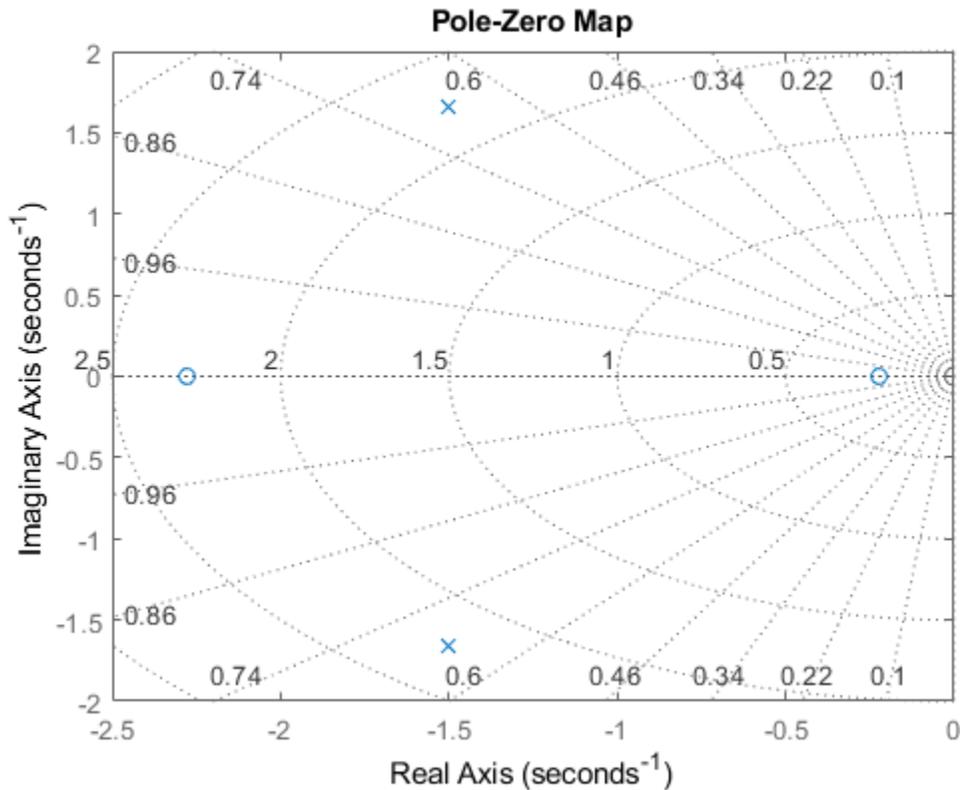
## Examples

### Pole-Zero Plot with Custom Plot Title

Plot the poles and zeros of the continuous-time system represented by the following transfer function:

$$sys(s) = \frac{2s^2 + 5s + 1}{s^2 + 3s + 5}$$

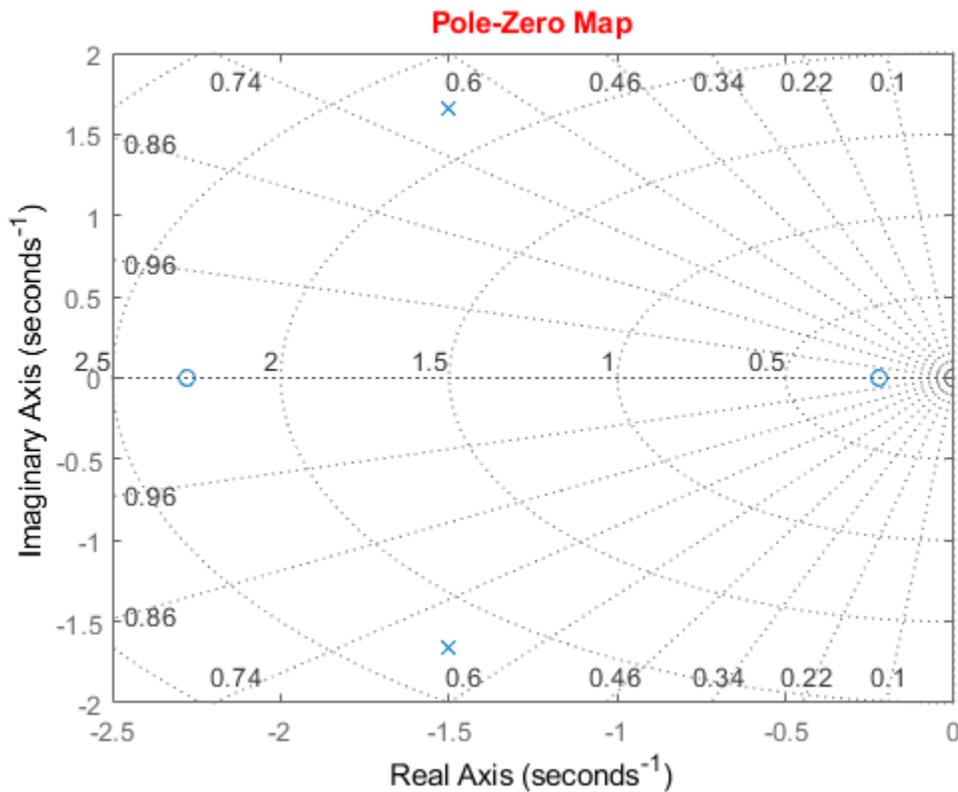
```
sys = tf([2 5 1],[1 3 5]);
h = pzplot(sys);
grid on
```



Turning on the grid displays lines of constant damping ratio ( $\zeta$ ) and lines of constant natural frequency ( $\omega_n$ ). This system has two real zeros, marked by o on the plot. The system also has a pair of complex poles, marked by x.

Change the color of the plot title. To do so, use the plot handle, h.

```
p = getoptions(h);
p.Title.Color = [1,0,0];
setoptions(h,p);
```



### Pole-Zero Plot of Multiple Models

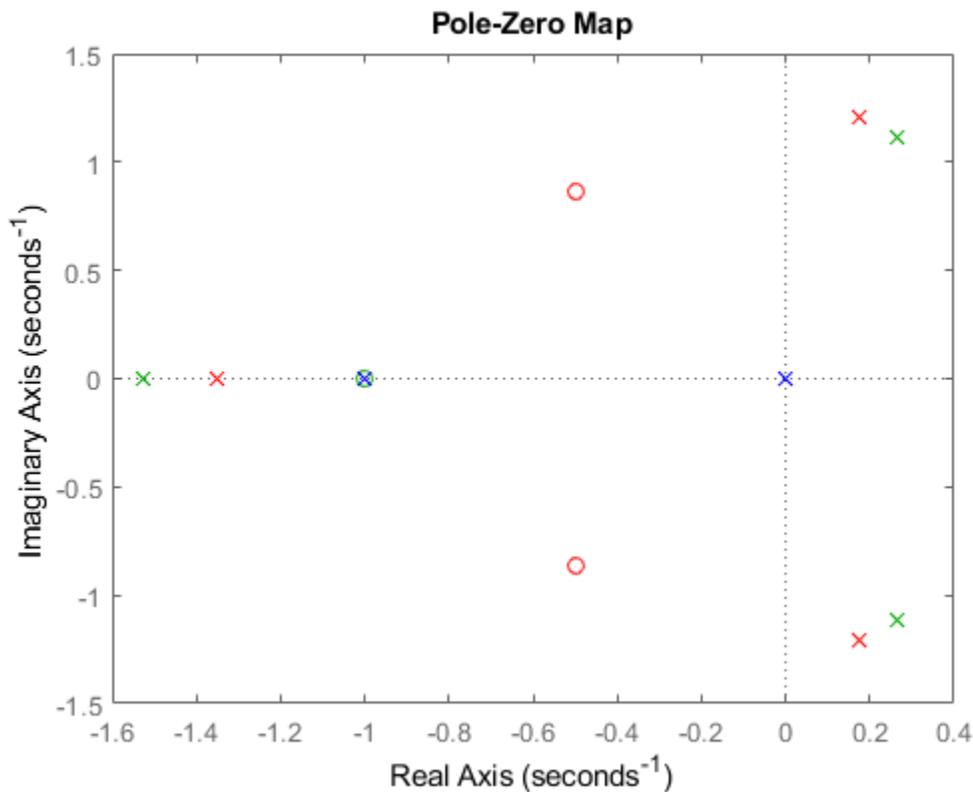
For this example, load a 3-by-1 array of transfer function models.

```
load('tfArrayMargin.mat','sys');  
size(sys)
```

```
3x1 array of transfer functions.  
Each model has 1 outputs and 1 inputs.
```

Plot the poles and zeros of the model array. Define the colors for each model. For this example, use red for the first model, green for the second and blue for the third model in the array.

```
pzplot(sys(:,:,1),'r',sys(:,:,2),'g',sys(:,:,3),'b');
```



### Pole-Zero Plot with Custom Options

Plot the poles and zeros of the continuous-time system represented by the following transfer function with a custom option set:

$$\text{sys}(s) = \frac{2s^2 + 5s + 1}{s^2 + 3s + 5}$$

Create the custom option set using `pzoptions`.

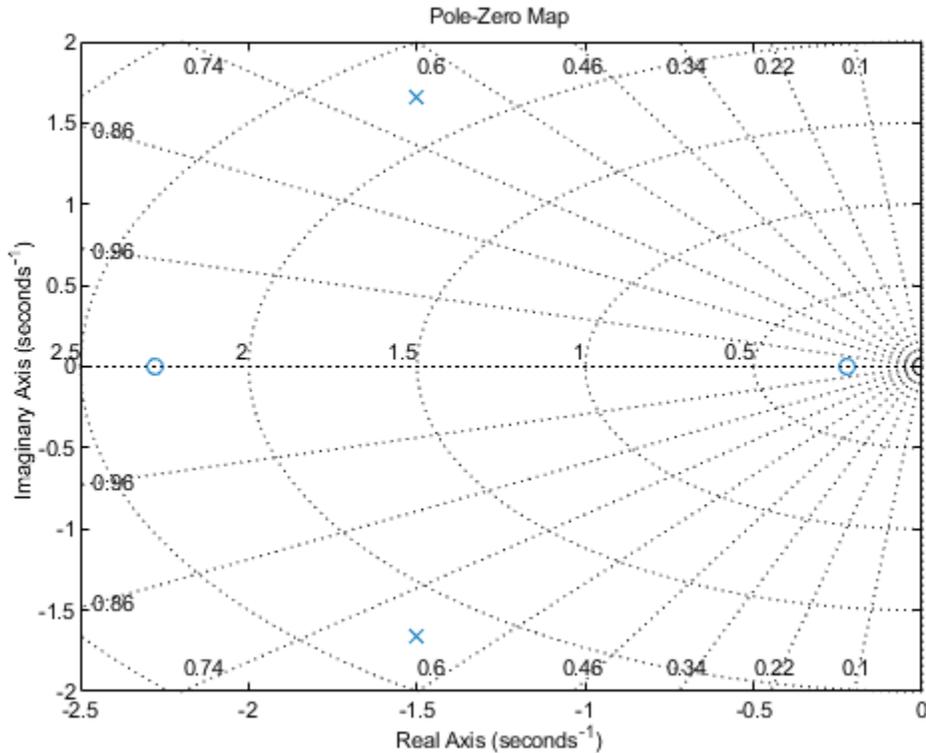
```
plotoptions = pzoptions;
```

For this example, specify the grid to be visible.

```
plotoptions.Grid = 'on';
```

Use the specified options to create a pole-zero map of the transfer function.

```
h = pzplot(tf([2 5 1],[1 3 5]),plotoptions);
```



Turning on the grid displays lines of constant damping ratio (zeta) and lines of constant natural frequency ( $\omega_n$ ). This system has two real zeros, marked by o on the plot. The system also has a pair of complex poles, marked by x.

## Input Arguments

**sys** — Dynamic system

dynamic system model | model array

Dynamic system, specified as a SISO or MIMO dynamic system model, or an array of SISO or MIMO dynamic system models. Dynamic systems that you can use include continuous-time or discrete-time numeric LTI models such as `tf`, `zpk`, or `ss` models.

If `sys` is a generalized state-space model `genss` or an uncertain state-space model `uss`, `pzplot` returns the poles and transmission of the current or nominal value of `sys`. If `sys` is an array of models, `pzplot` plots the poles and zeros of each model in the array on the same diagram.

### **ax — Axes handle**

axes object

Axes handle, specified as an axes object. If you do not specify the axes object, then `pzplot` uses the current axes `gca` to plot the poles and zeros of the system.

### **plotoptions — Pole-zero plot options**

options object

Pole-zero plot options, specified as an options object. See `pzoptions` for a list of available plot options.

## **Output Arguments**

### **h — Pole-zero plot options handle**

scalar

Pole-zero plot options handle, returned as a scalar. Use `h` to query and modify properties of your pole-zero plot. You can use this handle to customize the plot with the `getoptions` and `setoptions` commands.

## **Tips**

- Use `sgrid` or `zgrid` to plot lines of constant damping ratio and natural frequency in the *s*- or *z*-plane.

## **See Also**

`getoptions` | `iopzplot` | `pzmap` | `pzoptions` | `setoptions`

**Topics**

“Ways to Customize Plots” (Control System Toolbox)

**Introduced before R2006a**

## rarmax

(To be removed) Estimate recursively parameters of ARMAX or ARMA models

---

**Note** rarmax will be removed in a future release. Use recursiveARMA or recursiveARMAX instead.

---

## Syntax

```
thm = rarmax(z,nn,adm,adg)
```

```
[thm,yhat,P,phi,psi] = rarmax(z,nn,adm,adg,th0,P0,phi0,psi0)
```

## Description

The parameters of the ARMAX model structure

$$A(q)y(t) = B(q)u(t - nk) + C(q)e(t)$$

are estimated using a recursive prediction error method.

The input-output data is contained in `z`, which is either an `iddata` object or a matrix  $z = \begin{bmatrix} y & u \end{bmatrix}$  where `y` and `u` are column vectors. `nn` is given as

```
nn = [na nb nc nk]
```

where `na`, `nb`, and `nc` are the orders of the ARMAX model, and `nk` is the delay. Specifically,

$$na: A(q) = 1 + a_1q^{-1} + \dots + a_{na}q^{-na}$$

$$nb: B(q) = b_1 + b_2q^{-1} + \dots + b_{nb}q^{-nb+1}$$

$$nc: C(q) = 1 + c_1q^{-1} + \dots + c_{nc}q^{-nc}$$

See “What Are Polynomial Models?” for more information.

If  $z$  represents a time series  $y$  and  $nn = [na \ nc]$ , `rarmax` estimates the parameters of an ARMA model for  $y$ .

$$A(q)y(t) = C(q)e(t)$$

Only single-input, single-output models are handled by `rarmax`. Use `rpem` for the multiple-input case.

The estimated parameters are returned in the matrix `thm`. The  $k$ th row of `thm` contains the parameters associated with time  $k$ ; that is, they are based on the data in the rows up to and including row  $k$  in  $z$ . Each row of `thm` contains the estimated parameters in the following order:

$$\text{thm}(k, :) = [a_1, a_2, \dots, a_n, b_1, \dots, b_n, c_1, \dots, c_n]$$

`yhat` is the predicted value of the output, according to the current model; that is, row  $k$  of `yhat` contains the predicted value of  $y(k)$  based on all past data.

The actual algorithm is selected with the two arguments `adm` and `adg`. These are described under `rarx`.

The input argument `th0` contains the initial value of the parameters, a row vector consistent with the rows of `thm`. The default value of `th0` is all zeros.

The arguments `P0` and `P` are the initial and final values, respectively, of the scaled covariance matrix of the parameters. See `rarx`. The default value of `P0` is  $10^4$  times the unit matrix. The arguments `phi0`, `psi0`, `phi`, and `psi` contain initial and final values of the data vector and the gradient vector, respectively. The sizes of these depend on the chosen model orders. The normal choice of `phi0` and `psi0` is to use the outputs from a previous call to `rarmax` with the same model orders. (This call could be a dummy call with default input arguments.) The default values of `phi0` and `psi0` are all zeros.

Note that the function requires that the delay  $nk$  be larger than  $\theta$ . If you want  $nk = \theta$ , shift the input sequence appropriately and use  $nk = 1$ .

## Examples

Compute and plot, as functions of time, the four parameters in a second-order ARMA model of a time series given in the vector  $y$ . The forgetting factor algorithm with a forgetting factor of 0.98 is applied.

```
thm = rarmax(y,[2 2], 'ff', 0.98);  
plot(thm)
```

## Algorithms

The general recursive prediction error algorithm (11.44), (11.47) through (11.49) of Ljung (1999) is implemented. See “Recursive Algorithms for Online Parameter Estimation” for more information.

## See Also

[nkshift](#) | [recursiveARMA](#) | [recursiveARMAX](#) | [rpem](#) | [rplr](#)

## Topics

“Recursive Algorithms for Online Parameter Estimation”

**Introduced before R2006a**

## rarx

(To be removed) Estimate parameters of ARX or AR models recursively

---

**Note** rarx will be removed in a future release. Use recursiveAR or recursiveARX instead.

---

## Syntax

```
thm = rarx(z, nn, adm, adg)
[thm, yhat, P, phi] = rarx(z, nn, adm, adg, th0, P0, phi0)
```

## Description

`thm = rarx(z, nn, adm, adg)` estimates the parameters `thm` of single-output ARX model from input-output data `z` and model orders `nn` using the algorithm specified by `adm` and `adg`. If `z` is a time series `y` and `nn = na`, `rarx` estimates the parameters of a single-output AR model.

`[thm, yhat, P, phi] = rarx(z, nn, adm, adg, th0, P0, phi0)` estimates the parameters `thm`, the predicted output `yhat`, final values of the scaled covariance matrix of the parameters `P`, and final values of the data vector `phi` of single-output ARX model from input-output data `z` and model orders `nn` using the algorithm specified by `adm` and `adg`. If `z` is a time series `y` and `nn = na`, `rarx` estimates the parameters of a single-output AR model.

## Input Arguments

`z`

Name of the matrix `iddata` object that represents the input-output data or a matrix `z = [y u]`, where `y` and `u` are column vectors.

For multiple-input models, the `u` matrix contains each input as a column vector:

$u = [u_1 \dots u_n]$

**nn**

For input-output models, specifies the structure of the ARX model as:

$nn = [n_a \ n_b \ n_k]$

where  $n_a$  and  $n_b$  are the orders of the ARX model, and  $n_k$  is the delay.

For multiple-input models,  $n_b$  and  $n_k$  are row vectors that define orders and delays for each input.

For time-series models,  $nn = n_a$ , where  $n_a$  is the order of the AR model.

---

**Note** The delay  $n_k$  must be larger than 0. If you want  $n_k = 0$ , shift the input sequence appropriately and use  $n_k = 1$  (see `nkshift`).

---

**adm and adg**

$adm = 'ff'$  and  $adg = lam$  specify the *forgetting factor* algorithm with the forgetting factor  $\lambda = lam$ . This algorithm is also known as recursive least squares (RLS). In this case, the matrix  $P$  has the following interpretation:  $R_2/2 * P$  is approximately equal to the covariance matrix of the estimated parameters.  $R_2$  is the variance of the innovations (the true prediction errors  $e(t)$ ).

$adm = 'ug'$  and  $adg = gam$  specify the *unnormalized gradient* algorithm with gain  $gamma = gam$ . This algorithm is also known as the normalized least mean squares (LMS).

$adm = 'ng'$  and  $adg = gam$  specify the *normalized gradient* or normalized least mean squares (NLMS) algorithm. In these cases,  $P$  is not applicable.

$adm = 'kf'$  and  $adg = R1$  specify the *Kalman filter based* algorithm with  $R_2=1$  and  $R_1 = R1$ . If the variance of the innovations  $e(t)$  is not unity but  $R_2$ ; then  $R_2 * P$  is the covariance matrix of the parameter estimates, while  $R_1 = R1 / R_2$  is the covariance matrix of the parameter changes.

**th0**

Initial value of the parameters in a row vector, consistent with the rows of **thm**.

Default: All zeros.

**P0**

Initial values of the scaled covariance matrix of the parameters.

Default:  $10^4$  times the identity matrix.

**phi0**

The argument **phi0** contains the initial values of the data vector:

$$\varphi(t) = [y(t-1), \dots, y(t-na), u(t-1), \dots, u(t-nb-nk+1)]$$

If  $z = [y(1), u(1); \dots; y(N), u(N)]$ ,  $\text{phi0} = \varphi(1)$  and  $\text{phi} = \varphi(N)$ . For online use of **rarx**, use **phi0**, **th0**, and **P0** as the previous outputs **phi**, **thm** (last row), and **P**.

Default: All zeros.

## Output Arguments

**thm**

Estimated parameters of the model. The  $k$ th row of **thm** contains the parameters associated with time  $k$ ; that is, the estimate parameters are based on the data in rows up to and including row  $k$  in **z**. Each row of **thm** contains the estimated parameters in the following order:

$$\text{thm}(k, :) = [a_1, a_2, \dots, a_n, b_1, \dots, b_n]$$

For a multiple-input model, the  $b$  are grouped by input. For example, the  $b$  parameters associated with the first input are listed first, and the  $b$  parameters associated with the second input are listed next.

**yhat**

Predicted value of the output, according to the current model; that is, row  $k$  of **yhat** contains the predicted value of  $y(k)$  based on all past data.

**P**

Final values of the scaled covariance matrix of the parameters.

**phi**

**phi** contains the final values of the data vector:

$$\varphi(t) = [y(t-1), \dots, y(t-na), u(t-1), \dots, u(t-nb-nk+1)]$$

## Examples

Adaptive noise canceling: The signal  $y$  contains a component that originates from a known signal  $r$ . Remove this component by recursively estimating the system that relates  $r$  to  $y$  using a sixth-order FIR model and the NLMS algorithm.

```
z = [y r];
[thm,noise] = rarx(z,[0 6 1],'ng',0.1);
% noise is the adaptive estimate of the noise
% component of y
plot(y-noise)
```

If this is an online application, you can plot the best estimate of the signal  $y - \text{noise}$  at the same time as the data  $y$  and  $u$  become available, use the following code:

```
phi = zeros(6,1);
P=1000*eye(6);
th = zeros(1,6);
axis([0 100 -2 2]);
plot(0,0,'*'), hold on
% Use a while loop
while ~abort
[y,r,abort] = readAD(time);
[th,ns,P,phi] = rarx([y r],'ff',0.98,th,P,phi);
plot(time,y-ns,'*')
time = time + Dt
end
```

This example uses a forgetting factor algorithm with a forgetting factor of 0.98. `readAD` is a function that reads the value of an A/D converter at the indicated time instant.

## Definitions

### ARX Model Structure

The general ARX model structure is:

$$A(q)y(t) = B(q)u(t - nk) + e(t)$$

The orders of the ARX model are:

$$na: A(q) = 1 + a_1q^{-1} + \dots + a_{na}q^{-na}$$

$$nb: B(q) = b_1 + b_2q^{-1} + \dots + b_{nb}q^{-nb+1}$$

Models with several inputs are defined, as follows:

$$A(q)y(t) = B_1(q)u_1(t-nk_1) + \dots + B_{nu}(q)u_{nu}(t-nk_{nu}) + e(t)$$

## See Also

`nkshift` | `recursiveAR` | `recursiveARX` | `rpem` | `rplr`

## Topics

“Recursive Algorithms for Online Parameter Estimation”

**Introduced before R2006a**

## rbj

(To be removed) Estimate recursively parameters of Box-Jenkins models

---

**Note** rbj will be removed in a future release. Use recursiveBJ instead.

---

## Syntax

thm = rbj(z, nn, adm, adg)

[thm, yhat, P, phi, psi] = rbj(z, nn, adm, adg, th0, P0, phi0, psi0)

## Description

The parameters of the Box-Jenkins model structure

$$y(t) = \frac{B(q)}{F(q)} u(t-nk) + \frac{C(q)}{D(q)} e(t)$$

are estimated using a recursive prediction error method.

The input-output data is contained in z, which is either an `iddata` object or a matrix  $z = [y \ u]$  where y and u are column vectors. nn is given as

nn = [nb nc nd nf nk]

where nb, nc, nd, and nf are the orders of the Box-Jenkins model, and nk is the delay. Specifically,

$$nb: B(q) = b_1 + b_2q^{-1} + \dots + b_{nb}q^{-nb+1}$$

$$nc: C(q) = 1 + c_1q^{-1} + \dots + c_{nc}q^{-nc}$$

$$nd: D(q) = 1 + d_1q^{-1} + \dots + d_{nd}q^{-nd}$$

$$nf: F(q) = 1 + f_1q^{-1} + \dots + f_{nf}q^{-nf}$$

See “What Are Polynomial Models?” for more information.

Only single-input, single-output models are handled by `rbj`. Use `rpem` for the multiple-input case.

The estimated parameters are returned in the matrix `thm`. The  $k$ th row of `thm` contains the parameters associated with time  $k$ ; that is, they are based on the data in the rows up to and including row  $k$  in `z`. Each row of `thm` contains the estimated parameters in the following order:

```
thm(k,:) = [b1,...,bnb,c1,...,cnc,d1,...,dnd,f1,...,fnf]
```

`yhat` is the predicted value of the output, according to the current model; that is, row  $k$  of `yhat` contains the predicted value of  $y(k)$  based on all past data.

The actual algorithm is selected with the two arguments `adm` and `adg`. These are described under `rarx`.

The input argument `th0` contains the initial value of the parameters, a row vector consistent with the rows of `thm`. The default value of `th0` is all zeros.

The arguments `P0` and `P` are the initial and final values, respectively, of the scaled covariance matrix of the parameters. See `rarx`. The default value of `P0` is  $10^4$  times the unit matrix. The arguments `phi0`, `psi0`, `phi`, and `psi` contain initial and final values of the data vector and the gradient vector, respectively. The sizes of these depend on the chosen model orders. The normal choice of `phi0` and `psi0` is to use the outputs from a previous call to `rbj` with the same model orders. (This call could be a dummy call with default input arguments.) The default values of `phi0` and `psi0` are all zeros.

Note that the function requires that the delay  $n_k$  be larger than  $\theta$ . If you want  $n_k = \theta$ , shift the input sequence appropriately and use  $n_k = 1$ .

## Algorithms

The general recursive prediction error algorithm (11.44) of Ljung (1999) is implemented. See also “Recursive Algorithms for Online Parameter Estimation”.

## See Also

`nkshift` | `recursiveBJ` | `rpem` | `rplr`

## **Topics**

“Recursive Algorithms for Online Parameter Estimation”

**Introduced before R2006a**

## realdata

Determine whether `iddata` is based on real-valued signals

### Syntax

```
realdata(data)
```

### Description

`realdata` returns 1 if

- `data` contains only real-valued signals.
- `data` contains frequency-domain signals, obtained by Fourier transformation of real-valued signals.

Otherwise `realdata` returns 0.

### Examples

#### Determine if Data is Based on Real-Valued Signals

Load data.

```
load iddata1
```

Transform the data to frequency domain.

```
zf = fft(z1);
```

Determine if the time-domain data values are real.

```
isreal(z1)
```

```
ans = 1
```

Determine if the transformed data values are real.

```
isreal(zf)
```

```
ans = 0
```

Determine if the data is based on real-valued signals.

```
realdata(zf)
```

```
ans = logical  
     1
```

Add negative frequencies to `zf` and rerun the command.

```
zf = complex(zf);  
realdata(zf)
```

```
ans = logical  
     1
```

The command still returns 1.

**Introduced before R2006a**

## recursiveAR

Create System object for online parameter estimation of AR model

Use the `recursiveAR` command for parameter estimation with real-time data. If all data necessary for estimation is available at once, and you are estimating a time-invariant model, use the offline estimation command, `ar`.

### Syntax

```
obj = recursiveAR
obj = recursiveAR(na)
obj = recursiveAR(na,A0)
obj = recursiveAR( ____,Name,Value)
```

### Description

`obj = recursiveAR` creates a System object for online parameter estimation of a default single output AR model structure on page 1-1342. The default model structure has a polynomial of order 1 and initial polynomial coefficient value `eps`.

After creating the object, use the `step` command to update model parameter estimates using recursive estimation algorithms and real-time data.

`obj = recursiveAR(na)` specifies the polynomial order of the AR model to be estimated.

`obj = recursiveAR(na,A0)` specifies the polynomial order and initial values of the polynomial coefficients.

`obj = recursiveAR( ____,Name,Value)` specifies additional attributes of the AR model structure and recursive estimation algorithm using one or more `Name,Value` pair arguments.

## Object Description

recursiveAR creates a System object for online parameter estimation of single output AR models using a recursive estimation algorithm.

A System object is a specialized MATLAB object designed specifically for implementing and simulating dynamic systems with inputs that change over time. System objects use internal states to store past behavior, which is used in the next computational step.

After you create a System object, you use commands to process data or obtain information from or about the object. System objects use a minimum of two commands to process data — a constructor to create the object and the `step` command to update object parameters using real-time data. This separation of declaration from execution lets you create multiple, persistent, reusable objects, each with different settings.

You can use the following commands with the online estimation System objects in System Identification Toolbox:

Command	Description
<code>step</code>	Update model parameter estimates using recursive estimation algorithms and real-time data.  <code>step</code> puts the object into a locked state. In a locked state, you cannot change any nontunable properties or input specifications, such as model order, data type, or estimation algorithm. During execution, you can only change tunable properties.
<code>release</code>	Unlock the System object. Use this command to enable setting of nontunable parameters.
<code>reset</code>	Reset the internal states of a locked System object to the initial values, and leave the object locked.

Command	Description
<code>clone</code>	Create another System object with the same object property values.  Do not create additional objects using syntax <code>obj2 = obj</code> . Any changes made to the properties of the new object created this way ( <code>obj2</code> ) also change the properties of the original object ( <code>obj</code> ).
<code>isLocked</code>	Query locked status for input attributes and nontunable properties of the System object.

Use the `recursiveAR` command to create an online estimation System object. Then estimate the AR model parameter, `A`, and output using the `step` command with output data `y`.

```
[A,EstimatedOutput] = step(obj,y)
```

For `recursiveAR` object properties, see “Properties” on page 1-1333.

## Examples

### Estimate AR Model Online

Create a System object™ for online parameter estimation of an AR model using recursive estimation algorithms.

```
obj = recursiveAR;
```

The AR model has a default structure with polynomial of order 1 and initial polynomial coefficient values, `eps`.

Load the time-series estimation data. In this example, use a static data set for illustration.

```
load iddata9 z9;  
output = z9.y;
```

Estimate AR model parameters online using `step`.

```
for i = 1:numel(output)
[A,EstimatedOutput] = step(obj,output(i));
end
```

View the current estimated values of polynomial A coefficients.

```
obj.A
```

```
ans = 1×2
```

```
1.0000 -0.9592
```

View the current covariance estimate of the parameters.

```
obj.ParameterCovariance
```

```
ans = 1.6204e-04
```

View the current estimated output.

```
EstimatedOutput
```

```
EstimatedOutput = 0.7830
```

### **Create Online Estimation System Object for AR Model With Known Polynomial Order**

Specify AR model polynomial order.

```
na = 2;
```

Create a System object™ for online estimation of an AR model with the specified polynomial order.

```
obj = recursiveAR(na);
```

## Create Online Estimation System Object for AR Model With Known Initial Parameters

Specify AR model order.

```
na = 2;
```

Create a System object for online estimation of AR model with known initial polynomial coefficients.

```
A0 = [1 0.5 0.3];  
obj = recursiveAR(na,A0);
```

Specify the initial parameter covariance.

```
obj.InitialParameterCovariance = 0.1;
```

`InitialParameterCovariance` represents the uncertainty in your guess for the initial parameters. Typically, the default `InitialParameterCovariance` (10000) is too large relative to the parameter values. This results in initial guesses being given less importance during estimation. If you have confidence in the initial parameter guesses, specify a smaller initial parameter covariance.

## Specify Estimation Method for Online Estimation of AR Model

Create a System object that uses the normalized gradient algorithm for online parameter estimation of an AR model.

```
obj = recursiveAR(2,'EstimationMethod','NormalizedGradient');
```

## Input Arguments

### **na** — Model order

positive integer

Model order of the polynomial  $A(q)$  of an AR model on page 1-1342, specified as a positive integer.

**A0 — Initial value of polynomial coefficients**

row vector of real values | []

Initial value of coefficients of the polynomial  $A(q)$ , specified as a 1-by-(na+1) row vector of real values with 1 as the first element. Specify the coefficients in order of ascending powers of  $q^{-1}$ .

Specifying as [], uses the default value of eps for the polynomial coefficients.

---

**Note** If the initial parameter values are much smaller than `InitialParameterCovariance`, these initial values are given less importance during estimation. Specify a smaller initial parameter covariance if you have high confidence in the initial parameter values. This statement applies only for infinite-history estimation. Finite-history estimation does not use `InitialParameterCovariance`.

---

**Name-Value Pair Arguments**

Specify optional comma-separated pairs of `Name, Value` arguments. `Name` is the argument name and `Value` is the corresponding value. `Name` must appear inside quotes. You can specify several name and value pair arguments in any order as `Name1, Value1, ..., NameN, ValueN`.

Use `Name, Value` arguments to specify writable properties on page 1-1333 of recursiveAR System object during object creation. For example, `obj = recursiveAR(2, 'EstimationMethod', 'Gradient')` creates a System object to estimate an AR model using the 'Gradient' recursive estimation algorithm.

**Properties**

recursiveAR System object properties consist of read-only and writable properties. The writable properties are tunable and nontunable properties. The nontunable properties cannot be changed when the object is locked, that is, after you use the `step` command.

Use `Name, Value` arguments to specify writable properties of recursiveAR objects during object creation. After object creation, use dot notation to modify the tunable properties.

```
obj = recursiveAR;
obj.ForgettingFactor = 0.99;
```

## A

Estimated coefficients of polynomial  $A(q)$ , returned as a row vector of real values specified in order of ascending powers of  $q^{-1}$ .

A is a read-only property and is initially empty after you create the object. It is populated after you use the `step` command for online parameter estimation.

### **InitialA**

Initial values for the coefficients of polynomial  $A(q)$  of order `na`, specified as a row vector of length `na+1`, with 1 as the first element. Specify the coefficients in order of ascending powers of  $q^{-1}$ .

If the initial guesses are much smaller than the default `InitialParameterCovariance`, 10000, the initial guesses are given less importance during estimation. In that case, specify a smaller initial parameter covariance.

`InitialA` is a tunable property. You can change it when the object is in a locked state.

**Default:** [1 eps]

### **InitialOutputs**

Initial values of the outputs buffer in finite-history estimation, specified as `0` or as a  $(W + na)$ -by-1 vector, where  $W$  is the window length and  $na$  is the polynomial order you specify during object construction.

The `InitialOutputs` property provides a means of controlling the initial behavior of the algorithm.

When `InitialOutputs` is set to `0`, the object populates the buffer with zeros.

If the initial buffer is set to `0` or does not contain enough information, you see a warning message during the initial phase of your estimation. The warning should clear after a few cycles. The number of cycles it takes for sufficient information to be buffered depends upon the order of your polynomials and your input delays. If the warning persists, you should evaluate the content of your signals.

Specify `InitialOutputs` only when `History` is `Finite`.

`InitialOutputs` is a tunable property. You can change `InitialOutputs` when the object is in a locked state.

**Default:** 0

### ParameterCovariance

Estimated covariance  $P$  of the parameters, returned as an  $N$ -by- $N$  symmetric positive-definite matrix.  $N$  is the number of parameters to be estimated. The software computes  $P$  assuming that the residuals (difference between estimated and measured outputs) are white noise, and the variance of these residuals is 1.

ParameterCovariance is applicable only when EstimationMethod is 'ForgettingFactor' or 'KalmanFilter' or when History is Finite.

The interpretation of  $P$  depends on your settings for the History and EstimationMethod properties.

- If History is Infinite, then your EstimationMethod selection results in one of the following:
  - 'ForgettingFactor' —  $(R_2/2)P$  is approximately equal to the covariance matrix of the estimated parameters, where  $R_2$  is the true variance of the residuals.
  - 'KalmanFilter' —  $R_2P$  is the covariance matrix of the estimated parameters, and  $R_1/R_2$  is the covariance matrix of the parameter changes. Here,  $R_1$  is the covariance matrix that you specify in ProcessNoiseCovariance.
- If History is Finite (sliding-window estimation) —  $R_2P$  is the covariance of the estimated parameters. The sliding-window algorithm does not use this covariance in the parameter-estimation process. However, the algorithm does compute the covariance for output so that you can use it for statistical evaluation.

ParameterCovariance is a read-only property and is initially empty after you create the object. It is populated after you use the `step` command for online parameter estimation.

### InitialParameterCovariance

Covariance of the initial parameter estimates, specified as one of the following:

- Real positive scalar,  $\alpha$  — Covariance matrix is an  $N$ -by- $N$  diagonal matrix, with  $\alpha$  as the diagonal elements.  $N$  is the number of parameters to be estimated.
- Vector of real positive scalars,  $[\alpha_1, \dots, \alpha_N]$  — Covariance matrix is an  $N$ -by- $N$  diagonal matrix, with  $[\alpha_1, \dots, \alpha_N]$  as the diagonal elements.
- $N$ -by- $N$  symmetric positive-definite matrix.

`InitialParameterCovariance` represents the uncertainty in the initial parameter estimates. For large values of `InitialParameterCovariance`, less importance is placed on the initial parameter values and more on the measured data during beginning of estimation using `step`.

Use only when `EstimationMethod` is `'ForgettingFactor'` or `'KalmanFilter'`.

`InitialParameterCovariance` is a tunable property. You can change it when the object is in a locked state.

**Default:** 10000

## **EstimationMethod**

Recursive estimation algorithm used for online estimation of model parameters, specified as one of the following values:

- `'ForgettingFactor'` — Algorithm used for parameter estimation
- `'KalmanFilter'` — Algorithm used for parameter estimation
- `'NormalizedGradient'` — Algorithm used for parameter estimation
- `'Gradient'` — Unnormalized gradient algorithm used for parameter estimation

Forgetting factor and Kalman filter algorithms are more computationally intensive than gradient and unnormalized gradient methods. However, they have better convergence properties. For information about these algorithms, see “Recursive Algorithms for Online Parameter Estimation”.

These methods all use an infinite data history, and are available only when `History` is `'Infinite'`.

`EstimationMethod` is a nontunable property. You cannot change it during execution, that is, after the object is locked using the `step` command.

**Default:** Forgetting Factor

## **ForgettingFactor**

Forgetting factor,  $\lambda$ , relevant for parameter estimation, specified as a scalar in the range (0,1].

Suppose that the system remains approximately constant over  $T_0$  samples. You can choose  $\lambda$  such that:

$$T_0 = \frac{1}{1-\lambda}$$

- Setting  $\lambda = 1$  corresponds to “no forgetting” and estimating constant coefficients.
- Setting  $\lambda < 1$  implies that past measurements are less significant for parameter estimation and can be “forgotten”. Set  $\lambda < 1$  to estimate time-varying coefficients.

Typical choices of  $\lambda$  are in the range [0.98 0.995].

Use only when EstimationMethod is 'ForgettingFactor'.

ForgettingFactor is a tunable property. You can change it when the object is in a locked state.

**Default:** 1

### EnableAdapation

Enable or disable parameter estimation, specified as one of the following:

- `true` or `1`— The `step` command estimates the parameter values for that time step and updates the parameter values.
- `false` or `0` — The `step` command does not update the parameters for that time step and instead outputs the last estimated value. You can use this option when your system enters a mode where the parameter values do not vary with time.

---

**Note** If you set `EnableAdapation` to `false`, you must still execute the `step` command. Do not skip `step` to keep parameter values constant, because parameter estimation depends on current and past I/O measurements. `step` ensures past I/O data is stored, even when it does not update the parameters.

---

EnableAdapation is a tunable property. You can change it when the object is in a locked state.

**Default:** `true`

### Data Type

Floating point precision of parameters, specified as one of the following values:

- 'double' — Double-precision floating point
- 'single' — Single-precision floating point

Setting `DataType` to 'single' saves memory, but leads to loss of precision. Specify `DataType` based on the precision required by the target processor where you will deploy generated code.

`DataType` is a nontunable property. It can only be set during object construction using `Name, Value` arguments and cannot be changed afterward.

**Default:** 'double'

### **ProcessNoiseCovariance**

Covariance matrix of parameter variations, specified as one of the following:

- Real nonnegative scalar,  $\alpha$  — Covariance matrix is an  $N$ -by- $N$  diagonal matrix, with  $\alpha$  as the diagonal elements.
- Vector of real nonnegative scalars,  $[\alpha_1, \dots, \alpha_N]$  — Covariance matrix is an  $N$ -by- $N$  diagonal matrix, with  $[\alpha_1, \dots, \alpha_N]$  as the diagonal elements.
- $N$ -by- $N$  symmetric positive semidefinite matrix.

$N$  is the number of parameters to be estimated.

`ProcessNoiseCovariance` is applicable when `EstimationMethod` is 'KalmanFilter'.

Kalman filter algorithm treats the parameters as states of a dynamic system and estimates these parameters using a Kalman filter. `ProcessNoiseCovariance` is the covariance of the process noise acting on these parameters. Zero values in the noise covariance matrix correspond to estimating constant coefficients. Values larger than 0 correspond to time-varying parameters. Use large values for rapidly changing parameters. However, the larger values result in noisier parameter estimates.

`ProcessNoiseCovariance` is a tunable property. You can change it when the object is in a locked state.

**Default:** 0.1

## AdaptationGain

Adaptation gain,  $\gamma$ , used in gradient recursive estimation algorithms, specified as a positive scalar.

AdaptationGain is applicable when EstimationMethod is 'Gradient' or 'NormalizedGradient'.

Specify a large value for AdaptationGain when your measurements have a high signal-to-noise ratio.

AdaptationGain is a tunable property. You can change it when the object is in a locked state.

**Default:** 1

## NormalizationBias

Bias in adaptation gain scaling used in the 'NormalizedGradient' method, specified as a nonnegative scalar.

NormalizationBias is applicable when EstimationMethod is 'NormalizedGradient'.

The normalized gradient algorithm divides the adaptation gain at each step by the square of the two-norm of the gradient vector. If the gradient is close to zero, this can cause jumps in the estimated parameters. NormalizationBias is the term introduced in the denominator to prevent these jumps. Increase NormalizationBias if you observe jumps in estimated parameters.

NormalizationBias is a tunable property. You can change it when the object is in a locked state.

**Default:** eps

## History

Data history type defining which type of recursive algorithm you use, specified as:

- 'Infinite' — Use an algorithm that aims to minimize the error between the observed and predicted outputs for all time steps from the beginning of the simulation.
- 'Finite' — Use an algorithm that aims to minimize the error between the observed and predicted outputs for a finite number of past time steps.

Algorithms with infinite history aim to produce parameter estimates that explain all data since the start of the simulation. These algorithms still use a fixed amount of memory that does not grow over time. The object provides multiple algorithms of the 'Infinite' History type. Specifying this option activates the EstimationMethod property with which you specify an algorithm.

Algorithms with finite history aim to produce parameter estimates that explain only a finite number of past data samples. This method is also called sliding-window estimation. The object provides one algorithm of the 'Finite' type. Specifying this option activates the WindowLength property that sizes the window.

For more information on recursive estimation methods, see “Recursive Algorithms for Online Parameter Estimation”.

History is a nontunable property. It can be set only during object construction using Name, Value arguments and cannot be changed afterward.

**Default:** 'Infinite'

### **WindowLength**

Window size determining the number of time samples to use for the sliding-window estimation method, specified as a positive integer. Specify WindowLength only when History is Finite.

Choose a window size that balances estimation performance with computational and memory burden. Sizing factors include the number and time variance of the parameters in your model. Always specify **Window Length** in samples, even if you are using frame-based input processing.

WindowLength must be greater than or equal to the number of estimated parameters.

Suitable window length is independent of whether you are using sample-based or frame-based input processing (see InputProcessing). However, when using frame-based processing, your window length must be greater than or equal to the number of samples (time steps) contained in the frame.

WindowLength is a nontunable property. It can be set only during object construction using Name, Value arguments and cannot be changed afterward.

**Default:** 200

## InputProcessing

Option for sample-based or frame-based input processing, specified as a character vector or string.

- **Sample-based** processing operates on signals streamed one sample at a time.
- **Frame-based** processing operates on signals containing samples from multiple time steps. Many machine sensor interfaces package multiple samples and transmit these samples together in frames. **Frame-based** processing allows you to input this data directly without having to first unpack it.

Your `InputProcessing` specification impacts the dimensions for the input and output signals when using the `step` command:

```
[theta,EstimatedOutput] = step(obj,y)
```

- **Sample-based**
  - `y` and `EstimatedOutput` are scalars.
- **Frame-based** with  $M$  samples per frame
  - `y` and `EstimatedOutput` are  $M$ -by-1 vectors.

`InputProcessing` is a nontunable property. It can be set only during object construction using `Name, Value` arguments and cannot be changed afterward.

**Default:** 'Sample-based'

## Output Arguments

### **obj** — System object for online parameter estimation of AR model

recursiveAR System object

System object for online parameter estimation of AR model, returned as a `recursiveAR` System object. This object is created using the specified model orders and properties. Use `step` command to estimate the coefficients of the AR model polynomials. You can then access the estimated coefficients and parameter covariance using dot notation. For example, type `obj.A` to view the estimated  $A$  polynomial coefficients.

## Definitions

### AR Model Structure

The AR model structure is:

$$A(q)y(t) = e(t)$$

where,

$$A(q) = 1 + \alpha_1 q^{-1} + \dots + \alpha_{n_a} q^{-n_a}$$

Here,

- $y(t)$ — Output at time  $t$ . Data is a time series that has no input channels and one output channel.
- $n_a$  — Number of  $A$  polynomial coefficients.
- $e(t)$  — White-noise disturbance value at time  $t$ .
- $q^{-1}$  — Time-shift operator.

## Tips

- Starting in R2016b, instead of using the `step` command to update model parameter estimates, you can call the System object with input arguments, as if it were a function. For example, `[A,EstimatedOutput] = step(obj,y)` and `[A,EstimatedOutput] = obj(y)` perform equivalent operations.

## Extended Capabilities

### C/C++ Code Generation

Generate C and C++ code using MATLAB® Coder™.

Usage notes and limitations:

- For Simulink based workflows, use Recursive Polynomial Model Estimator.
- For limitations, see “Generate Code for Online Parameter Estimation in MATLAB”.

Supports MATLAB Function block: No

## See Also

Recursive Polynomial Model Estimator | ar | clone | isLocked | recursiveARMA | recursiveARMAX | recursiveARX | recursiveBJ | recursiveLS | recursiveOE | release | reset | step

## Topics

“Perform Online Parameter Estimation at the Command Line”

“Validate Online Parameter Estimation at the Command Line”

“What Is Online Estimation?”

“Recursive Algorithms for Online Parameter Estimation”

**Introduced in R2015b**

## recursiveARMA

Create System object for online parameter estimation of ARMA model

Use `recursiveARMA` command for parameter estimation with real-time data. If all data necessary for estimation is available at once, and you are estimating a time-invariant model, use the offline estimation command, `armax`.

### Syntax

```
obj = recursiveARMA
obj = recursiveARMA(Orders)
obj = recursiveARMA(Orders,A0,C0)
obj = recursiveARMA( ____,Name,Value)
```

### Description

`obj = recursiveARMA` creates a System object for online parameter estimation of a default single output ARMA model structure on page 1-1356. The default model structure has polynomials of order 1 and initial polynomial coefficient values `eps`.

After creating the object, use the `step` command to update model parameter estimates using recursive estimation algorithms and real-time data.

`obj = recursiveARMA(Orders)` specifies the polynomial orders of the ARMA model to be estimated.

`obj = recursiveARMA(Orders,A0,C0)` specifies the polynomial orders and initial values of the polynomial coefficients. Specify initial values to potentially avoid local minima during estimation. If the initial values are small compared to the default `InitialParameterCovariance` property value, and you have confidence in your initial values, also specify a smaller `InitialParameterCovariance`.

`obj = recursiveARMA( ____,Name,Value)` specifies additional attributes of the ARMA model structure and recursive estimation algorithm using one or more `Name,Value` pair arguments.

## Object Description

recursiveARMA creates a System object for online parameter estimation of single output ARMA models using a recursive estimation algorithm.

A System object is a specialized MATLAB object designed specifically for implementing and simulating dynamic systems with inputs that change over time. System objects use internal states to store past behavior, which is used in the next computational step.

After you create a System object, you use commands to process data or obtain information from or about the object. System objects use a minimum of two commands to process data — a constructor to create the object and the `step` command to update object parameters using real-time data. This separation of declaration from execution lets you create multiple, persistent, reusable objects, each with different settings.

You can use the following commands with the online estimation System objects in System Identification Toolbox:

Command	Description
<code>step</code>	Update model parameter estimates using recursive estimation algorithms and real-time data.  <code>step</code> puts the object into a locked state. In a locked state, you cannot change any nontunable properties or input specifications, such as model order, data type, or estimation algorithm. During execution, you can only change tunable properties.
<code>release</code>	Unlock the System object. Use this command to enable setting of nontunable parameters.
<code>reset</code>	Reset the internal states of a locked System object to the initial values, and leave the object locked.

Command	Description
<code>clone</code>	Create another System object with the same object property values.  Do not create additional objects using syntax <code>obj2 = obj</code> . Any changes made to the properties of the new object created this way ( <code>obj2</code> ) also change the properties of the original object ( <code>obj</code> ).
<code>isLocked</code>	Query locked status for input attributes and nontunable properties of the System object.

Use the `recursiveARMA` command to create an online estimation System object. Then estimate the ARMA model parameters (A and C) and output using the `step` command with output data `y`.

```
[A,C,EstimatedOutput] = step(obj,y)
```

For `recursiveARMA` object properties, see “Properties” on page 1-1350.

## Examples

### Estimate ARMA Model Online

Create a System object for online parameter estimation of an ARMA model.

```
obj = recursiveARMA;
```

The ARMA model has a default structure with polynomials of order 1 and initial polynomial coefficient values, `eps`.

Load the time-series estimation data. In this example, use a static data set for illustration.

```
load iddata9 z9;  
output = z9.y;
```

Estimate ARMA model parameters online using `step`.

```
for i = 1:numel(output)
[A,C,EstimatedOutput] = step(obj,output(i));
end
```

View the current estimated values of polynomial C coefficients.

```
obj.C
```

```
ans = 1×2
```

```
1.0000    0.2315
```

View the current covariance estimate of the parameters.

```
obj.ParameterCovariance
```

```
ans = 2×2
```

```
10-3 ×
```

```
0.6372    -0.0257
-0.0257    0.0017
```

View the current estimated output.

```
EstimatedOutput
```

```
EstimatedOutput = 11.8121
```

### Create Online Estimation System Object for ARMA Model With Known Orders

Specify ARMA model orders.

```
na = 2;
```

```
nc = 1;
```

Create a System object for online estimation of an ARMA model with the specified orders.

```
obj = recursiveARMA([na nc]);
```

## Create Online Estimation System Object for ARMA Model With Known Initial Parameters

Specify ARMA model orders.

```
na = 2;  
nc = 1;
```

Create a System object for online estimation of ARMA model with known initial polynomial coefficients.

```
A0 = [1 0.5 0.3];  
C0 = [1 0.7];  
obj = recursiveARMA([na nc],A0,C0);
```

Specify the initial parameter covariance.

```
obj.InitialParameterCovariance = 0.1;
```

`InitialParameterCovariance` represents the uncertainty in your guess for the initial parameters. Typically, the default `InitialParameterCovariance` (10000) is too large relative to the parameter values. This results in initial guesses being given less importance during estimation. If you have confidence in the initial parameter guesses, specify a smaller initial parameter covariance.

## Specify Estimation Method for Online Estimation of ARMA Model

Create a System object that uses the unnormalized gradient algorithm for online parameter estimation of an ARMA model.

```
obj = recursiveARMA([2 1], 'EstimationMethod', 'Gradient');
```

## Input Arguments

### Orders — Model orders

1-by-2 vector of integers

Model orders of an ARMA model on page 1-1356, specified as a 1-by-2 vector of integers, `[na nc]`.

- `na` — Order of the polynomial  $A(q)$ , specified as a nonnegative integer.
- `nc` — Order of the polynomial  $C(q)$ , specified as a nonnegative integer.

### **A0, C0 — Initial value of polynomial coefficients**

row vectors of real values | []

Initial value of polynomial coefficients, specified as row vectors of real values with elements in order of ascending powers of  $q^{-1}$ .

- `A0` — Initial guess for the coefficients of the polynomial  $A(q)$ , specified as a 1-by-(`na`+1) vector with 1 as the first element.
- `C0` — Initial guess for the coefficients of the polynomial  $C(q)$ , specified as a 1-by-(`nc`+1) vector with 1 as the first element.

The coefficients in `C0` must define a stable discrete-time polynomial with roots within a unit disk. For example,

```
C0 = [1 0.5 0.5];
all(abs(roots(C0))<1)
```

```
ans =
```

```
1
```

Specifying as [], uses the default value of `eps` for the polynomial coefficients.

---

**Note** If the initial guesses are much smaller than the default `InitialParameterCovariance`, 10000, the initial guesses are given less importance during estimation. In that case, specify a smaller initial parameter covariance.

---

## **Name-Value Pair Arguments**

Specify optional comma-separated pairs of `Name`, `Value` arguments. `Name` is the argument name and `Value` is the corresponding value. `Name` must appear inside quotes. You can specify several name and value pair arguments in any order as `Name1, Value1, ..., NameN, ValueN`.

Use `Name`, `Value` arguments to specify writable properties on page 1-1350 of recursiveARMA System object during object creation. For example, `obj = recursiveARMA([2 1], 'EstimationMethod', 'Gradient')` creates a System object to estimate an ARMA model using the 'Gradient' recursive estimation algorithm.

## Properties

recursiveARMA System object properties consist of read-only and writable properties. The writable properties are tunable and nontunable properties. The nontunable properties cannot be changed when the object is locked, that is, after you use the `step` command.

Use `Name`, `Value` arguments to specify writable properties of recursiveARMA objects during object creation. After object creation, use dot notation to modify the tunable properties.

```
obj = recursiveARMA;  
obj.ForgettingFactor = 0.99;
```

### A

Estimated coefficients of polynomial  $A(q)$ , returned as a row vector of real values specified in order of ascending powers of  $q^{-1}$ .

A is a read-only property and is initially empty after you create the object. It is populated after you use the `step` command for online parameter estimation.

### C

Estimated coefficients of polynomial  $C(q)$ , returned as a vector of real values specified in order of ascending powers of  $q^{-1}$ .

C is a read-only property and is initially empty after you create the object. It is populated after you use the `step` command for online parameter estimation.

### InitialA

Initial values for the coefficients of polynomial  $A(q)$  of order `na`, specified as a row vector of length `na+1`, with 1 as the first element. Specify the coefficients in order of ascending powers of  $q^{-1}$ .

If the initial guesses are much smaller than the default `InitialParameterCovariance`, 10000, the initial guesses are given less importance during estimation. In that case, specify a smaller initial parameter covariance.

`InitialA` is a tunable property. You can change it when the object is in a locked state.

**Default:** [1 eps]

## InitialC

Initial values for the coefficients of polynomial  $C(q)$  of order  $nc$ , specified as a row vector of length  $nc+1$ , with 1 as the first element. Specify the coefficients in order of ascending powers of  $q^{-1}$ .

The coefficients in `InitialC` must define a stable discrete-time polynomial with roots within a unit circle. For example,

```
InitialC = [1 0.5 0.5];
all(abs(roots(InitialC))<1)
```

```
ans =
```

```
1
```

If the initial guesses are much smaller than the default `InitialParameterCovariance`, 10000, the initial guesses are given less importance during estimation. In that case, specify a smaller initial parameter covariance.

`InitialC` is a tunable property. You can change it when the object is in a locked state.

**Default:** [1 eps]

## ParameterCovariance

Estimated covariance  $P$  of the parameters, returned as an  $N$ -by- $N$  symmetric positive-definite matrix.  $N$  is the number of parameters to be estimated. The software computes  $P$  assuming that the residuals (difference between estimated and measured outputs) are white noise, and the variance of these residuals is 1. `ParameterCovariance` is applicable only when `EstimationMethod` is 'ForgettingFactor' or 'KalmanFilter'.

The interpretation of  $P$  depends on the estimation method:

- 'ForgettingFactor' —  $R_2/2 * P$  is approximately equal to the covariance matrix of the estimated parameters, where  $R_2$  is the true variance of the residuals.
- 'KalmanFilter' —  $R_2 * P$  is the covariance matrix of the estimated parameters, and  $R_1 / R_2$  is the covariance matrix of the parameter changes. Where,  $R_1$  is the covariance matrix that you specify in `ProcessNoiseCovariance`.

`ParameterCovariance` is a read-only property and is initially empty after you create the object. It is populated after you use the `step` command for online parameter estimation.

## **InitialParameterCovariance**

Covariance of the initial parameter estimates, specified as one of the following:

- Real positive scalar,  $\alpha$  — Covariance matrix is an  $N$ -by- $N$  diagonal matrix, with  $\alpha$  as the diagonal elements.  $N$  is the number of parameters to be estimated.
- Vector of real positive scalars,  $[\alpha_1, \dots, \alpha_N]$  — Covariance matrix is an  $N$ -by- $N$  diagonal matrix, with  $[\alpha_1, \dots, \alpha_N]$  as the diagonal elements.
- $N$ -by- $N$  symmetric positive-definite matrix.

`InitialParameterCovariance` represents the uncertainty in the initial parameter estimates. For large values of `InitialParameterCovariance`, less importance is placed on the initial parameter values and more on the measured data during beginning of estimation using `step`.

Use only when `EstimationMethod` is 'ForgettingFactor' or 'KalmanFilter'.

`InitialParameterCovariance` is a tunable property. You can change it when the object is in a locked state.

**Default:** 10000

## **EstimationMethod**

Recursive estimation algorithm used for online estimation of model parameters, specified as one of the following values:

- 'ForgettingFactor' — Algorithm used for parameter estimation
- 'KalmanFilter' — Algorithm used for parameter estimation
- 'NormalizedGradient' — Algorithm used for parameter estimation
- 'Gradient' — Unnormalized gradient algorithm used for parameter estimation

Forgetting factor and Kalman filter algorithms are more computationally intensive than gradient and unnormalized gradient methods. However, they have better convergence properties. For information about these algorithms, see “Recursive Algorithms for Online Parameter Estimation”.

`EstimationMethod` is a nontunable property. You cannot change it during execution, that is after the object is locked using the `step` command. If you want to deploy code using MATLAB Coder, `EstimationMethod` can only be assigned once.

**Default:** 'ForgettingFactor'

### ForgettingFactor

Forgetting factor,  $\lambda$ , relevant for parameter estimation, specified as a scalar in the range (0,1].

Suppose that the system remains approximately constant over  $T_0$  samples. You can choose  $\lambda$  such that:

$$T_0 = \frac{1}{1-\lambda}$$

- Setting  $\lambda = 1$  corresponds to “no forgetting” and estimating constant coefficients.
- Setting  $\lambda < 1$  implies that past measurements are less significant for parameter estimation and can be “forgotten”. Set  $\lambda < 1$  to estimate time-varying coefficients.

Typical choices of  $\lambda$  are in the range [0.98 0.995].

Use only when EstimationMethod is 'ForgettingFactor'.

ForgettingFactor is a tunable property. You can change it when the object is in a locked state.

**Default:** 1

### EnableAdaption

Enable or disable parameter estimation, specified as one of the following:

- `true` or `1`— The `step` command estimates the parameter values for that time step and updates the parameter values.
- `false` or `0` — The `step` command does not update the parameters for that time step and instead outputs the last estimated value. You can use this option when your system enters a mode where the parameter values do not vary with time.

---

**Note** If you set `EnableAdaption` to `false`, you must still execute the `step` command. Do not skip `step` to keep parameter values constant, because parameter estimation depends on current and past I/O measurements. `step` ensures past I/O data is stored, even when it does not update the parameters.

---

EnableAdapation is a tunable property. You can change it when the object is in a locked state.

**Default:** true

## **DataType**

Floating point precision of parameters, specified as one of the following values:

- 'double' — Double-precision floating point
- 'single' — Single-precision floating point

Setting `DataType` to 'single' saves memory, but leads to loss of precision. Specify `DataType` based on the precision required by the target processor where you will deploy generated code.

`DataType` is a nontunable property. It can only be set during object construction using `Name, Value` arguments and cannot be changed afterward.

**Default:** 'double'

## **ProcessNoiseCovariance**

Covariance matrix of parameter variations, specified as one of the following:

- Real nonnegative scalar,  $\alpha$  — Covariance matrix is an  $N$ -by- $N$  diagonal matrix, with  $\alpha$  as the diagonal elements.
- Vector of real nonnegative scalars,  $[\alpha_1, \dots, \alpha_N]$  — Covariance matrix is an  $N$ -by- $N$  diagonal matrix, with  $[\alpha_1, \dots, \alpha_N]$  as the diagonal elements.
- $N$ -by- $N$  symmetric positive semidefinite matrix.

$N$  is the number of parameters to be estimated.

`ProcessNoiseCovariance` is applicable when `EstimationMethod` is 'KalmanFilter'.

Kalman filter algorithm treats the parameters as states of a dynamic system and estimates these parameters using a Kalman filter. `ProcessNoiseCovariance` is the covariance of the process noise acting on these parameters. Zero values in the noise covariance matrix correspond to estimating constant coefficients. Values larger than 0 correspond to time-varying parameters. Use large values for rapidly changing parameters. However, the larger values result in noisier parameter estimates.

ProcessNoiseCovariance is a tunable property. You can change it when the object is in a locked state.

**Default:** 0.1

### **AdaptationGain**

Adaptation gain,  $\gamma$ , used in gradient recursive estimation algorithms, specified as a positive scalar.

AdaptationGain is applicable when EstimationMethod is 'Gradient' or 'NormalizedGradient'.

Specify a large value for AdaptationGain when your measurements have a high signal-to-noise ratio.

AdaptationGain is a tunable property. You can change it when the object is in a locked state.

**Default:** 1

### **NormalizationBias**

Bias in adaptation gain scaling used in the 'NormalizedGradient' method, specified as a nonnegative scalar.

NormalizationBias is applicable when EstimationMethod is 'NormalizedGradient'.

The normalized gradient algorithm divides the adaptation gain at each step by the square of the two-norm of the gradient vector. If the gradient is close to zero, this can cause jumps in the estimated parameters. NormalizationBias is the term introduced in the denominator to prevent these jumps. Increase NormalizationBias if you observe jumps in estimated parameters.

NormalizationBias is a tunable property. You can change it when the object is in a locked state.

**Default:** eps

## Output Arguments

### **obj** — System object for online parameter estimation of ARMA model

`recursiveARMA` System object

System object for online parameter estimation of ARMA model, returned as a `recursiveARMA` System object. This object is created using the specified model orders and properties. Use `step` command to estimate the coefficients of the ARMA model polynomials. You can then access the estimated coefficients and parameter covariance using dot notation. For example, type `obj.A` to view the estimated *A* polynomial coefficients.

## Definitions

### ARMA Model Structure

The ARMA model structure is:

$$A(q)y(t) = C(q)e(t)$$

where,

$$A(q) = 1 + a_1q^{-1} + \dots + a_{n_a}q^{-n_a}$$

$$C(q) = 1 + c_1q^{-1} + \dots + c_{n_c}q^{-n_c}$$

Here,

- $y(t)$ — Output at time  $t$ . Data is a time series that has no input channels and one output channel.
- $n_a$  — Number of *A* polynomial coefficients
- $n_c$  — Number of *C* polynomial coefficients
- $e(t)$  — White-noise disturbance value at time  $t$
- $q^{-1}$  — Time-shift operator

## Tips

- Starting in R2016b, instead of using the `step` command to update model parameter estimates, you can call the System object with input arguments, as if it were a function. For example, `[A,C,EstimatedOutput] = step(obj,y)` and `[A,C,EstimatedOutput] = obj(y)` perform equivalent operations.

## Extended Capabilities

### C/C++ Code Generation

Generate C and C++ code using MATLAB® Coder™.

Usage notes and limitations:

- For Simulink-based workflows, use Recursive Polynomial Model Estimator.
- For limitations, see “Generate Code for Online Parameter Estimation in MATLAB”.

Supports MATLAB Function block: No

### See Also

[Recursive Polynomial Model Estimator](#) | [armax](#) | [clone](#) | [isLocked](#) | [recursiveAR](#) | [recursiveARMAX](#) | [recursiveARX](#) | [recursiveBJ](#) | [recursiveLS](#) | [recursiveOE](#) | [release](#) | [reset](#) | [step](#)

### Topics

“Perform Online Parameter Estimation at the Command Line”

“Validate Online Parameter Estimation at the Command Line”

“What Is Online Estimation?”

“Recursive Algorithms for Online Parameter Estimation”

**Introduced in R2015b**

## recursiveARMAX

Create System object for online parameter estimation of ARMAX model

Use `recursiveARMAX` command for parameter estimation with real-time data. If all data necessary for estimation is available at once, and you are estimating a time-invariant model, use the offline estimation command, `armax`.

### Syntax

```
obj = recursiveARMAX  
obj = recursiveARMAX(Orders)  
obj = recursiveARMAX(Orders,A0,B0,C0)  
obj = recursiveARMAX( ____,Name,Value)
```

### Description

`obj = recursiveARMAX` creates a System object for online parameter estimation of default single-input single-output (SISO) ARMAX model structure on page 1-1371. The default model structure has polynomials of order 1 and initial polynomial coefficient values `eps`.

After creating the object, use the `step` command to update model parameter estimates using recursive estimation algorithms and real-time data.

`obj = recursiveARMAX(Orders)` specifies the polynomial orders of the ARMAX model to be estimated.

`obj = recursiveARMAX(Orders,A0,B0,C0)` specifies the polynomial orders and initial values of the polynomial coefficients. Specify initial values to potentially avoid local minima during estimation. If the initial values are small compared to the default `InitialParameterCovariance` property value, and you have confidence in your initial values, also specify a smaller `InitialParameterCovariance`.

`obj = recursiveARMAX( ____,Name,Value)` specifies additional attributes of the ARMAX model structure and recursive estimation algorithm using one or more `Name,Value` pair arguments.

## Object Description

recursiveARMAX creates a System object for online parameter estimation of SISO ARMAX models using a recursive estimation algorithm.

A System object is a specialized MATLAB object designed specifically for implementing and simulating dynamic systems with inputs that change over time. System objects use internal states to store past behavior, which is used in the next computational step.

After you create a System object, you use commands to process data or obtain information from or about the object. System objects use a minimum of two commands to process data — a constructor to create the object and the `step` command to update object parameters using real-time data. This separation of declaration from execution lets you create multiple, persistent, reusable objects, each with different settings.

You can use the following commands with the online estimation System objects in System Identification Toolbox:

Command	Description
<code>step</code>	Update model parameter estimates using recursive estimation algorithms and real-time data.  <code>step</code> puts the object into a locked state. In a locked state, you cannot change any nontunable properties or input specifications, such as model order, data type, or estimation algorithm. During execution, you can only change tunable properties.
<code>release</code>	Unlock the System object. Use this command to enable setting of nontunable parameters.
<code>reset</code>	Reset the internal states of a locked System object to the initial values, and leave the object locked.

Command	Description
<code>clone</code>	Create another System object with the same object property values.  Do not create additional objects using syntax <code>obj2 = obj</code> . Any changes made to the properties of the new object created this way ( <code>obj2</code> ) also change the properties of the original object ( <code>obj</code> ).
<code>isLocked</code>	Query locked status for input attributes and nontunable properties of the System object.

Use the `recursiveARMAX` command to create an online estimation System object. Then estimate the ARMAX model parameters (A, B, and C) and output using the `step` command with incoming input and output data, `u`, and `y`.

```
[A,B,C,EstimatedOutput] = step(obj,y,u)
```

For `recursiveARMAX` object properties, see “Properties” on page 1-1364.

## Examples

### Estimate an ARMAX Model Online

Create a System object for online parameter estimation of an ARMAX model.

```
obj = recursiveARMAX;
```

The ARMAX model has a default structure with polynomials of order 1 and initial polynomial coefficient values, `eps`.

Load the estimation data. In this example, use a static data set for illustration.

```
load iddata1 z1;  
output = z1.y;  
input = z1.u;
```

Estimate ARMAX model parameters online using `step`.

```

for i = 1:numel(input)
[A,B,C,EstimatedOutput] = step(obj,output(i),input(i));
end

```

View the current estimated values of polynomial A coefficients.

```
obj.A
```

```
ans = 1x2
```

```
1.0000 -0.8298
```

View the current covariance estimate of the parameters.

```
obj.ParameterCovariance
```

```
ans = 3x3
```

```

0.0001    0.0001    0.0001
0.0001    0.0032    0.0000
0.0001    0.0000    0.0001

```

View the current estimated output.

```
EstimatedOutput
```

```
EstimatedOutput = -4.5595
```

### Create System Object for ARMAX Model With Known Polynomial Orders

Specify ARMAX model orders and delays.

```

na = 1;
nb = 2;
nc = 1;
nk = 1;

```

Create a System object for online estimation of ARMAX model with the specified orders and delays.

```
obj = recursiveARMAX([na nb nc nk]);
```

## Create Online Estimation System Object for ARMAX Model With Known Initial Parameters

Specify ARMAX model orders and delays.

```
na = 1;  
nb = 2;  
nc = 1;  
nk = 1;
```

Create a System object for online estimation of ARMAX model with known initial polynomial coefficients.

```
A0 = [1 0.5];  
B0 = [0 1 1];  
C0 = [1 0.5];  
obj = recursiveARMAX([na nb nc nk],A0,B0,C0);
```

Specify the initial parameter covariance.

```
obj.InitialParameterCovariance = 0.1;
```

`InitialParameterCovariance` represents the uncertainty in your guess for the initial parameters. Typically, the default `InitialParameterCovariance` (10000) is too large relative to the parameter values. This results in initial guesses being given less importance during estimation. If you have confidence in the initial parameter guesses, specify a smaller initial parameter covariance.

## Specify Estimation Method for Online Estimation of ARMAX Model

Create a System object that uses the Kalman filter algorithm for online parameter estimation of an ARMAX model.

```
obj = recursiveARMAX([1 2 1 1], 'EstimationMethod', 'KalmanFilter');  
obj.ProcessNoiseCovariance = 0.01;
```

The `ProcessNoiseCovariance` property of `obj` is applicable only when the Kalman filter algorithm is used for estimation.

## Input Arguments

### Orders — Model orders and delays

1-by-4 vector of integers

Model orders and delays of an ARMAX model on page 1-1371, specified as a 1-by-4 vector of integers, [na nb nc nk].

- **na** — Order of the polynomial  $A(q)$ , specified as a nonnegative integer. na represents the number of poles in your system.
- **nb** — Order of the polynomial  $B(q) + 1$ , specified as a positive integer. nb represents the number of zeroes in your system plus 1.
- **nc** — Order of the polynomial  $C(q)$ , specified as a nonnegative integer.
- **nk** — Input-output delay, specified as a nonnegative integer. nk is number of input samples that occur before the input affects the output. nk is expressed as fixed leading zeros of the  $B$  polynomial.

### A0, B0, C0 — Initial value of polynomial coefficients

row vectors of real values | []

Initial value of polynomial coefficients, specified as row vectors of real values with elements in order of ascending powers of  $q^{-1}$ .

- **A0** — Initial guess for the coefficients of the polynomial  $A(q)$ , specified as a 1-by-(na +1) vector with 1 as the first element.
- **B0** — Initial guess for the coefficients of the polynomial  $B(q)$ , specified as a 1-by-(nb +nk) vector with nk leading zeros.
- **C0** — Initial guess for the coefficients of the polynomial  $C(q)$ , specified as a 1-by-(nc +1) vector with 1 as the first element.

The coefficients in C0 must define a stable discrete-time polynomial with roots within a unit disk. For example,

```
C0 = [1 0.5 0.5];
all(abs(roots(C0))<1)
```

```
ans =
```

```
1
```

Specifying as [], uses the default value of eps for the polynomial coefficients.

---

**Note** If the initial guesses are much smaller than the default `InitialParameterCovariance`, 10000, the initial guesses are given less importance during estimation. In that case, specify a smaller initial parameter covariance.

---

## Name-Value Pair Arguments

Specify optional comma-separated pairs of `Name`, `Value` arguments. `Name` is the argument name and `Value` is the corresponding value. `Name` must appear inside quotes. You can specify several name and value pair arguments in any order as `Name1, Value1, ..., NameN, ValueN`.

Use `Name`, `Value` arguments to specify writable properties on page 1-1364 of `recursiveARMAX` System object during object creation. For example, `obj = recursiveARMAX([2 2 1 1], 'EstimationMethod', 'Gradient')` creates a System object to estimate an ARMAX model using the 'Gradient' recursive estimation algorithm.

## Properties

`recursiveARMAX` System object properties consist of read-only and writable properties. The writable properties are tunable and nontunable properties. The nontunable properties cannot be changed when the object is locked, that is, after you use the `step` command.

Use `Name`, `Value` arguments to specify writable properties of `recursiveARMAX` objects during object creation. After object creation, use dot notation to modify the tunable properties.

```
obj = recursiveARMAX;  
obj.ForgettingFactor = 0.99;
```

### A

Estimated coefficients of polynomial  $A(q)$ , returned as a row vector of real values specified in order of ascending powers of  $q^{-1}$ .

`A` is a read-only property and is initially empty after you create the object. It is populated after you use the `step` command for online parameter estimation.

**B**

Estimated coefficients of polynomial  $B(q)$ , returned as a vector of real values specified in order of ascending powers of  $q^{-1}$ .

**B** is a read-only property and is initially empty after you create the object. It is populated after you use the `step` command for online parameter estimation.

**C**

Estimated coefficients of polynomial  $C(q)$ , returned as a vector of real values specified in order of ascending powers of  $q^{-1}$ .

**C** is a read-only property and is initially empty after you create the object. It is populated after you use the `step` command for online parameter estimation.

**InitialA**

Initial values for the coefficients of polynomial  $A(q)$  of order `na`, specified as a row vector of length `na+1`, with 1 as the first element. Specify the coefficients in order of ascending powers of  $q^{-1}$ .

If the initial guesses are much smaller than the default `InitialParameterCovariance`, 10000, the initial guesses are given less importance during estimation. In that case, specify a smaller initial parameter covariance.

`InitialA` is a tunable property. You can change it when the object is in a locked state.

**Default:** [1 eps]

**InitialB**

Initial values for the coefficients of polynomial  $B(q)$  of order `nb-1`, specified as a row vector of length `nb+nk`, with `nk` leading zeros. `nk` is the input-output delay. Specify the coefficients in order of ascending powers of  $q^{-1}$ .

If the initial guesses are much smaller than the default `InitialParameterCovariance`, 10000, the initial guesses are given less importance during estimation. In that case, specify a smaller initial parameter covariance.

`InitialB` is a tunable property. You can change it when the object is in a locked state.

**Default:** [0 eps]

## InitialC

Initial values for the coefficients of polynomial  $C(q)$  of order  $nc$ , specified as a row vector of length  $nc+1$ , with 1 as the first element. Specify the coefficients in order of ascending powers of  $q^{-1}$ .

The coefficients in `InitialC` must define a stable discrete-time polynomial with roots within a unit circle. For example,

```
InitialC = [1 0.5 0.5];  
all(abs(roots(InitialC))<1)
```

```
ans =
```

```
1
```

If the initial guesses are much smaller than the default `InitialParameterCovariance`, 10000, the initial guesses are given less importance during estimation. In that case, specify a smaller initial parameter covariance.

`InitialC` is a tunable property. You can change it when the object is in a locked state.

**Default:** [1 eps]

## ParameterCovariance

Estimated covariance  $P$  of the parameters, returned as an  $N$ -by- $N$  symmetric positive-definite matrix.  $N$  is the number of parameters to be estimated. The software computes  $P$  assuming that the residuals (difference between estimated and measured outputs) are white noise, and the variance of these residuals is 1. `ParameterCovariance` is applicable only when `EstimationMethod` is 'ForgettingFactor' or 'KalmanFilter'.

The interpretation of  $P$  depends on the estimation method:

- 'ForgettingFactor' —  $R_2/2 * P$  is approximately equal to the covariance matrix of the estimated parameters, where  $R_2$  is the true variance of the residuals.
- 'KalmanFilter' —  $R_2 * P$  is the covariance matrix of the estimated parameters, and  $R_1 / R_2$  is the covariance matrix of the parameter changes. Where,  $R_1$  is the covariance matrix that you specify in `ProcessNoiseCovariance`.

`ParameterCovariance` is a read-only property and is initially empty after you create the object. It is populated after you use the `step` command for online parameter estimation.

## InitialParameterCovariance

Covariance of the initial parameter estimates, specified as one of the following:

- Real positive scalar,  $\alpha$  — Covariance matrix is an  $N$ -by- $N$  diagonal matrix, with  $\alpha$  as the diagonal elements.  $N$  is the number of parameters to be estimated.
- Vector of real positive scalars,  $[\alpha_1, \dots, \alpha_N]$  — Covariance matrix is an  $N$ -by- $N$  diagonal matrix, with  $[\alpha_1, \dots, \alpha_N]$  as the diagonal elements.
- $N$ -by- $N$  symmetric positive-definite matrix.

`InitialParameterCovariance` represents the uncertainty in the initial parameter estimates. For large values of `InitialParameterCovariance`, less importance is placed on the initial parameter values and more on the measured data during beginning of estimation using `step`.

Use only when `EstimationMethod` is 'ForgettingFactor' or 'KalmanFilter'.

`InitialParameterCovariance` is a tunable property. You can change it when the object is in a locked state.

**Default:** 10000

## EstimationMethod

Recursive estimation algorithm used for online estimation of model parameters, specified as one of the following values:

- 'ForgettingFactor' — Algorithm used for parameter estimation
- 'KalmanFilter' — Algorithm used for parameter estimation
- 'NormalizedGradient' — Algorithm used for parameter estimation
- 'Gradient' — Unnormalized gradient algorithm used for parameter estimation

Forgetting factor and Kalman filter algorithms are more computationally intensive than gradient and unnormalized gradient methods. However, they have better convergence properties. For information about these algorithms, see “Recursive Algorithms for Online Parameter Estimation”.

`EstimationMethod` is a nontunable property. You cannot change it during execution, that is after the object is locked using the `step` command. If you want to deploy code using MATLAB Coder, `EstimationMethod` can only be assigned once.

**Default:** 'ForgettingFactor'

### **ForgettingFactor**

Forgetting factor,  $\lambda$ , relevant for parameter estimation, specified as a scalar in the range (0,1].

Suppose that the system remains approximately constant over  $T_0$  samples. You can choose  $\lambda$  such that:

$$T_0 = \frac{1}{1 - \lambda}$$

- Setting  $\lambda = 1$  corresponds to “no forgetting” and estimating constant coefficients.
- Setting  $\lambda < 1$  implies that past measurements are less significant for parameter estimation and can be “forgotten”. Set  $\lambda < 1$  to estimate time-varying coefficients.

Typical choices of  $\lambda$  are in the range [0.98 0.995].

Use only when EstimationMethod is 'ForgettingFactor'.

ForgettingFactor is a tunable property. You can change it when the object is in a locked state.

**Default:** 1

### **EnableAdaption**

Enable or disable parameter estimation, specified as one of the following:

- `true` or `1`— The `step` command estimates the parameter values for that time step and updates the parameter values.
- `false` or `0` — The `step` command does not update the parameters for that time step and instead outputs the last estimated value. You can use this option when your system enters a mode where the parameter values do not vary with time.

---

**Note** If you set `EnableAdaption` to `false`, you must still execute the `step` command. Do not skip `step` to keep parameter values constant, because parameter estimation depends on current and past I/O measurements. `step` ensures past I/O data is stored, even when it does not update the parameters.

---

EnableAdapation is a tunable property. You can change it when the object is in a locked state.

**Default:** true

### **DataType**

Floating point precision of parameters, specified as one of the following values:

- 'double' — Double-precision floating point
- 'single' — Single-precision floating point

Setting `DataType` to 'single' saves memory, but leads to loss of precision. Specify `DataType` based on the precision required by the target processor where you will deploy generated code.

`DataType` is a nontunable property. It can only be set during object construction using `Name, Value` arguments and cannot be changed afterward.

**Default:** 'double'

### **ProcessNoiseCovariance**

Covariance matrix of parameter variations, specified as one of the following:

- Real nonnegative scalar,  $\alpha$  — Covariance matrix is an  $N$ -by- $N$  diagonal matrix, with  $\alpha$  as the diagonal elements.
- Vector of real nonnegative scalars,  $[\alpha_1, \dots, \alpha_N]$  — Covariance matrix is an  $N$ -by- $N$  diagonal matrix, with  $[\alpha_1, \dots, \alpha_N]$  as the diagonal elements.
- $N$ -by- $N$  symmetric positive semidefinite matrix.

$N$  is the number of parameters to be estimated.

`ProcessNoiseCovariance` is applicable when `EstimationMethod` is 'KalmanFilter'.

Kalman filter algorithm treats the parameters as states of a dynamic system and estimates these parameters using a Kalman filter. `ProcessNoiseCovariance` is the covariance of the process noise acting on these parameters. Zero values in the noise covariance matrix correspond to estimating constant coefficients. Values larger than 0 correspond to time-varying parameters. Use large values for rapidly changing parameters. However, the larger values result in noisier parameter estimates.

ProcessNoiseCovariance is a tunable property. You can change it when the object is in a locked state.

**Default:** 0.1

### **AdaptationGain**

Adaptation gain,  $\gamma$ , used in gradient recursive estimation algorithms, specified as a positive scalar.

AdaptationGain is applicable when EstimationMethod is 'Gradient' or 'NormalizedGradient'.

Specify a large value for AdaptationGain when your measurements have a high signal-to-noise ratio.

AdaptationGain is a tunable property. You can change it when the object is in a locked state.

**Default:** 1

### **NormalizationBias**

Bias in adaptation gain scaling used in the 'NormalizedGradient' method, specified as a nonnegative scalar.

NormalizationBias is applicable when EstimationMethod is 'NormalizedGradient'.

The normalized gradient algorithm divides the adaptation gain at each step by the square of the two-norm of the gradient vector. If the gradient is close to zero, this can cause jumps in the estimated parameters. NormalizationBias is the term introduced in the denominator to prevent these jumps. Increase NormalizationBias if you observe jumps in estimated parameters.

NormalizationBias is a tunable property. You can change it when the object is in a locked state.

**Default:** eps

## Output Arguments

**obj** — System object for online parameter estimation of ARMAX model  
recursiveARMAX System object

System object for online parameter estimation of ARMAX model, returned as a recursiveARMAX System object. This object is created using the specified model orders and properties. Use `step` command to estimate the coefficients of the ARMAX model polynomials. You can then access the estimated coefficients and parameter covariance using dot notation. For example, type `obj.A` to view the estimated A polynomial coefficients.

## Definitions

### ARMAX Model Structure

The ARMAX model structure is:

$$y(t) + a_1 y(t-1) + \dots + a_{n_a} y(t-n_a) = b_1 u(t-n_k) + \dots + b_{n_b} u(t-n_k-n_b+1) + c_1 e(t-1) + \dots + c_{n_c} e(t-n_c) + e(t)$$

A more compact way to write the difference equation is:

$$A(q)y(t) = B(q)u(t-n_k) + C(q)e(t)$$

where,

- $y(t)$  — Output at time  $t$ .
- $n_a$  — Number of poles.
- $n_b$  — Number of zeroes plus 1.
- $n_c$  — Number of  $C$  coefficients.

- $n_k$  — Number of input samples that occur before the input affects the output, also called the *dead time* in the system.
- $y(t-1)\dots y(t-n_a)$  — Previous outputs on which the current output depends.
- $u(t-n_k)\dots u(t-n_k-n_b+1)$  — Previous and delayed inputs on which the current output depends.
- $e(t-1)\dots e(t-n_c)$  — White-noise disturbance value.

The parameters  $n_a$ ,  $n_b$ , and  $n_c$  are the orders of the ARMAX model, and  $n_k$  is the delay.  $q$  is the delay operator. Specifically,

$$A(q) = 1 + a_1q^{-1} + \dots + a_{n_a}q^{-n_a}$$

$$B(q) = b_1 + b_2q^{-1} + \dots + b_{n_b}q^{-n_b+1}$$

$$C(q) = 1 + c_1q^{-1} + \dots + c_{n_c}q^{-n_c}$$

## Tips

- Starting in R2016b, instead of using the `step` command to update model parameter estimates, you can call the System object with input arguments, as if it were a function. For example, `[A,B,C,EstimatedOutput] = step(obj,y,u)` and `[A,B,C,EstimatedOutput] = obj(y,u)` perform equivalent operations.

## Extended Capabilities

### C/C++ Code Generation

Generate C and C++ code using MATLAB® Coder™.

Usage notes and limitations:

- For Simulink-based workflows, use Recursive Polynomial Model Estimator.
- For limitations, see “Generate Code for Online Parameter Estimation in MATLAB”.

Supports MATLAB Function block: No

## See Also

Recursive Polynomial Model Estimator | `armax` | `clone` | `isLocked` | `recursiveAR` | `recursiveARMA` | `recursiveARX` | `recursiveBJ` | `recursiveLS` | `recursiveOE` | `release` | `reset` | `step`

## Topics

“Perform Online Parameter Estimation at the Command Line”

“Validate Online Parameter Estimation at the Command Line”

“What Is Online Estimation?”

“Recursive Algorithms for Online Parameter Estimation”

**Introduced in R2015b**

## recursiveARX

Create System object for online parameter estimation of ARX model

Use `recursiveARX` command for parameter estimation with real-time data. If all data necessary for estimation is available at once, and you are estimating a time-invariant model, use the offline estimation command, `arx`.

### Syntax

```
obj = recursiveARX
obj = recursiveARX(Orders)
obj = recursiveARX(Orders,A0,B0)
obj = recursiveARX( ____,Name,Value)
```

### Description

`obj = recursiveARX` creates a System object for online parameter estimation of a default ARX model structure on page 1-1390. The default model structure has polynomials of order 1 and initial polynomial coefficient values `eps`.

After creating the object, use the `step` command to update model parameter estimates using recursive estimation algorithms and real-time data.

`obj = recursiveARX(Orders)` specifies the polynomial orders of the ARX model to be estimated.

`obj = recursiveARX(Orders,A0,B0)` specifies the polynomial orders and initial values of the polynomial coefficients.

`obj = recursiveARX( ____,Name,Value)` specifies additional attributes of the ARX model structure and recursive estimation algorithm using one or more `Name,Value` pair arguments.

## Object Description

recursiveARX creates a System object for online parameter estimation of single-input single-output (SISO) or multiple-input single-output (MISO) ARX models using a recursive estimation algorithm.

A System object is a specialized MATLAB object designed specifically for implementing and simulating dynamic systems with inputs that change over time. System objects use internal states to store past behavior, which is used in the next computational step.

After you create a System object, you use commands to process data or obtain information from or about the object. System objects use a minimum of two commands to process data — a constructor to create the object and the `step` command to update object parameters using real-time data. This separation of declaration from execution lets you create multiple, persistent, reusable objects, each with different settings.

You can use the following commands with the online estimation System objects in System Identification Toolbox:

Command	Description
<code>step</code>	Update model parameter estimates using recursive estimation algorithms and real-time data.  <code>step</code> puts the object into a locked state. In a locked state, you cannot change any nontunable properties or input specifications, such as model order, data type, or estimation algorithm. During execution, you can only change tunable properties.
<code>release</code>	Unlock the System object. Use this command to enable setting of nontunable parameters.
<code>reset</code>	Reset the internal states of a locked System object to the initial values, and leave the object locked.

Command	Description
<code>clone</code>	Create another System object with the same object property values.  Do not create additional objects using syntax <code>obj2 = obj</code> . Any changes made to the properties of the new object created this way ( <code>obj2</code> ) also change the properties of the original object ( <code>obj</code> ).
<code>isLocked</code>	Query locked status for input attributes and nontunable properties of the System object.

Use the `recursiveARX` command to create an online estimation System object. Then estimate the ARX model parameters (A and B) and output using the `step` command with incoming input and output data, `u` and `y`.

```
[A,B,EstimatedOutput] = step(obj,y,u)
```

For `recursiveARX` object properties, see “Properties” on page 1-1380.

## Examples

### Estimate a SISO ARX Model Online

Create a System object for online parameter estimation of a SISO ARX model.

```
obj = recursiveARX;
```

The ARX model has a default structure with polynomials of order 1 and initial polynomial coefficient values, `eps`.

Load the estimation data. In this example, use a static data set for illustration.

```
load iddata1 z1;  
output = z1.y;  
input = z1.u;
```

Estimate ARX model parameters online using `step`.

```
for i = 1:numel(input)
[A,B,EstimatedOutput] = step(obj,output(i),input(i));
end
```

View the current estimated values of polynomial B coefficients.

```
obj.B
```

```
ans = 1×2
      0      0.7974
```

View the current covariance estimate of the parameters.

```
obj.ParameterCovariance
```

```
ans = 2×2
      0.0002      0.0001
      0.0001      0.0034
```

View the current estimated output.

```
EstimatedOutput
```

```
EstimatedOutput = -4.7766
```

### Create System Object for SISO ARX Model With Known Initial Parameters

Specify ARX model orders and delays.

```
na = 1;
nb = 2;
nk = 1;
```

Create a System object for online estimation of SISO ARX model with known initial polynomial coefficients.

```
A0 = [1 0.5];
B0 = [0 1 1];
obj = recursiveARX([na nb nk],A0,B0);
```

Specify the initial parameter covariance.

```
obj.InitialParameterCovariance = 0.1;
```

`InitialParameterCovariance` represents the uncertainty in your guess for the initial parameters. Typically, the default `InitialParameterCovariance` (10000) is too large relative to the parameter values. This results in initial guesses being given less importance during estimation. If you have confidence in the initial parameter guesses, specify a smaller initial parameter covariance.

## Create System Object for MISO ARX Model With Known Initial Parameters

Specify orders and delays for ARX model with two inputs and one output.

```
na = 1;  
nb = [2 1];  
nk = [1 3];
```

`nb` and `nk` are specified as row vectors of length equal to number of inputs,  $Nu$ .

Specify initial polynomial coefficients.

```
A0 = [1 0.5];  
B0 = [0 1 1 0; 0 0 0 0.8];
```

`B0` has  $Nu$  rows and  $\max(nb+nk)$  columns. The  $i$ -th row corresponds to  $i$ -th input and is specified as having  $nk(i)$  zeros, followed by  $nb(i)$  initial values. Values after  $nb(i) + nk(i)$  are ignored.

Create a System object for online estimation of ARX model with known initial polynomial coefficients.

```
obj = recursiveARX([na nb nk],A0,B0);
```

## Specify Estimation Method for Online Estimation of ARX Model

Create a System object that uses the normalized gradient algorithm for online parameter estimation of an ARX model.

```
obj = recursiveARX([1 2 1], 'EstimationMethod', 'NormalizedGradient');
```

## Input Arguments

### Orders — Model orders and delays

1-by-3 vector of integers | 1-by-3 vector of vectors

Model orders and delays of an ARX model on page 1-1390, specified as a 1-by-3 vector of integers or vectors, [na nb nk].

- **na** — Order of the polynomial  $A(q)$ , specified as a nonnegative integer.
- **nb** — Order of the polynomial  $B(q) + 1$ , specified as 1-by- $Nu$  vector of positive integers.  $Nu$  is the number of inputs.

For MISO models, there are as many  $B(q)$  polynomials as the number of inputs.  $nb(i)$  is the order of  $i$ th polynomial  $B_i(q)+1$  for the  $i$ th input.

- **nk** — Input-output delay, specified as a 1-by- $Nu$  vector of nonnegative integers.  $Nu$  is the number of inputs.

For MISO models, there are as many  $B(q)$  polynomials as the number of inputs.  $nk(i)$  is the input-output delay time corresponding to the  $i$ th input.

### A0, B0 — Initial value of polynomial coefficients

row vector and matrix of real values | [ ]

Initial value of coefficients of  $A(q)$  and  $B(q)$  polynomials, specified as row vector and matrix or real values, respectively. Specify the elements in order of ascending powers of  $q^{-1}$ .

- **A0** — Initial value for the coefficients of the polynomial  $A(q)$ , specified as a 1-by-(na+1) row vector with 1 as the first element.
- **B0** — Initial value for the coefficients of the polynomial  $B(q)$ , specified as  $Nu$ -by- $\max(nb+nk)$  matrix.  $Nu$  is the number of inputs.

For MISO models, there are as many  $B(q)$  polynomials as the number of inputs. The  $i$ th row of **B0** corresponds to the  $i$ th input and must contain  $nk(i)$  leading zeros, followed by  $nb(i)$  initial parameter values. Entries beyond  $nk(i)+nb(i)$  are ignored.

na, nb, and nk are the Orders of the model.

Specifying as `[]`, uses the default value of `eps` for the polynomial coefficients.

If the initial parameter values are much smaller than `InitialParameterCovariance`, these initial values are given less importance during estimation. Specify a smaller initial parameter covariance if you have high confidence in the initial parameter values. This statement applies only for infinite-history estimation. Finite-history estimation does not use `InitialParameterCovariance`.

## Name-Value Pair Arguments

Specify optional comma-separated pairs of `Name`, `Value` arguments. `Name` is the argument name and `Value` is the corresponding value. `Name` must appear inside quotes. You can specify several name and value pair arguments in any order as `Name1, Value1, ..., NameN, ValueN`.

Use `Name`, `Value` arguments to specify writable properties on page 1-1380 of recursiveARX System object during object creation. For example, `obj = recursiveARX([2 2 1], 'EstimationMethod', 'Gradient')` creates a System object to estimate an ARX model using the 'Gradient' recursive estimation algorithm.

## Properties

recursiveARX System object properties consist of read-only and writable properties. The writable properties are tunable and nontunable properties. The nontunable properties cannot be changed when the object is locked, that is, after you use the `step` command.

Use `Name`, `Value` arguments to specify writable properties of recursiveARX objects during object creation. After object creation, use dot notation to modify the tunable properties.

```
obj = recursiveARX;  
obj.ForgettingFactor = 0.99;
```

### A

Estimated coefficients of polynomial  $A(q)$ , returned as a row vector of real values specified in order of ascending powers of  $q^{-1}$ .

$A$  is a read-only property and is initially empty after you create the object. It is populated after you use the `step` command for online parameter estimation.

## **B**

Estimated coefficients of polynomial  $B(q)$ , returned as a  $Nu$ -by- $\max(\text{nb}+\text{nk})$  matrix of real values.  $Nu$  is the number of inputs.

The  $i$ th row of  $B$  corresponds to the  $i$ th input and contains  $\text{nk}(i)$  leading zeros, followed by  $\text{nb}(i)$  estimated parameters, specified in order of ascending powers of  $q^{-1}$ . Ignore zero entries beyond  $\text{nk}(i)+\text{nb}(i)$ .

$B$  is a read-only property and is initially empty after you create the object. It is populated after you use the `step` command for online parameter estimation.

## **InitialA**

Initial values for the coefficients of polynomial  $A(q)$  of order  $\text{na}$ , specified as a row vector of length  $\text{na}+1$ , with 1 as the first element. Specify the coefficients in order of ascending powers of  $q^{-1}$ .

If the initial guesses are much smaller than the default `InitialParameterCovariance`, 10000, the initial guesses are given less importance during estimation. In that case, specify a smaller initial parameter covariance.

`InitialA` is a tunable property. You can change it when the object is in a locked state.

**Default:** [1 eps]

## **InitialB**

Initial values for the coefficients of polynomial  $B(q)$ , specified as an  $Nu$ -by- $\max(\text{nb}+\text{nk})$  matrix.  $Nu$  is the number of inputs.

For MISO models, there are as many  $B(q)$  polynomials as the number of inputs. The  $i$ th row of  $B_0$  corresponds to the  $i$ th input and must contain  $\text{nk}(i)$  zeros, followed by  $\text{nb}(i)$  initial parameter values. Entries beyond  $\text{nk}(i)+\text{nb}(i)$  are ignored.

If the initial guesses are much smaller than the default `InitialParameterCovariance`, 10000, the initial guesses are given less importance during estimation. In that case, specify a smaller initial parameter covariance.

`InitialB` is a tunable property. You can change it when the object is in a locked state.

**Default:** [0 eps]

### **InitialOutputs**

Initial values of the measured outputs buffer in finite-history estimation, specified as 0 or as a  $(W+na)$ -by-1 vector, where  $W$  is the window length and  $na$  is the order of the polynomial  $A(q)$  that you specify when constructing the object.

The `InitialOutputs` property provides a means of controlling the initial behavior of the algorithm.

When `InitialOutputs` is set to 0, the object populates the buffer with zeros.

If the initial buffer is set to 0 or does not contain enough information, you see a warning message during the initial phase of your estimation. The warning should clear after a few cycles. The number of cycles it takes for sufficient information to be buffered depends upon the order of your polynomials and your input delays. If the warning persists, you should evaluate the content of your signals.

Specify `InitialOutputs` only when `History` is `Finite`.

`InitialOutputs` is a tunable property. You can change `InitialOutputs` when the object is in a locked state.

**Default:** 0

### **InitialInputs**

Initial values of the inputs in the finite history window, specified as 0 or as a  $(W-1+\max(nb)+\max(nk))$ -by- $nu$  matrix, where  $W$  is the window length and  $nu$  is the number of inputs.  $nb$  is the vector of  $B(q)$  polynomial orders and  $nk$  is vector of input delays that you specify when constructing the `recursiveARX` object.

The `InitialInputs` property provides a means of controlling the initial behavior of the algorithm.

When the `InitialInputs` is set to 0, the object populates the buffer with zeros.

If the initial buffer is set to 0 or does not contain enough information, you see a warning message during the initial phase of your estimation. The warning should clear after a few cycles. The number of cycles it takes for sufficient information to be buffered depends upon the order of your polynomials and your input delays. If the warning persists, you should evaluate the content of your signals.

Specify `InitialInputs` only when `History` is `Finite`.

`InitialInputs` is a tunable property. You can change `InitialInputs` when the object is in a locked state.

**Default:** 0

### ParameterCovariance

Estimated covariance  $P$  of the parameters, returned as an  $N$ -by- $N$  symmetric positive-definite matrix.  $N$  is the number of parameters to be estimated. The software computes  $P$  assuming that the residuals (difference between estimated and measured outputs) are white noise, and the variance of these residuals is 1.

`ParameterCovariance` is applicable only when `EstimationMethod` is `'ForgettingFactor'` or `'KalmanFilter'` or when `History` is `Finite`.

The interpretation of  $P$  depends on your settings for the `History` and `EstimationMethod` properties.

- If `History` is `Infinite`, then your `EstimationMethod` selection results in one of the following:
  - `'ForgettingFactor'` —  $(R_2/2)P$  is approximately equal to the covariance matrix of the estimated parameters, where  $R_2$  is the true variance of the residuals.
  - `'KalmanFilter'` —  $R_2P$  is the covariance matrix of the estimated parameters, and  $R_1/R_2$  is the covariance matrix of the parameter changes. Here,  $R_1$  is the covariance matrix that you specify in `ProcessNoiseCovariance`.
- If `History` is `Finite` (sliding-window estimation) —  $R_2P$  is the covariance of the estimated parameters. The sliding-window algorithm does not use this covariance in the parameter-estimation process. However, the algorithm does compute the covariance for output so that you can use it for statistical evaluation.

`ParameterCovariance` is a read-only property and is initially empty after you create the object. It is populated after you use the `step` command for online parameter estimation.

### InitialParameterCovariance

Covariance of the initial parameter estimates, specified as one of the following:

- Real positive scalar,  $\alpha$  — Covariance matrix is an  $N$ -by- $N$  diagonal matrix, with  $\alpha$  as the diagonal elements.  $N$  is the number of parameters to be estimated.

- Vector of real positive scalars,  $[\alpha_1, \dots, \alpha_N]$  — Covariance matrix is an  $N$ -by- $N$  diagonal matrix, with  $[\alpha_1, \dots, \alpha_N]$  as the diagonal elements.
- $N$ -by- $N$  symmetric positive-definite matrix.

`InitialParameterCovariance` represents the uncertainty in the initial parameter estimates. For large values of `InitialParameterCovariance`, less importance is placed on the initial parameter values and more on the measured data during beginning of estimation using `step`.

Use only when `EstimationMethod` is 'ForgettingFactor' or 'KalmanFilter'.

`InitialParameterCovariance` is a tunable property. You can change it when the object is in a locked state.

**Default:** 10000

## **EstimationMethod**

Recursive estimation algorithm used for online estimation of model parameters, specified as one of the following values:

- 'ForgettingFactor' — Algorithm used for parameter estimation
- 'KalmanFilter' — Algorithm used for parameter estimation
- 'NormalizedGradient' — Algorithm used for parameter estimation
- 'Gradient' — Unnormalized gradient algorithm used for parameter estimation

Forgetting factor and Kalman filter algorithms are more computationally intensive than gradient and unnormalized gradient methods. However, they have better convergence properties. For information about these algorithms, see “Recursive Algorithms for Online Parameter Estimation”.

These methods all use an infinite data history, and are available only when `History` is 'Infinite'.

`EstimationMethod` is a nontunable property. You cannot change it during execution, that is, after the object is locked using the `step` command.

**Default:** Forgetting Factor

## ForgettingFactor

Forgetting factor,  $\lambda$ , relevant for parameter estimation, specified as a scalar in the range (0,1].

Suppose that the system remains approximately constant over  $T_0$  samples. You can choose  $\lambda$  such that:

$$T_0 = \frac{1}{1 - \lambda}$$

- Setting  $\lambda = 1$  corresponds to “no forgetting” and estimating constant coefficients.
- Setting  $\lambda < 1$  implies that past measurements are less significant for parameter estimation and can be “forgotten”. Set  $\lambda < 1$  to estimate time-varying coefficients.

Typical choices of  $\lambda$  are in the range [0.98 0.995].

Use only when EstimationMethod is 'ForgettingFactor'.

ForgettingFactor is a tunable property. You can change it when the object is in a locked state.

**Default:** 1

## EnableAdapation

Enable or disable parameter estimation, specified as one of the following:

- `true` or `1`— The `step` command estimates the parameter values for that time step and updates the parameter values.
- `false` or `0` — The `step` command does not update the parameters for that time step and instead outputs the last estimated value. You can use this option when your system enters a mode where the parameter values do not vary with time.

---

**Note** If you set `EnableAdapation` to `false`, you must still execute the `step` command. Do not skip `step` to keep parameter values constant, because parameter estimation depends on current and past I/O measurements. `step` ensures past I/O data is stored, even when it does not update the parameters.

---

`EnableAdapation` is a tunable property. You can change it when the object is in a locked state.

**Default:** true

### **DataType**

Floating point precision of parameters, specified as one of the following values:

- 'double' — Double-precision floating point
- 'single' — Single-precision floating point

Setting `DataType` to 'single' saves memory, but leads to loss of precision. Specify `DataType` based on the precision required by the target processor where you will deploy generated code.

`DataType` is a nontunable property. It can only be set during object construction using `Name, Value` arguments and cannot be changed afterward.

**Default:** 'double'

### **ProcessNoiseCovariance**

Covariance matrix of parameter variations, specified as one of the following:

- Real nonnegative scalar,  $\alpha$  — Covariance matrix is an  $N$ -by- $N$  diagonal matrix, with  $\alpha$  as the diagonal elements.
- Vector of real nonnegative scalars,  $[\alpha_1, \dots, \alpha_N]$  — Covariance matrix is an  $N$ -by- $N$  diagonal matrix, with  $[\alpha_1, \dots, \alpha_N]$  as the diagonal elements.
- $N$ -by- $N$  symmetric positive semidefinite matrix.

$N$  is the number of parameters to be estimated.

`ProcessNoiseCovariance` is applicable when `EstimationMethod` is 'KalmanFilter'.

Kalman filter algorithm treats the parameters as states of a dynamic system and estimates these parameters using a Kalman filter. `ProcessNoiseCovariance` is the covariance of the process noise acting on these parameters. Zero values in the noise covariance matrix correspond to estimating constant coefficients. Values larger than 0 correspond to time-varying parameters. Use large values for rapidly changing parameters. However, the larger values result in noisier parameter estimates.

`ProcessNoiseCovariance` is a tunable property. You can change it when the object is in a locked state.

**Default:** 0.1

### **AdaptationGain**

Adaptation gain,  $\gamma$ , used in gradient recursive estimation algorithms, specified as a positive scalar.

AdaptationGain is applicable when EstimationMethod is 'Gradient' or 'NormalizedGradient'.

Specify a large value for AdaptationGain when your measurements have a high signal-to-noise ratio.

AdaptationGain is a tunable property. You can change it when the object is in a locked state.

**Default:** 1

### **NormalizationBias**

Bias in adaptation gain scaling used in the 'NormalizedGradient' method, specified as a nonnegative scalar.

NormalizationBias is applicable when EstimationMethod is 'NormalizedGradient'.

The normalized gradient algorithm divides the adaptation gain at each step by the square of the two-norm of the gradient vector. If the gradient is close to zero, this can cause jumps in the estimated parameters. NormalizationBias is the term introduced in the denominator to prevent these jumps. Increase NormalizationBias if you observe jumps in estimated parameters.

NormalizationBias is a tunable property. You can change it when the object is in a locked state.

**Default:** eps

### **History**

Data history type defining which type of recursive algorithm you use, specified as:

- 'Infinite' — Use an algorithm that aims to minimize the error between the observed and predicted outputs for all time steps from the beginning of the simulation.

- `'Finite'` — Use an algorithm that aims to minimize the error between the observed and predicted outputs for a finite number of past time steps.

Algorithms with infinite history aim to produce parameter estimates that explain all data since the start of the simulation. These algorithms still use a fixed amount of memory that does not grow over time. The object provides multiple algorithms of the `'Infinite'` `History` type. Specifying this option activates the `EstimationMethod` property with which you specify an algorithm.

Algorithms with finite history aim to produce parameter estimates that explain only a finite number of past data samples. This method is also called sliding-window estimation. The object provides one algorithm of the `'Finite'` type. Specifying this option activates the `WindowLength` property that sizes the window.

For more information on recursive estimation methods, see “Recursive Algorithms for Online Parameter Estimation”.

`History` is a nontunable property. It can be set only during object construction using `Name, Value` arguments and cannot be changed afterward.

**Default:** `'Infinite'`

### **WindowLength**

Window size determining the number of time samples to use for the sliding-window estimation method, specified as a positive integer. Specify `WindowLength` only when `History` is `Finite`.

Choose a window size that balances estimation performance with computational and memory burden. Sizing factors include the number and time variance of the parameters in your model. Always specify **Window Length** in samples, even if you are using frame-based input processing.

`WindowLength` must be greater than or equal to the number of estimated parameters.

Suitable window length is independent of whether you are using sample-based or frame-based input processing (see `InputProcessing`). However, when using frame-based processing, your window length must be greater than or equal to the number of samples (time steps) contained in the frame.

`WindowLength` is a nontunable property. It can be set only during object construction using `Name, Value` arguments and cannot be changed afterward.

**Default:** 200

## InputProcessing

Option for sample-based or frame-based input processing, specified as a character vector or string.

- **Sample-based** processing operates on signals streamed one sample at a time.
- **Frame-based** processing operates on signals containing samples from multiple time steps. Many machine sensor interfaces package multiple samples and transmit these samples together in frames. **Frame-based** processing allows you to input this data directly without having to first unpack it.

Your `InputProcessing` specification impacts the dimensions for the input and output signals when using the `step` command:

```
[theta,EstimatedOutput] = step(obj,y,u)
```

- **Sample-based**
  - `y` and `EstimatedOutput` are scalars.
  - `u` is a 1-by- $Nu$  vector, where  $Nu$  is the number of inputs.
- **Frame-based with  $M$  samples per frame**
  - `y` and `EstimatedOutput` are  $M$ -by-1 vectors.
  - `u` is an  $M$ -by- $Nu$  matrix.

`InputProcessing` is a nontunable property. It can be set only during object construction using `Name,Value` arguments and cannot be changed afterward.

**Default:** 'Sample-based'

## Output Arguments

**obj** — System object for online parameter estimation of ARX model

recursiveARX System object

System object for online parameter estimation of ARX model, returned as a recursiveARX System object. This object is created using the specified model orders and properties. Use `step` command to estimate the coefficients of the ARX model

polynomials. You can then access the estimated coefficients and parameter covariance using dot notation. For example, type `obj.A` to view the estimated  $A$  polynomial coefficients.

## Definitions

### ARX Model Structure

The ARX model structure is :

$$y(t) + a_1 y(t-1) + \dots + a_{n_a} y(t-n_a) = b_1 u(t-n_k) + \dots + b_{n_b} u(t-n_b-n_k+1) + e(t)$$

The parameters  $n_a$  and  $n_b$  are the orders of the ARX model, and  $n_k$  is the delay.

- $y(t)$  — Output at time  $t$ .
- $n_a$  — Number of poles.
- $n_b$  — Number of zeroes plus 1.
- $n_k$  — Number of input samples that occur before the input affects the output, also called the *dead time* in the system.
- $y(t-1) \dots y(t-n_a)$  — Previous outputs on which the current output depends.
- $u(t-n_k) \dots u(t-n_k-n_b+1)$  — Previous and delayed inputs on which the current output depends.
- $e(t)$  — White-noise disturbance value.

A more compact way to write the difference equation is

$$A(q)y(t) = B(q)u(t-n_k) + e(t)$$

$q$  is the delay operator. Specifically,

$$A(q) = 1 + a_1 q^{-1} + \dots + a_{n_a} q^{-n_a}$$

$$B(q) = b_1 + b_2q^{-1} + \dots + b_{n_b}q^{-n_b+1}$$

## Tips

- Starting in R2016b, instead of using the `step` command to update model parameter estimates, you can call the System object with input arguments, as if it were a function. For example, `[A,B,EstimatedOutput] = step(obj,y,u)` and `[A,B,EstimatedOutput] = obj(y,u)` perform equivalent operations.

## Extended Capabilities

### C/C++ Code Generation

Generate C and C++ code using MATLAB® Coder™.

Usage notes and limitations:

- For Simulink based workflows, use Recursive Polynomial Model Estimator.
- For limitations, see “Generate Code for Online Parameter Estimation in MATLAB”.

Supports MATLAB Function block: No

## See Also

Recursive Polynomial Model Estimator | `arx` | `clone` | `isLocked` | `recursiveAR` | `recursiveARMA` | `recursiveARMAX` | `recursiveBJ` | `recursiveLS` | `recursiveOE` | `release` | `reset` | `step`

## Topics

“Perform Online Parameter Estimation at the Command Line”

“Validate Online Parameter Estimation at the Command Line”

“Online ARX Parameter Estimation for Tracking Time-Varying System Dynamics”

“What Is Online Estimation?”

“Recursive Algorithms for Online Parameter Estimation”

**Introduced in R2015b**

## recursiveBJ

Create System object for online parameter estimation of Box-Jenkins polynomial model

Use `recursiveBJ` command for parameter estimation with real-time data. If all data necessary for estimation is available at once, and you are estimating a time-invariant model, use the offline estimation command, `bj`.

### Syntax

```
obj = recursiveBJ
obj = recursiveBJ(Orders)
obj = recursiveBJ(Orders,B0,C0,D0,F0)
obj = recursiveBJ( ____,Name,Value)
```

### Description

`obj = recursiveBJ` creates a System object for online parameter estimation of a default single-input single-output (SISO) Box-Jenkins polynomial model structure on page 1-1407. The default model structure has polynomials of order 1 and initial polynomial coefficient values `eps`.

After creating the object, use the `step` command to update model parameter estimates using recursive estimation algorithms and real-time data.

`obj = recursiveBJ(Orders)` specifies the polynomial orders of the Box-Jenkins model to be estimated.

`obj = recursiveBJ(Orders,B0,C0,D0,F0)` specifies the polynomial orders and initial values of the polynomial coefficients. Specify initial values to potentially avoid local minima during estimation. If the initial values are small compared to the default `InitialParameterCovariance` property value, and you have confidence in your initial values, also specify a smaller `InitialParameterCovariance`.

`obj = recursiveBJ( ____,Name,Value)` specifies additional attributes of the Box-Jenkins model structure and recursive estimation algorithm using one or more `Name,Value` pair arguments.

## Object Description

`recursiveBJ` creates a System object for online parameter estimation of SISO Box-Jenkins polynomial models using a recursive estimation algorithm.

A System object is a specialized MATLAB object designed specifically for implementing and simulating dynamic systems with inputs that change over time. System objects use internal states to store past behavior, which is used in the next computational step.

After you create a System object, you use commands to process data or obtain information from or about the object. System objects use a minimum of two commands to process data — a constructor to create the object and the `step` command to update object parameters using real-time data. This separation of declaration from execution lets you create multiple, persistent, reusable objects, each with different settings.

You can use the following commands with the online estimation System objects in System Identification Toolbox:

Command	Description
<code>step</code>	Update model parameter estimates using recursive estimation algorithms and real-time data.  <code>step</code> puts the object into a locked state. In a locked state, you cannot change any nontunable properties or input specifications, such as model order, data type, or estimation algorithm. During execution, you can only change tunable properties.
<code>release</code>	Unlock the System object. Use this command to enable setting of nontunable parameters.
<code>reset</code>	Reset the internal states of a locked System object to the initial values, and leave the object locked.

Command	Description
clone	Create another System object with the same object property values.  Do not create additional objects using syntax <code>obj2 = obj</code> . Any changes made to the properties of the new object created this way ( <code>obj2</code> ) also change the properties of the original object ( <code>obj</code> ).
isLocked	Query locked status for input attributes and nontunable properties of the System object.

Use the `recursiveBJ` command to create an online estimation System object. Then estimate the Box-Jenkins polynomial model parameters (B, C, D, and F) and output using the `step` command with incoming input and output data, `u` and `y`.

```
[B,C,D,F,EstimatedOutput] = step(obj,y,u)
```

For `recursiveBJ` object properties, see “Properties” on page 1-1400.

## Examples

### Estimate Box-Jenkins Polynomial Model Online

Create a System object for online parameter estimation of a Box-Jenkins polynomial model.

```
obj = recursiveBJ;
```

The Box-Jenkins model has a default structure with polynomials of order 1 and initial polynomial coefficient values, `eps`.

Load the estimation data. In this example, use a static data set for illustration.

```
load iddata1 z1;
output = z1.y;
input = z1.u;
```

Estimate Box-Jenkins model parameters online using `step`.

```
for i = 1:numel(input)
[B,C,D,F,EstimatedOutput] = step(obj,output(i),input(i));
end
```

View the current estimated values of polynomial D coefficients.

```
obj.D
```

```
ans = 1x2
```

```
1.0000 -0.6876
```

View the current covariance estimate of the parameters.

```
obj.ParameterCovariance
```

```
ans = 4x4
```

```
0.0020 -0.0004 -0.0001 0.0002
-0.0004 0.0007 0.0006 -0.0001
-0.0001 0.0006 0.0007 -0.0000
0.0002 -0.0001 -0.0000 0.0001
```

View the current estimated output.

```
EstimatedOutput
```

```
EstimatedOutput = -4.1905
```

## Create System Object for Box-Jenkins Model With Known Orders and Delays

Specify Box-Jenkins polynomial model orders and delays.

```
nb = 1;
nc = 1;
nd = 2;
nf = 1;
nk = 1;
```

Create a System object for online estimation of Box-Jenkins model with the specified orders and delays.

```
obj = recursiveBJ([nb nc nd nf nk]);
```

### Create System Object for Box-Jenkins Model With Known Initial Parameters

Specify Box-Jenkins polynomial model orders and delays.

```
nb = 1;  
nc = 1;  
nd = 1;  
nf = 2;  
nk = 1;
```

Create a System object for online estimation of Box-Jenkins model with known initial polynomial coefficients.

```
B0 = [0 1];  
C0 = [1 0.5];  
D0 = [1 0.9];  
F0 = [1 0.7 0.8];  
obj = recursiveBJ([nb nc nd nf nk],B0,C0,D0,F0);
```

Specify the initial parameter covariance.

```
obj.InitialParameterCovariance = 0.1;
```

`InitialParameterCovariance` represents the uncertainty in your guess for the initial parameters. Typically, the default `InitialParameterCovariance` (10000) is too large relative to the parameter values. This results in initial guesses being given less importance during estimation. If you have confidence in the initial parameter guesses, specify a smaller initial parameter covariance.

### Specify Estimation Method for Online Estimation of Box-Jenkins Model

Create a System object that uses the normalized gradient algorithm for online parameter estimation of a Box-Jenkins model.

```
obj = recursiveBJ([1 1 1 2 1], 'EstimationMethod', 'NormalizedGradient');
```

## Input Arguments

### Orders — Model orders and delays

1-by-5 vector of integers

Model orders and delays of a Box-Jenkins polynomial model on page 1-1407, specified as a 1-by-5 vector of integers, [nb nc nd nf nk].

- nb — Order of the polynomial  $B(q) + 1$ , specified as a positive integer.
- nc — Order of the polynomial  $C(q)$ , specified as a nonnegative integer.
- nd — Order of the polynomial  $D(q)$ , specified as a nonnegative integer.
- nf — Order of the polynomial  $F(q)$ , specified as a nonnegative integer.
- nk — Input-output delay, specified as a positive integer. nk is number of input samples that occur before the input affects the output. nk is expressed as fixed leading zeros of the  $B$  polynomial.

### B0, C0, D0, F0 — Initial value of polynomial coefficients

row vectors of real values | []

Initial value of polynomial coefficients, specified as row vectors of real values with elements in order of ascending powers of  $q^{-1}$ .

- B0 — Initial guess for the coefficients of the polynomial  $B(q)$ , specified as a 1-by-(nb +nk) vector with nk leading zeros.
- C0 — Initial guess for the coefficients of the polynomial  $C(q)$ , specified as a 1-by-(nc +1) vector with 1 as the first element.

The coefficients in C0 must define a stable discrete-time polynomial with roots within a unit disk. For example,

```
C0 = [1 0.5 0.5];  
all(abs(roots(C0))<1)
```

```
ans =
```

```
1
```

- D0 — Initial guess for the coefficients of the polynomial  $D(q)$ , specified as a 1-by-(nd +1) vector with 1 as the first element.

The coefficients in D0 must define a stable discrete-time polynomial with roots within a unit disk. For example,

```
D0 = [1 0.9 0.8];
all(abs(roots(D0))<1)
```

```
ans =
```

```
1
```

- **F0** — Initial guess for the coefficients of the polynomial  $F(q)$ , specified as a 1-by-(nf +1) vector with 1 as the first element.

The coefficients in **F0** must define a stable discrete-time polynomial with roots within a unit disk. For example,

```
F0 = [1 0.5 0.5];
all(abs(roots(F0))<1)
```

```
ans =
```

```
1
```

Specifying as `[]`, uses the default value of `eps` for the polynomial coefficients.

---

**Note** If the initial guesses are much smaller than the default `InitialParameterCovariance`, 10000, the initial guesses are given less importance during estimation. In that case, specify a smaller initial parameter covariance.

---

## Name-Value Pair Arguments

Specify optional comma-separated pairs of `Name`, `Value` arguments. `Name` is the argument name and `Value` is the corresponding value. `Name` must appear inside quotes. You can specify several name and value pair arguments in any order as `Name1, Value1, ..., NameN, ValueN`.

Use `Name`, `Value` arguments to specify writable properties on page 1-1400 of recursiveBJ System object during object creation. For example, `obj = recursiveBJ([1 1 1 2 1], 'EstimationMethod', 'Gradient')` creates a System object to estimate a Box-Jenkins polynomial model using the 'Gradient' recursive estimation algorithm.

## Properties

`recursiveBJ` System object properties consist of read-only and writable properties. The writable properties are tunable and nontunable properties. The nontunable properties cannot be changed when the object is locked, that is, after you use the `step` command.

Use `Name`, `Value` arguments to specify writable properties of `recursiveBJ` objects during object creation. After object creation, use dot notation to modify the tunable properties.

```
obj = recursiveBJ;  
obj.ForgettingFactor = 0.99;
```

### **B**

Estimated coefficients of polynomial  $B(q)$ , returned as a vector of real values specified in order of ascending powers of  $q^{-1}$ .

`B` is a read-only property and is initially empty after you create the object. It is populated after you use the `step` command for online parameter estimation.

### **C**

Estimated coefficients of polynomial  $C(q)$ , returned as a vector of real values specified in order of ascending powers of  $q^{-1}$ .

`C` is a read-only property and is initially empty after you create the object. It is populated after you use the `step` command for online parameter estimation.

### **D**

Estimated coefficients of polynomial  $D(q)$ , returned as a vector of real values specified in order of ascending powers of  $q^{-1}$ .

`D` is a read-only property and is initially empty after you create the object. It is populated after you use the `step` command for online parameter estimation.

### **F**

Estimated coefficients of polynomial  $F(q)$ , returned as a vector of real values specified in order of ascending powers of  $q^{-1}$ .

F is a read-only property and is initially empty after you create the object. It is populated after you use the `step` command for online parameter estimation.

### **InitialB**

Initial values for the coefficients of polynomial  $B(q)$  of order `nb-1`, specified as a row vector of length `nb+nk`, with `nk` leading zeros. `nk` is the input-output delay. Specify the coefficients in order of ascending powers of  $q^{-1}$ .

If the initial guesses are much smaller than the default `InitialParameterCovariance`, 10000, the initial guesses are given less importance during estimation. In that case, specify a smaller initial parameter covariance.

`InitialB` is a tunable property. You can change it when the object is in a locked state.

**Default:** `[0 eps]`

### **InitialC**

Initial values for the coefficients of polynomial  $C(q)$  of order `nc`, specified as a row vector of length `nc+1`, with 1 as the first element. Specify the coefficients in order of ascending powers of  $q^{-1}$ .

The coefficients in `InitialC` must define a stable discrete-time polynomial with roots within a unit circle. For example,

```
InitialC = [1 0.5 0.5];
all(abs(roots(InitialC))<1)
```

```
ans =
```

```
1
```

If the initial guesses are much smaller than the default `InitialParameterCovariance`, 10000, the initial guesses are given less importance during estimation. In that case, specify a smaller initial parameter covariance.

`InitialC` is a tunable property. You can change it when the object is in a locked state.

**Default:** `[1 eps]`

## InitialD

Initial values for the coefficients of polynomial  $D(q)$  of order  $nd$ , specified as a row vector of length  $nd+1$ , with 1 as the first element. Specify the coefficients in order of ascending powers of  $q^{-1}$ .

The coefficients in `InitialD` must define a stable discrete-time polynomial with roots within a unit circle. For example,

```
InitialD = [1 0.9 0.8];  
all(abs(roots(InitialD))<1)
```

```
ans =
```

```
1
```

If the initial guesses are much smaller than the default `InitialParameterCovariance`, 10000, the initial guesses are given less importance during estimation. In that case, specify a smaller initial parameter covariance.

`InitialD` is a tunable property. You can change it when the object is in a locked state.

**Default:** [1 eps]

## InitialF

Initial values for the coefficients of polynomial  $F(q)$  of order  $nf$ , specified as a row vector of length  $nf+1$ , with 1 as the first element. Specify the coefficients in order of ascending powers of  $q^{-1}$ .

The coefficients in `InitialF` must define a stable discrete-time polynomial with roots within a unit circle. For example,

```
InitialF = [1 0.9 0.8];  
all(abs(roots(InitialF))<1)
```

```
ans =
```

```
1
```

If the initial guesses are much smaller than the default `InitialParameterCovariance`, 10000, the initial guesses are given less importance during estimation. In that case, specify a smaller initial parameter covariance.

InitialF is a tunable property. You can change it when the object is in a locked state.

**Default:** [1 eps]

### ParameterCovariance

Estimated covariance  $P$  of the parameters, returned as an  $N$ -by- $N$  symmetric positive-definite matrix.  $N$  is the number of parameters to be estimated. The software computes  $P$  assuming that the residuals (difference between estimated and measured outputs) are white noise, and the variance of these residuals is 1. ParameterCovariance is applicable only when EstimationMethod is 'ForgettingFactor' or 'KalmanFilter'.

The interpretation of  $P$  depends on the estimation method:

- 'ForgettingFactor' —  $R_2/2 * P$  is approximately equal to the covariance matrix of the estimated parameters, where  $R_2$  is the true variance of the residuals.
- 'KalmanFilter' —  $R_2 * P$  is the covariance matrix of the estimated parameters, and  $R_1 / R_2$  is the covariance matrix of the parameter changes. Where,  $R_1$  is the covariance matrix that you specify in ProcessNoiseCovariance.

ParameterCovariance is a read-only property and is initially empty after you create the object. It is populated after you use the `step` command for online parameter estimation.

### InitialParameterCovariance

Covariance of the initial parameter estimates, specified as one of the following:

- Real positive scalar,  $\alpha$  — Covariance matrix is an  $N$ -by- $N$  diagonal matrix, with  $\alpha$  as the diagonal elements.  $N$  is the number of parameters to be estimated.
- Vector of real positive scalars,  $[\alpha_1, \dots, \alpha_N]$  — Covariance matrix is an  $N$ -by- $N$  diagonal matrix, with  $[\alpha_1, \dots, \alpha_N]$  as the diagonal elements.
- $N$ -by- $N$  symmetric positive-definite matrix.

InitialParameterCovariance represents the uncertainty in the initial parameter estimates. For large values of InitialParameterCovariance, less importance is placed on the initial parameter values and more on the measured data during beginning of estimation using `step`.

Use only when EstimationMethod is 'ForgettingFactor' or 'KalmanFilter'.

`InitialParameterCovariance` is a tunable property. You can change it when the object is in a locked state.

**Default:** 10000

### **EstimationMethod**

Recursive estimation algorithm used for online estimation of model parameters, specified as one of the following values:

- 'ForgettingFactor' — Algorithm used for parameter estimation
- 'KalmanFilter' — Algorithm used for parameter estimation
- 'NormalizedGradient' — Algorithm used for parameter estimation
- 'Gradient' — Unnormalized gradient algorithm used for parameter estimation

Forgetting factor and Kalman filter algorithms are more computationally intensive than gradient and unnormalized gradient methods. However, they have better convergence properties. For information about these algorithms, see “Recursive Algorithms for Online Parameter Estimation”.

`EstimationMethod` is a nontunable property. You cannot change it during execution, that is after the object is locked using the `step` command. If you want to deploy code using MATLAB Coder, `EstimationMethod` can only be assigned once.

**Default:** 'ForgettingFactor'

### **ForgettingFactor**

Forgetting factor,  $\lambda$ , relevant for parameter estimation, specified as a scalar in the range (0,1].

Suppose that the system remains approximately constant over  $T_0$  samples. You can choose  $\lambda$  such that:

$$T_0 = \frac{1}{1 - \lambda}$$

- Setting  $\lambda = 1$  corresponds to “no forgetting” and estimating constant coefficients.
- Setting  $\lambda < 1$  implies that past measurements are less significant for parameter estimation and can be “forgotten”. Set  $\lambda < 1$  to estimate time-varying coefficients.

Typical choices of  $\lambda$  are in the range [0.98 0.995].

Use only when EstimationMethod is 'ForgettingFactor'.

ForgettingFactor is a tunable property. You can change it when the object is in a locked state.

**Default:** 1

### EnableAdapation

Enable or disable parameter estimation, specified as one of the following:

- `true` or `1`— The `step` command estimates the parameter values for that time step and updates the parameter values.
- `false` or `0` — The `step` command does not update the parameters for that time step and instead outputs the last estimated value. You can use this option when your system enters a mode where the parameter values do not vary with time.

---

**Note** If you set `EnableAdapation` to `false`, you must still execute the `step` command. Do not skip `step` to keep parameter values constant, because parameter estimation depends on current and past I/O measurements. `step` ensures past I/O data is stored, even when it does not update the parameters.

---

`EnableAdapation` is a tunable property. You can change it when the object is in a locked state.

**Default:** `true`

### DataType

Floating point precision of parameters, specified as one of the following values:

- `'double'` — Double-precision floating point
- `'single'` — Single-precision floating point

Setting `DataType` to `'single'` saves memory, but leads to loss of precision. Specify `DataType` based on the precision required by the target processor where you will deploy generated code.

`DataType` is a nontunable property. It can only be set during object construction using `Name, Value` arguments and cannot be changed afterward.

**Default:** 'double'

### **ProcessNoiseCovariance**

Covariance matrix of parameter variations, specified as one of the following:

- Real nonnegative scalar,  $\alpha$  — Covariance matrix is an  $N$ -by- $N$  diagonal matrix, with  $\alpha$  as the diagonal elements.
- Vector of real nonnegative scalars,  $[\alpha_1, \dots, \alpha_N]$  — Covariance matrix is an  $N$ -by- $N$  diagonal matrix, with  $[\alpha_1, \dots, \alpha_N]$  as the diagonal elements.
- $N$ -by- $N$  symmetric positive semidefinite matrix.

$N$  is the number of parameters to be estimated.

ProcessNoiseCovariance is applicable when EstimationMethod is 'KalmanFilter'.

Kalman filter algorithm treats the parameters as states of a dynamic system and estimates these parameters using a Kalman filter. ProcessNoiseCovariance is the covariance of the process noise acting on these parameters. Zero values in the noise covariance matrix correspond to estimating constant coefficients. Values larger than 0 correspond to time-varying parameters. Use large values for rapidly changing parameters. However, the larger values result in noisier parameter estimates.

ProcessNoiseCovariance is a tunable property. You can change it when the object is in a locked state.

**Default:** 0.1

### **AdaptationGain**

Adaptation gain,  $\gamma$ , used in gradient recursive estimation algorithms, specified as a positive scalar.

AdaptationGain is applicable when EstimationMethod is 'Gradient' or 'NormalizedGradient'.

Specify a large value for AdaptationGain when your measurements have a high signal-to-noise ratio.

AdaptationGain is a tunable property. You can change it when the object is in a locked state.

**Default:** 1

### **NormalizationBias**

Bias in adaptation gain scaling used in the 'NormalizedGradient' method, specified as a nonnegative scalar.

NormalizationBias is applicable when EstimationMethod is 'NormalizedGradient'.

The normalized gradient algorithm divides the adaptation gain at each step by the square of the two-norm of the gradient vector. If the gradient is close to zero, this can cause jumps in the estimated parameters. NormalizationBias is the term introduced in the denominator to prevent these jumps. Increase NormalizationBias if you observe jumps in estimated parameters.

NormalizationBias is a tunable property. You can change it when the object is in a locked state.

**Default:** eps

## **Output Arguments**

**obj** — System object for online parameter estimation of Box-Jenkins polynomial model

recursiveBJ System object

System object for online parameter estimation of Box-Jenkins polynomial model, returned as a recursiveBJ System object. This object is created using the specified model orders and properties. Use `step` command to estimate the coefficients of the Box-Jenkins model polynomials. You can then access the estimated coefficients and parameter covariance using dot notation. For example, type `obj.F` to view the estimated  $F$  polynomial coefficients.

## **Definitions**

### **Box-Jenkins Polynomial Model Structure**

The general Box-Jenkins model structure is:

$$y(t) = \sum_{i=1}^{nu} \frac{B_i(q)}{F_i(q)} u_i(t - nk_i) + \frac{C(q)}{D(q)} e(t)$$

where  $nu$  is the number of input channels.

The orders of Box-Jenkins model are defined as follows:

$$nb: B(q) = b_1 + b_2q^{-1} + \dots + b_{nb}q^{-nb+1}$$

$$nc: C(q) = 1 + c_1q^{-1} + \dots + c_{nc}q^{-nc}$$

$$nd: D(q) = 1 + d_1q^{-1} + \dots + d_{nd}q^{-nd}$$

$$nf: F(q) = 1 + f_1q^{-1} + \dots + f_{nf}q^{-nf}$$

## Tips

- Starting in R2016b, instead of using the `step` command to update model parameter estimates, you can call the System object with input arguments, as if it were a function. For example, `[B,C,D,F,EstimatedOutput] = step(obj,y,u)` and `[B,C,D,F,EstimatedOutput] = obj(y,u)` perform equivalent operations.

## Extended Capabilities

### C/C++ Code Generation

Generate C and C++ code using MATLAB® Coder™.

Usage notes and limitations:

- For Simulink-based workflows, use Recursive Polynomial Model Estimator.
- For limitations, see “Generate Code for Online Parameter Estimation in MATLAB”.

Supports MATLAB Function block: No

## See Also

Recursive Polynomial Model Estimator | `bj` | `clone` | `isLocked` | `recursiveAR` | `recursiveARMA` | `recursiveARMAX` | `recursiveARX` | `recursiveLS` | `recursiveOE` | `release` | `reset` | `step`

## Topics

[“Perform Online Parameter Estimation at the Command Line”](#)

[“Validate Online Parameter Estimation at the Command Line”](#)

[“What Is Online Estimation?”](#)

[“Recursive Algorithms for Online Parameter Estimation”](#)

**Introduced in R2015b**

## recursiveLS

Create System object for online parameter estimation using recursive least squares algorithm

Use the `recursiveLS` command for parameter estimation with real-time data. If all data necessary for estimation is available at once and you are estimating a time-invariant model, use `mldivide`, `\`.

### Syntax

```
obj = recursiveLS
obj = recursiveLS(Np)
obj = recursiveLS(Np,theta0)
obj = recursiveLS( ____,Name,Value)
```

### Description

`obj = recursiveLS` creates a System object for online parameter estimation of a default single output system that is linear in estimated parameters. Such a system can be represented as:

$$y(t) = H(t)\theta(t) + e(t).$$

Here,  $y$  is the output,  $\theta$  are the parameters,  $H$  are the regressors, and  $e$  is the white-noise disturbance. The default system has one parameter with initial parameter value 1.

After creating the object, use the `step` command to update model parameter estimates using recursive least squares algorithms and real-time data. Alternatively, you can call the object directly. For more information, see “Tips” on page 1-1429.

`obj = recursiveLS(Np)` also specifies the number of parameters to be estimated.

`obj = recursiveLS(Np,theta0)` also specifies the number of parameters and initial values of the parameters.

`obj = recursiveLS( ____,Name,Value)` specifies additional attributes of the system and recursive estimation algorithm using one or more `Name, Value` pair arguments.

## Object Description

recursiveLS creates a System object for online parameter estimation of a single output system that is linear in its parameters.

A System object is a specialized MATLAB object designed specifically for implementing and simulating dynamic systems with inputs that change over time. System objects use internal states to store past behavior, which is used in the next computational step.

After you create a System object, you use commands to process data or obtain information from or about the object. System objects use a minimum of two commands to process data — a constructor to create the object and the `step` command to update object parameters using real-time data. This separation of declaration from execution lets you create multiple, persistent, reusable objects, each with different settings.

You can use the following commands with the online estimation System objects in System Identification Toolbox:

Command	Description
<code>step</code>	Update model parameter estimates using recursive estimation algorithms and real-time data.  <code>step</code> puts the object into a locked state. In a locked state, you cannot change any nontunable properties or input specifications, such as model order, data type, or estimation algorithm. During execution, you can only change tunable properties.
<code>release</code>	Unlock the System object. Use this command to enable setting of nontunable parameters.
<code>reset</code>	Reset the internal states of a locked System object to the initial values, and leave the object locked.

Command	Description
clone	Create another System object with the same object property values.  Do not create additional objects using syntax <code>obj2 = obj</code> . Any changes made to the properties of the new object created this way ( <code>obj2</code> ) also change the properties of the original object ( <code>obj</code> ).
isLocked	Query locked status for input attributes and nontunable properties of the System object.

Use the `recursiveLS` command to create an online estimation System object. Then estimate the system parameters (`theta`) and output using the `step` command with regressors and incoming output data, `H` and `y`.

```
[theta,EstimatedOutput] = step(obj,y,H)
```

For `recursiveLS` object properties, see “Properties” on page 1-1420.

## Examples

### Create System Object for Online Estimation Using Recursive Least Squares Algorithm

```
obj = recursiveLS
```

```
obj =  
    recursiveLS with properties:  
  
        NumberOfParameters: 1  
            Parameters: []  
        InitialParameters: 1  
        ParameterCovariance: []  
    InitialParameterCovariance: 10000  
        EstimationMethod: 'ForgettingFactor'  
        ForgettingFactor: 1  
        EnableAdaptation: true  
        History: 'Infinite'
```

```
InputProcessing: 'Sample-based'
DataType: 'double'
```

### Estimate Parameters of System Using Recursive Least Squares Algorithm

The system has two parameters and is represented as:

$$y(t) = a_1 u(t) + a_2 u(t-1)$$

Here,

- $u$  and  $y$  are the real-time input and output data, respectively.
- $u(t)$  and  $u(t-1)$  are the regressors,  $H$ , of the system.
- $a_1$  and  $a_2$  are the parameters,  $\theta$ , of the system.

Create a System object for online estimation using the recursive least squares algorithm.

```
obj = recursiveLS(2);
```

Load the estimation data, which for this example is a static data set.

```
load iddata3
input = z3.u;
output = z3.y;
```

Create a variable to store  $u(t-1)$ . This variable is updated at each time step.

```
oldInput = 0;
```

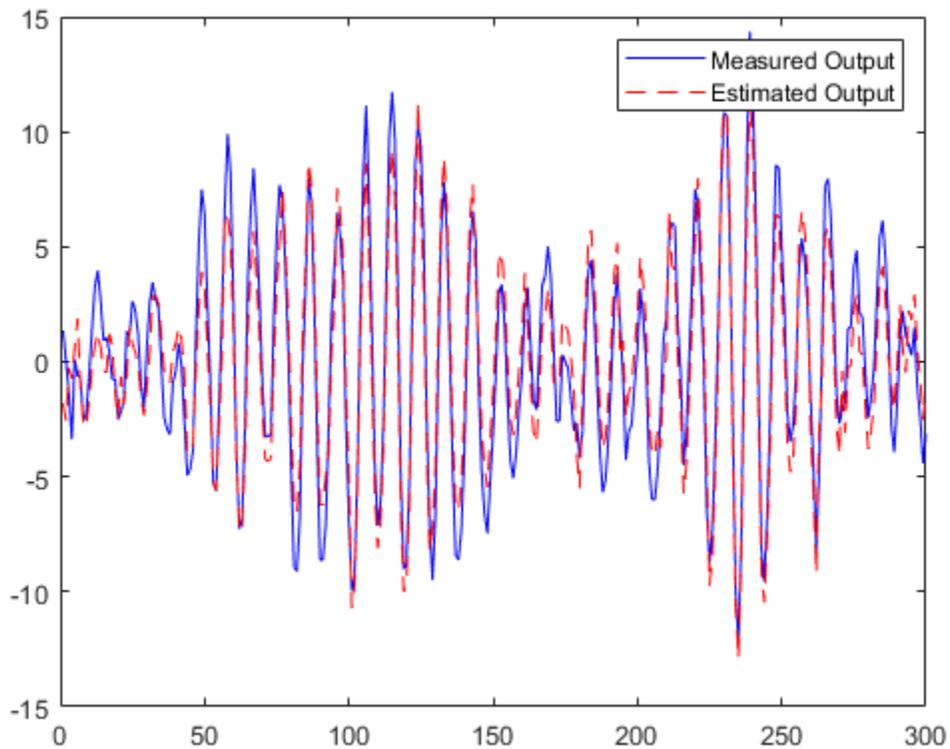
Estimate the parameters and output using `step` and input-output data, maintaining the current regressor pair in  $H$ . Invoke the `step` function implicitly by calling the `obj` system object with input arguments.

```
for i = 1:numel(input)
    H = [input(i) oldInput];
    [theta, EstimatedOutput] = obj(output(i),H);
    estimatedOut(i) = EstimatedOutput;
```

```
    theta_est(i,:) = theta;  
    oldInput = input(i);  
end
```

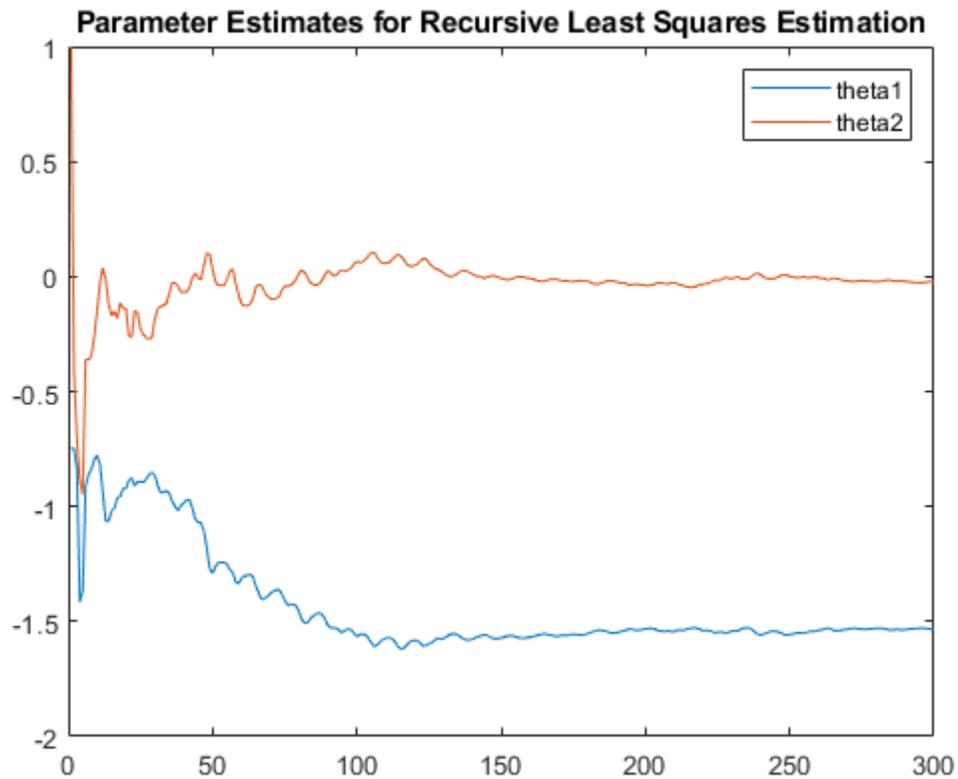
Plot the measured and estimated output data.

```
numSample = 1:numel(input);  
plot(numSample,output,'b',numSample,estimatedOut,'r--');  
legend('Measured Output','Estimated Output');
```



Plot the parameters.

```
plot(numSample,theta_est(:,1),numSample,theta_est(:,2))  
title('Parameter Estimates for Recursive Least Squares Estimation')  
legend("theta1", "theta2")
```



View the final estimates.

```
theta_final = theta
```

```
theta_final = 2x1
```

```
-1.5322
```

```
-0.0235
```

## Use Frame-Based Data for Recursive Least Squares Estimation

Use frame-based signals with the `recursiveLS` command. Machine interfaces often provide sensor data in frames containing multiple samples, rather than in individual samples. The `recursiveLS` object accepts these frames directly when you set `InputProcessing` to `Frame-based`.

The object uses the same estimation algorithms for sample-based and frame-based input processing. The estimation results are identical. There are some special considerations, however, for working with frame-based inputs.

This example is the frame-based version of the sample-based `recursiveLS` example in “Estimate Parameters of System Using Recursive Least Squares Algorithm” on page 1-1413.

The system has two parameters and is represented as:

$$y(t) = a_1 u(t) + a_2 u(t-1)$$

Here,

- $u$  and  $y$  are the real-time input and output data, respectively.
- $u(t)$  and  $u(t-1)$  are the regressors,  $H$ , of the system.
- $a_1$  and  $a_2$  are the parameters,  $\theta$ , of the system.

Create a System object for online estimation using the recursive least squares algorithm.

```
obj_f = recursiveLS(2, 'InputProcessing', 'Frame-Based');
```

Load the data, which contains input and output time series signals. Each signal consists of 30 frames and each frame contains ten individual time samples.

```
load iddata3_frames input_sig_frame output_sig_frame
input = input_sig_frame.data;
output = output_sig_frame.data;
numframes = size(input,3)

numframes = 30

mframe = size(input,1)
```

```
mframe = 10
```

Initialize the regressor frame, which for a given frame, is of the form

$$H_f = \begin{bmatrix} u_1 & u_0 \\ u_2 & u_1 \\ \vdots & \vdots \\ u_{10} & u_9 \end{bmatrix},$$

where the most recent point in the frame is  $u_{10}$ .

```
Hframe = zeros(10,2);
```

For this first-order example, the regressor frame includes one point from the previous frame. Initialize this point.

```
oldInput = 0;
```

Estimate the parameters and output using `step` and input-output data, maintaining the current regressor frame in `Hframe`.

- The input and output arrays have three dimensions. The third dimension is the frame index, and the first two dimensions represent the contents of individual frames.
- Use the `circshift` function to populate the second column of `Hframe` with the past input value for each regressor pair by shifting the input vector by one position.
- Populate the `Hframe` element holding the oldest value, `Hframe(1,2)`, with the regressor value stored from the previous frame.
- Invoke the `step` function implicitly by calling the `obj` system object with input arguments. The `step` function is compatible with frames, so no loop function within the frame is necessary.
- Save the most recent input value to use for the next frame calculation.

```
EstimatedOutput = zeros(10,1,30);
```

```
theta = zeros(2,30);
```

```
for i = 1:numframes
```

```
    Hframe = [input(:,:,i) circshift(input(:,:,i),1)];
```

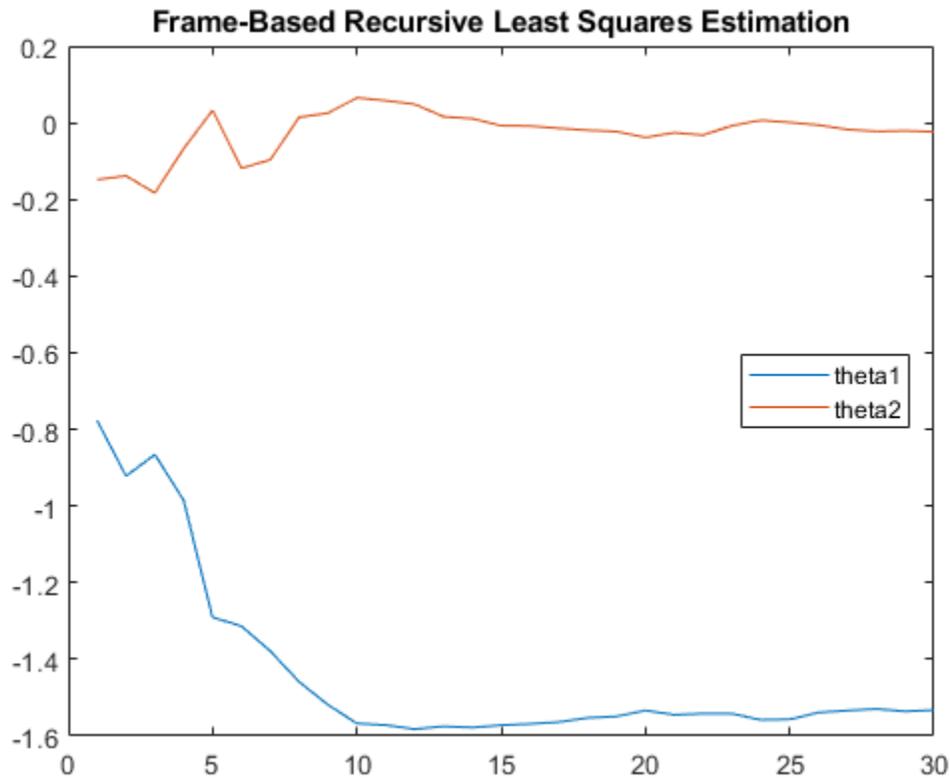
```
    Hframe(1,2) = oldInput;
```

```
    [theta(:,i), EstimatedOutput(:,:,i)] = obj_f(output(:,:,i),Hframe);
```

```
        oldInput = input(10,:,i);  
    end
```

Plot the parameters.

```
theta1 = theta(1,:);  
theta2 = theta(2,:);  
iframe = 1:numframes;  
plot(iframe,theta1,iframe,theta2)  
title('Frame-Based Recursive Least Squares Estimation')  
legend('theta1','theta2','location','best')
```



View the final estimates.

```
theta_final = theta(:,numframes)
```

```
theta_final = 2x1
    -1.5322
    -0.0235
```

The final estimates are identical to the sample-based estimation.

### Specify Initial Parameters for Online Estimation Using Recursive Least Squares Algorithm

Create System object for online parameter estimation using recursive least squares algorithm of a system with two parameters and known initial parameter values.

```
obj = recursiveLS(2,[0.8 1], 'InitialParameterCovariance', 0.1);
```

`InitialParameterCovariance` represents the uncertainty in your guess for the initial parameters. Typically, the default `InitialParameterCovariance` (10000) is too large relative to the parameter values. This results in initial guesses being given less importance during estimation. If you have confidence in the initial parameter guesses, specify a smaller initial parameter covariance.

## Input Arguments

### **Np** — Number of parameters

positive integer

Number of parameters in the system, specified as a positive integer.

### **theta0** — Initial value of parameters

scalar | vector of real values

Initial value of parameters, specified as one of the following:

- Scalar — All the parameters have the same initial value.
- Vector of real values of length `Np`— The  $i$ th parameter has initial value `theta0(i)`.

The default initial value for all parameters is 1.

---

**Note** If the initial parameter values are much smaller than `InitialParameterCovariance`, these initial values are given less importance during estimation. Specify a smaller initial parameter covariance if you have high confidence in the initial parameter values. This statement applies only for infinite-history estimation. Finite-history estimation does not use `InitialParameterCovariance`.

---

## Name-Value Pair Arguments

Specify optional comma-separated pairs of `Name`, `Value` arguments. `Name` is the argument name and `Value` is the corresponding value. `Name` must appear inside quotes. You can specify several name and value pair arguments in any order as `Name1, Value1, . . . , NameN, ValueN`.

Use `Name`, `Value` arguments to specify writable properties on page 1-1420 of `recursiveLS` System object during object creation. For example, `obj = recursiveLS(2, 'EstimationMethod', 'Gradient')` creates a System object to estimate the system parameters using the 'Gradient' recursive estimation algorithm.

## Properties

`recursiveLS` System object properties consist of read-only and writable properties. The writable properties are tunable and nontunable properties. The nontunable properties cannot be changed when the object is locked, that is, after you use the `step` command.

Use `Name`, `Value` arguments to specify writable properties of `recursiveLS` objects during object creation. After object creation, use dot notation to modify the tunable properties.

```
obj = recursiveLS;  
obj.ForgettingFactor = 0.99;
```

### NumberOfParameters

Number of parameters to be estimated, returned as a positive integer.

`NumberOfParameters` is a read-only property. If `Np` is specified during object construction, `NumberOfParameters` takes the value assigned to `Np`.

**Default:** 1

## Parameters

Estimated parameters, returned as a column vector of real values.

Parameters is a read-only property and is initially empty after you create the object. It is populated after you use the `step` command for online parameter estimation.

## InitialParameters

Initial values of parameters, specified as one of the following:

- Scalar — All the parameters have the same initial value.
- Vector of real values of length  $N_p$ — The  $i$ th parameter has initial value `InitialParameters(i)`.

If the initial parameter values are much smaller than `InitialParameterCovariance`, these initial values are given less importance during estimation. Specify a smaller initial parameter covariance if you have high confidence in initial parameter values. This statement applies only for infinite-history estimation. Finite-history estimation does not use `InitialParameterCovariance`.

`InitialParameters` is a tunable property. You can change `InitialParameters` when the object is in a locked state.

### Default: 1

## InitialOutputs

Initial values of the outputs buffer in finite-history estimation, specified as  $\emptyset$  or as a  $W$ -by-1 vector, where  $W$  is the window length.

The `InitialOutputs` property provides a means of controlling the initial behavior of the algorithm.

When `InitialOutputs` is set to  $\emptyset$ , the object populates the buffer with zeros.

If the initial buffer is set to  $\emptyset$  or does not contain enough information, you see a warning message during the initial phase of your estimation. The warning should clear after a few cycles. The number of cycles it takes for sufficient information to be buffered depends upon the order of your polynomials and your input delays. If the warning persists, you should evaluate the content of your signals.

Specify `InitialOutputs` only when `History` is `Finite`.

`InitialOutputs` is a tunable property. You can change `InitialOutputs` when the object is in a locked state.

**Default:** 0

## **InitialRegressors**

Initial values of the regressors buffer in finite-history estimation, specified as 0 or as a  $W$ -by- $N_p$  matrix, where  $W$  is the window length and  $N_p$  is the number of parameters.

The `InitialRegressors` property provides a means of controlling the initial behavior of the algorithm.

When the `InitialRegressors` is set to 0, the object populates the buffer with zeros.

If the initial buffer is set to 0 or does not contain enough information, you see a warning message during the initial phase of your estimation. The warning should clear after a few cycles. The number of cycles it takes for sufficient information to be buffered depends upon the order of your polynomials and your input delays. If the warning persists, you should evaluate the content of your signals.

Specify `InitialRegressors` only when `History` is `Finite`.

`InitialRegressors` is a tunable property. You can change `InitialRegressors` when the object is in a locked state.

**Default:** 0

## **ParameterCovariance**

Estimated covariance  $P$  of the parameters, returned as an  $N$ -by- $N$  symmetric positive-definite matrix.  $N$  is the number of parameters to be estimated. The software computes  $P$  assuming that the residuals (difference between estimated and measured outputs) are white noise, and the variance of these residuals is 1.

`ParameterCovariance` is applicable only when `EstimationMethod` is 'ForgettingFactor' or 'KalmanFilter' or when `History` is `Finite`.

The interpretation of  $P$  depends on your settings for the `History` and `EstimationMethod` properties.

- If `History` is `Infinite`, then your `EstimationMethod` selection results in one of the following:

- 'ForgettingFactor' —  $(R_2/2)P$  is approximately equal to the covariance matrix of the estimated parameters, where  $R_2$  is the true variance of the residuals.
- 'KalmanFilter' —  $R_2P$  is the covariance matrix of the estimated parameters, and  $R_1/R_2$  is the covariance matrix of the parameter changes. Here,  $R_1$  is the covariance matrix that you specify in `ProcessNoiseCovariance`.
- If `History` is `Finite` (sliding-window estimation) —  $R_2P$  is the covariance of the estimated parameters. The sliding-window algorithm does not use this covariance in the parameter-estimation process. However, the algorithm does compute the covariance for output so that you can use it for statistical evaluation.

`ParameterCovariance` is a read-only property and is initially empty after you create the object. It is populated after you use the `step` command for online parameter estimation.

### **InitialParameterCovariance**

Covariance of the initial parameter estimates, specified as one of the following:

- Real positive scalar,  $\alpha$  — Covariance matrix is an  $N$ -by- $N$  diagonal matrix, with  $\alpha$  as the diagonal elements.  $N$  is the number of parameters to be estimated.
- Vector of real positive scalars,  $[\alpha_1, \dots, \alpha_N]$  — Covariance matrix is an  $N$ -by- $N$  diagonal matrix, with  $[\alpha_1, \dots, \alpha_N]$  as the diagonal elements.
- $N$ -by- $N$  symmetric positive-definite matrix.

`InitialParameterCovariance` represents the uncertainty in the initial parameter estimates. For large values of `InitialParameterCovariance`, less importance is placed on the initial parameter values and more on the measured data during beginning of estimation using `step`.

Use only when `EstimationMethod` is 'ForgettingFactor' or 'KalmanFilter'.

`InitialParameterCovariance` is a tunable property. You can change it when the object is in a locked state.

**Default:** 10000

### **EstimationMethod**

Recursive least squares estimation algorithm used for online estimation of model parameters, specified as one of the following values:

- 'ForgettingFactor' — Algorithm used for parameter estimation
- 'KalmanFilter' — Algorithm used for parameter estimation
- 'NormalizedGradient' — Algorithm used for parameter estimation
- 'Gradient' — Unnormalized gradient algorithm used for parameter estimation

Forgetting factor and Kalman filter algorithms are more computationally intensive than gradient and unnormalized gradient methods. However, they have better convergence properties. For information about these algorithms, see “Recursive Algorithms for Online Parameter Estimation”.

These methods all use an infinite data history, and are available only when `History` is 'Infinite'.

`EstimationMethod` is a nontunable property. You cannot change it during execution, that is, after the object is locked using the `step` command.

**Default:** Forgetting Factor

### ForgettingFactor

Forgetting factor,  $\lambda$ , relevant for parameter estimation, specified as a scalar in the range (0,1].

Suppose that the system remains approximately constant over  $T_0$  samples. You can choose  $\lambda$  such that:

$$T_0 = \frac{1}{1 - \lambda}$$

- Setting  $\lambda = 1$  corresponds to “no forgetting” and estimating constant coefficients.
- Setting  $\lambda < 1$  implies that past measurements are less significant for parameter estimation and can be “forgotten”. Set  $\lambda < 1$  to estimate time-varying coefficients.

Typical choices of  $\lambda$  are in the range [0.98 0.995].

Use only when `EstimationMethod` is 'ForgettingFactor'.

`ForgettingFactor` is a tunable property. You can change it when the object is in a locked state.

**Default:** 1

## EnableAdapation

Enable or disable parameter estimation, specified as one of the following:

- `true` or `1`— The `step` command estimates the parameter values for that time step and updates the parameter values.
- `false` or `0` — The `step` command does not update the parameters for that time step and instead outputs the last estimated value. You can use this option when your system enters a mode where the parameter values do not vary with time.

---

**Note** If you set `EnableAdapation` to `false`, you must still execute the `step` command. Do not skip `step` to keep parameter values constant, because parameter estimation depends on current and past I/O measurements. `step` ensures past I/O data is stored, even when it does not update the parameters.

---

`EnableAdapation` is a tunable property. You can change it when the object is in a locked state.

**Default:** `true`

## DataType

Floating point precision of parameters, specified as one of the following values:

- `'double'` — Double-precision floating point
- `'single'` — Single-precision floating point

Setting `DataType` to `'single'` saves memory, but leads to loss of precision. Specify `DataType` based on the precision required by the target processor where you will deploy generated code.

`DataType` is a nontunable property. It can only be set during object construction using `Name, Value` arguments and cannot be changed afterward.

**Default:** `'double'`

## ProcessNoiseCovariance

Covariance matrix of parameter variations, specified as one of the following:

- Real nonnegative scalar,  $\alpha$  — Covariance matrix is an  $N$ -by- $N$  diagonal matrix, with  $\alpha$  as the diagonal elements.

- Vector of real nonnegative scalars,  $[\alpha_1, \dots, \alpha_N]$  — Covariance matrix is an  $N$ -by- $N$  diagonal matrix, with  $[\alpha_1, \dots, \alpha_N]$  as the diagonal elements.
- $N$ -by- $N$  symmetric positive semidefinite matrix.

$N$  is the number of parameters to be estimated.

`ProcessNoiseCovariance` is applicable when `EstimationMethod` is `'KalmanFilter'`.

Kalman filter algorithm treats the parameters as states of a dynamic system and estimates these parameters using a Kalman filter. `ProcessNoiseCovariance` is the covariance of the process noise acting on these parameters. Zero values in the noise covariance matrix correspond to estimating constant coefficients. Values larger than 0 correspond to time-varying parameters. Use large values for rapidly changing parameters. However, the larger values result in noisier parameter estimates.

`ProcessNoiseCovariance` is a tunable property. You can change it when the object is in a locked state.

**Default:** 0.1

### **AdaptationGain**

Adaptation gain,  $\gamma$ , used in gradient recursive estimation algorithms, specified as a positive scalar.

`AdaptationGain` is applicable when `EstimationMethod` is `'Gradient'` or `'NormalizedGradient'`.

Specify a large value for `AdaptationGain` when your measurements have a high signal-to-noise ratio.

`AdaptationGain` is a tunable property. You can change it when the object is in a locked state.

**Default:** 1

### **NormalizationBias**

Bias in adaptation gain scaling used in the `'NormalizedGradient'` method, specified as a nonnegative scalar.

`NormalizationBias` is applicable when `EstimationMethod` is `'NormalizedGradient'`.

The normalized gradient algorithm divides the adaptation gain at each step by the square of the two-norm of the gradient vector. If the gradient is close to zero, this can cause jumps in the estimated parameters. `NormalizationBias` is the term introduced in the denominator to prevent these jumps. Increase `NormalizationBias` if you observe jumps in estimated parameters.

`NormalizationBias` is a tunable property. You can change it when the object is in a locked state.

**Default:** `eps`

### History

Data history type defining which type of recursive algorithm you use, specified as:

- `'Infinite'` — Use an algorithm that aims to minimize the error between the observed and predicted outputs for all time steps from the beginning of the simulation.
- `'Finite'` — Use an algorithm that aims to minimize the error between the observed and predicted outputs for a finite number of past time steps.

Algorithms with infinite history aim to produce parameter estimates that explain all data since the start of the simulation. These algorithms still use a fixed amount of memory that does not grow over time. The object provides multiple algorithms of the `'Infinite'` `History` type. Specifying this option activates the `EstimationMethod` property with which you specify an algorithm.

Algorithms with finite history aim to produce parameter estimates that explain only a finite number of past data samples. This method is also called sliding-window estimation. The object provides one algorithm of the `'Finite'` type. Specifying this option activates the `WindowLength` property that sizes the window.

For more information on recursive estimation methods, see “Recursive Algorithms for Online Parameter Estimation”.

`History` is a nontunable property. It can be set only during object construction using `Name, Value` arguments and cannot be changed afterward.

**Default:** `'Infinite'`

## WindowLength

Window size determining the number of time samples to use for the sliding-window estimation method, specified as a positive integer. Specify `WindowLength` only when `History` is `Finite`.

Choose a window size that balances estimation performance with computational and memory burden. Sizing factors include the number and time variance of the parameters in your model. Always specify **Window Length** in samples, even if you are using frame-based input processing.

`WindowLength` must be greater than or equal to the number of estimated parameters.

Suitable window length is independent of whether you are using sample-based or frame-based input processing (see `InputProcessing`). However, when using frame-based processing, your window length must be greater than or equal to the number of samples (time steps) contained in the frame.

`WindowLength` is a nontunable property. It can be set only during object construction using `Name, Value` arguments and cannot be changed afterward.

**Default:** 200

## InputProcessing

Option for sample-based or frame-based input processing, specified as a character vector or string.

- **Sample-based** processing operates on signals streamed one sample at a time.
- **Frame-based** processing operates on signals containing samples from multiple time steps. Many machine sensor interfaces package multiple samples and transmit these samples together in frames. **Frame-based** processing allows you to input this data directly without having to first unpack it.

Your `InputProcessing` specification impacts the dimensions for the input and output signals when using the `step` command:

```
[theta, EstimatedOutput] = step(obj, y, H)
```

- **Sample-based**
  - `y` and `EstimatedOutput` are scalars.

- $H$  is a 1-by- $N_p$  vector, where  $N_p$  is the number of parameters.
- • Frame-based with  $M$  samples per frame
  - $y$  and `EstimatedOutput` are  $M$ -by-1 vectors.
  - $H$  is an  $M$ -by- $N_p$  matrix.

`InputProcessing` is a nontunable property. It can be set only during object construction using `Name`, `Value` arguments and cannot be changed afterward.

**Default:** 'Sample-based'

## Output Arguments

### **obj** — System object for online parameter estimation

recursiveLS System object

System object for online parameter estimation, returned as a recursiveLS System object. Use `step` command to estimate the parameters of the system. You can then access the estimated parameters and parameter covariance using dot notation. For example, type `obj.Parameters` to view the estimated parameters.

## Tips

- Starting in R2016b, instead of using the `step` command to update model parameter estimates, you can call the System object with input arguments, as if it were a function. For example, `[theta,EstimatedOutput] = step(obj,y,H)` and `[theta,EstimatedOutput] = obj(y,H)` perform equivalent operations.

## Extended Capabilities

### **C/C++ Code Generation**

Generate C and C++ code using MATLAB® Coder™.

Usage notes and limitations:

- For Simulink based workflows, use Recursive Least Squares Estimator.
- For limitations, see “Generate Code for Online Parameter Estimation in MATLAB”.

Supports MATLAB Function block: No

### See Also

[Recursive Least Squares Estimator](#) | [clone](#) | [isLocked](#) | [mldivide](#) | [recursiveAR](#) | [recursiveARMA](#) | [recursiveARMAX](#) | [recursiveARX](#) | [recursiveBJ](#) | [recursiveOE](#) | [release](#) | [reset](#) | [step](#)

### Topics

[“Perform Online Parameter Estimation at the Command Line”](#)  
[“Validate Online Parameter Estimation at the Command Line”](#)  
[“Line Fitting with Online Recursive Least Squares Estimation”](#)  
[“What Is Online Estimation?”](#)  
[“Recursive Algorithms for Online Parameter Estimation”](#)

### Introduced in R2015b

## recursiveOE

Create System object for online parameter estimation of Output-Error polynomial model

Use `recursiveOE` command for parameter estimation with real-time data. If all data necessary for estimation is available at once, and you are estimating a time-invariant model, use the offline estimation command, `oe`.

### Syntax

```
obj = recursiveOE
obj = recursiveOE(Orders)
obj = recursiveOE(Orders,B0,F0)
obj = recursiveOE( ____,Name,Value)
```

### Description

`obj = recursiveOE` creates a System object for online parameter estimation of a default single-input-single output (SISO) Output-Error model structure on page 1-1443. The default model structure has polynomials of order 1 and initial polynomial coefficient values `eps`.

After creating the object, use the `step` command to update model parameter estimates using recursive estimation algorithms and real-time data.

`obj = recursiveOE(Orders)` specifies the polynomial orders of the Output-Error model to be estimated.

`obj = recursiveOE(Orders,B0,F0)` specifies the polynomial orders and initial values of the polynomial coefficients. Specify initial values to potentially avoid local minima during estimation. If the initial values are small compared to the default `InitialParameterCovariance` property value, and you have confidence in your initial values, also specify a smaller `InitialParameterCovariance`.

`obj = recursiveOE( ____,Name,Value)` specifies additional attributes of the Output-Error model structure and recursive estimation algorithm using one or more `Name,Value` pair arguments.

## Object Description

`recursiveOE` creates a System object for online parameter estimation of SISO Output-Error polynomial models using a recursive estimation algorithm.

A System object is a specialized MATLAB object designed specifically for implementing and simulating dynamic systems with inputs that change over time. System objects use internal states to store past behavior, which is used in the next computational step.

After you create a System object, you use commands to process data or obtain information from or about the object. System objects use a minimum of two commands to process data — a constructor to create the object and the `step` command to update object parameters using real-time data. This separation of declaration from execution lets you create multiple, persistent, reusable objects, each with different settings.

You can use the following commands with the online estimation System objects in System Identification Toolbox:

Command	Description
<code>step</code>	Update model parameter estimates using recursive estimation algorithms and real-time data.  <code>step</code> puts the object into a locked state. In a locked state, you cannot change any nontunable properties or input specifications, such as model order, data type, or estimation algorithm. During execution, you can only change tunable properties.
<code>release</code>	Unlock the System object. Use this command to enable setting of nontunable parameters.
<code>reset</code>	Reset the internal states of a locked System object to the initial values, and leave the object locked.

Command	Description
clone	Create another System object with the same object property values.  Do not create additional objects using syntax <code>obj2 = obj</code> . Any changes made to the properties of the new object created this way ( <code>obj2</code> ) also change the properties of the original object ( <code>obj</code> ).
isLocked	Query locked status for input attributes and nontunable properties of the System object.

Use the `recursiveOE` command to create an online estimation System object. Then estimate the Output-Error polynomial model parameters (B and F) and output using the `step` command with incoming input and output data, `u` and `y`.

```
[B,F,EstimatedOutput] = step(obj,y,u)
```

For `recursiveOE` object properties, see “Properties” on page 1-1437.

## Examples

### Estimate Output-Error Polynomial Model Online

Create a System object for online parameter estimation of a Output-Error polynomial model using recursive estimation algorithms.

```
obj = recursiveOE;
```

The Output-Error model has a default structure with polynomials of order 1 and initial polynomial coefficient values, `eps`.

Load the estimation data. In this example, use a static data set for illustration.

```
load iddata1 z1;
output = z1.y;
input = z1.u;
```

Estimate Output-Error model parameters online using `step`.

```
for i = 1:numel(input)
[B,F,EstimatedOutput] = step(obj,output(i),input(i));
end
```

View the current estimated values of polynomial F coefficients.

```
obj.F
```

```
ans = 1×2
```

```
1.0000 -0.7618
```

View the current covariance estimate of the parameters.

```
obj.ParameterCovariance
```

```
ans = 2×2
```

```
0.0024 0.0002
0.0002 0.0001
```

View the current estimated output.

```
EstimatedOutput
```

```
EstimatedOutput = -6.6140
```

## Create System Object for Output-Error Model With Known Orders and Delays

Specify Output-Error polynomial model orders and delays.

```
nb = 1;
nf = 2;
nk = 1;
```

Create a System object for online estimation of Output-Error polynomial model with the specified orders and delays.

```
obj = recursiveOE([nb nf nk]);
```

## Create System Object for Output-Error Model With Known Initial Parameters

Specify Output-Error polynomial model orders and delays.

```
nb = 1;  
nf = 2;  
nk = 1;
```

Create a System object for online estimation of Output-Error model with known initial polynomial coefficients.

```
B0 = [0 1];  
F0 = [1 0.7 0.8];  
obj = recursiveOE([nb nf nk],B0,F0);
```

Specify the initial parameter covariance.

```
obj.InitialParameterCovariance = 0.1;
```

`InitialParameterCovariance` represents the uncertainty in your guess for the initial parameters. Typically, the default `InitialParameterCovariance` (10000) is too large relative to the parameter values. This results in initial guesses being given less importance during estimation. If you have confidence in the initial parameter guesses, specify a smaller initial parameter covariance.

## Specify Estimation Method for Online Estimation of Output-Error Model

Create a System object that uses the unnormalized gradient algorithm for online parameter estimation of an Output-Error model.

```
obj = recursiveOE([1 2 1], 'EstimationMethod', 'Gradient');
```

## Input Arguments

### Orders — Model orders and delays

1-by-3 vector of integers

Model orders and delays of a Output-Error polynomial model on page 1-1443, specified as a 1-by-3 vector of integers, `[nb nf nk]`.

- `nb` — Order of the polynomial  $B(q) + 1$ , specified as a positive integer.
- `nf` — Order of the polynomial  $F(q)$ , specified as a nonnegative integer.
- `nk` — Input-output delay, specified as a positive integer. `nk` is number of input samples that occur before the input affects the output. `nk` is expressed as fixed leading zeros of the  $B$  polynomial.

## **B0, F0 — Initial value of polynomial coefficients**

row vectors of real values | []

Initial value of polynomial coefficients, specified as row vectors of real values with elements in order of ascending powers of  $q^{-1}$ .

- `B0` — Initial guess for the coefficients of the polynomial  $B(q)$ , specified as a 1-by-(`nb` + `nk`) vector with `nk` leading zeros.
- `F0` — Initial guess for the coefficients of the polynomial  $F(q)$ , specified as a 1-by-(`nf` + 1) vector with 1 as the first element.

The coefficients in `F0` must define a stable discrete-time polynomial with roots within a unit disk. For example,

```
F0 = [1 0.5 0.5];  
all(abs(roots(F0))<1)
```

```
ans =
```

```
1
```

Specifying as [], uses the default value of `eps` for the polynomial coefficients.

---

**Note** If the initial guesses are much smaller than the default `InitialParameterCovariance`, 10000, the initial guesses are given less importance during estimation. In that case, specify a smaller initial parameter covariance.

---

## **Name-Value Pair Arguments**

Specify optional comma-separated pairs of `Name`, `Value` arguments. `Name` is the argument name and `Value` is the corresponding value. `Name` must appear inside quotes. You can specify several name and value pair arguments in any order as `Name1, Value1, ..., NameN, ValueN`.

Use `Name`, `Value` arguments to specify writable properties on page 1-1437 of recursiveOE System object during object creation. For example, `obj = recursiveOE([1 2 1], 'EstimationMethod', 'Gradient')` creates a System object to estimate a Output-Error polynomial model using the 'Gradient' recursive estimation algorithm.

## Properties

recursiveOE System object properties consist of read-only and writable properties. The writable properties are tunable and nontunable properties. The nontunable properties cannot be changed when the object is locked, that is, after you use the `step` command.

Use `Name`, `Value` arguments to specify writable properties of recursiveOE objects during object creation. After object creation, use dot notation to modify the tunable properties.

```
obj = recursiveOE;  
obj.ForgettingFactor = 0.99;
```

### **B**

Estimated coefficients of polynomial  $B(q)$ , returned as a vector of real values specified in order of ascending powers of  $q^{-1}$ .

**B** is a read-only property and is initially empty after you create the object. It is populated after you use the `step` command for online parameter estimation.

### **F**

Estimated coefficients of polynomial  $F(q)$ , returned as a vector of real values specified in order of ascending powers of  $q^{-1}$ .

**F** is a read-only property and is initially empty after you create the object. It is populated after you use the `step` command for online parameter estimation.

### **InitialB**

Initial values for the coefficients of polynomial  $B(q)$  of order `nb-1`, specified as a row vector of length `nb+nk`, with `nk` leading zeros. `nk` is the input-output delay. Specify the coefficients in order of ascending powers of  $q^{-1}$ .

If the initial guesses are much smaller than the default `InitialParameterCovariance`, 10000, the initial guesses are given less importance during estimation. In that case, specify a smaller initial parameter covariance.

`InitialB` is a tunable property. You can change it when the object is in a locked state.

**Default:** [0 eps]

### **InitialF**

Initial values for the coefficients of polynomial  $F(q)$  of order `nf`, specified as a row vector of length `nf+1`, with 1 as the first element. Specify the coefficients in order of ascending powers of  $q^{-1}$ .

The coefficients in `InitialF` must define a stable discrete-time polynomial with roots within a unit circle. For example,

```
InitialF = [1 0.9 0.8];  
all(abs(roots(InitialF))<1)
```

```
ans =
```

```
1
```

If the initial guesses are much smaller than the default `InitialParameterCovariance`, 10000, the initial guesses are given less importance during estimation. In that case, specify a smaller initial parameter covariance.

`InitialF` is a tunable property. You can change it when the object is in a locked state.

**Default:** [1 eps]

### **ParameterCovariance**

Estimated covariance  $P$  of the parameters, returned as an  $N$ -by- $N$  symmetric positive-definite matrix.  $N$  is the number of parameters to be estimated. The software computes  $P$  assuming that the residuals (difference between estimated and measured outputs) are white noise, and the variance of these residuals is 1. `ParameterCovariance` is applicable only when `EstimationMethod` is 'ForgettingFactor' or 'KalmanFilter'.

The interpretation of  $P$  depends on the estimation method:

- 'ForgettingFactor' —  $R_2/2 * P$  is approximately equal to the covariance matrix of the estimated parameters, where  $R_2$  is the true variance of the residuals.
- 'KalmanFilter' —  $R_2 * P$  is the covariance matrix of the estimated parameters, and  $R_1 / R_2$  is the covariance matrix of the parameter changes. Where,  $R_1$  is the covariance matrix that you specify in `ProcessNoiseCovariance`.

`ParameterCovariance` is a read-only property and is initially empty after you create the object. It is populated after you use the `step` command for online parameter estimation.

### **InitialParameterCovariance**

Covariance of the initial parameter estimates, specified as one of the following:

- Real positive scalar,  $\alpha$  — Covariance matrix is an  $N$ -by- $N$  diagonal matrix, with  $\alpha$  as the diagonal elements.  $N$  is the number of parameters to be estimated.
- Vector of real positive scalars,  $[\alpha_1, \dots, \alpha_N]$  — Covariance matrix is an  $N$ -by- $N$  diagonal matrix, with  $[\alpha_1, \dots, \alpha_N]$  as the diagonal elements.
- $N$ -by- $N$  symmetric positive-definite matrix.

`InitialParameterCovariance` represents the uncertainty in the initial parameter estimates. For large values of `InitialParameterCovariance`, less importance is placed on the initial parameter values and more on the measured data during beginning of estimation using `step`.

Use only when `EstimationMethod` is 'ForgettingFactor' or 'KalmanFilter'.

`InitialParameterCovariance` is a tunable property. You can change it when the object is in a locked state.

**Default:** 10000

### **EstimationMethod**

Recursive estimation algorithm used for online estimation of model parameters, specified as one of the following values:

- 'ForgettingFactor' — Algorithm used for parameter estimation
- 'KalmanFilter' — Algorithm used for parameter estimation
- 'NormalizedGradient' — Algorithm used for parameter estimation
- 'Gradient' — Unnormalized gradient algorithm used for parameter estimation

Forgetting factor and Kalman filter algorithms are more computationally intensive than gradient and unnormalized gradient methods. However, they have better convergence properties. For information about these algorithms, see “Recursive Algorithms for Online Parameter Estimation”.

`EstimationMethod` is a nontunable property. You cannot change it during execution, that is after the object is locked using the `step` command. If you want to deploy code using MATLAB Coder, `EstimationMethod` can only be assigned once.

**Default:** 'ForgettingFactor'

### **ForgettingFactor**

Forgetting factor,  $\lambda$ , relevant for parameter estimation, specified as a scalar in the range (0,1].

Suppose that the system remains approximately constant over  $T_0$  samples. You can choose  $\lambda$  such that:

$$T_0 = \frac{1}{1 - \lambda}$$

- Setting  $\lambda = 1$  corresponds to “no forgetting” and estimating constant coefficients.
- Setting  $\lambda < 1$  implies that past measurements are less significant for parameter estimation and can be “forgotten”. Set  $\lambda < 1$  to estimate time-varying coefficients.

Typical choices of  $\lambda$  are in the range [0.98 0.995].

Use only when `EstimationMethod` is 'ForgettingFactor'.

`ForgettingFactor` is a tunable property. You can change it when the object is in a locked state.

**Default:** 1

### **EnableAdaption**

Enable or disable parameter estimation, specified as one of the following:

- `true` or 1— The `step` command estimates the parameter values for that time step and updates the parameter values.

- `false` or `0` — The `step` command does not update the parameters for that time step and instead outputs the last estimated value. You can use this option when your system enters a mode where the parameter values do not vary with time.

---

**Note** If you set `EnableAdapation` to `false`, you must still execute the `step` command. Do not skip `step` to keep parameter values constant, because parameter estimation depends on current and past I/O measurements. `step` ensures past I/O data is stored, even when it does not update the parameters.

---

`EnableAdapation` is a tunable property. You can change it when the object is in a locked state.

**Default:** `true`

### **DataType**

Floating point precision of parameters, specified as one of the following values:

- `'double'` — Double-precision floating point
- `'single'` — Single-precision floating point

Setting `DataType` to `'single'` saves memory, but leads to loss of precision. Specify `DataType` based on the precision required by the target processor where you will deploy generated code.

`DataType` is a nontunable property. It can only be set during object construction using `Name, Value` arguments and cannot be changed afterward.

**Default:** `'double'`

### **ProcessNoiseCovariance**

Covariance matrix of parameter variations, specified as one of the following:

- Real nonnegative scalar,  $\alpha$  — Covariance matrix is an  $N$ -by- $N$  diagonal matrix, with  $\alpha$  as the diagonal elements.
- Vector of real nonnegative scalars,  $[\alpha_1, \dots, \alpha_N]$  — Covariance matrix is an  $N$ -by- $N$  diagonal matrix, with  $[\alpha_1, \dots, \alpha_N]$  as the diagonal elements.
- $N$ -by- $N$  symmetric positive semidefinite matrix.

$N$  is the number of parameters to be estimated.

`ProcessNoiseCovariance` is applicable when `EstimationMethod` is `'KalmanFilter'`.

Kalman filter algorithm treats the parameters as states of a dynamic system and estimates these parameters using a Kalman filter. `ProcessNoiseCovariance` is the covariance of the process noise acting on these parameters. Zero values in the noise covariance matrix correspond to estimating constant coefficients. Values larger than 0 correspond to time-varying parameters. Use large values for rapidly changing parameters. However, the larger values result in noisier parameter estimates.

`ProcessNoiseCovariance` is a tunable property. You can change it when the object is in a locked state.

**Default:** 0.1

### **AdaptationGain**

Adaptation gain,  $\gamma$ , used in gradient recursive estimation algorithms, specified as a positive scalar.

`AdaptationGain` is applicable when `EstimationMethod` is `'Gradient'` or `'NormalizedGradient'`.

Specify a large value for `AdaptationGain` when your measurements have a high signal-to-noise ratio.

`AdaptationGain` is a tunable property. You can change it when the object is in a locked state.

**Default:** 1

### **NormalizationBias**

Bias in adaptation gain scaling used in the `'NormalizedGradient'` method, specified as a nonnegative scalar.

`NormalizationBias` is applicable when `EstimationMethod` is `'NormalizedGradient'`.

The normalized gradient algorithm divides the adaptation gain at each step by the square of the two-norm of the gradient vector. If the gradient is close to zero, this can cause jumps in the estimated parameters. `NormalizationBias` is the term introduced in the

denominator to prevent these jumps. Increase `NormalizationBias` if you observe jumps in estimated parameters.

`NormalizationBias` is a tunable property. You can change it when the object is in a locked state.

**Default:** `eps`

## Output Arguments

**obj** — System object for online parameter estimation of Output-Error polynomial model

`recursiveOE` System object

System object for online parameter estimation of SISO Output-Error polynomial model, returned as a `recursiveOE` System object. This object is created using the specified model orders and properties. Use `step` command to estimate the coefficients of the Output-Error model polynomials. You can then access the estimated coefficients and parameter covariance using dot notation. For example, type `obj.B` to view the estimated  $B$  polynomial coefficients.

## Definitions

### Output-Error Model Structure

The general Output-Error model structure is:

$$y(t) = \frac{B(q)}{F(q)} u(t - n_k) + e(t)$$

The orders of the Output-Error model are:

$$\begin{aligned} nb: \quad B(q) &= b_1 + b_2 q^{-1} + \dots + b_{nb} q^{-nb+1} \\ nf: \quad F(q) &= 1 + f_1 q^{-1} + \dots + f_{nf} q^{-nf} \end{aligned}$$

## Tips

- Starting in R2016b, instead of using the `step` command to update model parameter estimates, you can call the System object with input arguments, as if it were a function. For example, `[B,F,EstimatedOutput] = step(obj,y,u)` and `[B,F,EstimatedOutput] = obj(y,u)` perform equivalent operations.

## Extended Capabilities

### C/C++ Code Generation

Generate C and C++ code using MATLAB® Coder™.

Usage notes and limitations:

- For Simulink-based workflows, use Recursive Polynomial Model Estimator.
- For limitations, see “Generate Code for Online Parameter Estimation in MATLAB”.

Supports MATLAB Function block: No

### See Also

[Recursive Polynomial Model Estimator](#) | [clone](#) | [isLocked](#) | [oe](#) | [recursiveAR](#) | [recursiveARMA](#) | [recursiveARMAX](#) | [recursiveARX](#) | [recursiveBJ](#) | [recursiveLS](#) | [release](#) | [reset](#) | [step](#)

### Topics

“Perform Online Parameter Estimation at the Command Line”

“Validate Online Parameter Estimation at the Command Line”

“What Is Online Estimation?”

“Recursive Algorithms for Online Parameter Estimation”

**Introduced in R2015b**

# release

Unlock online parameter estimation System object

## Syntax

```
release(obj)
```

## Description

`release(obj)` unlocks the online parameter estimation System object, `obj`. Use `release` to change nontunable properties of the object.

---

**Note** You can use `release` on a System object in code generated using MATLAB Coder, but once you release its resources, you cannot use that System object again.

---

## Examples

### Unlock Online Estimation System Object

Create a System object™ for online estimation of an ARMAX model with default properties.

```
obj = recursiveARMAX;
```

Estimate model parameters online using `step` and input-output data.

```
[A,B,C,EstimatedOutput] = step(obj,1,1);
```

`step` puts the object in a locked state.

```
L = isLocked(obj)
```

```
L = logical
    1
```

Unlock the object.

```
release(obj)
```

Check the locked status of the object.

```
L = isLocked(obj)
```

```
L = logical
    0
```

## Input Arguments

### **obj** — System object for online parameter estimation

`recursiveAR` object | `recursiveARMA` object | `recursiveARX` object | `recursiveARMAX` object | `recursiveOE` object | `recursiveBJ` object | `recursiveLS` object

System object for online parameter estimation, created using one of the following commands:

- `recursiveAR`
- `recursiveARMA`
- `recursiveARX`
- `recursiveARMAX`
- `recursiveOE`
- `recursiveBJ`
- `recursiveLS`

## See Also

`clone` | `isLocked` | `recursiveAR` | `recursiveARMA` | `recursiveARMAX` | `recursiveARX` | `recursiveBJ` | `recursiveLS` | `recursiveOE` | `reset` | `step`

## **Topics**

“Perform Online Parameter Estimation at the Command Line”

“Validate Online Parameter Estimation at the Command Line”

“What Is Online Estimation?”

**Introduced in R2015b**

## repsys

Replicate and tile models

### Syntax

```
rsys = repsys(sys,[M N])  
rsys = repsys(sys,N)  
rsys = repsys(sys,[M N S1,...,Sk])
```

### Description

`rsys = repsys(sys,[M N])` replicates the model `sys` into an M-by-N tiling pattern. The resulting model `rsys` has `size(sys,1)*M` outputs and `size(sys,2)*N` inputs.

`rsys = repsys(sys,N)` creates an N-by-N tiling.

`rsys = repsys(sys,[M N S1,...,Sk])` replicates and tiles `sys` along both I/O and array dimensions to produce a model array. The indices `S` specify the array dimensions. The size of the array is `[size(sys,1)*M, size(sys,2)*N, size(sys,3)*S1, ...]`.

### Input Arguments

#### **sys**

Model to replicate.

#### **M**

Number of replications of `sys` along the output dimension.

#### **N**

Number of replications of `sys` along the input dimension.

**S**

Numbers of replications of `sys` along array dimensions.

**Output Arguments****rsys**

Model having `size(sys,1)*M` outputs and `size(sys,2)*N` inputs.

If you provide array dimensions `S1, ..., Sk`, `rsys` is an array of dynamic systems which each have `size(sys,1)*M` outputs and `size(sys,2)*N` inputs. The size of `rsys` is `[size(sys,1)*M, size(sys,2)*N, size(sys,3)*S1, ...]`.

**Examples****Replicate SISO Transfer Function to Create MIMO Transfer Function**

Create a single-input single-output (SISO) transfer function.

```
sys = tf(2,[1 3])
```

```
sys =
```

```
      2
-----
s + 3
```

Continuous-time transfer function.

Replicate the SISO transfer function to create a MIMO transfer function that has three inputs and two outputs.

```
rsys = repsys(sys,[2 3])
```

```
rsys =
```

```
From input 1 to output...
      2
1:  -----
```

$$\begin{array}{r} s + 3 \\ 2: \frac{\quad}{s + 3} \end{array}$$

From input 2 to output...

$$\begin{array}{r} 2 \\ 1: \frac{\quad}{s + 3} \end{array}$$

$$\begin{array}{r} 2 \\ 2: \frac{\quad}{s + 3} \end{array}$$

From input 3 to output...

$$\begin{array}{r} 2 \\ 1: \frac{\quad}{s + 3} \end{array}$$

$$\begin{array}{r} 2 \\ 2: \frac{\quad}{s + 3} \end{array}$$

Continuous-time transfer function.

Alternatively, you can obtain the MIMO transfer function as follows:

```
rsys = [sys sys sys; sys sys sys];
```

## Replicate SISO Transfer Function to Create Array of Transfer Functions

Create a SISO transfer function.

```
sys = tf(2,[1 3]);
```

Replicate the transfer function into a 3-by-4 array of two-input, one-output transfer functions.

```
rsys = repsys(sys,[1 2 3 4]);
```

Check the size of rsys.

```
size(rsys)
```

```
3x4 array of transfer functions.  
Each model has 1 outputs and 2 inputs.
```

## Tips

`rsys = repsys(sys,N)` produces the same result as `rsys = repsys(sys,[N N])`.  
To produce a diagonal tiling, use `rsys = sys*eye(N)`.

## See Also

`append`

**Introduced in R2010b**

## resample

Resample time-domain data by decimation or interpolation (requires Signal Processing Toolbox software)

### Syntax

```
resample(data,P,Q)  
resample(data,P,Q,order)
```

### Description

`resample(data,P,Q)` resamples data such that the data is interpolated by a factor  $P$  and then decimated by a factor  $Q$ . `resample(z,1,Q)` results in decimation by a factor  $Q$ .

`resample(data,P,Q,order)` filters the data by applying a filter of specified order before interpolation and decimation.

### Input Arguments

`data`

Name of time-domain `iddata` object. Can be input-output or time-series data.

Data must be sampled at equal time intervals.

`P, Q`

Integers that specify the resampling factor, such that the new sample time is  $Q/P$  times the original one.

$(Q/P) > 1$  results in decimation and  $(Q/P) < 1$  results in interpolation.

`order`

Order of the filters applied before interpolation and decimation.

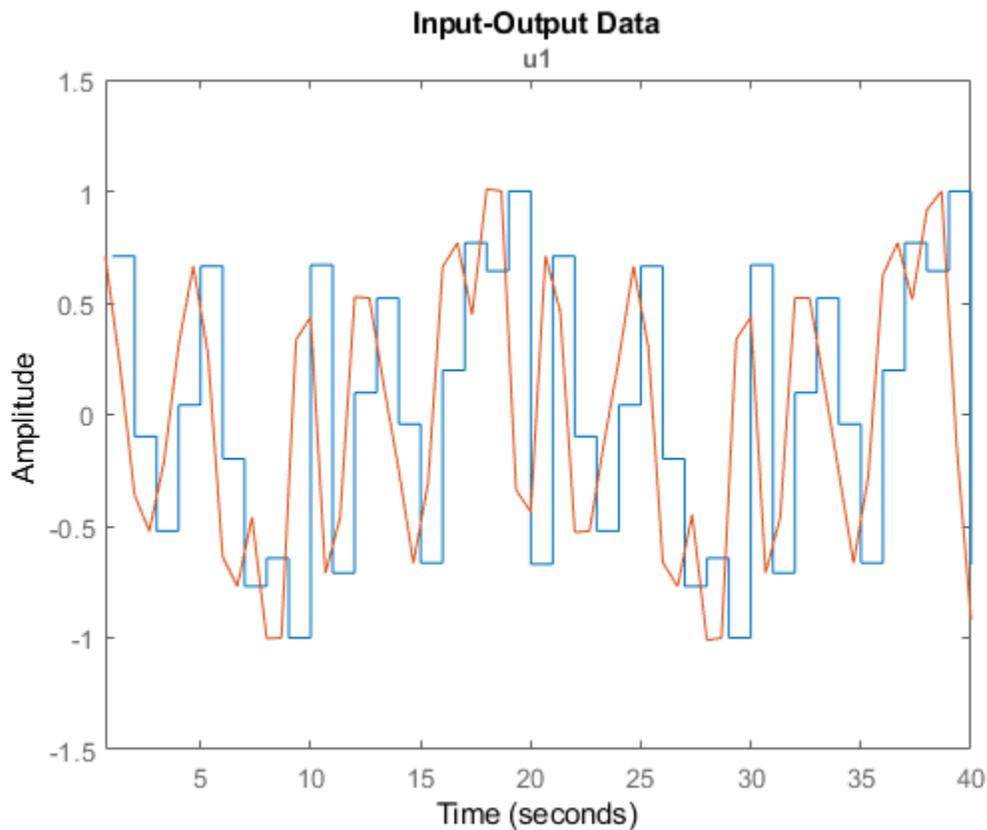
Default: 10

## Examples

### Resample Time-Domain Data

Increase the sampling rate of data by a factor of 1.5 and compare the resampled and the original data signals.

```
u = idinput([20 1 2], 'sine', [], [], [5 10 1]);  
u = iddata([], u, 1);  
plot(u)  
ur = resample(u, 3, 2);  
plot(u, ur)
```



## Algorithms

If you have installed the Signal Processing Toolbox software, `resample` calls the Signal Processing Toolbox `resample` function. The algorithm takes into account the intersample characteristics of the input signal, as described by `data.InterSample`.

## See Also

`idresamp`

**Introduced before R2006a**

## reset

Reset online parameter estimation System object

## Syntax

```
reset(obj)
```

## Description

`reset(obj)` resets the states of a locked online parameter estimation System object, `obj`, to initial values and leaves the object locked. The states of the object are the estimated parameters and parameter covariance. Use `reset` if you are not satisfied with the estimation or if your system changes modes.

## Examples

### Reset Online Estimation System Object

Create a System object for online estimation of an Output-Error model.

```
obj = recursiveOE('InitialB',[0 0.5],'InitialF',[1 0.8],...  
    'InitialParameterCovariance',0.1);
```

Load the estimation data. For this example, use a static data set for illustration.

```
load iddata1 z1;  
output = z1.y;  
input = z1.u;
```

Estimate model parameters online using `step`.

```
for i = 1:numel(input)  
    [B,F,EstimatedOutput] = step(obj,output(i),input(i));  
end
```

View the object properties.

obj

```
obj =  
  recursiveOE with properties:  
  
          B: [0 2.0014]  
          F: [1 -0.7639]  
    InitialB: [0 0.5000]  
    InitialF: [1 0.8000]  
  ParameterCovariance: [2x2 double]  
InitialParameterCovariance: [2x2 double]  
    EstimationMethod: 'ForgettingFactor'  
    ForgettingFactor: 1  
    EnableAdaptation: true  
    InputProcessing: 'Sample-based'  
    DataType: 'double'
```

Reset the System object.

reset(obj)

The estimated parameters, B and F, and parameter covariance, ParameterCovariance are reset to the initial values.

obj

```
obj =  
  recursiveOE with properties:  
  
          B: [0 0.5000]  
          F: [1 0.8000]  
    InitialB: [0 0.5000]  
    InitialF: [1 0.8000]  
  ParameterCovariance: [2x2 double]  
InitialParameterCovariance: [2x2 double]  
    EstimationMethod: 'ForgettingFactor'  
    ForgettingFactor: 1  
    EnableAdaptation: true  
    InputProcessing: 'Sample-based'  
    DataType: 'double'
```

## Input Arguments

### **obj** — System object for online parameter estimation

recursiveAR object | recursiveARMA object | recursiveARX object | recursiveARMAX object | recursiveOE object | recursiveBJ object | recursiveLS object

System object for online parameter estimation, created using one of the following commands:

Online Estimation System Object	Estimated Parameters
recursiveAR	A — Reset to InitialA
recursiveARMA	A — Reset to InitialA C — Reset to InitialC
recursiveARX	A — Reset to InitialA B — Reset to InitialB
recursiveARMAX	A — Reset to InitialA B — Reset to InitialB C — Reset to InitialC
recursiveOE	B — Reset to InitialB F — Reset to InitialF
recursiveBJ	B — Reset to InitialB C — Reset to InitialC D — Reset to InitialD F — Reset to InitialF
recursiveLS	Parameters — Reset to InitialParameters

When EstimationMethod property of obj is 'ForgettingFactor' or 'KalmanFilter', the ParameterCovariance property of obj is reset to the value of InitialParameterCovariance.

## See Also

clone | isLocked | recursiveAR | recursiveARMA | recursiveARMAX | recursiveARX | recursiveBJ | recursiveLS | recursiveOE | release | step

## Topics

“Perform Online Parameter Estimation at the Command Line”

“Validate Online Parameter Estimation at the Command Line”

“What Is Online Estimation?”

**Introduced in R2015b**

## reshape

Change shape of model array

### Syntax

```
sys = reshape(sys,s1,s2,...,sk)
sys = reshape(sys,[s1 s2 ... sk])
```

### Description

`sys = reshape(sys,s1,s2,...,sk)` (or, equivalently, `sys = reshape(sys,[s1 s2 ... sk])`) reshapes the LTI array `sys` into an `s1`-by-`s2`-by-...-by-`sk` model array. With either syntax, there must be `s1*s2*...*sk` models in `sys` to begin with.

### Examples

#### Change Shape of Model Array

Generate a 2-by-3 array of SISO models with four states each.

```
sys = rss(4,1,1,2,3);
size(sys)
```

2x3 array of state-space models.  
Each model has 1 outputs, 1 inputs, and 4 states.

Change the shape of the array to create a 6-by-1 model array.

```
sys1 = reshape(sys,6,1);
size(sys1)
```

6x1 array of state-space models.  
Each model has 1 outputs, 1 inputs, and 4 states.

## See Also

`ndims` | `size`

**Introduced before R2006a**

## resid

Compute and test residuals

### Syntax

```
resid(Data,sys)
resid(Data,sys,Linespec)
resid(Data,sys1,...,sysn)
resid(Data,sys1,Linespec1,...,sysn,Linespecn)
```

```
resid( ____,Options)
```

```
resid( ____,Type)
```

```
[E,R] = resid(Data,sys)
```

### Description

`resid(Data,sys)` computes the 1-step-ahead prediction errors (residuals) for an identified model, `sys`, and plots residual-input dynamics as one of the following, depending on the data in `Data`:

- For time-domain data, `resid` plots the auto-correlation of the residuals and the cross-correlation of the residuals with the input signals. The correlations are generated for lags -25 to 25. To specify a different maximum lag value, use `residOptions`. The 99% confidence region marking statistically insignificant correlations displays as a shaded region around the X-axis.
- For frequency-domain data, `resid` plots a bode plot of the frequency response from the input signals to the residuals. The 99% confidence region marking statistically insignificant response is shown as a region around the X-axis.

To change display options, right-click the plot to access the context menu. For more details about the menu, see “Tips” on page 1-1473.

`resid(Data,sys,Linespec)` sets the line style, marker symbol, and color.

`resid(Data,sys1,...,sysn)` computes and plots the residual of multiple identified models `sys1,...,sysn`.

`resid(Data,sys1,Linespec1,...,sysn,Linespecn)` sets the line style, marker symbol, and color for each system.

`resid( ____,Options)` specifies additional residual calculation options. Use `Options` with any of the previous syntaxes.

`resid( ____,Type)` specifies the plot type. Use `Type` with any of the previous syntaxes.

`[E,R] = resid(Data,sys)` returns the calculated residuals, `E`, and residual correlations, `R`. No plot is generated.

## Examples

### Plot Model Residual Correlations

Load time-domain data.

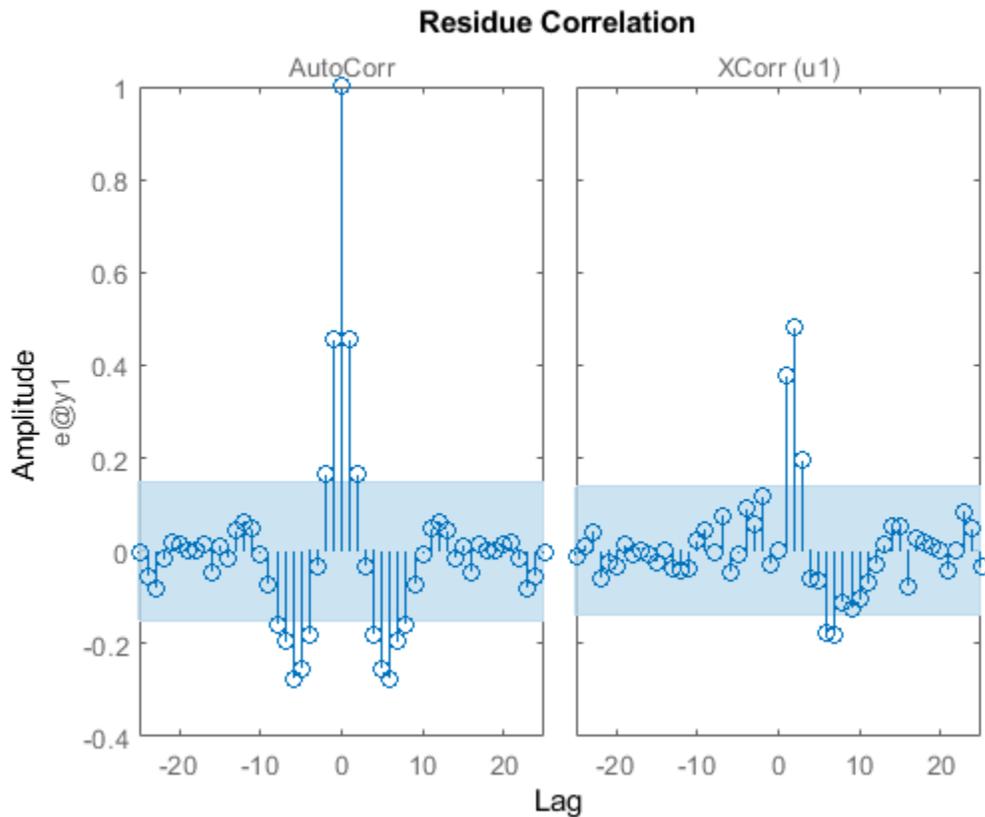
```
load iddata1
data = z1;
```

Estimate an ARX model.

```
sys = arx(data,[1 1 0]);
```

Plot the autocorrelation of the residuals and cross-correlation between the residuals and the inputs.

```
resid(data,sys)
```



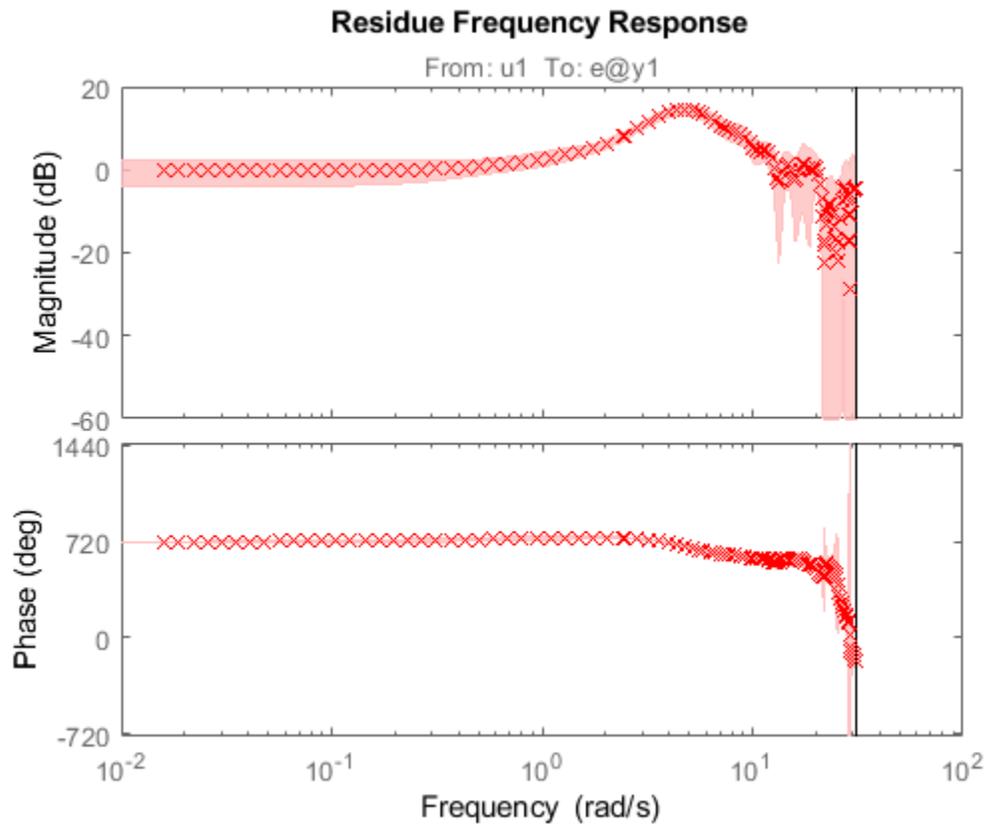
The correlations are calculated until the default maximum lag, 25. The 99% confidence region marking statistically insignificant correlations displays as a shaded region around the X-axis.

Convert data to frequency domain.

```
data2 = fft(data);
```

Compute the residuals for identified model, `sys`, and the frequency-domain data. Plot the residual response using red crosses.

```
resid(data2,sys, 'rx')
```



For frequency-domain data, `resid` plots the Bode plot showing frequency response from the input to the residuals.

### Compare the Residuals for Multiple Identified Models

Load time-domain data.

```
load iddata1
```

Estimate an ARX model.

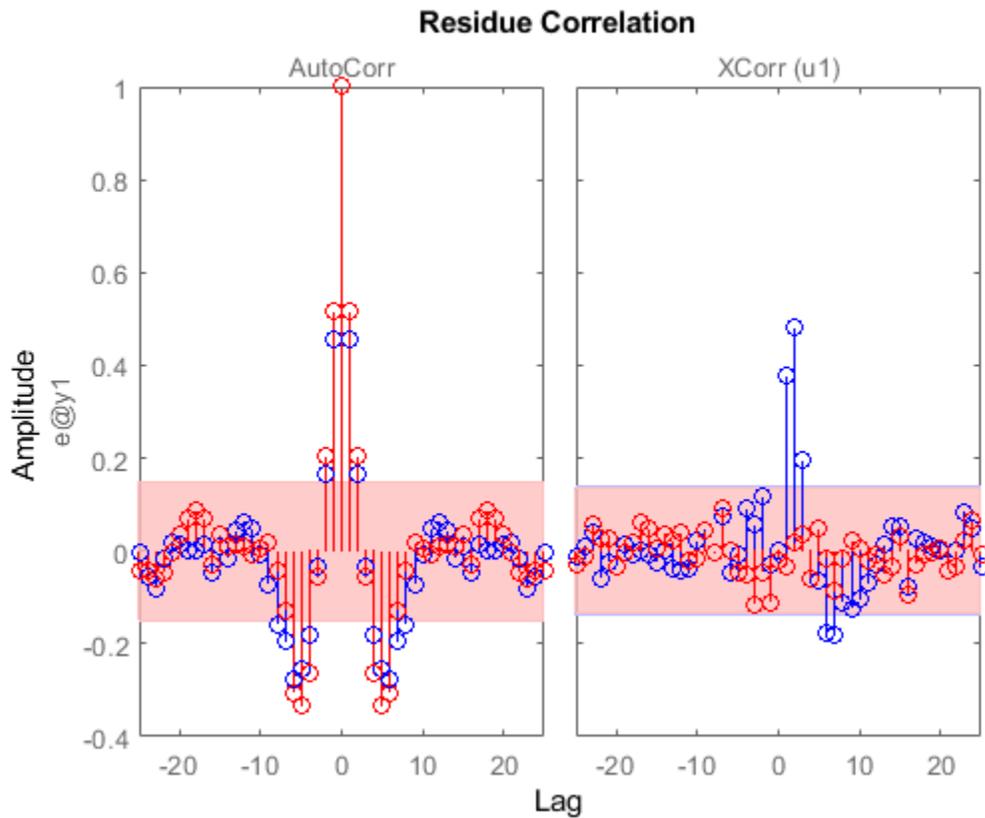
```
sys1 = arx(z1,[1 1 0]);
```

Estimate a transfer function model.

```
sys2 = tfest(z1,2);
```

Plot the correlations of the residuals.

```
resid(z1,sys1,'b',sys2,'r')
```



The cross-correlation between residuals of `sys2` and the inputs lie in the 99% confidence band for all lags.

### Specify Maximum Lag for Residual Impulse Response Calculations

Load time-domain data.

```
load iddata1
```

Estimate an ARX model.

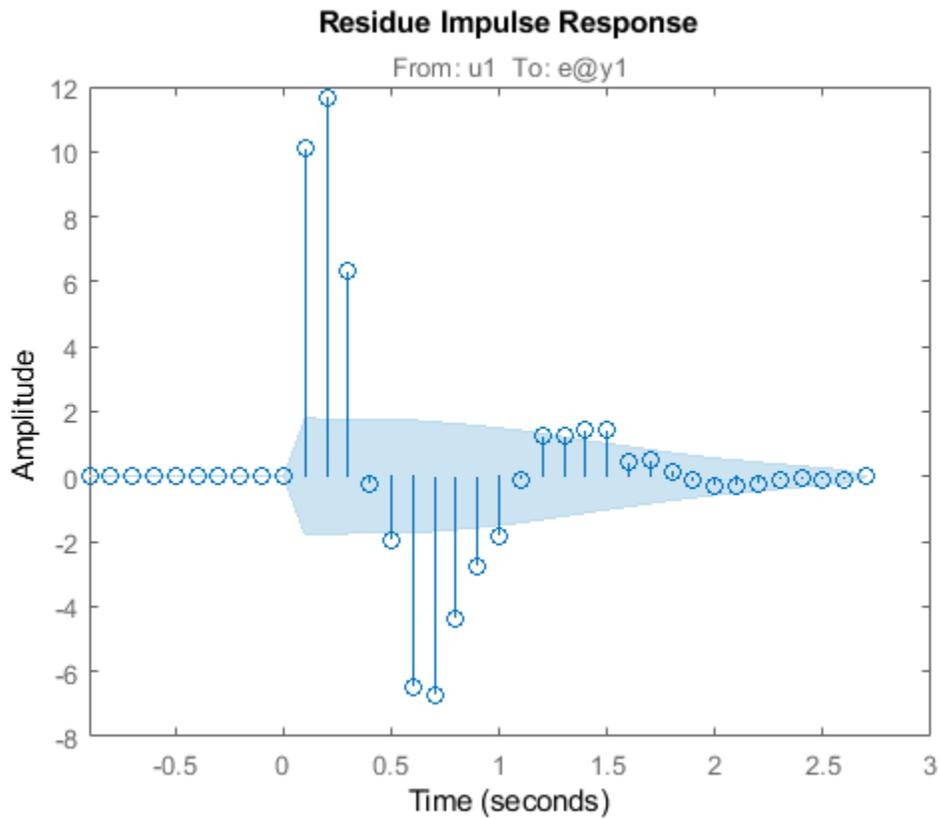
```
sys = arx(z1,[1 1 0]);
```

Specify the maximum lag for residual correlation calculations.

```
opt = residOptions('MaxLag',35);
```

Plot the impulse response from the input to the residuals.

```
resid(z1,sys,opt,'ir')
```



### Calculate Residuals for a MIMO System

Load time-domain data.

```
load iddata7
```

The data is a two-input, single-output dataset.

Estimate an ARX model.

```
sys = tfest(z7,2);
```

Calculate the residuals and their autocorrelations and cross-correlations with inputs.

```
[E,R] = resid(z7,sys);
```

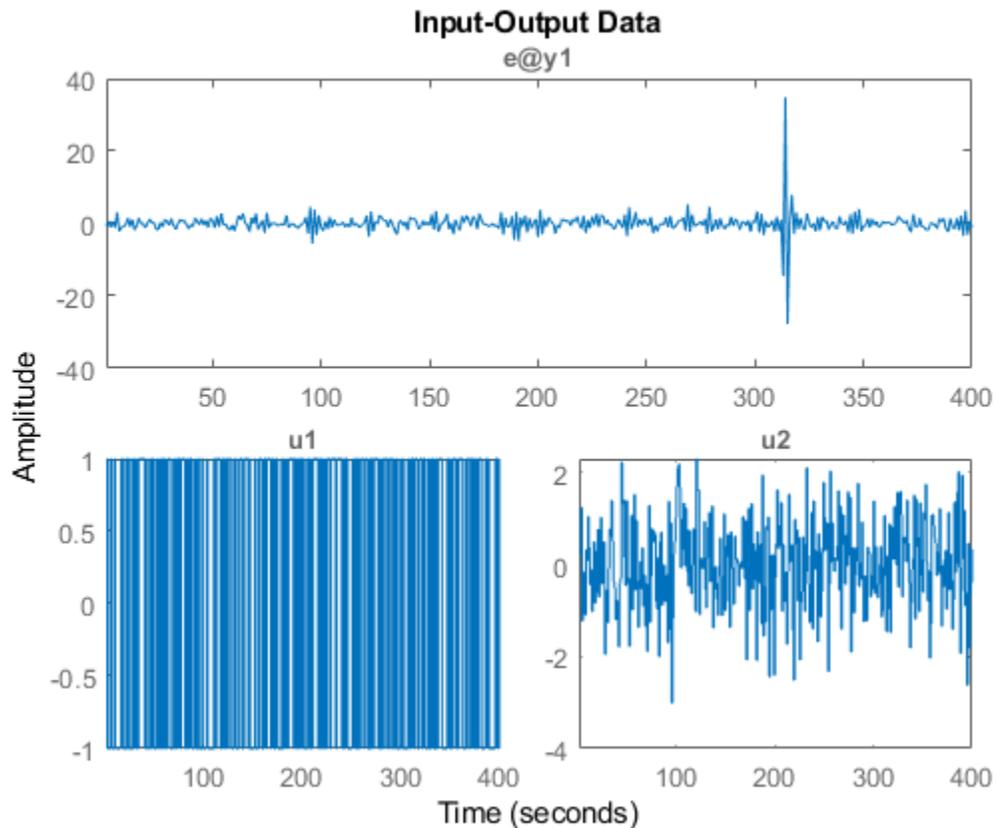
R is a 26-by-3-by-3 matrix of correlations. For example,

- $R(:,1,1)$  is the autocorrelation of the residuals until lag 25.
- $R(:,1,2)$  is the cross-correlation of the residuals with the first input, until lag 25.

E is an iddata object with the residuals as output data and the inputs of validation data (z7) as input data. You can use E to identify error models and analyze the error dynamics.

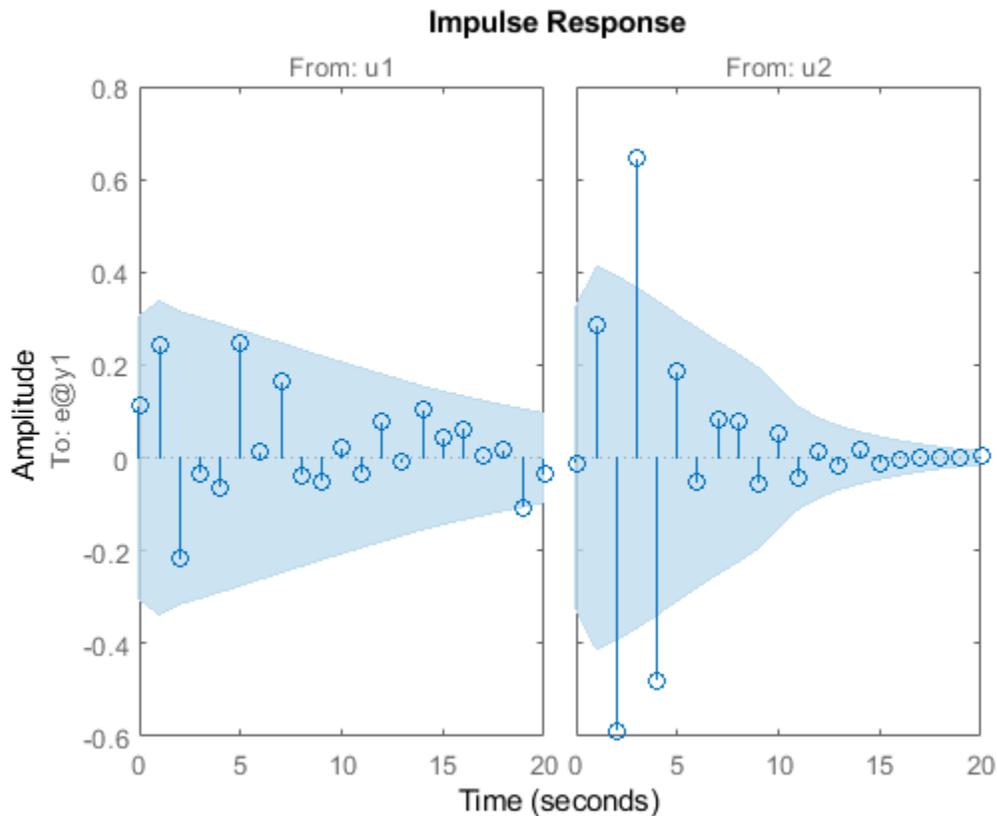
Plot the error.

```
plot(E)
```



Estimate impulse response between inputs and residuals. Plot them with a 3 standard deviation confidence region.

```
I = impulseest(E);  
showConfidence(impulseplot(I,20),3)
```



## Input Arguments

**Data — Validation data**

iddata object

Validation input-output data, specified as an `iddata` object. Data can have multiple input-output channels. When `sys` is linear, Data is time-domain or frequency-domain. When `sys` is nonlinear, Data is time-domain.

### **sys — System used for computing residuals**

identified linear or nonlinear model

System used for computing residuals, specified as an identified linear or nonlinear model.

Example: `idpoly`

### **Linespec — Line style, marker symbol, and color**

character vector

Line style, marker symbol, and color, specified as a character vector. For more information, see `plot` in the MATLAB documentation. When `Type` is specified as `'corr'`, only the line style is used.

Example: `'Linespec', 'kx'`

### **Options — Residual analysis options**

`residOptions` option set

Residual analysis options, specified as an `residOptions` option set.

### **Type — Plot type**

`'corr' | 'ir' | 'fr'`

Plot type, specified as one of the following values:

- `'corr'` — Plots the autocorrelation of the residuals,  $e$ , and the cross-correlation of the residuals with the input signals,  $u$ . The correlations are generated for lags -25 to 25. Use `residOptions` to specify a different maximum lag value. The 99% confidence region marking statistically insignificant correlations is also shown as a shaded region around the X-axis. The computation of the confidence region is done assuming  $e$  to be white and independent of  $u$ .  
  
`'corr'` is default for time-domain data. This plot type is not available for frequency-domain data.
- `'ir'` — Plots the impulse response up to lag 25 of a system from the input to the residuals. The `impulseest` command first estimates the impulse response model with  $e$  as output data and  $u$  as inputs. Then `impulseest` calculates the impulse response

of the estimated model. The 99% confidence region marking statistically insignificant response displays as a shaded region. A low magnitude indicates a reliable model.

This plot type is not available for frequency-domain data.

- 'fr' — The frequency response from the input to the residuals (based on a high-order FIR model) is shown as a Bode plot. The 99% confidence region marking statistically insignificant response displays as a shaded region. A low magnitude in the frequency range of interest indicates a reliable model.

'fr' is default for frequency-domain data.

## Output Arguments

### **E — Model residuals**

`iddata` object

Model residuals, returned as an `iddata` object. The residuals are stored in `E.OutputData`, and the inputs are stored in `E.InputData`. Use `E` to build models that describe the dynamics from the inputs to the residuals. The dynamics are negligible if `sys` is a reliable identified model.

### **R — Correlations of the residuals**

matrix of doubles | []

Correlations of the residuals, returned as one of the following:

- Matrix of doubles — For time-domain-data

$R$  is a matrix of size  $M+1$ -by- $(ny+nu)$ -by- $(ny+nu)$ . Where,  $M$  is the maximum lag specified in `Options`,  $ny$  is the number of outputs, and  $nu$  is the number of inputs. The default value of  $M$  is 25.

At each lag  $k$  ( $k = 0:M$ ),  $R(k, i, j)$  is the expectation value,  $\langle Z(t, i) \cdot Z(t+k-1, j) \rangle$ . Here,  $Z = [E.OutputData, E.InputData]$ .

For example, for a two-output, single-input model,  $Z = [e1, e2, u1]$ . Where,  $e1$  is the residual of the first output,  $e2$  is the residual of the second output, and  $u1$  is the input.  $R$  is a 26-by-3-by-3 matrix, where:

- $R(5, 1, 2) = \langle e1(t) \cdot e2(t+4) \rangle$  is the cross-correlation at lag 4 between  $e1$  and  $e2$ .

- $R(5, 1, 3) = \langle e1(t) \cdot u1(t+4) \rangle$  is the cross-correlation at lag 4 between  $e1$  and  $u1$ .
- $R(5, 1, 1)$ ,  $R(5, 2, 2)$ ,  $R(5, 3, 3)$  are the autocorrelations at lag 4 for  $e1$ ,  $e2$ , and  $u1$ , respectively.
- [ ] — For frequency-domain data

## Tips

- Right-clicking the plot opens the context menu, where you can access the following options:
  - **Systems** — Select systems to view the residual correlation or response plots. By default, all systems are plotted.
  - **Show Confidence Region** — View the 99% confidence region marking statistically insignificant correlations. Applicable only for the correlation plots.
  - **Data Experiment** — For multi-experiment data only. Toggle between data from different experiments.
  - **Characteristics** — View data characteristics. Not applicable for correlation plots.
    - **Peak Response** — View peak response of the data.
    - **Confidence Region** — View the 99% confidence region marking statistically insignificant response.
  - **Show** — Applicable only for frequency-response plots.
    - **Magnitude** — View magnitude of frequency response.
    - **Phase** — View phase of frequency response.
  - **I/O Grouping** — For datasets containing more than one input or output channel. Select grouping of input and output channels on the plot. Not applicable for correlation plots.
    - **None** — Plot input-output channels in their own separate axes.
    - **All** — Group all input channels together and all output channels together.
  - **I/O Selector** — For datasets containing more than one input or output channel. Select a subset of the input and output channels to plot. By default, all output channels are plotted.

- **Grid** — Add grids to the plot.
- **Normalize** — Normalize the y-scale of all data in the plot. Not applicable for frequency-response data.
- **Full View** — Return to full view. By default, the plot is scaled to full view.
- **Initial Condition** — Specify handling of initial conditions.

Specify as one of the following:

- **Estimate** — Treat the initial conditions as estimation parameters.
- **Zero** — Set all initial conditions to zero.
- **Absorb delays and estimate** — Absorb nonzero delays into the model coefficients and treat the initial conditions as estimation parameters. Use this option for discrete-time models only.
- **Properties** — Open the Property Editor dialog box to customize plot attributes.

## References

- [1] Ljung, L. *System Identification: Theory for the User*. Upper Saddle River, NJ: Prentice-Hall PTR, 1999, Section 16.6.

## See Also

`compare` | `predict` | `residOptions` | `sim` | `simsd`

## Topics

“What Is Residual Analysis?”

**Introduced before R2006a**

# residOptions

Option set for resid

## Syntax

```
opt = residOptions  
opt = residOptions(Name,Value)
```

## Description

`opt = residOptions` creates the default option set for `resid`. Use dot notation to customize the option set, if needed.

`opt = residOptions(Name,Value)` creates an option set with options specified by one or more `Name,Value` pair arguments. The options that you do not specify retain their default value.

## Examples

### Create and Modify Default Option Set for Residual Analysis

Create a default option set for `resid`.

```
opt = residOptions;
```

Specify the maximum lag for residual correlation calculations.

```
opt.MaxLag = 35;
```

### Specify Options for Residual Analysis

Create an option set for `resid` that specifies initial condition as zero.

```
opt = residOptions('InitialCondition','z');
```

## Input Arguments

### Name-Value Pair Arguments

Specify optional comma-separated pairs of `Name`, `Value` arguments. `Name` is the argument name and `Value` is the corresponding value. `Name` must appear inside quotes. You can specify several name and value pair arguments in any order as `Name1,Value1,...,NameN,ValueN`.

Example: `residOptions('InitialCondition','e')`

#### **MaxLag** — Maximum positive lag

25 (default) | positive integer

Maximum positive lag for residual correlation and impulse response calculations, specified as the comma-separated pair consisting of `'MaxLag'` and a positive integer.

#### **InitialCondition** — Handling of initial conditions

'e' (default) | 'z' | 'd' | column vector | matrix | structure | idpar object x00bj

Handling of initial conditions, specified as the comma-separated pair consisting of `'InitialCondition'` and one of the following values:

- `'z'` — Zero initial conditions.
- `'e'` — Estimate initial conditions such that the prediction error for observed output is minimized.

For nonlinear grey-box models, only those initial states `i` that are designated as free in the model (`sys.InitialStates(i).Fixed = false`) are estimated. To estimate all the states of the model, first specify all the `Nx` states of the `idnlgrey` model `sys` as free.

```
for i = 1:Nx
    sys.InitialStates(i).Fixed = false;
end
```

Similarly, to fix all the initial states to values specified in `sys.InitialStates`, first specify all the states as fixed in the `sys.InitialStates` property of the nonlinear grey-box model.

- 'd' — Similar to 'e', but absorbs nonzero delays into the model coefficients. The delays are first converted to explicit model states, and the initial values of those states are also estimated and returned.

Use this option for linear models only.

- Vector or Matrix — Initial guess for state values, specified as a numerical column vector of length equal to the number of states. For multi-experiment data, specify a matrix with  $N_e$  columns, where  $N_e$  is the number of experiments. Otherwise, use a column vector to specify the same initial conditions for all experiments. Use this option for state-space (`idss` and `idgrey`) and nonlinear models (`idnlarx`, `idnlhw`, and `idnlgrey`) only.
- Structure with the following fields, which contain the historical input and output values for a time interval immediately before the start time of the data used by `resid`:

Field	Description
Input	Input history, specified as a matrix with $N_u$ columns, where $N_u$ is the number of input channels. For time series models, use <code>[ ]</code> . The number of rows must be greater than or equal to the model order.
Output	Output history, specified as a matrix with $N_y$ columns, where $N_y$ is the number of output channels. The number of rows must be greater than or equal to the model order.

For multi-experiment data, configure the initial conditions separately for each experiment by specifying `InitialCondition` as a structure array with  $N_e$  elements. To specify the same initial conditions for all experiments, use a single structure.

The software uses `data2state` to map the historical data to states. If your model is not `idss`, `idgrey`, `idnlgrey`, or `idnlarx`, the software first converts the model to its state-space representation and then maps the data to states. If conversion of your model to `idss` is not possible, the estimated states are returned empty.

- `x0obj` — Specification object created using `idpar`. Use this object for discrete-time state-space (`idss` and `idgrey`) and nonlinear grey-box (`idnlgrey`) models only. Use `x0obj` to impose constraints on the initial states by fixing their value or specifying minimum or maximum bounds.

## **InputOffset — Removal of offset from time-domain input data during estimation**

[ ] (default) | vector of positive integers | matrix

Removal of offset from time-domain input data during estimation, specified as the comma-separated pair consisting of 'InputOffset' and one of the following:

- A column vector of positive integers of length  $Nu$ , where  $Nu$  is the number of inputs.
- [ ] — Indicates no offset.
- $Nu$ -by- $Ne$  matrix — For multi-experiment data, specify InputOffset as an  $Nu$ -by- $Ne$  matrix.  $Nu$  is the number of inputs, and  $Ne$  is the number of experiments.

Each entry specified by InputOffset is subtracted from the corresponding input data.

## **OutputOffset — Removal of offset from time-domain output data during estimation**

[ ] (default) | vector | matrix

Removal of offset from time-domain output data during estimation, specified as the comma-separated pair consisting of 'OutputOffset' and one of the following:

- A column vector of length  $Ny$ , where  $Ny$  is the number of outputs.
- [ ] — Indicates no offset.
- $Ny$ -by- $Ne$  matrix — For multi-experiment data, specify OutputOffset as a  $Ny$ -by- $Ne$  matrix.  $Ny$  is the number of outputs, and  $Ne$  is the number of experiments.

Each entry specified by OutputOffset is subtracted from the corresponding output data.

## **OutputWeight — Weight of output for initial condition estimation**

[ ] (default) | 'noise' | matrix

Weight of output for initial condition estimation, specified as the comma-separated pair consisting of 'OutputWeight' and one of the following:

- [ ] — No weighting is used. This option is the same as using `eye(Ny)` for the output weight.  $Ny$  is the number of outputs.
- 'noise' — Inverse of the noise variance stored with the model.
- Matrix of doubles — A positive semidefinite matrix of dimension  $Ny$ -by- $Ny$ .  $Ny$  is the number of outputs.

## Output Arguments

**opt** — Option set for resid  
residOptions option set

Option set for resid, returned as an residOptions option set.

## See Also

resid

**Introduced in R2016a**

## retrend

Add offsets or trends to data signals

### Syntax

```
data = retrend(data_d,T)
```

### Description

`data = retrend(data_d,T)` returns a data object `data` by adding the trend information `T` to each signal in `data_d`. `data_d` is a time-domain `iddata` object. `T` is an `TrendInfo` object.

### Examples

#### Retrend Simulated Model Output

Subtract means from input-output signals, estimate a linear model, and retrend the simulated output.

Load SISO data containing vectors `u2` and `y2`.

```
load dryer2
```

Create a data object with sample time of 0.08 seconds.

```
data = iddata(y2,u2,0.08);
```

Remove the mean from the data.

```
[data_d,T] = detrend(data,0);
```

Estimate a linear ARX model.

```
m = arx(data_d,[2 2 1]);
```

Simulate the model output with zero initial states.

```
y_sim = sim(m,data_d(:,[],:));
```

Retrend the simulated model output.

```
y_tot = retrend(y_sim,T);
```

## See Also

[TrendInfo](#) | [detrrend](#) | [getTrend](#)

## Topics

“Handling Offsets and Trends in Data”

**Introduced in R2009a**

## roe

(To be removed) Estimate recursively output-error models (IIR-filters)

---

**Note** roe will be removed in a future release. Use `recursiveOE` instead.

---

## Syntax

```
thm = roe(z,nn,adm,adg)
```

```
[thm,yhat,P,phi,psi] = roe(z,nn,adm,adg,th0,P0,phi0,psi0)
```

## Description

The parameters of the output-error model structure

$$y(t) = \frac{B(q)}{F(q)} u(t - n_k) + e(t)$$

are estimated using a recursive prediction error method.

The input-output data are contained in `z`, which is either an `iddata` object or a matrix `z = [y u]` where `y` and `u` are column vectors. `nn` is given as

```
nn = [nb nf nk]
```

where `nb` and `nf` are the orders of the output-error model, and `nk` is the delay. Specifically,

$$\begin{aligned}nb: B(q) &= b_1 + b_2 q^{-1} + \dots + b_{nb} q^{-nb+1} \\nf: F(q) &= 1 + f_1 q^{-1} + \dots + f_{nf} q^{-nf}\end{aligned}$$

See “What Are Polynomial Models?” for more information.

Only single-input, single-output models are handled by `roe`. Use `rpem` for the multiple-input case.

The estimated parameters are returned in the matrix `thm`. The  $k$ th row of `thm` contains the parameters associated with time  $k$ ; that is, they are based on the data in the rows up to and including row  $k$  in `z`.

Each row of `thm` contains the estimated parameters in the following order.

```
thm(k,:) = [b1, ..., bnb, f1, ..., fnf]
```

`yhat` is the predicted value of the output, according to the current model; that is, row  $k$  of `yhat` contains the predicted value of  $y(k)$  based on all past data.

The actual algorithm is selected with the two arguments `adg` and `adm`. These are described under `rarx`.

The input argument `th0` contains the initial value of the parameters, a row vector consistent with the rows of `thm`. The default value of `th0` is all zeros.

The arguments `P0` and `P` are the initial and final values, respectively, of the scaled covariance matrix of the parameters. The default value of `P0` is  $10^4$  times the unit matrix. The arguments `phi0`, `psi0`, `phi`, and `psi` contain initial and final values of the data vector and the gradient vector, respectively. The sizes of these depend on the chosen model orders. The normal choice of `phi0` and `psi0` is to use the outputs from a previous call to `roe` with the same model orders. (This call could be a dummy call with default input arguments.) The default values of `phi0` and `psi0` are all zeros.

Note that the function requires that the delay  $n_k$  be larger than  $\theta$ . If you want  $n_k = \theta$ , shift the input sequence appropriately and use  $n_k = 1$ .

## Algorithms

The general recursive prediction error algorithm (11.44) of Ljung (1999) is implemented. See also "Recursive Algorithms for Online Parameter Estimation".

## See Also

`nkshift` | `recursiveOE` | `rpem` | `rplr`

**Topics**

“Recursive Algorithms for Online Parameter Estimation”

**Introduced before R2006a**

## rpem

Estimate general input-output models using recursive prediction-error minimization method

rpem is not compatible with MATLAB Coder or MATLAB Compiler™. For the special cases of ARX, AR, ARMA, ARMAX, Box-Jenkins, and Output-Error models, use `recursiveARX`, `recursiveAR`, `recursiveARMA`, `recursiveARMAX`, `recursiveBJ`, and `recursiveOE`, respectively.

## Syntax

```
t hm = rpem(z,nn,adm,adg)
```

```
[t hm,yhat,P,phi,psi] = rpem(z,nn,adm,adg,th0,P0,phi0,psi0)
```

## Description

The parameters of the general linear model structure

$$A(q)y(t) = \frac{B_1(q)}{F_1(q)}u_1(t - nk_1) + \dots + \frac{B_{nu}(q)}{F_{nu}(q)}u_{nu}(t - nk_{nu}) + \frac{C(q)}{D(q)}e(t)$$

are estimated using a recursive prediction error method.

The input-output data is contained in `z`, which is either an `iddata` object or a matrix `z = [y u]` where `y` and `u` are column vectors. (In the multiple-input case, `u` contains one column for each input.) `nn` is given as

```
nn = [na nb nc nd nf nk]
```

where `na`, `nb`, `nc`, `nd`, and `nf` are the orders of the model, and `nk` is the delay. For multiple-input systems, `nb`, `nf`, and `nk` are row vectors giving the orders and delays of each input. See “What Are Polynomial Models?” for an exact definition of the orders.

The estimated parameters are returned in the matrix `t hm`. The `k`th row of `t hm` contains the parameters associated with time `k`; that is, they are based on the data in the rows up

to and including row  $k$  in  $z$ . Each row of  $thm$  contains the estimated parameters in the following order.

```
thm(k,:) = [a1,a2,...,ana,b1,...,bnb,...  
           c1,...,cnc,d1,...,dnd,f1,...,fnf]
```

For multiple-input systems, the  $B$  part in the above expression is repeated for each input before the  $C$  part begins, and the  $F$  part is also repeated for each input. This is the same ordering as in `m.par`.

$yhat$  is the predicted value of the output, according to the current model; that is, row  $k$  of  $yhat$  contains the predicted value of  $y(k)$  based on all past data.

The actual algorithm is selected with the two arguments `adg` and `adm`:

- `adm = 'ff'` and `adg = lam` specify the *forgetting factor* algorithm with the forgetting factor  $\lambda = lam$ . This algorithm is also known as recursive least squares (RLS). In this case, the matrix  $P$  has the following interpretation:  $R_2/2 * P$  is approximately equal to the covariance matrix of the estimated parameters.  $R_2$  is the variance of the innovations (the true prediction errors  $e(t)$ ).

`adm = 'ug'` and `adg = gam` specify the *unnormalized gradient* algorithm with gain  $gamma = gam$ . This algorithm is also known as the normalized least mean squares (LMS).

`adm = 'ng'` and `adg = gam` specify the *normalized gradient* or normalized least mean squares (NLMS) algorithm. In these cases,  $P$  is not applicable.

`adm = 'kf'` and `adg = R1` specify the *Kalman filter based* algorithm with  $R_2 = 1$  and  $R_1 = R1$ . If the variance of the innovations  $e(t)$  is not unity but  $R_2$ ; then  $R_2 * P$  is the covariance matrix of the parameter estimates, while  $R_1 = R1 / R_2$  is the covariance matrix of the parameter changes.

The input argument `th0` contains the initial value of the parameters, a row vector consistent with the rows of `thm`. The default value of `th0` is all zeros.

The arguments `P0` and `P` are the initial and final values, respectively, of the scaled covariance matrix of the parameters. The default value of `P0` is  $10^4$  times the unit matrix. The arguments `phi0`, `psi0`, `phi`, and `psi` contain initial and final values of the data vector and the gradient vector, respectively. The sizes of these depend on the chosen model orders. The normal choice of `phi0` and `psi0` is to use the outputs from a previous call to `rpem` with the same model orders. (This call could be a dummy call with default input arguments.) The default values of `phi0` and `psi0` are all zeros.

Note that the function requires that the delay  $n_k$  be larger than 0. If you want  $n_k = 0$ , shift the input sequence appropriately and use  $n_k = 1$ .

## Examples

### Estimate Model Parameters Using Recursive Prediction-Error Minimization

Specify the order and delays of a polynomial model structure.

```
na = 2;
nb = 1;
nc = 1;
nd = 1;
nf = 0;
nk = 1;
```

Load the estimation data.

```
load iddata1 z1
```

Estimate the parameters using forgetting factor algorithm with forgetting factor 0.99.

```
EstimatedParameters = rpem(z1,[na nb nc nd nf nk],'ff',0.99);
```

Get the last set of estimated parameters.

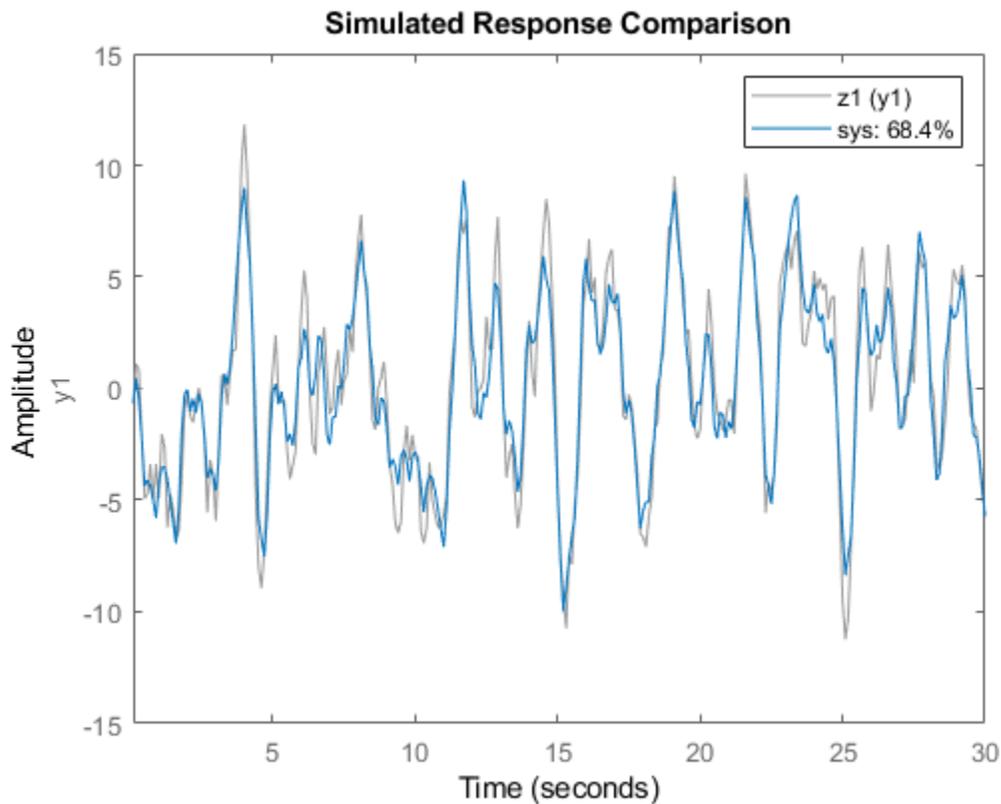
```
p = EstimatedParameters(end,:);
```

Construct a polynomial model with the estimated parameters.

```
sys = idpoly([1 p(1:na)],... % A polynomial
             [zeros(1,nk) p(na+1:na+nb)],... % B polynomial
             [1 p(na+nb+1:na+nb+nc)],... % C polynomial
             [1 p(na+nb+nc+1:na+nb+nc+nd)]); % D polynomial
sys.Ts = z1.Ts;
```

Compare the estimated output with measured data.

```
compare(z1,sys);
```



## Algorithms

The general recursive prediction error algorithm (11.44) of Ljung (1999) is implemented. See also "Recursive Algorithms for Online Parameter Estimation".

## See Also

[nkshift](#) | [recursiveAR](#) | [recursiveARMA](#) | [recursiveARMAX](#) | [recursiveARX](#) | [recursiveBJ](#) | [recursiveOE](#) | [rplr](#)

## **Topics**

“What Is Online Estimation?”

“Recursive Algorithms for Online Parameter Estimation”

**Introduced before R2006a**

## rplr

Estimate general input-output models using recursive pseudolinear regression method  
rplr is not compatible with MATLAB Coder or MATLAB Compiler.

## Syntax

```
thm = rplr(z,nn,adm,adg)
```

```
[thm,yhat,P,phi] = rplr(z,nn,adm,adg,th0,P0,phi0)
```

## Description

This is a direct alternative to `rpem` and has essentially the same syntax. See `rpem` for an explanation of the input and output arguments.

`rplr` differs from `rpem` in that it uses another gradient approximation. See Section 11.5 in Ljung (1999). Several of the special cases are well-known algorithms.

When applied to ARMAX models, ( $nn = [na \ nb \ nc \ 0 \ 0 \ nk]$ ), `rplr` gives the extended least squares (ELS) method. When applied to the output-error structure ( $nn = [0 \ nb \ 0 \ 0 \ nf \ nk]$ ), the method is known as HARF in the `adm = 'ff'` case and SHARF in the `adm = 'ng'` case.

`rplr` can also be applied to the time-series case in which an ARMA model is estimated with:

```
z = y  
nn = [na nc]
```

## Examples

### Estimate Output-Error Model Parameters Using Recursive Pseudolinear Regression

Specify the order and delays of an Output-Error model structure.

```
na = 0;  
nb = 2;  
nc = 0;  
nd = 0;  
nf = 2;  
nk = 1;
```

Load the estimation data.

```
load iddata1 z1
```

Estimate the parameters using forgetting factor algorithm, with forgetting factor 0.99.

```
EstimatedParameters = rplr(z1,[na nb nc nd nf nk], 'ff',0.99);
```

## References

For more information about HARF and SHARF, see Chapter 11 in Ljung (1999).

## See Also

nkshift | recursiveAR | recursiveARMA | recursiveARMAX | recursiveARX | recursiveBJ | recursiveOE | rpem

## Topics

“What Is Online Estimation?”

“Recursive Algorithms for Online Parameter Estimation”

**Introduced before R2006a**

## rsample

Random sampling of linear identified systems

### Syntax

```
sys_array = rsample(sys,N)  
sys_array = rsample(sys,N,sd)
```

### Description

`sys_array = rsample(sys,N)` creates `N` random samples of the identified linear system, `sys`. `sys_array` contains systems with the same structure as `sys`, whose parameters are perturbed about their nominal values, based on the parameter covariance.

`sys_array = rsample(sys,N,sd)` specifies the standard deviation level, `sd`, for perturbing the parameters of `sys`.

### Input Arguments

#### **sys**

Identifiable system.

#### **N**

Number of samples to be generated.

**Default:** 10

#### **sd**

Standard deviation level for perturbing the identifiable parameters of `sys`.

**Default:** 1

## Output Arguments

### **sys\_array**

Array of random samples of `sys`.

If `sys` is an array of models, then the size of `sys_array` is equal to `[size(sys) N]`. There are `N` randomized samples for each model in `sys`.

The parameters of the samples in `sys_array` vary from the original identifiable model within 1 standard deviation of their nominal values.

## Examples

### **Create Random Samples of Estimated Model**

Estimate a third-order, discrete-time, state-space model.

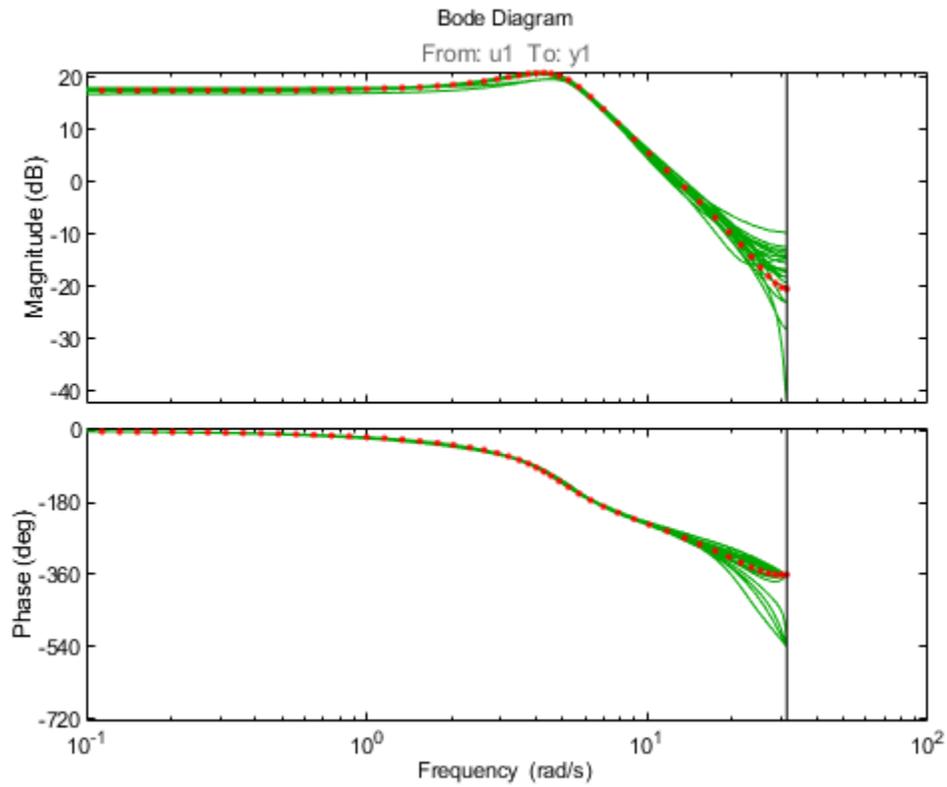
```
load iddata2 z2;  
sys = n4sid(z2,3);
```

Randomly sample the estimated model.

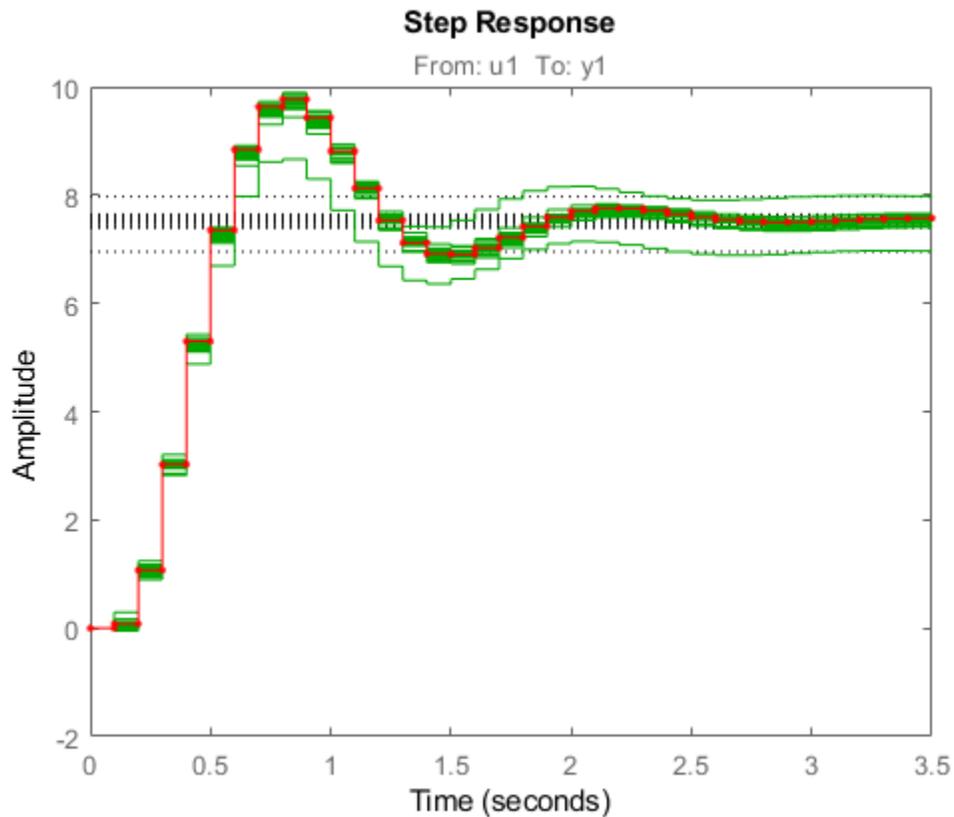
```
N = 20;  
sys_array = rsample(sys,N);
```

Analyze the uncertainty in time (step) and frequency (Bode) responses.

```
opt = bodeoptions;  
opt.PhaseMatching = 'on';  
figure;  
bodeplot(sys_array,'g',sys,'r.',opt)
```



```
figure;  
stepplot(sys_array, 'g', sys, 'r.-')
```



### Specify Standard Deviation Level for Parameter Perturbation

Estimate a third-order, discrete-time, state-space model.

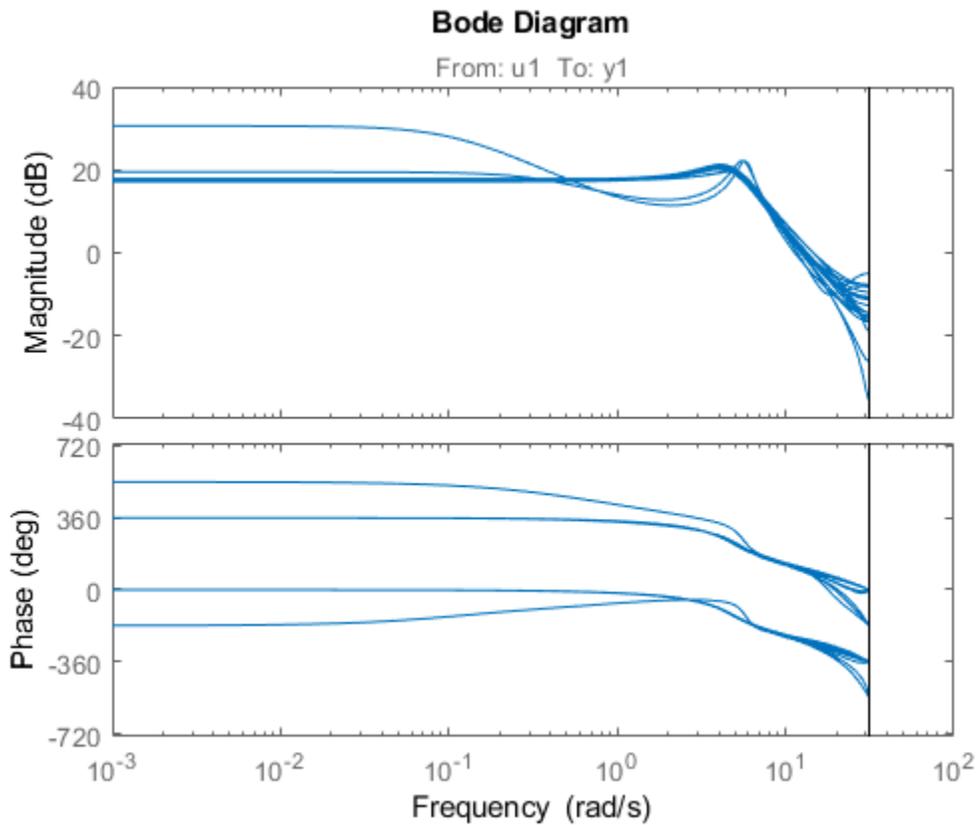
```
load iddata2 z2;  
sys = n4sid(z2,3);
```

Randomly sample the estimated model. Specify the standard deviation level for perturbing the model parameters.

```
N = 20;  
sd = 2;  
sys_array = rsample(sys,N,sd);
```

Analyze the model uncertainty.

```
figure;  
bode(sys_array);
```



### Compare Frequency Response Confidence Regions for Sampled ARMAX Model

Estimate an ARMAX model.

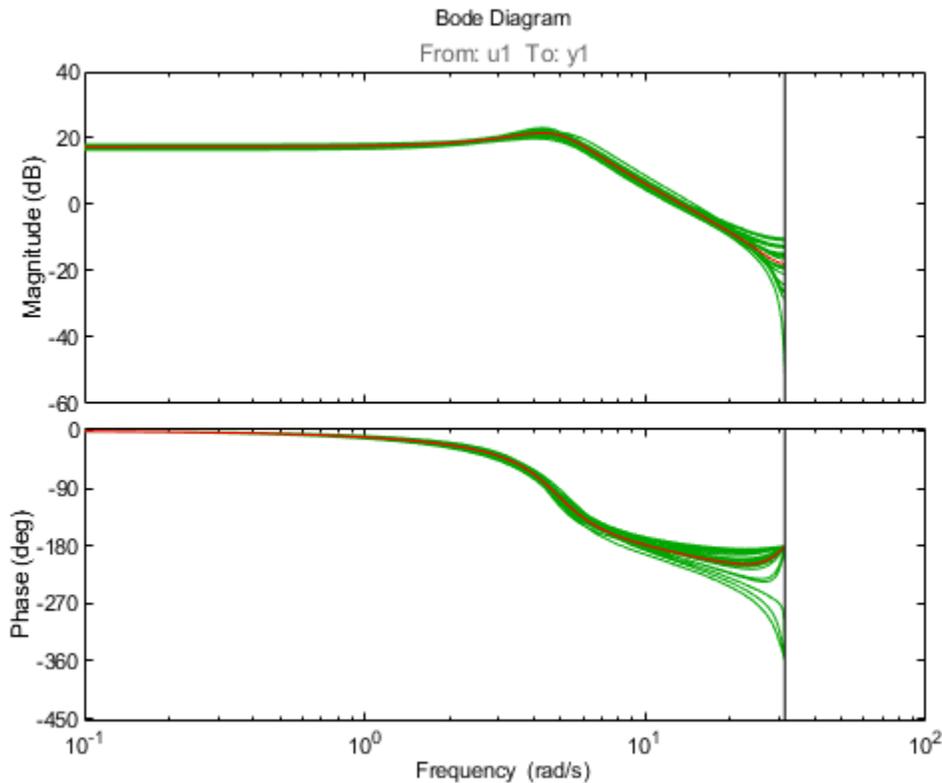
```
load iddata1 z1
sys = armax(z1,[2 2 2 1]);
```

Randomly sample the ARMAX model. Perturb the model parameters up to 2 standard deviations.

```
N = 20;
sd = 2;
sys_array = rsample(sys,N,sd);
```

Compare the frequency response confidence region corresponding to 2 standard deviations (asymptotic estimate) with the model array response.

```
opt = bodeoptions; opt.PhaseMatching = 'on';
opt.ConfidenceRegionNumberSD = 2;
bodeplot(sys_array,'g',sys,'r',opt)
```



To view the confidence region, right click the plot, and choose **Characteristics > Confidence Region**.

## Tips

- For systems with large parameter uncertainties, the randomized systems may contain unstable elements. These unstable elements may make it difficult to analyze the properties of the identified system. Execution of analysis commands, such as `step`, `bode`, `sim`, etc., on such systems can produce unreliable results. Instead, use a dedicated Monte-Carlo analysis command, such as `simsd`.

## See Also

`bode` | `init` | `iopzmap` | `noise2meas` | `noisecnv` | `simsd` | `step` | `translatecov`

**Introduced in R2012a**

## saturation

Create a saturation nonlinearity estimator object

### Syntax

```
NL = saturation  
NL = saturation('LinearInterval', [a,b])
```

### Description

`NL = saturation` creates a default saturation nonlinearity estimator object for estimating Hammerstein-Wiener models. The linear interval is set to `[NaN NaN]`. The initial value of the linear interval is determined from the estimation data range during estimation using `nllhw`. Use dot notation to customize the object properties, if needed.

`NL = saturation('LinearInterval', [a,b])` creates a saturation nonlinearity estimator object initialized with linear interval, `[a,b]`.

Alternatively, use `NL = saturation([a,b])`.

### Object Description

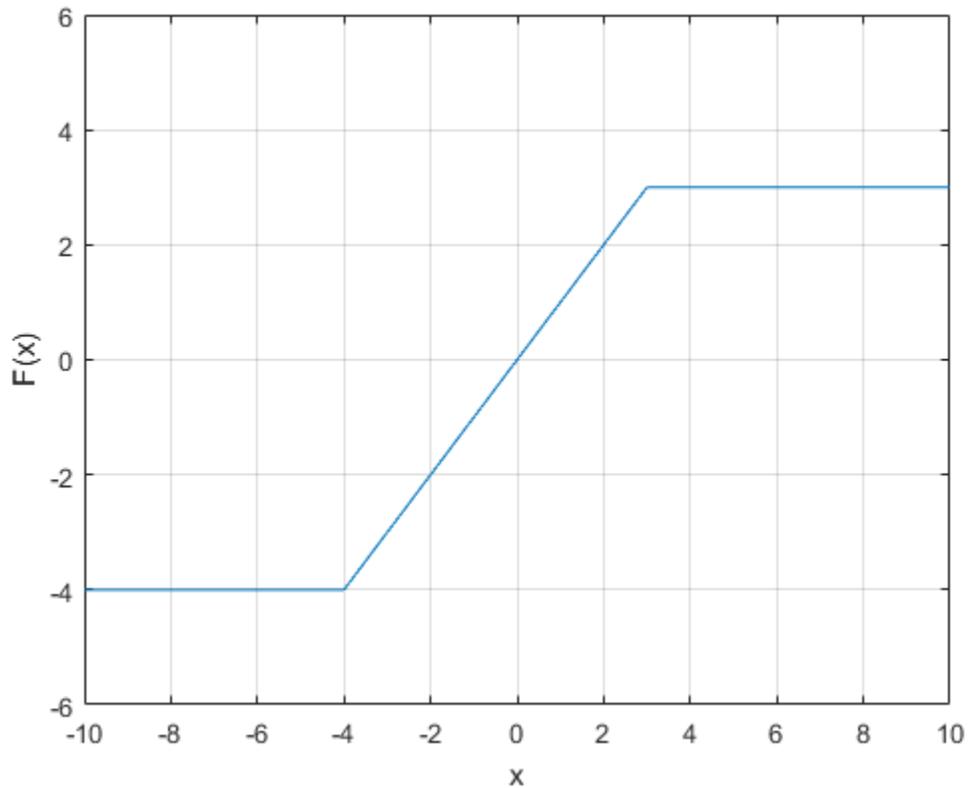
`saturation` is an object that stores the saturation nonlinearity estimator for estimating Hammerstein-Wiener models.

Use `saturation` to define a nonlinear function  $y = F(x, \theta)$ , where  $y$  and  $x$  are scalars, and  $\theta$  represents the parameters  $a$  and  $b$  that define the linear interval, `[a,b]`.

The saturation nonlinearity function has the following characteristics:

$$\begin{array}{ll} a \leq x < b & F(x) = x \\ a > x & F(x) = a \\ b \leq x & F(x) = b \end{array}$$

For example, in the following plot, the linear interval is  $[-4, 3]$ .



The value  $F(x)$  is computed by `evaluate(NL, x)`, where `NL` is the saturation object.

For saturation object properties, see “Properties” on page 1-1505.

## Examples

### Create a Default Saturation Nonlinearity Estimator

```
NL = saturation;
```

Specify the linear interval.

```
NL.LinearInterval = [-4,5];
```

## Estimate a Hammerstein Model with Saturation

Load data.

```
load twotankdata;  
z = iddata(y,u,0.2,'Name','Two tank system');  
z1 = z(1:1000);
```

Create a saturation object with lower limit of 0 and upper limit of 5.

```
InputNL = saturation('LinearInterval',[0 5]);
```

Estimate model with no output nonlinearity.

```
m = nlhw(z1,[2 3 0],InputNL,[]);
```

## Estimate MIMO Hammerstein-Wiener Model

Load the estimation data.

```
load motorizedcamera;
```

Create an iddata object.

```
z = iddata(y,u,0.02,'Name','Motorized Camera','TimeUnit','s');
```

z is an iddata object with 6 inputs and 2 outputs.

Specify the model orders and delays.

```
Orders = [ones(2,6),ones(2,6),ones(2,6)];
```

Specify the same nonlinearity estimator for each input channel.

```
InputNL = saturation;
```

Specify different nonlinearity estimators for each output channel.

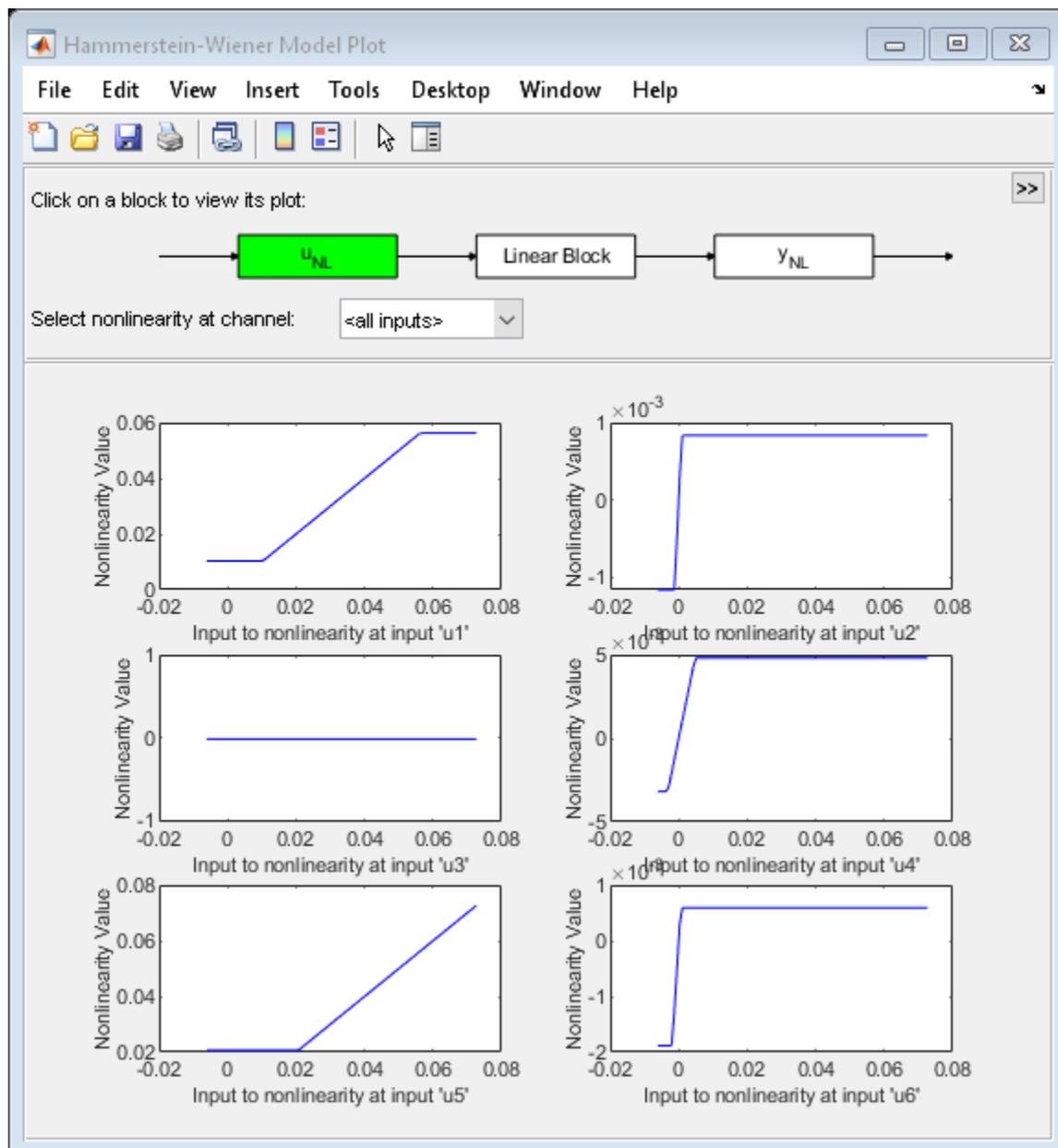
```
OutputNL = [deadzone,wavenet];
```

Estimate the Hammerstein-Wiener model.

```
sys = nlhw(z,Orders,InputNL,OutputNL);
```

To see the shape of the estimated input and output nonlinearities, plot the nonlinearities.

```
plot(sys)
```



Click on the input and output nonlinearity blocks on the top of the plot to see the nonlinearities.

## Input Arguments

### **[a, b] — Linear interval**

[NaN NaN] (default) | 2-element row vector

Linear interval of the saturation, specified as a 2-element row vector of doubles.

The saturation nonlinearity is initialized at the interval [a, b]. The interval values are adjusted to the estimation data by `n_lhw`. To remove the lower limit, set a to `-Inf`. The lower limit is not adjusted during estimation. To remove the upper limit, set b to `Inf`. The upper limit is not adjusted during estimation.

When the interval is [NaN NaN], the initial value of the linear interval is determined from the estimation data range during estimation using `n_lhw`.

Example: [-2 1]

## Properties

### **LinearInterval**

Linear interval of the saturation, specified as a 2-element row vector of doubles.

**Default:** [NaN NaN]

## Output Arguments

### **NL — Saturation nonlinearity estimator object**

saturation object

Saturation nonlinearity estimator object, returned as a saturation object.

## **See Also**

n|hw

**Introduced in R2007a**

# segment

Segment data and estimate models for each segment

segment is not compatible with MATLAB Coder or MATLAB Compiler.

## Syntax

```
segm = segment(z,nn)
```

```
[segm,V,thm,R2e] = segment(z,nn,R2,q,R1,M,th0,P0,ll,mu)
```

## Description

segment builds models of AR, ARX, or ARMAX/ARMA type,

$$A(q)y(t) = B(q)u(t - nk) + C(q)e(t)$$

assuming that the model parameters are piecewise constant over time. It results in a model that has split the data record into segments over which the model remains constant. The function models signals and systems that might undergo abrupt changes.

The input-output data is contained in `z`, which is either an `iddata` object or a matrix  $z = [y \ u]$  where `y` and `u` are column vectors. If the system has several inputs, `u` has the corresponding number of columns.

The argument `nn` defines the model order. For the ARMAX model

```
nn = [na nb nc nk];
```

where `na`, `nb`, and `nc` are the orders of the corresponding polynomials. See “What Are Polynomial Models?”. Moreover, `nk` is the delay. If the model has several inputs, `nb` and `nk` are row vectors, giving the orders and delays for each input.

For an ARX model (`nc = 0`) enter

```
nn = [na nb nk];
```

For an ARMA model of a time series

```
z = y;  
nn = [na nc];
```

and for an AR model

```
nn = na;
```

The output argument `segm` is a matrix, where the  $k$ th row contains the parameters corresponding to time  $k$ . This is analogous to output estimates returned by the `recursiveARX` and `recursiveARMAX` estimators. The output argument `thm` of `segment` contains the corresponding model parameters that have not yet been segmented. Each row of `thm` contains the parameter estimates at the corresponding time instant. These estimates are formed by weighting together the parameters of  $M$  (default: 5) different time-varying models, with the participating models changing at every time step. Consider `segment` as an alternative to the online estimation commands when you are not interested in continuously tracking the changes in parameters of a single model, but need to detect abrupt changes in the system dynamics.

The output argument  $V$  contains the sum of the squared prediction errors of the segmented model. It is a measure of how successful the segmentation has been.

The input argument  $R2$  is the assumed variance of the innovations  $e(t)$  in the model. The default value of  $R2$ ,  $R2 = []$ , is that it is estimated. Then the output argument  $R2e$  is a vector whose  $k$ th element contains the estimate of  $R2$  at time  $k$ .

The argument  $q$  is the probability that the model exhibits an abrupt change at any given time. The default value is  $0.01$ .

$R1$  is the assumed covariance matrix of the parameter jumps when they occur. The default value is the identity matrix with dimension equal to the number of estimated parameters.

$M$  is the number of parallel models used in the algorithm (see below). Its default value is 5.

$th0$  is the initial value of the parameters. Its default is zero.  $P0$  is the initial covariance matrix of the parameters. The default is 10 times the identity matrix.

$ll$  is the guaranteed life of each of the models. That is, any created candidate model is not abolished until after at least  $ll$  time steps. The default is  $ll = 1$ .  $\mu$  is a forgetting parameter that is used in the scheme that estimates  $R2$ . The default is  $0.97$ .

The most critical parameter for you to choose is  $R2$ . It is usually more robust to have a reasonable guess of  $R2$  than to estimate it. Typically, you need to try different values of  $R2$  and evaluate the results. (See the example below.)  $\text{sqrt}(R2)$  corresponds to a change in the value  $y(t)$  that is normal, giving no indication that the system or the input might have changed.

## Examples

### Divide Sinusoid into Segments with Constant Levels

Create a sinusoid for the simulated model output.

```
y = sin([1:50]/3)';
```

Specify the input signal to be constant at 1.

```
u = ones(size(y));
```

Specify the estimated noise variance for the model.

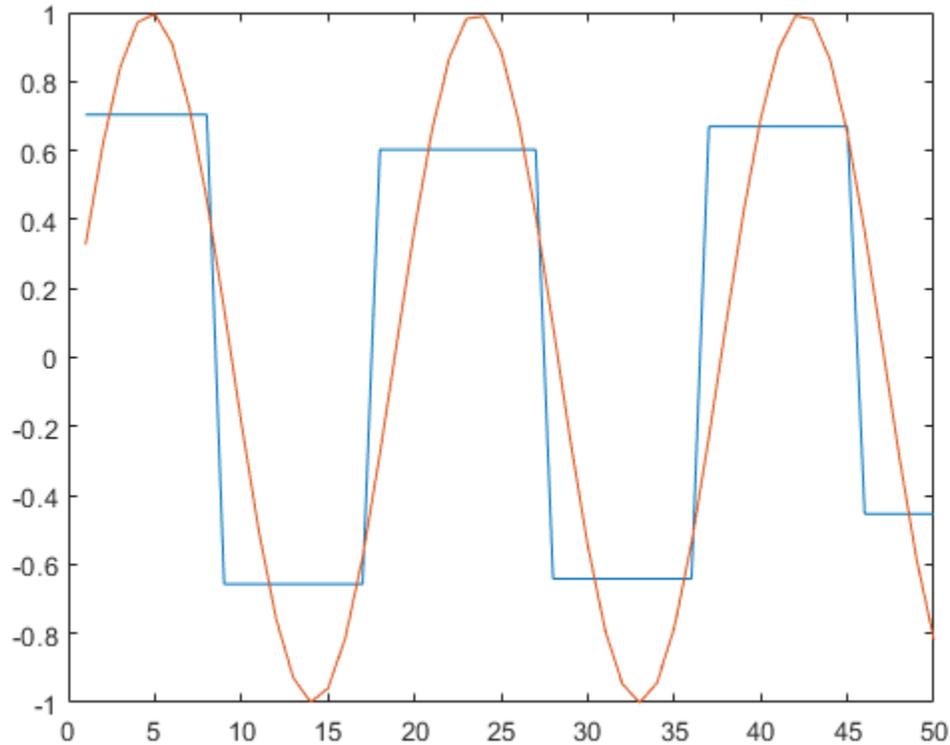
```
R2 = 0.1;
```

Segment the signal and estimate an ARX model for each segment. Use the simple model  $y(t) = b_1 u(t-1)$ , where  $b_1$  is the model parameter describing the piecewise constant level of the estimated output,  $y(t)$ .

```
segm = segment([y,u],[0 1 1],R2);
```

Examine the result.

```
plot([segm,y])
```

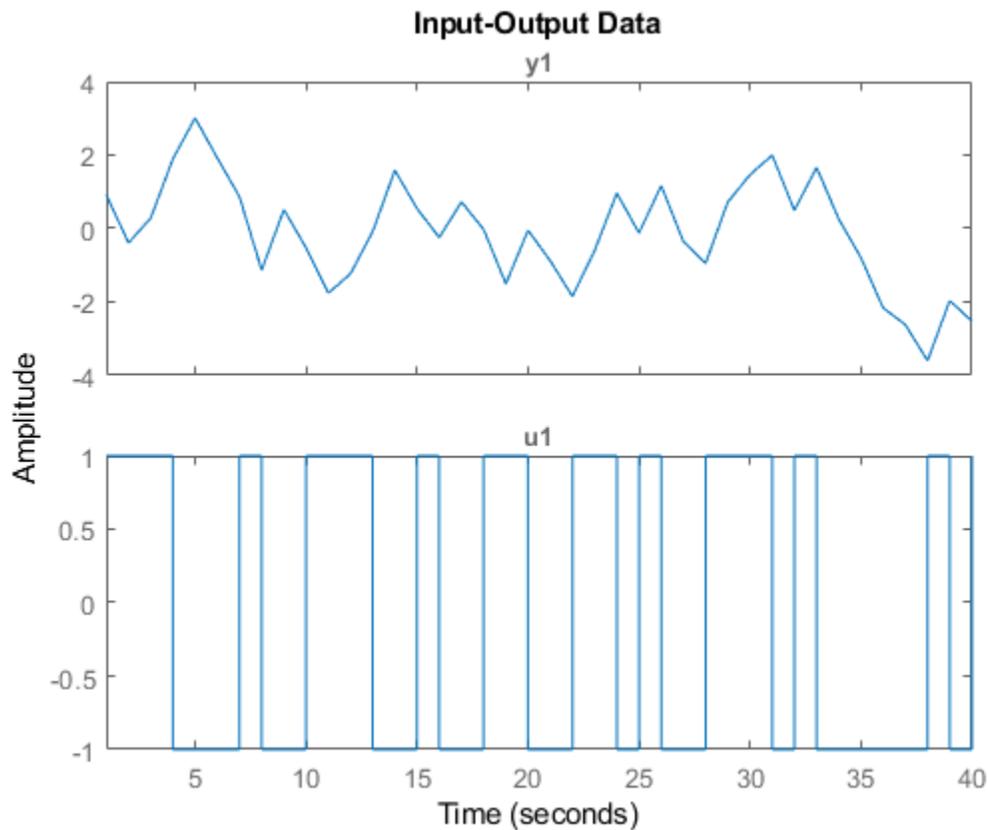


Vary the value of R2 to change the estimated noise variance. Decreasing R2 increases the number of segments produced for this model.

## Model Abrupt Change In Time Delay Using Segmentation

Load and plot the estimation data.

```
load iddemo6m.mat z
z = iddata(z(:,1),z(:,2));
plot(z)
```



This data contains a change in time delay from 2 to 1, which is difficult to detect by examining the data.

Specify the model orders to estimate an ARX model of the form:

$$y(t) + ay(t-1) = b_1u(t-1) + b_2u(t-2)$$

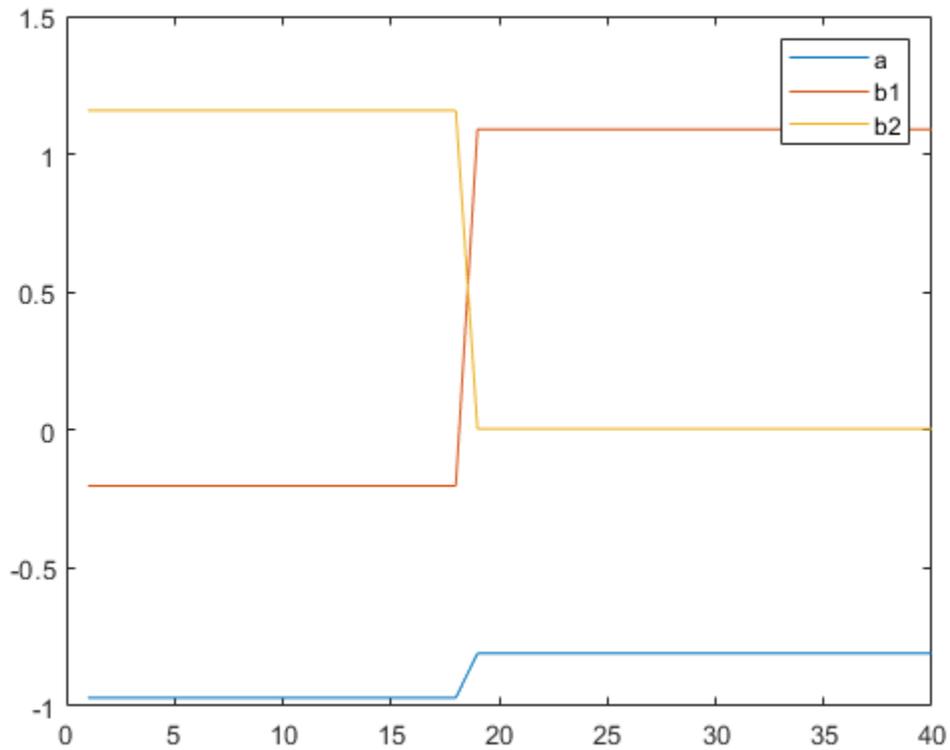
`nn = [1 2 1];`

Segment the data and estimate ARX models for each segment. Specify an estimated noise variance of 0.1.

`seg = segment(z,nn,0.1);`

Examine the parameters of the segmented model.

```
plot(seg)  
legend('a', 'b1', 'b2');
```



The data has been divided into two segments, as indicated by the change in model parameters around sample number 19. The increase in b1, along with a corresponding decrease in b2, shows the change in model delay.

## Algorithms

The algorithm is based on  $M$  parallel models, each recursively estimated by an algorithm of Kalman filter type. Each model is updated independently, and its posterior probability is computed. The time-varying estimate  $\hat{t}_{hm}$  is formed by weighting together the  $M$  different models with weights equal to their posterior probability. At each time step the model (among those that have lived at least  $ll$  samples) that has the lowest posterior probability is abolished. A new model is started, assuming that the system parameters have changed, with probability  $q$ , a random jump from the most likely among the models. The covariance matrix of the parameter change is set to  $R1$ .

After all the data are examined, the surviving model with the highest posterior probability is tracked back and the time instances where it jumped are marked. This defines the different segments of the data. (If no models had been abolished in the algorithm, this would have been the maximum likelihood estimates of the jump instances.) The segmented model  $segm$  is then formed by smoothing the parameter estimate, assuming that the jump instances are correct. In other words, the last estimate over a segment is chosen to represent the whole segment.

## See Also

### Topics

“Data Segmentation”

**Introduced before R2006a**

## selstruc

Select model order for single-output ARX models

### Syntax

```
nn = selstruc(v)
```

```
[nn,vmod] = selstruc(v,c)
```

### Description

---

**Note** Use `selstruc` for single-output systems only. `selstruc` supports both single-input and multiple-input systems.

---

`selstruc` is a function to help choose a model structure (order) from the information contained in the matrix `v` obtained as the output from `arxstruc` or `ivstruc`.

The default value of `c` is `'plot'`. The plot shows the percentage of the output variance that is not explained by the model as a function of the number of parameters used. Each value shows the best fit for that number of parameters. By clicking in the plot you can examine which orders are of interest. When you click **Select**, the variable `nn` is exported to the workspace as the optimal model structure for your choice of number of parameters. Several choices can be made.

`c = 'aic'` gives no plots, but returns in `nn` the structure that minimizes

$$\begin{aligned} V_{\text{mod}} &= \log\left(V\left(1 + \frac{2d}{N}\right)\right) \\ &= \log(V) + \frac{2d}{N}, N \gg d \end{aligned}$$

where  $V$  is the loss function,  $d$  is the total number of parameters in the structure in

question, and  $N$  is the number of data points used for the estimation.  $\log(V) + \frac{2d}{N}$  is the Akaike's Information Criterion (AIC). See `aic` for more details.

`c = 'mdl'` returns in `nn` the structure that minimizes Rissanen's Minimum Description Length (MDL) criterion.

$$V_{\text{mod}} = V\left(1 + \frac{d \log(N)}{N}\right)$$

When `c` equals a numerical value, the structure that minimizes  $V_{\text{mod}} = V\left(1 + \frac{cd}{N}\right)$  is selected.

The output argument `vmod` has the same format as `v`, but it contains the logarithms of the accordingly modified criteria.

## Examples

### Generate Model-Order Combinations and Estimate ARX Model Using IV Method

Create estimation and validation data sets

```
load iddata1;
ze = z1(1:150);
zv = z1(151:300);
```

Generate model-order combinations for estimation, specifying ranges for model orders and delays.

```
NN = struc(1:3,1:2,2:4);
```

Estimate ARX models using the instrumental variable method, and compute the loss function for each model order combination.

```
V = ivstruc(ze,zv,NN);
```

Select the model order with the best fit to the validation data.

```
order = selstruc(V,0);
```

Estimate an ARX model of selected order.

```
M = iv4(ze,order);
```

## Generate Model-Order Combinations and Estimate Multi-Input ARX Model

Create estimation and validation data sets.

```
load co2data;  
Ts = 0.5; % Sample time is 0.5 min  
ze = iddata(Output_exp1,Input_exp1,Ts);  
zv = iddata(Output_exp2,Input_exp2,Ts);
```

Generate model-order combinations for:

- na = 2:4
- nb = 2:5 for the first input, and 1 or 4 for the second input.
- nk = 1:4 for the first input, and 0 for the second input.

```
NN = struc(2:4,2:5,[1 4],1:4,0);
```

Estimate an ARX model for each model order combination.

```
V = arxstruc(ze,zv,NN);
```

Select the model order with the best fit to the validation data.

```
order = selstruc(V,0);
```

Estimate an ARX model of selected order.

```
M = arx(ze,order);
```

**Introduced before R2006a**

## set

Set or modify model properties

### Syntax

```
set(sys, 'Property', Value)
set(sys, 'Property1', Value1, 'Property2', Value2, ...)
sysnew = set( ___ )
set(sys, 'Property')
```

### Description

`set` is used to set or modify the properties of a dynamic system model using property name/property value pairs.

`set(sys, 'Property', Value)` assigns the value `Value` to the property of the model `sys`. 'Property' can be the full property name (for example, 'UserData') or any unambiguous case-insensitive abbreviation (for example, 'user'). The specified property must be compatible with the model type. For example, if `sys` is a transfer function, `Variable` is a valid property but `StateName` is not. For a complete list of available system properties for any linear model type, see the reference page for that model type. This syntax is equivalent to `sys.Property = Value`.

`set(sys, 'Property1', Value1, 'Property2', Value2, ...)` sets multiple property values with a single statement. Each property name/property value pair updates one particular property.

`sysnew = set( ___ )` returns the modified dynamic system model, and can be used with any of the previous syntaxes.

`set(sys, 'Property')` displays help for the property specified by 'Property'.

### Examples

### Specify Model Properties

Create a SISO state-space model with matrices  $A$ ,  $B$ ,  $C$ , and  $D$  equal to 1, 2, 3, and 4, respectively.

```
sys = ss(1,2,3,4);
```

Modify the properties of the model. Add an input delay of 0.1 second, label the input as torque, and set the  $D$  matrix to 0.

```
set(sys, 'InputDelay',0.1, 'InputName', 'torque', 'D',0);
```

View the model properties, and verify the changes.

```
get(sys)

        A: 1
        B: 2
        C: 3
        D: 0
        E: []
    Scaled: 0
    StateName: {''}
    StateUnit: {''}
InternalDelay: [0x1 double]
    InputDelay: 0.1000
    OutputDelay: 0
        Ts: 0
    TimeUnit: 'seconds'
    InputName: {'torque'}
    InputUnit: {''}
    InputGroup: [1x1 struct]
    OutputName: {''}
    OutputUnit: {''}
    OutputGroup: [1x1 struct]
    Notes: [0x1 string]
    UserData: []
        Name: ''
    SamplingGrid: [1x1 struct]
```

## Tips

For discrete-time transfer functions, the convention used to represent the numerator and denominator depends on the choice of variable (see `tf` for details). Like `tf`, the syntax for `set` changes to remain consistent with the choice of variable. For example, if the `Variable` property is set to `'z'` (the default),

```
set(h, 'num', [1 2], 'den', [1 3 4])
```

produces the transfer function

$$h(z) = \frac{z+2}{z^2+3z+4}$$

However, if you change the `Variable` to `'z^-1'` by

```
set(h, 'Variable', 'z^-1'),
```

the same command

```
set(h, 'num', [1 2], 'den', [1 3 4])
```

now interprets the row vectors `[1 2]` and `[1 3 4]` as the polynomials  $1 + 2z^{-1}$  and  $1 + 3z^{-1} + 4z^{-2}$  and produces:

$$\bar{h}(z^{-1}) = \frac{1+2z^{-1}}{1+3z^{-1}+4z^{-2}} = zh(z)$$

---

**Note** Because the resulting transfer functions are different, make sure to use the convention consistent with your choice of variable.

---

## See Also

`frd` | `get` | `idfrd` | `idgrey` | `idnlarx` | `idnlgrey` | `idnlhw` | `idpoly` | `idproc` | `idss` | `idtf` | `ss` | `tf` | `zpk`

**Introduced before R2006a**

## setcov

Set parameter covariance data in identified model

## Syntax

```
sys = setcov(sys0,cov)
```

## Description

`sys = setcov(sys0,cov)` sets the parameter covariance of identified model `sys0` as `cov`.

The model parameter covariance is calculated and stored automatically when a model is estimated. Therefore, you do not need to set the parameter covariance explicitly for estimated models. Use this function for analysis, such as to study how the parameter covariance affects the response of a model obtained by explicit construction.

## Input Arguments

### **sys0**

Identified model.

Identified model, specified as an `idtf`, `idss`, `idgrey`, `idpoly`, `idproc`, or `idnlgrey` model. You cannot set the covariance for nonlinear black-box models (`idnlrx` and `idnlhw`).

### **cov**

Parameter covariance matrix.

`cov` is one of the following:

- an  $np$ -by- $np$  semi-positive definite symmetric matrix, where  $np$  is equal to the number of parameters of `sys0`.

- a structure with the following fields that describe the parameter covariance in a factored form:
  - R — usually the Cholesky factor of inverse of covariance.
  - T — transformation matrix.
  - Free — logical vector of length  $np$  indicating if a parameter is free. Here  $np$  is equal to the number of parameters of `sys0`.

$$\text{cov}(\text{Free}, \text{Free}) = \text{T} * \text{inv}(\text{R}' * \text{R}) * \text{T}'.$$

**Default:**

## Output Arguments

**sys**

Identified model.

The values of all the properties of `sys` are the same as those in `sys0`, except for the parameter covariance values which are modified as specified by `cov`.

## Examples

### Set Raw Covariance Data for Identified Model

Create a transfer function model for the following system:

$$\text{sys0} = \frac{4}{s^2 + 2s + 1}$$

```
sys0 = idtf(4, [1 2 1]);
np = nparams(sys0);
```

`sys0` contains `np` model parameters.

Specify the covariance values for the denominator parameters only.

```
cov = zeros(np);  
den_index = 2:3;  
cov(den_index,den_index) = diag([0.04 0.001]);
```

cov is a covariance matrix with nonzero entries for the denominator parameters.

Set the covariance for `sys0`.

```
sys = setcov(sys0,cov);
```

### See Also

[getcov](#) | [rsample](#) | [setpvec](#) | [sim](#)

**Introduced in R2012a**

# setinit

Set initial states of `idnlgrey` model object

## Syntax

```
model = setinit(model,Property,Values)
```

## Description

`model = setinit(model,Property,Values)` sets the values of the `Property` field of the `InitialStates` model property. `Property` can be `'Name'`, `'Unit'`, `'Value'`, `'Minimum'`, `'Maximum'`, and `'Fixed'`.

## Input Arguments

`model`

Name of the `idnlgrey` model object.

`Property`

Name of the `InitialStates` model property field, such as `'Name'`, `'Unit'`, `'Value'`, `'Minimum'`, `'Maximum'`, and `'Fixed'`.

`Values`

Values of the specified property `Property`. `Values` are an `Nx-by-1` cell array of values, where `Nx` is the number of states.

## See Also

`getinit` | `getpar` | `idnlgrey` | `setpar`

**Introduced in R2007a**

## setoptions

Set plot options for response plot

### Syntax

```
setoptions(h, PlotOpts)
setoptions(h, 'Property1', 'value1', ...)
setoptions(h, PlotOpts, 'Property1', 'value1', ...)
```

### Description

`setoptions(h, PlotOpts)` sets preferences for response plot using the plot handle. `h` is the plot handle, `PlotOpts` is a plot options handle containing information about plot options.

There are two ways to create a plot options handle:

- Use `getoptions`, which accepts a plot handle and returns a plot options handle.  
`p = getoptions(h)`
- Create a default plot options handle using one of the following commands:
  - `bodeoptions` — Bode plots
  - `hsvoptions` — Hankel singular values plots
  - `nicholsoptions` — Nichols plots
  - `nyquistoptions` — Nyquist plots
  - `pzoptions` — Pole/zero plots
  - `sigmaoptions` — Sigma plots
  - `timeoptions` — Time plots (step, initial, impulse, etc.)

For example,

```
p = bodeoptions
```

returns a plot options handle for Bode plots.

`setoptions(h, 'Property1', 'value1', ...)` assigns values to property pairs instead of using `PlotOpts`. To find out what properties and values are available for a particular plot, type `help <function>options`. For example, for Bode plots type

```
help bodeoptions
```

For a list of the properties and values available for each plot type, see “Properties and Values Reference” (Control System Toolbox).

`setoptions(h, PlotOpts, 'Property1', 'value1', ...)` first assigns plot properties as defined in `@PlotOptions`, and then overrides any properties governed by the specified property/value pairs.

## Examples

### Change Frequency Units in Response Plot

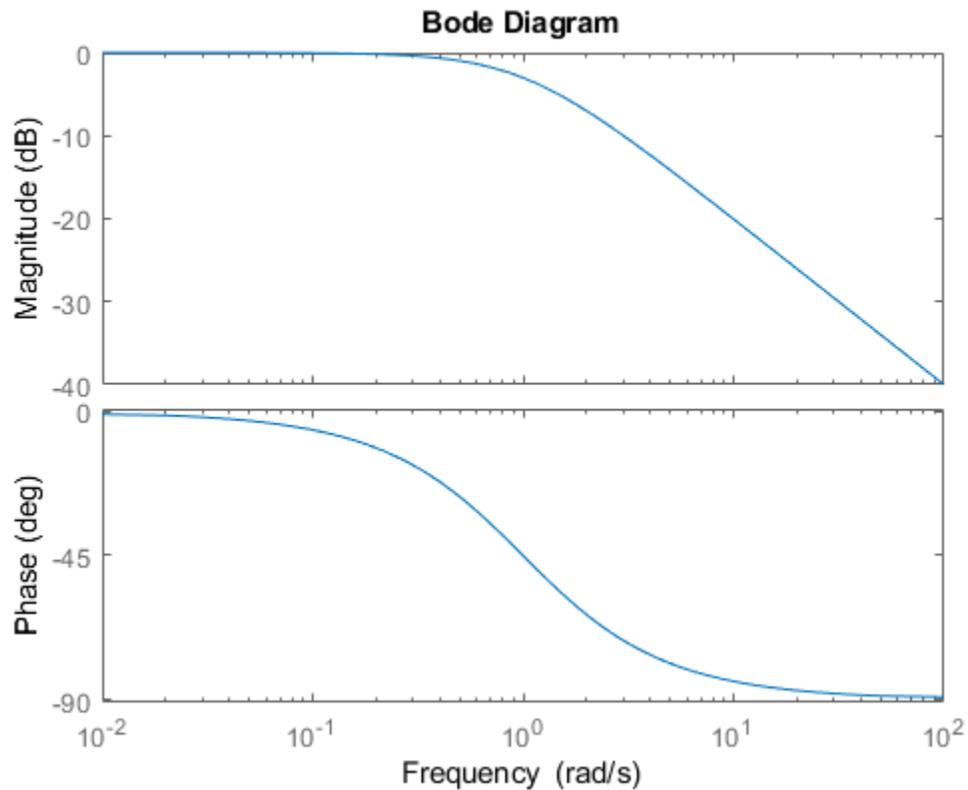
Create the following continuous-time transfer function:

$$H(s) = \frac{1}{s + 1}$$

```
sys = tf(1,[1 1]);
```

Create a Bode plot with plot handle `h`.

```
h = bodeplot(sys);
```



Create a plot options handle p.

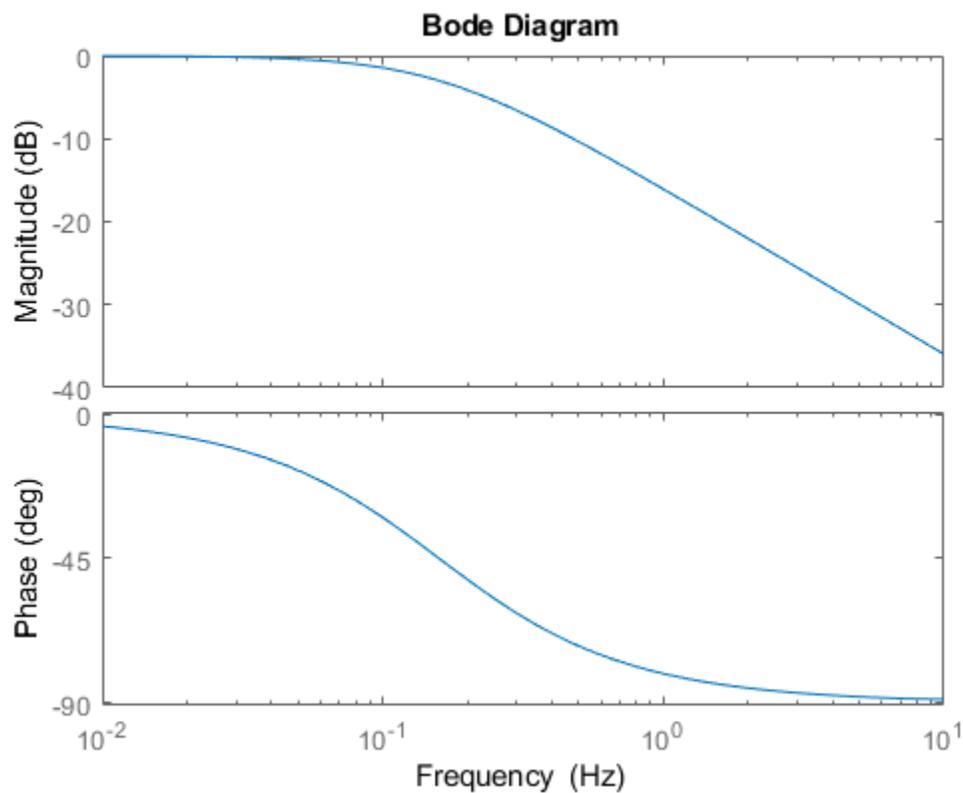
```
p = getoptions(h);
```

Change frequency units of the plot to Hz.

```
p.FreqUnits = 'Hz';
```

Apply the plot options to the Bode plot.

```
setoptions(h,p);
```



Alternatively, use `setoptions(h, 'FrequencyUnits', 'Hz')`.

## See Also

`getoptions`

**Introduced in R2012a**

## idParametric/setpar

Set attributes such as values and bounds of linear model parameters

### Syntax

```
sys1 = setpar(sys,'value',value)
sys1 = setpar(sys,'free',free)
sys1 = setpar(sys,'bounds',bounds)
sys1 = setpar(sys,'label',label)
```

### Description

`sys1 = setpar(sys,'value',value)` sets the parameter values of the model `sys`. For model arrays, use `setpar` separately on each model in the array.

`sys1 = setpar(sys,'free',free)` sets the free or fixed status of the parameters.

`sys1 = setpar(sys,'bounds',bounds)` sets the minimum and maximum bounds on the parameters.

`sys1 = setpar(sys,'label',label)` sets the labels for the parameters.

### Examples

#### Assign Model Parameter Values

Estimate an ARMAX model.

```
load iddata8;
init_data = z8(1:100);
na = 1;
nb = [1 1 1];
nc = 1;
nk = [0 0 0];
sys = armax(init_data,[na nb nc nk]);
```

Set the parameter values.

```
sys = setpar(sys, 'value', [0.5 0.1 0.3 0.02 0.5]);
```

To view the values, type `val = getpar(sys, 'value')`.

### Fix or Free Model Parameters

Construct a process model.

```
m = idproc('P2DUZI');
m.Kp = 1;
m.Tw = 100;
m.Zeta = .3;
m.Tz = 10;
m.Td = 0.4;
```

Set the free status of the parameters.

```
m = setpar(m, 'free', [1 1 1 1 0]);
```

Here, you set Tz to be a fixed parameter.

To check the free status of Tz, type `m.Structure.Tz`.

### Set Minimum and Maximum Bounds on Parameters

Estimate an ARMAX model.

```
load iddata8;
init_data = z8(1:100);
na = 1;
nb = [1 1 1];
nc = 1;
nk = [0 0 0];
sys = armax(init_data, [na nb nc nk]);
```

Set the minimum and maximum bounds for the parameters. Each row represents the bounds for a single parameter. The first value in each row specifies the minimum bound and the second value specifies the maximum bound.

```
sys = setpar(sys, 'bounds', [0 1; 1 1.5; 0 2; 0.5 1; 0 1]);
```

## Assign Default Labels to Parameters

Estimate an ARMAX model.

```
load iddata8;
init_data = z8(1:100);
na = 1;
nb = [1 1 1];
nc = 1;
nk = [0 0 0];
sys = armax(init_data, [na nb nc nk]);
```

Assign default labels to model parameters.

```
sys = setpar(sys, 'label', 'default');
```

View the default labels.

```
getpar(sys, 'label')
ans = 5x1 cell array
    {'A1(1)'}
    {'B0(1)'}
    {'B0(2)'}
    {'B0(3)'}
    {'C1' }
```

## Input Arguments

### **sys** — Identified linear model

idss | idproc | idgrey | idtf | idpoly

Identified linear model, specified as an idss, idproc, idgrey, idtf, or idpoly model object.

### **value** — Parameter values

vector of doubles

Parameter values, specified as a double vector of length `nparams(sys)`.

**free — Free or fixed status of parameters**

vector of logical values

Free or fixed status of parameters, specified as a logical vector of length `nparams(sys)`.

**bounds — Minimum and maximum bounds on parameters**

matrix of doubles

Minimum and maximum bounds on parameters, specified as a double matrix of size `nparams(sys)-by-2`. The first column specifies the minimum bound and the second column the maximum bound.

**label — Parameter labels**

cell array of character vectors | 'default'

Parameter labels, specified as a cell array of character vectors. The cell array is of length `nparams(sys)`. For example, {'a1', 'a3'}, if `nparams(sys)` is two.

Use 'default' to assign default labels, A1, A2, ..., B1, B2, ..., to the parameters.

## Output Arguments

**sys1 — Model with specified values of parameter attributes**

idss | idproc | idgrey | idtf | idpoly

Model with specified values of parameter attributes. The model `sys` you specify as the input to `setpar` gets updated with the specified parameter attribute values.

## See Also

getpar | setcov | setpvec

**Introduced in R2013b**

## setpar

Set initial parameter values of `idnlgrey` model object

### Syntax

```
setpar(model,property,values)
```

### Input Arguments

`model`

Name of the `idnlgrey` model object.

`property`

Name of the Parameters model property field, such as 'Name', 'Unit', 'Value', 'Minimum', or 'Maximum'.

Default: 'Value'.

`values`

Values of the specified property `Property`. `values` are an `Np`-by-1 cell array of values, where `Np` is the number of parameters.

### Description

`setpar(model,property,values)` sets the model parameter values in the `property` field of the Parameters model property. `property` can be 'Name', 'Unit', 'Value', 'Minimum', and 'Maximum'.

### See Also

`getinit` | `getpar` | `idnlgrey` | `setinit`

**Introduced in R2007a**

# setPolyFormat

Specify format for  $B$  and  $F$  polynomials of multi-input polynomial model

## Syntax

```
modelOut = setPolyFormat(modelIn,'double')  
modelOut = setPolyFormat(modelIn,'cell')
```

## Description

`modelOut = setPolyFormat(modelIn,'double')` converts the  $B$  and  $F$  polynomials of a multi-input polynomial model, `modelIn`, to double matrices.

By default, the  $B$  and  $F$  polynomials of an `idpoly` model are cell arrays. For MATLAB scripts written before R2012a, convert the cell arrays to double matrices for backward compatibility using this syntax. For example:

```
model = arx(data,[3 2 2 1 1]);  
model = setPolyFormat(model,'double');
```

`modelOut = setPolyFormat(modelIn,'cell')` converts the  $B$  and  $F$  polynomials of `modelIn` to cell arrays.

MATLAB data files saved before R2012a store `idpoly` models with their  $B$  and  $F$  polynomials represented as double matrices. If these models were previously set to operate in backward-compatibility mode, they are not converted to use cell arrays when loaded. Convert these models to use cell arrays using this syntax. For example:

```
load polyData.mat model;  
model = setPolyFormat(model,'cell');
```

## Examples

### Convert $B$ and $F$ Polynomials of a Multi-Input ARX Model to Double Matrices

Load estimation data.

```
load iddata8;
```

Estimate the model.

```
m1 = arx(z8,[3 [2 2 1] [1 1 1]]);
```

Convert the  $b$  and  $f$  polynomials to use double matrices.

```
m2 = setPolyFormat(m1,'double');
```

Extract pole and zero information from the model using matrix syntax.

```
Poles1 = roots(m2.F(1,:));  
Zeros1 = roots(m2.B(1,:));
```

## Input Arguments

### **modelIn** — Polynomial model

*idpoly* object

Polynomial model, specified as an *idpoly* object. The  $B$  and  $F$  polynomials of *modelIn* are either:

- Cell arrays with  $N_u$  elements, where  $N_u$  is the number of model inputs, with each element containing a double vector. This configuration is the default.
- Double matrices with  $N_u$  rows. This configuration applies to backward-compatible *idpoly* models stored in MATLAB data files before R2012a.

---

**Note** `setPolyFormat` only supports multi-input, single-output models. Specifying *modelIn* as a:

- Multi-output model generates an error.
  - Single-input, single-output model has no effect. The  $B$  and  $F$  polynomials remain as double vectors.
-

## Output Arguments

### **modelOut** — Polynomial model

idpoly object

Polynomial model, returned as an idpoly object.

To access the b and f polynomials of modelOut, use:

- Matrix syntax after using `modelOut = setPolyFormat(modelIn, 'double')`. For example:

```
modelOut.B(1,:);
```

- Cell array syntax after using `modelOut = setPolyFormat(modelIn, 'cell')`. For example:

```
modelOut.B{1};
```

After using `modelOut = setPolyFormat(modelIn, 'cell')`, you can resave the converted model in cell array format. For example:

```
save polyNew.mat modelOut;
```

## Tips

- To verify the current format of the B and F polynomials for a given idpoly model, enter:

```
class(model.B)
```

If the model uses double matrices, the displayed result is:

```
ans =
```

```
double
```

Otherwise, for cell arrays, the result is:

```
ans =
```

```
cell
```

## **See Also**

get | idpoly | polydata | set | tfdata

## **Topics**

“Extracting Numerical Model Data”

**Introduced in R2010a**

# setpvec

Modify value of model parameters

## Syntax

```
sys = setpvec(sys0,par)
sys = setpvec(sys0,par,'free')
```

## Description

`sys = setpvec(sys0,par)` modifies the value of the parameters of the identified model `sys0` to the value specified by `par`.

`par` must be of length `nparams(sys0)`. `nparams(sys0)` returns a count of all the parameters of `sys0`.

`sys = setpvec(sys0,par,'free')` modifies the value of all the free estimation parameters of `sys0` to the value specified by `par`.

`par` must be of length `nparams(sys0,'free')`. `nparams(sys0,'free')` returns a count of all the free parameters of `sys0`. For `idnlarx` and `idnlhw` models, all parameters are treated as free.

## Input Arguments

### **sys0**

Identified model, specified as an `idtf`, `idss`, `idgrey`, `idpoly`, `idproc`, `idnlarx`, `idnlhw`, or `idnlgrey` object.

### **par**

Replacement value for the parameters of the identified model `sys0`.

For the syntax `sys = setpvec(sys0, par)`, `par` must be of length `nparams(sys0)`. `nparams(sys0)` returns a count of all the parameters of `sys0`.

For the syntax `sys = setpvec(sys0, par, 'free')`, `par` must be of length `nparams(sys0, 'free')`. `nparams(sys0, 'free')` returns a count of all the free parameters of `sys0`.

Use `NaN` to denote unknown parameter values.

If `sys0` is an array of models, then specify `par` as a cell array with an entry corresponding to each model in `sys0`.

## Output Arguments

### `sys`

Identified model obtained from `sys0` by updating the values of the specified parameters.

## Examples

### Modify Parameter Values of Transfer Function Model

Construct a transfer function model.

```
sys0 = idtf(1,[1 2]);
```

Define a parameter vector and use it to set the model parameters. The second parameter is set to `NaN`, indicating that its value is unknown.

```
par = [1;NaN;0];  
sys = setpvec(sys0,par);
```

The constructed model, `sys`, can be used to initialize a model estimation.

### Modify Free Parameter Values of Transfer Function Model

Construct a transfer function model.

```
sys0 = idtf([1 0],[1 2 0]);
```

Set the first three parameters of `sys0` as free parameters.

```
sys0 = setpar(sys0, 'free', [1 1 1 0 0]);
```

Define a parameter vector and use it to set the free model parameters.

```
par = [1;2;1];  
sys = setpvec(sys0,par, 'free');
```

## See Also

[getpvec](#) | [nparams](#) | [setcov](#)

**Introduced in R2012a**

## sgrid

Generate s-plane grid of constant damping factors and natural frequencies

### Syntax

```
sgrid  
sgrid(z,wn)
```

### Description

`sgrid` generates, for pole-zero and root locus plots, a grid of constant damping factors from zero to one in steps of 0.1 and natural frequencies from zero to 10 rad/sec in steps of one rad/sec, and plots the grid over the current axis. If the current axis contains a continuous s-plane root locus diagram or pole-zero map, `sgrid` draws the grid over the plot.

`sgrid(z,wn)` plots a grid of constant damping factor and natural frequency lines for the damping factors and natural frequencies in the vectors `z` and `wn`, respectively. If the current axis contains a continuous s-plane root locus diagram or pole-zero map, `sgrid(z,wn)` draws the grid over the plot.

Alternatively, you can select **Grid** from the right-click menu to generate the same s-plane grid.

### Examples

#### Generate S-Plane Grid on Root Locus Plot

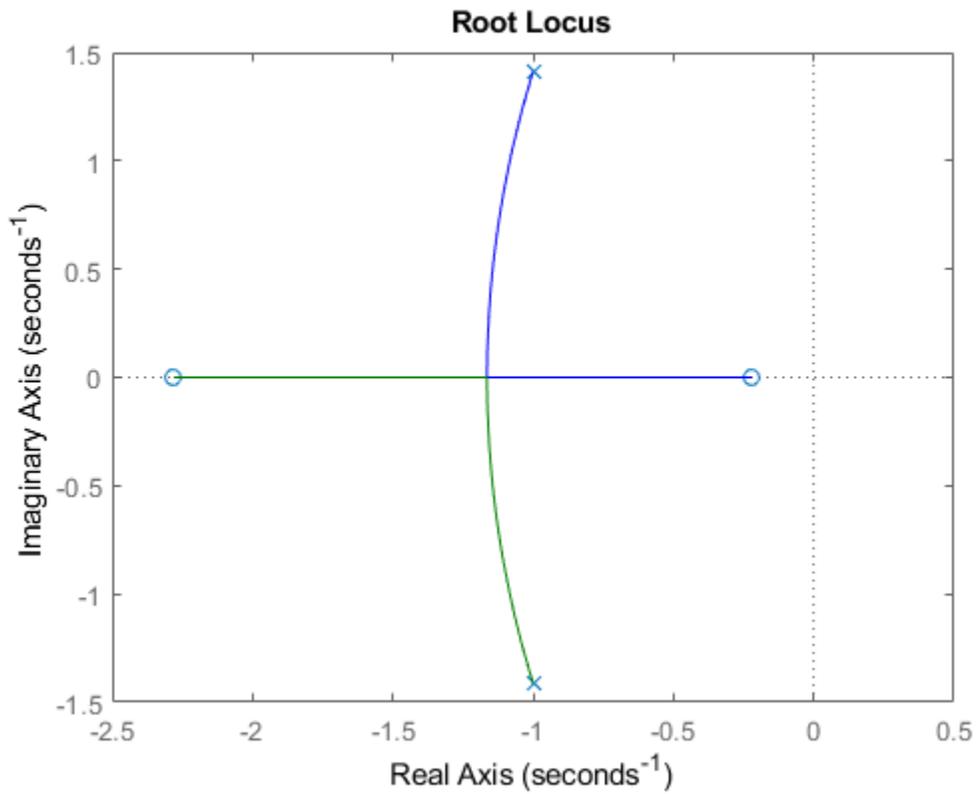
Create the following continuous-time transfer function:

$$H(s) = \frac{2s^2 + 5s + 1}{s^2 + 2s + 3}$$

```
H = tf([2 5 1],[1 2 3]);
```

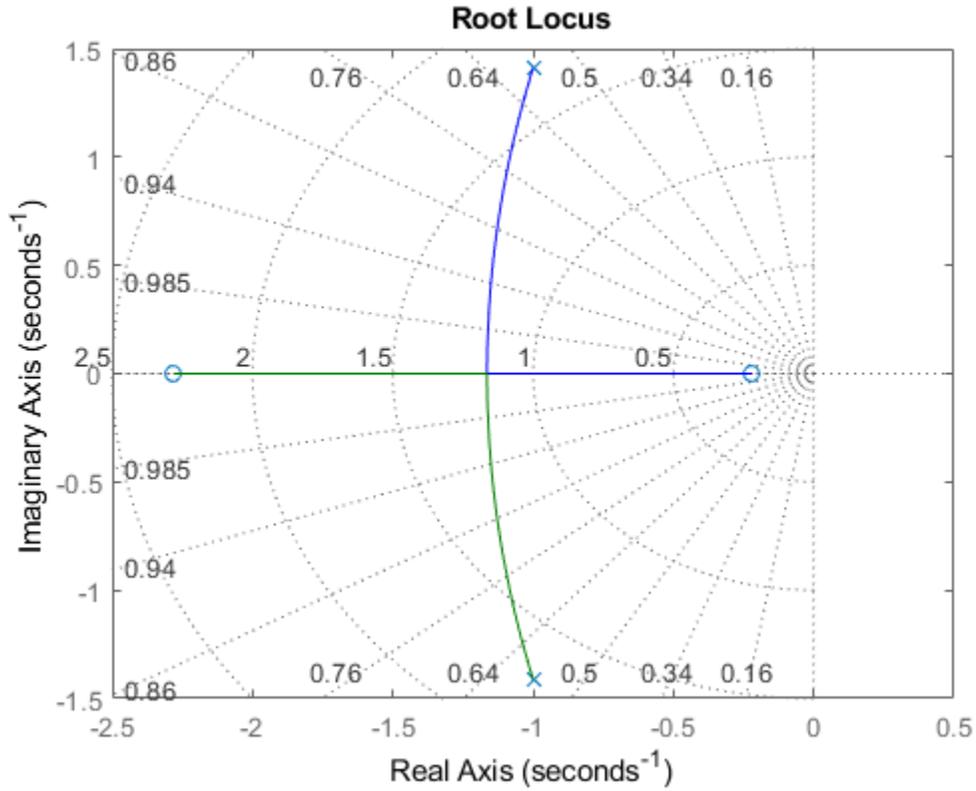
Plot the root locus of the transfer function.

```
rlocus(H)
```



Plot s-plane grid lines on the root locus.

```
sgrid
```



## See Also

[pzmap](#) | [rlocus](#) | [zgrid](#)

Introduced in R2012a

# showConfidence

Display confidence regions on response plots for identified models

## Syntax

```
showConfidence(plot_handle)
showConfidence(plot_handle, sd)
```

## Description

`showConfidence(plot_handle)` displays the confidence region on the response plot, with handle `plot_handle`, for an identified model.

`showConfidence(plot_handle, sd)` displays the confidence region for `sd` standard deviations.

## Input Arguments

### **plot\_handle**

Response plot handle.

`plot_handle` is the handle for the response plot of an identified model on which the confidence region is displayed. It is obtained as an output of one of the following plot commands: `bodeplot`, `stepplot`, `impzplot`, `nyquistplot`, or `iopzplot`.

### **sd**

Standard deviation of the confidence region. A common choice is 3 standard deviations, which gives 99.7% significance.

**Default:** `getoptions(plot_handle, 'ConfidenceRegionNumberSD')`

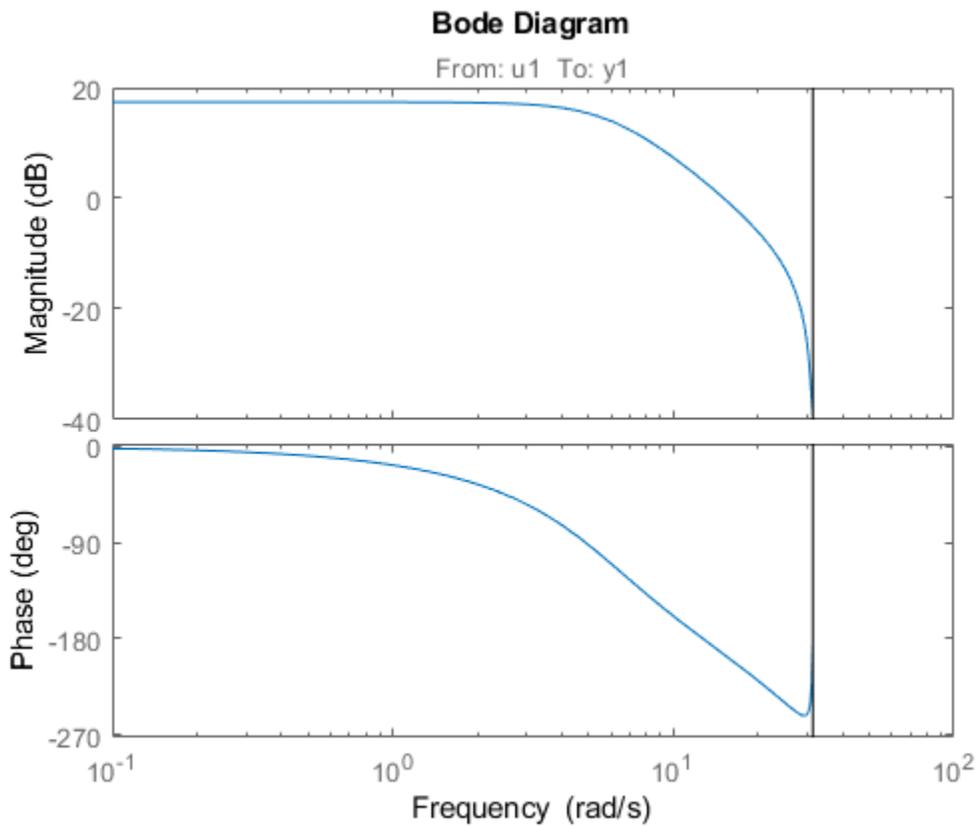
## Examples

### View Confidence Region for Identified Model

Show the confidence bounds on the bode plot of an identified ARX model.

Obtain identified model and plot its bode response.

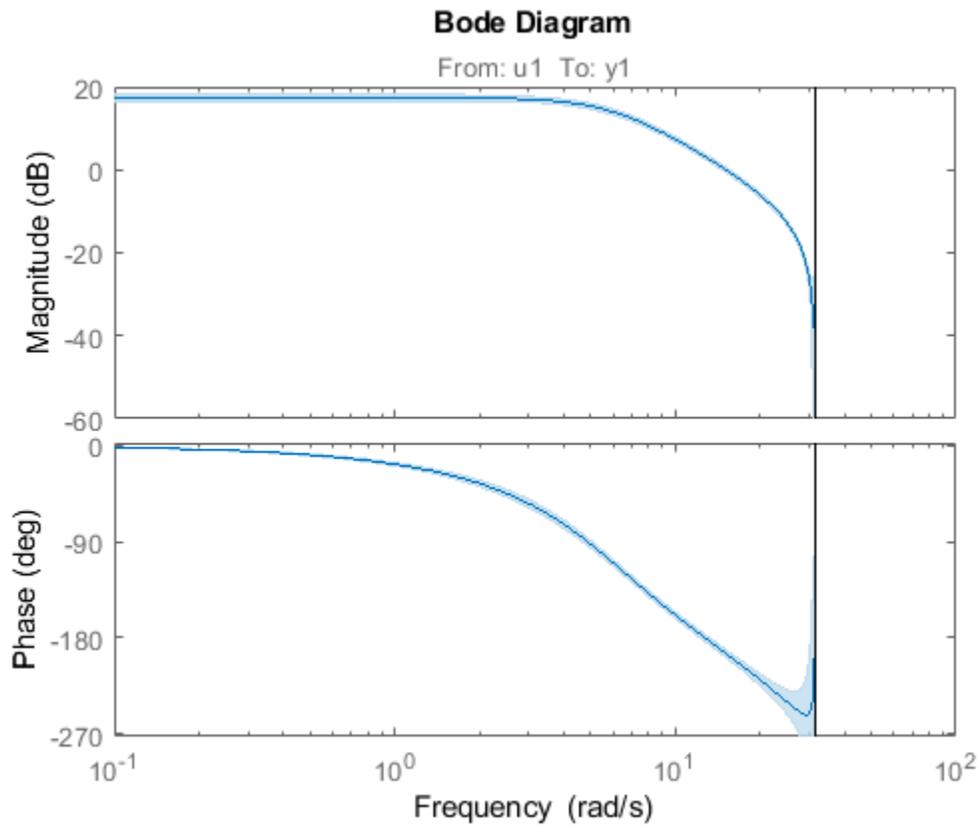
```
load iddata1 z1
sys = arx(z1, [2 2 1]);
h = bodeplot(sys);
```



`z1` is an `iddata` object that contains time domain system response data. `sys` is an `idpoly` model containing the identified polynomial model. `h` is the plot handle for the bode response plot of `sys`.

Show the confidence bounds for `sys`.

```
showConfidence(h);
```



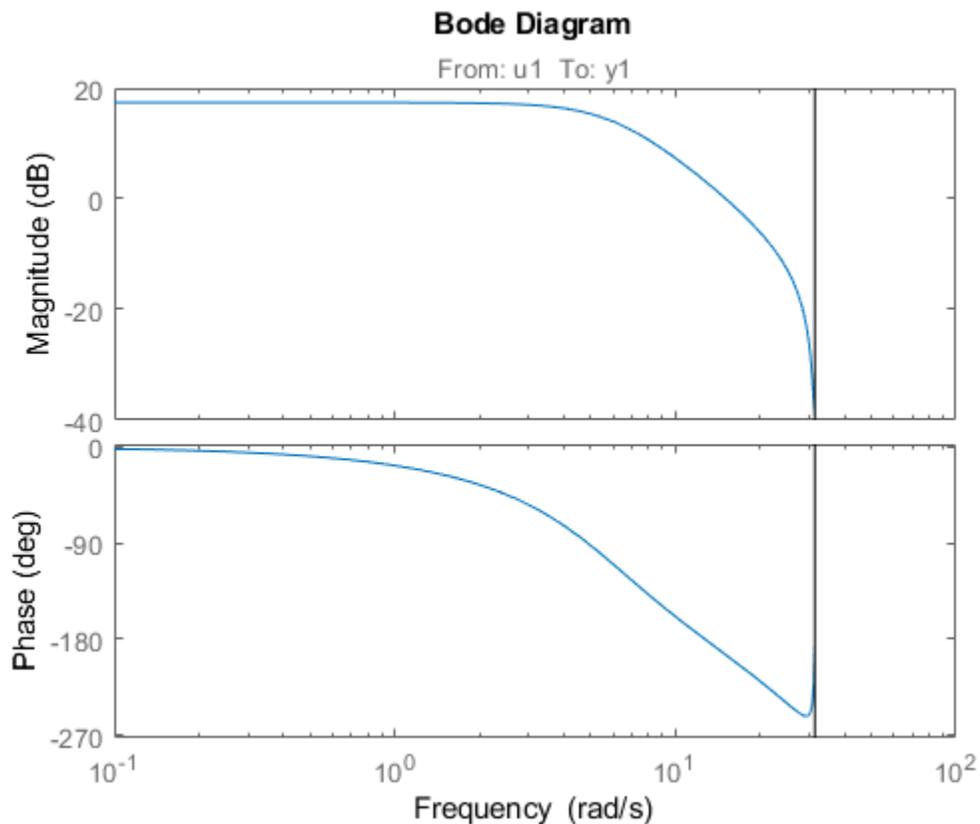
This plot depicts the confidence region for 1 standard deviation.

### Specify the Standard Deviation of the Confidence Region

Show the confidence bounds on the bode plot of an identified ARX model.

Obtain identified model and plot its bode response.

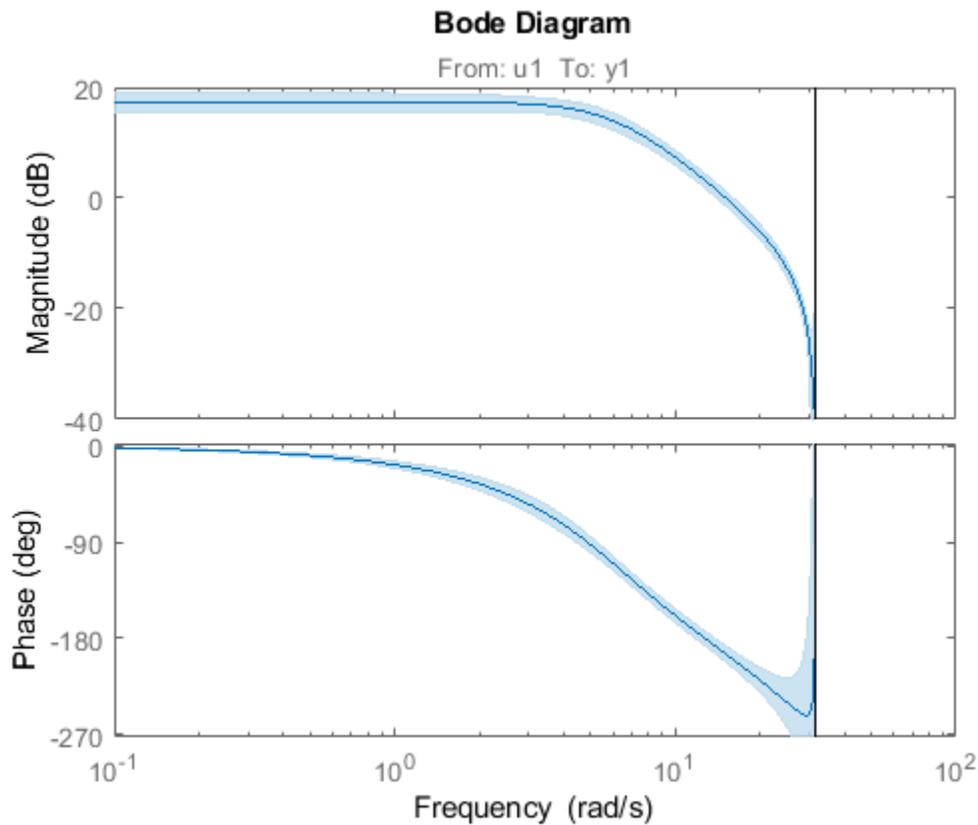
```
load iddata1 z1
sys = arx(z1, [2 2 1]);
h = bodeplot(sys);
```



`z1` is an `iddata` object that contains time domain system response data. `sys` is an `idpoly` model containing the identified polynomial model. `h` is the plot handle for the bode response plot of `sys`.

Show the confidence bounds for `sys` using 2 standard deviations.

```
sd = 2;  
showConfidence(h,sd);
```



`sd` specifies the number of standard deviations for the confidence region displayed on the plot.

## Alternatives

You can interactively turn on the confidence region display on a response plot. Right-click the response plot, and select **Characteristics > Confidence Region**.

## **See Also**

bodeplot | impulseplot | iopzplot | nyquistplot | stepplot

**Introduced in R2012a**

# sigmoidnet

Class representing sigmoid network nonlinearity estimator for nonlinear ARX and Hammerstein-Wiener models

## Syntax

```
s=sigmoidnet('NumberOfUnits',N)
s=sigmoidnet(Property1,Value1,...PropertyN,ValueN)
```

## Description

`sigmoidnet` is an object that stores the sigmoid network nonlinear estimator for estimating nonlinear ARX and Hammerstein-Wiener models.

You can use the constructor to create the nonlinearity object, as follows:

`s=sigmoidnet('NumberOfUnits',N)` creates a sigmoid nonlinearity estimator object with `N` terms in the sigmoid expansion.

`s=sigmoidnet(Property1,Value1,...PropertyN,ValueN)` creates a sigmoid nonlinearity estimator object specified by properties in “sigmoidnet Properties” on page 1-1549.

Use `evaluate(s,x)` to compute the value of the function defined by the `sigmoidnet` object `s` at `x`.

## sigmoidnet Properties

You can include property-value pairs in the constructor to specify the object.

After creating the object, you can use `get` or dot notation to access the object property values. For example:

```
% List all property values
get(s)
```

```
% Get value of NumberOfUnits property  
s.NumberOfUnits
```

You can also use the `set` function to set the value of particular properties. For example:

```
set(s, 'LinearTerm', 'on')
```

The first argument to `set` must be the name of a MATLAB variable.

Property Name	Description
NumberOfUnits	Integer specifies the number of nonlinearity units in the expansion. Default=10.  For example: <code>sigmoidnet(H, 'NumberOfUnits', 5)</code>
LinearTerm	Can have the following values: <ul style="list-style-type: none"><li>• 'on'—Estimates the vector <math>L</math> in the expansion.</li><li>• 'off'—Fixes the vector <math>L</math> to zero.</li></ul> For example: <code>sigmoidnet(H, 'LinearTerm', 'on')</code>

Property Name	Description
Parameters	<p>A structure containing the parameters in the nonlinear expansion, as follows:</p> <ul style="list-style-type: none"> <li>• <b>RegressorMean</b>: 1-by-m vector containing the means of <math>x</math> in estimation data, <math>r</math>.</li> <li>• <b>NonLinearSubspace</b>: m-by-q matrix containing <math>Q</math>.</li> <li>• <b>LinearSubspace</b>: m-by-p matrix containing <math>P</math>.</li> <li>• <b>LinearCoef</b>: p-by-1 vector <math>L</math>.</li> <li>• <b>Dilation</b>: q-by-n matrix containing the values <math>b_n</math>.</li> <li>• <b>Translation</b>: 1-by-n vector containing the values <math>c_n</math>.</li> <li>• <b>OutputCoef</b>: n-by-1 vector containing the values <math>a_n</math>.</li> <li>• <b>OutputOffset</b>: scalar <math>d</math>.</li> </ul> <p>Typically, the values of this structure are set by estimating a model with a <code>sigmoidnet</code> nonlinearity.</p>

## Examples

Use `sigmoidnet` to specify the nonlinear estimator in nonlinear ARX and Hammerstein-Wiener models. For example:

```
m=nlarx(Data,Orders,sigmoidnet('num',5));
```

## Tips

Use `sigmoidnet` to define a nonlinear function  $y = F(x)$ , where  $y$  is scalar and  $x$  is an  $m$ -dimensional row vector. The sigmoid network function is based on the following expansion:

$$F(x) = (x-r)PL + a_1 f((x-r)Qb_1 + c_1) + \dots + a_n f((x-r)Qb_n + c_n) + d$$

where  $f$  is the sigmoid function, given by the following equation:

$$f(z) = \frac{1}{e^{-z} + 1}.$$

$P$  and  $Q$  are  $m$ -by- $p$  and  $m$ -by- $q$  projection matrices. The projection matrices  $P$  and  $Q$  are determined by principal component analysis of estimation data. Usually,  $p=m$ . If the components of  $x$  in the estimation data are linearly dependent, then  $p < m$ . The number of columns of  $Q$ ,  $q$ , corresponds to the number of components of  $x$  used in the sigmoid function.

When used in a nonlinear ARX model,  $q$  is equal to the size of the `NonlinearRegressors` property of the `idnlarx` object. When used in a Hammerstein-Wiener model,  $m=q=1$  and  $Q$  is a scalar.

$r$  is a  $1$ -by- $m$  vector and represents the mean value of the regressor vector computed from estimation data.

$d$ ,  $a$ , and  $c$  are scalars.

$L$  is a  $p$ -by- $1$  vector.

$b$  are  $q$ -by- $1$  vectors.

## Algorithms

`sigmoidnet` uses an iterative search technique for estimating parameters.

## See Also

`nlarx` | `nlnhw`

**Introduced in R2007a**

# sim

Simulate response of identified model

## Syntax

```
y = sim(sys, udata)
y = sim(sys, udata, opt)

[y, y_sd] = sim( ___ )
[y, y_sd, x] = sim( ___ )
[y, y_sd, x, x_sd] = sim( ___ )

sim( ___ )
```

## Description

`y = sim(sys, udata)` returns the simulated response of an identified model using the input data, `udata`. By default, zero initial conditions are used for all model types except `idnlgrey`, in which case the initial conditions stored internally in the model are used.

`y = sim(sys, udata, opt)` uses the option set, `opt`, to configure the simulation option, including the specification of initial conditions.

`[y, y_sd] = sim( ___ )` returns the estimated standard deviation, `y_sd`, of the simulated response.

`[y, y_sd, x] = sim( ___ )` returns the state trajectory, `x`, for state-space models.

`[y, y_sd, x, x_sd] = sim( ___ )` returns the standard deviation of the state trajectory, `x_sd`, for state-space models.

`sim( ___ )` plots the simulated response of the identified model.

## Examples

## Simulate State-Space Model Using Input Data

Load the estimation data.

```
load iddata2 z2
```

Estimate a third-order state-space model.

```
sys = ssest(z2,3);
```

Simulate the identified model using the input channels from the estimation data.

```
y = sim(sys,z2);
```

## Add Noise to Simulated Model Response

Load the data, and obtain the identified model.

```
load iddata2 z2  
sys = n4sid(z2,3);
```

sys is a third-order state-space model estimated using a subspace method.

Create a simulation option set to add noise to the simulated model response.

```
opt1 = simOptions('AddNoise',true);
```

Simulate the model.

```
y = sim(sys,z2,opt1);
```

Default Gaussian white noise is filtered by the noise transfer function of the model and added to the simulated model response.

You can also add your own noise signal, e, using the NoiseData option.

```
e = randn(length(z2.u),1);  
opt2 = simOptions('AddNoise',true,'NoiseData',e);
```

Simulate the model.

```
y = sim(sys,z2,opt2);
```

## Simulate Model Using Initial Conditions Obtained During Estimation

Load data.

```
load iddata1 z1
```

Specify the estimation option to estimate the initial state.

```
estimOpt = ssestOptions('InitialState','estimate');
```

Estimate a state-space model, and return the value of the estimated initial state.

```
[sys,x0] = ssest(z1,2,estimOpt);
```

Specify initial conditions for simulation

```
simOpt = simOptions('InitialCondition',x0);
```

Simulate the model, and obtain the model response and standard deviation.

```
[y,y_sd] = sim(sys,z1,simOpt);
```

## Estimate Standard Deviation and State Trajectory for State-Space Models

Load estimation data, and estimate a state-space model.

```
load iddata1 z1  
sys = ssest(z1,2);
```

Return the standard deviation and state trajectory.

```
[y,y_sd,x] = sim(sys,z1);
```

## Estimate State Trajectory and Standard Deviations of Simulated Response

Load estimation data, and estimate a state-space model.

```
load iddata1 z1  
sys = ssest(z1,2);
```

Create a simulation option set, and specify the initial states.

```
opt = simOptions('InitialCondition',[1;2]);
```

Specify the covariance of the initial states.

```
opt.X0Covariance = [0.1 0; 0 0.1];
```

Calculate the standard deviations of simulated response, `y_sd`, and state trajectory, `x_sd`.

```
[y,y_sd,x,x_sd] = sim(sys,z1,opt);
```

## Plot Simulated Model Response

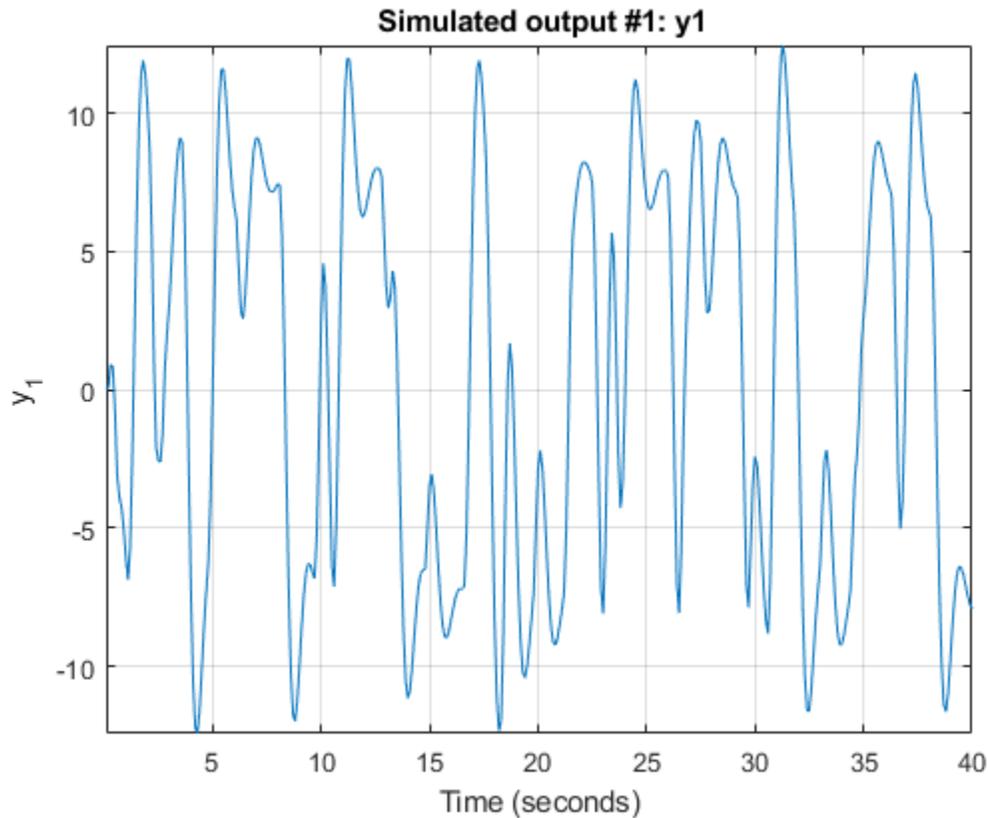
Obtain the identified model.

```
load iddata2 z2  
sys = tfest(z2,3);
```

`sys` is an `idtf` model that encapsulates the third-order transfer function estimated for the measured data `z2`.

Simulate the model.

```
sim(sys,z2)
```



### Simulate Nonlinear ARX Model

Simulate a single-input single-output nonlinear ARX model around a known equilibrium point, with an input level of 1 and output level of 10.

Load the sample data.

```
load iddata2
```

Estimate a nonlinear ARX model from the data.

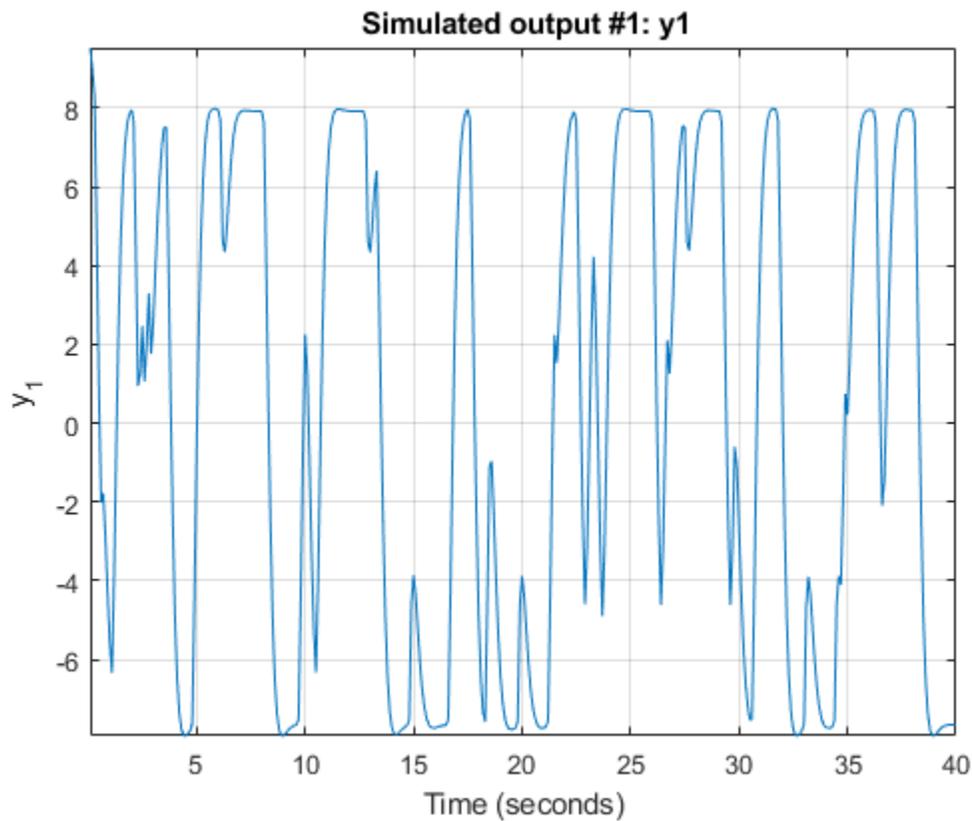
```
M = nlarx(z2,[2 2 1], 'treepartition');
```

Estimate current states of model based on past data. Specify as many past samples as there are lags in the input and output variables (2 here).

```
x0 = data2state(M,struct('Input',ones(2,1),'Output',10*ones(2,1)));
```

Simulate the model using the initial states returned by `data2state`.

```
opt = simOptions('InitialCondition',x0);  
sim(M,z2,opt)
```



## Continue from End of Previous Simulation

Continue the simulation of a nonlinear ARX model from the end of a previous simulation run.

Estimate a nonlinear ARX model from data.

```
load iddata2
M = nlarx(z2,[2 2 1],'treepartition');
```

Simulate the model using the first half of the input data z2. Start the simulation from zero initial states.

```
u1 = z2(1:200,[]);
opt1 = simOptions('InitialCondition','zero');
ys1 = sim(M,u1,opt1);
```

Start another simulation using the second half of the input data z2. Use the same states of the model from the end of the first simulation.

```
u2 = z2(201:end,[]);
```

To set the initial states for the second simulation correctly, package input u1 and output ys1 from the first simulation into one iddata object. Pass this data as initial conditions for the next simulation.

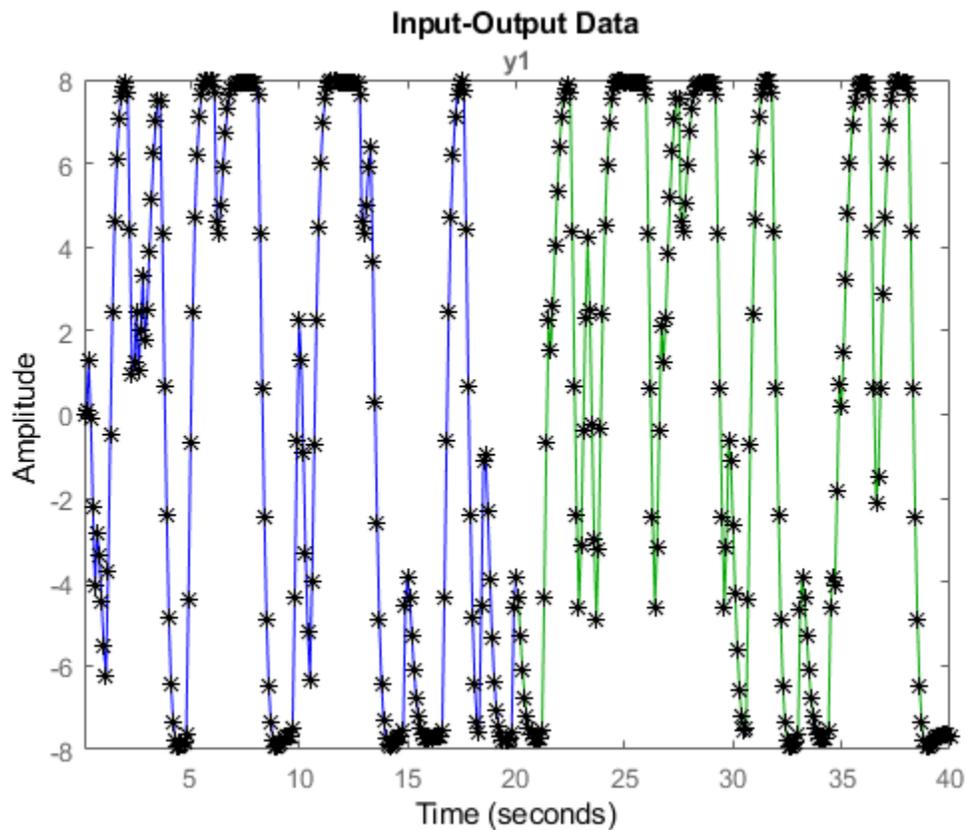
```
firstSimData = [ys1,u1];
opt2 = simOptions('InitialCondition',firstSimData);
ys2 = sim(M,u2,opt2);
```

Verify the two simulations by comparing to a complete simulation using all the input data z2. First, extract the whole set of input data.

```
uTotal = z2(:,[]);
opt3 = simOptions('InitialCondition','zero');
ysTotal = sim(M,uTotal,opt3);
```

Plot the three responses ys1, ys2 and ysTotal. ys1 should be equal to first half of ysTotal. ys2 should be equal to the second half of ysTotal.

```
plot(ys1,'b',ys2,'g',ysTotal,'k*')
```



The plot shows that the three responses `ys1`, `ys2`, and `ysTotal` overlap as expected.

### Match Model Response to Output Data

Estimate initial states of model `M` such that, the response best matches the output in data set `z2`.

Load the sample data.

```
load iddata2;
```

Estimate a nonlinear ARX model from the data.

```
M = nlarx(z2,[4 3 2],wavenet('NumberOfUnits',20));
```

Estimate the initial states of M to best fit z2.y in the simulated response.

```
x0 = findstates(M,z2,Inf);
```

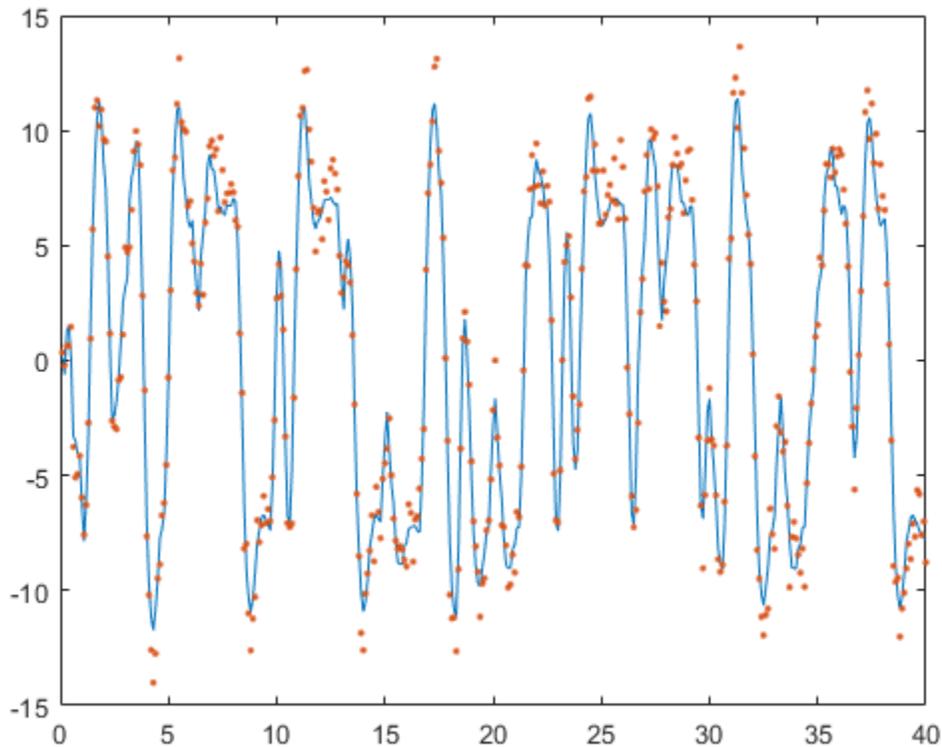
Simulate the model.

```
opt = simOptions('InitialCondition',x0);
```

```
ysim = sim(M,z2.u,opt);
```

Compare the simulated model output ysim with the output signal in z2.

```
time = z2.SamplingInstants;  
plot(time,ysim,time,z2.y,'.')
```



## Simulate Model Near Steady State with Known Input and Unknown Output

Start simulation of a model near steady state, where the input is known to be 1, but the output is unknown.

Load the sample data.

```
load iddata2
```

Estimate a nonlinear ARX model from the data.

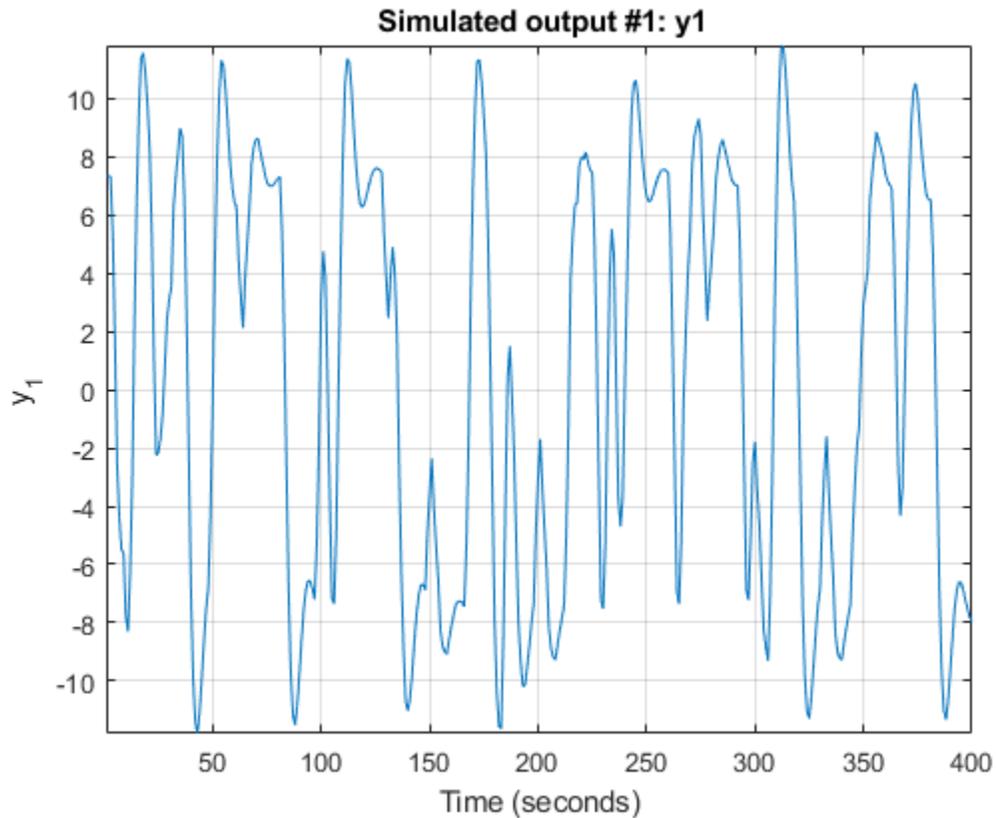
```
M = nlarx(z2,[4 3 2], 'wavenet');
```

Determine equilibrium state values for input 1 and unknown target output.

```
x0 = findop(M, 'steady', 1, NaN);
```

Simulate the model using initial states  $x_0$ .

```
opt = simOptions('InitialCondition', x0);  
sim(M, z2.u, opt)
```



### Simulate Hammerstein-Wiener Model at Steady-State Operating Point

Load the sample data.

```
load iddata2
```

Create a Hammerstein-Wiener model.

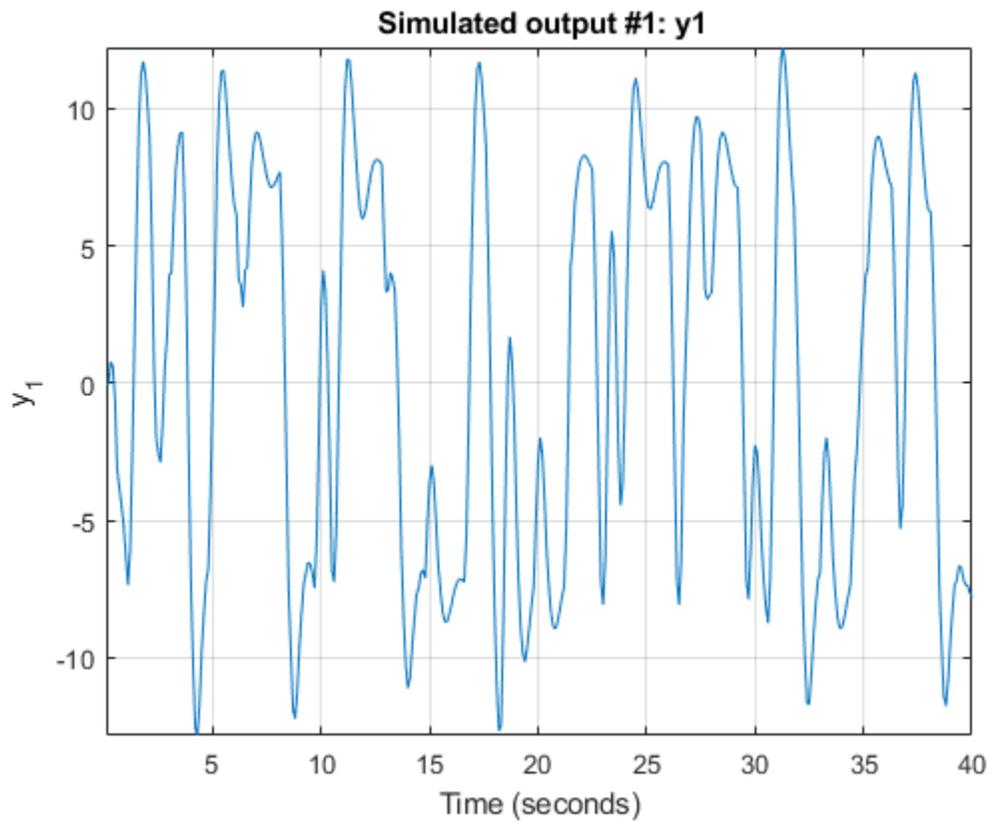
```
M = nlhw(z2,[4 3 2],[],'pwlinear');
```

Compute steady-state operating point values corresponding to an input level of 1 and an unknown output level.

```
x0 = findop(M, 'steady', 1, NaN);
```

Simulate the model using the estimated initial states.

```
opt = simOptions('InitialCondition', x0);  
sim(M, z2.u)
```



## Simulate Time Series Model

Load time series data, and estimate an AR model using the least-squares approach.

---

```
load iddata9 z9
sys = ar(z9,6,'ls');
```

For time series data, specify the desired simulation length,  $N = 200$  using an  $N$ -by-0 input data set.

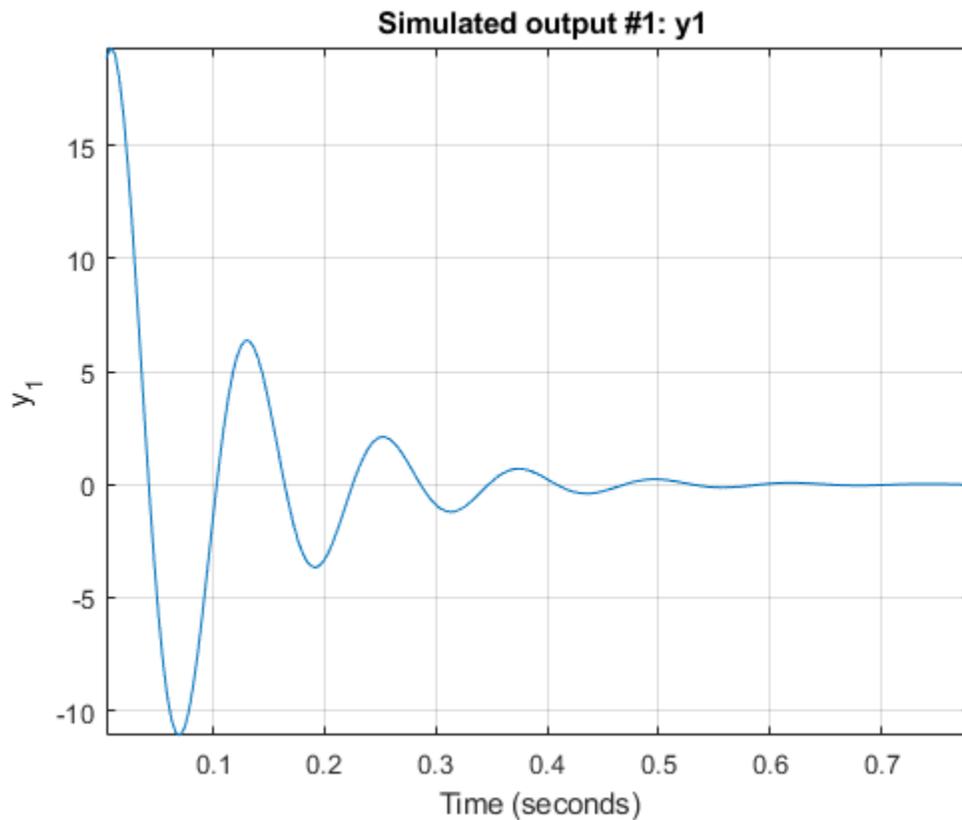
```
data = iddata([],zeros(200,0),z9.Ts);
```

Set the initial conditions to use the initial samples of the time series as historical output samples.

```
IC = struct('Input',[],'Output',z9.y(1:6));
opt = simOptions('InitialCondition',IC);
```

Simulate the model.

```
sim(sys,data,opt)
```



### **Understand Use of Historical Data for Model Simulation**

Use historical input-output data as a proxy for initial conditions when simulating your model. You first simulate using the `sim` command and specify the historical data using the `simOptions` option set. You then reproduce the simulated output by manually mapping the historical data to initial states.

Load a two-input, one-output data set.

```
load iddata7 z7
```

Identify a fifth-order state-space model using the data.

```
sys = n4sid(z7,5);
```

Split the data set into two parts.

```
zA = z7(1:15);  
zB = z7(16:end);
```

Simulate the model using the input signal in zB.

```
uSim = zB;
```

Simulation requires initial conditions. The signal values in zA are the historical data, that is, they are the input and output values for the time immediately preceding data in zB. Use zA as a proxy for the required initial conditions.

```
I0 = struct('Input',zA.InputData,'Output',zA.OutputData);  
opt = simOptions('InitialCondition',I0);
```

Simulate the model.

```
ysim = sim(sys,uSim,opt);
```

Now reproduce the output by manually mapping the historical data to initial states of sys. To do so, use the `data2state` command.

```
xf = data2state(sys,zA);
```

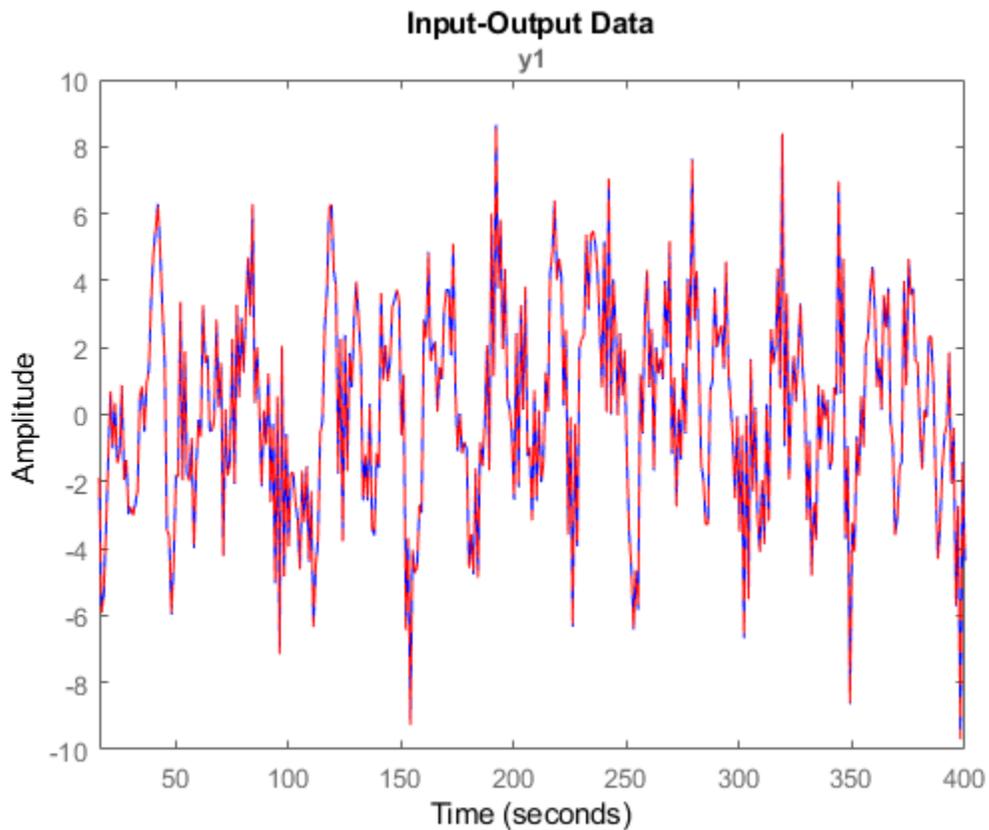
xf contains the state values of sys at the time instant immediately after the most recent data sample in zA.

Simulate the system using xf as the initial states.

```
opt2 = simOptions('InitialCondition',xf);  
ysim2 = sim(sys,uSim,opt2);
```

Plot the output of the `sim` command `ysim` and the manually computed results `ysim2`.

```
plot(ysim,'b',ysim2,'--r')
```



`ysim2` is the same as `ysim`.

## Input Arguments

### **sys** — Identified model

identified linear model | identified nonlinear model

Identified model, specified as one of the following model objects:

	<b>Model Type</b>	<b>Model Object</b>
<b>Identified Linear Model</b>	Polynomial model	idpoly
	Process model	idproc
	State-space model	idss
	Transfer function model	idtf
	Linear grey-box model	idgrey
<b>Identified Nonlinear Model</b>	Nonlinear ARX model	idnlarx
	Nonlinear Hammerstein-Wiener model	idnlhw
	Nonlinear grey-box model	idnlgrey

### **udata — Simulation input data**

iddata object | matrix

Simulation input data, specified as an `iddata` object or a matrix. `sim` uses the input channels from this object as the simulation inputs. For time-domain simulation of discrete-time systems, you can also specify `udata` as a matrix with columns that correspond to each input channel.

If `sys` is a linear model, you can use either time-domain or frequency-domain data. If `sys` is a nonlinear model, you can only use time-domain data.

If `sys` is a time-series model, that is a model with no inputs, specify `udata` as an  $N_s$ -by-0 signal, where  $N_s$  is the wanted number of simulation output samples. For example, to simulate 100 output samples, specify `udata` as follows.

```
udata = iddata([],zeros(100,0),Ts);
```

If you do not have data from an experiment, use `idinput` to generate signals with various characteristics.

## **opt — Simulation options**

`simOptions` option set

Simulation options, specified as a `simOptions` option set for setting the following options:

- Initial conditions
- Input/output offsets
- Additive noise

## **Output Arguments**

### **y — Simulated response**

`iddata` object | matrix

Simulated response for `sys`, returned as an `iddata` object or matrix, depending on how you specify `udata`. For example, if `udata` is an `iddata` object, then so is `y`.

If `udata` represents time-domain data, then `y` is the simulated response for the time vector corresponding to `udata`.

If `udata` represents frequency-domain data,  $U(\omega)$ , then `y` contains the Fourier transform of the corresponding sampled time-domain output signal. This signal is the product of the frequency response of `sys`,  $G(\omega)$ , and  $U(\omega)$ .

For multi-experiment data, `y` is a corresponding multi-experiment `iddata` object.

### **y\_sd — Estimated standard deviation**

double matrix

Estimated standard deviation of the simulated response for linear models or nonlinear grey-box models, returned as an  $N_s$ -by- $N_y$  matrix, where  $N_s$  is the number of samples and  $N_y$  is the number of outputs. The software computes the standard deviation by taking into account the model parameter covariance, initial state covariance, and additive noise covariance. The additive noise covariance is stored in the `NoiseVariance` property of the model.

`y_sd` is derived using first order sensitivity considerations (Gauss approximation formula).

For nonlinear models, `y_sd` is `[]`.

### **x** — Estimated state trajectory for state-space models

matrix | `[]`

Estimated state trajectory for state-space models, returned as an  $N_s$ -by- $N_x$  matrix, where  $N_s$  is the number of samples and  $N_x$  is the number of states.

`x` is only relevant if `sys` is an `idss`, `idgrey`, or `idnlgrey` model. If `sys` is not a state-space model, `x` is returned as `[]`.

### **x\_sd** — Estimated standard deviation of state trajectory

matrix | `[]`

Estimated standard deviation of state trajectory for state-space models, returned as an  $N_s$ -by- $N_x$  matrix, where  $N_s$  is the number of samples and  $N_x$  is the number of states. The software computes the standard deviation by taking into account the model parameter covariance, initial state covariance, and additive noise covariance. The additive noise covariance is stored in the `NoiseVariance` property of the model.

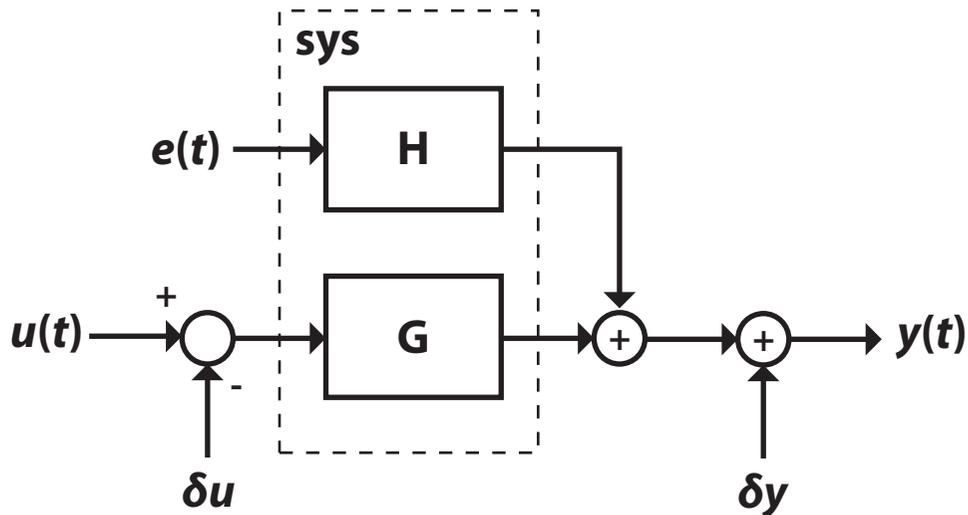
`x_sd` is only relevant if `sys` is an `idss`, `idgrey`, or `idnlgrey` model. If `sys` is not a state-space model, `x_sd` is returned as `[]`.

## Tips

- When the initial conditions of the estimated model and the system that measured the validation data set are different, the simulated and measured responses may also differ, especially at the beginning of the response. To minimize this difference, estimate the initial state values using `findstates` and use the estimated values to set the `InitialCondition` option using `simOptions`. For an example, see “Match Model Response to Output Data” on page 1-1560.

## Algorithms

*Simulation* means computing the model response using input data and initial conditions. `sim` simulates the following system:



Here,

- $u(t)$  is the simulation input data, `udata`.
- $y(t)$  is the simulated output response.
- $G$  is the transfer function from the input to the output and is defined in `sys`. The simulation initial conditions, as specified using `simOptions`, set the initial state of  $G$ .
- $e(t)$  is an optional noise signal. Add noise to your simulation by creating a `simOptions` option set, and setting the `AddNoise` option to `true`. Additionally, you can change the default noise signal by specifying the `NoiseData` option.
- $H$  is the noise transfer function and is defined in `sys`.
- $\delta u$  is an optional input offset subtracted from the input signal,  $u(t)$ , before the input is used to simulate the model. Specify an input offset by setting the `InputOffset` option using `simOptions`.
- $\delta y$  is an optional output offset added to the output response,  $y(t)$ , after simulation. Specify an output offset by setting the `OutputOffset` option using `simOptions`.

For more information on specifying simulation initial conditions, input and output offsets, and noise signal data, see `simOptions`. For multiexperiment data, you can specify these options separately for each experiment.

## Alternatives

- Use `simsd` for a Monte-Carlo method of computing the standard deviation of the response.
- `sim` extends `lsim` to facilitate additional features relevant to identified models:
  - Simulation of nonlinear models
  - Simulation with additive noise
  - Incorporation of signal offsets
  - Computation of response standard deviation (linear models only)
  - Frequency-domain simulation (linear models only)
  - Simulations using different intersample behavior for different inputs

To obtain the simulated response without any of the preceding operations, use `lsim`.

## See Also

`compare` | `findstates` | `forecast` | `idinput` | `lsim` | `predict` | `simOptions` | `simsd` | `step`

## Topics

“Simulate and Predict Identified Model Output”

“Simulation and Prediction at the Command Line”

**Introduced before R2006a**

## simOptions

Option set for sim

### Syntax

```
opt = simOptions  
opt = simOptions(Name,Value)
```

### Description

`opt = simOptions` creates the default option set for `sim`.

`opt = simOptions(Name,Value)` creates an option set with the options specified by one or more `Name,Value` pair arguments.

### Examples

#### Create Default Option Set for Model Simulation

```
opt = simOptions;
```

#### Specify Options for Model Simulation

Create an option set for `sim` specifying the following options.

- Zero initial conditions
- Input offset of 5 for the second input of a two-input model

```
opt = simOptions('InitialCondition','z','InputOffset',[0; 5]);
```

### Add Noise to Simulation Output

Create noise data for a simulation with 500 input data samples and two outputs.

```
noiseData = randn(500,2);
```

Create a default option set.

```
opt = simOptions;
```

Modify the option set to add the noise data.

```
opt.AddNoise = true;  
opt.NoiseData = noiseData;
```

### Use Historical Data to Specify Initial Conditions for Model Simulation

Use historical input-output data as a proxy for initial conditions when simulating your model.

Load a two-input, one-output data set.

```
load iddata7 z7
```

Identify a fifth-order state-space model using the data.

```
sys = n4sid(z7, 5);
```

Split the data set into two parts.

```
zA = z7(1:15);  
zB = z7(16:end);
```

Simulate the model using the input signal in zB.

```
uSim = zB;
```

Simulation requires initial conditions. The signal values in zA are the historical data, that is, they are the input and output values for the time immediately preceding data in zB. Use zA as a proxy for the required initial conditions.

```
I0 = struct('Input',zA.InputData,'Output',zA.OutputData);  
opt = simOptions('InitialCondition',I0);
```

Simulate the model.

```
ysim = sim(sys,uSim,opt);
```

To understand how the past data is mapped to the initial states of the model, see “Understand Use of Historical Data for Model Simulation” on page 1-1566.

## Input Arguments

### Name-Value Pair Arguments

Specify optional comma-separated pairs of `Name`, `Value` arguments. `Name` is the argument name and `Value` is the corresponding value. `Name` must appear inside quotes. You can specify several name and value pair arguments in any order as `Name1,Value1,...,NameN,ValueN`.

Example: `'AddNoise',true','InputOffset',[5;0]` adds default Gaussian white noise to the response model and specifies an input offset of 5 for the first of two model inputs.

### InitialCondition — Simulation initial conditions

`[]` (default) | column vector | matrix | structure | structure array | `'model'`

Simulation initial conditions, specified as one of the following:

- `'z'` — Zero initial conditions.
- Numerical column vector of initial states with length equal to the model order.

For multi-experiment data, specify a matrix with  $N_e$  columns, where  $N_e$  is the number of experiments, to configure the initial conditions separately for each experiment. Otherwise, use a column vector to specify the same initial conditions for all experiments.

Use this option for state-space models (`idss` and `idgrey`) only.

- Structure with the following fields, which contain the historical input and output values for a time interval immediately before the start time of the data used in the simulation:

Field	Description
Input	Input history, specified as a matrix with $N_u$ columns, where $N_u$ is the number of input channels. For time-series models, use <code>[]</code> . The number of rows must be greater than or equal to the model order.
Output	Output history, specified as a matrix with $N_y$ columns, where $N_y$ is the number of output channels. The number of rows must be greater than or equal to the model order.

For an example, see “Use Historical Data to Specify Initial Conditions for Model Simulation” on page 1-1575.

For multi-experiment data, configure the initial conditions separately for each experiment by specifying `InitialCondition` as a structure array with  $N_e$  elements. To specify the same initial conditions for all experiments, use a single structure.

The software uses `data2state` to map the historical data to states. If your model is not `idss`, `idgrey`, `idnlgrey`, or `idnlrx`, the software first converts the model to its state-space representation and then maps the data to states. If conversion of your model to `idss` is not possible, the estimated states are returned empty.

- `'model'` — Use this option for `idnlgrey` models only. The software sets the initial states to the values specified in the `sys.InitialStates` property of the model `sys`.
- `[]` — Corresponds to zero initial conditions for all models except `idnlgrey`. For `idnlgrey` models, the software treats `[]` as `'model'` and specifies the initial states as `sys.InitialStates`.

### **X0Covariance — Covariance of initial states vector**

`[]` (default) | matrix

Covariance of initial states vector, specified as one of the following:

- Positive definite matrix of size  $N_x$ -by- $N_x$ , where  $N_x$  is the model order.

For multi-experiment data, specify as an  $N_x$ -by- $N_x$ -by- $N_e$  matrix, where  $N_e$  is the number of experiments.

- `[]` — No uncertainty in the initial states.

Use this option only for state-space models (`idss` and `idgrey`) when `'InitialCondition'` is specified as a column vector. Use this option to account for initial condition uncertainty when computing the standard deviation of the simulated response of a model.

## **InputOffset — Input signal offset**

[ ] (default) | column vector | matrix

Input signal offset, specified as a column vector of length  $N_u$ . Use [ ] if there are no input offsets. Each element of `InputOffset` is subtracted from the corresponding input data before the input is used to simulate the model.

For multiexperiment data, specify `InputOffset` as:

- An  $N_u$ -by- $N_e$  matrix to set offsets separately for each experiment.
- A column vector of length  $N_u$  to apply the same offset for all experiments.

## **OutputOffset — Output signal offset**

[ ] (default) | column vector | matrix

Output signal offset, specified as a column vector of length  $N_y$ . Use [ ] if there are no output offsets. Each element of `OutputOffset` is added to the corresponding simulated output response of the model.

For multiexperiment data, specify `OutputOffset` as:

- An  $N_y$ -by- $N_e$  matrix to set offsets separately for each experiment.
- A column vector of length  $N_y$  to apply the same offset for all experiments.

## **AddNoise — Noise addition toggle**

false (default) | true

Noise addition toggle, specified as a logical value indicating whether to add noise to the response model.

## **NoiseData — Noise signal data**

[ ] (default) | matrix | cell array of matrices

Noise signal data specified as one of the following:

- [ ] — Default Gaussian white noise.
- Matrix with  $N_s$  rows and  $N_y$  columns, where  $N_s$  is the number of input data samples, and  $N_y$  is the number of outputs. Each matrix entry is scaled according to `NoiseVariance` property of the simulated model and added to the corresponding output data point. To set `NoiseData` at a level that is consistent with the model, use white noise with zero mean and a unit covariance matrix.

- Cell array of  $N_e$  matrices, where  $N_e$  is the number of experiments for multiexperiment data. Use a cell array to set the `NoiseData` separately for each experiment, otherwise set the same noise signal for all experiments using a matrix.

`NoiseData` is the noise signal,  $e(t)$ , for the model

$$y(t) = Gu(t) + He(t).$$

Here,  $G$  is the transfer function from the input,  $u(t)$ , to the output,  $y(t)$ , and  $H$  is the noise transfer function.

`NoiseData` is used for simulation only when `AddNoise` is true.

## Output Arguments

**opt** — Option set for `sim` command

`simOptions` option set

Option set for `sim` command, returned as a `simOptions` option set.

## See Also

`sim`

**Introduced in R2012a**

## **simsd**

Simulate linear models with uncertainty using Monte Carlo method

`simsd` simulates linear models using the Monte Carlo method. The command performs multiple simulations using different values of the uncertain parameters of the model, and different realizations of additive noise and simulation initial conditions. `simsd` uses Monte Carlo techniques to generate response uncertainty, whereas `sim` generates the uncertainty using the Gauss Approximation Formula.

## **Syntax**

```
simsd(sys, udata)
simsd(sys, udata, N)
simsd(sys, udata, N, opt)
```

```
y = simsd(____)
[y, y_sd] = simsd(____)
```

## **Description**

`simsd(sys, udata)` simulates and plots the response of 10 perturbed realizations of the identified model `sys`. Simulation input data `udata` is used to compute the simulated response.

The parameters of the perturbed realizations of `sys` are consistent with the parameter covariance of the original model, `sys`. If `sys` does not contain parameter covariance information, the 10 simulated responses are identical. For information about how the parameter covariance information is used to generate the perturbed models, see “Generating Perturbations of Identified Model” on page 1-1591.

`simsd(sys, udata, N)` simulates and plots the response of `N` perturbed realizations of the identified model `sys`.

`simsd(sys, udata, N, opt)` simulates the system response using the simulation behavior specified in the option set `opt`. Use `opt` to specify uncertainties in the initial conditions and include the effect of additive disturbances.

The simulated responses are all identical if `sys` does not contain parameter covariance information, and you do not specify additive noise or covariance values for initial states. You specify these values in the `AddNoise` and `X0Covariance` options of `opt`.

`y = simsd( ___ )` returns the `N` simulation results in `y` as a cell array. No simulated response plot is produced. Use with any of the input argument combinations in the previous syntaxes.

`[y,y_sd] = simsd( ___ )` also returns the estimated standard deviation `y_sd` for the simulated response.

## Examples

### Simulate State-Space Model Using Monte Carlo Method

Load the estimation data.

```
load iddata1 z1
```

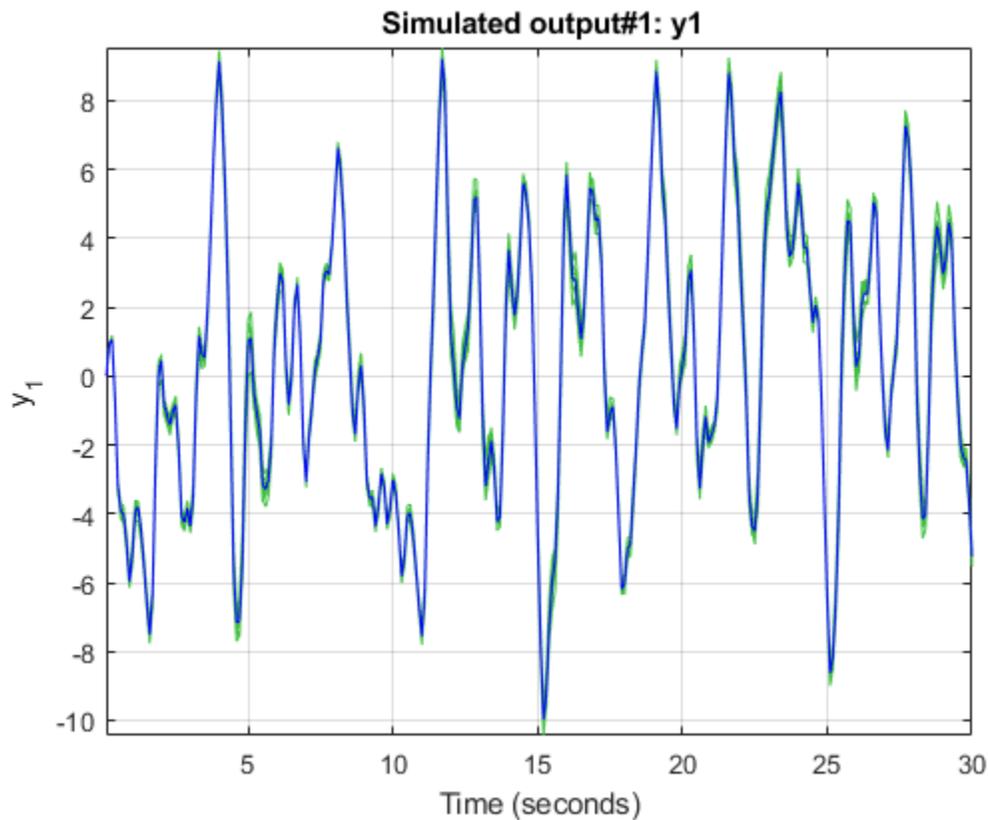
`z1` is an `iddata` object that stores the input-output estimation data.

Estimate a third-order state-space model.

```
sys = ssest(z1,3);
```

Simulate the response of the estimated model using the Monte Carlo method and input estimation data, and plot the response.

```
simsd(sys,z1);
```



The blue line plots the simulated response of the original nominal model  $\text{sys}$ . The green lines plot the simulated response of 10 perturbed realizations of  $\text{sys}$ .

### **Simulate Estimated Model Using Monte Carlo Method**

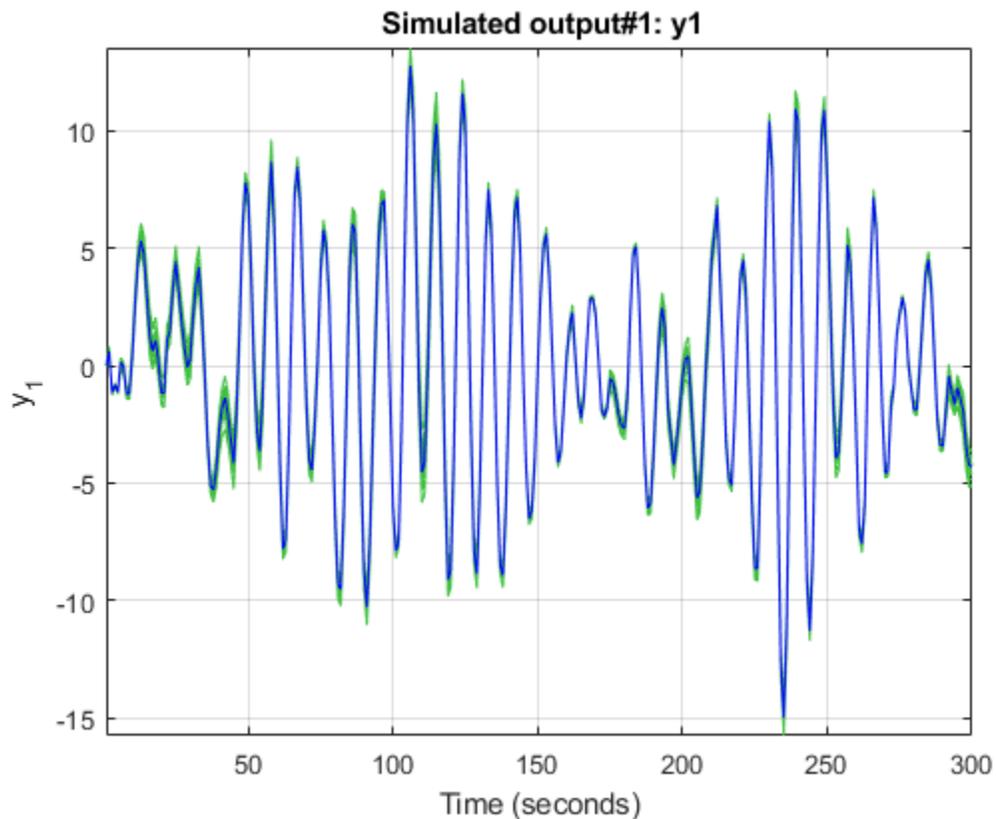
Simulate an estimated model using the Monte Carlo method for a specified number of model perturbations.

Estimate a second-order state-space model using estimation data. Obtain  $\text{sys}$  in the observability canonical form.

```
load iddata3 z3
sys = ssest(z3,2,'Form','canonical');
```

Compute the simulated response of the estimated model using the Monte Carlo method, and plot the responses. Specify the number of random model perturbations as 20.

```
N = 20;
simsd(sys,z3,N)
```



The blue line plots the simulated response of the original nominal model `sys`. The green lines plot the simulated response of the 20 perturbed realizations of `sys`.

You can also obtain the simulated response for each perturbation of `sys`. No plot is generated when you use this syntax.

```
y = simsd(sys,z3,N);
```

y is the simulated response, returned as a cell array of N+1 elements. y{1} contains the nominal response for sys. The remaining elements contain the simulated response for the N perturbed realizations.

## Simulate Time Series Model Using Monte Carlo Method

Load time series data.

```
load iddata9 z9
```

z9 is an iddata object with 200 output data samples and no inputs.

Estimate a sixth-order AR model using the least-squares algorithm.

```
sys = ar(z9,6,'ls');
```

For time series data, specify the desired simulation length, Ns = 200 using an Ns-by-0 input data set.

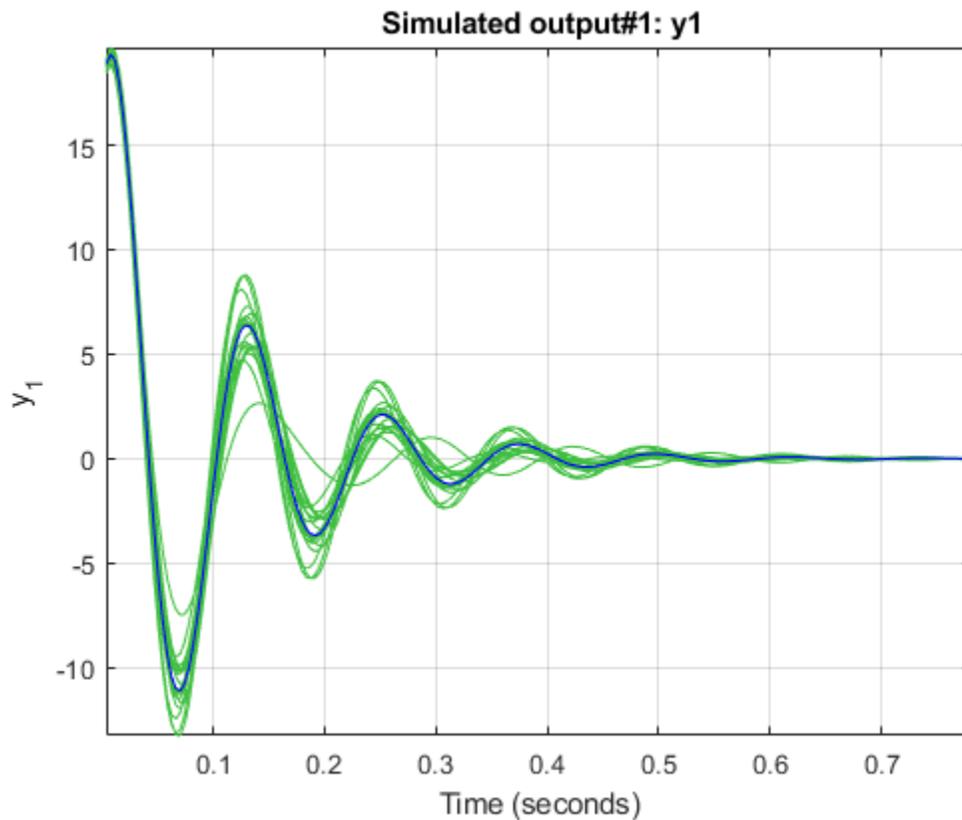
```
data = iddata([],zeros(200,0),z9.Ts);
```

Set the initial conditions to use the initial samples of the time series as historical output samples. The past data is mapped to the initial states of each perturbed system individually.

```
IC = struct('Input',[],'Output',z9.y(1:6));  
opt = simsdOptions('InitialCondition',IC);
```

Simulate the model using Monte Carlo method and specified initial conditions. Specify the number of random model perturbations as 20.

```
simsd(sys,data,20,opt)
```



The blue line plots the simulated response of the original nominal model `sys`. The green lines plot the simulated response of the 20 perturbed realizations of `sys`.

### Study Effect of Initial Condition Uncertainty on Model Response

Load data, and split it into estimation and simulation data.

```
load iddata3
ze = z3(1:200);
zs = z3(201:256);
```

Estimate a second-order state-space model `sys` using estimation data. Specify that no parameter covariance data is generated. Obtain `sys` in the observability canonical form.

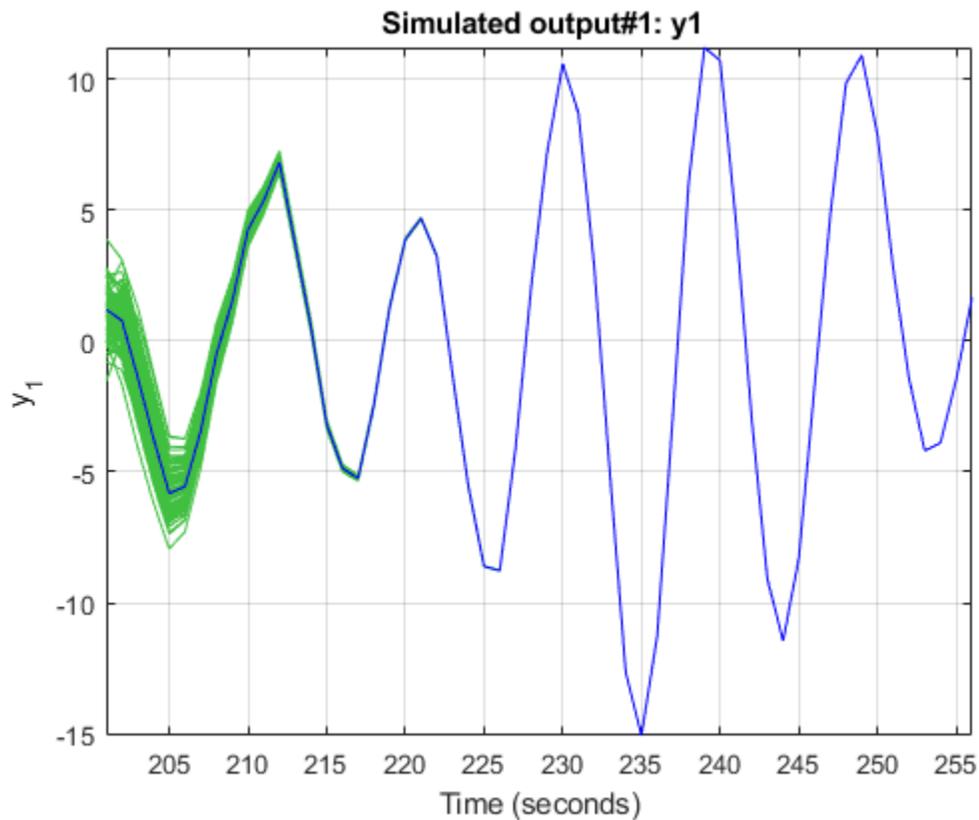
```
opt = ssestOptions('EstimateCovariance',false);  
sys = ssest(ze,2,'Form','canonical',opt);
```

Set the initial conditions for simulating the estimated model. Specify initial state values `x0` for the two states and also the covariance of initial state values `x0Cov`. The covariance is specified as a 2-by-2 matrix because there are two states.

```
x0 = [1.2; -2.4];  
x0Cov = [0.86 -0.39; -0.39 1.42];  
opt = simsdOptions('InitialCondition',x0,'X0Covariance',x0Cov);
```

Simulate the model using Monte Carlo method and specified initial conditions. Specify the number of random model perturbations as 100.

```
simsd(sys,zsim,100,opt)
```



The blue line plots the simulated response of the original nominal model `sys`. The green lines plot the simulated response of the 100 perturbed realizations of `sys`. The software uses a different realization of the initial states to simulate each perturbed model. Initial states are drawn from a Gaussian distribution with mean `InitialCondition` and covariance `X0Covariance`.

### Study Effect of Additive Disturbance on Response Uncertainty

Load the estimation data.

```
load iddata1 z1
```

`z1` is an `iddata` object that stores 300 input-output estimation data samples.

Estimate a second-order state-space model using the estimation data.

```
sys = ssest(z1,2);
```

Create a default option set for `simsd`, and modify the option set to add noise.

```
opt = simsdOptions;  
opt.AddNoise = true;
```

Compute the simulated response of the estimated model using the Monte Carlo method. Specify the number of random model perturbations as 20, and simulate the model using the specified option set.

```
[y,y_sd] = simsd(sys,z1,20,opt);
```

`y` is the simulated response, returned as a cell array of 21 elements. `y{1}` contains the nominal, noise-free response for `sys`. The remaining elements contain the simulated response for the 20 perturbed realizations of `sys` with additive disturbances added to each response.

`y_sd` is the estimated standard deviation of simulated response, returned as an `iddata` object with no inputs. The standard deviations are computed from the 21 simulated outputs. To access the standard deviation, use `y_sd.OutputData`.

## Input Arguments

### **sys** — Model to be simulated

parametric linear identified model

Model to be simulated, specified as one of the following parametric linear identified models: `idtf`, `idproc`, `idpoly`, `idss`, or `idgrey`.

To generate the set of simulated responses, the software perturbs the parameters of `sys` in a way that is consistent with the parameter covariance information. Use `getcov` to examine the parameter uncertainty for `sys`. For information about how the perturbed models are generated from `sys`, see `rsample`.

The simulated responses are all identical if `sys` does not contain parameter covariance information and you do not specify additive noise or covariance values for initial states. You specify these values in the `AddNoise` and `X0Covariance` options of `opt`.

### **udata — Simulation input data**

`iddata` object | matrix

Simulation input data, specified as one of the following:

- `iddata` object — Input data can be either time-domain or frequency-domain. The software uses only the input channels of the `iddata` object.

If `sys` is a time series model, that is, a model with no inputs, specify `udata` as an  $N_s$ -by-0 signal, where  $N_s$  is the wanted number of simulation output samples for each of the  $N$  perturbed realizations of `sys`. For example, to simulate 100 output samples, specify `udata` as follows.

```
udata = iddata([],zeros(100,0),Ts);
```

For an example, see “Simulate Time Series Model Using Monte Carlo Method” on page 1-1584.

- matrix — For simulation of discrete-time systems using time-domain data only. Columns of the matrix correspond to each input channel.

If you do not have data from an experiment, use `idinput` to generate signals with various characteristics.

### **N — Number of perturbed realizations**

10 (default) | positive integer

Number of perturbed realizations of `sys` to be simulated, specified as a positive integer.

### **opt — Simulation options**

`simsdOptions` option set

Simulation options for simulating models using Monte Carlo methods, specified as a `simsdOptions` option set. You can use this option set to specify:

- Input and output signal offsets — Specify an offset to remove from the input signal and an offset to add to the response of `sys`.
- Initial condition handling — Specify initial conditions for simulation and their covariance. For state-space and linear grey-box models (`idss` and `idgrey`), if you

want to simulate the effect of uncertainty in initial states, set the `InitialCondition` option to a double vector, and specify its covariance using the `X0Covariance` option. For an example, see “Study Effect of Initial Condition Uncertainty on Model Response” on page 1-1585.

- Addition of noise to simulated data — If you want to include the influence of additive disturbances, specify the `AddNoise` option as `true`. For an example, see “Study Effect of Additive Disturbance on Response Uncertainty” on page 1-1587.

## Output Arguments

### **y** — Simulated response

cell array

Simulated response, returned as a cell array of  $N+1$  elements. `y{1}` contains the nominal response for `sys`. The remaining elements contain the simulated response for the  $N$  perturbed realizations.

The command performs multiple simulations using different values of the uncertain parameters of the model, and different realizations of additive noise and simulation initial conditions. Thus, the simulated responses are all identical if `sys` does not contain parameter covariance information and you do not specify additive noise and covariance values for initial states in `opt`.

### **y\_sd** — Estimated standard deviation of simulated response

`iddata` object

Estimated standard deviation of simulated response, returned as an `iddata` object. The standard deviation is computed as the sample standard deviation of the `y` ensemble:

$$y\_sd = \sqrt{\frac{1}{N} \sum_{i=2}^{N+1} (y\{1\} - y\{i\})^2}$$

Here `y{1}` is the nominal response for `sys`, and `y{i}` ( $i = 2:N+1$ ) are the simulated responses for the  $N$  perturbed realizations of `sys`.

## Definitions

### Generating Perturbations of Identified Model

The software generates  $N$  perturbations of the identified model  $\text{sys}$  and then simulates the response of each of these perturbations. The parameters of the perturbed realizations of  $\text{sys}$  are consistent with the parameter covariance of the original model  $\text{sys}$ . The parameter covariance of  $\text{sys}$  gives information about the distribution of the parameters. However, for some parameter values, the resulting perturbed systems can be unstable. To reduce the probability of generation of unrealistic systems, the software prescales the parameter covariance.

If  $\Delta p$  is the parameter covariance for the parameters  $p$  of  $\text{sys}$ , then the simulated output  $f(p+\Delta p)$  of a perturbed model as a first-order approximation is:

$$f(p + \Delta p) = f(p) + \frac{\partial f}{\partial p} \Delta p$$

The `simsd` command first scales  $\Delta p$  by a scaling factor  $s$  (approximately 0.1%) to generate perturbed systems with parameters  $(p+s\Delta p)$ . The command then computes  $f(p+s\Delta p)$ , the simulated response of these perturbed systems. Where,

$$f(p + s\Delta p) = f(p) + s \frac{\partial f}{\partial p} \Delta p$$

The command then computes the simulated response  $f(p+\Delta p)$  as:

$$f(p + \Delta p) = f(p) + \frac{1}{s}(f(p + s\Delta p) - f(p))$$

---

**Note** This scaling is not applied to the free delays of `idproc` or `idtf` models.

---

If you specify the `AddNoise` option of `simsdOptions` as `true`, the software adds different realizations of the noise sequence to the noise-free responses of the perturbed system. The realizations of the noise sequence are consistent with the noise component of the model.

For state-space models, if you specify the covariance of initial state values in `X0Covariance` option of `simsdOptions`, different realizations of the initial states are used to simulate each perturbed model. Initial states are drawn from a Gaussian distribution with mean `InitialCondition` and covariance `X0Covariance`.

### **See Also**

`getcov` | `rsample` | `showConfidence` | `sim` | `simsdOptions`

**Introduced before R2006a**

# simsdOptions

Option set for `simsd`

## Syntax

```
opt = simsdOptions  
opt = simsdOptions(Name,Value)
```

## Description

`opt = simsdOptions` creates the default option set for `simsd`.

`opt = simsdOptions(Name,Value)` creates an option set with the options specified by one or more `Name,Value` pair arguments.

## Examples

### Create Default Option Set for Uncertain Model Simulation

```
opt = simsdOptions;
```

### Specify Options for Uncertain Model Simulation

Create an option set for `simsd` specifying the following options.

- Zero initial conditions
- Input offset of 5 for the second input of a two-input model

```
opt = simsdOptions('InitialCondition','z','InputOffset',[0; 5]);
```

## Add Noise to Uncertain Simulation Output

Create a default option set.

```
opt = simsdOptions;
```

Modify the option set to add noise to the data.

```
opt.AddNoise = true;
```

When you use this option set and `simsd` command to simulate the response of a model `sys`. The command returns the perturbed realizations of `sys` with additive disturbances added to each response.

## Input Arguments

### Name-Value Pair Arguments

Specify optional comma-separated pairs of `Name`, `Value` arguments. `Name` is the argument name and `Value` is the corresponding value. `Name` must appear inside quotes. You can specify several name and value pair arguments in any order as `Name1,Value1,...,NameN,ValueN`.

Example: `opt = simsdOptions('AddNoise',true,'InputOffset',[5;0])` adds default Gaussian white noise to the response model, and specifies an input offset of 5 for the first of two model inputs.

### InitialCondition – Simulation initial conditions

'z' (default) | column vector | matrix | structure | structure array

Simulation initial conditions, specified as one of the following:

- 'z' — Zero initial conditions.
- Numerical column vector  $X_0$  of initial states with length equal to the model order.

For multi-experiment data, specify a matrix with  $N_e$  columns, where  $N_e$  is the number of experiments, to configure the initial conditions separately for each experiment. Otherwise, use a column vector to specify the same initial conditions for all experiments.

Use this option for state-space models (`idss` and `idgrey`) only. You can also specify the covariance of the initial state vector in `X0Covariance`.

- Structure with the following fields, which contain the historical input and output values for a time interval immediately before the start time of the data used in the simulation:

Field	Description
Input	Input history, specified as a matrix with $Nu$ columns, where $Nu$ is the number of input channels. For time-series models, use <code>[]</code> . The number of rows must be greater than or equal to the model order.
Output	Output history, specified as a matrix with $Ny$ columns, where $Ny$ is the number of output channels. The number of rows must be greater than or equal to the model order.

For multi-experiment data, you can configure the initial conditions separately for each experiment by specifying `InitialCondition` as a structure array with  $Ne$  elements. Otherwise, use a single structure to specify the same initial conditions for all experiments.

The software uses `data2state` to map the historical data to states. If your model is not `idss` or `idgrey`, the software first converts the model to its state-space representation and then maps the data to states. If conversion of your model to `idss` is not possible, the estimated states are returned empty.

### **X0Covariance — Covariance of initial states vector**

`[]` (default) | matrix

Covariance of initial states vector, specified as one of the following:

- Positive definite matrix of size  $Nx$ -by- $Nx$ , where  $Nx$  is the model order.

For multi-experiment data, specify as an  $Nx$ -by- $Nx$ -by- $Ne$  matrix, where  $Ne$  is the number of experiments. For the  $k^{th}$  experiment, `X0Covariance(:, :, k)` specifies the covariance of initial states `X0(:, k)`.

- `[]` — No uncertainty in the initial states.

Use this option for state-space models (`idss` and `idgrey`) when `'InitialCondition'` is specified as a numerical column vector `X0`. When you specify this option, the software uses a different realization of the initial states to simulate each perturbed model. Initial

states are drawn from a Gaussian distribution with mean `InitialCondition` and covariance `X0Covariance`.

### **InputOffset — Input signal offset**

`[]` (default) | column vector | matrix

Input signal offset, specified as a column vector of length  $Nu$ . Use `[]` if there are no input offsets. Each element of `InputOffset` is subtracted from the corresponding input data before the input is used to simulate the model.

For multiexperiment data, specify `InputOffset` as:

- An  $Nu$ -by- $Ne$  matrix to set offsets separately for each experiment.
- A column vector of length  $Nu$  to apply the same offset for all experiments.

### **OutputOffset — Output signal offset**

`[]` (default) | column vector | matrix

Output signal offset, specified as a column vector of length  $Ny$ . Use `[]` if there are no output offsets. Each element of `OutputOffset` is added to the corresponding simulated output response of the model.

For multiexperiment data, specify `OutputOffset` as:

- An  $Ny$ -by- $Ne$  matrix to set offsets separately for each experiment.
- A column vector of length  $Ny$  to apply the same offset for all experiments.

### **AddNoise — Noise addition toggle**

`false` (default) | `true`

Noise addition toggle, specified as a logical value indicating whether to add noise to the response model. Set `NoiseModel` to `true` to study the effect of additive disturbances on the response. A different realization of the noise sequence, consistent with the noise component of the perturbed system, is added to the noise-free response of that system.

## **Output Arguments**

### **opt — Option set for `simsd` command**

`simsdOptions` option set

Option set for `simsd` command, returned as a `simsdOptions` option set.

## **See Also**

simsd

**Introduced in R2012a**

## size

Query output/input/array dimensions of input-output model and number of frequencies of FRD model

### Syntax

```
size(sys)
d = size(sys)
Ny = size(sys,1)
Nu = size(sys,2)
Sk = size(sys,2+k)
Nf = size(sys,'frequency')
```

### Description

When invoked without output arguments, `size(sys)` returns a description of type and the input-output dimensions of `sys`. If `sys` is a model array, the array size is also described. For identified models, the number of free parameters is also displayed. The lengths of the array dimensions are also included in the response to `size` when `sys` is a model array.

`d = size(sys)` returns:

- The row vector `d = [Ny Nu]` for a single dynamic model `sys` with `Ny` outputs and `Nu` inputs
- The row vector `d = [Ny Nu S1 S2 ... Sp]` for an `S1-by-S2-by-...-by-Sp` array of dynamic models with `Ny` outputs and `Nu` inputs

`Ny = size(sys,1)` returns the number of outputs of `sys`.

`Nu = size(sys,2)` returns the number of inputs of `sys`.

`Sk = size(sys,2+k)` returns the length of the `k`-th array dimension when `sys` is a model array.

`Nf = size(sys, 'frequency')` returns the number of frequencies when `sys` is a frequency response data model. This is the same as the length of `sys.frequency`.

## Examples

### Query Dimensions of Model Array

Create a 3-by-1 model array of random state-space models with 3 outputs, 2 inputs, and 5 states.

```
sys = rss(5,3,2,3);
```

Verify the size of the model array.

```
size(sys)
```

```
3x1 array of state-space models.
```

```
Each model has 3 outputs, 2 inputs, and 5 states.
```

### Query Dimensions of Identified Model

Create a 2-input 2-output continuous-time process model with identifiable parameters.

```
type = {'p1d', 'p2'; 'p3uz', 'p0'};  
sys = idproc(type);
```

Each element of the `type` cell array describes the model structure for the corresponding input-output pair.

Query the input-output dimensions and number of free parameters in the model.

```
size(sys)
```

```
Process model with 2 outputs, 2 inputs and 12 free parameters.
```

## See Also

`isempty` | `issiso` | `ndims` | `nparams`

**Introduced before R2006a**

## spa

Estimate frequency response with fixed frequency resolution using spectral analysis

## Syntax

```
G = spa(data)
G = spa(data,winSize,freq)
G = spa(data,winSize,freq,MaxSize)
```

## Description

`G = spa(data)` estimates frequency response (with uncertainty) and noise spectrum from time- or frequency-domain data. `data` is an `iddata` or `idfrd` object and can be complex valued. `G` is as an `idfrd` object. For time-series `data`, `G` is the estimated spectrum and standard deviation.

Information about the estimation results and options used is stored in the model's `Report` property. `Report` has the following fields:

- `Status` — Summary of the model status, which indicates whether the model was created by construction or obtained by estimation.
- `Method` — Estimation command used.
- `WindowSize` — Size of the Hann window.
- `DataUsed` — Attributes of the data used for estimation. Structure with the following fields:
  - `Name` — Name of the data set.
  - `Type` — Data type.
  - `Length` — Number of data samples.
  - `Ts` — Sample time.
  - `InterSample` — Input intersample behavior.
  - `InputOffset` — Offset removed from time-domain input data during estimation.

- `OutputOffset` — Offset removed from time-domain output data during estimation.

`G = spa(data,winSize,freq)` estimates frequency response at frequencies `freq`. `freq` is a row vector of values in `rad/TimeUnit`, where `TimeUnit` refers to the `TimeUnit` property of `data`. For discrete-time data, set `freq` within the Nyquist frequency bound. `winSize` is a scalar integer that sets the size of the Hann window.

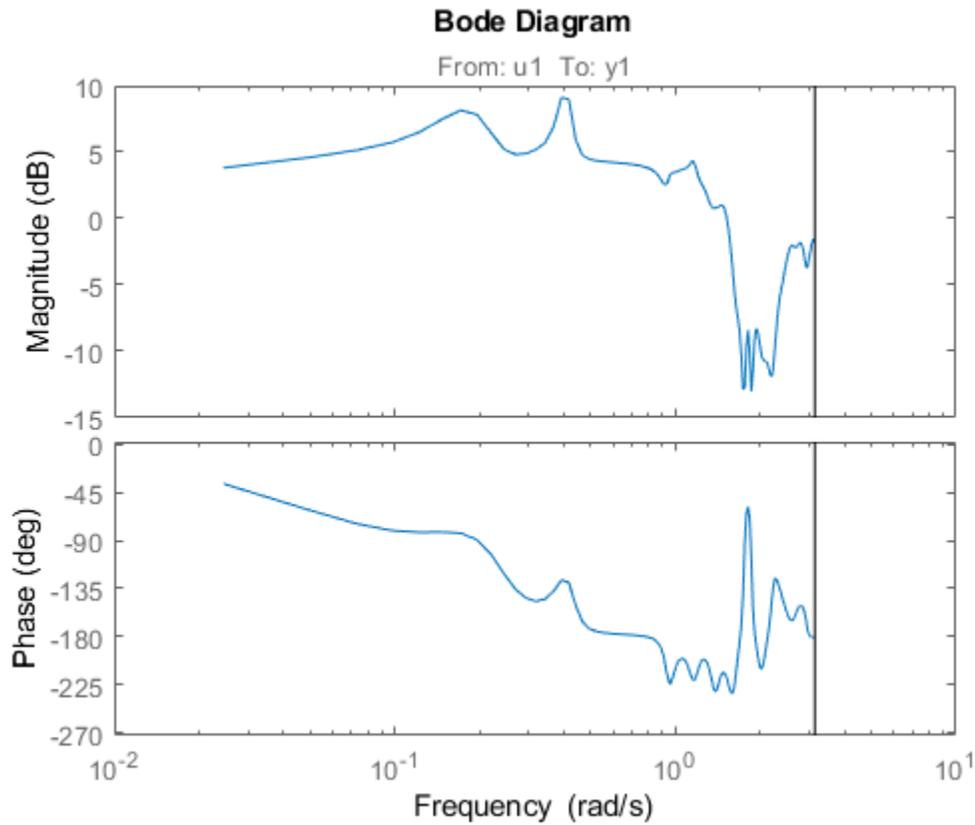
`G = spa(data,winSize,freq,MaxSize)` can improve computational performance using `MaxSize` to split the input-output data such that each segment contains fewer than `MaxSize` elements. `MaxSize` is a positive integer.

## Examples

### Estimate Frequency Response

Estimate frequency response with fixed resolution at 128 equally spaced, logarithmic frequency values between 0 (excluded) and  $\pi$ .

```
load iddata3;  
g = spa(z3);  
bode(g)
```



### Estimate Frequency Response at Specified Frequencies

Define the frequency vector.

```
w = logspace(-2,pi,128);
```

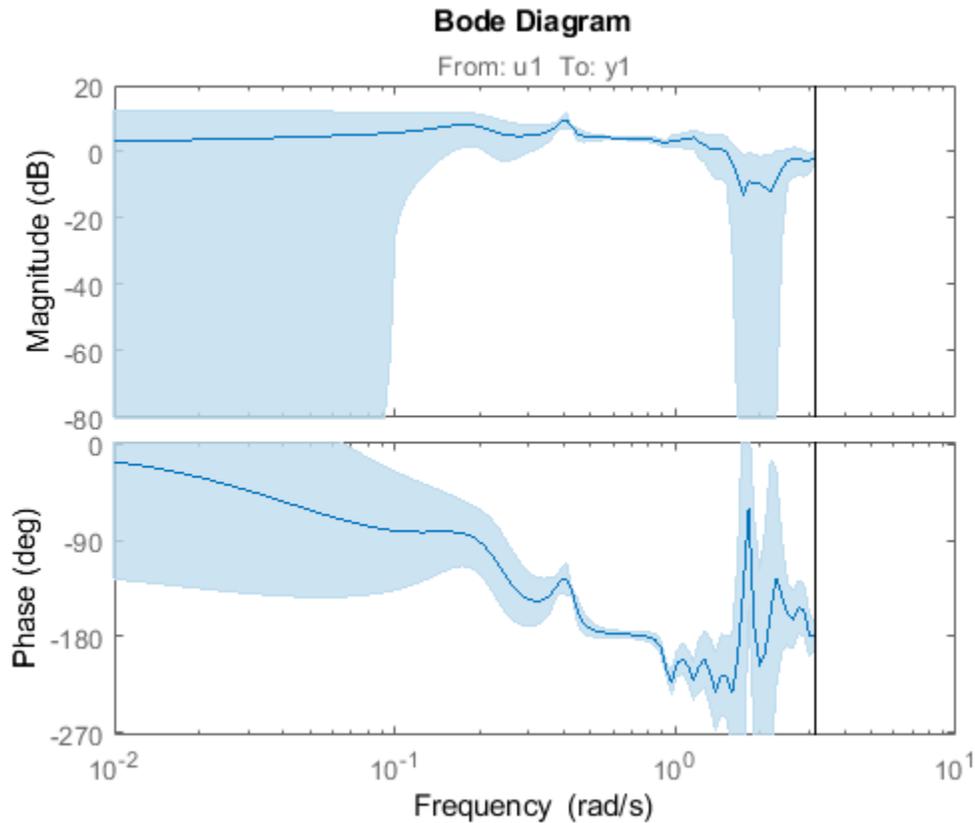
Compute the frequency response.

```
load iddata3;  
g = spa(z3,[],w);
```

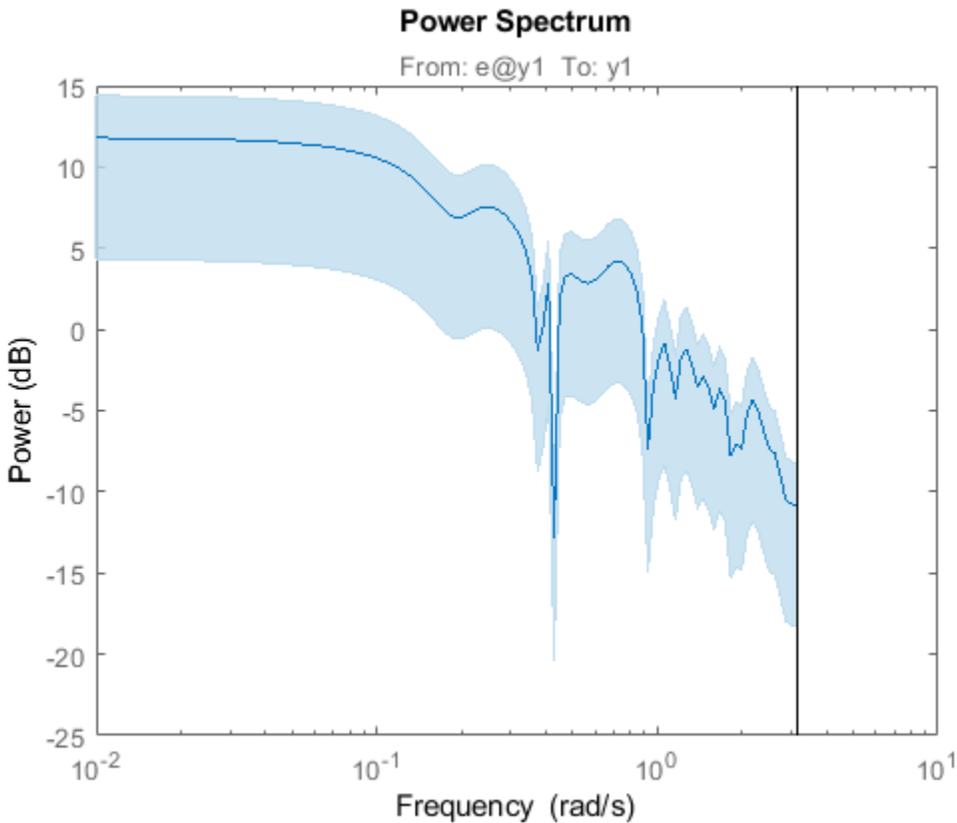
[ ] specifies the default lag window size.

Plot the Bode response and disturbance spectrum with confidence interval of 3 standard deviations.

```
h = bodeplot(g);  
showConfidence(h,3)
```



```
figure  
h = spectrumplot(g);  
showConfidence(h,3)
```



## Definitions

### Frequency Response Function

*Frequency response function* describes the steady-state response of a system to sinusoidal inputs. For a linear system, a sinusoidal input of a specific frequency results in an output that is also a sinusoid with the same frequency, but with a different amplitude and phase. The frequency response function describes the amplitude change and phase shift as a function of frequency.

To better understand the frequency response function, consider the following description of a linear, dynamic system:

$$y(t) = G(q)u(t) + v(t)$$

where  $u(t)$  and  $y(t)$  are the input and output signals, respectively.  $G(q)$  is called the transfer function of the system—it captures the system dynamics that take the input to the output. The notation  $G(q)u(t)$  represents the following operation:

$$G(q)u(t) = \sum_{k=1}^{\infty} g(k)u(t-k)$$

$q$  is the *shift operator*, defined by the following equation:

$$G(q) = \sum_{k=1}^{\infty} g(k)q^{-k} \quad q^{-1}u(t) = u(t-1)$$

$G(q)$  is the *frequency-response function*, which is evaluated on the unit circle,  $G(q=e^{i\omega})$ .

Together,  $G(q=e^{i\omega})$  and the output noise spectrum  $\hat{\Phi}_v(\omega)$  are the frequency-domain description of the system.

The frequency-response function estimated using the Blackman-Tukey approach is given by the following equation:

$$\hat{G}_N(e^{i\omega}) = \frac{\hat{\Phi}_{yu}(\omega)}{\hat{\Phi}_u(\omega)}$$

In this case,  $\hat{\phantom{x}}$  represents approximate quantities. For a derivation of this equation, see the chapter on nonparametric time- and frequency-domain methods in *System Identification: Theory for the User*, Second Edition, by Lennart Ljung, Prentice Hall PTR, 1999.

## Output Noise Spectrum

The output noise spectrum (spectrum of  $v(t)$ ) is given by the following equation:

$$\hat{\Phi}_v(\omega) = \hat{\Phi}_y(\omega) - \frac{|\hat{\Phi}_{yu}(\omega)|^2}{\hat{\Phi}_u(\omega)}$$

This equation for the noise spectrum is derived by assuming the linear relationship  $y(t) = G(q)u(t) + v(t)$ , that  $u(t)$  is independent of  $v(t)$ , and the following relationships between the spectra:

$$\Phi_y(\omega) = \left| G(e^{i\omega}) \right|^2 \Phi_u(\omega) + \Phi_v(\omega)$$

$$\Phi_{yu}(\omega) = G(e^{i\omega}) \Phi_u(\omega)$$

where the noise spectrum is given by the following equation:

$$\Phi_v(\omega) \equiv \sum_{\tau=-\infty}^{\infty} R_v(\tau) e^{-i\omega\tau}$$

$\hat{\Phi}_{yu}(\omega)$  is the output-input cross-spectrum and  $\hat{\Phi}_u(\omega)$  is the input spectrum.

Alternatively, the disturbance  $v(t)$  can be described as filtered white noise:

$$v(t) = H(q)e(t)$$

where  $e(t)$  is the white noise with variance  $\lambda$  and the noise power spectrum is given by the following equation:

$$\Phi_v(\omega) = \lambda \left| H(e^{i\omega}) \right|^2$$

## Algorithms

spa applies the Blackman-Tukey spectral analysis method by following these steps:

- 1** Computes the covariances and cross-covariance from  $u(t)$  and  $y(t)$ :

$$\hat{R}_y(\tau) = \frac{1}{N} \sum_{t=1}^N y(t+\tau)y(t)$$

$$\hat{R}_u(\tau) = \frac{1}{N} \sum_{t=1}^N u(t+\tau)u(t)$$

$$\hat{R}_{yu}(\tau) = \frac{1}{N} \sum_{t=1}^N y(t+\tau)u(t)$$

- 2** Computes the Fourier transforms of the covariances and the cross-covariance:

$$\hat{\Phi}_y(\omega) = \sum_{\tau=-M}^M \hat{R}_y(\tau)W_M(\tau)e^{-i\omega\tau}$$

$$\hat{\Phi}_u(\omega) = \sum_{\tau=-M}^M \hat{R}_u(\tau)W_M(\tau)e^{-i\omega\tau}$$

$$\hat{\Phi}_{yu}(\omega) = \sum_{\tau=-M}^M \hat{R}_{yu}(\tau)W_M(\tau)e^{-i\omega\tau}$$

where  $W_M(\tau)$  is the Hann window with a width (lag size) of  $M$ . You can specify  $M$  to control the frequency resolution of the estimate, which is approximately equal  $2\pi/M$  rad/sample time.

By default, this operation uses 128 equally spaced frequency values between 0 (excluded) and  $\pi$ , where  $w = [1:128]/128*\pi/T_s$  and  $T_s$  is the sample time of that data set. The default lag size of the Hann window is  $M = \min(\text{length}(\text{data})/10, 30)$ . For default frequencies, uses fast Fourier transforms (FFT)—which is more efficient than for user-defined frequencies.

---

**Note**  $M = \gamma$  is in Table 6.1 of Ljung (1999). Standard deviations are on pages 184 and 188 in Ljung (1999).

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- 3**

Compute the frequency-response function  $\hat{G}_N(e^{i\omega})$  and the output noise spectrum  $\hat{\Phi}_v(\omega)$ .

$$\hat{G}_N(e^{i\omega}) = \frac{\hat{\Phi}_{yu}(\omega)}{\hat{\Phi}_u(\omega)}$$

$$\Phi_v(\omega) \equiv \sum_{\tau=-\infty}^{\infty} R_v(\tau) e^{-i\omega\tau}$$

spectrum is the spectrum matrix for both the output and the input channels. That is, if  $z = [\text{data.OutputData}, \text{data.InputData}]$ , `spectrum` contains as spectrum data the matrix-valued power spectrum of  $z$ .

$$S = \sum_{m=-M}^M E z(t+m) z(t)' W_M(T_s) \exp(-i\omega m)$$

' is a complex-conjugate transpose.

## References

Ljung, L. *System Identification: Theory for the User*, Second Ed., Prentice Hall PTR, 1999.

## See Also

`bode` | `etfe` | `freqresp` | `idfrd` | `spafdr` | `spectrum`

## Topics

“Estimate Impulse-Response Models at the Command Line”  
 “Spectrum Normalization”

**Introduced before R2006a**

## spafdr

Estimate frequency response and spectrum using spectral analysis with frequency-dependent resolution

### Syntax

```
g = spafdr(data)
g = spafdr(data,Resol,w)
```

### Description

`g = spafdr(data)` estimates the input-to-output frequency response  $G(\omega)$  and noise spectrum  $\Phi_v$  of the general linear model

$$y(t) = G(q)u(t) + v(t)$$

where  $\Phi_v(\omega)$  is the spectrum of  $v(t)$ . `data` contains the output-input data as an `iddata` object. The data can be complex valued, and either time or frequency domain. It can also be an `idfrd` object containing frequency-response data. `g` is an `idfrd` object with the

estimate of  $G(e^{i\omega})$  at the frequencies  $\omega$  specified by row vector `w`. `g` also includes information about the spectrum estimate of  $\Phi_v(\omega)$  at the same frequencies. Both results are returned with estimated covariances, included in `g`. The normalization of the spectrum is the same as described in `spa`.

Information about the estimation results and options used is stored in the model's `Report` property. `Report` has the following fields:

- `Status` — Summary of the model status, which indicates whether the model was created by construction or obtained by estimation.
- `Method` — Estimation command used.
- `WindowSize` — Frequency resolution.
- `DataUsed` — Attributes of the data used for estimation. Structure with the following fields:

- Name — Name of the data set.
- Type — Data type.
- Length — Number of data samples.
- Ts — Sample time.
- InterSample — Input intersample behavior.
- InputOffset — Offset removed from time-domain input data during estimation.
- OutputOffset — Offset removed from time-domain output data during estimation.

`g = spafdr(data,Resol,w)` specifies frequencies and frequency resolution.

## Frequencies

The frequency variable `w` is either specified as a row vector of frequencies in rad/TimeUnit, where TimeUnit refers to the TimeUnit property of data, or as a cell array `{wmin,wmax}`. In the latter case the covered frequencies will be 50 logarithmically spaced points from `wmin` to `wmax`. You can change the number of points to `NP` by entering `{wmin,wmax,NP}`.

Omitting `w` or entering it as an empty matrix gives the default value, which is 100 logarithmically spaced frequencies between the smallest and largest frequency in data.

For time-domain data, the default range goes from  $\frac{2\pi}{NT_s}$  to  $\frac{\pi}{T_s}$ , where  $T_s$  is the sample time of data and  $N$  is the number of data points.

## Resolution

The argument `Resol` defines the frequency resolution of the estimates. The resolution (measured in rad/TimeUnit) is the size of the smallest detail in the frequency function and the spectrum that is resolved by the estimate. The resolution is a tradeoff between obtaining estimates with fine, reliable details, and suffering from spurious, random effects: The finer the resolution, the higher the variance in the estimate. `Resol` can be entered as a scalar (measured in rad/TimeUnit), which defines the resolution over the whole frequency interval. It can also be entered as a row vector of the same length as `w`. Then `Resol(k)` is the local, frequency-dependent resolution around frequency `w(k)`.

The default value of `Resol`, obtained by omitting it or entering it as the empty matrix, is `Resol(k) = 2(w(k+1)-w(k))`, adjusted upwards, so that a reasonable estimate is

guaranteed. In all cases, the resolution is returned in the variable `g.Report.WindowSize`.

### Algorithms

If the data is given in the time domain, it is first converted to the frequency domain. Then averages of  $Y(\omega)\text{Conj}(U(\omega))$  and  $U(\omega)\text{Conj}(U(\omega))$  are formed over the frequency ranges  $\omega$ , corresponding to the desired resolution around the frequency in question. The ratio of these averages is then formed for the frequency-function estimate, and corresponding expressions define the noise spectrum estimate.

### See Also

`bode` | `etfe` | `freqresp` | `idfrd` | `nyquist` | `spa` | `spectrum`

**Introduced before R2006a**

# spectrum

Output power spectrum of time series models

## Syntax

```
spectrum(sys)
spectrum(sys,{wmin, wmax})
spectrum(sys,w)
spectrum(sys1,...,sysN,w)
ps = spectrum(sys,w)
[ps,w] = spectrum(sys)
[ps,w,sdps] = spectrum(sys)
```

## Description

`spectrum(sys)` creates an output power spectrum plot of the identified time series model `sys`. The frequency range and number of points are chosen automatically.

`sys` is a time series model, which represents the system:

$$y(t) = He(t)$$

Where,  $e(t)$  is a Gaussian white noise and  $y(t)$  is the observed output.

`spectrum` plots  $\text{abs}(H'H)$ , scaled by the variance of  $e(t)$  and the sample time.

If `sys` is an input-output model, it represents the system:

$$y(t) = Gu(t) + He(t)$$

Where,  $u(t)$  is the measured input,  $e(t)$  is a Gaussian white noise and  $y(t)$  is the observed output.

In this case, `spectrum` plots the spectrum of the disturbance component  $He(t)$ .

`spectrum(sys, {wmin, wmax})` creates a spectrum plot for frequencies ranging from `wmin` to `wmax`.

`spectrum(sys, w)` creates a spectrum plot using the frequencies specified in the vector `w`.

`spectrum(sys1, ..., sysN, w)` creates a spectrum plot of several identified models on a single plot. The `w` argument is optional.

You can specify a color, line style and marker for each model. For example:

```
spectrum(sys1, 'r', sys2, 'y--', sys3, 'gx');
```

`ps = spectrum(sys, w)` returns the power spectrum amplitude of `sys` for the specified frequencies, `w`. No plot is drawn on the screen.

`[ps, w] = spectrum(sys)` returns the frequency vector, `w`, for which the output power spectrum is plotted.

`[ps, w, sdps] = spectrum(sys)` returns the estimated standard deviations of the power spectrum.

For discrete-time models with sample time `Ts`, `spectrum` uses the transformation  $z = \exp(j*w*Ts)$  to map the unit circle to the real frequency axis. The spectrum is only plotted for frequencies smaller than the Nyquist frequency  $\pi/Ts$ , and the default value 1 (time unit) is assumed when `Ts` is unspecified.

## Input Arguments

### **sys**

Identified model.

If `sys` is a time series model, it represents the system:

$$y(t) = He(t)$$

Where,  $e(t)$  is a Gaussian white noise and  $y(t)$  is the observed output.

If `sys` is an input-output model, it represents the system:

$$y(t) = Gu(t) + He(t)$$

Where,  $u(t)$  is the measured input,  $e(t)$  is a Gaussian white noise and  $y(t)$  is the observed output.

### **wmin**

Minimum frequency of the frequency range for which the output power spectrum is plotted.

Specify `wmin` in rad/TimeUnit, where `TimeUnit` is `sys.TimeUnit`.

### **wmax**

Maximum frequency of the frequency range for which the output power spectrum is plotted.

Specify `wmax` in rad/TimeUnit, where `TimeUnit` is `sys.TimeUnit`.

### **w**

Frequencies for which the output power spectrum is plotted.

Specify `w` in rad/TimeUnit, where `TimeUnit` is `sys.TimeUnit`.

### **sys1, ..., sysN**

Identified systems for which the output power spectrum is plotted.

## **Output Arguments**

### **ps**

Power spectrum amplitude.

If `sys` has  $N_y$  outputs, then `ps` is an array of size  $[N_y \ N_y \ \text{length}(w)]$ . Where `ps(:, :, k)` corresponds to the power spectrum for the frequency at `w(k)`.

For amplitude values in dB, type `psdb = 10*log10(ps)`.

## w

Frequency vector for which the output power spectrum is plotted.

## sdps

Estimated standard deviation of the power spectrum.

## Examples

### Plot Noise Spectrum of SISO Linear Identified Model

Load the estimation data.

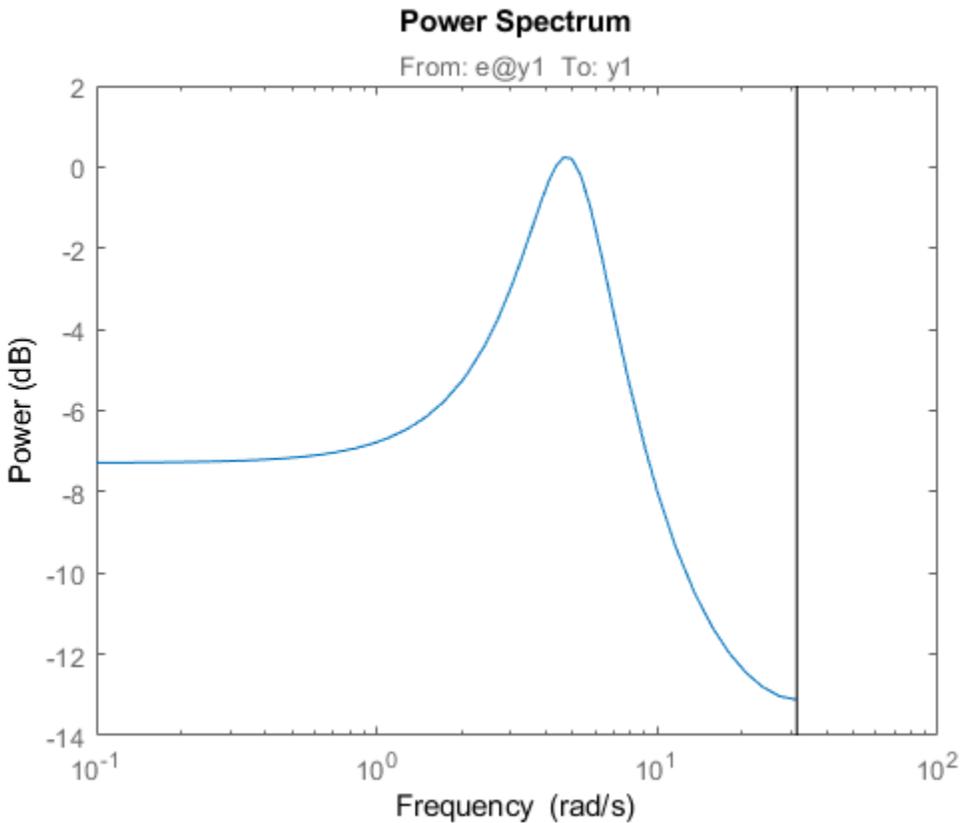
```
load iddata1 z1;
```

Estimate a single-input single-output state-space model.

```
sys = n4sid(z1,2);
```

Plot the noise spectrum for the model.

```
spectrum(sys);
```



### Plot Output Spectrum of Time-Series Model

Load the time-series estimation data.

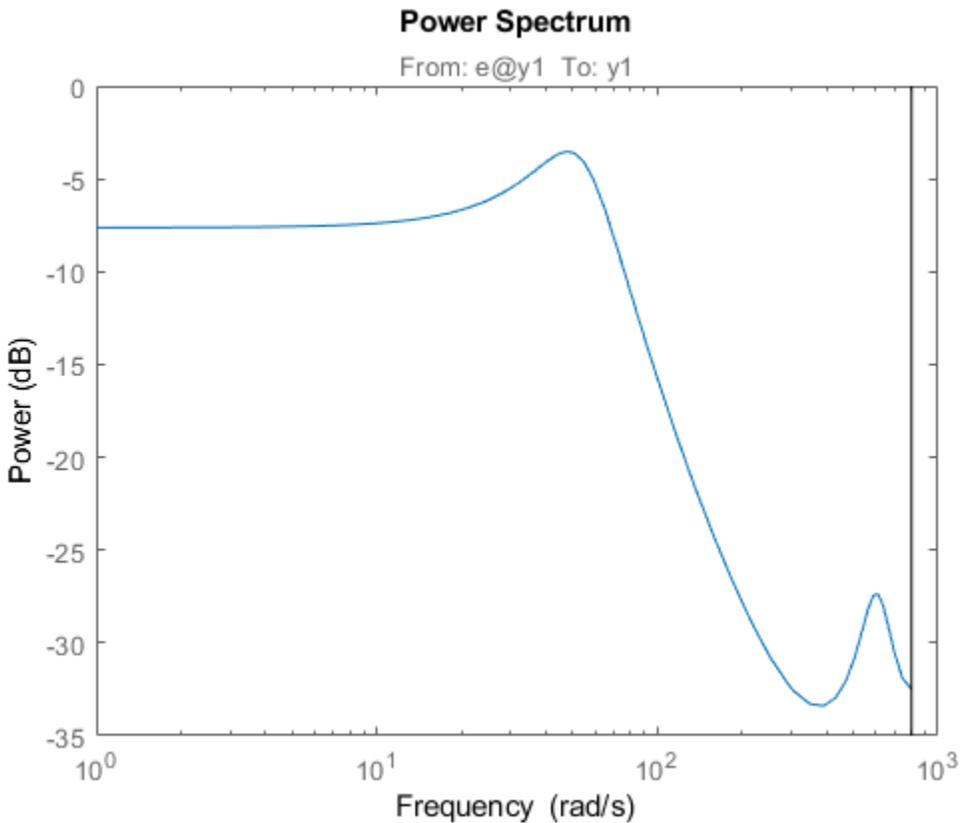
```
load iddata9 z9
```

Estimate a fourth-order AR model using a least-squares approach.

```
sys = ar(z9,4,'ls');
```

Plot the output spectrum of the model.

```
spectrum(sys);
```

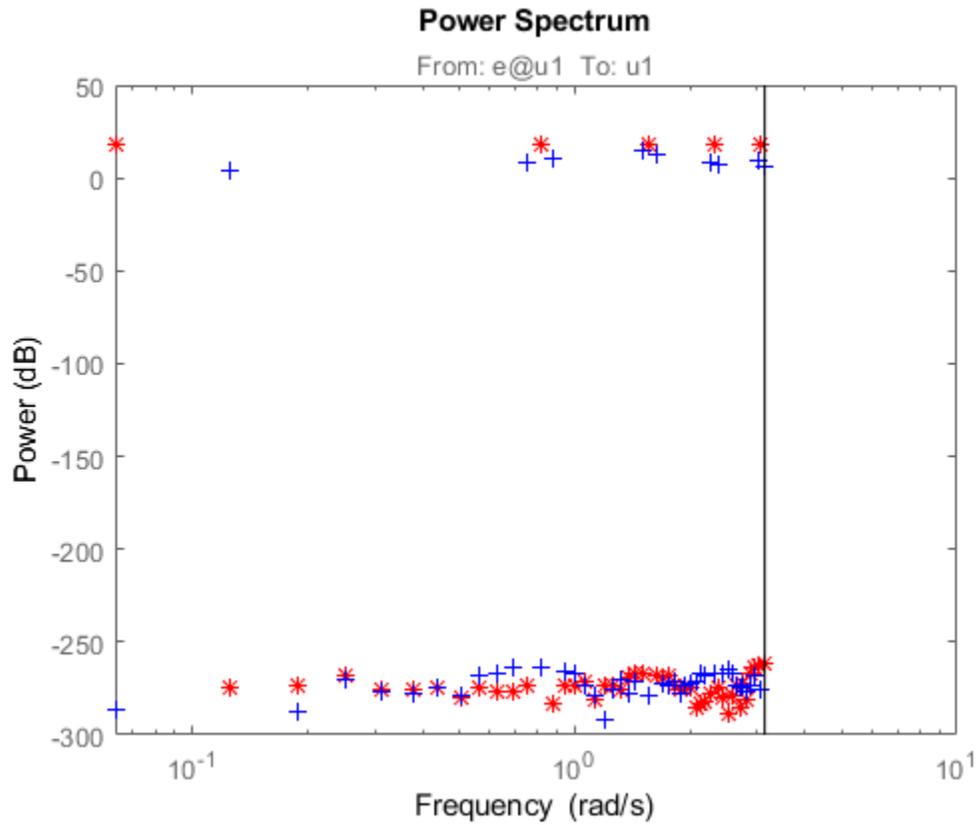


### Compare Spectrum of Sinusoid Signal with Square

Create an input consisting of five sinusoids spread over the whole frequency interval. Compare the spectrum of this signal with that of its square. The frequency splitting (the square having spectral support at other frequencies) reveals the nonlinearity involved.

```
u = idinput([100 1 20], 'sine', [], [], [5 10 1]);  
u = iddata([], u, 1, 'per', 100);  
u2 = u.u.^2;
```

```
u2 = iddata([],u2,1,'per',100);  
spectrum(etfe(u),'r*',etfe(u2),'+')
```



## See Also

[ar](#) | [armax](#) | [arx](#) | [bode](#) | [forecast](#) | [freqresp](#) | [nlarx](#)

**Introduced in R2012a**

## spectrumsoptions

Option set for `spectrumplot`

### Syntax

```
opt = spectrumsoptions  
opt = spectrumsoptions('identpref')
```

### Description

`opt = spectrumsoptions` creates the default option set for `spectrumplot`. Use dot notation to customize the option set, if needed.

`opt = spectrumsoptions('identpref')` initializes the plot options with the System Identification Toolbox preferences. Use this syntax to change a few plot options but otherwise use your toolbox preferences.

### Examples

#### Specify Options for Spectrum Plot

Specify the plot options.

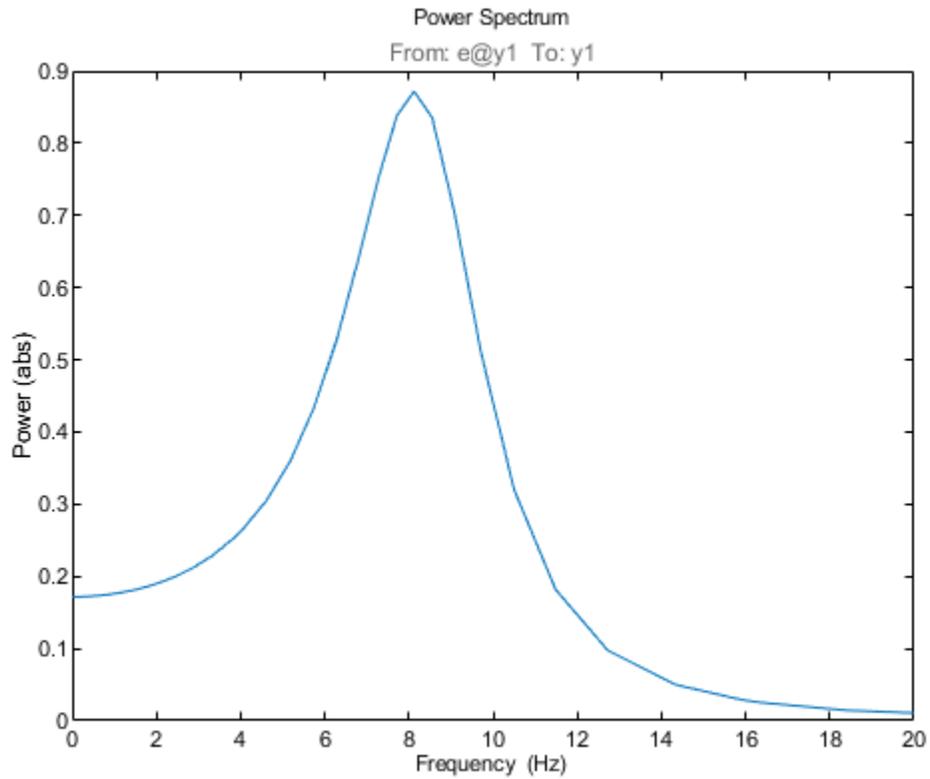
```
plot_options = spectrumsoptions;  
plot_options.FreqUnits = 'Hz';  
plot_options.FreqScale = 'linear';  
plot_options.Xlim = {[0 20]};  
plot_options.MagUnits = 'abs';
```

Estimate an AR model.

```
load iddata9 z9  
sys = ar(z9,4);
```

Plot the output spectrum for the model.

```
spectrumplot(sys,plot_options);
```



### Initialize Plot Options Using Toolbox Preferences

```
opt = spectrumoptions('identpref');
```

## Output Arguments

**opt** — Option set for `spectrumplot`  
`spectrumoptions` option set

Option set containing the specified options for `spectrumplot`.

Field	Description
Title, XLabel, YLabel	<p>Text and style for axes labels and plot title, specified as a structure array with the following fields:</p> <ul style="list-style-type: none"> <li>• <b>String</b> — Title and axes label text, specified as a character vector. <b>Default Title:</b> 'Power Spectrum' <b>Default XLabel:</b> 'Frequency' <b>Default YLabel:</b> 'Power'</li> <li>• <b>FontSize</b> — Font size, specified as data type scalar. <b>Default:</b> 8</li> <li>• <b>FontWeight</b> — Thickness of text, specified as one of the following values: 'Normal'   'Bold' <b>Default:</b> 'Normal'</li> <li>• <b>Font Angle</b> — Text character angle, specified as one of the following values: 'Normal'   'Italic' <b>Default:</b> 'Normal'</li> <li>• <b>Color</b> — Color of text, specified as vector of RGB values between 0 to 1. <b>Default:</b> [0, 0, 0]</li> <li>• <b>Interpreter</b> — Interpretation of text characters, specified as one of the following values: 'tex'   'latex'   'none' <b>Default:</b> 'tex'</li> </ul>

Field	Description
TickLabel	<p>Tick label style, specified as a structure array with the following fields:</p> <ul style="list-style-type: none"> <li>• <b>FontSize</b> — Font size, specified as data type scalar. <b>Default:</b> 8</li> <li>• <b>FontWeight</b> — Thickness of text, specified as one of the following values: 'Normal'   'Bold' <b>Default:</b> 'Normal'</li> <li>• <b>Font Angle</b> — Text character angle, specified as one of the following values: 'Normal'   'Italic' <b>Default:</b> 'Normal'</li> <li>• <b>Color</b> — Color of text, specified as vector of RGB values between 0 to 1   character vector of color name   'none'. For example, for yellow color, specify as one of the following: [1 1 0], 'yellow', or 'y'. <b>Default:</b> [0,0,0]</li> </ul>
Grid	<p>Show or hide the grid, specified as one of the following values: 'off'   'on'</p> <p><b>Default:</b> 'off'</p>
GridColor	<p>Color of the grid lines, specified as one of the following: vector of RGB values in the range [0,1]   character vector of color name   'none'. For example, for yellow color, specify as one of the following: [1 1 0], 'yellow', or 'y'.</p> <p><b>Default:</b> [0.15,0.15,0.15]</p>

Field	Description
XLimMode, YLimMode	<p>Axes limit modes, specified as one of the following values:</p> <ul style="list-style-type: none"><li>• 'auto' — The axes limits are based on the data plotted</li><li>• 'manual' — The values are explicitly set with Xlim, Ylim</li></ul> <p><b>Default:</b> 'auto'</p>
XLim, YLim	<p>Axes limits, specified as an array of the form [min,max]</p>
IOGrouping	<p>Grouping of input-output pairs in the plot, specified as one of the following values: 'none'   'inputs'   'outputs'   'all'</p> <p><b>Default:</b> 'none'</p>

Field	Description
InputLabels, OutputLabels	<p>Input and output label styles, specified as a structure array with the following fields:</p> <ul style="list-style-type: none"> <li>• <b>FontSize</b> — Font size, specified as data type scalar. <b>Default:</b> 8</li> <li>• <b>FontWeight</b> — Thickness of text, specified as one of the following values: 'Normal'   'Bold' <b>Default:</b> 'Normal'</li> <li>• <b>Font Angle</b> — Text character angle, specified as one of the following values: 'Normal'   'Italic' <b>Default:</b> 'Normal'</li> <li>• <b>Color</b> — Color of text, specified as a vector of RGB values between 0 to 1   character vector of color name   'none'. For example, for yellow color, specify as one of the following: [1 1 0], 'yellow', or 'y'. <b>Default:</b> [0.4,0.4,0.4]</li> <li>• <b>Interpreter</b> — Interpretation of text characters, specified as one of the following values: 'tex'   'latex'   'none' <b>Default:</b> 'tex'</li> </ul>
InputVisible, OutputVisible	<p>Visibility of input and output channels, specified as one of the following values: 'off'   'on'</p> <p><b>Default:</b> 'on'</p>
ConfidenceRegionNumberSD	<p>Number of standard deviations to use to plot the response confidence region.</p> <p><b>Default:</b> 1</p>

Field	Description
FreqUnits	<p>Frequency units, specified as one of the following values:</p> <ul style="list-style-type: none"> <li>• 'Hz'</li> <li>• 'rad/second'</li> <li>• 'rpm'</li> <li>• 'kHz'</li> <li>• 'MHz'</li> <li>• 'GHz'</li> <li>• 'rad/nanosecond'</li> <li>• 'rad/microsecond'</li> <li>• 'rad/millisecond'</li> <li>• 'rad/minute'</li> <li>• 'rad/hour'</li> <li>• 'rad/day'</li> <li>• 'rad/week'</li> <li>• 'rad/month'</li> <li>• 'rad/year'</li> <li>• 'cycles/nanosecond'</li> <li>• 'cycles/microsecond'</li> <li>• 'cycles/millisecond'</li> <li>• 'cycles/hour'</li> <li>• 'cycles/day'</li> <li>• 'cycles/week'</li> <li>• 'cycles/month'</li> <li>• 'cycles/year'</li> </ul> <p><b>Default:</b> 'rad/s'</p> <p>You can also specify 'auto', which uses frequency units rad/TimeUnit relative to system time units specified in the</p>

<b>Field</b>	<b>Description</b>
	TimeUnit property. For multiple systems with different time units, the units of the first system are used.
FreqScale	Frequency scale, specified as one of the following values: 'linear'   'log' <b>Default:</b> 'log'
MagUnits	Magnitude units, specified as one of the following values: 'dB'   'abs' <b>Default:</b> 'dB'
MagScale	Magnitude scale, specified as one of the following values: 'linear'   'log' <b>Default:</b> 'linear'
MagLowerLimMode	Enables a lower magnitude limit, specified as one of the following values: 'auto'   'manual' <b>Default:</b> 'auto'
MagLowerLim	Lower magnitude limit, specified as data type double.

## See Also

getoptions | identpref | setoptions | spectrumplot

**Introduced in R2012a**

# spectrumplot

Plot disturbance spectrum of linear identified models

## Syntax

```
spectrumplot(sys)
spectrumplot(sys,line_spec)
spectrumplot(sys1,line_spec1,...,sysN,line_specN)
spectrumplot(ax, ___)
spectrumplot(___,plot_options)
spectrumplot(sys,w)
h = spectrumplot(___)
```

## Description

`spectrumplot(sys)` plots the disturbance spectrum of the model, `sys`. The software chooses the number of points on the plot and the plot frequency range.

If `sys` is a time-series model, its disturbance spectrum is the same as the model output spectrum. You generally use this function with time-series models.

`spectrumplot(sys,line_spec)` uses `line_spec` to specify the line type, marker symbol, and color.

`spectrumplot(sys1,line_spec1,...,sysN,line_specN)` plots the disturbance spectrum for one or more models on the same axes.

You can mix `sys,line_spec` pairs with `sys` models as in `spectrumplot(sys1,sys2,line_spec2,sys3)`. `spectrumplot` automatically chooses colors and line styles in the order specified by the `ColorOrder` and `LineStyleOrder` properties of the current axes.

`spectrumplot(ax, ___)` plots into the axes with handle `ax`. All input arguments described for the previous syntaxes also apply here.

`spectrogram( ____, plot_options )` uses `plot_options` to specify options such as plot title, frequency units, etc. All input arguments described for the previous syntaxes also apply here.

`spectrogram(sys, w)` uses `w` to specify the plot frequencies.

- If `w` is specified as a 2-element cell array, `{wmin, wmax}`, the plot spans the frequency range `{wmin, wmax}`.
- If `w` is specified as vector, the spectrum is plotted for the specified frequencies.

Specify `w` as `radians/time_unit`, where `time_unit` must equal `sys.TimeUnit`.

`h = spectrogram( ____, )` returns the handle to the spectrum plot. You use the handle to customize the plot. All input arguments described for the previous syntaxes also apply here.

## Input Arguments

### **sys**

Identified linear model.

### **Default:**

### **line\_spec**

Line style, marker, and color of both the line and marker, specified as a character vector. For example, `'b'`, `'b+:'`.

For more information, see [Chart Line](#) .

### **ax**

Plot axes handle.

Specify as a double-precision value.

You can obtain the current axes handle by using the function, `gca`.

### **plot\_options**

Plot customization options.

Specify as a plot options object.

You use the command, `spectrumoptions`, to create `plot_options`. For more information, type `help spectrumoptions`.

**w**

Frequency range.

Specify in radians/time\_unit, where time\_unit must equal `sys.TimeUnit`.

## Output Arguments

**h**

Plot handle for spectrum plot, returned as a double-precision value.

## Examples

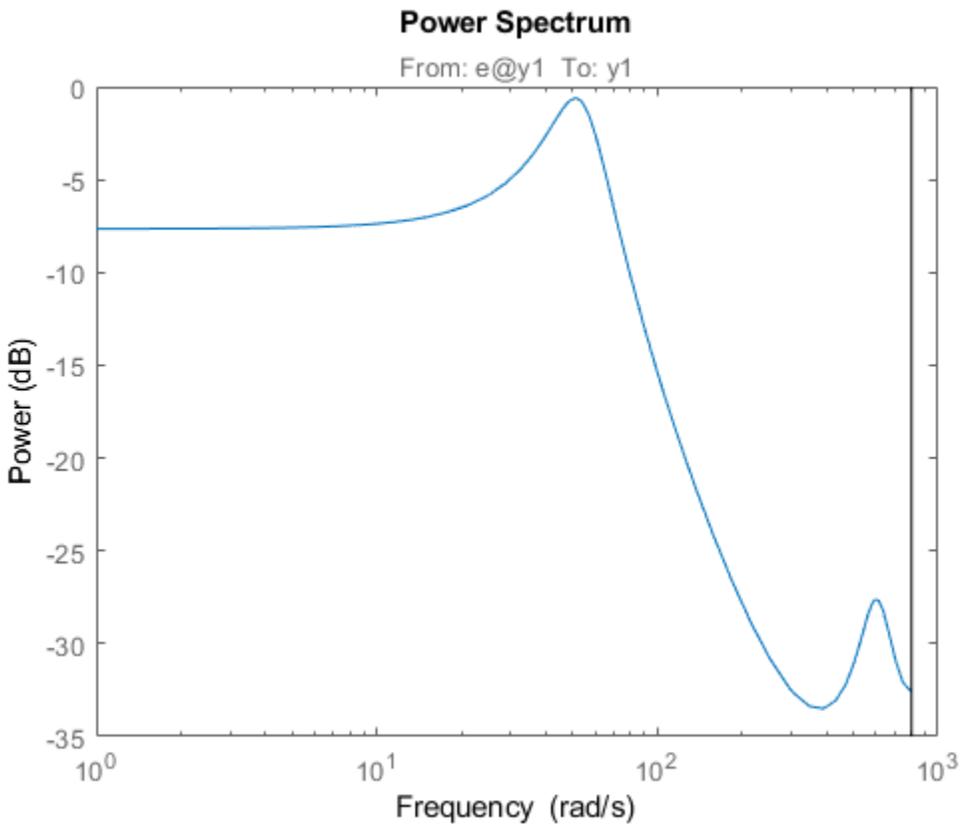
### Plot Model Output Spectrum for Identified Model

Obtain the identified model.

```
load iddata9 z9  
sys = ar(z9,4);
```

Plot the output spectrum for the model.

```
spectrumplot(sys);
```



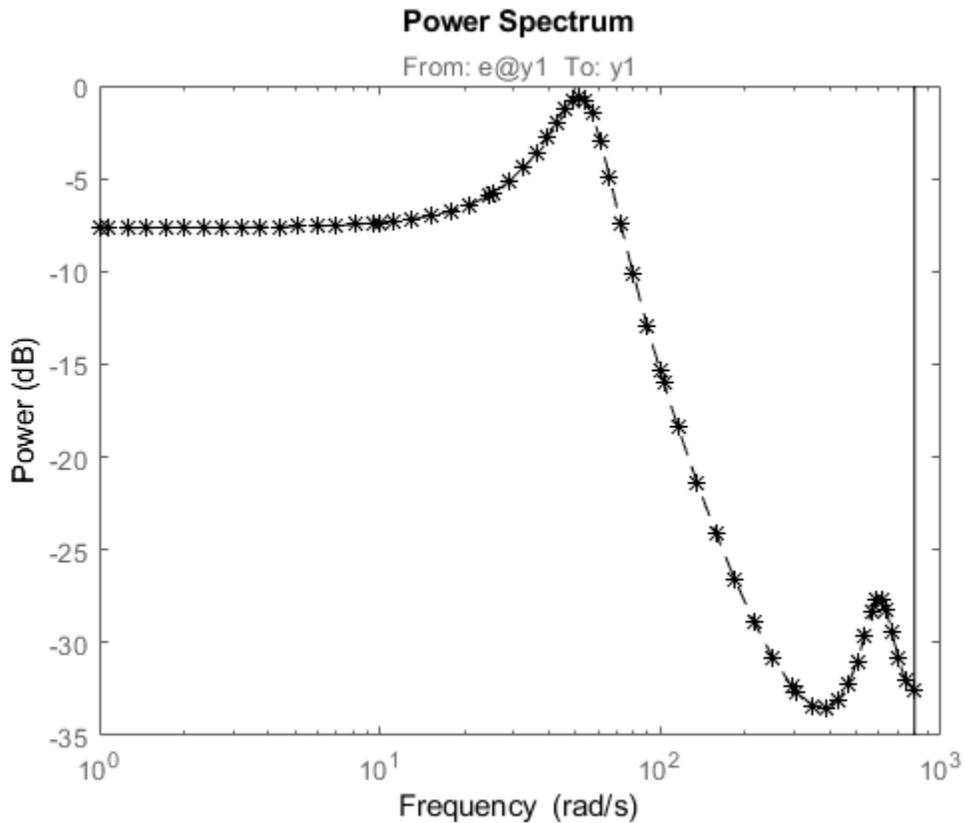
### Specify Line Width and Marker Style on Spectrum Plot

Obtain the identified model.

```
load iddata9 z9  
sys = ar(z9,4);
```

Specify the line width and marker style for the spectrum plot.

```
spectrumplot(sys, 'k*--');
```



'k\*--', specifies a dashed line (--) that is black (k), with star markers (\*).

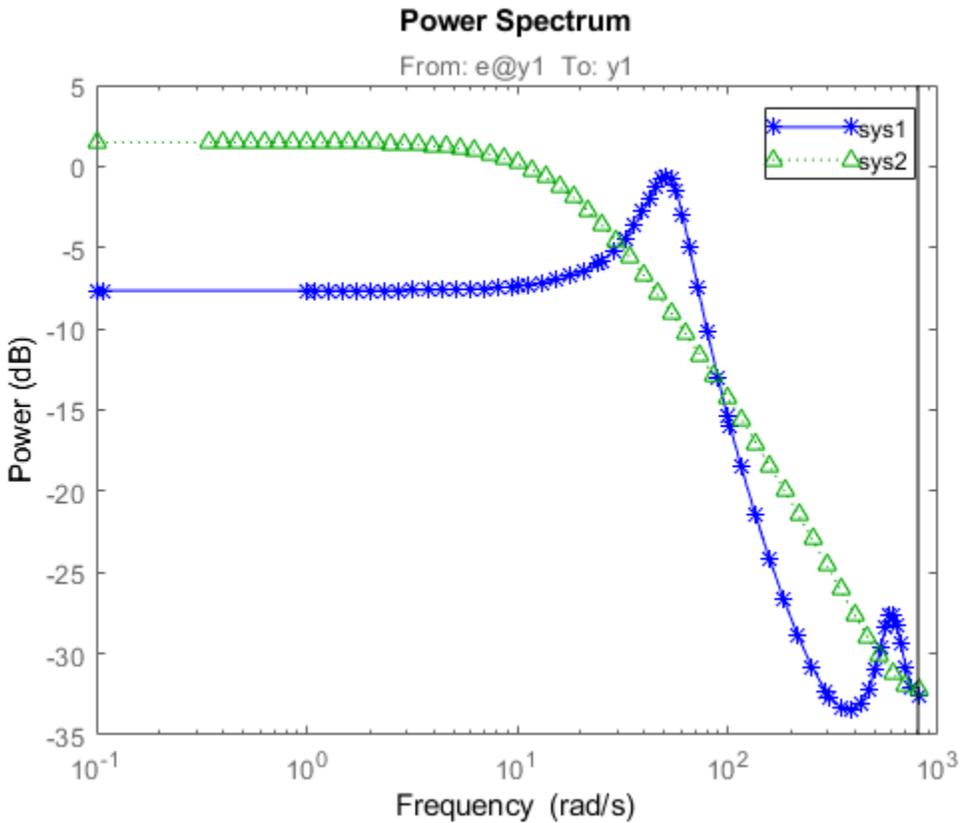
### Plot Multiple Models on the Same Axes

Obtain multiple identified models.

```
load iddata9 z9
sys1 = ar(z9,4);
sys2 = ar(z9,2);
```

Plot the output spectrum for both models.

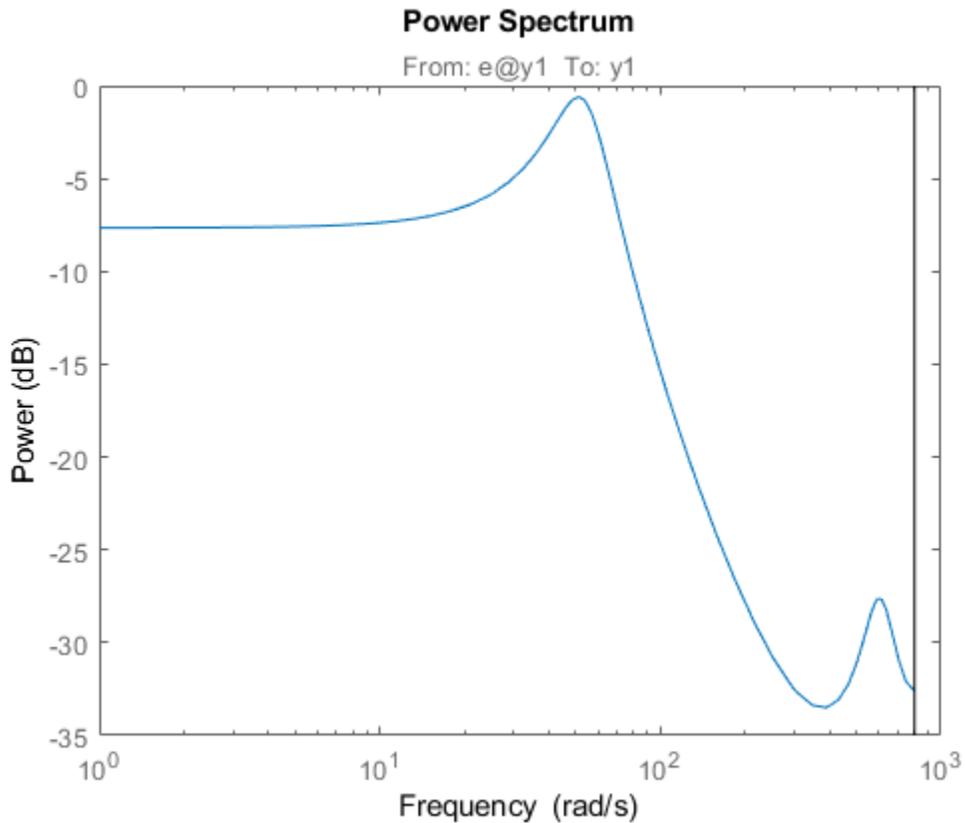
```
spectrumplot(sys1, 'b*-', sys2, 'g^:');  
legend('sys1', 'sys2');
```



### Specify Plot Axes for Spectrum Plot

Obtain the axes handle for a plot.

```
load iddata9 z9  
sys1 = ar(z9,4);  
spectrumplot(sys1);
```



```
ax = gca;
```

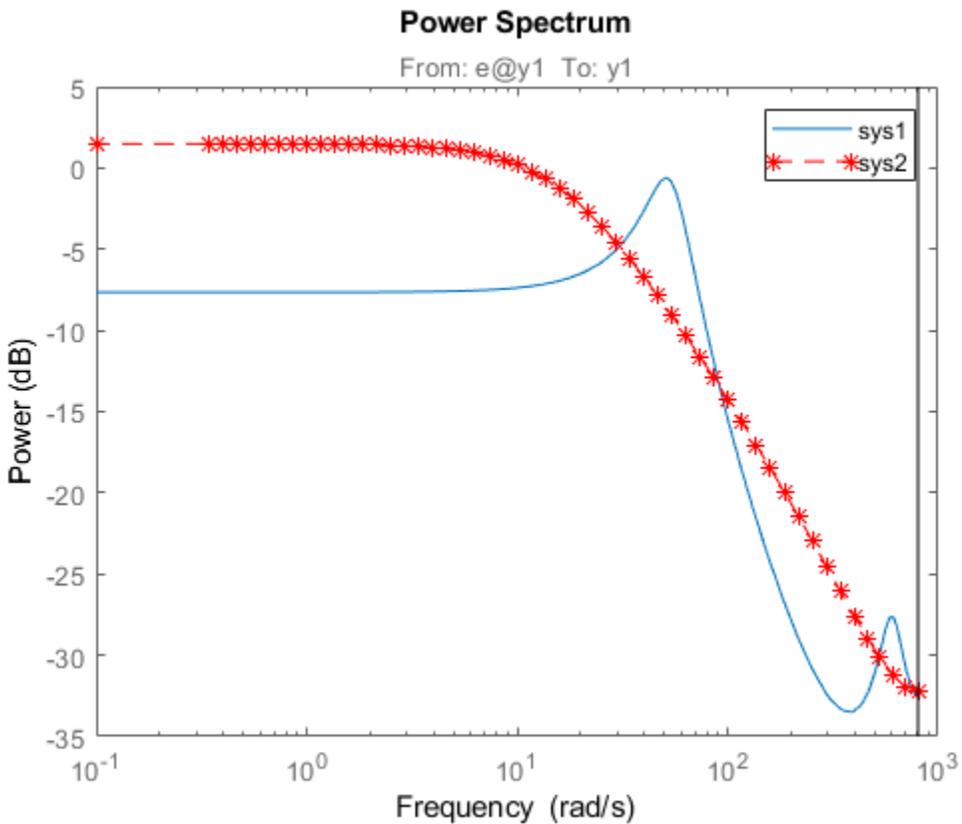
ax is the handle for the spectrum plot axes.

Plot the output spectrum for another model on the specified axes.

```
sys2 = ar(z9,2);
```

```
hold on;  
spectrumplot(ax,sys2,'r*--');
```

```
legend('sys1','sys2');
```



### Specify Options for Spectrum Plot

Specify the plot options.

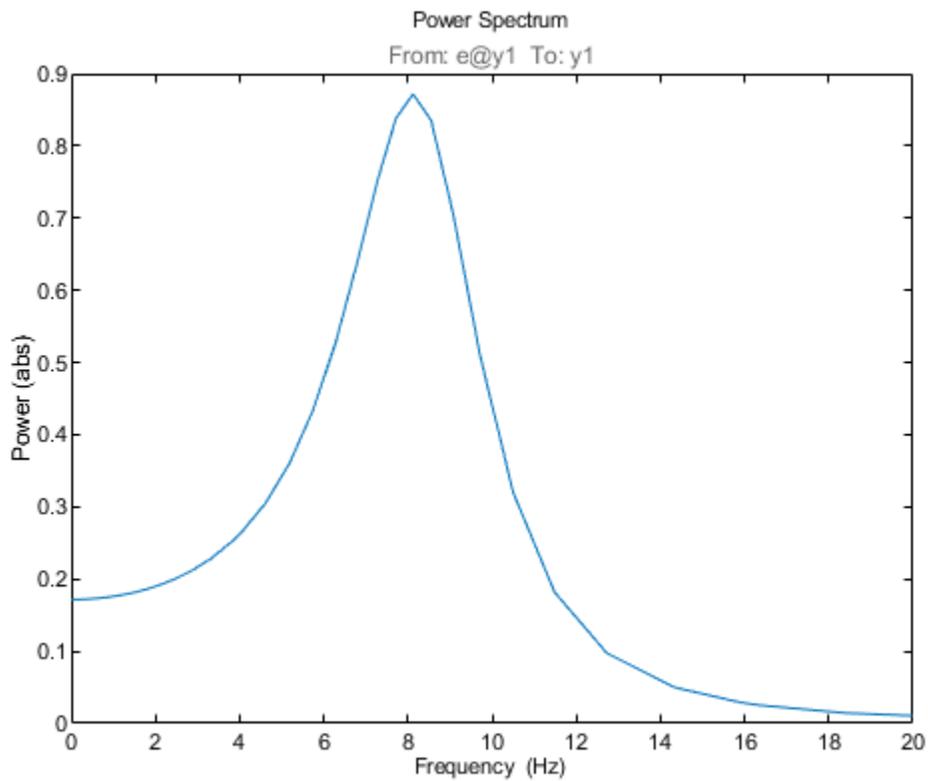
```
plot_options = spectrumoptions;  
plot_options.FreqUnits = 'Hz';  
plot_options.FreqScale = 'linear';  
plot_options.Xlim = {[0 20]};  
plot_options.MagUnits = 'abs';
```

Estimate an AR model.

```
load iddata9 z9  
sys = ar(z9,4);
```

Plot the output spectrum for the model.

```
spectrumplot(sys,plot_options);
```



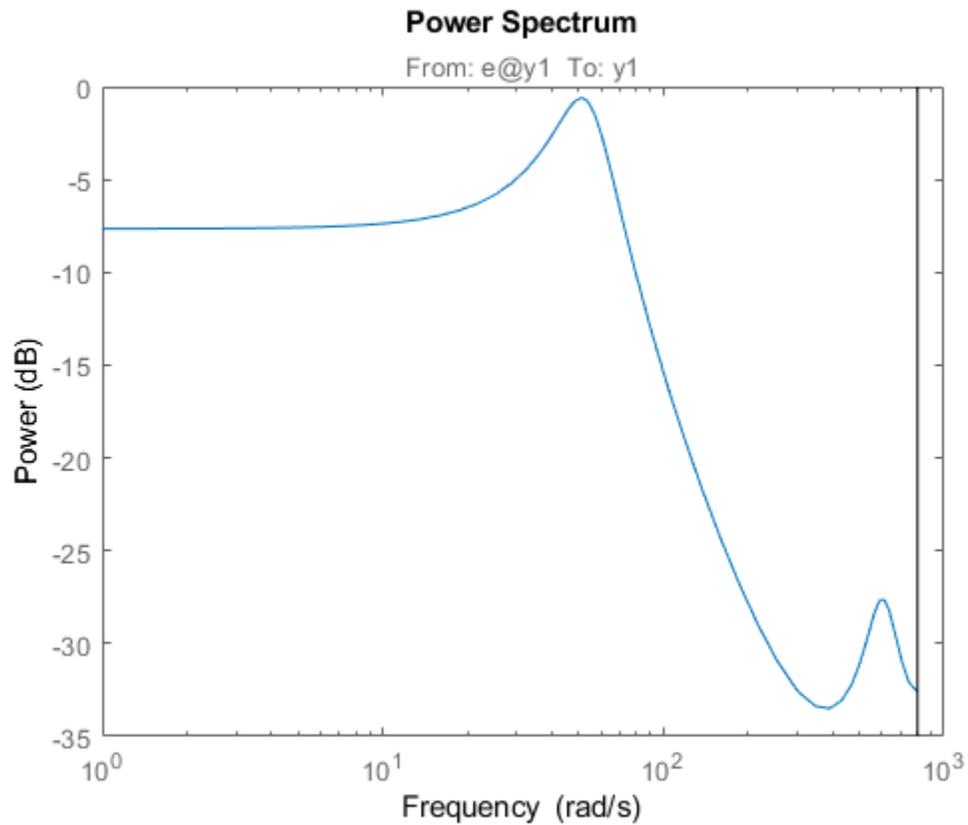
### Specify Spectrum Plot Frequency Range

Obtain the identified model.

```
load iddata9 z9  
sys = ar(z9,4);
```

Specify the frequency range for the output spectrum plot for the model.

```
spectrumplot(sys, {1, 1000});
```



The 2-element cell array `{1, 1000}` specifies the frequency range from 1 rad/s to 1000 rad/s.

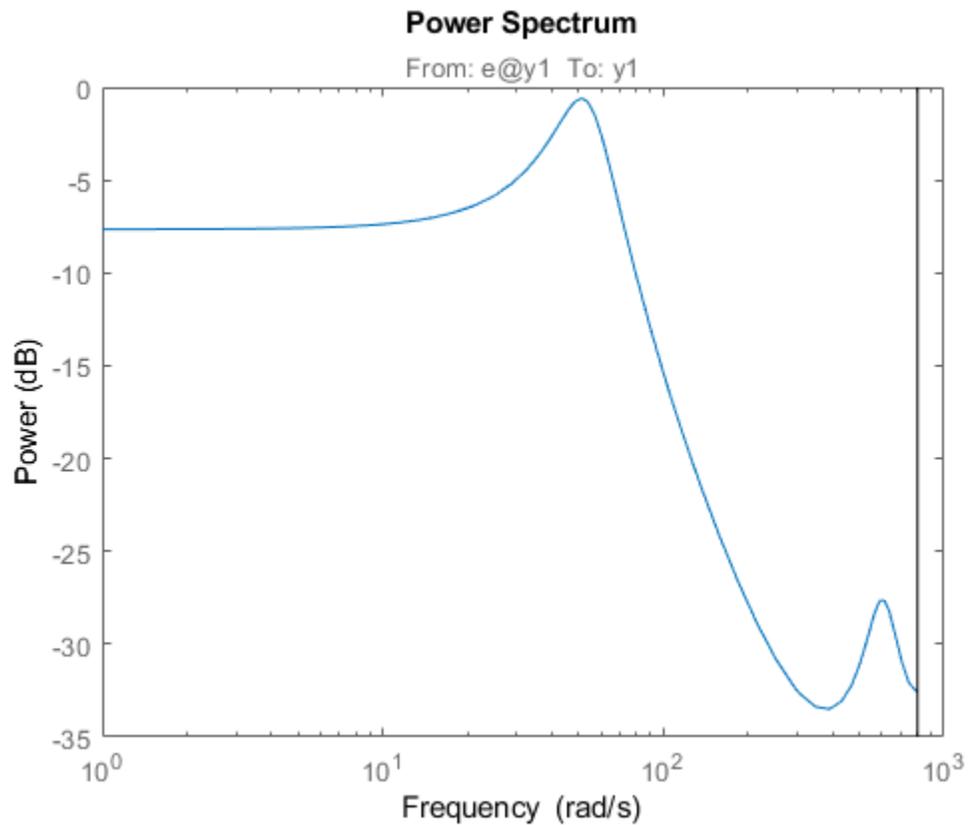
## Get Plot Handle for Spectrum Plot Customization

Obtain the identified model.

```
load iddata9 z9  
sys = ar(z9,4);
```

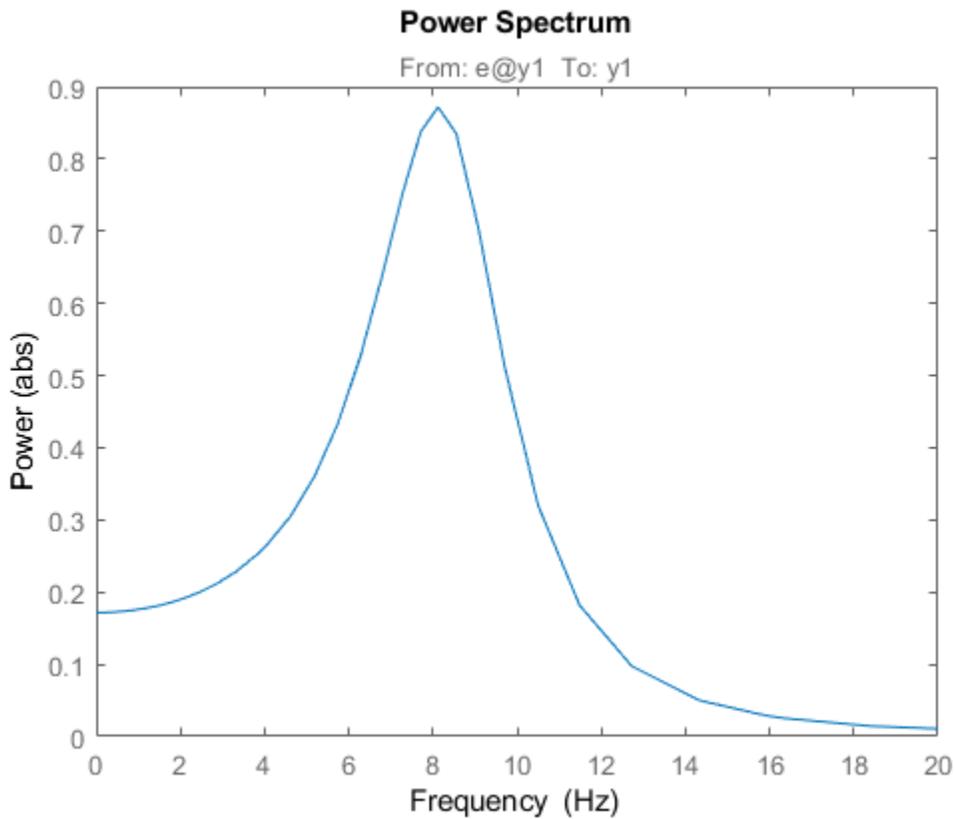
Get the plot handle for the model spectrum plot.

```
h = spectrumplot(sys);
```



(Optional) Specify the plot options, using the plot handle.

```
setoptions(h, 'FreqUnits', 'Hz', 'FreqScale', 'linear', 'Xlim', {[0 20]}, 'MagUnits', 'abs');
```



## See Also

[Axes](#) | [Chart Line](#) | [getoptions](#) | [setoptions](#) | [showConfidence](#) | [spectrum](#) | [spectrumoptions](#)

**Introduced in R2012b**

## ss2ss

State coordinate transformation for state-space model

### Syntax

`sysT = ss2ss(sys,T)`

### Description

Given a state-space model `sys` with equations

$$\begin{aligned}\dot{x} &= Ax + Bu \\ y &= Cx + Du\end{aligned}$$

or the innovations form used by the identified state-space (IDSS) models:

$$\begin{aligned}\frac{dx}{dt} &= Ax + Bu + Ke \\ y &= Cx + Du + e\end{aligned}$$

(or their discrete-time counterpart), `ss2ss` performs the similarity transformation  $\bar{x} = Tx$  on the state vector  $x$  and produces the equivalent state-space model `sysT` with equations.

$$\begin{aligned}\dot{\bar{x}} &= TAT^{-1}\bar{x} + TBu \\ y &= CT^{-1}\bar{x} + Du\end{aligned}$$

or, in the case of an IDSS model:

$$\begin{aligned}\dot{\bar{x}} &= TAT^{-1}\bar{x} + TBu + TKe \\ y &= CT^{-1}\bar{x} + Du + e\end{aligned}$$

`sysT = ss2ss(sys,T)` returns the transformed state-space model `sysT` given `sys` and the state coordinate transformation `T`. The model `sys` must be in state-space form and the matrix `T` must be invertible. `ss2ss` is applicable to both continuous- and discrete-time models.

## Examples

Perform a similarity transform to improve the conditioning of the `A` matrix.

```
T = balance(sys.A)
sysb = ss2ss(sys,inv(T))
```

## See Also

`balreal` | `canon`

**Introduced before R2006a**

## ssdata

Access state-space model data

### Syntax

```
[a,b,c,d] = ssdata(sys)
[a,b,c,d,Ts] = ssdata(sys)
```

### Description

`[a,b,c,d] = ssdata(sys)` extracts the matrix (or multidimensional array) data A, B, C, D from the state-space model (LTI array) `sys`. If `sys` is a transfer function or zero-pole-gain model (LTI array), it is first converted to state space. See `ss` for more information on the format of state-space model data.

If `sys` appears in descriptor form (nonempty E matrix), an equivalent explicit form is first derived.

If `sys` has internal delays, A, B, C, D are obtained by first setting all internal delays to zero (creating a zero-order Padé approximation). For some systems, setting delays to zero creates singular algebraic loops, which result in either improper or ill-defined, zero-delay approximations. For these systems, `ssdata` cannot display the matrices and returns an error. This error does not imply a problem with the model `sys` itself.

For generalized state-space (`genss`) models, `ssdata` returns the state-space models evaluated at the current, nominal value of all control design blocks. To access the dependency of a `genss` model on its static control design blocks, use the A, B, C, and D properties of the model.

`[a,b,c,d,Ts] = ssdata(sys)` also returns the sample time `Ts`.

You can access the remaining LTI properties of `sys` with `get` or by direct referencing. For example:

```
sys.statename
```

For arrays of state-space models with variable numbers of states, use the syntax:

```
[a,b,c,d] = ssdata(sys,'cell')
```

to extract the state-space matrices of each model as separate cells in the cell arrays **a**, **b**, **c**, and **d**.

### **See Also**

[dssdata](#) | [get](#) | [getdelaymodel](#) | [idssdata](#) | [set](#) | [ss](#) | [tfdata](#) | [zpkdata](#)

**Introduced before R2006a**

## ssest

Estimate state-space model using time or frequency domain data

### Syntax

```
sys = ssest(data,nx)
sys = ssest(data,nx,Name,Value)
sys = ssest( ___,opt)
```

```
sys = ssest(data,init_sys)
sys = ssest(data,init_sys,opt)
```

```
[sys,x0] = ssest( ___ )
```

### Description

`sys = ssest(data,nx)` estimates a state-space model, `sys`, using time- or frequency-domain data, `data`. `sys` is a state-space model of order `nx` and represents:

$$\begin{aligned}\dot{x}(t) &= Ax(t) + Bu(t) + Ke(t) \\ y(t) &= Cx(t) + Du(t) + e(t)\end{aligned}$$

$A$ ,  $B$ ,  $C$ ,  $D$ , and  $K$  are state-space matrices.  $u(t)$  is the input,  $y(t)$  is the output,  $e(t)$  is the disturbance and  $x(t)$  is the vector of `nx` states.

All the entries of  $A$ ,  $B$ ,  $C$ , and  $K$  are free estimable parameters by default.  $D$  is fixed to zero by default, meaning that there is no feedthrough, except for static systems (`nx=0`).

`sys = ssest(data,nx,Name,Value)` estimates the model using the additional options specified by one or more `Name,Value` pair arguments. Use the `Form`, `Feedthrough` and `DisturbanceModel` name-value pair arguments to modify the default behavior of the  $A$ ,  $B$ ,  $C$ ,  $D$ , and  $K$  matrices.

`sys = ssest( ___,opt)` estimates the model using an option set, `opt`, that specifies options such as estimation objective, handling of initial conditions and numerical search method used for estimation.

`sys = ssest(data,init_sys)` estimates a state-space model using the linear system `init_sys` to configure the initial parameterization.

`sys = ssest(data,init_sys,opt)` estimates the model using an option set, `opt`.

`[sys,x0] = ssest( ___ )` returns the value of initial states computed during estimation.

## Input Arguments

### **data**

Estimation data.

For time-domain estimation, `data` must be an `iddata` object containing the input and output signal values.

For frequency-domain estimation, `data` can be one of the following:

- Recorded frequency response data (`frd` or `idfrd`)
- `iddata` object with its properties specified as follows:
  - `InputData` — Fourier transform of the input signal
  - `OutputData` — Fourier transform of the output signal
  - `Domain` — 'Frequency'

### **nx**

Order of estimated model.

Specify `nx` as a positive integer. `nx` may be a scalar or a vector. If `nx` is a vector, then `ssest` creates a plot which you can use to choose a suitable model order. The plot shows the Hankel singular values for models of different orders. States with relatively small Hankel singular values can be safely discarded. A default choice is suggested in the plot.

### **opt**

Estimation options.

`opt` is an options set, created using `ssestOptions`, that specifies options including:

- Estimation objective
- Handling of initial conditions
- Numerical search method used for estimation

If `opt` is not specified and `init_sys` is a previously estimated `idss` model, the options from `init_sys.Report.OptionsUsed` are used.

### **init\_sys**

Linear system that configures the initial parameterization of `sys`.

You obtain `init_sys` by either performing an estimation using measured data or by direct construction.

If `init_sys` is an `idss` model, `ssest` uses the parameter values of `init_sys` as the initial guess for estimating `sys`. For information on how to specify `idss`, see “Estimate State-Space Models with Structured Parameterization”. Constraints on the parameters of `init_sys`, such as fixed coefficients and minimum/maximum bounds are honored in estimating `sys`.

Use the `Structure` property of `init_sys` to configure initial guesses and constraints for the  $A$ ,  $B$ ,  $C$ ,  $D$ , and  $K$  matrices. For example:

- To specify an initial guess for the  $A$  matrix of `init_sys`, set `init_sys.Structure.A.Value` as the initial guess.
- To specify constraints for the  $B$  matrix of `init_sys`:
  - Set `init_sys.Structure.B.Minimum` to the minimum  $B$  matrix value
  - Set `init_sys.Structure.B.Maximum` to the maximum  $B$  matrix value
  - Set `init_sys.Structure.B.Free` to indicate if entries of the  $B$  matrix are free parameters for estimation

If `init_sys` is not a state-space (`idss`) model, the software first converts `init_sys` to an `idss` model. `ssest` uses the parameters of the resulting model as the initial guess for estimation.

If `opt` is not specified, and `init_sys` was obtained by estimation, then the estimation options from `init_sys.Report.OptionsUsed` are used.

## Name-Value Pair Arguments

Specify optional comma-separated pairs of `Name`, `Value` arguments. `Name` is the argument name and `Value` is the corresponding value. `Name` must appear inside quotes. You can specify several name and value pair arguments in any order as `Name1, Value1, . . . , NameN, ValueN`.

### **Ts** — Sample time

0 (continuous-time) (default) | sample time of data (`data.Ts`) | positive scalar

Sample time, specified as a positive scalar.

For continuous-time models, use `Ts = 0`. For discrete-time models, specify `Ts` as a positive scalar whose value is equal to the data sample time.

### **InputDelay** — Input delays

0 (default) | scalar | vector

Input delay for each input channel, specified as a numeric vector. For continuous-time systems, specify input delays in the time unit stored in the `TimeUnit` property. For discrete-time systems, specify input delays in integer multiples of the sample time `Ts`. For example, `InputDelay = 3` means a delay of three sampling periods.

For a system with `Nu` inputs, set `InputDelay` to an `Nu`-by-1 vector. Each entry of this vector is a numerical value that represents the input delay for the corresponding input channel.

You can also set `InputDelay` to a scalar value to apply the same delay to all channels.

### **Form** — Type of canonical form

'free' (default) | 'modal' | 'companion' | 'canonical'

Type of canonical form of `sys`, specified as one of the following values:

- 'modal' — Obtain `sys` in modal form on page 1-1663.
- 'companion' — Obtain `sys` in companion form on page 1-1663.
- 'free' — All entries of the `A`, `B` and `C` matrices are treated as free.
- 'canonical' — Obtain `sys` in the observability canonical form [1].

Use the `Form`, `Feedthrough` and `DisturbanceModel` name-value pair arguments to modify the default behavior of the `A`, `B`, `C`, `D`, and `K` matrices.

For more information, see “Estimate State-Space Models with Canonical Parameterization”.

### **Feedthrough — Direct feedthrough from input to output**

0 (default) | 1 | logical vector

Direct feedthrough from input to output, specified as a logical vector of length  $Nu$ , where  $Nu$  is the number of inputs. If `Feedthrough` is specified as a logical scalar, it is applied to all the inputs.

Use the `Form`, `Feedthrough` and `DisturbanceModel` name-value pair arguments to modify the default behavior of the  $A$ ,  $B$ ,  $C$ ,  $D$ , and  $K$  matrices.

### **DisturbanceModel — Specify whether to estimate the $K$ matrix**

'estimate' (default) | 'none'

Specify whether to estimate the  $K$  matrix which specifies the noise component, specified as one of the following values:

- 'none' — Noise component is not estimated. The value of the  $K$  matrix is fixed to zero value.
- 'estimate' — The  $K$  matrix is treated as a free parameter.

`DisturbanceModel` must be 'none' when using frequency-domain data.

Use the `Form`, `Feedthrough` and `DisturbanceModel` name-value pair arguments to modify the default behavior of the  $A$ ,  $B$ ,  $C$ ,  $D$ , and  $K$  matrices.

## **Output Arguments**

### **sys**

Identified state-space model, returned as a `idss` model. This model is created using the specified model orders, delays, and estimation options.

Information about the estimation results and options used is stored in the `Report` property of the model. `Report` has the following fields:

<b>Report Field</b>	<b>Description</b>
Status	Summary of the model status, which indicates whether the model was created by construction or obtained by estimation.
Method	Estimation command used.
Initial State	<p>How initial states were handled during estimation, returned as one of the following values:</p> <ul style="list-style-type: none"> <li>• 'zero' — The initial state is set to zero.</li> <li>• 'estimate' — The initial state is treated as an independent estimation parameter.</li> <li>• 'backcast' — The initial state is estimated using the best least squares fit.</li> <li>• Column vector of length <math>N_x</math>, where <math>N_x</math> is the number of states. For multi-experiment data, a matrix with <math>N_e</math> columns, where <math>N_e</math> is the number of experiments.</li> <li>• Parametric initial condition object (<code>x0obj</code>) created using <code>idpar</code>. Only for discrete-time state-space models.</li> </ul> <p>This field is especially useful when the <code>InitialState</code> option in the estimation option set is 'auto'.</p>
N4Weight	<p>Weighting scheme used for singular-value decomposition by the N4SID algorithm, returned as one of the following values:</p> <ul style="list-style-type: none"> <li>• 'MOESP' — Uses the MOESP algorithm.</li> <li>• 'CVA' — Uses the Canonical Variable Algorithm.</li> <li>• 'SSARX' — A subspace identification method that uses an ARX estimation based algorithm to compute the weighting.</li> </ul> <p>This option is especially useful when the <code>N4Weight</code> option in the estimation option set is 'auto'.</p>
N4Horizon	Forward and backward prediction horizons used by the N4SID algorithm, returned as a row vector with three elements — <code>[r sy su]</code> , where <code>r</code> is the maximum forward prediction horizon. <code>sy</code> is the number of past outputs, and <code>su</code> is the number of past inputs that are used for the predictions.

Report Field	Description																		
Fit	<p>Quantitative assessment of the estimation, returned as a structure. See “Loss Function and Model Quality Metrics” for more information on these quality metrics. The structure has the following fields:</p> <table border="1" data-bbox="387 456 1332 951"> <thead> <tr> <th data-bbox="387 456 520 499">Field</th> <th data-bbox="520 456 1332 499">Description</th> </tr> </thead> <tbody> <tr> <td data-bbox="387 499 520 609">FitPercent</td> <td data-bbox="520 499 1332 609">Normalized root mean squared error (NRMSE) measure of how well the response of the model fits the estimation data, expressed as a percentage.</td> </tr> <tr> <td data-bbox="387 609 520 652">LossFcn</td> <td data-bbox="520 609 1332 652">Value of the loss function when the estimation completes.</td> </tr> <tr> <td data-bbox="387 652 520 730">MSE</td> <td data-bbox="520 652 1332 730">Mean squared error (MSE) measure of how well the response of the model fits the estimation data.</td> </tr> <tr> <td data-bbox="387 730 520 774">FPE</td> <td data-bbox="520 730 1332 774">Final prediction error for the model.</td> </tr> <tr> <td data-bbox="387 774 520 817">AIC</td> <td data-bbox="520 774 1332 817">Raw Akaike Information Criteria (AIC) measure of model quality.</td> </tr> <tr> <td data-bbox="387 817 520 861">AICc</td> <td data-bbox="520 817 1332 861">Small sample-size corrected AIC.</td> </tr> <tr> <td data-bbox="387 861 520 904">nAIC</td> <td data-bbox="520 861 1332 904">Normalized AIC.</td> </tr> <tr> <td data-bbox="387 904 520 951">BIC</td> <td data-bbox="520 904 1332 951">Bayesian Information Criteria (BIC).</td> </tr> </tbody> </table>	Field	Description	FitPercent	Normalized root mean squared error (NRMSE) measure of how well the response of the model fits the estimation data, expressed as a percentage.	LossFcn	Value of the loss function when the estimation completes.	MSE	Mean squared error (MSE) measure of how well the response of the model fits the estimation data.	FPE	Final prediction error for the model.	AIC	Raw Akaike Information Criteria (AIC) measure of model quality.	AICc	Small sample-size corrected AIC.	nAIC	Normalized AIC.	BIC	Bayesian Information Criteria (BIC).
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Parameters	Estimated values of model parameters.																		
OptionsUsed	Option set used for estimation. If no custom options were configured, this is a set of default options. See <code>ssestOptions</code> for more information.																		
RandState	State of the random number stream at the start of estimation. Empty, [], if randomization was not used during estimation. For more information, see <code>rng</code> in the MATLAB documentation.																		

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Report Field	Description	
Termination	Termination conditions for the iterative search used for prediction error minimization. Structure with the following fields:	
	Field	Description
	WhyStop	Reason for terminating the numerical search.
	Iterations	Number of search iterations performed by the estimation algorithm.
	FirstOrderOptimality	$\infty$ -norm of the gradient search vector when the search algorithm terminates.
	FcnCount	Number of times the objective function was called.
	UpdateNorm	Norm of the gradient search vector in the last iteration. Omitted when the search method is 'lsqnonlin' or 'fmincon'.
	LastImprovement	Criterion improvement in the last iteration, expressed as a percentage. Omitted when the search method is 'lsqnonlin' or 'fmincon'.
	Algorithm	Algorithm used by 'lsqnonlin' or 'fmincon' search method. Omitted when other search methods are used.
For estimation methods that do not require numerical search optimization, the Termination field is omitted.		

For more information on using Report, see “Estimation Report”.

### **x0**

Initial states computed during the estimation.

If data contains multiple experiments, then x0 is an array with each column corresponding to an experiment.

This value is also stored in the Parameters field of the model's Report property.

## Examples

### Determine Optimal Estimated Model Order

Obtain measured input-output data.

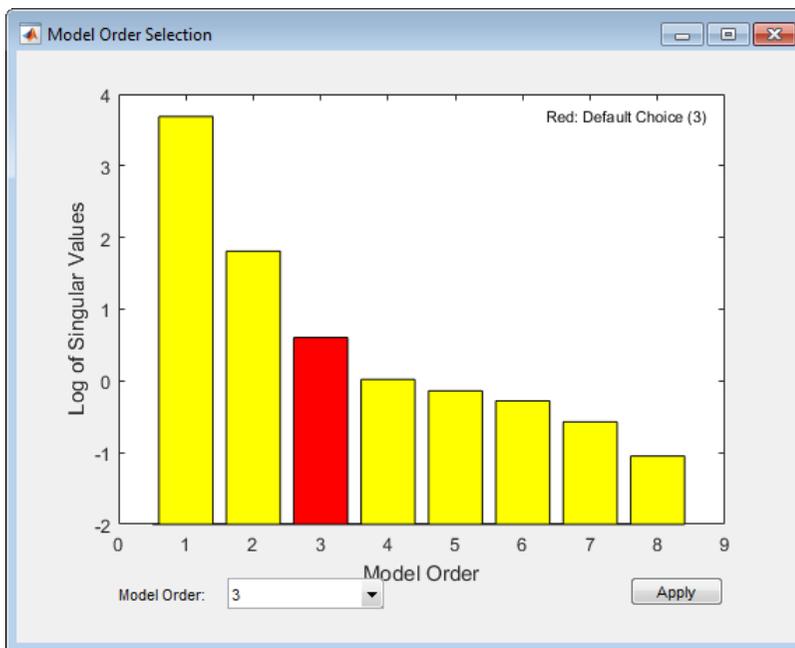
```
load icEngine.mat;  
data = iddata(y,u,0.04);
```

data is an iddata object containing 1500 input-output data samples. The data sample time is 0.04 seconds.

Estimate a state-space model for measured input-output data. Determine the optimal model order within a given model order range.

```
nx = 1:10;  
sys = ssest(data,nx);
```

A plot that shows the Hankel singular values (SVD) for models of the orders specified by nx appears.



States with relatively small Hankel singular values can be safely discarded. The suggested default order choice is 3.

Select the model order in the **Model Order** drop-down list and click **Apply**.

## Identify State-Space Model With Input Delay

Load time-domain system response data.

```
load iddata7 z7;
```

Identify a fourth-order state-space model of the data. Specify a known delay of 2 seconds for the first input and 0 seconds for the second input.

```
nx = 4;
sys = ssest(z7(1:300),nx,'InputDelay',[2;0]);
```

## Estimate State-Space Model Using Regularization

Obtain a regularized 5th order state-space model for a 2nd order system from a narrow bandwidth signal.

Load estimation data.

```
load regularizationExampleData eData;
```

Create the transfer function model used for generating the estimation data (true system).

```
trueSys = idtf([0.02008 0.04017 0.02008],[1 -1.561 0.6414],1);
```

Estimate an unregularized state-space model.

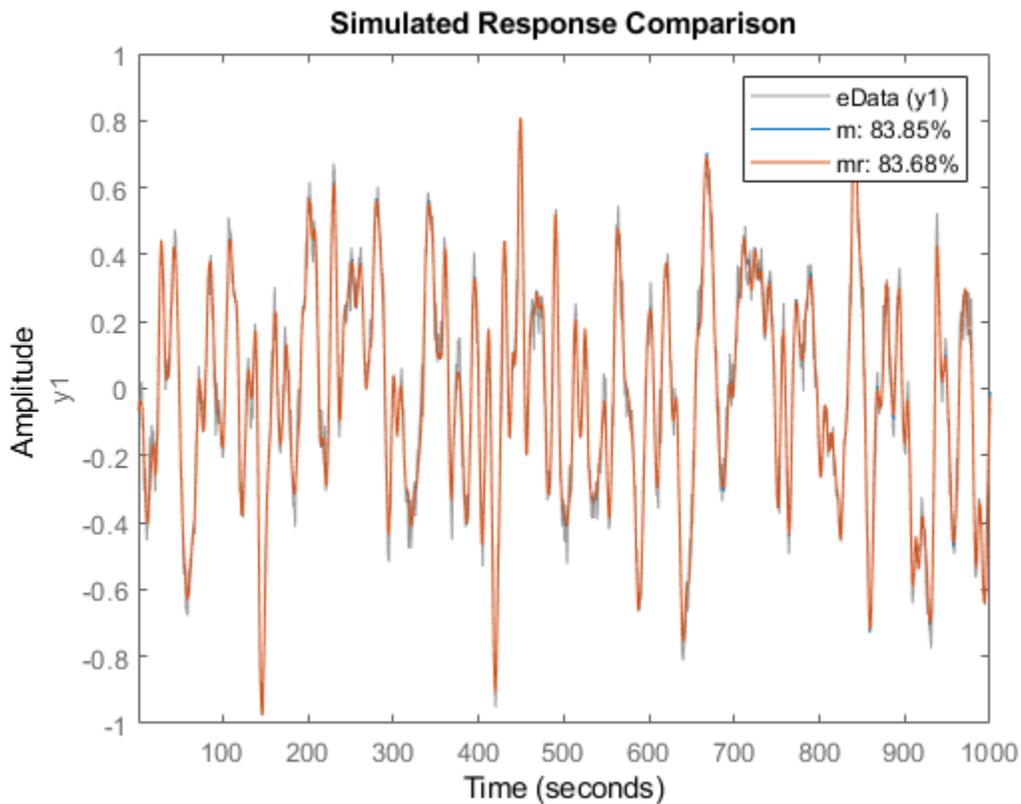
```
opt = ssestOptions('SearchMethod','lm');
m = ssest(eData,5,'form','modal','DisturbanceModel','none','Ts',eData.Ts,opt);
```

Estimate a regularized state-space model.

```
opt.Regularization.Lambda = 10;
mr = ssest(eData,5,'form','modal','DisturbanceModel','none','Ts',eData.Ts,opt);
```

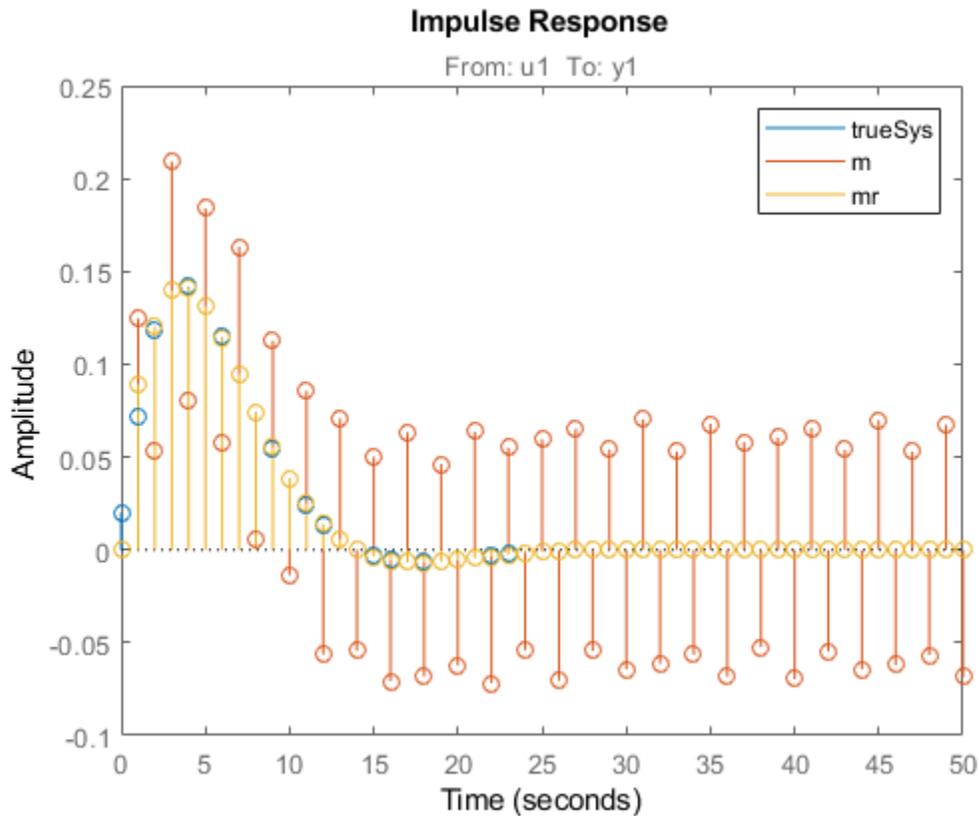
Compare the model outputs with the estimation data.

```
compare(eData,m,mr);
```



Compare the model impulse responses.

```
impulse(trueSys,m,mr,50);  
legend('trueSys','m','mr');
```



## Estimate Partially Known State-Space Model Using Structured Estimation

Estimate a state-space model of measured input-output data. Configure the parameter constraints and initial values for estimation using a state-space model.

Create an `idss` model to specify the initial parameterization for estimation.

```
A = blkdiag([-0.1 0.4; -0.4 -0.1],[-1 5; -5 -1]);
B = [1; zeros(3,1)];
C = [1 1 1 1];
D = 0;
```

```
K = zeros(4,1);
x0 = [0.1 0.1 0.1 0.1];
Ts = 0;
init_sys = idss(A,B,C,D,K,x0,Ts);
```

Setting all entries of `K` to `0` creates an `idss` model with no state disturbance element.

Use the `Structure` property to fix the values of some of the model parameters. Configure the model so that `B` and `K` are fixed, and only the nonzero entries of `A` are estimable.

```
init_sys.Structure.A.Free = (A~=0);
init_sys.Structure.B.Free = false;
init_sys.Structure.K.Free = false;
```

The entries in `init_sys.Structure.A.Free` determine whether the corresponding entries in `init_sys.A` are free (true) or fixed (false).

Load the measured data and estimate a state-space model using the parameter constraints and initial values specified by `init_sys`.

```
load iddata2 z2;
sys = ssest(z2,init_sys);
```

The estimated parameters of `sys` satisfy the constraints specified by `init_sys`.

## Model Order Reduction by Estimation

Consider the Simulink model `idF14Model`. Linearizing this model gives a ninth-order model. However, the dynamics of the model can be captured, without compromising the fit quality too much, using a lower-order model.

Obtain the linearized model.

```
load_system('idF14Model');
io = getlinio('idF14Model');
sys_lin = linearize('idF14Model',io);
```

`sys_lin` is a ninth-order state-space model with two outputs and one input.

Simulate the step response of the linearized model, and use the data to create an `iddata` object.

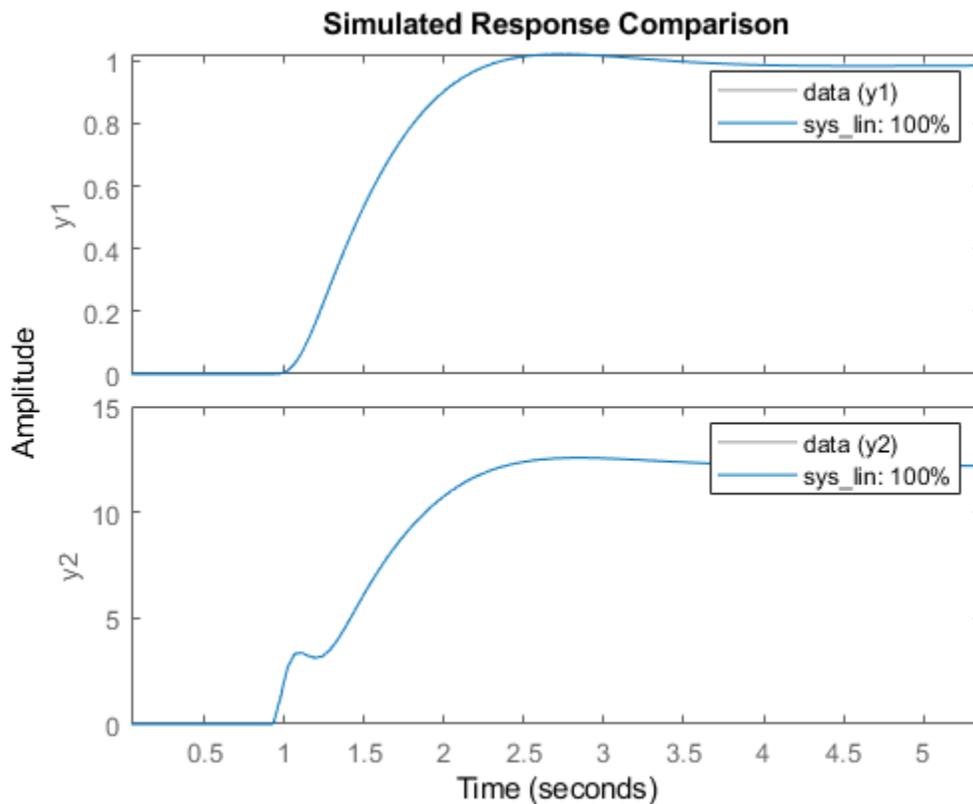
```
Ts = 0.0444;  
t = (0:Ts:4.44)';  
y = step(sys_lin,t);
```

```
data = iddata([zeros(20,2);y],[zeros(20,1); ones(101,1)],Ts);
```

data is an iddata object that encapsulates the step response of sys\_lin.

Compare the data to the model linearization.

```
compare(data,sys_lin);
```



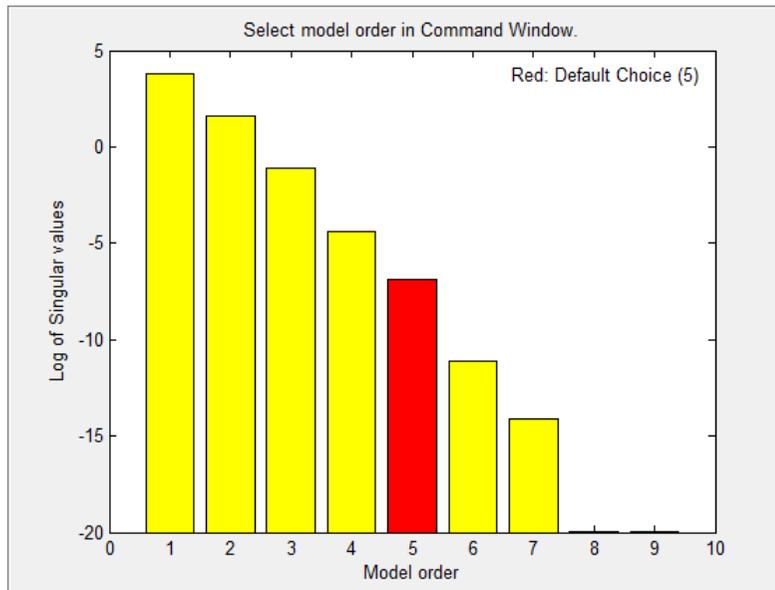
Because the data was obtained by simulating the linearized model, there is a 100% match between the data and model linearization response.

Identify a state-space model with a reduced order that adequately fits the data.

Determine an optimal model order.

```
nx = 1:9;  
sys1 = ssest(data,nx,'DisturbanceModel','none');
```

A plot showing the Hankel singular values (SVD) for models of the orders specified by `nx` appears.

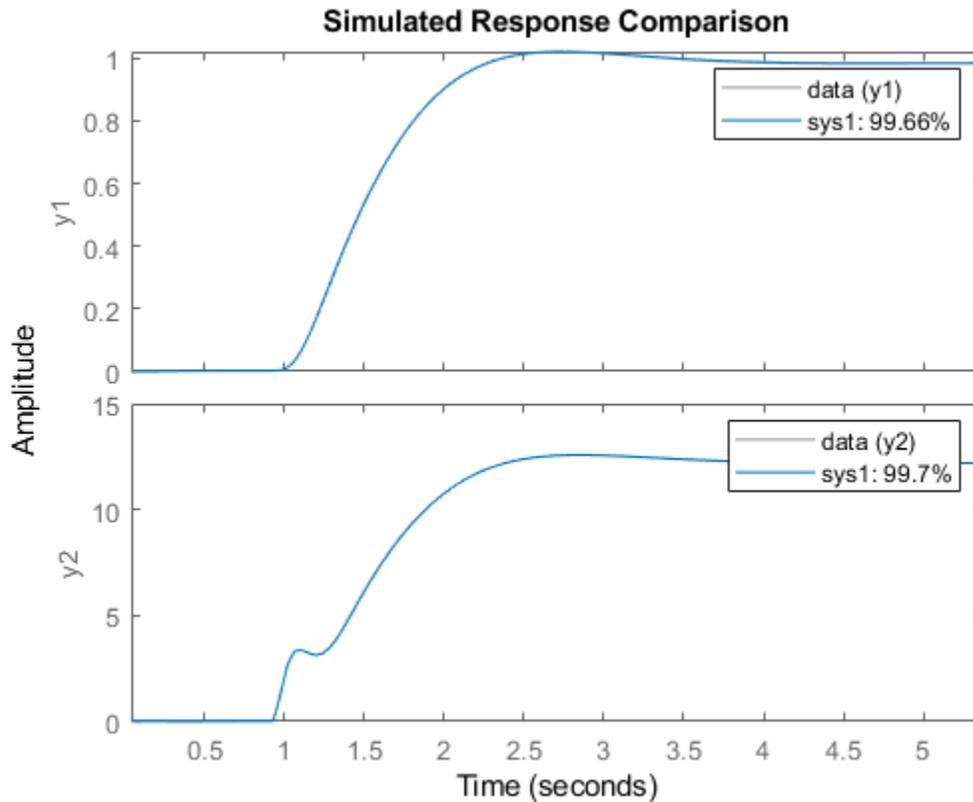


States with relatively small Hankel singular values can be safely discarded. The plot suggests using a fifth-order model.

At the MATLAB command prompt, select the model order for the estimated state-space model. Specify the model order as 5, or press **Enter** to use the default order value.

Compare the data to the estimated model.

```
compare(data,sys1);
```



sys1 provides a 98.4% fit for the first output and a 97.7% fit for the second output.

Examine the stopping condition for the search algorithm.

```
sys1.Report.Termination.WhyStop
```

```
ans =  
'Maximum number of iterations reached.'
```

Create an estimation options set that specifies the 'lm' search method and allows a maximum of 50 search iterations.

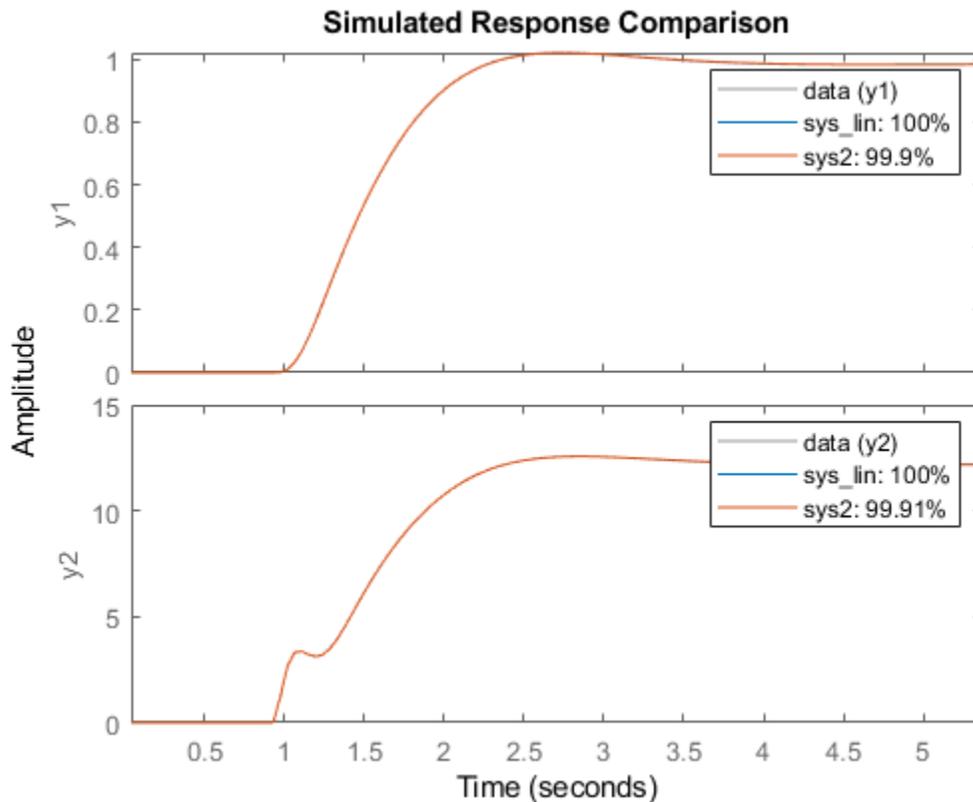
```
opt = ssestOptions('SearchMethod','lm');  
opt.SearchOptions.MaxIterations = 50;  
opt.Display = 'on';
```

Identify a state-space model using the estimation option set and `sys1` as the estimation initialization model.

```
sys2 = ssest(data,sys1,opt);
```

Compare the response of the linearized and the estimated models.

```
compare(data,sys_lin,sys2);
```



`sys2` provides a 99% fit for the first output and a 98% fit for the second output while using 4 less states than `sys_lin`.

## Definitions

### Modal Form

In modal form,  $A$  is a block-diagonal matrix. The block size is typically 1-by-1 for real eigenvalues and 2-by-2 for complex eigenvalues. However, if there are repeated eigenvalues or clusters of nearby eigenvalues, the block size can be larger.

For example, for a system with eigenvalues  $(\lambda_1, \sigma \pm j\omega, \lambda_2)$ , the modal  $A$  matrix is of the form

$$\begin{bmatrix} \lambda_1 & 0 & 0 & 0 \\ 0 & \sigma & \omega & 0 \\ 0 & -\omega & \sigma & 0 \\ 0 & 0 & 0 & \lambda_2 \end{bmatrix}$$

### Companion Form

In the companion realization, the characteristic polynomial of the system appears explicitly in the right-most column of the  $A$  matrix.

For a system with characteristic polynomial

$$p(s) = s^n + \alpha_1 s^{n-1} + \dots + \alpha_{n-1} s + \alpha_n$$

the corresponding companion  $A$  matrix is

$$A = \begin{bmatrix} 0 & 0 & \dots & \dots & 0 & -\alpha_n \\ 1 & 0 & 0 & \dots & 0 & -\alpha_{n-1} \\ 0 & 1 & 0 & \dots & \vdots & \vdots \\ \vdots & 0 & \dots & \dots & \vdots & \vdots \\ 0 & \dots & \dots & 1 & 0 & -\alpha_2 \\ 0 & \dots & \dots & 0 & 1 & -\alpha_1 \end{bmatrix}$$

The companion transformation requires that the system be controllable from the first input. The companion form is poorly conditioned for most state-space computations; avoid using it when possible.

## Algorithms

`ssest` initializes the parameter estimates using a noniterative subspace approach. It then refines the parameter values using the prediction error minimization approach. See `pem` for more information.

## References

- [1] Ljung, L. *System Identification: Theory For the User*, Second Edition, Upper Saddle River, N.J.: Prentice Hall, 1999.

## Extended Capabilities

### Automatic Parallel Support

Accelerate code by automatically running computation in parallel using Parallel Computing Toolbox™.

Parallel computing support is available for estimation using the `lsqnonlin` search method (requires Optimization Toolbox). To enable parallel computing, use `ssestOptions`, set `SearchMethod` to `'lsqnonlin'`, and set `SearchOptions.Advanced.UseParallel` to `true`.

For example:

```
opt = ssestOptions;  
opt.SearchMethod = 'lsqnonlin';  
opt.SearchOptions.Advanced.UseParallel = true;
```

## See Also

canon | iddata | idfrd | idgrey | idss | n4sid | pem | polyest | procest |  
ssestOptions | ssregest | tfest

## Topics

*“Estimate State-Space Models at the Command Line”*

*“Estimate State-Space Models with Free-Parameterization”*

*“Estimate State-Space Models with Canonical Parameterization”*

*“Estimate State-Space Models with Structured Parameterization”*

*“What Are State-Space Models?”*

*“Supported State-Space Parameterizations”*

*“State-Space Model Estimation Methods”*

*“Regularized Estimates of Model Parameters”*

*“Estimating Models Using Frequency-Domain Data”*

## Introduced in R2012a

## ssestOptions

Option set for ssest

### Syntax

```
opt = ssestOptions  
opt = ssestOptions(Name,Value)
```

### Description

`opt = ssestOptions` creates the default option set for `ssest`.

`opt = ssestOptions(Name,Value)` creates an option set with the options specified by one or more `Name,Value` pair arguments.

### Input Arguments

#### Name-Value Pair Arguments

Specify optional comma-separated pairs of `Name,Value` arguments. `Name` is the argument name and `Value` is the corresponding value. `Name` must appear inside quotes. You can specify several name and value pair arguments in any order as `Name1,Value1,...,NameN,ValueN`.

#### InitializeMethod — Algorithm used to initialize the state-space parameters

```
'auto' (default) | 'n4sid' | 'lsrf'
```

Algorithm used to initialize the state-space parameter values for `ssest`, specified as one of the following values:

- `'auto'` — `ssest` selects automatically:
  - `lsrf`, if the system is non-MIMO, the data is frequency-domain, and the state-space parameters are real-valued.

- `n4sid` otherwise (time-domain, MIMO, or with complex-valued state-space parameters).
- `'n4sid'` — Subspace state-space estimation approach — can be used with all systems (see `n4sid`).
- `'lsrf'` — Least-squares rational function estimation-based approach [7] (see “Continuous-Time Transfer Function Estimation Using Continuous-Time Frequency-Domain Data” on page 1-1796) — can provide higher-accuracy results for non-MIMO frequency-domain systems with real-valued state-space parameters, but cannot be used for any other systems (time-domain, MIMO, or with complex-valued state-space parameters).

### InitialState — Handling of initial states

`'auto'` (default) | `'zero'` | `'estimate'` | `'backcast'` | vector | parametric initial condition object (`x0obj`)

Handling of initial states during estimation, specified as one of the following values:

- `'zero'` — The initial state is set to zero.
- `'estimate'` — The initial state is treated as an independent estimation parameter.
- `'backcast'` — The initial state is estimated using the best least squares fit.
- `'auto'` — `ssest` chooses the initial state handling method, based on the estimation data. The possible initial state handling methods are `'zero'`, `'estimate'` and `'backcast'`.
- Vector of doubles — Specify a column vector of length  $N_x$ , where  $N_x$  is the number of states. For multi-experiment data, specify a matrix with  $N_e$  columns, where  $N_e$  is the number of experiments. The specified values are treated as fixed values during the estimation process.
- Parametric initial condition object (`x0obj`) — Specify initial conditions by using `idpar` to create a parametric initial condition object. You can specify minimum/maximum bounds and fix the values of specific states using the parametric initial condition object. The free entries of `x0obj` are estimated together with the `idss` model parameters.

Use this option only for discrete-time state-space models.

### N4Weight — Weighting scheme used for singular-value decomposition by the N4SID algorithm

`'auto'` (default) | `'MOESP'` | `'CVA'` | `'SSARX'`

Weighting scheme used for singular-value decomposition by the N4SID algorithm, specified as one of the following values:

- 'MOESP' — Uses the MOESP algorithm by Verhaegen [2].
- 'CVA' — Uses the Canonical Variable Algorithm by Larimore [1].
- 'SSARX' — A subspace identification method that uses an ARX estimation based algorithm to compute the weighting.

Specifying this option allows unbiased estimates when using data that is collected in closed-loop operation. For more information about the algorithm, see [6].

- 'auto' — The estimating function chooses between the MOESP and CVA algorithms.

### **N4Horizon — Forward- and backward-prediction horizons used by the N4SID algorithm**

'auto' (default) | vector [r sy su] | k-by-3 matrix

Forward and backward prediction horizons used by the N4SID algorithm, specified as one of the following values:

- A row vector with three elements — [r sy su], where r is the maximum forward prediction horizon. The algorithm uses up to r step-ahead predictors. sy is the number of past outputs, and su is the number of past inputs that are used for the predictions. See pages 209 and 210 in [4] for more information. These numbers can have a substantial influence on the quality of the resulting model, and there are no simple rules for choosing them. Making 'N4Horizon' a k-by-3 matrix means that each row of 'N4Horizon' is tried, and the value that gives the best (prediction) fit to data is selected. k is the number of guesses of [r sy su] combinations. If you specify N4Horizon as a single column, r = sy = su is used.
- 'auto' — The software uses an Akaike Information Criterion (AIC) for the selection of sy and su.

### **Focus — Error to be minimized**

'prediction' (default) | 'simulation'

Error to be minimized in the loss function during estimation, specified as the comma-separated pair consisting of 'Focus' and one of the following values:

- 'prediction' — The one-step ahead prediction error between measured and predicted outputs is minimized during estimation. As a result, the estimation focuses on producing a good predictor model.

- 'simulation' — The simulation error between measured and simulated outputs is minimized during estimation. As a result, the estimation focuses on making a good fit for simulation of model response with the current inputs.

The `Focus` option can be interpreted as a weighting filter in the loss function. For more information, see “Loss Function and Model Quality Metrics”.

### **WeightingFilter — Weighting prefilter**

[ ] (default) | vector | matrix | cell array | linear system

Weighting prefilter applied to the loss function to be minimized during estimation. To understand the effect of `WeightingFilter` on the loss function, see “Loss Function and Model Quality Metrics”.

Specify `WeightingFilter` as one of the following values:

- [ ] — No weighting prefilter is used.
- Passbands — Specify a row vector or matrix containing frequency values that define desired passbands. You select a frequency band where the fit between estimated model and estimation data is optimized. For example, `[wl,wh]` where `wl` and `wh` represent lower and upper limits of a passband. For a matrix with several rows defining frequency passbands, `[w1l,w1h;w2l,w2h;w3l,w3h;...]`, the estimation algorithm uses the union of the frequency ranges to define the estimation passband.

Passbands are expressed in `rad/TimeUnit` for time-domain data and in `FrequencyUnit` for frequency-domain data, where `TimeUnit` and `FrequencyUnit` are the time and frequency units of the estimation data.

- SISO filter — Specify a single-input-single-output (SISO) linear filter in one of the following ways:
  - A SISO LTI model
  - `{A,B,C,D}` format, which specifies the state-space matrices of a filter with the same sample time as estimation data.
  - `{numerator,denominator}` format, which specifies the numerator and denominator of the filter as a transfer function with same sample time as estimation data.

This option calculates the weighting function as a product of the filter and the input spectrum to estimate the transfer function.

- Weighting vector — Applicable for frequency-domain data only. Specify a column vector of weights. This vector must have the same length as the frequency vector of

the data set, `Data.Frequency`. Each input and output response in the data is multiplied by the corresponding weight at that frequency.

- `'invsqrt'` — Applicable for frequency-domain data only, with `InitializeMethod` set to `'lsrf'` only. Uses  $1/\sqrt{|G(\omega)|}$  as the weighting filter, where  $G(\omega)$  is the complex frequency-response data. Use this option for capturing relatively low amplitude dynamics in data.
- `'inv'` — Applicable for frequency-domain data only, with `InitializeMethod` set to `'lsrf'` only. Uses  $1/|G(\omega)|$  as the weighting filter. Similarly to `'invsqrt'`, this option captures relatively low-amplitude dynamics in data. Use it when `'invsqrt'` weighting produces an estimate that is missing dynamics in the low-amplitude regions. `'inv'` is more sensitive to noise than `'invsqrt'`.

### **EnforceStability — Control whether to enforce stability of model**

`false` (default) | `true`

Control whether to enforce stability of estimated model, specified as the comma-separated pair consisting of `'EnforceStability'` and either `true` or `false`.

Data Types: `logical`

### **EstimateCovariance — Control whether to generate parameter covariance data**

`true` (default) | `false`

Controls whether parameter covariance data is generated, specified as `true` or `false`.

If `EstimateCovariance` is `true`, then use `getcov` to fetch the covariance matrix from the estimated model.

### **Display — Specify whether to display the estimation progress**

`'off'` (default) | `'on'`

Specify whether to display the estimation progress, specified as one of the following values:

- `'on'` — Information on model structure and estimation results are displayed in a progress-viewer window.
- `'off'` — No progress or results information is displayed.

### **InputOffset — Removal of offset from time-domain input data during estimation**

[ ] (default) | vector of positive integers | matrix

Removal of offset from time-domain input data during estimation, specified as the comma-separated pair consisting of 'InputOffset' and one of the following:

- A column vector of positive integers of length  $Nu$ , where  $Nu$  is the number of inputs.
- [ ] — Indicates no offset.
- $Nu$ -by- $Ne$  matrix — For multi-experiment data, specify InputOffset as an  $Nu$ -by- $Ne$  matrix.  $Nu$  is the number of inputs, and  $Ne$  is the number of experiments.

Each entry specified by InputOffset is subtracted from the corresponding input data.

### **OutputOffset — Removal of offset from time-domain output data during estimation**

[ ] (default) | vector | matrix

Removal of offset from time-domain output data during estimation, specified as the comma-separated pair consisting of 'OutputOffset' and one of the following:

- A column vector of length  $Ny$ , where  $Ny$  is the number of outputs.
- [ ] — Indicates no offset.
- $Ny$ -by- $Ne$  matrix — For multi-experiment data, specify OutputOffset as a  $Ny$ -by- $Ne$  matrix.  $Ny$  is the number of outputs, and  $Ne$  is the number of experiments.

Each entry specified by OutputOffset is subtracted from the corresponding output data.

### **OutputWeight — Weighting of prediction errors in multi-output estimations**

[ ] (default) | 'noise' | positive semidefinite symmetric matrix

Weighting of prediction errors in multi-output estimations, specified as one of the following values:

- 'noise' — Minimize  $\det(E^*E / N)$ , where  $E$  represents the prediction error and  $N$  is the number of data samples. This choice is optimal in a statistical sense and leads to maximum likelihood estimates if nothing is known about the variance of the noise. It uses the inverse of the estimated noise variance as the weighting function.

**Note** OutputWeight must not be 'noise' if SearchMethod is 'lsqnonlin'.

- Positive semidefinite symmetric matrix ( $W$ ) — Minimize the trace of the weighted prediction error matrix  $\text{trace}(E' * E * W / N)$  where:
  - $E$  is the matrix of prediction errors, with one column for each output, and  $W$  is the positive semidefinite symmetric matrix of size equal to the number of outputs. Use  $W$  to specify the relative importance of outputs in multiple-output models, or the reliability of corresponding data.
  - $N$  is the number of data samples.
- `[]` — The software chooses between the 'noise' or using the identity matrix for  $W$ .

This option is relevant for only multi-output models.

## **Regularization — Options for regularized estimation of model parameters** structure

Options for regularized estimation of model parameters. For more information on regularization, see “Regularized Estimates of Model Parameters”.

Regularization is a structure with the following fields:

- **Lambda** — Constant that determines the bias versus variance tradeoff.

Specify a positive scalar to add the regularization term to the estimation cost.

The default value of zero implies no regularization.

**Default:** 0

- **R** — Weighting matrix.

Specify a vector of nonnegative numbers or a square positive semi-definite matrix. The length must be equal to the number of free parameters of the model.

For black-box models, using the default value is recommended. For structured and grey-box models, you can also specify a vector of `np` positive numbers such that each entry denotes the confidence in the value of the associated parameter.

The default value of 1 implies a value of `eye(npfree)`, where `npfree` is the number of free parameters.

**Default:** 1

- **Nominal** — The nominal value towards which the free parameters are pulled during estimation.

The default value of zero implies that the parameter values are pulled towards zero. If you are refining a model, you can set the value to `'model'` to pull the parameters towards the parameter values of the initial model. The initial parameter values must be finite for this setting to work.

**Default:** 0

### **SearchMethod** — Numerical search method used for iterative parameter estimation

`'auto'` (default) | `'gn'` | `'gna'` | `'lm'` | `'grad'` | `'lsqnonlin'` | `'fmincon'`

Numerical search method used for iterative parameter estimation, specified as the comma-separated pair consisting of `'SearchMethod'` and one of the following:

- `'auto'` — A combination of the line search algorithms, `'gn'`, `'lm'`, `'gna'`, and `'grad'` methods is tried in sequence at each iteration. The first descent direction leading to a reduction in estimation cost is used.
- `'gn'` — Subspace Gauss-Newton least squares search. Singular values of the Jacobian matrix less than  $GnPinvConstant * eps * \max(\text{size}(J)) * \text{norm}(J)$  are discarded when computing the search direction.  $J$  is the Jacobian matrix. The Hessian matrix is approximated as  $J^T J$ . If there is no improvement in this direction, the function tries the gradient direction.
- `'gna'` — Adaptive subspace Gauss-Newton search. Eigenvalues less than  $gamma * \max(sv)$  of the Hessian are ignored, where  $sv$  contains the singular values of the Hessian. The Gauss-Newton direction is computed in the remaining subspace.  $gamma$  has the initial value `InitialGnaTolerance` (see `Advanced in 'SearchOptions'` for more information). This value is increased by the factor `LMStep` each time the search fails to find a lower value of the criterion in fewer than five bisections. This value is decreased by the factor  $2 * LMStep$  each time a search is successful without any bisections.
- `'lm'` — Levenberg-Marquardt least squares search, where the next parameter value is  $-pinv(H+d*I) * grad$  from the previous one.  $H$  is the Hessian,  $I$  is the identity matrix, and  $grad$  is the gradient.  $d$  is a number that is increased until a lower value of the criterion is found.
- `'grad'` — Steepest descent least squares search.
- `'lsqnonlin'` — Trust-region-reflective algorithm of `lsqnonlin`. Requires Optimization Toolbox software.

- 'fmincon' — Constrained nonlinear solvers. You can use the sequential quadratic programming (SQP) and trust-region-reflective algorithms of the fmincon solver. If you have Optimization Toolbox software, you can also use the interior-point and active-set algorithms of the fmincon solver. Specify the algorithm in the SearchOptions.Algorithm option. The fmincon algorithms may result in improved estimation results in the following scenarios:
  - Constrained minimization problems when there are bounds imposed on the model parameters.
  - Model structures where the loss function is a nonlinear or non smooth function of the parameters.
  - Multi-output model estimation. A determinant loss function is minimized by default for multi-output model estimation. fmincon algorithms are able to minimize such loss functions directly. The other search methods such as 'lm' and 'gn' minimize the determinant loss function by alternately estimating the noise variance and reducing the loss value for a given noise variance value. Hence, the fmincon algorithms can offer better efficiency and accuracy for multi-output model estimations.

### **SearchOptions — Option set for the search algorithm**

search option set

Option set for the search algorithm, specified as the comma-separated pair consisting of 'SearchOptions' and a search option set with fields that depend on the value of SearchMethod.

**SearchOptions Structure When SearchMethod is Specified as 'gn', 'gna', 'lm', 'grad', or 'auto'**

Field Name	Description	Default
Tolerance	Minimum percentage difference between the current value of the loss function and its expected improvement after the next iteration, specified as a positive scalar. When the percentage of expected improvement is less than <code>Tolerance</code> , the iterations stop. The estimate of the expected loss-function improvement at the next iteration is based on the Gauss-Newton vector computed for the current parameter value.	0.01
MaxIterations	<p>Maximum number of iterations during loss-function minimization, specified as a positive integer. The iterations stop when <code>MaxIterations</code> is reached or another stopping criterion is satisfied, such as <code>Tolerance</code>.</p> <p>Setting <code>MaxIterations = 0</code> returns the result of the start-up procedure.</p> <p>Use <code>sys.Report.Termination.Iterations</code> to get the actual number of iterations during an estimation, where <code>sys</code> is an <code>idtf</code> model.</p>	20

Field Name	Description	Default
Advanced	Advanced search settings, specified as a structure with the following fields:	
<b>Field Name</b>	<b>Description</b>	<b>Default</b>
GnPinvConstant	Jacobian matrix singular value threshold, specified as a positive scalar. Singular values of the Jacobian matrix that are smaller than $\text{GnPinvConstant} * \max(\text{size}(J) * \text{norm}(J) * \text{eps})$ are discarded when computing the search direction. Applicable when SearchMethod is 'gn'.	10000
InitialGnaTolerance	Initial value of $\gamma$ , specified as a positive scalar. Applicable when SearchMethod is 'gna'.	0.0001
LMStartValue	Starting value of search-direction length $d$ in the Levenberg-Marquardt method, specified as a positive scalar. Applicable when SearchMethod is 'lm'.	0.001
LMStep	Size of the Levenberg-Marquardt step, specified as a positive integer. The next value of the search-direction length $d$ in the Levenberg-Marquardt method is LMStep times the previous one. Applicable when SearchMethod is 'lm'.	2
MaxBisections	Maximum number of bisections used for line search along the search direction, specified as a positive integer.	25
MaxFunctionEvaluations	Maximum number of calls to the model file, specified as a positive integer. Iterations stop if the number of calls to the model file exceeds this value.	Inf
MinParameterChange	Smallest parameter update allowed per iteration, specified as a nonnegative scalar.	0
RelativeImprovement	Relative improvement threshold, specified as a nonnegative scalar. Iterations stop if the relative	0

Field Name	Description		Default
	<b>Field Name</b>	<b>Description</b>	<b>Default</b>
		improvement of the criterion function is less than this value.	
	StepReduction	<p>Step reduction factor, specified as a positive scalar that is greater than 1. The suggested parameter update is reduced by the factor <code>StepReduction</code> after each try. This reduction continues until <code>MaxBisections</code> tries are completed or a lower value of the criterion function is obtained.</p> <p><code>StepReduction</code> is not applicable for <code>SearchMethod 'lm'</code> (Levenberg-Marquardt method).</p>	2

**SearchOptions Structure When SearchMethod is Specified as 'lsqnonlin'**

<b>Field Name</b>	<b>Description</b>	<b>Default</b>
FunctionTolerance	Termination tolerance on the loss function that the software minimizes to determine the estimated parameter values, specified as a positive scalar.  The value of FunctionTolerance is the same as that of <code>opt.SearchOptions.Advanced.TolFun</code> .	1e-5
StepTolerance	Termination tolerance on the estimated parameter values, specified as a positive scalar.  The value of StepTolerance is the same as that of <code>opt.SearchOptions.Advanced.TolX</code> .	1e-6
MaxIterations	Maximum number of iterations during loss-function minimization, specified as a positive integer. The iterations stop when MaxIterations is reached or another stopping criterion is satisfied, such as FunctionTolerance.  The value of MaxIterations is the same as that of <code>opt.SearchOptions.Advanced.MaxIter</code> .	20
Advanced	Advanced search settings, specified as an option set for <code>lsqnonlin</code> .  For more information, see the Optimization Options table in "Optimization Options" (Optimization Toolbox).	Use <code>optimset('lsqnonlin')</code> to create a default option set.

### SearchOptions Structure When SearchMethod is Specified as 'fmincon'

Field Name	Description	Default
Algorithm	<p>fmincon optimization algorithm, specified as one of the following:</p> <ul style="list-style-type: none"> <li>'sqp' — Sequential quadratic programming algorithm. The algorithm satisfies bounds at all iterations, and it can recover from NaN or Inf results. It is not a large-scale algorithm. For more information, see “Large-Scale vs. Medium-Scale Algorithms” (Optimization Toolbox).</li> <li>'trust-region-reflective' — Subspace trust-region method based on the interior-reflective Newton method. It is a large-scale algorithm.</li> <li>'interior-point' — Large-scale algorithm that requires Optimization Toolbox software. The algorithm satisfies bounds at all iterations, and it can recover from NaN or Inf results.</li> <li>'active-set' — Requires Optimization Toolbox software. The</li> </ul>	'sqp'

Field Name	Description	Default
	<p>algorithm can take large steps, which adds speed. It is not a large-scale algorithm.</p> <p>For more information about the algorithms, see “Constrained Nonlinear Optimization Algorithms” (Optimization Toolbox) and “Choosing the Algorithm” (Optimization Toolbox).</p>	
FunctionTolerance	Termination tolerance on the loss function that the software minimizes to determine the estimated parameter values, specified as a positive scalar.	1e-6
StepTolerance	Termination tolerance on the estimated parameter values, specified as a positive scalar.	1e-6
MaxIterations	Maximum number of iterations during loss function minimization, specified as a positive integer. The iterations stop when MaxIterations is reached or another stopping criterion is satisfied, such as FunctionTolerance.	100

**Advanced – Additional advanced options**  
structure

Additional advanced options, specified as a structure with the following fields:

- `ErrorThreshold` — Specifies when to adjust the weight of large errors from quadratic to linear.

Errors larger than `ErrorThreshold` times the estimated standard deviation have a linear weight in the loss function. The standard deviation is estimated robustly as the median of the absolute deviations from the median of the prediction errors, divided by 0.7. For more information on robust norm choices, see section 15.2 of [4].

`ErrorThreshold = 0` disables robustification and leads to a purely quadratic loss function. When estimating with frequency-domain data, the software sets `ErrorThreshold` to zero. For time-domain data that contains outliers, try setting `ErrorThreshold` to 1.6.

**Default:** 0

- `MaxSize` — Specifies the maximum number of elements in a segment when input-output data is split into segments.

`MaxSize` must be a positive integer.

**Default:** 250000

- `StabilityThreshold` — Specifies thresholds for stability tests.

`StabilityThreshold` is a structure with the following fields:

- `s` — Specifies the location of the right-most pole to test the stability of continuous-time models. A model is considered stable when its right-most pole is to the left of `s`.

**Default:** 0

- `z` — Specifies the maximum distance of all poles from the origin to test stability of discrete-time models. A model is considered stable if all poles are within the distance `z` from the origin.

**Default:** `1+sqrt(eps)`

- `AutoInitThreshold` — Specifies when to automatically estimate the initial conditions.

The initial condition is estimated when

$$\frac{\|y_{p,z} - y_{meas}\|}{\|y_{p,e} - y_{meas}\|} > \text{AutoInitThreshold}$$

- $y_{meas}$  is the measured output.
- $y_{p,z}$  is the predicted output of a model estimated using zero initial states.
- $y_{p,e}$  is the predicted output of a model estimated using estimated initial states.

Applicable when `InitialState` is 'auto'.

**Default:** 1.05

- DDC — Specifies if the Data Driven Coordinates algorithm [5] is used to estimate freely parameterized state-space models.

Specify DDC as one of the following values:

- 'on' — The free parameters are projected to a reduced space of identifiable parameters using the Data Driven Coordinates algorithm.
- 'off' — All the entries of  $A$ ,  $B$ , and  $C$  updated directly using the chosen `SearchMethod`.

**Default:** 'on'

## Output Arguments

**opt** — Option set for `ssest`

`ssestOptions` option set

Option set for `ssest`, returned as an `ssestOptions` option set.

## Examples

**Create Default Option Set for State Space Estimation**

```
opt = ssestOptions;
```

## Specify Options for State Space Estimation

Create an option set for `ssest` using the 'backcast' algorithm to initialize the state and set the `Display` to 'on'.

```
opt = ssestOptions('InitialState','backcast','Display','on');
```

Alternatively, use dot notation to set the values of `opt`.

```
opt = ssestOptions;  
opt.InitialState = 'backcast';  
opt.Display = 'on';
```

## References

- [1] Larimore, W.E. "Canonical variate analysis in identification, filtering and adaptive control." *Proceedings of the 29th IEEE Conference on Decision and Control*, pp. 596-604, 1990.
- [2] Verhaegen, M. "Identification of the deterministic part of MIMO state space models." *Automatica*, Vol. 30, No. 1, 1994, pp. 61-74.
- [3] Wills, Adrian, B. Ninness, and S. Gibson. "On Gradient-Based Search for Multivariable System Estimates." *Proceedings of the 16th IFAC World Congress, Prague, Czech Republic, July 3-8, 2005*. Oxford, UK: Elsevier Ltd., 2005.
- [4] Ljung, L. *System Identification: Theory for the User*. Upper Saddle River, NJ: Prentice-Hall PTR, 1999.
- [5] McKelvey, T., A. Helmersson, and T. Ribarits. "Data driven local coordinates for multivariable linear systems and their application to system identification." *Automatica*, Volume 40, No. 9, 2004, pp. 1629-1635.
- [6] Jansson, M. "Subspace identification and ARX modeling." *13th IFAC Symposium on System Identification*, Rotterdam, The Netherlands, 2003.
- [7] Ozdemir, A. A., and S. Gumossoy. "Transfer Function Estimation in System identification Toolbox via Vector Fitting." *Proceedings of the 20th World Congress of the International Federation of Automatic Control*. Toulouse, France, July 2017.

## **See Also**

ssest

## **Topics**

“Loss Function and Model Quality Metrics”

**Introduced in R2012a**

# ssform

Quick configuration of state-space model structure

## Syntax

```
sys1 = ssform(sys,Name,Value)
```

## Description

`sys1 = ssform(sys,Name,Value)` specifies the type of parameterization and whether feedthrough and disturbance dynamics are present for the state-space model `sys` using one or more `Name,Value` pair arguments.

## Input Arguments

### `sys`

State-space model

### Default:

## Name-Value Pair Arguments

Specify comma-separated pairs of `Name,Value` arguments, where `Name` is the argument name and `Value` is the corresponding value. `Name` must appear inside single quotes ( ' '). You can specify several name and value pair arguments in any order as `Name1,Value1,...,NameN,ValueN`.

### Form

Specify structure of A, B and C matrices as one of the following values:

- 'free'

All entries of A, B, C are set free

- 'companion'

Companion form of the model where the characteristic polynomial appears in the far-right column of the state matrix  $A$

- 'modal'

Modal decomposition form, where the state matrix  $A$  is block diagonal. Each block corresponds to a real or complex-conjugate pair of poles.

You cannot use this value for models with repeated poles.

- 'canonical'

Observability canonical form of  $A$ ,  $B$ , and  $C$  matrices, as described in [1].

### **Default:**

### **Feedthrough**

Specify whether the model has direct feedthrough from the input  $u(t)$  to the output  $y(t)$ , (whether the elements of the matrix  $D$  are nonzero).

Must be a logical vector (`true` or `false`) of length equal to the number of inputs ( $Nu$ ).

`Feedthrough(i) = false` sets `sys.Structure.D.Value(:,i)` to zero and `sys.Structure.D.Free(:,i)` to `false`.

`Feedthrough(i) = true` sets `sys.Structure.D.Free(:,i)` to `true`.

---

**Note** Specifying this option for a previously estimated model causes the model parameter covariance information to be lost. Use `translatecov` to recompute the covariance.

---

### **Default:**

### **DisturbanceModel**

Specify whether to estimate the noise component of the model, specified as one of the following values:

- 'none'

The value of the  $K$  matrix is fixed to zero.

- 'estimate'

The K matrix is treated as a free parameter

---

**Note** Specifying this option for a previously estimated model causes the model parameter covariance information to be lost. Use `translatecov` to recompute the covariance.

---

**Default:**

## Output Arguments

**sys1**

State-space model with configured parameterization, feedthrough, and disturbance dynamics

## Examples

### Convert State-Space Model to Canonical Form

Create a state-space model.

```
rng('default');  
A = randn(2) - 2*eye(2);  
B = randn(2,1);  
C = randn(1,2);  
D = 0;  
K = randn(2,1);  
model = idss(A,B,C,D,K,'Ts',0);
```

The state-space model has free parameterization and no feedthrough.

Convert the model to observability canonical form.

```
model1 = ssform(model,'Form','canonical');
```

## Estimate State-Space Model Parameters in Canonical Form with Feedthrough

Load the estimation data.

```
load iddata1 z1;
```

Create a state-space model.

```
rng('default');  
A = randn(2) - 2*eye(2);  
B = randn(2,1);  
C = randn(1,2);  
D = 0;  
K = randn(2,1);  
model = idss(A,B,C,D,K,'Ts',0);
```

The state-space model has free parameterization and no feedthrough.

Convert the model to observability canonical form and specify to estimate its feedthrough behavior.

```
model1 = ssform(model,'Form','canonical','Feedthrough', true);
```

Estimate the parameters of the model.

```
model2 = ssest(z1,model1);
```

## Alternatives

Use the `Structure` property of an `idss` model to specify the parameterization, feedthrough, and disturbance dynamics by modifying the `Value` and `Free` attributes of the `A`, `B`, `C`, `D` and `K` parameters.

## References

- [1] Ljung, L. *System Identification: Theory For the User*, Second Edition, Appendix 4A, pp 132-134, Upper Saddle River, N.J: Prentice Hall, 1999.

## See Also

`idss` | `n4sid` | `ssest`

## Topics

“Estimate State-Space Models at the Command Line”

“Supported State-Space Parameterizations”

**Introduced in R2012b**

## ssregest

Estimate state-space model by reduction of regularized ARX model

### Syntax

```
sys = ssregest(data,nx)
sys = ssregest(data,nx,Name,Value)
sys = ssregest( ___,opt)
```

```
[sys,x0] = ssregest( ___)
```

### Description

`sys = ssregest(data,nx)` estimates a state-space model by reduction of a regularized ARX model.

`sys = ssregest(data,nx,Name,Value)` specifies additional options using one or more `Name,Value` pair arguments.

`sys = ssregest( ___,opt)` specifies estimation options that configure the estimation objective, ARX orders, and order reduction options. This syntax can include any of the input argument combinations in the previous syntaxes.

`[sys,x0] = ssregest( ___)` returns the value of initial states computed during estimation. This syntax can include any of the input argument combinations in the previous syntaxes.

### Examples

#### Estimate State-Space Model by Reduction of Regularized ARX Model

Load estimation data.

```
load iddata2 z2;
```

z2 is an `iddata` object that contains time-domain system response data.

Identify a third-order state-space model.

```
sys = ssregest(z2,3);
```

### **Estimate State-Space Model With Input Delay**

Load estimation data.

```
load iddata2 z2
```

Estimate a third-order state-space model with input delay.

```
sys = ssregest(z2,3,'InputDelay',2);
```

### **Configure the ARX Orders and Estimation Focus**

Load estimation data.

```
load iddata2 z2;
```

Specify the order of the regularized ARX model used by the software during estimation. Also, set the estimation focus to simulation.

```
opt = ssregestOptions('ARXOrder',[100 100 1],'Focus','simulation');
```

Identify a third-order state-space model.

```
sys = ssregest(z2,3,opt);
```

### **Return Initial State Values Computed During Estimation**

Load estimation data.

```
load iddata2 z2;
```

Obtain the initial state values when identifying a third-order state-space model.

```
[sys,x0] = ssregest(z2,3);
```

### **Compare Regularized State-Space Models Estimated Using Impulse Response and Reduction of ARX Models**

Load data.

```
load regularizationExampleData eData;
```

Create a transfer function model used for generating the estimation data (true system).

```
trueSys = idtf([0.02008 0.04017 0.02008],[1 -1.561 0.6414],1);
```

Obtain regularized impulse response (FIR) model.

```
opt = impulseestOptions('RegularizationKernel','DC');  
m0 = impulseest(eData,70,opt);
```

Convert the model into a state-space model and reduce the model order.

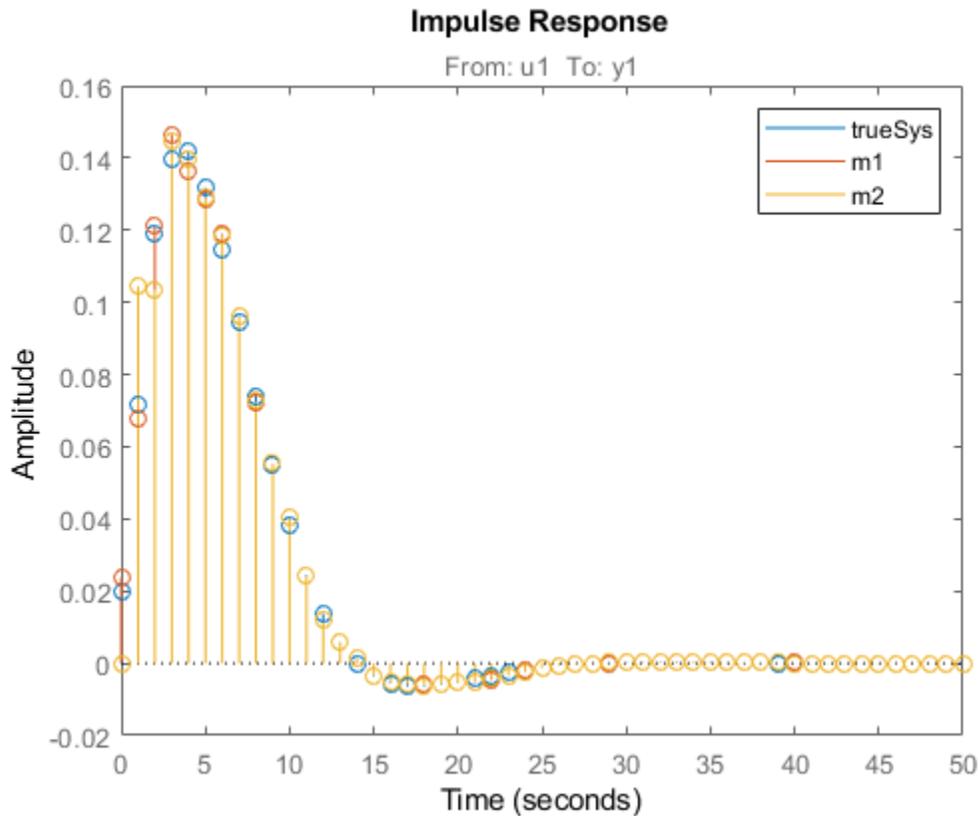
```
m1 = balred(idss(m0),15);
```

Obtain a second state-space model using regularized reduction of an ARX model.

```
m2 = ssregest(eData,15);
```

Compare the impulse responses of the true system and the estimated models.

```
impulse(trueSys,m1,m2,50);  
legend('trueSys','m1','m2');
```



## Input Arguments

### data — Estimation data

`iddata` | `idfrd` | `frd`

Estimation data, specified as an `iddata`, `idfrd` or `frd` object.

For time-domain estimation, data must be an `iddata` object containing the input and output signal values.

For frequency-domain estimation, data can be one of the following:

- Recorded frequency response data (`frd` or `idfrd`)
- `iddata` object with its properties specified as follows:
  - `InputData` — Fourier transform of the input signal
  - `OutputData` — Fourier transform of the output signal
  - `Domain` — 'Frequency'

The sample time `Ts` of the `iddata` object must be nonzero.

### **`nx` — Order of estimated model**

positive scalar | positive vector | 'best'

Order of the estimated model, specified as a positive scalar or vector.

If `nx` is a vector, then `ssregest` creates a plot which you can use to choose a suitable model order. The plot shows the Hankel singular values for models of chosen values in the vector. States with relatively small Hankel singular values can be safely discarded. A default choice is suggested in the plot.

You can also specify `nx = 'best'`, as in `ssregest(data, 'best')`, in which case the optimal order is chosen automatically in the 1:10 range.

### **`opt` — Options set for `ssregest`**

`ssregestOptions` options set

Estimation options for `ssregest`, specified as an options set you create using `ssregestOptions`.

## **Name-Value Pair Arguments**

Specify optional comma-separated pairs of `Name`, `Value` arguments. `Name` is the argument name and `Value` is the corresponding value. `Name` must appear inside quotes. You can specify several name and value pair arguments in any order as `Name1, Value1, ..., NameN, ValueN`.

Example: `sys = ssregest(z2, 3, 'InputDelay', 2)` specifies a delay of 2 sampling periods.

### **`Ts` — Sample time**

sample time of data (`data.Ts`) (default) | positive scalar | 0

Sample time of the model, specified as 0 or equal to the sample time of data.

For continuous-time models, use  $T_s = 0$ . For discrete-time models, specify  $T_s$  as a positive scalar whose value is equal to the data sample time.

### **InputDelay — Input delays**

0 (default) | scalar | vector

Input delay for each input channel, specified as a numeric vector. For continuous-time systems, specify input delays in the time unit stored in the `TimeUnit` property. For discrete-time systems, specify input delays in integer multiples of the sample time  $T_s$ . For example, `InputDelay = 3` means a delay of three sampling periods.

For a system with  $N_u$  inputs, set `InputDelay` to an  $N_u$ -by-1 vector. Each entry of this vector is a numerical value that represents the input delay for the corresponding input channel.

You can also set `InputDelay` to a scalar value to apply the same delay to all channels.

### **Form — Type of canonical form**

'free' (default) | 'modal' | 'companion' | 'canonical'

Type of canonical form of `sys`, specified as one of the following values:

- 'modal' — Obtain `sys` in modal form on page 1-1699.
- 'companion' — Obtain `sys` in companion form on page 1-1699.
- 'free' — All entries of the  $A$ ,  $B$  and  $C$  matrices are treated as free.
- 'canonical' — Obtain `sys` in the observability canonical form [1].

Use the `Form`, `Feedthrough` and `DisturbanceModel` name-value pair arguments to modify the default behavior of the  $A$ ,  $B$ ,  $C$ ,  $D$ , and  $K$  matrices.

### **Feedthrough — Direct feedthrough from input to output**

0 (default) | 1 | logical vector

Direct feedthrough from input to output, specified as a logical vector of length  $N_u$ , where  $N_u$  is the number of inputs. If `Feedthrough` is specified as a logical scalar, it is applied to all the inputs.

Use the `Form`, `Feedthrough` and `DisturbanceModel` name-value pair arguments to modify the default behavior of the  $A$ ,  $B$ ,  $C$ ,  $D$ , and  $K$  matrices.

## **DisturbanceModel — Specify whether to estimate the $K$ matrix**

'estimate' (default) | 'none'

Specify whether to estimate the  $K$  matrix which specifies the noise component, specified as one of the following values:

- 'none' — Noise component is not estimated. The value of the  $K$  matrix is fixed to zero value.
- 'estimate' — The  $K$  matrix is treated as a free parameter.

`DisturbanceModel` must be 'none' when using frequency-domain data.

Use the `Form`, `Feedthrough` and `DisturbanceModel` name-value pair arguments to modify the default behavior of the  $A$ ,  $B$ ,  $C$ ,  $D$ , and  $K$  matrices.

## **Output Arguments**

### **sys — Estimated state-space model**

`idss`

Estimated state-space model of order  $n_x$ , returned as an `idss` model object. The model represents:

$$\dot{x}(t) = Ax(t) + Bu(t) + Ke(t)$$

$$y(t) = Cx(t) + Du(t) + e(t)$$

$A$ ,  $B$ ,  $C$ ,  $D$ , and  $K$  are state-space matrices.  $u(t)$  is the input,  $y(t)$  is the output,  $e(t)$  is the disturbance and  $x(t)$  is the vector of  $n_x$  states.

All the entries of  $A$ ,  $B$ ,  $C$ , and  $K$  are free estimable parameters by default.  $D$  is fixed to zero by default, meaning that there is no feedthrough, except for static systems ( $n_x=0$ ).

Information about the estimation results and options used is stored in the `Report` property of the model. `Report` has the following fields:

Report Field	Description																		
Status	Summary of the model status, which indicates whether the model was created by construction or obtained by estimation.																		
Method	Estimation command used.																		
Initial State	<p>Handling of initial states during estimation, returned as one of the following values:</p> <ul style="list-style-type: none"> <li>• 'zero' — The initial state was set to zero.</li> <li>• 'estimate' — The initial state was treated as an independent estimation parameter.</li> </ul> <p>This field is especially useful when the InitialState option in the estimation option set is 'auto'.</p>																		
ARXOrder	ARX model orders, returned as a matrix of nonnegative integers [na nb nk].																		
Fit	<p>Quantitative assessment of the estimation, returned as a structure. See “Loss Function and Model Quality Metrics” for more information on these quality metrics. The structure has the following fields:</p> <table border="1" data-bbox="387 951 1338 1451"> <thead> <tr> <th data-bbox="387 951 516 994">Field</th> <th data-bbox="516 951 1338 994">Description</th> </tr> </thead> <tbody> <tr> <td data-bbox="387 994 516 1107">FitPercent</td> <td data-bbox="516 994 1338 1107">Normalized root mean squared error (NRMSE) measure of how well the response of the model fits the estimation data, expressed as a percentage.</td> </tr> <tr> <td data-bbox="387 1107 516 1150">LossFcn</td> <td data-bbox="516 1107 1338 1150">Value of the loss function when the estimation completes.</td> </tr> <tr> <td data-bbox="387 1150 516 1225">MSE</td> <td data-bbox="516 1150 1338 1225">Mean squared error (MSE) measure of how well the response of the model fits the estimation data.</td> </tr> <tr> <td data-bbox="387 1225 516 1269">FPE</td> <td data-bbox="516 1225 1338 1269">Final prediction error for the model.</td> </tr> <tr> <td data-bbox="387 1269 516 1312">AIC</td> <td data-bbox="516 1269 1338 1312">Raw Akaike Information Criteria (AIC) measure of model quality.</td> </tr> <tr> <td data-bbox="387 1312 516 1355">AICc</td> <td data-bbox="516 1312 1338 1355">Small sample-size corrected AIC.</td> </tr> <tr> <td data-bbox="387 1355 516 1399">nAIC</td> <td data-bbox="516 1355 1338 1399">Normalized AIC.</td> </tr> <tr> <td data-bbox="387 1399 516 1451">BIC</td> <td data-bbox="516 1399 1338 1451">Bayesian Information Criteria (BIC).</td> </tr> </tbody> </table>	Field	Description	FitPercent	Normalized root mean squared error (NRMSE) measure of how well the response of the model fits the estimation data, expressed as a percentage.	LossFcn	Value of the loss function when the estimation completes.	MSE	Mean squared error (MSE) measure of how well the response of the model fits the estimation data.	FPE	Final prediction error for the model.	AIC	Raw Akaike Information Criteria (AIC) measure of model quality.	AICc	Small sample-size corrected AIC.	nAIC	Normalized AIC.	BIC	Bayesian Information Criteria (BIC).
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BIC	Bayesian Information Criteria (BIC).																		
Parameters	Estimated values of model parameters.																		

Report Field	Description																
OptionsUsed	Option set used for estimation. If no custom options were configured, this is a set of default options. See <code>ssregestOptions</code> for more information.																
RandState	State of the random number stream at the start of estimation. Empty, <code>[]</code> , if randomization was not used during estimation. For more information, see <code>rng</code> in the MATLAB documentation.																
DataUsed	Attributes of the data used for estimation, returned as a structure with the following fields: <table border="1" data-bbox="387 604 1332 1296"> <thead> <tr> <th>Field</th> <th>Description</th> </tr> </thead> <tbody> <tr> <td>Name</td> <td>Name of the data set.</td> </tr> <tr> <td>Type</td> <td>Data type.</td> </tr> <tr> <td>Length</td> <td>Number of data samples.</td> </tr> <tr> <td>Ts</td> <td>Sample time.</td> </tr> <tr> <td>InterSample</td> <td>Input intersample behavior, returned as one of the following values: <ul style="list-style-type: none"> <li>'zoh' — Zero-order hold maintains a piecewise-constant input signal between samples.</li> <li>'foh' — First-order hold maintains a piecewise-linear input signal between samples.</li> <li>'bl' — Band-limited behavior specifies that the continuous-time input signal has zero power above the Nyquist frequency.</li> </ul> </td> </tr> <tr> <td>InputOffset</td> <td>Offset removed from time-domain input data during estimation. For nonlinear models, it is <code>[]</code>.</td> </tr> <tr> <td>OutputOffset</td> <td>Offset removed from time-domain output data during estimation. For nonlinear models, it is <code>[]</code>.</td> </tr> </tbody> </table>	Field	Description	Name	Name of the data set.	Type	Data type.	Length	Number of data samples.	Ts	Sample time.	InterSample	Input intersample behavior, returned as one of the following values: <ul style="list-style-type: none"> <li>'zoh' — Zero-order hold maintains a piecewise-constant input signal between samples.</li> <li>'foh' — First-order hold maintains a piecewise-linear input signal between samples.</li> <li>'bl' — Band-limited behavior specifies that the continuous-time input signal has zero power above the Nyquist frequency.</li> </ul>	InputOffset	Offset removed from time-domain input data during estimation. For nonlinear models, it is <code>[]</code> .	OutputOffset	Offset removed from time-domain output data during estimation. For nonlinear models, it is <code>[]</code> .
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For more information on using `Report`, see “Estimation Report”.

### **x0 — Initial states computed during estimation**

scalar | matrix

Initial states computed during estimation, returned as a scalar. If `data` contains multiple experiments, then `x0` is a matrix with each column corresponding to an experiment.

This value is also stored in the Parameters field of the model's Report property.

## Definitions

### Modal Form

In modal form,  $A$  is a block-diagonal matrix. The block size is typically 1-by-1 for real eigenvalues and 2-by-2 for complex eigenvalues. However, if there are repeated eigenvalues or clusters of nearby eigenvalues, the block size can be larger.

For example, for a system with eigenvalues  $(\lambda_1, \sigma \pm j\omega, \lambda_2)$ , the modal  $A$  matrix is of the form

$$\begin{bmatrix} \lambda_1 & 0 & 0 & 0 \\ 0 & \sigma & \omega & 0 \\ 0 & -\omega & \sigma & 0 \\ 0 & 0 & 0 & \lambda_2 \end{bmatrix}$$

### Companion Form

In the companion realization, the characteristic polynomial of the system appears explicitly in the right-most column of the  $A$  matrix.

For a system with characteristic polynomial

$$p(s) = s^n + \alpha_1 s^{n-1} + \dots + \alpha_{n-1} s + \alpha_n$$

the corresponding companion  $A$  matrix is

$$A = \begin{bmatrix} 0 & 0 & \dots & \dots & 0 & -\alpha_n \\ 1 & 0 & 0 & \dots & 0 & -\alpha_{n-1} \\ 0 & 1 & 0 & \dots & \dots & \vdots \\ \vdots & 0 & \dots & \dots & \vdots & \vdots \\ 0 & \dots & \dots & 1 & 0 & -\alpha_2 \\ 0 & \dots & \dots & 0 & 1 & -\alpha_1 \end{bmatrix}$$

The companion transformation requires that the system be controllable from the first input. The companion form is poorly conditioned for most state-space computations; avoid using it when possible.

### Tips

- `ssregest` function provides improved accuracy than `n4sid` for short, noisy data sets.
- For some problems, the quality of fit using `n4sid` is sensitive to options, such as `N4Horizon`, whose values can be difficult to determine. In comparison, the quality of fit with `ssregest` is less sensitive to its options, which makes `ssregest` simpler to use.

### Algorithms

`ssregest` estimates a regularized ARX model and converts the ARX model to a state-space model. The software then uses balanced model reduction techniques to reduce the state-space model to the specified order.

### References

[1] Ljung, L. *System Identification: Theory For the User*, Second Edition, Appendix 4A, pp 132-134, Upper Saddle River, N.J: Prentice Hall, 1999.

### See Also

`arx` | `arxRegul` | `balred` | `n4sid` | `ssest` | `ssregestOptions`

### Topics

“Regularized Estimates of Model Parameters”

**Introduced in R2014a**

# ssregestOptions

Option set for `ssregest`

## Syntax

```
options = ssregestOptions  
options = ssregestOptions(Name,Value)
```

## Description

`options = ssregestOptions` creates a default option set for `ssregest`.

`options = ssregestOptions(Name,Value)` specifies additional options using one or more `Name,Value` pair arguments.

## Examples

### Create Default Option Set for State-Space Estimation Using Reduction of Regularized ARX Model

```
options = ssregestOptions;
```

### Specify Options for State-Space Estimation Using Reduction of Regularized ARX Model

Create an option set for `ssregest` that fixes the value of the initial states to `'zero'`. Also, set the `Display` to `'on'`.

```
opt = ssregestOptions('InitialState','zero','Display','on');
```

Alternatively, use dot notation to set the values of `opt`.

```
opt = ssregestOptions;  
opt.InitialState = 'zero';  
opt.Display = 'on';
```

## Input Arguments

### Name-Value Pair Arguments

Specify optional comma-separated pairs of `Name`, `Value` arguments. `Name` is the argument name and `Value` is the corresponding value. `Name` must appear inside quotes. You can specify several name and value pair arguments in any order as `Name1, Value1, ..., NameN, ValueN`.

Example: `opt = ssregestOptions('InitialState', 'zero')` fixes the value of the initial states to zero.

#### **InitialState** — Handling of initial states

'estimate' (default) | 'zero'

Handling of initial states during estimation, specified as one of the following values:

- 'zero' — The initial state is set to zero.
- 'estimate' — The initial state is treated as an independent estimation parameter.

#### **ARXOrder** — ARX model orders

'auto' (default) | matrix of nonnegative integers

ARX model orders, specified as a matrix of nonnegative integers [`na` `nb` `nk`]. The `max(ARXOrder)+1` must be greater than the desired state-space model order (number of states). If you specify a value, it is recommended that you use a large value for `nb` order. To learn more about ARX model orders, see `arx`.

#### **RegularizationKernel** — Regularizing kernel

'TC' (default) | 'SE' | 'SS' | 'HF' | 'DI' | 'DC'

Regularizing kernel used for regularized estimates of the underlying ARX model, specified as one of the following values:

- 'TC' — Tuned and correlated kernel

- 'SE' — Squared exponential kernel
- 'SS' — Stable spline kernel
- 'HF' — High frequency stable spline kernel
- 'DI' — Diagonal kernel
- 'DC' — Diagonal and correlated kernel

For more information, see [1].

### Reduction — Options for model order reduction

structure

Options for model order reduction, specified as a structure with the following fields:

- StateElimMethod

State elimination method. Specifies how to eliminate the weakly coupled states (states with smallest Hankel singular values). Specified as one of the following values:

'MatchDC'	Discards the specified states and alters the remaining states to preserve the DC gain.
'Truncate'	Discards the specified states without altering the remaining states. This method tends to product a better approximation in the frequency domain, but the DC gains are not guaranteed to match.

**Default:** 'Truncate'

- AbsTol, RelTol

Absolute and relative error tolerance for stable/unstable decomposition. Positive scalar values. For an input model  $G$  with unstable poles, the reduction algorithm of `ssregest` first extracts the stable dynamics by computing the stable/unstable decomposition  $G \rightarrow GS + GU$ . The `AbsTol` and `RelTol` tolerances control the accuracy of this decomposition by ensuring that the frequency responses of  $G$  and  $GS + GU$  differ by no more than  $\text{AbsTol} + \text{RelTol} \cdot \text{abs}(G)$ . Increasing these tolerances helps separate nearby stable and unstable modes at the expense of accuracy. See `stabsep` for more information.

**Default:** `AbsTol = 0; RelTol = 1e-8`

- Offset

Offset for the stable/unstable boundary. Positive scalar value. In the stable/unstable decomposition, the stable term includes only poles satisfying

- $\text{Re}(s) < -\text{Offset} * \max(1, |\text{Im}(s)|)$  (Continuous time)
- $|z| < 1 - \text{Offset}$  (Discrete time)

Increase the value of `Offset` to treat poles close to the stability boundary as unstable.

**Default:** 1e-8

### Focus — Error to be minimized

'prediction' (default) | 'simulation'

Error to be minimized in the loss function during estimation, specified as the comma-separated pair consisting of 'Focus' and one of the following values:

- 'prediction' — The one-step ahead prediction error between measured and predicted outputs is minimized during estimation. As a result, the estimation focuses on producing a good predictor model.
- 'simulation' — The simulation error between measured and simulated outputs is minimized during estimation. As a result, the estimation focuses on making a good fit for simulation of model response with the current inputs.

The `Focus` option can be interpreted as a weighting filter in the loss function. For more information, see “Loss Function and Model Quality Metrics”.

### WeightingFilter — Weighting prefilter

[] (default) | vector | matrix | cell array | linear system

Weighting prefilter applied to the loss function to be minimized during estimation. To understand the effect of `WeightingFilter` on the loss function, see “Loss Function and Model Quality Metrics”.

Specify `WeightingFilter` as one of the following values:

- [] — No weighting prefilter is used.
- Passbands — Specify a row vector or matrix containing frequency values that define desired passbands. You select a frequency band where the fit between estimated model and estimation data is optimized. For example, `[wl, wh]` where `wl` and `wh` represent lower and upper limits of a passband. For a matrix with several rows defining frequency passbands, `[w1l, w1h; w2l, w2h; w3l, w3h; . . .]`, the estimation algorithm uses the union of the frequency ranges to define the estimation passband.

Passbands are expressed in `rad/TimeUnit` for time-domain data and in `FrequencyUnit` for frequency-domain data, where `TimeUnit` and `FrequencyUnit` are the time and frequency units of the estimation data.

- SISO filter — Specify a single-input-single-output (SISO) linear filter in one of the following ways:
  - A SISO LTI model
  - `{A, B, C, D}` format, which specifies the state-space matrices of a filter with the same sample time as estimation data.
  - `{numerator, denominator}` format, which specifies the numerator and denominator of the filter as a transfer function with same sample time as estimation data.

This option calculates the weighting function as a product of the filter and the input spectrum to estimate the transfer function.

- Weighting vector — Applicable for frequency-domain data only. Specify a column vector of weights. This vector must have the same length as the frequency vector of the data set, `Data.Frequency`. Each input and output response in the data is multiplied by the corresponding weight at that frequency.

### **EstimateCovariance — Control whether to generate parameter covariance data**

`true` (default) | `false`

Controls whether parameter covariance data is generated, specified as `true` or `false`.

If `EstimateCovariance` is `true`, then use `getcov` to fetch the covariance matrix from the estimated model.

### **Display — Specify whether to display the estimation progress**

`'off'` (default) | `'on'`

Specify whether to display the estimation progress, specified as one of the following values:

- `'on'` — Information on model structure and estimation results are displayed in a progress-viewer window.
- `'off'` — No progress or results information is displayed.

### **InputOffset — Removal of offset from time-domain input data during estimation**

`[]` (default) | vector of positive integers | matrix

Removal of offset from time-domain input data during estimation, specified as the comma-separated pair consisting of 'InputOffset' and one of the following:

- A column vector of positive integers of length  $Nu$ , where  $Nu$  is the number of inputs.
- [] — Indicates no offset.
- $Nu$ -by- $Ne$  matrix — For multi-experiment data, specify InputOffset as an  $Nu$ -by- $Ne$  matrix.  $Nu$  is the number of inputs, and  $Ne$  is the number of experiments.

Each entry specified by InputOffset is subtracted from the corresponding input data.

### **OutputOffset — Removal of offset from time-domain output data during estimation**

[] (default) | vector | matrix

Removal of offset from time-domain output data during estimation, specified as the comma-separated pair consisting of 'OutputOffset' and one of the following:

- A column vector of length  $Ny$ , where  $Ny$  is the number of outputs.
- [] — Indicates no offset.
- $Ny$ -by- $Ne$  matrix — For multi-experiment data, specify OutputOffset as a  $Ny$ -by- $Ne$  matrix.  $Ny$  is the number of outputs, and  $Ne$  is the number of experiments.

Each entry specified by OutputOffset is subtracted from the corresponding output data.

### **OutputWeight — Weight of prediction errors in multi-output estimation**

[] (default) | positive semidefinite, symmetric matrix

Weight of prediction errors in multi-output estimation, specified as one of the following values:

- Positive semidefinite, symmetric matrix ( $W$ ). The software minimizes the trace of the weighted prediction error matrix  $\text{trace}(E' * E * W / N)$  where:
  - $E$  is the matrix of prediction errors, with one column for each output, and  $W$  is the positive semidefinite, symmetric matrix of size equal to the number of outputs. Use  $W$  to specify the relative importance of outputs in multiple-output models, or the reliability of corresponding data.
  - $N$  is the number of data samples.
- [] — No weighting is used. Specifying as [] is the same as  $\text{eye}(Ny)$ , where  $Ny$  is the number of outputs.

This option is relevant only for multi-output models.

### **Advanced — Advanced estimation options**

structure

Advanced options for regularized estimation, specified as a structure with the following fields:

- **MaxSize** — Maximum allowable size of Jacobian matrices formed during estimation, specified as a large positive number.

**Default:** 250e3

- **SearchMethod** — Search method for estimating regularization parameters, specified as one of the following values:
  - `'gn'`: Quasi-Newton line search.
  - `'fmincon'`: Trust-region-reflective constrained minimizer. In general, `'fmincon'` is better than `'gn'` for handling bounds on regularization parameters that are imposed automatically during estimation.

**Default:** `'fmincon'`

## **Output Arguments**

### **options — Option set for ssregest**

ssregestOptions options set

Estimation options for `ssregest`, returned as an `ssregestOptions` option set.

## **References**

- [1] T. Chen, H. Ohlsson, and L. Ljung. “On the Estimation of Transfer Functions, Regularizations and Gaussian Processes - Revisited”, *Automatica*, Volume 48, August 2012.

## **See Also**

ssregest

**Topics**

“Loss Function and Model Quality Metrics”

**Introduced in R2014a**

# stack

Build model array by stacking models or model arrays along array dimensions

## Syntax

```
sys = stack(arraydim,sys1,sys2,...)
```

## Description

`sys = stack(arraydim,sys1,sys2,...)` produces an array of dynamic system models `sys` by stacking (concatenating) the models (or arrays) `sys1,sys2,...` along the array dimension `arraydim`. All models must have the same number of inputs and outputs (the same I/O dimensions), but the number of states can vary. The I/O dimensions are not counted in the array dimensions. For more information about model arrays and array dimensions, see “Model Arrays” (Control System Toolbox).

For arrays of state-space models with variable order, you cannot use the dot operator (e.g., `sys.A`) to access arrays. Use the syntax

```
[A,B,C,D] = ssdata(sys,'cell')
```

to extract the state-space matrices of each model as separate cells in the cell arrays `A`, `B`, `C`, and `D`.

## Examples

### Example 1

If `sys1` and `sys2` are two models:

- `stack(1,sys1,sys2)` produces a 2-by-1 model array.
- `stack(2,sys1,sys2)` produces a 1-by-2 model array.
- `stack(3,sys1,sys2)` produces a 1-by-1-by-2 model array.

## Example 2

Stack identified state-space models derived from the same estimation data and compare their bode responses.

```
load iddata1 z1
sysc = cell(1,5);
opt = ssestOptions('Focus','simulation');
for i = 1:5
    sysc{i} = ssest(z1,i-1,opt);
end
sysArray = stack(1, sysc{:});
bode(sysArray);
```

**Introduced in R2012a**

# step

Step response plot of dynamic system; step response data

## Syntax

```
step(sys)
step(sys,Tfinal)
step(sys,t)
step(sys1,sys2,...,sysN)
step(sys1,sys2,...,sysN,Tfinal)
step(sys1,sys2,...,sysN,t)
y = step(sys,t)
[y,t] = step(sys)
[y,t] = step(sys,Tfinal)
[y,t,x] = step(sys)
[y,t,x,ysd] = step(sys)
[y,...] = step(sys,...,options)
```

## Description

`step` calculates the step response of a dynamic system. For the state-space case, zero initial state is assumed. When it is invoked with no output arguments, this function plots the step response on the screen.

`step(sys)` plots the step response of an arbitrary dynamic system model, `sys`. This model can be continuous- or discrete-time, and SISO or MIMO. The step response of multi-input systems is the collection of step responses for each input channel. The duration of simulation is determined automatically, based on the system poles and zeros.

`step(sys,Tfinal)` simulates the step response from  $t = 0$  to the final time  $t = T_{\text{final}}$ . Express `Tfinal` in the system time units, specified in the `TimeUnit` property of `sys`. For discrete-time systems with unspecified sample time ( $T_s = -1$ ), `step` interprets `Tfinal` as the number of sampling periods to simulate.

`step(sys,t)` uses the user-supplied time vector `t` for simulation. Express `t` in the system time units, specified in the `TimeUnit` property of `sys`. For discrete-time models, `t`

should be of the form  $T_i:T_s:T_f$ , where  $T_s$  is the sample time. For continuous-time models,  $t$  should be of the form  $T_i:dt:T_f$ , where  $dt$  becomes the sample time of a discrete approximation to the continuous system (see “Algorithms” on page 1-1721). The `step` command always applies the step input at  $t=0$ , regardless of  $T_i$ .

To plot the step response of several models `sys1, ..., sysN` on a single figure, use

```
step(sys1,sys2, ..., sysN)
step(sys1,sys2, ..., sysN,Tfinal)
step(sys1,sys2, ..., sysN,t)
```

All of the systems plotted on a single plot must have the same number of inputs and outputs. You can, however, plot a mix of continuous- and discrete-time systems on a single plot. This syntax is useful to compare the step responses of multiple systems.

You can also specify a distinctive color, linestyle, marker, or all three for each system. For example,

```
step(sys1, 'y: ', sys2, 'g--')
```

plots the step response of `sys1` with a dotted yellow line and the step response of `sys2` with a green dashed line.

When invoked with output arguments:

```
y = step(sys,t)
[y,t] = step(sys)
[y,t] = step(sys,Tfinal)
[y,t,x] = step(sys)
```

`step` returns the output response  $y$ , the time vector  $t$  used for simulation (if not supplied as an input argument), and the state trajectories  $x$  (for state-space models only). No plot generates on the screen. For single-input systems,  $y$  has as many rows as time samples (length of  $t$ ), and as many columns as outputs. In the multi-input case, the step responses of each input channel are stacked up along the third dimension of  $y$ . The dimensions of  $y$  are then

*(length of t) × (number of outputs) × (number of inputs)*

and  $y(:, :, j)$  gives the response to a unit step command injected in the  $j$ th input channel. Similarly, the dimensions of  $x$  are

$$(length\ of\ t) \times (number\ of\ states) \times (number\ of\ inputs)$$

For identified models (see `idlti` and `idn1model`) `[y,t,x,ysd] = step(sys)` also computes the standard deviation `ysd` of the response `y` (`ysd` is empty if `sys` does not contain parameter covariance information).

`[y,...] = step(sys,...,options)` specifies additional options for computing the step response, such as the step amplitude or input offset. Use `stepDataOptions` to create the option set `options`.

## Examples

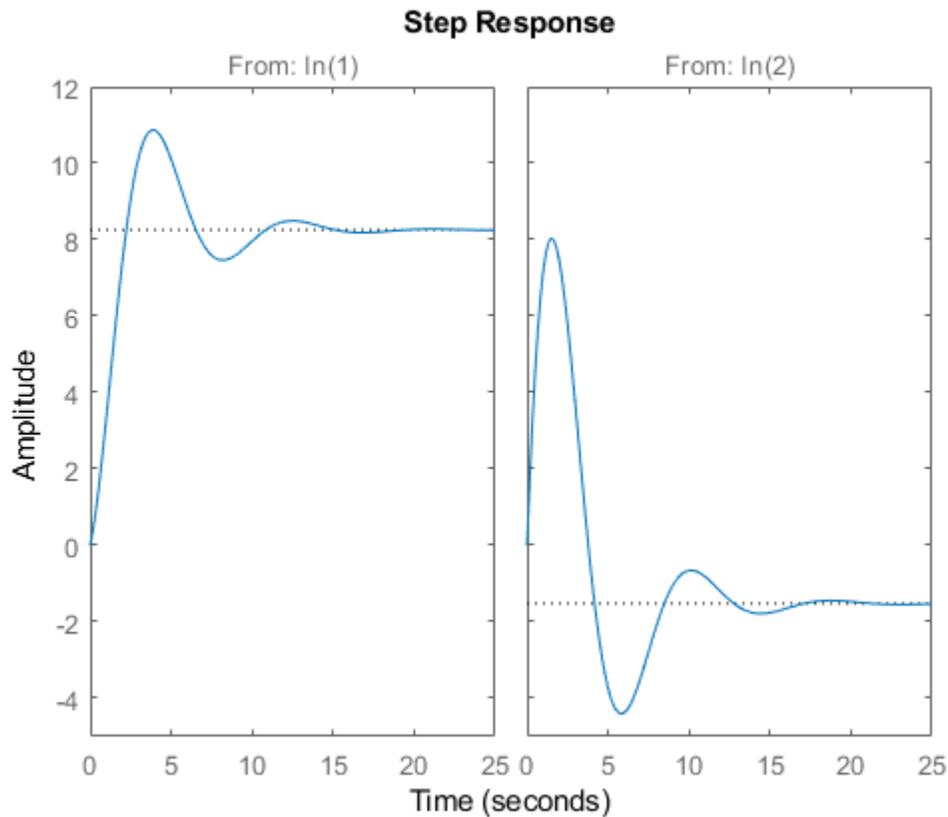
### Step Response Plot of Dynamic System

Plot the step response of the following second-order state-space model:

$$\begin{bmatrix} \dot{x}_1 \\ \dot{x}_2 \end{bmatrix} = \begin{bmatrix} -0.5572 & -0.7814 \\ 0.7814 & 0 \end{bmatrix} \begin{bmatrix} x_1 \\ x_2 \end{bmatrix} + \begin{bmatrix} 1 & -1 \\ 0 & 2 \end{bmatrix} \begin{bmatrix} u_1 \\ u_2 \end{bmatrix}$$

$$y = \begin{bmatrix} 1.9691 & 6.4493 \end{bmatrix} \begin{bmatrix} x_1 \\ x_2 \end{bmatrix}$$

```
a = [-0.5572, -0.7814; 0.7814, 0];
b = [1, -1; 0, 2];
c = [1.9691, 6.4493];
sys = ss(a, b, c, 0);
step(sys)
```

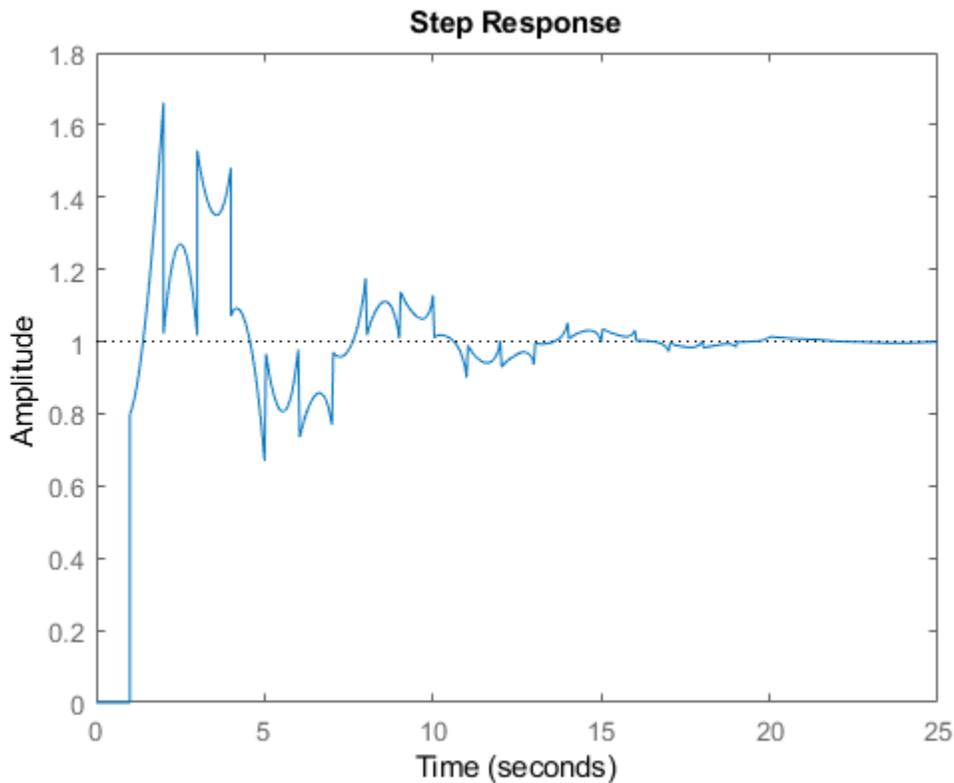


The left plot shows the step response of the first input channel, and the right plot shows the step response of the second input channel.

### Step Response Plot of Feedback Loop with Delay

Create a feedback loop with delay and plot its step response.

```
s = tf('s');  
G = exp(-s) * (0.8*s^2+s+2)/(s^2+s);  
T = feedback(ss(G),1);  
step(T)
```



The system step response displayed is chaotic. The step response of systems with internal delays may exhibit odd behavior, such as recurring jumps. Such behavior is a feature of the system and not software anomalies.

### Step Responses of Identified Models with Confidence Regions

Compare the step response of a parametric identified model to a non-parametric (empirical) model. Also view their  $3\sigma$  confidence regions.

Load the data.

```
load iddata1 z1
```

Estimate a parametric model.

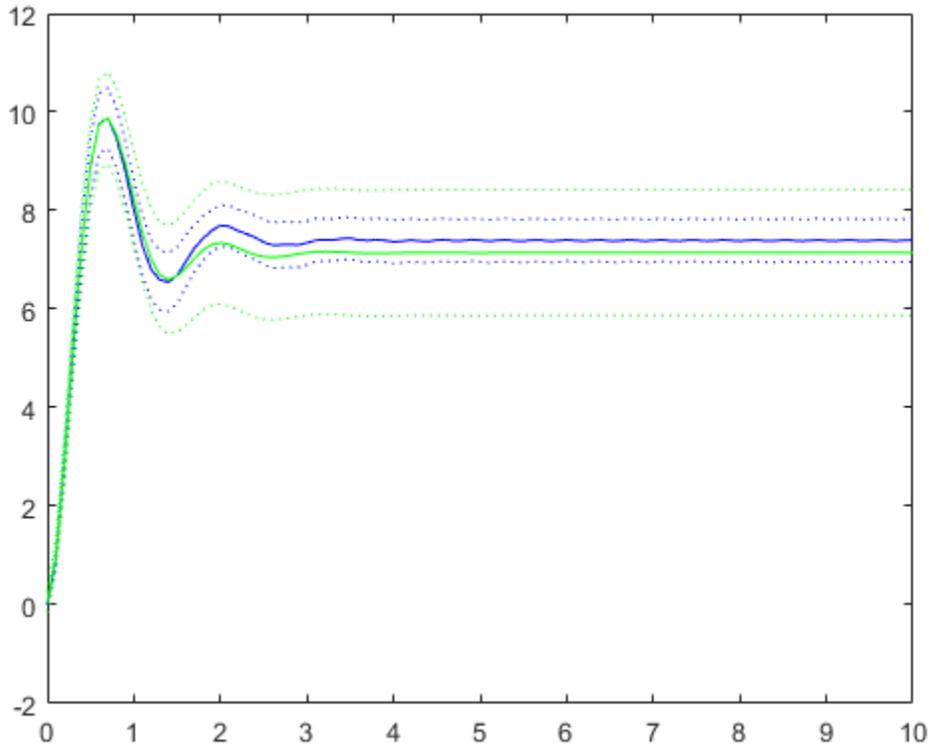
```
sys1 = ssest(z1,4);
```

Estimate a non-parametric model.

```
sys2 = impulseest(z1);
```

Plot the step responses for comparison.

```
t = (0:0.1:10)';  
[y1, ~, ~, ysd1] = step(sys1,t);  
[y2, ~, ~, ysd2] = step(sys2,t);  
plot(t, y1, 'b', t, y1+3*ysd1, 'b:', t, y1-3*ysd1, 'b:')  
hold on  
plot(t, y2, 'g', t, y2+3*ysd2, 'g:', t, y2-3*ysd2, 'g:')
```



### Validate Linearization of Identified Nonlinear ARX Model

Validate the linearization of a nonlinear ARX model by comparing the small amplitude step responses of the linear and nonlinear models.

Load the data.

```
load iddata2 z2;
```

Estimate a nonlinear ARX model.

```
nlsys = nlarx(z2,[4 3 10], 'tree', 'custom', {'sin(y1(t-2)*u1(t))+y1(t-2)*u1(t)+u1(t).*u1(t)'});
```

Determine an equilibrium operating point for `nlsys` corresponding to a steady-state input value of 1.

```
u0 = 1;  
[X,~,r] = findop(nlsys, 'steady', 1);  
y0 = r.SignalLevels.Output;
```

Obtain a linear approximation of `nlsys` at this operating point.

```
sys = linearize(nlsys,u0,X);
```

Validate the usefulness of `sys` by comparing its small-amplitude step response to that of `nlsys`.

The nonlinear system `nlsys` is operating at an equilibrium level dictated by  $(u_0, y_0)$ . Introduce a step perturbation of size 0.1 about this steady-state and compute the corresponding response.

```
opt = stepDataOptions;  
opt.InputOffset = u0;  
opt.StepAmplitude = 0.1;  
t = (0:0.1:10)';  
ynl = step(nlsys, t, opt);
```

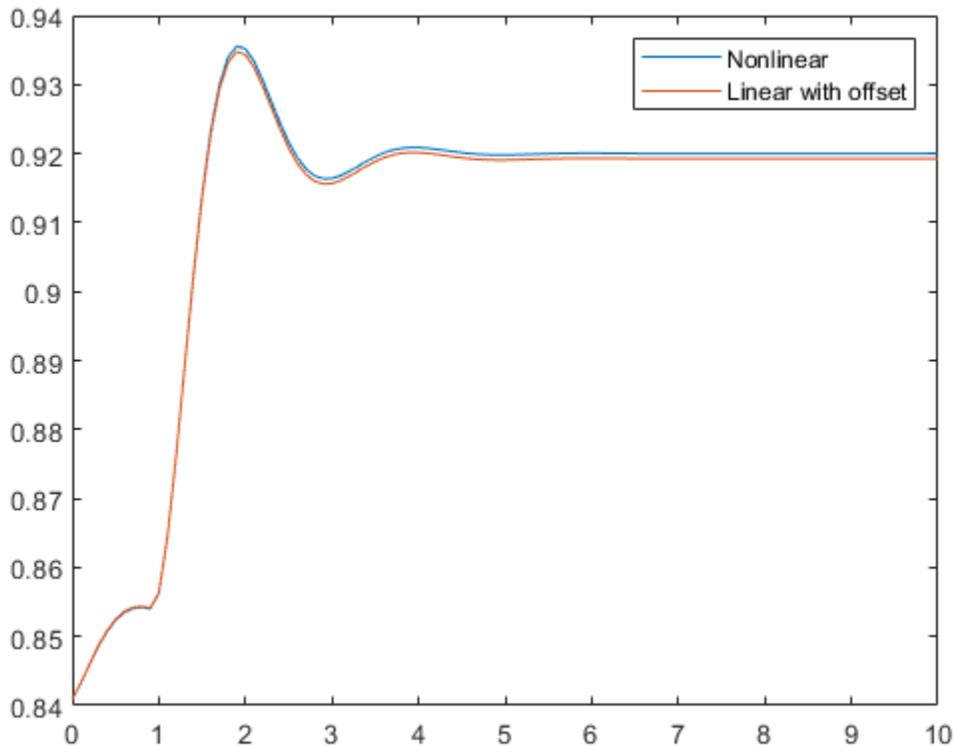
The linear system `sys` expresses the relationship between the perturbations in input to the corresponding perturbation in output. It is unaware of the nonlinear system's equilibrium values.

Plot the step response of the linear system.

```
opt = stepDataOptions;  
opt.StepAmplitude = 0.1;  
yl = step(sys, t, opt);
```

Add the steady-state offset,  $y_0$ , to the response of the linear system and plot the responses.

```
plot(t, ynl, t, yl+y0)  
legend('Nonlinear', 'Linear with offset')
```



### Step Response of Identified Time-Series Model

Compute the step response of an identified time-series model.

A time-series model, also called a signal model, is one without measured input signals. The step plot of this model uses its (unmeasured) noise channel as the input channel to which the step signal is applied.

Load the data.

```
load iddata9;
```

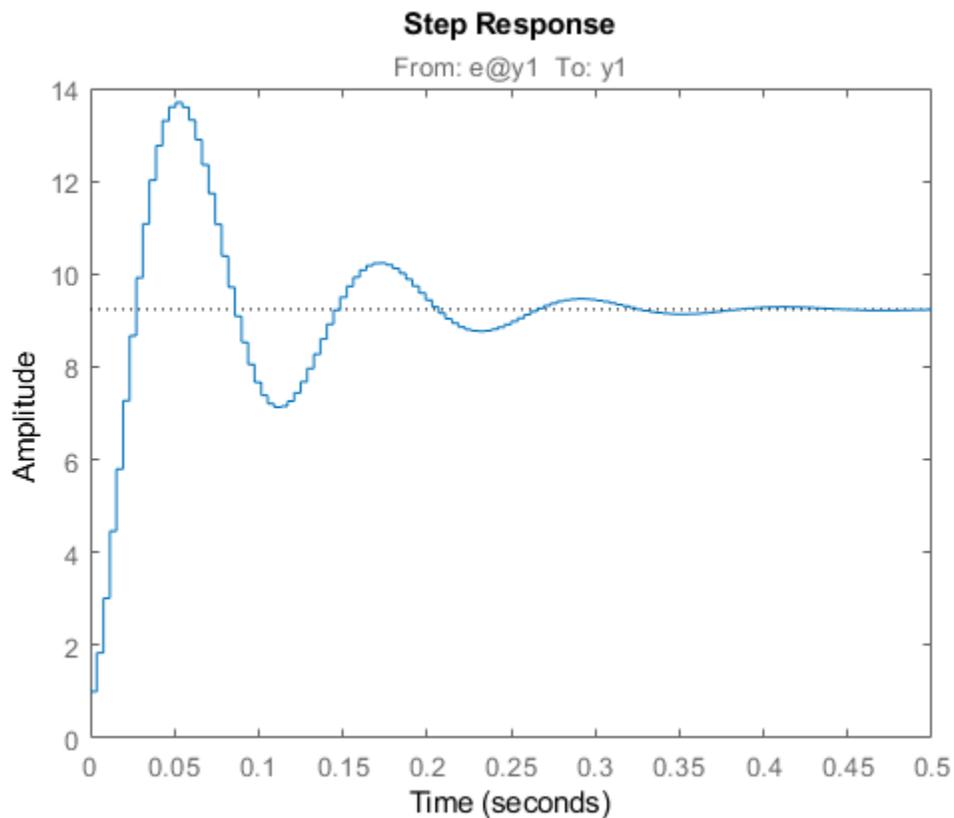
Estimate a time-series model.

```
sys = ar(z9, 4);
```

ys is a model of the form  $A y(t) = e(t)$ , where  $e(t)$  represents the noise channel. For computation of step response,  $e(t)$  is treated as an input channel, and is named e@y1.

Plot the step response.

```
step(sys)
```



## Tips

You can change the properties of your plot, for example the units. For information on the ways to change properties of your plots, see “Ways to Customize Plots” (Control System Toolbox).

## Algorithms

Continuous-time models without internal delays are converted to state space and discretized using zero-order hold on the inputs. The sample time is chosen automatically based on the system dynamics, except when a time vector  $t = 0:dt:Tf$  is supplied ( $dt$  is then used as sampling period). The resulting simulation time steps  $t$  are equisampled with spacing  $dt$ .

For systems with internal delays, Control System Toolbox software uses variable step solvers. As a result, the time steps  $t$  are not equisampled.

## References

- [1] L.F. Shampine and P. Gahinet, "Delay-differential-algebraic equations in control theory," *Applied Numerical Mathematics*, Vol. 56, Issues 3-4, pp. 574-588.

## See Also

**Linear System Analyzer** | `impulse` | `initial` | `lsim` | `stepDataOptions`

**Introduced before R2006a**

## step

Update model parameters and output online using recursive estimation algorithm

## Syntax

```
[EstimatedParameters,EstimatedOutput] = step(obj,y,InputData)
```

## Description

[EstimatedParameters,EstimatedOutput] = step(obj,y,InputData) updates parameters and output of the model specified in System object, obj, using measured output, y, and input data.

step puts the object into a locked state. In a locked state you cannot change any nontunable properties of the object, such as model order, data type, or estimation algorithm.

The EstimatedParameters and InputData depend on the online estimation System object:

- recursiveAR — step returns the estimated polynomial  $A(q)$  coefficients of a single output AR model using time-series output data.  
[A,EstimatedOutput] = step(obj,y)
- recursiveARMA — step returns the estimated polynomial  $A(q)$  and  $C(q)$  coefficients of a single output ARMA model using time-series output data, y.  
[A,C,EstimatedOutput] = step(obj,y)
- recursiveARX — step returns the estimated polynomial  $A(q)$  and  $B(q)$  coefficients of a SISO or MISO ARX model using measured input and output data, u and y, respectively.  
[A,B,EstimatedOutput] = step(obj,y,u).
- recursiveARMAX — step returns the estimated polynomial  $A(q)$ ,  $B(q)$ , and  $C(q)$  coefficients of a SISO ARMAX model using measured input and output data, u and y, respectively.  
[A,B,C,EstimatedOutput] = step(obj,y,u).

- `recursiveOE` — `step` returns the estimated polynomial  $B(q)$ , and  $F(q)$  coefficients of a SISO Output-Error polynomial model using measured input and output data,  $u$  and  $y$ , respectively.  
`[B,F,EstimatedOutput] = step(obj,y,u).`
- `recursiveBJ` — `step` returns the estimated polynomial  $B(q)$ ,  $C(q)$ ,  $D(q)$ , and  $F(q)$  coefficients of a SISO Box-Jenkins polynomial model using measured input and output data,  $u$  and  $y$ , respectively.  
`[B,C,D,F,EstimatedOutput] = step(obj,y,u).`
- `recursiveLS` — `step` returns the estimated system parameters,  $\theta$ , of a single output system that is linear in estimated parameters, using regressors  $H$  and output data  $y$ .  
`[theta,EstimatedOutput] = step(obj,y,H).`

## Examples

### Estimate an ARMAX Model Online

Create a System object for online parameter estimation of an ARMAX model.

```
obj = recursiveARMAX;
```

The ARMAX model has a default structure with polynomials of order 1 and initial polynomial coefficient values, `eps`.

Load the estimation data. In this example, use a static data set for illustration.

```
load iddata1 z1;
output = z1.y;
input = z1.u;
```

Estimate ARMAX model parameters online using `step`.

```
for i = 1:numel(input)
[A,B,C,EstimatedOutput] = step(obj,output(i),input(i));
end
```

View the current estimated values of polynomial A coefficients.

```
obj.A
```

```
ans = 1x2
```

```
1.0000 -0.8298
```

View the current covariance estimate of the parameters.

```
obj.ParameterCovariance
```

```
ans = 3×3
```

```
0.0001 0.0001 0.0001
0.0001 0.0032 0.0000
0.0001 0.0000 0.0001
```

View the current estimated output.

```
EstimatedOutput
```

```
EstimatedOutput = -4.5595
```

## **Tune Recursive Estimation Algorithm Properties During Online Parameter Estimation**

Create a System object for online parameter estimation of an ARMAX model.

```
obj = recursiveARMAX;
```

The ARMAX model has a default structure with polynomials of order 1 and initial polynomial coefficient values, `eps`.

Load the estimation data. In this example, use a static data set for illustration.

```
load iddata1 z1;
output = z1.y;
input = z1.u;
dataSize = numel(input);
```

Estimate ARMAX model parameters online using the default recursive estimation algorithm, Forgetting Factor. Change the `ForgettingFactor` property during online parameter estimation.

```
for i = 1:dataSize
    if i == dataSize/2
```

```

        obj.ForgettingFactor = 0.98;
    end
    [A,B,C,EstimatedOutput] = step(obj,output(i),input(i));
end

```

## Estimate Parameters of System Using Recursive Least Squares Algorithm

The system has two parameters and is represented as:

$$y(t) = a_1 u(t) + a_2 u(t-1)$$

Here,

- $u$  and  $y$  are the real-time input and output data, respectively.
- $u(t)$  and  $u(t-1)$  are the regressors,  $H$ , of the system.
- $a_1$  and  $a_2$  are the parameters,  $\theta$ , of the system.

Create a System object for online estimation using the recursive least squares algorithm.

```
obj = recursiveLS(2);
```

Load the estimation data, which for this example is a static data set.

```
load iddata3
input = z3.u;
output = z3.y;
```

Create a variable to store  $u(t-1)$ . This variable is updated at each time step.

```
oldInput = 0;
```

Estimate the parameters and output using `step` and input-output data, maintaining the current regressor pair in  $H$ . Invoke the `step` function implicitly by calling the `obj` system object with input arguments.

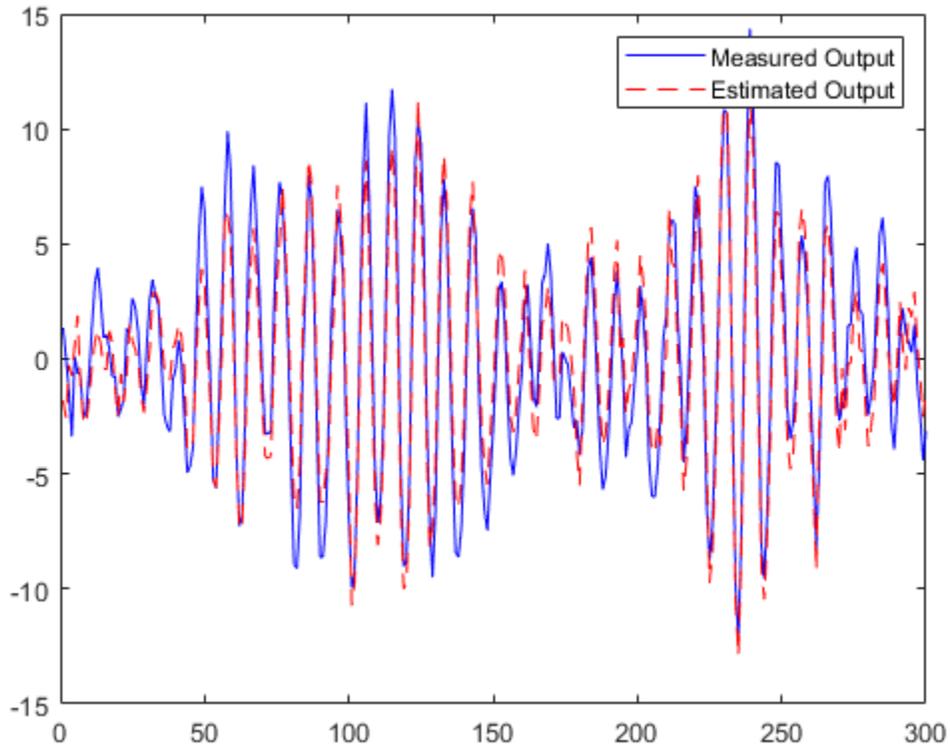
```
for i = 1:numel(input)
    H = [input(i) oldInput];
    [theta, EstimatedOutput] = obj(output(i),H);
end

```

```
    estimatedOut(i)= EstimatedOutput;  
    theta_est(i,:) = theta;  
    oldInput = input(i);  
end
```

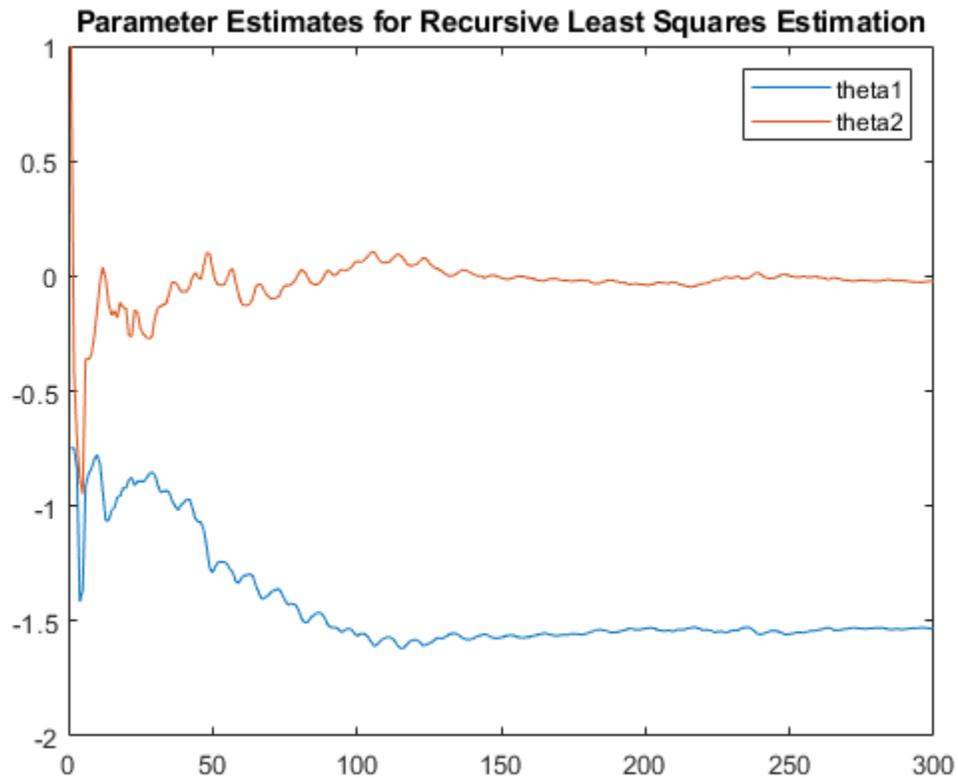
Plot the measured and estimated output data.

```
numSample = 1:numel(input);  
plot(numSample,output,'b',numSample,estimatedOut,'r--');  
legend('Measured Output','Estimated Output');
```



Plot the parameters.

```
plot(numSample, theta_est(:,1), numSample, theta_est(:,2))  
title('Parameter Estimates for Recursive Least Squares Estimation')  
legend("theta1", "theta2")
```



View the final estimates.

```
theta_final = theta
```

```
theta_final = 2×1
```

```
-1.5322
```

```
-0.0235
```

## Input Arguments

### **obj** — System object for online parameter estimation

recursiveAR object | recursiveARMA object | recursiveARX object | recursiveARMAX object | recursiveOE object | recursiveBJ object | recursiveLS object

System object for online parameter estimation, created using one of the following commands:

- recursiveAR
- recursiveARMA
- recursiveARX
- recursiveARMAX
- recursiveOE
- recursiveBJ
- recursiveLS

The `step` command updates parameters of the model using the recursive estimation algorithm specified in `obj` and the incoming input-output data.

### **y** — Output data

real scalar

Output data acquired in real time, specified as a real scalar.

### **InputData** — Input data

scalar | vector of real values

Input data acquired in real time, specified as a scalar or vector of real values depending on the type of System object.

<b>System object</b>	<b>Model Type</b>	<b>InputData</b>
recursiveAR	Time-series	Not Applicable
recursiveARMA	Time-series	Not Applicable
recursiveARX	SISO ARX	Real scalar

System object	Model Type	InputData
	MISO ARX with $N$ inputs	Column vector of length $N$ , specified as real values
recursiveARMAX	SISO	Real scalar
recursiveOE	SISO	Real scalar
recursiveBJ	SISO	Real scalar
recursiveLS	Single output system with $Np$ system parameters	Regressors, $H$ , specified as a vector of real values of length $Np$

## Output Arguments

### EstimatedParameters — Estimated model parameters

vector of real values for each parameter

Estimated model parameters, returned as vectors of real values. The number of estimated parameters, and so the step syntax, depend on the type of System object:

Online Estimation System Object	Estimated Parameters
recursiveAR	Polynomial $A(q)$ coefficients
recursiveARMA	Polynomials $A(q)$ and $C(q)$ coefficients
recursiveARX	Polynomials $A(q)$ and $B(q)$ coefficients
recursiveARMAX	Polynomials $A(q)$ , $B(q)$ , and $C(q)$ coefficients
recursiveOE	Polynomials $B(q)$ and $F(q)$
recursiveBJ	Polynomials $B(q)$ , $C(q)$ , $D(q)$ , and $F(q)$ coefficients
recursiveLS	System parameters, $\theta$

### EstimatedOutput — Estimated output

real scalar

Estimated output, returned as a real scalar. The output is estimated using input-output estimation data, current parameter values, and recursive estimation algorithm specified in obj.

## Tips

- Starting in R2016b, instead of using the `step` command to update model parameter estimates, you can call the System object with input arguments, as if it were a function. For example, `[A,EstimatedOutput] = step(obj,y)` and `[A,EstimatedOutput] = obj(y)` perform equivalent operations.

## See Also

`clone` | `isLocked` | `recursiveAR` | `recursiveARMA` | `recursiveARMAX` | `recursiveARX` | `recursiveBJ` | `recursiveLS` | `recursiveOE` | `release` | `reset`

## Topics

“Perform Online Parameter Estimation at the Command Line”

“Validate Online Parameter Estimation at the Command Line”

“Online ARX Parameter Estimation for Tracking Time-Varying System Dynamics”

“Line Fitting with Online Recursive Least Squares Estimation”

“What Is Online Estimation?”

## Introduced in R2015b

# stepDataOptions

Options set for step

## Syntax

```
opt = stepDataOptions  
opt = stepDataOptions(Name, Value)
```

## Description

`opt = stepDataOptions` creates the default options for `step`.

`opt = stepDataOptions(Name, Value)` creates an options set with the options specified by one or more `Name, Value` pair arguments.

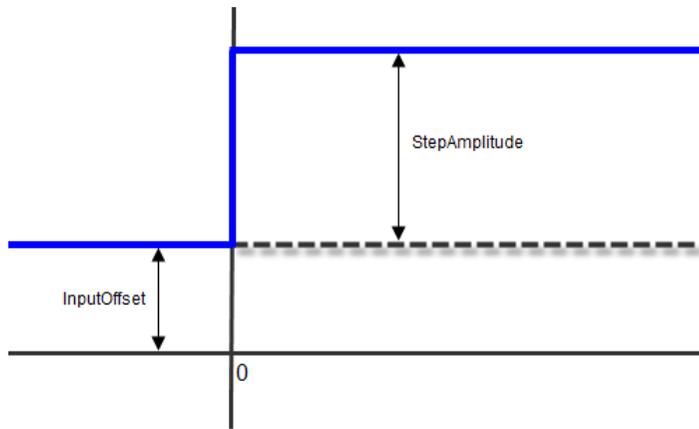
## Input Arguments

### Name-Value Pair Arguments

Specify optional comma-separated pairs of `Name, Value` arguments. `Name` is the argument name and `Value` is the corresponding value. `Name` must appear inside quotes. You can specify several name and value pair arguments in any order as `Name1, Value1, ..., NameN, ValueN`.

### InputOffset

Input signal level for all time  $t < 0$ , as shown in the next figure.



**Default:** 0

## **StepAmplitude**

Change of input signal level which occurs at time  $t = 0$ , as shown in the previous figure.

**Default:** 1

## **Output Arguments**

### **opt**

Option set containing the specified options for `step`.

## **Examples**

### **Specify Input Offset and Step Amplitude Level for Step Response**

Create a transfer function model.

```
sys = tf(1,[1,1]);
```

Create an option set for `step` to specify input offset and step amplitude level.

```
opt = stepDataOptions('InputOffset',-1,'StepAmplitude',2);
```

Calculate the step response using the specified options.

```
[y,t] = step(sys,opt);
```

## See Also

step

**Introduced in R2012a**

## stepinfo

Rise time, settling time, and other step-response characteristics

### Syntax

```
S = stepinfo(sys)
```

```
S = stepinfo(y,t)
```

```
S = stepinfo(y,t,yfinal)
```

```
S = stepinfo( ____, 'SettlingTimeThreshold', ST)
```

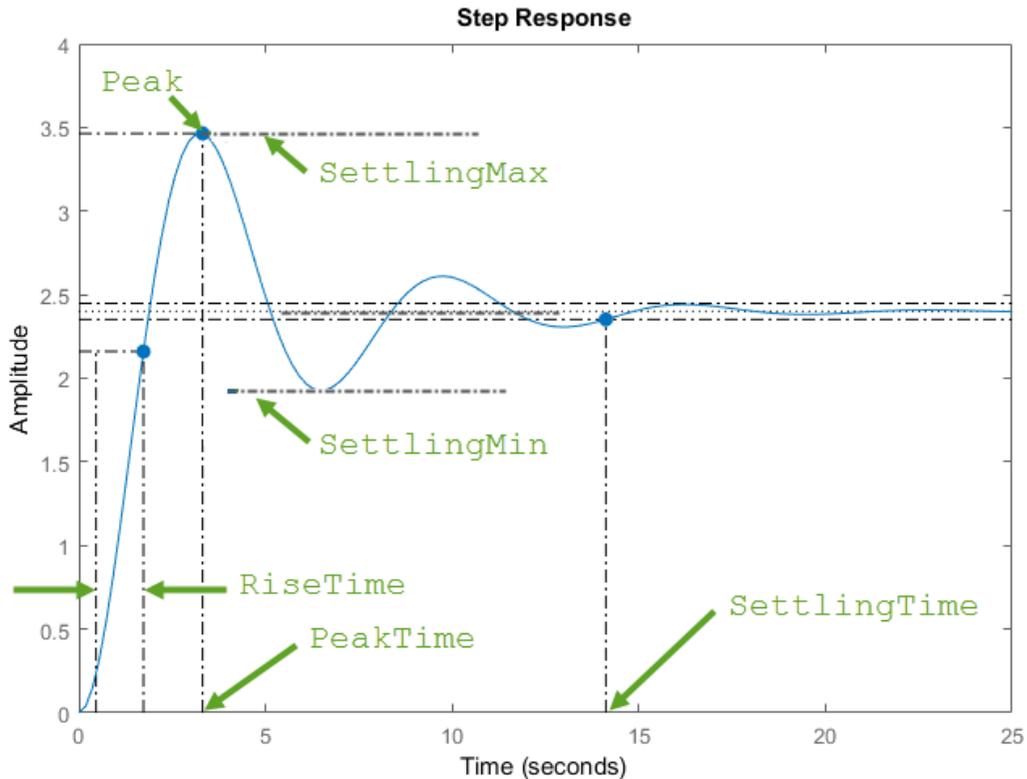
```
S = stepinfo( ____, 'RiseTimeLimits', RT)
```

### Description

`S = stepinfo(sys)` computes the step-response characteristics for a dynamic system model `sys`. The function returns the characteristics in a structure containing the fields:

- **RiseTime** — Time it takes for the response to rise from 10% to 90% of the steady-state response.
- **SettlingTime** — Time it takes for the error  $|y(t) - y_{final}|$  between the response  $y(t)$  and the steady-state response  $y_{final}$  to fall to within 2% of  $y_{final}$ .
- **SettlingMin** — Minimum value of  $y(t)$  once the response has risen.
- **SettlingMax** — Maximum value of  $y(t)$  once the response has risen.
- **Overshoot** — Percentage overshoot, relative to  $y_{final}$ .
- **Undershoot** — Percentage undershoot.
- **Peak** — Peak absolute value of  $y(t)$
- **PeakTime** — Time at which the peak value occurs.

The following figure illustrates some of these quantities on a typical second-order response.



Using this syntax requires a Control System Toolbox license.

`S = stepinfo(y,t)` computes step-response characteristics from an array of step-response data `y` and corresponding time vector `t`. For SISO system responses, `y` is a vector with the same number of entries as `t`. For MIMO response data, `y` is an array containing the responses of each I/O channel. This syntax uses the last value in `y` (or the last value in each channel's corresponding response data) as the steady-state value for computing characteristics that depend on that value.

`S = stepinfo(y,t,yfinal)` computes step-response characteristics relative to the steady-state value `yfinal`. This syntax is useful when you know that the expected steady-state system response differs from the last value in `y` for reasons such as measurement noise.

For SISO responses,  $t$  and  $y$  are vectors with the same length  $NS$ . For systems with  $NU$  inputs and  $NY$  outputs, you can specify  $y$  as an  $NS$ -by- $NY$ -by- $NU$  array (see `step`) and  $y_{final}$  as an  $NY$ -by- $NU$  array. `stepinfo` then returns a  $NY$ -by- $NU$  structure array  $S$  of performance metrics for each I/O pair.

`S = stepinfo( ____, 'SettlingTimeThreshold', ST)` lets you specify the threshold  $ST$  used in the definition of settling time. The response has settled when the error  $|y(t) - y_{final}|$  becomes smaller than a fraction  $ST$  of its peak value. The default value is  $ST = 0.02$  (2%). You can use this syntax with any of the previous input-argument combinations.

`S = stepinfo( ____, 'RiseTimeLimits', RT)` lets you specify the lower and upper thresholds used in the definition of rise time. By default, the rise time is defined as the time the response takes to rise from 10 to 90% of the steady-state value ( $RT = [0.1 \ 0.9]$ ). The upper threshold  $RT(2)$  is also used to calculate `SettlingMin` and `SettlingMax`. These values are the minimum and maximum values of the response occurring after the response has reached the upper threshold. You can use this syntax with any of the previous input-argument combinations.

## Examples

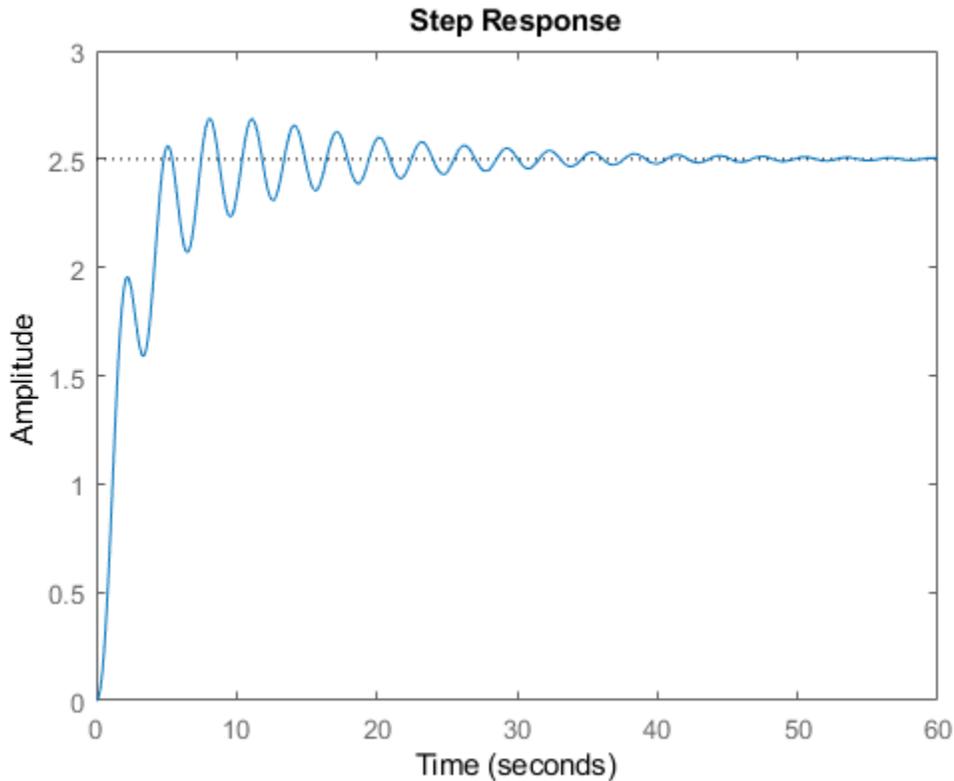
### Step-Response Characteristics of Dynamic System

Compute step-response characteristics such as rise time, settling time, and overshoot for a dynamic system model. For this example, use the continuous-time transfer function:

$$sys = \frac{s^2 + 5s + 5}{s^4 + 1.65s^3 + 5s^2 + 6.5s + 2}$$

Create the transfer function and examine its step response.

```
sys = tf([1 5 5],[1 1.65 5 6.5 2]);  
step(sys)
```



The plot shows that the response rises in a few seconds, and then rings down to a steady-state value of about 2.5. Compute the characteristics of this response using `stepinfo`.

```
S = stepinfo(sys)
```

```
S = struct with fields:  
    RiseTime: 3.8456  
    SettlingTime: 27.9762  
    SettlingMin: 2.0689  
    SettlingMax: 2.6873  
    Overshoot: 7.4915  
    Undershoot: 0  
    Peak: 2.6873
```

PeakTime: 8.0530

By default, the settling time is the time it takes for  $y(t) - y_{final}$  to fall below 2% of its peak value, where  $y(t)$  is the system response at time  $t$  and  $y_{final}$  is the steady-state response. The result `S.SettlingTime` shows that for `sys`, this condition occurs after about 28 seconds. The default definition of rise time is the time it takes for the response to go from 10% of its steady-state value to 90% of that value. `S.RiseTime` shows that for `sys`, this rise occurs in less than 4 seconds. The maximum overshoot is returned in `S.Overshoot`. For this system, the peak value `S.Peak`, which occurs at the time `S.PeakTime`, overshoots the steady-state value by about 7.5% of the steady-state value.

## Step-Response Characteristics of MIMO System

For a MIMO system, `stepinfo` returns a structure array in which each entry contains the response characteristics of the corresponding I/O channel of the system. For this example, use a two-output, two-input discrete-time system. Compute the step-response characteristics.

```
A = [0.68 -0.34; 0.34 0.68];  
B = [0.18 -0.05; 0.04 0.11];  
C = [0 -1.53; -1.12 -1.10];  
D = [0 0; 0.06 -0.37];  
sys = ss(A,B,C,D,0.2);
```

```
S = stepinfo(sys)
```

```
S = 2x2 struct array with fields:  
  RiseTime  
  SettlingTime  
  SettlingMin  
  SettlingMax  
  Overshoot  
  Undershoot  
  Peak  
  PeakTime
```

Access the response characteristics for a particular I/O channel by indexing into *S*. For instance, examine the response characteristics for the response from the first input to the second output of *sys*, corresponding to *S*(2,1).

*S*(2,1)

```
ans = struct with fields:
    RiseTime: 0.4000
    SettlingTime: 2.8000
    SettlingMin: -0.6724
    SettlingMax: -0.5188
    Overshoot: 24.6476
    Undershoot: 11.1224
    Peak: 0.6724
    PeakTime: 1
```

To access a particular value, use dot notation. For instance, extract the rise time of the (2,1) channel.

```
rt21 = S(2,1).RiseTime
rt21 = 0.4000
```

### Specify Definition of Settling Time or Rise Time

By default, `stepinfo` defines settling time as the time it takes for the error  $|y(t) - y_{final}|$  between the response  $y(t)$  and the steady-state response  $y_{final}$  to come within 2% of  $y_{final}$ . Also, `stepinfo` defines the rise time as the time it takes for the response to rise from 10% of  $y_{final}$  to 90% of  $y_{final}$ . You can change these definitions using `SettlingTimeThreshold` and `RiseTimeThreshold`. For this example, use the system given by:

$$sys = \frac{s^2 + 5s + 5}{s^4 + 1.65s^3 + 6.5s + 2}$$

Create the transfer function.

```
sys = tf([1 5 5],[1 1.65 5 6.5 2]);
```

Compute the time it takes for the error in the response of `sys` to reach 0.5% of the steady-state response. To do so, set `SettlingTimeThreshold` to 0.5%, or 0.005.

```
S1 = stepinfo(sys, 'SettlingTimeThreshold', 0.005);  
st1 = S1.SettlingTime
```

```
st1 = 46.1325
```

Compute the time it takes the response of `sys` to rise from 5% to 95% of the steady-state value. To do so, set `RiseTimeThreshold` to a vector containing those bounds.

```
S2 = stepinfo(sys, 'RiseTimeThreshold', [0.05 0.95]);  
rt2 = S2.RiseTime
```

```
rt2 = 4.1690
```

You can define both settling time and rise time in the same computation.

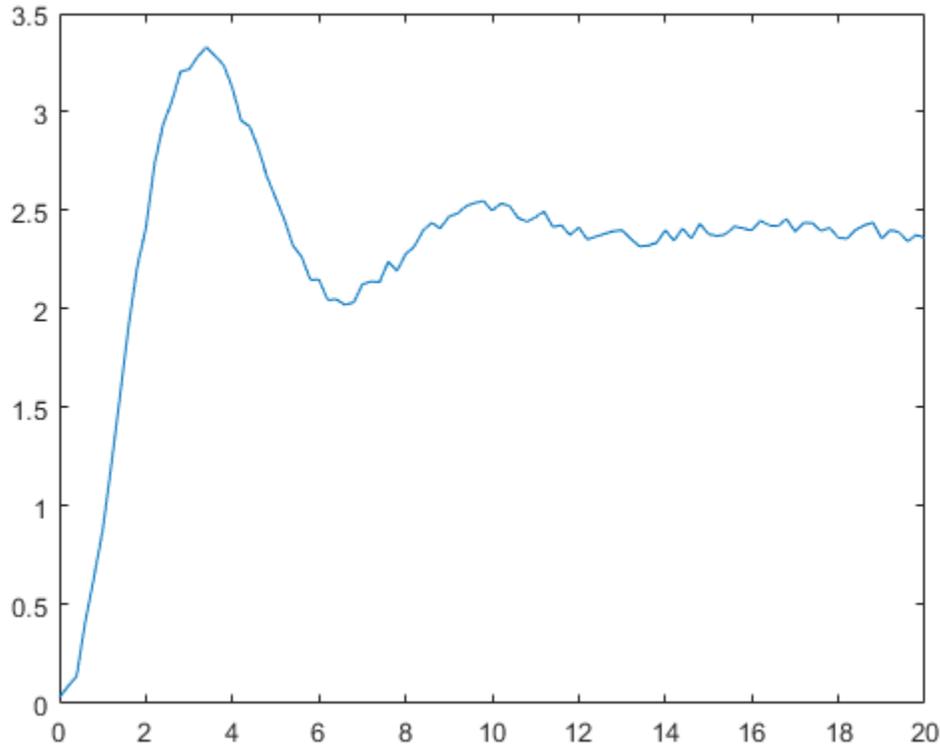
```
S3 = stepinfo(sys, 'SettlingTimeThreshold', 0.005, 'RiseTimeThreshold', [0.05 0.95])
```

```
S3 = struct with fields:  
    RiseTime: 4.1690  
    SettlingTime: 46.1325  
    SettlingMin: 2.0689  
    SettlingMax: 2.6873  
    Overshoot: 7.4915  
    Undershoot: 0  
    Peak: 2.6873  
    PeakTime: 8.0530
```

## Step-Response Characteristics from Response Data

You can extract step-response characteristics from step-response data even if you do not have a model of your system. For instance, suppose you have measured the response of your system to a step input, and saved the resulting response data in a vector `y` of response values at the times stored in another vector, `t`. Load the response data and examine it.

```
load StepInfoData t y  
plot(t,y)
```



Compute step-response characteristics from this response data using `stepinfo`. If you do not specify the steady-state response value `yfinal`, then `stepinfo` assumes that the last value in the response vector `y` is the steady-state response. Because there is some noise in the data, the last value in `y` is likely not the true steady-state response value. When you know what the steady-state value should be, you can provide it to `stepinfo`. For this example, suppose that the steady-state response is 2.4.

```
S1 = stepinfo(y,t,2.4)
```

```
S1 = struct with fields:  
    RiseTime: 1.2713  
    SettlingTime: 19.6478  
    SettlingMin: 2.0219  
    SettlingMax: 3.3302
```

```
Overshoot: 38.7575
Undershoot: 0
Peak: 3.3302
PeakTime: 3.4000
```

Because of the noise in the data, the default definition of the settling time is too stringent, resulting in an arbitrary value of almost 20 seconds. To allow for the noise, increase the settling-time threshold from the default 2% to 5%.

```
S2 = stepinfo(y,t,2.4,'SettlingTimeThreshold',0.05)
```

```
S2 = struct with fields:
  RiseTime: 1.2713
  SettlingTime: 10.4201
  SettlingMin: 2.0219
  SettlingMax: 3.3302
  Overshoot: 38.7575
  Undershoot: 0
  Peak: 3.3302
  PeakTime: 3.4000
```

## Input Arguments

### **sys** — Dynamic system

dynamic system model

Dynamic system, specified as a SISO or MIMO dynamic system model. Dynamic systems that you can use include:

- Continuous-time or discrete-time numeric LTI models, such as `tf`, `zpk`, or `ss` models.
- Generalized or uncertain LTI models such as `genss` or `uss` models. (Using uncertain models requires Robust Control Toolbox software.) For generalized models, `stepinfo` computes the step-response characteristics using the current value of tunable blocks and the nominal value of uncertain blocks.
- Identified LTI models, such as `idtf`, `idss`, or `idproc` models.

### **y** — Step-response data

vector | array

Step-response data, specified as:

- For SISO response data, a vector of length  $N_s$ , where  $N_s$  is the number of samples in the response data.
- For MIMO response data, an  $N_s$ -by- $N_y$ -by- $N_u$  array, where  $N_y$  is the number of system outputs, and  $N_u$  is the number of system inputs.

### **t** — Time vector

vector

Time vector corresponding to the response data in  $y$ , specified as a vector of length  $N_s$ .

### **yfinal** — Steady-state response

scalar | array

Steady-state response, specified as:

- For SISO response data, a scalar value.
- For MIMO response data, an  $N_y$ -by- $N_u$  array, where each entry provides the steady-state response value for the corresponding system channel.

If you do not provide `yfinal`, then `stepinfo` uses the last value in the corresponding channel of  $y$  as the steady-state response value.

### **ST** — Threshold for defining settling time

0.02 (default) | scalar between 0 and 1

Threshold for defining settling time, specified as a scalar value between 0 and 1. By default, `stepinfo` defines settling time as the time it takes for the error  $|y(t) - y_{final}|$  between the response  $y(t)$  and the steady-state response  $y_{final}$  to fall to within 2% of  $y_{final}$ . To change this definition, set `ST` to a different value. For instance, to set a threshold of 5%, set `ST` to 0.05.

### **RT** — Thresholds for defining rise time

[0.1 0.9] (default) | 2-element row vector

Threshold for defining rise time, specified as a 2-element row vector of nondecreasing values between 0 and 1. By default, `stepinfo` defines rise time as the time it takes for the response to rise from 10% to 90% of the steady-state value  $y_{final}$ . To change this definition, set `RT` to a different value. For instance, to define the rise time as the time it takes for the response to rise from 5% to 95% of the steady-state value, set `RT` to [0.05 0.95].

## Output Arguments

### **S** — Step-response characteristics

structure

Step-response characteristics of `sys`, returned as a structure containing the fields:

- `RiseTime` — Time it takes for the response to rise from 10% to 90% of the steady-state response.
- `SettlingTime` — Time it takes for the error  $|y(t) - y_{final}|$  between the response  $y(t)$  and the steady-state response  $y_{final}$  to fall to within 2% of  $y_{final}$ .
- `SettlingMin` — Minimum value of  $y(t)$  once the response has risen.
- `SettlingMax` — Maximum value of  $y(t)$  once the response has risen.
- `Overshoot` — Percentage overshoot, relative to  $y_{final}$ .
- `Undershoot` — Percentage undershoot.
- `Peak` — Peak absolute value of  $y(t)$
- `PeakTime` — Time at which the peak value occurs.

For MIMO models or responses data, `S` is a structure array in which each entry contains the step-response characteristics of the corresponding I/O channel. For instance, if you provide a 3-input, 3-output model or array of response data, then `S(2,3)` contains the characteristics of the response from the third input to the second output. For an example, see “Step-Response Characteristics of MIMO System” (Control System Toolbox).

If `sys` is unstable, then all step-response characteristics are NaN, except for `Peak` and `PeakTime`, which are Inf.

## See Also

`lsiminfo` | `step`

**Introduced in R2006a**

# stepplot

Plot step response and return plot handle

## Syntax

```
h = stepplot(sys)
stepplot(sys,Tfinal)
stepplot(sys,t)
stepplot(sys1,sys2,...,sysN)
stepplot(sys1,sys2,...,sysN,Tfinal)
stepplot(sys1,sys2,...,sysN,t)
stepplot(AX,...)
stepplot(..., plotoptions)
stepplot(..., dataoptions)
```

## Description

`h = stepplot(sys)` plots the step response of the dynamic system model `sys`. It also returns the plot handle `h`. You can use this handle to customize the plot with the `getoptions` and `setoptions` commands. Type

```
help timeoptions
```

for a list of available plot options.

For multiinput models, independent step commands are applied to each input channel. The time range and number of points are chosen automatically.

`stepplot(sys,Tfinal)` simulates the step response from  $t = 0$  to the final time  $t = T_{\text{final}}$ . Express `Tfinal` in the system time units, specified in the `TimeUnit` property of `sys`. For discrete-time systems with unspecified sample time ( $T_s = -1$ ), `stepplot` interprets `Tfinal` as the number of sampling intervals to simulate.

`stepplot(sys,t)` uses the user-supplied time vector `t` for simulation. Express `t` in the system time units, specified in the `TimeUnit` property of `sys`. For discrete-time models, `t` should be of the form `Ti:Ts:Tf`, where `Ts` is the sample time. For continuous-time

models,  $t$  should be of the form  $T_i:dt:T_f$ , where  $dt$  becomes the sample time of a discrete approximation to the continuous system (see `step`). The `stepplot` command always applies the step input at  $t=0$ , regardless of  $T_i$ .

To plot the step responses of multiple models `sys1,sys2,...` on a single plot, use:

```
stepplot(sys1,sys2,...,sysN)
```

```
stepplot(sys1,sys2,...,sysN,Tfinal)
```

```
stepplot(sys1,sys2,...,sysN,t)
```

You can also specify a color, line style, and marker for each system, as in

```
stepplot(sys1,'r',sys2,'y--',sys3,'gx')
```

`stepplot(AX,...)` plots into the axes with handle `AX`.

`stepplot(..., plotoptions)` customizes the plot appearance using the options set, `plotoptions`. Use `timeOptions` to create the options set.

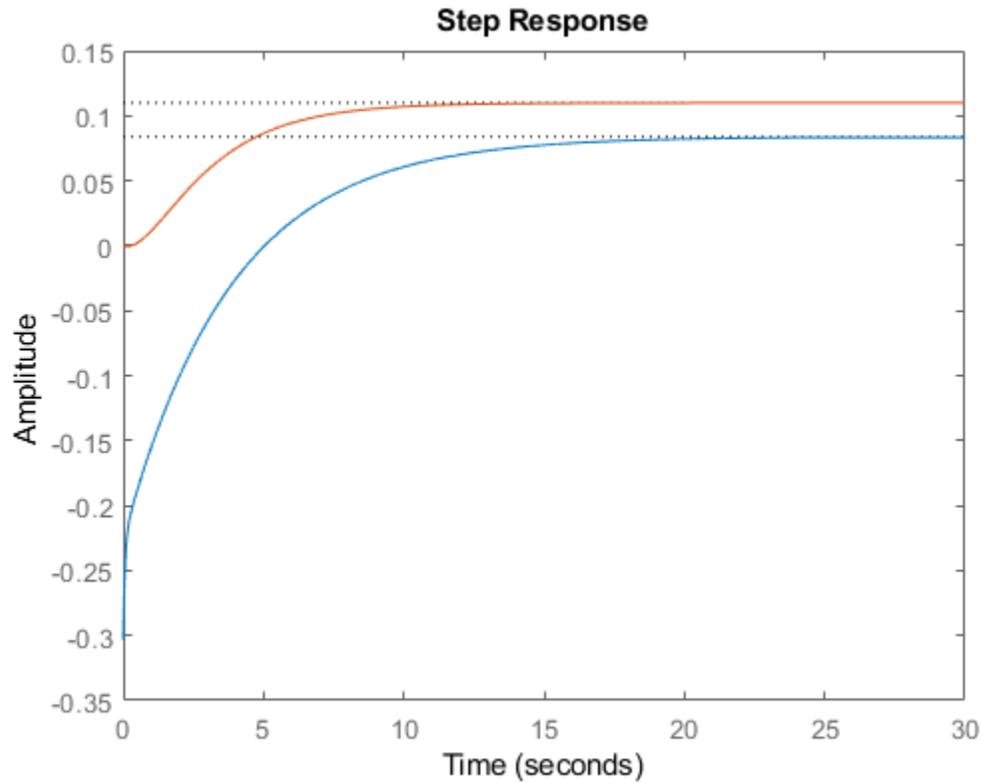
`stepplot(..., dataoptions)` specifies options such as the step amplitude and input offset using the options set, `dataoptions`. Use `stepDataOptions` to create the options set.

## Examples

### Display Normalized Response on Step Plot

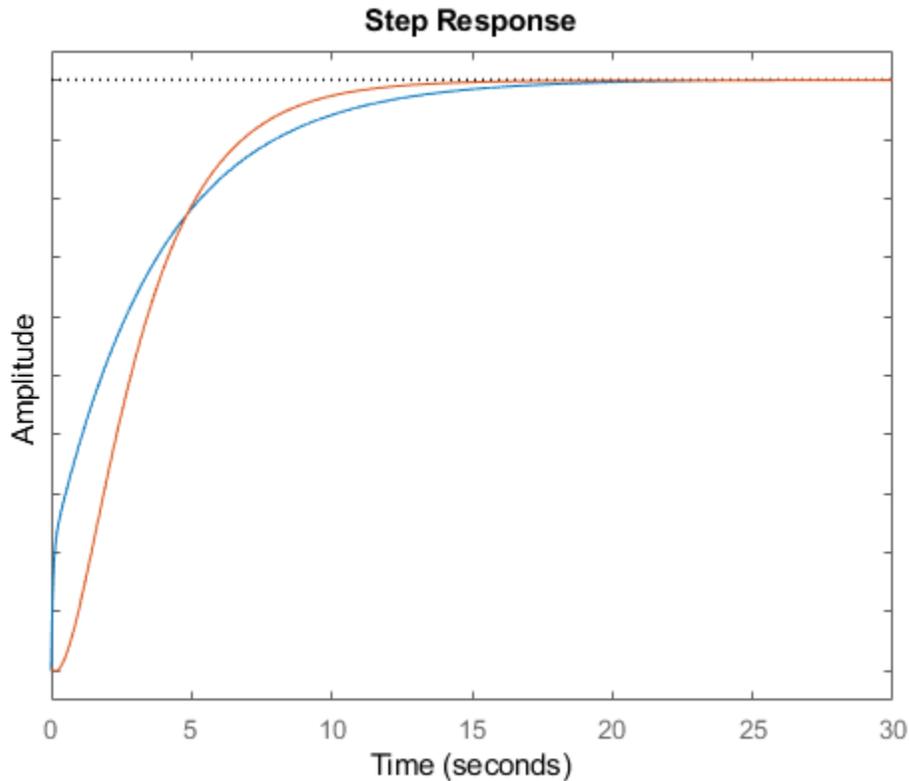
Generate a step response plot for two dynamic systems.

```
sys1 = rss(3);  
sys2 = rss(3);  
h = stepplot(sys1,sys2);
```



Each step response settles at a different steady-state value. Use the plot handle to normalize the plotted response.

```
setoptions(h, 'Normalize', 'on')
```



Now, the responses settle at the same value expressed in arbitrary units.

### **Plot Step Responses of Identified Models with Confidence Region**

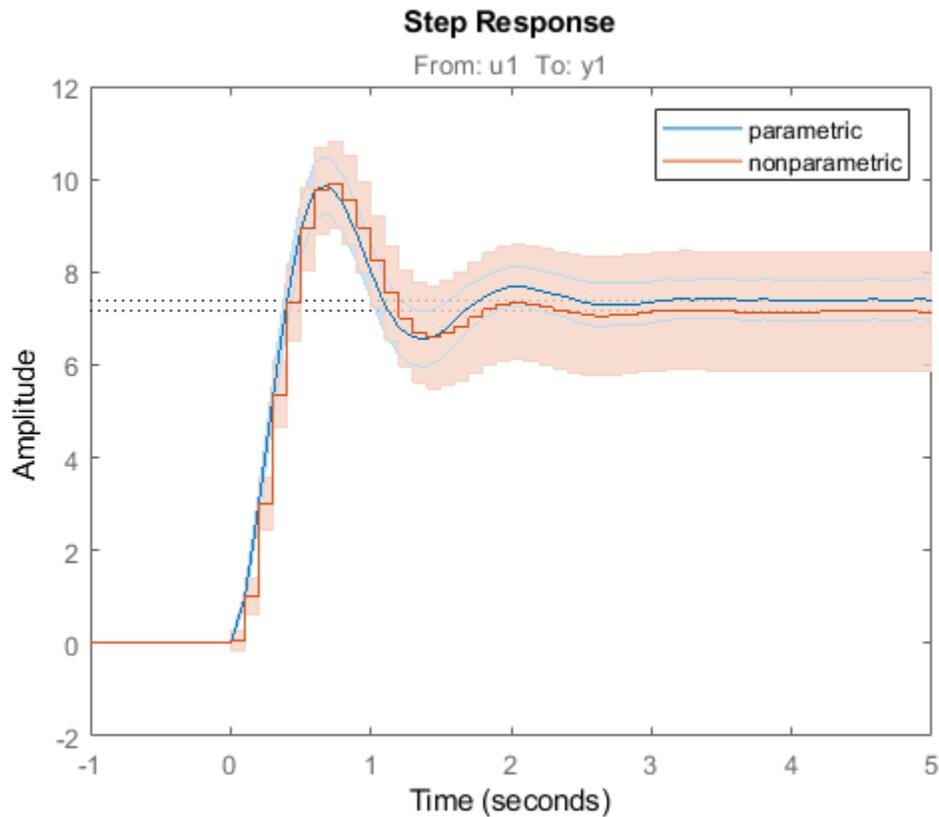
Compare the step response of a parametric identified model to a nonparametric (empirical) model, and view their 3- $\sigma$  confidence regions. (Identified models require System Identification Toolbox™ software.)

Identify a parametric and a nonparametric model from sample data.

```
load iddata1 z1
sys1 = ssest(z1,4);
sys2 = impulseest(z1);
```

Plot the step responses of both identified models. Use the plot handle to display the 3- $\sigma$  confidence regions.

```
t = -1:0.1:5;
h = stepplot(sys1,sys2,t);
showConfidence(h,3)
legend('parametric','nonparametric')
```



The nonparametric model sys2 shows higher uncertainty.

## Plot Step Response of Nonlinear Identified Model

Load data for estimating a nonlinear Hammerstein-Wiener model.

```
load(fullfile(matlabroot, 'toolbox', 'ident', 'iddemos', 'data', 'twotankdata'));  
z = iddata(y,u,0.2, 'Name', 'Two tank system');
```

z is an iddata object that stores the input-output estimation data.

Estimate a Hammerstein-Wiener Model of order [1 5 3] using the estimation data. Specify the input nonlinearity as piecewise linear and output nonlinearity as one-dimensional polynomial.

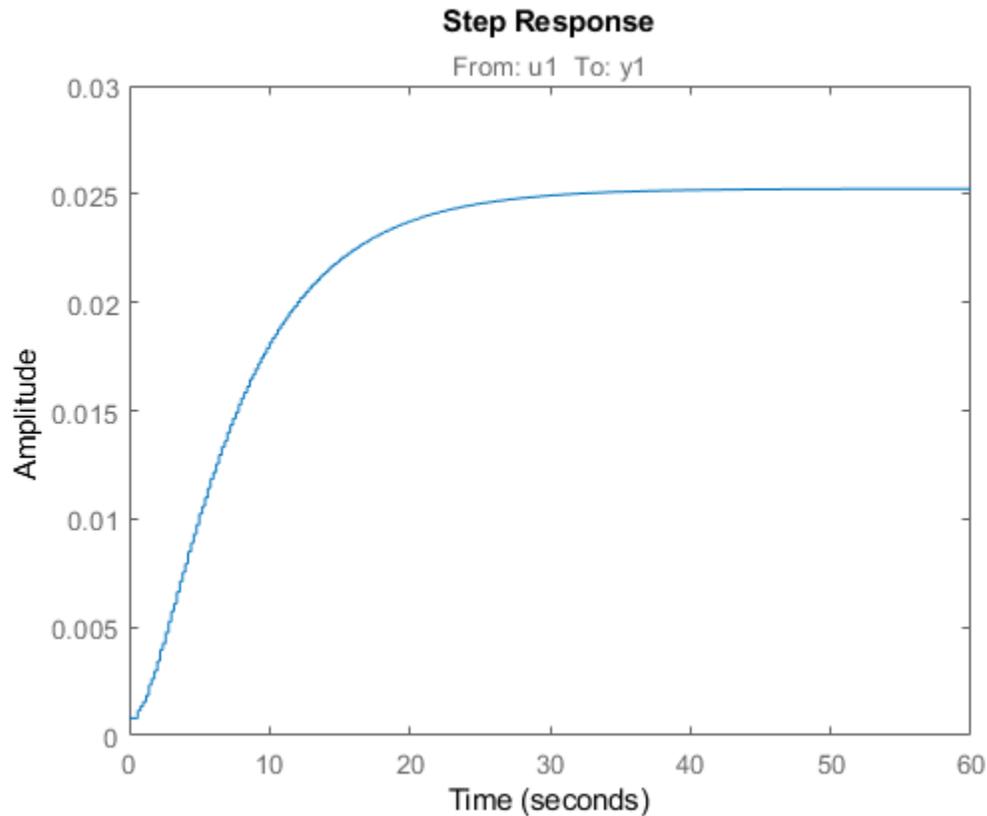
```
sys = nlhw(z, [1 5 3], pwlinear, poly1d);
```

Create an option set to specify input offset and step amplitude level.

```
opt = stepDataOptions('InputOffset', 2, 'StepAmplitude', 0.5);
```

Plot the step response until 60 seconds using the specified options.

```
stepplot(sys, 60, opt);
```



## Tips

You can change the properties of your plot, for example the units. For information on the ways to change properties of your plots, see “Ways to Customize Plots” (Control System Toolbox).

## See Also

`getoptions` | `setoptions` | `showConfidence` | `step`

**Introduced in R2012a**

## strseq

Create sequence of indexed character vectors

### Syntax

```
txtarray = strseq(TXT,INDICES)
```

### Description

`txtarray = strseq(TXT,INDICES)` creates a sequence of indexed character vectors in the cell array `txtarray` by appending the integer values `INDICES` to the character vector `TXT`.

---

**Note** You can use `strvec` to aid in system interconnection. For an example, see the `sumbk` reference page.

---

### Examples

#### Create a Cell Array of Indexed Text

Index the text 'e' with the numbers 1, 2, and 4.

```
txtarray = strseq('e',[1 2 4])
```

```
txtarray = 3x1 cell array
    {'e1'}
    {'e2'}
    {'e4'}
```

## **See Also**

connect | strcat

**Introduced in R2012a**

## struc

Generate model-order combinations for single-output ARX model estimation

### Syntax

```
nn = struc(na,nb,nk)  
nn = struc(na,nb_1,...,nb_nu,nk_1,...,nk_nu)
```

### Description

*nn* = struc(*na*,*nb*,*nk*) generates model-order combinations for single-input, single-output ARX model estimation. *na* and *nb* are row vectors that specify ranges of model orders. *nk* is a row vector that specifies a range of model delays. *nn* is a matrix that contains all combinations of the orders and delays.

*nn* = struc(*na*,*nb\_1*,...,*nb\_nu*,*nk\_1*,...,*nk\_nu*) generates model-order combinations for an ARX model with *nu* input channels.

### Examples

#### Generate Model-Order Combinations and Estimate ARX Model Using IV Method

Create estimation and validation data sets

```
load iddata1;  
ze = z1(1:150);  
zv = z1(151:300);
```

Generate model-order combinations for estimation, specifying ranges for model orders and delays.

```
NN = struc(1:3,1:2,2:4);
```

Estimate ARX models using the instrumental variable method, and compute the loss function for each model order combination.

```
V = ivstruc(ze,zv,NN);
```

Select the model order with the best fit to the validation data.

```
order = selstruc(V,0);
```

Estimate an ARX model of selected order.

```
M = iv4(ze,order);
```

## Generate Model-Order Combinations and Estimate Multi-Input ARX Model

Create estimation and validation data sets.

```
load co2data;  
Ts = 0.5; % Sample time is 0.5 min  
ze = iddata(Output_exp1,Input_exp1,Ts);  
zv = iddata(Output_exp2,Input_exp2,Ts);
```

Generate model-order combinations for:

- na = 2:4
- nb = 2:5 for the first input, and 1 or 4 for the second input.
- nk = 1:4 for the first input, and 0 for the second input.

```
NN = struc(2:4,2:5,[1 4],1:4,0);
```

Estimate an ARX model for each model order combination.

```
V = arxstruc(ze,zv,NN);
```

Select the model order with the best fit to the validation data.

```
order = selstruc(V,0);
```

Estimate an ARX model of selected order.

```
M = arx(ze,order);
```

## Tips

- Use with `arxstruc` or `ivstruc` to compute loss functions for ARX models, one for each model order combination returned by `struc`.

## See Also

`arxstruc` | `ivstruc` | `selstruc`

## Topics

“Estimating Model Orders Using an ARX Model Structure”

“Preliminary Step - Estimating Model Orders and Input Delays”

**Introduced before R2006a**

## System Identification

Identify models of dynamic systems from measured data

### Description

The **System Identification** app enables you to identify models of dynamic systems from measured input-output data. You can estimate both linear and nonlinear models and compare responses of different models.

Using this app you can:

- Import, plot, and preprocess measured input-output data.
- Estimate linear models such as transfer functions, process models, polynomial models, and state-space models using time-domain, time series, or frequency-domain data.
- Estimate nonlinear ARX and Hammerstein-Wiener models using time-domain data.
- Validate estimated models using independent data sets.
- Export estimated models for further analysis to MATLAB workspace or to the **Linear System Analyzer** app in Control System Toolbox.

For more information, in the **System Identification** app, select **Help > System Identification App Help**.

### Open the System Identification App

- MATLAB Toolstrip: On the **Apps** tab, under **Control System Design and Analysis**, click the app icon.
- MATLAB command prompt: Enter `systemIdentification`.

### Examples

- “Working with System Identification App”
- “Identify Linear Models Using System Identification App”

- “Identify Nonlinear Black-Box Models Using System Identification App”

## Programmatic Use

`systemIdentification` opens the **System Identification** app. If the app is already open, the command brings the app into focus.

`systemIdentification(sessionFile)` opens the app and loads a previously saved session file, `sessionFile`, on the MATLAB path. A session includes data sets and models in the app at the time of saving. If the app is already open, the command merges the contents of the session file with those already present in the app.

For example, `systemIdentification('mySession')` opens the app and loads the previously saved app session `mySession.sid`.

To save a session, in the **System Identification** app, select **File > Save session**. The session is saved to a file with a `.sid` extension.

`systemIdentification(sessionFile,path)` specifies the path to the session file. Use this syntax if the file is not on the MATLAB path.

For example, `systemIdentification('mySession','C:\matlab\work')` opens the app and loads the previously saved app session `mySession.sid` located at `C:\matlab\work`.

## See Also

### Functions

`arx` | `iddata` | `midprefs` | `nlarx` | `nllhw` | `polyest` | `procest` | `ssest` | `tfest`

### Topics

“Working with System Identification App”

“Identify Linear Models Using System Identification App”

“Identify Nonlinear Black-Box Models Using System Identification App”

**Introduced before R2006a**

## tfdata

Access transfer function data

### Syntax

```
[num,den] = tfdata(sys)
[num,den,Ts] = tfdata(sys)
[num,den,Ts,sdnum,sdden]=tfdata(sys)
[num,den,Ts,...]=tfdata(sys,J1,...,Jn)
```

### Description

`[num,den] = tfdata(sys)` returns the numerator(s) and denominator(s) of the transfer function for the TF, SS or ZPK model (or LTI array of TF, SS or ZPK models) `sys`. For single LTI models, the outputs `num` and `den` of `tfdata` are cell arrays with the following characteristics:

- `num` and `den` have as many rows as outputs and as many columns as inputs.
- The  $(i, j)$  entries `num{i, j}` and `den{i, j}` are row vectors specifying the numerator and denominator coefficients of the transfer function from input `j` to output `i`. These coefficients are ordered in *descending* powers of  $s$  or  $z$ .

For arrays `sys` of LTI models, `num` and `den` are multidimensional cell arrays with the same sizes as `sys`.

If `sys` is a state-space or zero-pole-gain model, it is first converted to transfer function form using `tf`. For more information on the format of transfer function model data, see the `tf` reference page.

For SISO transfer functions, the syntax

```
[num,den] = tfdata(sys,'v')
```

forces `tfdata` to return the numerator and denominator directly as row vectors rather than as cell arrays (see example below).

```
[num,den,Ts] = tfdata(sys)
```

 also returns the sample time `Ts`.

`[num,den,Ts,sdnum,sdden]=tfdata(sys)` also returns the uncertainties in the numerator and denominator coefficients of identified system `sys`. `sdnum{i,j}(k)` is the 1 standard uncertainty in the value `num{i,j}(k)` and `sdden{i,j}(k)` is the 1 standard uncertainty in the value `den{i,j}(k)`. If `sys` does not contain uncertainty information, `sdnum` and `sdden` are empty (`[]`).

`[num,den,Ts,...]=tfdata(sys,J1,...,Jn)` extracts the data for the `(J1,...,JN)` entry in the model array `sys`.

You can access the remaining LTI properties of `sys` with `get` or by direct referencing, for example,

```
sys.Ts
sys.variable
```

## Examples

### Example 1

Given the SISO transfer function

```
h = tf([1 1],[1 2 5])
```

you can extract the numerator and denominator coefficients by typing

```
[num,den] = tfdata(h,'v')
num =
     0     1     1
den =
     1     2     5
```

This syntax returns two row vectors.

If you turn `h` into a MIMO transfer function by typing

```
H = [h ; tf(1,[1 1])]
```

the command

```
[num,den] = tfdata(H)
```

now returns two cell arrays with the numerator/denominator data for each SISO entry. Use `celldisp` to visualize this data. Type

```
celldisp(num)
```

This command returns the numerator vectors of the entries of H.

```
num{1} =  
    0     1     1
```

```
num{2} =  
    0     1
```

Similarly, for the denominators, type

```
celldisp(den)  
den{1} =  
    1     2     5
```

```
den{2} =  
    1     1
```

## Example 2

Extract the numerator, denominator and their standard deviations for a 2-input, 1 output identified transfer function.

```
load iddata7
```

```
transfer function model
```

```
sys1 = tfest(z7, 2, 1, 'InputDelay',[1 0]);
```

```
an equivalent process model
```

```
sys2 = procest(z7, {'P2UZ', 'P2UZ'}, 'InputDelay',[1 0]);
```

```
[num1, den1, ~, dnum1, dden1] = tfdata(sys1);
```

```
[num2, den2, ~, dnum2, dden2] = tfdata(sys2);
```

## See Also

[get](#) | [ssdata](#) | [tf](#) | [zpkdata](#)

**Introduced before R2006a**

## tfest

Transfer function estimation

### Syntax

```
sys = tfest(data,np)
sys = tfest(data,np,nz)
sys = tfest(data,np,nz,iodelay)
sys = tfest( ____,Name,Value)
sys = tfest(data,init_sys)
sys = tfest( ____,opt)
```

### Description

`sys = tfest(data,np)` estimates a continuous-time transfer function, `sys`, using time- or frequency-domain data, `data`, and contains `np` poles. The number of zeros in the `sys` is `max(np-1,0)`.

`sys = tfest(data,np,nz)` estimates a transfer function containing `nz` zeros.

`sys = tfest(data,np,nz,iodelay)` estimates a transfer function with transport delay for input/output pairs `iodelay`.

`sys = tfest( ____,Name,Value)` uses additional options specified by one or more `Name,Value` pair arguments. All input arguments described for previous syntaxes also apply here.

`sys = tfest(data,init_sys)` uses the linear system `init_sys` to configure the initial parameterization of `sys`.

`sys = tfest( ____,opt)` specifies the estimation behavior using the option set `opt`. All input arguments described for previous syntaxes also apply here.

## Input Arguments

### data

Estimation data.

For time domain estimation, `data` is an `iddata` object containing the input and output signal values.

Time-series models, which are models that contain no measured inputs, cannot be estimated using `tfest`. Use `ar`, `arx` or `armax` for time-series models instead.

For frequency domain estimation, `data` can be one of the following:

- `frd` or `idfrd` object that represents recorded frequency response data:
  - Complex-values  $G(e^{i\omega})$ , for given frequencies  $\omega$
  - Amplitude  $|G|$  and phase shift  $\phi = \arg G$  values
- `iddata` object with its properties specified as follows:
  - `InputData` — Fourier transform of the input signal
  - `OutputData` — Fourier transform of the output signal
  - `Domain` — 'Frequency'

For multi-experiment data, the sample times and intersample behavior of all the experiments must match.

### np

Number of poles in the estimated transfer function.

`np` is a nonnegative number.

For systems that are multiple-input, or multiple-output, or both:

- To use the same number of poles for all the input/output pairs, specify `np` as a scalar.
- To use different number of poles for the input/output pairs, specify `np` as an  $ny$ -by- $nu$  matrix.  $ny$  is the number of outputs, and  $nu$  is the number of inputs.

**nz**

Number of zeros in the estimated transfer function.

`nz` is a nonnegative number.

For systems that are multiple-input, or multiple-output, or both:

- To use the same number of zeros for all the input/output pairs, specify `nz` as a scalar.
- To use a different number of zeros for the input/output pairs, specify `nz` as an  $N_y$ -by- $N_u$  matrix.  $N_y$  is the number of outputs, and  $N_u$  is the number of inputs.

For a continuous-time model, estimated using discrete-time data, set `nz`  $\leq$  `np`.

For discrete-time model estimation, specify `nz` as the number of zeros of the numerator polynomial of the transfer function. For example, `tfest(data,2,1,'Ts',data.Ts)`

estimates a transfer function of the form  $b_1z^{-1} / (1 + a_1z^{-1} + b_2z^{-2})$ , while `tfest(data,2,2,'Ts',data.Ts)` estimates  $(b_1z^{-1} + b_2z^{-2}) / (1 + a_1z^{-1} + b_2z^{-2})$ . Here  $z^{-1}$  is the Z-transform lag variable. For more information about discrete-time transfer functions, see “Discrete-Time Representation”. For an example, see “Estimate Discrete-Time Transfer Function” on page 1-1776.

**iodelay**

Transport delay.

For continuous-time systems, specify transport delays in the time unit stored in the `TimeUnit` property of `data`. For discrete-time systems, specify transport delays as integers denoting delay of a multiple of the sample time `Ts`.

For a MIMO system with  $N_y$  outputs and  $N_u$  inputs, set `iodelay` to an  $N_y$ -by- $N_u$  array. Each entry of this array is a numerical value that represents the transport delay for the corresponding input/output pair. You can also set `iodelay` to a scalar value to apply the same delay to all input/output pairs.

The specified values are treated as fixed delays.

`iodelay` must contain either nonnegative numbers or NaNs. Use NaN in the `iodelay` matrix to denote unknown transport delays.

Use `[]` or `0` to indicate that there is no transport delay.

## **opt**

Estimation options.

**opt** is an options set, created using `tfestOptions`, that specifies estimation options including:

- Estimation objective
- Handling of initial conditions
- Numerical search method to be used in estimation

## **init\_sys**

Linear system that configures the initial parameterization of `sys`.

You obtain `init_sys` by either performing an estimation using measured data or by direct construction.

If `init_sys` is an `idtf` model, `tfest` uses the parameters and constraints defined in `init_sys` as the initial guess for estimating `sys`. Use the `Structure` property of `init_sys` to configure initial guesses and constraints for the numerator, denominator, and transport lag. For example:

- To specify an initial guess for the numerator of `init_sys`, set `init_sys.Structure.Numerator.Value` to the initial guess.
- To specify constraints for the numerator of `init_sys`:
  - Set `init_sys.Structure.Numerator.Minimum` to the minimum numerator coefficient values
  - Set `init_sys.Structure.Numerator.Maximum` to the maximum numerator coefficient values
  - Set `init_sys.Structure.Numerator.Free` to indicate which numerator coefficients are free for estimation

If `init_sys` is not an `idtf` model, the software first converts `init_sys` to a transfer function. `tfest` uses the parameters of the resulting model as the initial guess for estimation.

If `opt` is not specified, and `init_sys` was obtained by estimation, then the estimation options from `init_sys.Report.OptionsUsed` are used.

## Name-Value Pair Arguments

Specify optional comma-separated pairs of `Name`, `Value` arguments. `Name` is the argument name and `Value` is the corresponding value. `Name` must appear inside quotes. You can specify several name and value pair arguments in any order as `Name1, Value1, . . . , NameN, ValueN`.

### **Ts**

Sample time.

Use the following values for `Ts`:

- `0` — Continuous-time model.
- `data.Ts` — Discrete-time model. In this case, `np` and `nz` refer to the number of roots of  $z^{-1}$  for the numerator and denominator polynomials.

**Default:** `0`

### **InputDelay**

Input delay for each input channel, specified as a scalar value or numeric vector. For continuous-time systems, specify input delays in the time unit stored in the `TimeUnit` property. For discrete-time systems, specify input delays in integer multiples of the sample time `Ts`. For example, `InputDelay = 3` means a delay of three sample times.

For a system with `Nu` inputs, set `InputDelay` to an `Nu`-by-1 vector. Each entry of this vector is a numerical value that represents the input delay for the corresponding input channel.

You can also set `InputDelay` to a scalar value to apply the same delay to all channels.

**Default:** `0`

### **Feedthrough**

Feedthrough for discrete-time transfer function, specified as an `Ny`-by-`Nu` logical matrix. `Ny` is the number of outputs, and `Nu` is the number of inputs. To use the same feedthrough for all input-output channels, specify `Feedthrough` as a scalar.

Consider a discrete-time model with two poles and three zeros:

$$H(z^{-1}) = \frac{b_0 + b_1z^{-1} + b_2z^{-2} + b_3z^{-3}}{1 + a_1z^{-1} + a_2z^{-2}}$$

When the model has direct feedthrough,  $b_0$  is a free parameter whose value is estimated along with the rest of the model parameters  $b_1$ ,  $b_2$ ,  $b_3$ ,  $a_1$ ,  $a_2$ . When the model has no feedthrough,  $b_0$  is fixed to zero. For an example, see “Estimate Discrete-Time Transfer Function With Feedthrough” on page 1-1777.

**Default:** false ( $N_y, N_u$ )

## Output Arguments

**sys**

Identified transfer function, returned as an `idtf` model. This model is created using the specified model orders, delays and estimation options.

Information about the estimation results and options used is stored in the `Report` property of the model. `Report` has the following fields:

Report Field	Description
Status	Summary of the model status, which indicates whether the model was created by construction or obtained by estimation.
Method	Estimation command used.
InitializeMethod	<p>Algorithm used to initialize the numerator and denominator for estimation of continuous-time transfer functions using time-domain data, returned as one of the following values:</p> <ul style="list-style-type: none"> <li>• 'iv' — Instrument Variable approach.</li> <li>• 'svf' — State Variable Filters approach.</li> <li>• 'gpmf' — Generalized Poisson Moment Functions approach.</li> <li>• 'n4sid' — Subspace state-space estimation approach.</li> </ul> <p>This field is especially useful to view the algorithm used when the <code>InitializeMethod</code> option in the estimation option set is 'all'.</p>

<b>Report Field</b>	<b>Description</b>
N4Weight	<p>Weighting matrices used in the singular-value decomposition step when InitializeMethod is 'n4sid', returned as one of the following values:</p> <ul style="list-style-type: none"> <li>• 'MOESP' — Uses the MOESP algorithm by Verhaegen.</li> <li>• 'CVA' — Uses the canonical variable algorithm (CVA) by Larimore.</li> <li>• 'SSARX' — A subspace identification method that uses an ARX estimation based algorithm to compute the weighting.</li> </ul> <p>This field is especially useful to view the weighting matrices used when the N4Weight option in the estimation option set is 'auto'.</p>
N4Horizon	<p>Forward and backward prediction horizons used when InitializeMethod is 'n4sid', returned as a row vector with three elements — [r sy su], where r is the maximum forward prediction horizon. sy is the number of past outputs, and su is the number of past inputs that are used for the predictions.</p>
InitialCondition	<p>Handling of initial conditions during model estimation, returned as one of the following values:</p> <ul style="list-style-type: none"> <li>• 'zero' — The initial conditions were set to zero.</li> <li>• 'estimate' — The initial conditions were treated as independent estimation parameters.</li> <li>• 'backcast' — The initial conditions were estimated using the best least squares fit.</li> </ul> <p>This field is especially useful to view how the initial conditions were handled when the InitialCondition option in the estimation option set is 'auto'.</p>

Report Field	Description																		
Fit	<p>Quantitative assessment of the estimation, returned as a structure. See “Loss Function and Model Quality Metrics” for more information on these quality metrics. The structure has the following fields:</p> <table border="1" data-bbox="387 456 1328 951"> <thead> <tr> <th data-bbox="387 456 516 499">Field</th> <th data-bbox="516 456 1328 499">Description</th> </tr> </thead> <tbody> <tr> <td data-bbox="387 499 516 609">FitPercent</td> <td data-bbox="516 499 1328 609">Normalized root mean squared error (NRMSE) measure of how well the response of the model fits the estimation data, expressed as a percentage.</td> </tr> <tr> <td data-bbox="387 609 516 652">LossFcn</td> <td data-bbox="516 609 1328 652">Value of the loss function when the estimation completes.</td> </tr> <tr> <td data-bbox="387 652 516 730">MSE</td> <td data-bbox="516 652 1328 730">Mean squared error (MSE) measure of how well the response of the model fits the estimation data.</td> </tr> <tr> <td data-bbox="387 730 516 774">FPE</td> <td data-bbox="516 730 1328 774">Final prediction error for the model.</td> </tr> <tr> <td data-bbox="387 774 516 817">AIC</td> <td data-bbox="516 774 1328 817">Raw Akaike Information Criteria (AIC) measure of model quality.</td> </tr> <tr> <td data-bbox="387 817 516 861">AICc</td> <td data-bbox="516 817 1328 861">Small sample-size corrected AIC.</td> </tr> <tr> <td data-bbox="387 861 516 904">nAIC</td> <td data-bbox="516 861 1328 904">Normalized AIC.</td> </tr> <tr> <td data-bbox="387 904 516 951">BIC</td> <td data-bbox="516 904 1328 951">Bayesian Information Criteria (BIC).</td> </tr> </tbody> </table>	Field	Description	FitPercent	Normalized root mean squared error (NRMSE) measure of how well the response of the model fits the estimation data, expressed as a percentage.	LossFcn	Value of the loss function when the estimation completes.	MSE	Mean squared error (MSE) measure of how well the response of the model fits the estimation data.	FPE	Final prediction error for the model.	AIC	Raw Akaike Information Criteria (AIC) measure of model quality.	AICc	Small sample-size corrected AIC.	nAIC	Normalized AIC.	BIC	Bayesian Information Criteria (BIC).
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Parameters	Estimated values of model parameters.																		
OptionsUsed	Option set used for estimation. If no custom options were configured, this is a set of default options. See <code>polyestOptions</code> for more information.																		
RandState	State of the random number stream at the start of estimation. Empty, <code>[]</code> , if randomization was not used during estimation. For more information, see <code>rng</code> in the MATLAB documentation.																		

Report Field	Description																
DataUsed	Attributes of the data used for estimation, returned as a structure with the following fields:																
	<table border="1"> <thead> <tr> <th data-bbox="385 425 516 465">Field</th> <th data-bbox="516 425 1335 465">Description</th> </tr> </thead> <tbody> <tr> <td data-bbox="385 465 516 512">Name</td> <td data-bbox="516 465 1335 512">Name of the data set.</td> </tr> <tr> <td data-bbox="385 512 516 558">Type</td> <td data-bbox="516 512 1335 558">Data type.</td> </tr> <tr> <td data-bbox="385 558 516 605">Length</td> <td data-bbox="516 558 1335 605">Number of data samples.</td> </tr> <tr> <td data-bbox="385 605 516 652">Ts</td> <td data-bbox="516 605 1335 652">Sample time.</td> </tr> <tr> <td data-bbox="385 652 516 956">InterSample</td> <td data-bbox="516 652 1335 956">                     Input intersample behavior, returned as one of the following values:                     <ul style="list-style-type: none"> <li>• 'zoh' — Zero-order hold maintains a piecewise-constant input signal between samples.</li> <li>• 'foh' — First-order hold maintains a piecewise-linear input signal between samples.</li> <li>• 'bl' — Band-limited behavior specifies that the continuous-time input signal has zero power above the Nyquist frequency.</li> </ul> </td> </tr> <tr> <td data-bbox="385 956 516 1038">InputOffset</td> <td data-bbox="516 956 1335 1038">Offset removed from time-domain input data during estimation. For nonlinear models, it is [ ].</td> </tr> <tr> <td data-bbox="385 1038 516 1112">OutputOffset</td> <td data-bbox="516 1038 1335 1112">Offset removed from time-domain output data during estimation. For nonlinear models, it is [ ].</td> </tr> </tbody> </table>	Field	Description	Name	Name of the data set.	Type	Data type.	Length	Number of data samples.	Ts	Sample time.	InterSample	Input intersample behavior, returned as one of the following values: <ul style="list-style-type: none"> <li>• 'zoh' — Zero-order hold maintains a piecewise-constant input signal between samples.</li> <li>• 'foh' — First-order hold maintains a piecewise-linear input signal between samples.</li> <li>• 'bl' — Band-limited behavior specifies that the continuous-time input signal has zero power above the Nyquist frequency.</li> </ul>	InputOffset	Offset removed from time-domain input data during estimation. For nonlinear models, it is [ ].	OutputOffset	Offset removed from time-domain output data during estimation. For nonlinear models, it is [ ].
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Report Field	Description	
Termination	Termination conditions for the iterative search used for prediction error minimization. Structure with the following fields:	
	Field	Description
	WhyStop	Reason for terminating the numerical search.
	Iterations	Number of search iterations performed by the estimation algorithm.
	FirstOrderOptimality	$\infty$ -norm of the gradient search vector when the search algorithm terminates.
	FcnCount	Number of times the objective function was called.
	UpdateNorm	Norm of the gradient search vector in the last iteration. Omitted when the search method is 'lsqnonlin' or 'fmincon'.
	LastImprovement	Criterion improvement in the last iteration, expressed as a percentage. Omitted when the search method is 'lsqnonlin' or 'fmincon'.
	Algorithm	Algorithm used by 'lsqnonlin' or 'fmincon' search method. Omitted when other search methods are used.
For estimation methods that do not require numerical search optimization, the Termination field is omitted.		

For more information on using Report, see “Estimation Report”.

## Examples

### Estimate Transfer Function Model By Specifying Number of Poles

Load time-domain system response data and use it to estimate a transfer function for the system.

```
load iddata1 z1;  
np = 2;  
sys = tfest(z1,np);
```

`z1` is an `iddata` object that contains time-domain, input-output data.

`np` specifies the number of poles in the estimated transfer function.

`sys` is an `idtf` model containing the estimated transfer function.

To see the numerator and denominator coefficients of the resulting estimated model `sys`, enter:

```
sys.Numerator
```

```
ans = 1×2
```

```
    2.4554    176.9856
```

```
sys.Denominator
```

```
ans = 1×3
```

```
    1.0000    3.1625   23.1631
```

To view the uncertainty in the estimates of the numerator and denominator and other information, use `tfdata`.

## Specify Number of Poles and Zeros in Estimated Transfer Function

Load time-domain system response data and use it to estimate a transfer function for the system.

```
load iddata2 z2;  
np = 2;  
nz = 1;  
sys = tfest(z2,np,nz);
```

`z2` is an `iddata` object that contains time-domain system response data.

`np` and `nz` specify the number of poles and zeros in the estimated transfer function, respectively.

`sys` is an `idtf` model containing the estimated transfer function.

### **Estimate Transfer Function Containing Known Transport Delay**

Load time-domain system response data and use it to estimate a transfer function for the system. Specify a known transport delay for the transfer function.

```
load iddata2 z2;  
np = 2;  
nz = 1;  
iodelay = 0.2;  
sys = tfest(z2,np,nz,iodelay);
```

`z2` is an `iddata` object that contains time-domain system response data.

`np` and `nz` specify the number of poles and zeros in the estimated transfer function, respectively.

`iodelay` specifies the transport delay for the estimated transfer function as 0.2 seconds.

`sys` is an `idtf` model containing the estimated transfer function, with `IODelay` property set to 0.2 seconds.

### **Estimate Transfer Function Containing Unknown Transport Delay**

Load time-domain system response data and use it to estimate a transfer function for the system. Specify an unknown transport delay for the transfer function.

```
load iddata2 z2;  
np = 2;  
nz = 1;  
iodelay = NaN;  
sys = tfest(z2,np,nz,iodelay);
```

`z2` is an `iddata` object that contains time-domain system response data.

`np` and `nz` specify the number of poles and zeros in the estimated transfer function, respectively.

`iodelay` specifies the transport delay for the estimated transfer function. `iodelay = NaN` denotes the transport delay as an unknown parameter to be estimated.

`sys` is an `idtf` model containing the estimated transfer function, whose `IODElay` property is estimated using data.

## Estimate Discrete-Time Transfer Function

Load time-domain system response data.

```
load iddata2 z2;
```

`z2` is an `iddata` object that contains the time-domain system response data.

Estimate a discrete-time transfer function with two poles and one zero. Specify the sample time as 0.1 sec and the transport delay as 2 sec.

```
np = 2;  
nz = 1;  
iodelay = 2;  
Ts = 0.1;  
sysd = tfest(z2,np,nz,iodelay,'Ts',Ts)
```

```
sysd =
```

```
From input "u1" to output "y1":  
1.8 z^-1  
z^(-2) * -----  
1 - 1.418 z^-1 + 0.6613 z^-2
```

```
Sample time: 0.1 seconds  
Discrete-time identified transfer function.
```

```
Parameterization:
```

```
Number of poles: 2   Number of zeros: 1
```

```
Number of free coefficients: 3
```

```
Use "tfdata", "getpvec", "getcov" for parameters and their uncertainties.
```

```
Status:
```

Estimated using TFEST on time domain data "z2".  
 Fit to estimation data: 80.26%  
 FPE: 2.095, MSE: 2.063

By default, the model has no feedthrough, and the numerator polynomial of the estimated transfer function has a zero leading coefficient  $b_0$ . To estimate  $b_0$ , specify the Feedthrough property during estimation.

### Estimate Discrete-Time Transfer Function With Feedthrough

Load the estimation data.

```
load iddata5 z5
```

First estimate a discrete-time transfer function model with two poles, one zero, and no feedthrough.

```
np = 2;  
nz = 1;  
model = tfest(z5,np,nz,'Ts',z5.Ts);
```

The estimated transfer function has the form:

$$H(z^{-1}) = \frac{b_1 z^{-1} + b_2 z^{-2}}{1 + a_1 z^{-1} + a_2 z^{-2}}$$

By default, the model has no feedthrough, and the numerator polynomial of the estimated transfer function has a zero leading coefficient  $b_0$ . To estimate  $b_0$ , specify the Feedthrough property during estimation.

```
model = tfest(z5,np,nz,'Ts',z5.Ts,'Feedthrough',true);
```

The numerator polynomial of the estimated transfer function now has a nonzero leading coefficient:

$$H(z^{-1}) = \frac{b_0 + b_1 z^{-1} + b_2 z^{-2}}{1 + a_1 z^{-1} + a_2 z^{-2}}$$

## Analyze the Origin of Delay in Measured Data

Compare two discrete-time models with and without feedthrough and transport delay.

If there is a delay from the measured input to output, it can be attributed to a lack of feedthrough or to a true transport delay. For discrete-time models, absence of feedthrough corresponds to a lag of 1 sample between the input and output. Estimating a model with `Feedthrough = false` and `IODelay = 0` thus produces a discrete-time system that is equivalent to a system with `Feedthrough = true` and `IODelay = 1`. Both systems show the same time- and frequency-domain responses, for example, on step and Bode plots. However, you get different results if you reduce these models using `balred` or convert them to their continuous-time representation. Therefore, you should check if the observed delay should be attributed to transport delay or to a lack of feedthrough.

Estimate a discrete-time model with no feedthrough.

```
load iddata1 z1
np = 2;
nz = 2;
model1 = tfest(z1,np,nz,'Ts',z1.Ts);
```

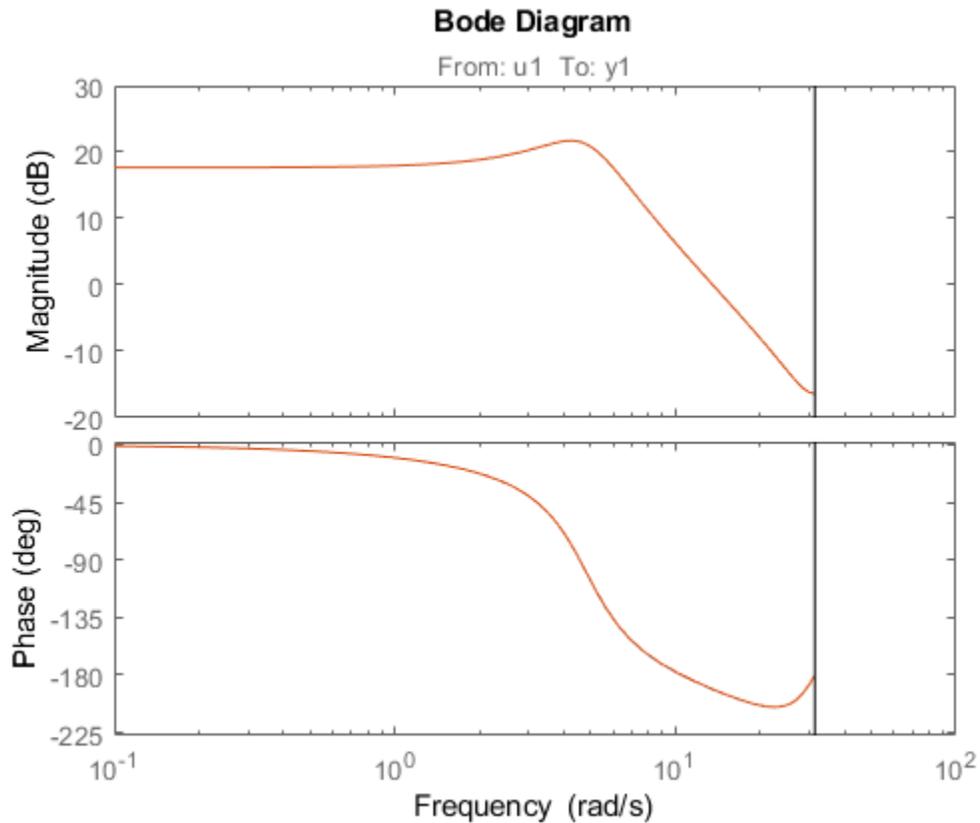
`model1` has a transport delay of 1 sample and its `IODelay` property is 0. Its numerator polynomial begins with  $z^{-1}$ .

Estimate another discrete-time model with feedthrough and 1 sample input-output delay.

```
model2 = tfest(z1,np,nz-1,1,'Ts',z1.Ts,'Feedthrough',true);
```

Compare the Bode response of the models.

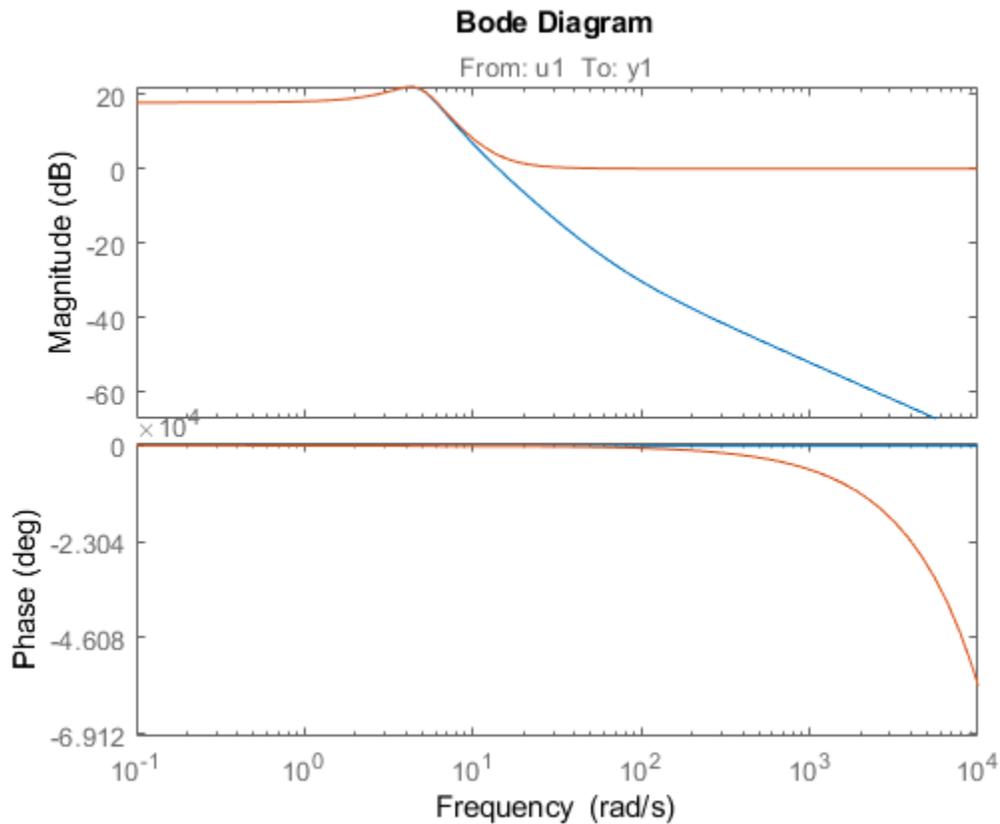
```
bode(model1,model2);
```



The equations for `model1` and `model2` are equivalent, but the transport delay of `model2` has been absorbed into the numerator of `model1`.

Convert the models to continuous time, and compare their Bode responses.

```
bode(d2c(model1),d2c(model2));
```



As the plot shows, the Bode responses of the two models do not match when you convert them to continuous time.

### **Estimate MISO Discrete-Time Transfer Function with Feedthrough and Delay Specifications for Individual Channels**

Estimate a 2-input, 1-output discrete-time transfer function with a delay of 2 samples on first input and zero seconds on the second input. Both inputs have no feedthrough.

Split data into estimation and validation data sets.

```
load iddata7 z7
ze = z7(1:300);
zv = z7(200:400);
```

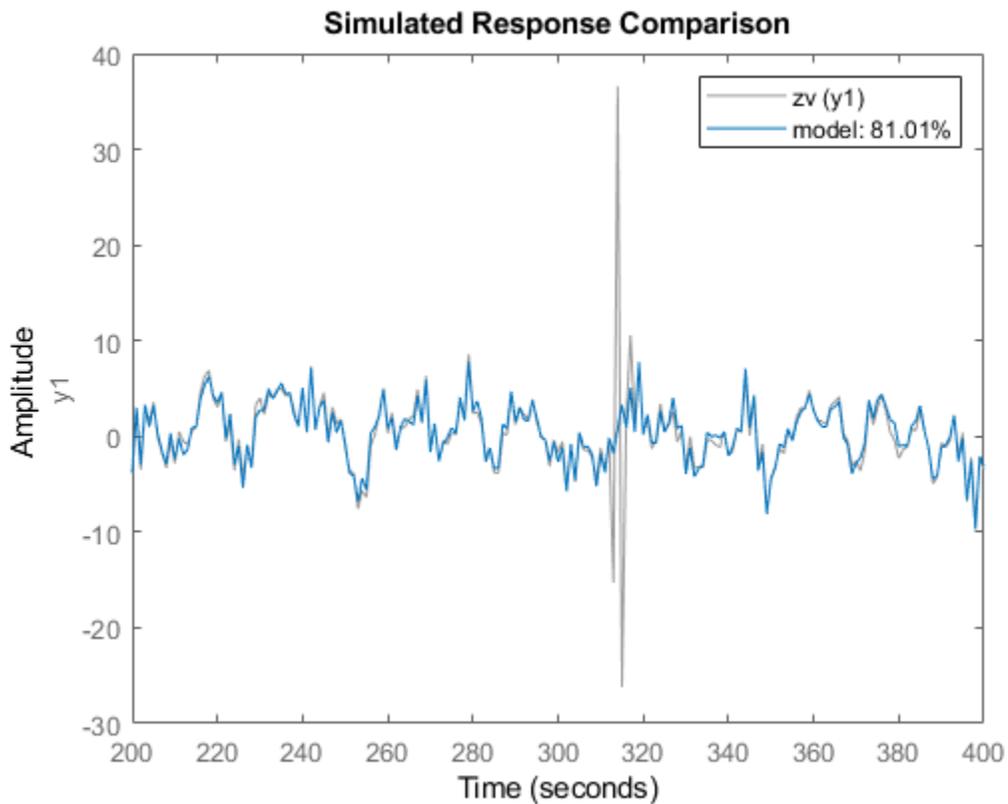
Estimate a 2-input, 1-output transfer function with 2 poles and 1 zero for each input-to-output transfer function.

```
Lag = [2;0];
Ft = [false,false];
model = tfest(ze,2,1,'Ts',z7.Ts,'Feedthrough',Ft,'InputDelay',Lag);
```

Choice of `Feedthrough` dictates whether the leading numerator coefficient is zero (no feedthrough) or not (nonzero feedthrough). Delays are expressed separately using `InputDelay` or `IODelay` property. This example uses `InputDelay` to express the delays.

Validate the estimated model. Exclude the data outliers for validation.

```
I = 1:201;
I(114:118) = [];
opt = compareOptions('Samples',I);
compare(zv,model,opt)
```



### Estimate Transfer Function Model Using Regularized Impulse Response Model

Identify a 15th order transfer function model by using regularized impulse response estimation

Load data.

```
load regularizationExampleData m0simdata;
```

Obtain regularized impulse response (FIR) model.

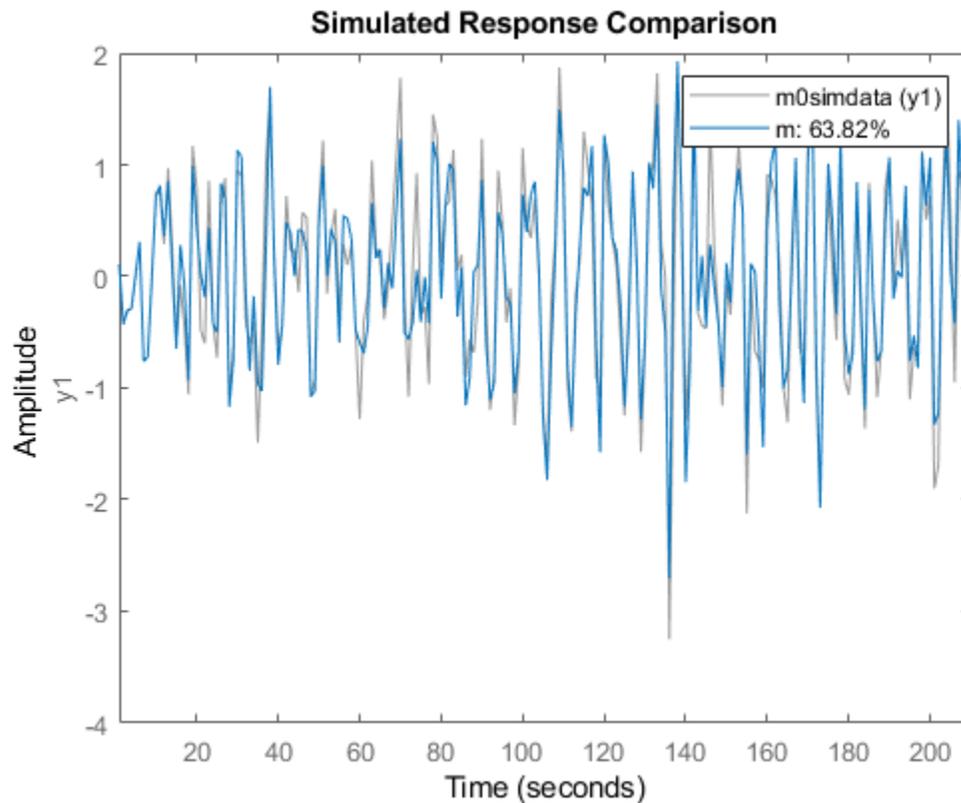
```
opt = impulseestOptions('RegularizationKernel','DC');  
m0 = impulseest(m0simdata,70,opt);
```

Convert model into a transfer function model after reducing order to 15.

```
m = idtf(balred(idss(m0),15));
```

Compare the model output with data.

```
compare(m0simdata,m);
```



## Estimate Transfer Function Using an Estimation Option Set

Create the option set for `tfest`.

```
opt = tfestOptions('InitializeMethod','n4sid','Display','on','SearchMethod','lsqnonlin
```

`opt` specifies that the initialization method as `'n4sid'`, and the search method as `'lsqnonlin'`. It also specifies that the loss-function values for each iteration be shown.

Load time-domain system response data and use it to estimate a transfer function for the system. Specify the estimation options using `opt`.

```
load iddata2 z2;  
np = 2;  
nz = 1;  
iodelay = 0.2;  
sysc = tfest(z2,np,nz,iodelay,opt);
```

`z2` is an `iddata` object that contains time-domain system response data.

`np` and `nz` specify the number of poles and zeros in the estimated transfer function, respectively.

`iodelay` specifies the transport delay for the estimated transfer function as 0.2 seconds.

`opt` specifies the estimation options.

`sys` is an `idtf` model containing the estimated transfer function.

## Specify Model Properties of the Estimated Transfer Function

Load time-domain system response data, and use it to estimate a transfer function for the system. Specify the input delay for the estimated transfer function.

```
load iddata2 z2;  
np = 2;  
nz = 1;  
input_delay = 0.2;  
sys = tfest(z2,np,nz,'InputDelay',input_delay);
```

`z2` is an `iddata` object that contains time-domain system response data.

`np` and `nz` specify the number of poles and zeros in the estimated transfer function, respectively.

`input_delay` specifies the input delay for the estimated transfer function as 0.2 seconds.

`sys` is an `idtf` model containing the estimated transfer function with an input delay of 0.2 seconds.

### Convert Frequency-Response Data into Transfer Function

This example requires a Control System Toolbox™ license.

Obtain frequency-response data.

For example, use `bode` to obtain the magnitude and phase response data for the following system:

$$H(s) = \frac{s + 0.2}{s^3 + 2s^2 + s + 1}$$

Use 100 frequency points, ranging from 0.1 rad/s to 10 rad/s, to obtain the frequency-response data. Use `frd` to create a frequency-response data object.

```
freq = logspace(-1,1,100);  
[mag,phase] = bode(tf([1 0.2],[1 2 1 1]),freq);  
data = frd(mag.*exp(1j*phase*pi/180),freq);
```

Estimate a transfer function using `data`.

```
np = 3;  
nz = 1;  
sys = tfest(data,np,nz);
```

`np` and `nz` specify the number of poles and zeros in the estimated transfer function, respectively.

`sys` is an `idtf` model containing the estimated transfer function.

## Estimate Transfer Function with Known Transport Delays for Multiple Inputs

Load time-domain system response data.

```
load co2data;  
Ts = 0.5;  
data = iddata(Output_exp1,Input_exp1,Ts);
```

data is an iddata object and has a sample rate of 0.5 seconds.

Specify the search method as gna. Also specify the maximum search iterations and input/output offsets.

```
opt = tfestOptions('SearchMethod','gna');  
opt.InputOffset = [170;50];  
opt.OutputOffset = mean(data.y(1:75));  
opt.SearchOptions.MaxIterations = 50;
```

opt is an estimation option set that specifies the search method as gna, with a maximum of 50 iterations. opt also specifies the input offset and the output offset.

Estimate a transfer function using the measured data and the estimation option set. Specify the transport delays from the inputs to the output.

```
np = 3;  
nz = 1;  
iodelay = [2 5];  
sys = tfest(data,np,nz,iodelay,opt);
```

iodelay specifies the input to output delay from the first and second inputs to the output as 2 seconds and 5 seconds, respectively.

sys is an idtf model containing the estimated transfer function.

## Estimate Transfer Function with Known and Unknown Transport Delays

Load time-domain system response data and use it to estimate a transfer function for the system. Specify the known and unknown transport delays.

```
load co2data;  
Ts = 0.5;  
data = iddata(Output_exp1,Input_exp1,Ts);
```

`data` is an `iddata` object and has a sample rate of 0.5 seconds.

Specify the search method as `gna`. Also specify the maximum search iterations and input/output offsets.

```
opt = tfestOptions('Display','on','SearchMethod','gna');
opt.InputOffset = [170; 50];
opt.OutputOffset = mean(data.y(1:75));
opt.SearchOptions.MaxIterations = 50;
```

`opt` is an estimation option set that specifies the search method as `gna`, with a maximum of 50 iterations. `opt` also specifies the input/output offsets.

Estimate the transfer function. Specify the unknown and known transport delays.

```
np = 3;
nz = 1;
iodelay = [2 nan];
sys = tfest(data,np,nz,iodelay,opt);
```

`iodelay` specifies the transport delay from the first input to the output as 2 seconds. Using `NaN` specifies the transport delay from the second input to the output as unknown.

`sys` is an `idtf` model containing the estimated transfer function.

### Estimate Transfer Function with Unknown, Constrained Transport Delays

Create a transfer function model with the expected numerator and denominator structure and delay constraints.

In this example, the experiment data consists of two inputs and one output. Both transport delays are unknown and have an identical upper bound. Additionally, the transfer functions from both inputs to the output are identical in structure.

```
init_sys = idtf(NaN(1,2),[1,NaN(1,3)],'IODelay',NaN);
init_sys.Structure(1).IODelay.Free = true;
init_sys.Structure(1).IODelay.Maximum = 7;
```

`init_sys` is an `idtf` model describing the structure of the transfer function from one input to the output. The transfer function consists of one zero, three poles and a transport delay. The use of `NaN` indicates unknown coefficients.

`init_sys.Structure(1).IODelay.Free = true` indicates that the transport delay is not fixed.

`init_sys.Structure(1).IODelay.Maximum = 7` sets the upper bound for the transport delay to 7 seconds.

Specify the transfer function from both inputs to the output.

```
init_sys = [init_sys,init_sys];
```

Load time-domain system response data and use it to estimate a transfer function.

```
load co2data;
Ts = 0.5;
data = iddata(Output_exp1,Input_exp1,Ts);
opt = tfestOptions('Display','on','SearchMethod','gna');
opt.InputOffset = [170;50];
opt.OutputOffset = mean(data.y(1:75));
opt.SearchOptions.MaxIterations = 50;
sys = tfest(data,init_sys,opt);
```

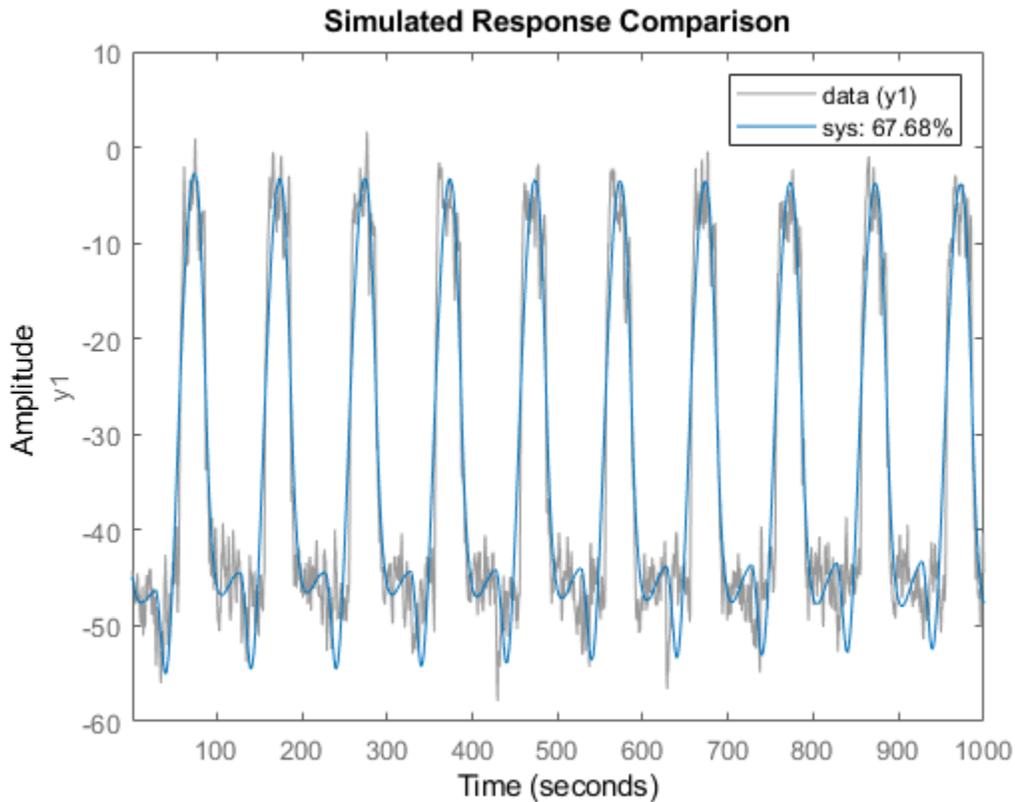
`data` is an `iddata` object and has a sample rate of 0.5 seconds.

`opt` is an estimation option set that specifies the search method as `gna`, with a maximum of 50 iterations. `opt` also specifies the input offset and the output offset.

`sys` is an `idtf` model containing the estimated transfer function.

Analyze the estimation result by comparison.

```
opt2 = compareOptions;
opt2.InputOffset = opt.InputOffset;
opt2.OutputOffset = opt.OutputOffset;
compare(data,sys,opt2)
```



### Estimate Transfer Function Containing Different Number of Poles for Input/Output Pairs

Estimate a multiple-input, single-output transfer function containing different number of poles for input/output pairs for given data.

This example requires a Control System Toolbox™ license.

Obtain frequency-response data.

For example, use `f rd` to create a frequency-response data model for the following system:

$$G = \begin{bmatrix} e^{-4s} \frac{s+2}{s^3+2s^2+4s+5} \\ e^{-0.6s} \frac{5}{s^4+2s^3+s^2+s} \end{bmatrix}$$

Use 100 frequency points, ranging from 0.01 rad/s to 100 rad/s, to obtain the frequency-response data.

```
G = tf([1 2],[5]},{[1 2 4 5],[1 2 1 1 0]},0,'IODelay',[4 0.6]);  
data = frd(G,logspace(-2,2,100));
```

`data` is an `frd` object containing the continuous-time frequency response for `G`.

Estimate a transfer function for `data`.

```
np = [3 4];  
nz = [1 0];  
iodelay = [4 0.6];  
sys = tfest(data,np,nz,iodelay);
```

`np` specifies the number of poles in the estimated transfer function. The first element of `np` indicates that the transfer function from the first input to the output contains 3 poles. Similarly, the second element of `np` indicates that the transfer function from the second input to the output contains 4 poles.

`nz` specifies the number of zeros in the estimated transfer function. The first element of `nz` indicates that the transfer function from the first input to the output contains 1 zero. Similarly, the second element of `np` indicates that the transfer function from the second input to the output does not contain any zeros.

`iodelay` specifies the transport delay from the first input to the output as 4 seconds. The transport delay from the second input to the output is specified as 0.6 seconds.

`sys` is an `idtf` model containing the estimated transfer function.

### **Estimate Transfer Function for Unstable System**

Estimate a transfer function describing an unstable system for given data.

Obtain frequency-response data.

For example, use `frd` to create frequency-response data model for the following system:

$$G = \begin{bmatrix} \frac{s+2}{s^3+2s^2+4s+5} \\ \frac{5}{s^4+2s^3+s^2+s+1} \end{bmatrix}$$

Use 100 frequency points, ranging from 0.01 rad/s to 100 rad/s, to obtain the frequency-response data.

```
G = idtf([1 2], 5), [1 2 4 5], [1 2 1 1 1]);
data = idfrd(G, logspace(-2, 2, 100));
```

`data` is an `idfrd` object containing the continuous-time frequency response for `G`.

Estimate a transfer function for `data`.

```
np = [3 4];
nz = [1 0];
sys = tfest(data, np, nz);
```

`np` specifies the number of poles in the estimated transfer function. The first element of `np` indicates that the transfer function from the first input to the output contains 3 poles. Similarly, the second element of `np` indicates that the transfer function from the second input to the output contains 4 poles.

`nz` specifies the number of zeros in the estimated transfer function. The first element of `nz` indicates that the transfer function from the first input to the output contains 1 zero. Similarly, the second element of `nz` indicates that the transfer function from the second input to the output does not contain any zeros.

`sys` is an `idtf` model containing the estimated transfer function.

```
pole(sys)
```

```
ans = 7×1 complex
```

```
-1.5260 + 0.0000i
-0.2370 + 1.7946i
-0.2370 - 1.7946i
-1.4656 + 0.0000i
-1.0000 + 0.0000i
0.2328 + 0.7926i
```

0.2328 - 0.7926i

sys is an unstable system as verified by the pole display.

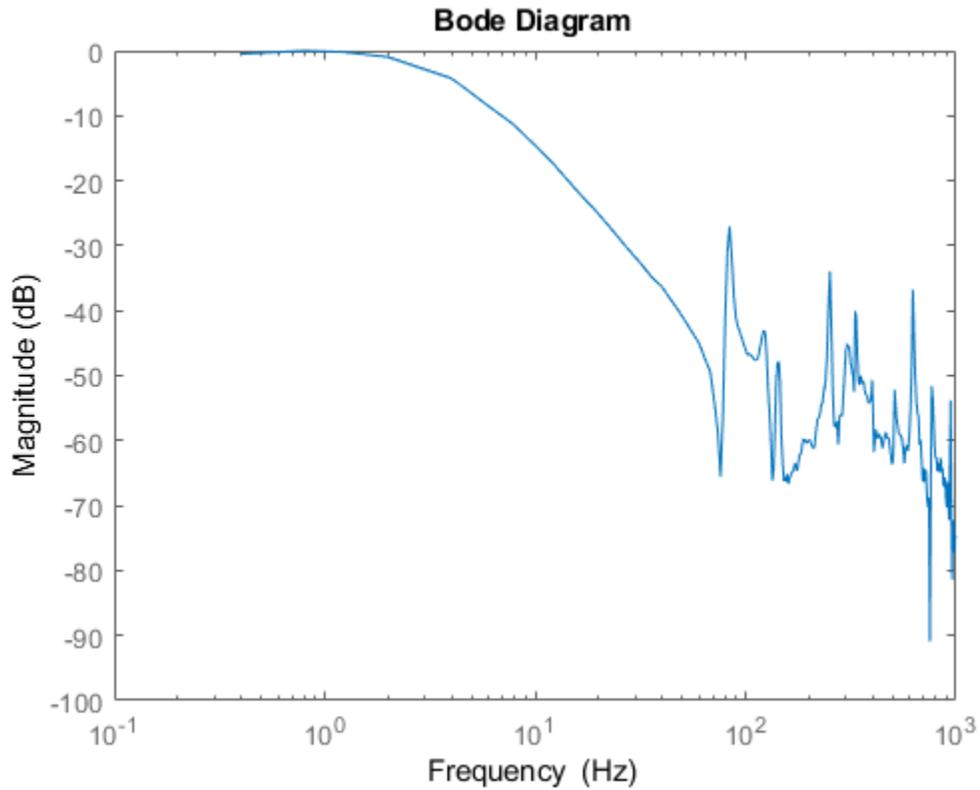
### **Estimate Transfer Function using High Modal Density Frequency Response Data**

Load the high density frequency response measurement data. The data corresponds to an unstable process maintained at equilibrium using feedback control.

```
load HighModalDensityData FRF f
```

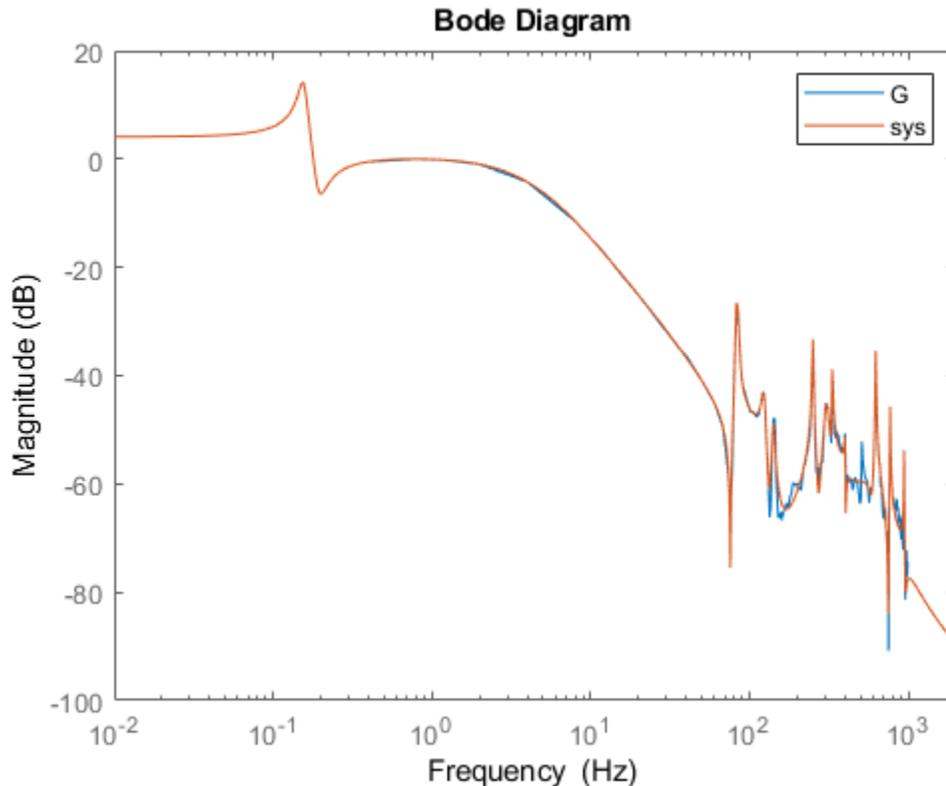
Package the data as an `idfrd` object for identification and, find the Bode magnitude response.

```
G = idfrd(permute(FRF,[2 3 1]),f,0,'FrequencyUnit','Hz');  
bodemag(G)
```



Estimate a transfer function with 32 poles and 32 zeros, and compare the Bode magnitude response.

```
sys = tfest(G,32,32);  
bodemag(G, sys)  
xlim([0.01,2e3])  
legend
```



## Definitions

## Compatibility

Starting in R2016b, a new algorithm is used for performing transfer function estimation from frequency-domain data. You are likely to see faster and more accurate results with the new algorithm, particularly for data with dynamics over a large range of frequencies and amplitudes. However, the estimation results may not match results from previous releases. To perform estimation using the previous estimation algorithm, append ' - R2016a ' to the syntax.

For example, suppose that you are estimating a transfer function model with `np` poles using frequency-domain data, `data`.

```
sys = tfest(data,np)
```

To use the previous estimation algorithm, use the following syntax.

```
sys = tfest(data,np, '-R2016a')
```

## Algorithms

The details of the estimation algorithms used by `tfest` vary depending on a variety of factors, including the sampling of the estimated model and the estimation data.

### Continuous-Time Transfer Function Estimation Using Time-Domain Data

The estimation algorithm initializes the estimable parameters using the method specified by the `InitializeMethod` estimation option. The default method is the Instrument Variable (IV) method.

The State-Variable Filters (SVF) approach and the Generalized Poisson Moment Functions (GPMF) approach to continuous-time parameter estimation use prefiltered data [1] [2].

The constant  $\frac{1}{\tau}$  in [1] and [2] corresponds to the initialization option (`InitializeOptions`) field `FilterTimeConstant`. IV is the simplified refined IV method and is called SRIVC in [3]. This method has a prefilter that is the denominator of the current model, initialized with SVF. This prefilter is iterated up to `MaxIterations` times, until the model change is less than `Tolerance`. `MaxIterations` and `Tolerance` are options that you can specify using the `InitializeOptions` structure. The 'n4sid' initialization option estimates a discrete-time model, using the N4SID estimation algorithm, that it transforms to continuous-time using `d2c`.

You use `tfestOptions` to create the option set used to estimate a transfer function.

The initialized parameters are updated using a nonlinear least-squares search method, specified by the `SearchMethod` estimation option. The objective of the search method is to minimize the weighted prediction error norm.

## Discrete-Time Transfer Function Estimation Using Time-Domain Data

For discrete-time data, `tfest` uses the same algorithm as `oe` to determine the numerator and denominator polynomial coefficients. In this algorithm, the initialization is performed using `arx`, followed by nonlinear least-squares search based updates to minimize a weighted prediction error norm.

## Continuous-Time Transfer Function Estimation Using Continuous-Time Frequency-Domain Data

The estimation algorithm performs the following tasks:

- 1 Perform a bilinear mapping to transform the domain (frequency grid) of the transfer function. For continuous-time models, the imaginary axis is transformed to the unit disk. For discrete-time models, the original domain unit disk is transformed to another unit disk.
- 2 Perform S-K iterations [4] to solve a nonlinear least-squares problem — Consider a multi-input single-output system. The nonlinear least-squares problem is to minimize the loss function:

$$\text{minimize}_{D, N_i} \sum_{k=1}^{n_f} \left| W(\omega_k) \left( y(\omega_k) - \sum_{i=1}^{n_u} \frac{N_i(\omega_k)}{D(\omega_k)} u_i(\omega_k) \right) \right|^2$$

Here  $W$  is a frequency-dependent weight that you specify.  $D$  is the denominator of the transfer function model that is to be estimated, and  $N_i$  is the numerator corresponding to the  $i$ th input.  $y$  and  $u$  are the measured output and input data, respectively.  $n_f$  and  $n_u$  are the number of frequencies and inputs, and  $w$  is the frequency. Rearranging the terms gives:

$$\text{minimize}_{D, N_i} \sum_{k=1}^{n_f} \left| \frac{W(\omega_k)}{D(\omega_k)} \left( D(\omega_k) y(\omega_k) - \sum_{i=1}^{n_u} N_i(\omega_k) u_i(\omega_k) \right) \right|^2$$

To perform the S-K iterations, the algorithm iteratively solves

$$\underset{D_m, N_{i,m}}{\text{minimize}} \sum_{k=1}^{n_f} \left| \frac{W(\omega_k)}{D_{m-1}(\omega_k)} \left( D_m(\omega_k) y(\omega_k) - \sum_{i=1}^{n_u} N_{i,m}(\omega_k) u_i(\omega_k) \right) \right|^2$$

where  $m$  is the current iteration, and  $D_{m-1}(\omega)$  is the denominator response identified at the previous iteration. Now each step of the iteration is a linear least-squares problem, where the identified parameters capture the responses  $D_m(\omega)$  and  $N_{i,m}(\omega)$  for  $i = 1, 2, \dots, n_u$ . The iteration is initialized by choosing  $D_0(\omega) = 1$ .

- The first iteration of the algorithm identifies  $D_1(\omega)$ . The  $D_1(\omega)$  and  $N_{i,1}(\omega)$  polynomials are expressed in monomial basis.
- The second and following iterations express the polynomials  $D_m(\omega)$  and  $N_{i,m}(\omega)$  in terms of orthogonal rational basis functions on the unit disk. These basis functions have the form:

$$B_{j,m}(\omega) = \frac{\left( \frac{\sqrt{1 - |\lambda_{j,m-1}|^2}}{q - \lambda_{j,m-1}} \right) \prod_{r=0}^{j-1} \frac{1 - (\lambda_{j,m-1})^* q(\omega)}{q(\omega) - \lambda_{r,m-1}}}{1}$$

Here  $\lambda_{j,m-1}$  is the  $j$ th pole that is identified at the previous step,  $m-1$ , of the iteration.  $\lambda_{j,m-1}^*$  is the complex conjugate of  $\lambda_{j,m-1}$ , and  $q$  is the frequency-domain variable on the unit disk.

- The algorithm runs for a maximum of 20 iterations. The iterations are terminated early if the relative change in the value of the loss function is less than 0.001 in the last three iterations.

If you specify bounds on transfer function coefficients, these bounds correspond to affine constraints on the identified parameters. If you only have equality constraints (fixed transfer function coefficients), the corresponding equality constrained least-squares problem is solved algebraically. To do so, the software computes an orthogonal basis for the null space of the equality constraint matrix, and then solves least-squares problem within this null space. If you have upper or lower bounds on transfer function coefficients, the corresponding inequality constrained least-squares problem is solved using interior-point methods.

- 3** Perform linear refinements — The S-K iterations, even when they converge, do not always yield a locally optimal solution. To find a critical point of the optimization problem that may yield a locally optimal solution, a second set of iterations are

performed. The critical points are solutions to a set of nonlinear equations. The algorithm searches for a critical point by successively constructing a linear approximation to the nonlinear equations and solving the resulting linear equations in the least-squares sense. The equations are:

- Equation for the  $j$ th denominator parameter:

$$0 = 2 \sum_{k=1}^{n_f} \operatorname{Re} \left\{ \frac{|W(\omega_k)|^2 B_j^*(\omega_k) \sum_{i=1}^{n_u} N_{i,m-1}^*(\omega_k) u_i^*(\omega_k)}{D_{m-1}^*(\omega_k) |D_{m-1}(\omega_k)|^2} \left( D_m(\omega_k) y(\omega_k) - \sum_{i=1}^{n_u} N_{i,m}(\omega_k) u_i(\omega_k) \right) \right\}$$

- Equation for the  $j$ th numerator parameter that corresponds to input  $l$ :

$$0 = -2 \sum_{k=1}^{n_f} \operatorname{Re} \left\{ \frac{|W(\omega_k)|^2 B_j^*(\omega_k) u_l^*(\omega_k)}{|D_{m-1}(\omega_k)|^2} \left( D_m(\omega_k) y(\omega_k) - \sum_{i=1}^{n_u} N_{i,m}(\omega_k) u_i(\omega_k) \right) \right\}$$

The first iteration is started with the best solution found for the numerators  $N_i$  and denominator  $D$  parameters during S-K iterations. Unlike S-K iterations, the basis functions  $B_j(\omega)$  are not changed at each iteration, the iterations are performed with the basis functions that yielded the best solution in the S-K iterations. As before, the algorithm runs for a maximum of 20 iterations. The iterations are terminated early if the relative change in the value of the loss function is less than 0.001 in the last three iterations.

If you specify bounds on transfer function coefficients, these bounds are incorporated into the necessary optimality conditions via generalized Lagrange multipliers. The resulting constrained linear least-squares problems are solved using the same methods explained in the S-K iterations step.

- 4 Return the transfer function parameters corresponding to the optimal solution — Both the S-K and linear refinement iteration steps do not guarantee an improvement in the loss function value. The algorithm tracks the best parameter value observed during these steps, and returns these values.
- 5 Invert the bilinear mapping performed in step 1.
- 6 Perform an iterative refinement of the transfer function parameters using the nonlinear least-squares search method specified in the `SearchMethod` estimation option. This step is implemented in the following situations:

- When you specify the `EnforceStability` estimation option as `true` (stability is requested), and the result of step 5 of this algorithm is an unstable model. The unstable poles are reflected inside the stability boundary and the resulting parameters are iteratively refined. For information about estimation options, see `tfestOptions`.
- When you add a regularization penalty to the loss function using the `Regularization` estimation option. For an example about regularization, see “Regularized Identification of Dynamic Systems”.
- You estimate a continuous-time model using discrete-time data (see “Discrete-Time Transfer Function Estimation Using Discrete-Time Frequency-Domain Data” on page 1-1799).
- You use frequency domain input-output data to identify a multi-input model.

If you are using the estimation algorithm from R2016a or earlier (see “Compatibility” on page 1-1794) for estimating a continuous-time model using continuous-time frequency-domain data, then for continuous-time data and fixed delays, the Output-Error algorithm is used for model estimation. For continuous-time data and free delays, the state-space estimation algorithm is used. In this algorithm, the model coefficients are initialized using the N4SID estimation method. This initialization is followed by nonlinear least-squares search based updates to minimize a weighted prediction error norm.

## Discrete-Time Transfer Function Estimation Using Discrete-Time Frequency-Domain Data

The estimation algorithm is the same as for continuous-time transfer function estimation using continuous-time frequency-domain data on page 1-1796, except discrete-time data is used.

If you are using the estimation algorithm from R2016a or earlier (see “Compatibility” on page 1-1794), the algorithm is the same as the algorithm for discrete-time transfer function estimation using time-domain data on page 1-1796.

---

**Note** The software does not support estimation of a discrete-time transfer function using continuous-time frequency-domain data.

---

## Continuous-Time Transfer Function Estimation Using Discrete-Time Frequency-Domain Data

`tfest` command first estimates a discrete-time model from the discrete-time data. The estimated model is then converted to a continuous-time model using the `d2c` command. The frequency response of the resulting continuous-time model is then computed over the frequency grid of the estimation data. A continuous-time model of the desired (user-specified) structure is then fit to this frequency response. The estimation algorithm for using the frequency-response data to obtain the continuous-time model is the same as that for continuous-time transfer function estimation using continuous-time data on page 1-1796.

If you are using the estimation algorithm from R2016a or earlier (see “Compatibility” on page 1-1794), the state-space estimation algorithm is used for estimating continuous-time models from discrete-time data. In this algorithm, the model coefficients are initialized using the N4SID estimation method. This initialization is followed by nonlinear least-squares search based updates to minimize a weighted prediction error norm.

### Delay Estimation

- When delay values are specified as `NaN`, they are estimated separate from the model numerator and denominator coefficients, using `delayest`. The delay values thus determined are treated as fixed values during the iterative update of the model using a nonlinear least-squares search method. Thus, the delay values are not iteratively updated.
- For an initial model, `init_sys`, with:
  - `init_sys.Structure.IODelay.Value` specified as finite values
  - `init_sys.Structure.IODelay.Free` specified as `true`

the initial delay values are left unchanged.

Estimation of delays is often a difficult problem. You should assess the presence and the value of a delay. To do so, use physical insight of the process being modeled and functions such as `arxstruc`, `delayest`, and `impulseest`. For an example of determining input delay, see “Model Structure Selection: Determining Model Order and Input Delay”.

## References

- [1] Garnier, H., M. Mensler, and A. Richard. "Continuous-time Model Identification From Sampled Data: Implementation Issues and Performance Evaluation." *International Journal of Control*, 2003, Vol. 76, Issue 13, pp 1337-1357.
- [2] Ljung, L. "Experiments With Identification of Continuous-Time Models." *Proceedings of the 15th IFAC Symposium on System Identification*. 2009.
- [3] Young, P. C. and A.J. Jakeman. "Refined instrumental variable methods of time-series analysis: Part III, extensions." *International Journal of Control* 31, 1980, pp 741-764.
- [4] Drmac, Z., S. Gugercin, and C. Beattie. "Quadrature-based vector fitting for discretized  $H_2$  approximation." *SIAM Journal on Scientific Computing*. Vol. 37, Numer 2, 2014, pp A625-A652.
- [5] Ozdemir, A. A., and S. Gumussoy. "Transfer Function Estimation in System Identification Toolbox via Vector Fitting." *Proceedings of the 20th World Congress of the International Federation of Automatic Control*. Toulouse, France, July 2017.

## Extended Capabilities

### Automatic Parallel Support

Accelerate code by automatically running computation in parallel using Parallel Computing Toolbox™.

Parallel computing support is available for estimation using the `lsqnonlin` search method (requires Optimization Toolbox). To enable parallel computing, use `tfestOptions`, set `SearchMethod` to `'lsqnonlin'`, and set `SearchOptions.Advanced.UseParallel` to `true`.

For example:

```
opt = tfestOptions;  
opt.SearchMethod = 'lsqnonlin';  
opt.SearchOptions.Advanced.UseParallel = true;
```

## See Also

ar | arx | bj | greyest | idtf | oe | polyest | procest | ssest | tfestOptions

## Topics

“Estimate Transfer Function Models at the Command Line”

“Estimate Transfer Function Models with Transport Delay to Fit Given Frequency-Response Data”

“Estimate Transfer Function Models With Prior Knowledge of Model Structure and Constraints”

“Troubleshoot Frequency-Domain Identification of Transfer Function Models”

“What are Transfer Function Models?”

“Regularized Estimates of Model Parameters”

“Estimating Models Using Frequency-Domain Data”

**Introduced in R2012a**

# tfestOptions

Option set for tfest

## Syntax

```
opt = tfestOptions  
opt = tfestOptions(Name,Value)
```

## Description

`opt = tfestOptions` creates the default option set for `tfest`.

`opt = tfestOptions(Name,Value)` creates an option set with the options specified by one or more `Name,Value` pair arguments.

## Input Arguments

### Name-Value Pair Arguments

Specify optional comma-separated pairs of `Name,Value` arguments. `Name` is the argument name and `Value` is the corresponding value. `Name` must appear inside quotes. You can specify several name and value pair arguments in any order as `Name1,Value1,...,NameN,ValueN`.

### InitializeMethod — Algorithm used to initialize the numerator and denominator

```
'iv' (default) | 'svf' | 'gpmf' | 'n4sid' | 'all'
```

Algorithm used to initialize the values of the numerator and denominator of the output of `tfest`, applicable only for estimation of continuous-time transfer functions using time-domain data, specified as one of the following values:

- `'iv'` — Instrument Variable approach.

- 'svf' — State Variable Filters approach.
- 'gpmf' — Generalized Poisson Moment Functions approach.
- 'n4sid' — Subspace state-space estimation approach.
- 'all' — Combination of all of the preceding approaches. The software tries all these methods and selects the method that yields the smallest value of prediction error norm.

## **InitializeOptions — Option set for the initialization algorithm**

structure

Option set for the initialization algorithm used to initialize the values of the numerator and denominator of the output of `tfest`, specified as a structure with the following fields:

- **N4Weight** — Calculates the weighting matrices used in the singular-value decomposition step of the 'n4sid' algorithm. Applicable when `InitializeMethod` is 'n4sid'.

**N4Weight** is specified as one of the following values:

- 'MOESP' — Uses the MOESP algorithm by Verhaegen.
- 'CVA' — Uses the canonical variable algorithm (CVA) by Larimore.
- 'SSARX' — A subspace identification method that uses an ARX estimation based algorithm to compute the weighting.

Specifying this option allows unbiased estimates when using data that is collected in closed-loop operation. For more information about the algorithm, see [6].

- 'auto' — The software automatically determines if the MOESP algorithm or the CVA algorithm should be used in the singular-value decomposition step.

**Default:** 'auto'

- **N4Horizon** — Determines the forward and backward prediction horizons used by the 'n4sid' algorithm. Applicable when `InitializeMethod` is 'n4sid'.

**N4Horizon** is a row vector with three elements:  $[r \ sy \ su]$ , where  $r$  is the maximum forward prediction horizon. The algorithm uses up to  $r$  step-ahead predictors.  $sy$  is the number of past outputs, and  $su$  is the number of past inputs that are used for the predictions. See pages 209 and 210 in [1] for more information. These numbers can have a substantial influence on the quality of the resulting model, and there are no simple rules for choosing them. Making 'N4Horizon' a  $k$ -by-3 matrix

means that each row of 'N4Horizon' is tried, and the value that gives the best (prediction) fit to data is selected. k is the number of guesses of [r sy su] combinations.

If N4Horizon = 'auto', the software uses an Akaike Information Criterion (AIC) for the selection of sy and su.

**Default:** 'auto'

- **FilterTimeConstant** — Time constant of the differentiating filter used by the iv, svf, and gpmf initialization methods (see [4] and [5]).

FilterTimeConstant specifies the cutoff frequency of the differentiating filter,  $F_{cutoff}$ , as:

$$F_{cutoff} = \frac{\text{FilterTimeConstant}}{T_s}$$

$T_s$  is the sample time of the estimation data.

Specify FilterTimeConstant as a positive number, typically less than 1. A good value of FilterTimeConstant is the ratio of  $T_s$  to the dominating time constant of the system.

**Default:** 0.1

- **MaxIterations** — Maximum number of iterations. Applicable when InitializeMethod is 'iv'.

**Default:** 30

- **Tolerance** — Convergence tolerance. Applicable when InitializeMethod is 'iv'.

**Default:** 0.01

### **InitialCondition — Handling of initial conditions**

'auto' (default) | 'zero' | 'estimate' | 'backcast'

Handling of initial conditions during estimation, specified as one of the following values:

- 'zero' — All initial conditions are taken as zero.
- 'estimate' — The necessary initial conditions are treated as estimation parameters.

- 'backcast' — The necessary initial conditions are estimated by a backcasting (backward filtering) process, described in [2].
- 'auto' — An automatic choice among the preceding options is made, guided by the data.

## **WeightingFilter — Weighting prefilter**

[ ] (default) | vector | matrix | cell array | linear system | 'inv' | 'invsqrt'

Weighting prefilter applied to the loss function to be minimized during estimation. To understand the effect of `WeightingFilter` on the loss function, see “Loss Function and Model Quality Metrics”.

Specify `WeightingFilter` as one of the following values:

- [ ] — No weighting prefilter is used.
- Passbands — Specify a row vector or matrix containing frequency values that define desired passbands. You select a frequency band where the fit between estimated model and estimation data is optimized. For example, `[wl, wh]` where `wl` and `wh` represent lower and upper limits of a passband. For a matrix with several rows defining frequency passbands, `[w1l, w1h; w2l, w2h; w3l, w3h; . . . ]`, the estimation algorithm uses the union of the frequency ranges to define the estimation passband.

Passbands are expressed in `rad/TimeUnit` for time-domain data and in `FrequencyUnit` for frequency-domain data, where `TimeUnit` and `FrequencyUnit` are the time and frequency units of the estimation data.

- SISO filter — Specify a single-input-single-output (SISO) linear filter in one of the following ways:
  - A SISO LTI model
  - `{A, B, C, D}` format, which specifies the state-space matrices of a filter with the same sample time as estimation data.
  - `{numerator, denominator}` format, which specifies the numerator and denominator of the filter as a transfer function with the same sample time as estimation data.

This option calculates the weighting function as a product of the filter and the input spectrum to estimate the transfer function.

- Weighting vector — Applicable for frequency-domain data only. Specify a column vector of weights. This vector must have the same length as the frequency vector of

the data set, `Data.Frequency`. Each input and output response in the data is multiplied by the corresponding weight at that frequency.

- `'inv'` — Applicable for estimation using frequency-response data only. Use  $1/|G(\omega)|$  as the weighting filter, where  $G(\omega)$  is the complex frequency-response data. Use this option for capturing relatively low amplitude dynamics in data, or for fitting data with high modal density. This option also makes it easier to specify channel-dependent weighting filters for MIMO frequency-response data.
- `'invsqrt'` — Applicable for estimation using frequency-response data only. Use  $1/\sqrt{|G(\omega)|}$  as the weighting filter. Use this option for capturing relatively low amplitude dynamics in data, or for fitting data with high modal density. This option also makes it easier to specify channel-dependent weighting filters for MIMO frequency-response data.

### **EnforceStability — Control whether to enforce stability of model**

`false` (default) | `true`

Control whether to enforce stability of estimated model, specified as the comma-separated pair consisting of `'EnforceStability'` and either `true` or `false`.

Use this option when estimating models using frequency-domain data. Models estimated using time-domain data are always stable.

Data Types: `logical`

### **EstimateCovariance — Control whether to generate parameter covariance data**

`true` (default) | `false`

Controls whether parameter covariance data is generated, specified as `true` or `false`.

If `EstimateCovariance` is `true`, then use `getcov` to fetch the covariance matrix from the estimated model.

### **Display — Specify whether to display the estimation progress**

`'off'` (default) | `'on'`

Specify whether to display the estimation progress, specified as one of the following values:

- `'on'` — Information on model structure and estimation results are displayed in a progress-viewer window.

- 'off' — No progress or results information is displayed.

## **InputOffset — Removal of offset from time-domain input data during estimation**

[ ] (default) | vector of positive integers | matrix

Removal of offset from time-domain input data during estimation, specified as the comma-separated pair consisting of 'InputOffset' and one of the following:

- A column vector of positive integers of length  $N_u$ , where  $N_u$  is the number of inputs.
- [ ] — Indicates no offset.
- $N_u$ -by- $N_e$  matrix — For multi-experiment data, specify InputOffset as an  $N_u$ -by- $N_e$  matrix.  $N_u$  is the number of inputs, and  $N_e$  is the number of experiments.

Each entry specified by InputOffset is subtracted from the corresponding input data.

## **OutputOffset — Removal of offset from time-domain output data during estimation**

[ ] (default) | vector | matrix

Removal of offset from time-domain output data during estimation, specified as the comma-separated pair consisting of 'OutputOffset' and one of the following:

- A column vector of length  $N_y$ , where  $N_y$  is the number of outputs.
- [ ] — Indicates no offset.
- $N_y$ -by- $N_e$  matrix — For multi-experiment data, specify OutputOffset as a  $N_y$ -by- $N_e$  matrix.  $N_y$  is the number of outputs, and  $N_e$  is the number of experiments.

Each entry specified by OutputOffset is subtracted from the corresponding output data.

## **OutputWeight — Weighting of prediction errors in multi-output estimations**

[ ] (default) | 'noise' | positive semidefinite symmetric matrix

Weighting of prediction errors in multi-output estimations, specified as one of the following values:

- 'noise' — Minimize  $\det(E^*E / N)$ , where  $E$  represents the prediction error and  $N$  is the number of data samples. This choice is optimal in a statistical sense and leads to maximum likelihood estimates if nothing is known about the variance of the noise. It uses the inverse of the estimated noise variance as the weighting function.

---

**Note** OutputWeight must not be 'noise' if SearchMethod is 'lsqnonlin'.

---

- Positive semidefinite symmetric matrix ( $W$ ) — Minimize the trace of the weighted prediction error matrix  $\text{trace}(E' * E * W / N)$  where:
  - $E$  is the matrix of prediction errors, with one column for each output, and  $W$  is the positive semidefinite symmetric matrix of size equal to the number of outputs. Use  $W$  to specify the relative importance of outputs in multiple-output models, or the reliability of corresponding data.
  - $N$  is the number of data samples.
- `[]` — The software chooses between the 'noise' or using the identity matrix for  $W$ .

This option is relevant for only multi-output models.

### Regularization — Options for regularized estimation of model parameters structure

Options for regularized estimation of model parameters. For more information on regularization, see “Regularized Estimates of Model Parameters”.

Regularization is a structure with the following fields:

- **Lambda** — Constant that determines the bias versus variance tradeoff.

Specify a positive scalar to add the regularization term to the estimation cost.

The default value of zero implies no regularization.

**Default:** 0

- **R** — Weighting matrix.

Specify a vector of nonnegative numbers or a square positive semi-definite matrix. The length must be equal to the number of free parameters of the model.

For black-box models, using the default value is recommended. For structured and grey-box models, you can also specify a vector of `np` positive numbers such that each entry denotes the confidence in the value of the associated parameter.

The default value of 1 implies a value of `eye(npfree)`, where `npfree` is the number of free parameters.

**Default:** 1

- **Nominal** — The nominal value towards which the free parameters are pulled during estimation.

The default value of zero implies that the parameter values are pulled towards zero. If you are refining a model, you can set the value to `'model'` to pull the parameters towards the parameter values of the initial model. The initial parameter values must be finite for this setting to work.

**Default:** 0

### **SearchMethod** — Numerical search method used for iterative parameter estimation

`'auto'` (default) | `'gn'` | `'gna'` | `'lm'` | `'grad'` | `'lsqnonlin'` | `'fmincon'`

Numerical search method used for iterative parameter estimation, specified as the comma-separated pair consisting of `'SearchMethod'` and one of the following:

- `'auto'` — A combination of the line search algorithms, `'gn'`, `'lm'`, `'gna'`, and `'grad'` methods is tried in sequence at each iteration. The first descent direction leading to a reduction in estimation cost is used.
- `'gn'` — Subspace Gauss-Newton least squares search. Singular values of the Jacobian matrix less than  $GnPinvConstant * eps * \max(\text{size}(J)) * \text{norm}(J)$  are discarded when computing the search direction.  $J$  is the Jacobian matrix. The Hessian matrix is approximated as  $J^T J$ . If there is no improvement in this direction, the function tries the gradient direction.
- `'gna'` — Adaptive subspace Gauss-Newton search. Eigenvalues less than  $\gamma * \max(sv)$  of the Hessian are ignored, where  $sv$  contains the singular values of the Hessian. The Gauss-Newton direction is computed in the remaining subspace.  $\gamma$  has the initial value `InitialGnaTolerance` (see `Advanced in 'SearchOptions'` for more information). This value is increased by the factor `LMStep` each time the search fails to find a lower value of the criterion in fewer than five bisections. This value is decreased by the factor  $2 * LMStep$  each time a search is successful without any bisections.
- `'lm'` — Levenberg-Marquardt least squares search, where the next parameter value is  $-\text{pinv}(H+d*I) * \text{grad}$  from the previous one.  $H$  is the Hessian,  $I$  is the identity matrix, and  $\text{grad}$  is the gradient.  $d$  is a number that is increased until a lower value of the criterion is found.
- `'grad'` — Steepest descent least squares search.
- `'lsqnonlin'` — Trust-region-reflective algorithm of `lsqnonlin`. Requires Optimization Toolbox software.

- 'fmincon' — Constrained nonlinear solvers. You can use the sequential quadratic programming (SQP) and trust-region-reflective algorithms of the fmincon solver. If you have Optimization Toolbox software, you can also use the interior-point and active-set algorithms of the fmincon solver. Specify the algorithm in the SearchOptions.Algorithm option. The fmincon algorithms may result in improved estimation results in the following scenarios:
  - Constrained minimization problems when there are bounds imposed on the model parameters.
  - Model structures where the loss function is a nonlinear or non smooth function of the parameters.
  - Multi-output model estimation. A determinant loss function is minimized by default for multi-output model estimation. fmincon algorithms are able to minimize such loss functions directly. The other search methods such as 'lm' and 'gn' minimize the determinant loss function by alternately estimating the noise variance and reducing the loss value for a given noise variance value. Hence, the fmincon algorithms can offer better efficiency and accuracy for multi-output model estimations.

### **SearchOptions — Option set for the search algorithm**

search option set

Option set for the search algorithm, specified as the comma-separated pair consisting of 'SearchOptions' and a search option set with fields that depend on the value of SearchMethod.

**SearchOptions Structure When SearchMethod is Specified as 'gn', 'gna', 'lm', 'grad', or 'auto'**

Field Name	Description	Default
Tolerance	Minimum percentage difference between the current value of the loss function and its expected improvement after the next iteration, specified as a positive scalar. When the percentage of expected improvement is less than <code>Tolerance</code> , the iterations stop. The estimate of the expected loss-function improvement at the next iteration is based on the Gauss-Newton vector computed for the current parameter value.	0.01
MaxIterations	<p>Maximum number of iterations during loss-function minimization, specified as a positive integer. The iterations stop when <code>MaxIterations</code> is reached or another stopping criterion is satisfied, such as <code>Tolerance</code>.</p> <p>Setting <code>MaxIterations = 0</code> returns the result of the start-up procedure.</p> <p>Use <code>sys.Report.Termination.Iterations</code> to get the actual number of iterations during an estimation, where <code>sys</code> is an <code>idtf</code> model.</p>	20

Field Name	Description	Default
Advanced	Advanced search settings, specified as a structure with the following fields:	
Field Name	Description	Default
GnPinvConstant	Jacobian matrix singular value threshold, specified as a positive scalar. Singular values of the Jacobian matrix that are smaller than $\text{GnPinvConstant} \times \max(\text{size}(J) \times \text{norm}(J) \times \text{eps})$ are discarded when computing the search direction. Applicable when SearchMethod is 'gn'.	10000
InitialGnaTolerance	Initial value of $\gamma$ , specified as a positive scalar. Applicable when SearchMethod is 'gna'.	0.0001
LMStartValue	Starting value of search-direction length $d$ in the Levenberg-Marquardt method, specified as a positive scalar. Applicable when SearchMethod is 'lm'.	0.001
LMStep	Size of the Levenberg-Marquardt step, specified as a positive integer. The next value of the search-direction length $d$ in the Levenberg-Marquardt method is LMStep times the previous one. Applicable when SearchMethod is 'lm'.	2
MaxBisections	Maximum number of bisections used for line search along the search direction, specified as a positive integer.	25
MaxFunctionEvaluations	Maximum number of calls to the model file, specified as a positive integer. Iterations stop if the number of calls to the model file exceeds this value.	Inf
MinParameterChange	Smallest parameter update allowed per iteration, specified as a nonnegative scalar.	0
RelativeImprovement	Relative improvement threshold, specified as a nonnegative scalar. Iterations stop if the relative	0

<b>Field Name</b>	<b>Description</b>		<b>Default</b>
	<b>Field Name</b>	<b>Description</b>	<b>Default</b>
		improvement of the criterion function is less than this value.	
	StepReduction	<p>Step reduction factor, specified as a positive scalar that is greater than 1. The suggested parameter update is reduced by the factor <code>StepReduction</code> after each try. This reduction continues until <code>MaxBisections</code> tries are completed or a lower value of the criterion function is obtained.</p> <p><code>StepReduction</code> is not applicable for SearchMethod 'lm' (Levenberg-Marquardt method).</p>	2

**SearchOptions Structure When SearchMethod is Specified as 'lsqnonlin'**

<b>Field Name</b>	<b>Description</b>	<b>Default</b>
FunctionTolerance	Termination tolerance on the loss function that the software minimizes to determine the estimated parameter values, specified as a positive scalar.  The value of FunctionTolerance is the same as that of <code>opt.SearchOptions.Advanced.TolFun</code> .	1e-5
StepTolerance	Termination tolerance on the estimated parameter values, specified as a positive scalar.  The value of StepTolerance is the same as that of <code>opt.SearchOptions.Advanced.TolX</code> .	1e-6
MaxIterations	Maximum number of iterations during loss-function minimization, specified as a positive integer. The iterations stop when MaxIterations is reached or another stopping criterion is satisfied, such as FunctionTolerance.  The value of MaxIterations is the same as that of <code>opt.SearchOptions.Advanced.MaxIter</code> .	20
Advanced	Advanced search settings, specified as an option set for <code>lsqnonlin</code> .  For more information, see the Optimization Options table in "Optimization Options" (Optimization Toolbox).	Use <code>optimset('lsqnonlin')</code> to create a default option set.

**SearchOptions Structure When SearchMethod is Specified as 'fmincon'**

Field Name	Description	Default
Algorithm	<p>fmincon optimization algorithm, specified as one of the following:</p> <ul style="list-style-type: none"> <li>• 'sqp' — Sequential quadratic programming algorithm. The algorithm satisfies bounds at all iterations, and it can recover from NaN or Inf results. It is not a large-scale algorithm. For more information, see “Large-Scale vs. Medium-Scale Algorithms” (Optimization Toolbox).</li> <li>• 'trust-region-reflective' — Subspace trust-region method based on the interior-reflective Newton method. It is a large-scale algorithm.</li> <li>• 'interior-point' — Large-scale algorithm that requires Optimization Toolbox software. The algorithm satisfies bounds at all iterations, and it can recover from NaN or Inf results.</li> <li>• 'active-set' — Requires Optimization Toolbox software. The</li> </ul>	'sqp'

Field Name	Description	Default
	<p>algorithm can take large steps, which adds speed. It is not a large-scale algorithm.</p> <p>For more information about the algorithms, see “Constrained Nonlinear Optimization Algorithms” (Optimization Toolbox) and “Choosing the Algorithm” (Optimization Toolbox).</p>	
FunctionTolerance	Termination tolerance on the loss function that the software minimizes to determine the estimated parameter values, specified as a positive scalar.	1e-6
StepTolerance	Termination tolerance on the estimated parameter values, specified as a positive scalar.	1e-6
MaxIterations	Maximum number of iterations during loss function minimization, specified as a positive integer. The iterations stop when MaxIterations is reached or another stopping criterion is satisfied, such as FunctionTolerance.	100

### Advanced – Additional advanced options structure

Additional advanced options, specified as a structure with the following fields:

- `ErrorThreshold` — Specifies when to adjust the weight of large errors from quadratic to linear.

Errors larger than `ErrorThreshold` times the estimated standard deviation have a linear weight in the loss function. The standard deviation is estimated robustly as the median of the absolute deviations from the median of the prediction errors, divided by 0.7. For more information on robust norm choices, see section 15.2 of [1].

`ErrorThreshold = 0` disables robustification and leads to a purely quadratic loss function. When estimating with frequency-domain data, the software sets `ErrorThreshold` to zero. For time-domain data that contains outliers, try setting `ErrorThreshold` to 1.6.

**Default:** 0

- `MaxSize` — Specifies the maximum number of elements in a segment when input-output data is split into segments.

`MaxSize` must be a positive, integer value.

**Default:** 250000

- `StabilityThreshold` — Specifies thresholds for stability tests.

`StabilityThreshold` is a structure with the following fields:

- `s` — Specifies the location of the right-most pole to test the stability of continuous-time models. A model is considered stable when its right-most pole is to the left of `s`.

**Default:** 0

- `z` — Specifies the maximum distance of all poles from the origin to test stability of discrete-time models. A model is considered stable if all poles are within the distance `z` from the origin.

**Default:** `1+sqrt(eps)`

- `AutoInitThreshold` — Specifies when to automatically estimate the initial conditions.

The initial condition is estimated when

$$\frac{\|y_{p,z} - y_{meas}\|}{\|y_{p,e} - y_{meas}\|} > \text{AutoInitThreshold}$$

- $y_{meas}$  is the measured output.
- $y_{p,z}$  is the predicted output of a model estimated using zero initial states.
- $y_{p,e}$  is the predicted output of a model estimated using estimated initial states.

Applicable when `InitialCondition` is 'auto'.

**Default:** 1.05

## Output Arguments

### **opt** — Option set for `tfest`

`tfestOptions` option set

Option set for `tfest`, returned as an `tfestOptions` option set.

## Examples

### **Create Default Options Set for Transfer Function Estimation**

```
opt = tfestOptions;
```

### **Specify Options for Transfer Function Estimation**

Create an options set for `tfest` using the 'n4sid' initialization algorithm and set the `Display` to 'on'.

```
opt = tfestOptions('InitializeMethod','n4sid','Display','on');
```

Alternatively, use dot notation to set the values of `opt`.

```
opt = tfestOptions;  
opt.InitializeMethod = 'n4sid';  
opt.Display = 'on';
```

## References

- [1] Ljung, L. *System Identification: Theory for the User*. Upper Saddle River, NJ: Prentice-Hall PTR, 1999.
- [2] Knudsen, T. "New method for estimating ARMAX models," *In Proceedings of 10th IFAC Symposium on System Identification, SYSID'94*, Copenhagen, Denmark, July 1994, Vol. 2, pp. 611-617.
- [3] Wills, Adrian, B. Ninness, and S. Gibson. "On Gradient-Based Search for Multivariable System Estimates." *Proceedings of the 16th IFAC World Congress, Prague, Czech Republic, July 3-8, 2005*. Oxford, UK: Elsevier Ltd., 2005.
- [4] Garnier, H., M. Mensler, and A. Richard. "Continuous-time Model Identification From Sampled Data: Implementation Issues and Performance Evaluation" *International Journal of Control*, 2003, Vol. 76, Issue 13, pp 1337-1357.
- [5] Ljung, L. "Experiments With Identification of Continuous-Time Models." *Proceedings of the 15th IFAC Symposium on System Identification*. 2009.
- [6] Jansson, M. "Subspace identification and ARX modeling." *13th IFAC Symposium on System Identification*, Rotterdam, The Netherlands, 2003.

## See Also

tfest

## Topics

"Loss Function and Model Quality Metrics"

**Introduced in R2012b**

# timeoptions

Create list of time plot options

## Syntax

```
P = timeoptions
P = timeoptions('cstprefs')
```

## Description

`P = timeoptions` returns a list of available options for time plots with default values set. You can use these options to customize the time value plot appearance from the command line.

`P = timeoptions('cstprefs')` initializes the plot options you selected in the Control System and System Identification Toolbox Preferences Editor. For more information about the editor, see “Toolbox Preferences Editor”.

This table summarizes the available time plot options.

Option	Description
Title, XLabel, YLabel	Label text and style
TickLabel	Tick label style
Grid	Show or hide the grid, specified as one of the following values: 'off'   'on' <b>Default:</b> 'off'
GridColor	Color of the grid lines, specified as one of the following: Vector of RGB values in the range [0,1]  character vector of color name   'none'. For example, for yellow color, specify as one of the following: [1 1 0], 'yellow', or 'y'. <b>Default:</b> [0.15,0.15,0.15]

<b>Option</b>	<b>Description</b>
XlimMode, YlimMode	Limit modes
Xlim, Ylim	Axes limits
IOGrouping	Grouping of input-output pairs, specified as one of the following values: 'none'   'inputs'   'outputs'   'all' <b>Default:</b> 'none'
InputLabelLabels, OutputLabels	Input and output label styles
InputVisible, OutputVisible	Visibility of input and output channels
Normalize	Normalize responses, specified as one of the following values: 'on'   'off' <b>Default:</b> 'off'
SettleTimeThreshold	Settling time threshold
RiseTimeLimits	Rise time limits

Option	Description
TimeUnits	<p>Time units, specified as one of the following values:</p> <ul style="list-style-type: none"><li>• 'nanoseconds'</li><li>• 'microseconds'</li><li>• 'milliseconds'</li><li>• 'seconds'</li><li>• 'minutes'</li><li>• 'hours'</li><li>• 'days'</li><li>• 'weeks'</li><li>• 'months'</li><li>• 'years'</li></ul> <p><b>Default:</b> 'seconds'</p> <p>You can also specify 'auto' which uses time units specified in the TimeUnit property of the input system. For multiple systems with different time units, the units of the first system is used.</p>

## Examples

### Plot Normalized Step Response

Create a default time options set.

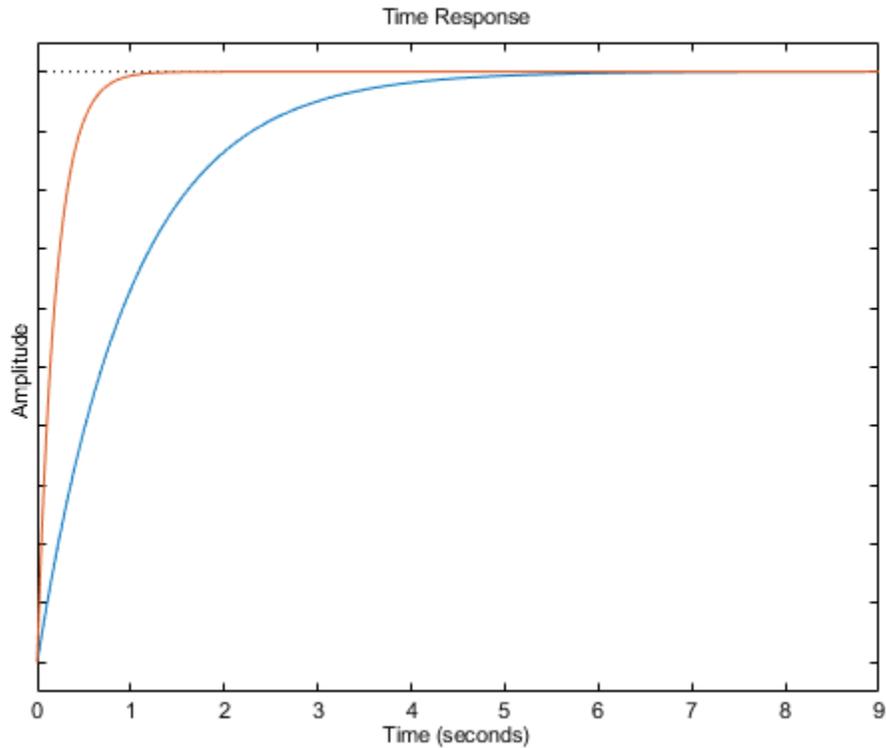
```
opt = timeoptions;
```

Enable plotting of normalized responses.

```
opt.Normalize = 'on';
```

Plot the step response of two transfer function models using the specified options.

```
sys1 = tf(10,[1,1]);  
sys2 = tf(5,[1,5]);  
stepplot(sys1,sys2,opt);
```



The plot shows the normalized step response for the two transfer function models.

## See Also

[getoptions](#) | [impzplot](#) | [initialplot](#) | [lsimplot](#) | [setoptions](#) | [stepplot](#)

**Introduced in R2012a**

# totaldelay

Total combined I/O delays for LTI model

## Syntax

```
td = totaldelay(sys)
```

## Description

`td = totaldelay(sys)` returns the total combined I/O delays for an LTI model `sys`. The matrix `td` combines contributions from the `InputDelay`, `OutputDelay`, and `ioDelayMatrix` properties.

Delays are expressed in seconds for continuous-time models, and as integer multiples of the sample period for discrete-time models. To obtain the delay times in seconds, multiply `td` by the sample time `sys.Ts`.

## Examples

### Compute Combined Input-Output Delay for Transfer Function

Create the transfer function model,  $1/s$ .

```
sys = tf(1,[1 0]);
```

Specify a 2 second input delay, and a 1.5 second output delay.

```
sys.InputDelay = 2;  
sys.OutputDelay = 1.5;
```

Compute the combined input-output delay for `sys`.

```
td = totaldelay(sys)
```

```
td = 3.5000
```

The resulting transfer function has the following form:

$$e^{-2s}x \frac{1}{s} e^{-1.5s} = e^{-3.5s} \frac{1}{s}$$

This result is equivalent to specifying an input-output delay of 3.5 seconds for the original transfer function,  $1/s$ .

## See Also

`absorbDelay` | `hasdelay`

**Introduced in R2012a**

# translatecov

Translate parameter covariance across model transformation operations

## Syntax

```
sys_new = translatecov(fcn,sys)
sys_new = translatecov(fcn,Input1,...,InputN)
```

## Description

`sys_new = translatecov(fcn,sys)` transforms `sys` into `sys_new = fcn(sys)`, and translates the parameter covariance of `sys` to the parameter covariance of the transformed model. `fcn` is a transformation function that you specify. The command computes the parameter covariance of `sys_new` by applying the Gauss Approximation formula. To view the translated parameter covariance, use `getcov`.

Applying model transformations directly does not always translate the parameter covariance of the original model to that of the transformed model. For example, `d2c(sys)` does not translate the parameter covariance of `sys`. In contrast, `translatecov(@(x)d2c(x),sys)` produces a transformed model that has the same coefficients as `d2c(sys)` and has the translated parameter covariance of `sys`.

`sys_new = translatecov(fcn,Input1,...,InputN)` returns the model `sys_new = fcn(Input1,...,InputN)` and its parameter covariance. At least one of the `N` inputs must be a linear model with parameter covariance information.

## Input Arguments

### **fcn**

Model transformation function, specified as a function handle.

For single-input functions, `sys_new = fcn(sys)`. The input to `fcn` must be an identified model with parameter covariance information. Typical single-input transformation operations include:

- Model-type conversion, such as `idpoly` and `idss`. For example, `fcn = @(x)idpoly(x)` converts the model `x` to a polynomial model.
- Sample time conversion, such as `c2d` and `d2c`. For example, `fcn = @(x)c2d(x,Ts)` converts the continuous-time model `x` to discrete-time using the specified sample time `Ts`.

For multi-input functions, `sys_new = fcn(Input1, . . . InputN)`. At least one of the input arguments `Input1, . . . , InputN` must be an identified model with parameter covariance information. Typical multi-input operations include merging and concatenation of multiple models. For example, `fcn = @(x,y)[x,y]` performs horizontal concatenation of the models `x` and `y`.

### **sys**

Linear model with parameter covariance information, specified as one of the following model types:

- `idtf`
- `idproc`
- `idss`
- `idpoly`
- `idgrey`

The model must contain parameter covariance information, that is `getcov(sys)` must be nonempty.

### **Input1, . . . , InputN**

Multiple input arguments to the translation function `fcn`. At least one of the `N` inputs must be a linear identified model with parameter covariance information. The other inputs can be any MATLAB data type. For an example, see “Translate Parameter Covariance to Closed-Loop Model” on page 1-1832.

## **Output Arguments**

### **sys\_new**

Model resulting from transformation operation. The model includes parameter covariance information.

## Examples

### Translate Parameter Covariance During Model Conversion

Convert an estimated transfer function model into state-space model while also translating the estimated parameter covariance.

Estimate a transfer function model.

```
load iddata1
sys1 = tfest(z1,2);
```

Convert the estimated model to state-space form while also translating the estimated parameter covariance.

```
sys2 = translatecov(@(x)idss(x),sys1);
```

If you convert the transfer function model to state-space form directly, the estimated parameter covariance is lost (the output of `getcov` is empty).

```
sys3 = idss(sys1);
getcov(sys3)
```

```
ans =
```

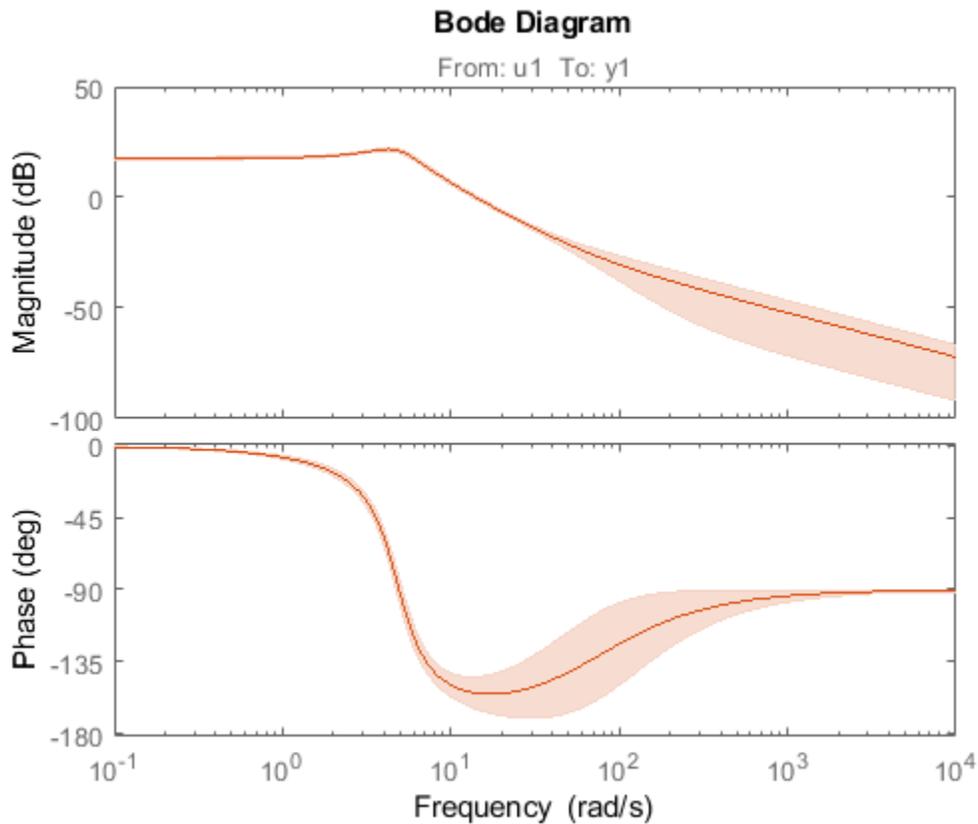
```
    []
```

View the parameter covariance in the estimated and converted models.

```
covsys1 = getcov(sys1);
covsys2 = getcov(sys2);
```

Compare the confidence regions.

```
h = bodeplot(sys1,sys2);
showConfidence(h,2);
```



The confidence bounds for sys1 overlaps with sys2.

### Translate Parameter Covariance During Model Concatenation

Concatenate 3 single-output models such that the covariance data from the 3 models combine to produce the covariance data for the resulting model.

Construct a state-space model.

```
a = [-1.1008 0.3733;0.3733 -0.9561];  
b = [0.7254 0.7147;-0.0631 -0.2050];
```

```
c = [-0.1241 0; 1.4897 0.6715; 1.4090 -1.2075];
d = [0 1.0347; 1.6302 0; 0.4889 0];
sys = idss(a,b,c,d,'Ts',0);
```

Generate multi-output estimation data.

```
t = (0:0.01:0.99)';
u = randn(100,2);
y = lsim(sys,u,t,'zoh');
y = y + rand(size(y))/10;
data = iddata(y,u,0.01);
```

Estimate a separate model for each output signal.

```
m1 = ssest(data(:,1,:),2,'feedthrough',true(1,2), 'DisturbanceModel', 'none');
m2 = ssest(data(:,2,:),2,'feedthrough',true(1,2), 'DisturbanceModel', 'none');
m3 = ssest(data(:,3,:),2,'feedthrough',true(1,2), 'DisturbanceModel', 'none');
```

Combine the estimated models while also translating the covariance information.

```
f = @(x,y,z)[x;y;z];
M2 = translatecov(f, m1, m2, m3);
```

The parameter covariance is not empty.

```
getcov(M2, 'factors')

ans = struct with fields:
    R: [36x36 double]
    T: [24x36 double]
    Free: [90x1 logical]
```

If you combine the estimated models into one 3-output model directly, the covariance information is lost (the output of `getcov` is empty).

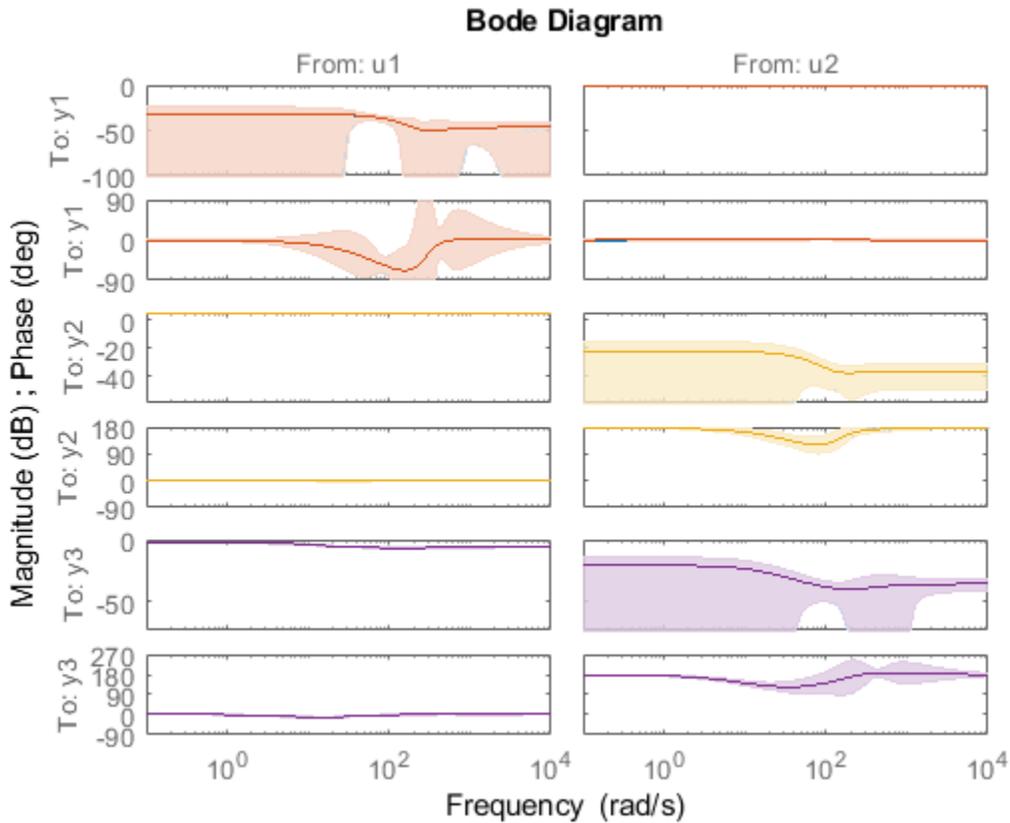
```
M1 = [m1;m2;m3];
getcov(M1)

ans =

    []
```

Compare the confidence bounds.

```
h = bodeplot(M2, m1, m2, m3);
showConfidence(h);
```



The confidence bounds for M2 overlap with those of m1, m2 and m3 models on their respective plot axes.

### Translate Parameter Covariance to Closed-Loop Model

Consider a closed-loop feedback model consisting of a plant and controller. Translate the parameter covariance of the plant to the closed-loop feedback model.

Estimate a plant as a fourth-order state-space model using estimation data z1.

```
load iddata1 z1  
Plant = ssest(z1,4);
```

Plant contains parameter covariance information.

Create a controller as a continuous-time zero-pole-gain model with zeros, poles, and gain equal to -2, -10, 5, respectively.

```
Controller = zp(-2, -10, 5);
```

Define a transformation function to generate the closed-loop feedback state-space model.

```
fcn = @(x,y) idss(feedback(x,y));
```

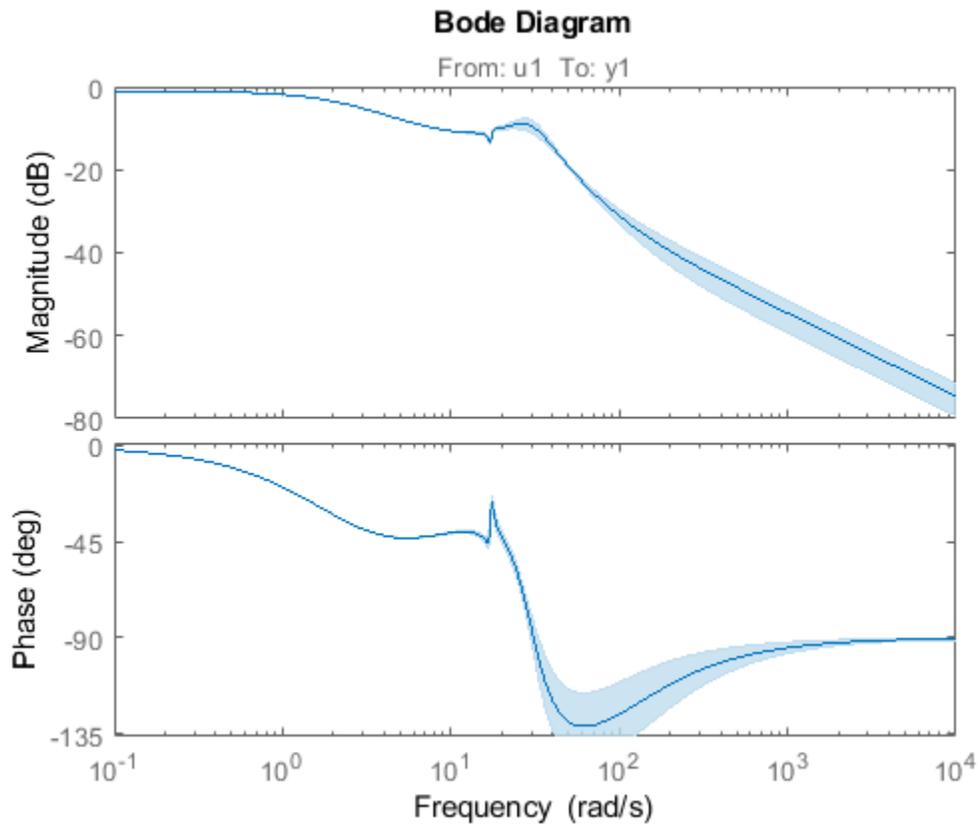
Translate the parameter covariance of the plant to the closed-loop feedback model.

```
sys_new = translatecov(fcn,Plant,Controller);
```

sys\_new contains the translated parameter covariance information.

Plot the frequency-response of the transformed model sys\_new, and view the confidence region of the response.

```
h = bodeplot(sys_new);  
showConfidence(h);
```



The plot shows the effect of the uncertainty in Plant on the closed-loop response.

## Tips

- If you obtained `sys` through estimation and have access to the estimation data, you can use zero-iteration update to recompute the parameter covariance. For example:

```
load iddata1
m = ssest(z1,4);
opt = ssestOptions
opt.SearchOptions.MaxIterations = 0;
m_new = ssest(z1,m2,opt)
```

You cannot run a zero-iteration update in the following cases:

- If `MaxIterations` option, which depends on the `SearchMethod` option, is not available.
- For some model and data types. For example, a continuous-time `idpoly` model using time-domain data.

## Algorithms

`translatecov` uses numerical perturbations of individual parameters of `sys` to compute the Jacobian of `fcn(sys)` parameters with respect to parameters of `sys`. `translatecov`

then applies Gauss Approximation formula  $cov\_new = J \times cov \times J^T$  to translate the covariance, where `J` is the Jacobian matrix. This operation can be slow for models containing a large number of free parameters.

## See Also

`getcov` | `getpvec` | `rsample` | `setcov`

## Topics

“What Is Model Covariance?”

“Types of Model Uncertainty Information”

**Introduced in R2012b**

## treepartition

Class representing binary-tree nonlinearity estimator for nonlinear ARX models

### Syntax

```
t=treepartition(Property1,Value1,...PropertyN,ValueN)  
t=treepartition('NumberOfUnits',N)
```

### Description

treepartition is an object that stores the binary-tree nonlinear estimator for estimating nonlinear ARX models. The object defines a nonlinear function  $y = F(x)$ , where  $F$  is a piecewise-linear (affine) function of  $x$ ,  $y$  is scalar, and  $x$  is a 1-by- $m$  vector. Compute the value of  $F$  using `evaluate(t,x)`, where  $t$  is the treepartition object at  $x$ .

### Construction

`t=treepartition(Property1,Value1,...PropertyN,ValueN)` creates a binary tree nonlinearity estimator object specified by properties in “treepartition Properties” on page 1-1837. The tree has the number of leaves equal to  $2^{(J+1)} - 1$ , where  $J$  is the number of nodes in the tree and set by the property `NumberOfUnits`. The default value of `NumberOfUnits` is computed automatically and sets an upper limit on the actual number of tree nodes used by the estimator.

`t=treepartition('NumberOfUnits',N)` creates a binary tree nonlinearity estimator object with  $N$  terms in the binary tree expansion (the number of nodes in the tree). When you estimate a model containing  $t$ , the value of the `NumberOfUnits` property,  $N$ , in  $t$  is automatically changed to show the actual number of leaves used—which is the largest integer of the form  $2^n - 1$  and less than or equal to  $N$ .

## treepartition Properties

You can include property-value pairs in the constructor to specify the object.

After creating the object, you can use `get` or dot notation to access the object property values. For example:

```
% List all property values
get(t)
% Get value of NumberOfUnits property
t.NumberOfUnits
```

You can also use the `set` function to set the value of particular properties. For example:

```
set(t, 'NumberOfUnits', 5)
```

The first argument to `set` must be the name of a MATLAB variable.

Property Name	Description
NumberOfUnits	<p>Integer specifies the number of nodes in the tree. Default= 'auto' selects the number of nodes from the data using the pruning algorithm.</p> <p>When you estimate a model containing a <code>treepartition</code> nonlinearity, the value of <code>NumberOfUnits</code> is automatically changed to show the actual number of leaves used—which is the largest integer of the form <math>2^n - 1</math> and less than or equal to <math>N</math> (the integer value of units you specify).</p> <p>For example:</p> <pre>treepartition('NumberOfUnits',5)</pre>

Property Name	Description
Parameters	<p>Structure containing the following fields:</p> <ul style="list-style-type: none"> <li>• <b>RegressorMean</b>: 1-by-m vector containing the means of <math>x</math> in estimation data, <math>r</math>.</li> <li>• <b>RegressorMinMax</b>: m-by-2 matrix containing the maximum and minimum estimation-data regressor values.</li> <li>• <b>OutputOffset</b>: scalar <math>d</math>.</li> <li>• <b>LinearCoef</b>: m-by-1 vector <math>L</math>.</li> <li>• <b>SampleLength</b>: Length of estimation data.</li> <li>• <b>NoiseVariance</b>: Estimated variance of the noise in estimation data.</li> <li>• <b>Tree</b>: A structure containing the following tree parameters: <ul style="list-style-type: none"> <li>• <b>TreeLevelPtr</b>: N-by-1 vector containing the levels <math>j</math> of each node.</li> <li>• <b>AncestorDescendantPtr</b>: N-by-3 matrix, such that the entry <math>(k, 1)</math> is the ancestor of node <math>k</math>, and entries <math>(k, 2)</math> and <math>(k, 3)</math> are the left and right descendants, respectively.</li> <li>• <b>LocalizingVectors</b>: N-by-(m+1) matrix, such that the <math>r</math>th row is <math>B_r</math>.</li> <li>• <b>LocalParVector</b>: N-by-(m+1) matrix, such that the <math>k</math>th row is <math>C_k</math>.</li> <li>• <b>LocalCovMatrix</b>: N-by-((m+1)m/2) matrix such that the <math>k</math>th row is the covariance matrix of <math>C_k</math>. <math>C_k</math> is reshaped as a row vector.</li> </ul> </li> </ul>

Property Name	Description
Options	<p>Structure containing the following fields that affect the initial model:</p> <ul style="list-style-type: none"> <li>• <b>FinestCell</b>: Integer or character vector specifying the minimum number of data points in the smallest partition. Default: 'auto', which computes the value from the data.</li> <li>• <b>Threshold</b>: Threshold parameter used by the adaptive pruning algorithm. Smaller threshold value corresponds to a shorter branch that is terminated by the active partition <math>D_a</math>. Higher threshold value results in a longer branch. Default: 1.0.</li> <li>• <b>Stabilizer</b>: Penalty parameter of the penalized least-squares algorithm used to compute local parameter vectors <math>C_k</math>. Higher stabilizer value improves stability, but may deteriorate the accuracy of the least-square estimate. Default: 1e-6.</li> </ul>

## Examples

Use `treepartition` to specify the nonlinear estimator in nonlinear ARX models. For example:

```
m=nlarx(Data,Orders,treepartition('num',5));
```

The following commands provide an example of using advanced `treepartition` options:

```
% Define the treepartition object.
t=treepartition('num',100);
% Set the Threshold, which is a field
% in the Options structure.
t.Options.Threshold=2;
% Estimate the nonlinear ARX model.
m=nlarx(Data,Orders,t);
```

## Algorithms

The mapping  $F$  is defined by a dyadic partition  $P$  of the  $x$ -space, such that on each partition element  $P_k$ ,  $F$  is a linear mapping. When  $x$  belongs to  $P_k$ ,  $F(x)$  is given by:

$$F(x) = d + xL + (1, x)C_k,$$

where  $L$  is 1-by- $m$  vector and  $d$  is a scalar common for all elements of partition.  $C_k$  is a 1-by- $(m+1)$  vector.

The mapping  $F$  and associated partition  $P$  of the  $x$ -space are computed as follows:

- 1 Given the value of  $J$ , a dyadic tree with  $J$  levels and  $N = 2^{J-1}$  nodes is initialized.
- 2 Each node at level  $1 < j < J$  has two descendants at level  $j + 1$  and one parent at level  $j - 1$ .
  - The root node at level 1 has two descendants.
  - Nodes at level  $J$  are terminating leaves of the tree and have one parent.
- 3 One partition element is associated to each node  $k$  of the tree.
  - The vector of coefficients  $C_k$  is computed using the observations on the corresponding partition element  $P_k$  by the penalized least-squares algorithm.
  - When the node  $k$  is not a terminating leaf, the partition element  $P_k$  is cut into two to obtain the partition elements of descendant nodes. The cut is defined by the half-spaces  $(1, x)B_k > 0$  or  $\leq 0$  (move to left or right descendant), where  $B_k$  is chosen to improve the stability of least-square computation on the partitions at the descendant nodes.
- 4 When the value of the mapping  $F$ , defined by the `treepartition` object, is computed at  $x$ , an adaptive algorithm selects the *active node*  $k$  of the tree on the branch of partitions which contain  $x$ .

When the `Focus` option in `nlarxOptions` is 'prediction', `treepartition` uses a noniterative technique for estimating parameters. Iterative refinements are not possible for models containing this nonlinearity estimator.

You cannot use `treepartition` when `Focus` is 'simulation' because this nonlinearity estimator is not differentiable. Minimization of simulation error requires differentiable nonlinear functions.

## See Also

`nlarx`

**Introduced in R2007a**

## TrendInfo

Offset and linear trend slope values for detrending data

### Description

TrendInfo class represents offset and linear trend information of input and output data. Constructing the corresponding object lets you:

- Compute and store mean values or best-fit linear trends of input and output data signals.
- Define specific offsets and trends to be removed from input-output data.

By storing offset and trend information, you can apply it to multiple data sets.

After estimating a linear model from detrended data, you can simulate the model at original operation conditions by adding the saved trend to the simulated output using `retrend`.

### Construction

For transient data, if you want to define a specific offset or trend to be removed from this data, create the TrendInfo object using `getTrend`. For example:

```
T = getTrend(data)
```

where `data` is the `iddata` object from which you will be removing the offset or linear trend, and `T` is the `TrendInfo` object. You must then assign specific offset and slope values as properties of this object before passing the object as an argument to `detrend`.

For steady-state data, if you want to detrend the data and store the trend information, use the `detrend` command with the `output` argument for storing trend information.

## Properties

After creating the object, you can use `get` or dot notation to access the object property values.

Property Name	Default	Description
DataName	''	Name of the <code>iddata</code> object from which trend information is derived (if any)
InputOffset	<code>zeros(1,nu)</code> , where <code>nu</code> is the number of inputs	<ul style="list-style-type: none"> <li>For transient data, the physical equilibrium offset you specify for each input signal.</li> <li>For steady-state data, the mean of input values. Computed automatically when detrending the data.</li> <li>If removing a linear trend from the input-output data, the value of the line at <code>t0</code>, where <code>t0</code> is the start time.</li> </ul> <p>For multiple experiment data, this is a cell array of size equal to the number of experiments in the data set.</p>
InputSlope	<code>zeros(1,nu)</code> , where <code>nu</code> is the number of inputs	<p>Slope of linear trend in input data, computed automatically when using the <code>detrend</code> command to remove the linear trend in the data.</p> <p>For multiple experiment data, this is a cell array of size equal to the number of experiments in the data set.</p>
OutputOffset	<code>zeros(1,ny)</code> , where <code>ny</code> is the number of outputs	<ul style="list-style-type: none"> <li>For transient data, the physical equilibrium offset you specify for each output signal</li> <li>For steady-state data, the mean of output values. Computed automatically when detrending the data.</li> <li>If removing a linear trend from the input-output data, the value of the line at <code>t0</code>, where <code>t0</code> is the start time.</li> </ul> <p>For multiple experiment data, this is a cell array of size equal to the number of experiments in the data set.</p>

Property Name	Default	Description
OutputSlope	<code>zeros(1,ny)</code> , where <code>ny</code> is the number of outputs	<p>Slope of linear trend in output data, computed automatically when using the <code>detrend</code> command to remove the linear trend in the data.</p> <p>For multiple experiment data, this is a cell array of size equal to the number of experiments in the data set.</p>

## Examples

### Remove Offsets From Data

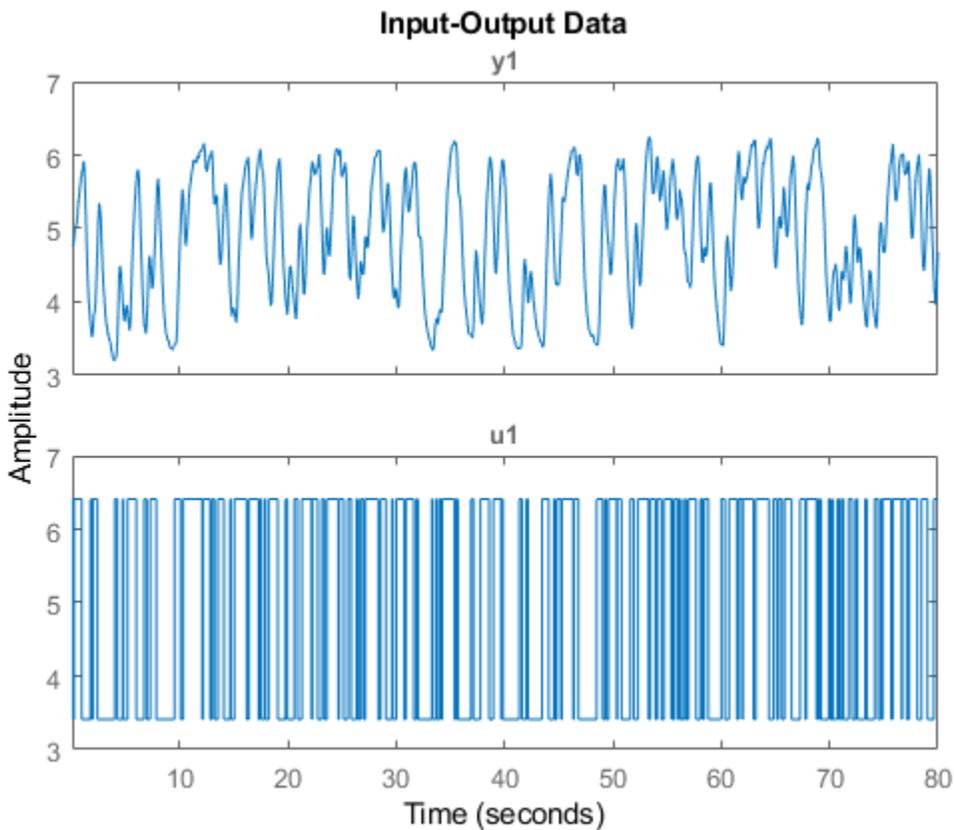
Remove specified offset from input and output signals.

Load SISO data containing vectors `u2` and `y2`.

```
load dryer2
```

Create a data object with sample time of 0.08 seconds and plot it.

```
data = iddata(y2,u2,0.08);  
plot(data)
```



The data has a nonzero mean value.

Store the data offset and trend information in a `TrendInfo` object.

```
T = getTrend(data);
```

Assign offset values to the `TrendInfo` object.

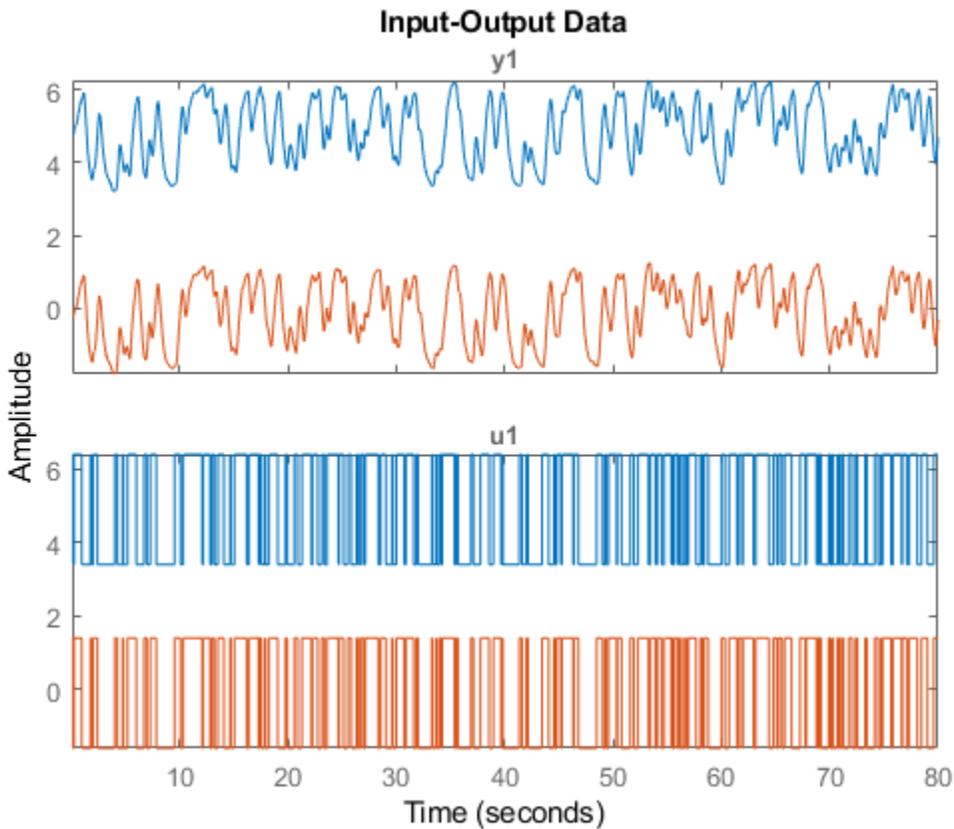
```
T.InputOffset = 5;  
T.OutputOffset = 5;
```

Subtract offset from the data.

```
data_d = detrend(data,T);
```

Plot the detrended data on the same plot.

```
hold on  
plot(data_d)
```



View the mean value removed from the data.

```
get(T)
```

```
ans = struct with fields:  
    DataName: 'data'  
    InputOffset: 5  
    OutputOffset: 5  
    InputSlope: 0
```

```
OutputSlope: 0
```

### Store Trend Information

Construct the TrendInfo object that stores trend information as part of data detrending.

Load SISO data containing vectors u2 and y2.

```
load dryer2
```

Create data object with sample time of 0.08 seconds.

```
data = iddata(y2,u2,0.08);
```

Detrend the mean from the data and store the mean as a TrendInfo object T.

```
[data_d,T] = detrend(data,0);
```

View the mean value removed from the data.

```
get(T)
```

```
ans = struct with fields:
  DataName: 'data'
  InputOffset: 5.0000
  OutputOffset: 4.8901
  InputSlope: 0
  OutputSlope: 0
```

## See Also

[detrend](#) | [getTrend](#) | [retrend](#)

## Topics

“Handling Offsets and Trends in Data”

**Introduced in R2009a**

## unitgain

Specify absence of nonlinearities for specific input or output channels in Hammerstein-Wiener models

### Syntax

```
unit=unitgain
```

### Description

`unit=unitgain` instantiates an object that specifies an identity mapping  $F(x)=x$  to exclude specific input and output channels from being affected by a nonlinearity in Hammerstein-Wiener models.

Use the `unitgain` object as an argument in the `n_lhw` estimator to set the corresponding channel nonlinearity to unit gain.

For example, for a two-input and one-output model, to exclude the second input from being affected by a nonlinearity, use the following syntax:

```
m = n_lhw(data,orders,['saturation' 'unitgain'],'deadzone')
```

In this case, the first input saturates and the output has an associated deadzone nonlinearity.

### unitgain Properties

`unitgain` does not have properties.

### Examples

For example, for a one-input and one-output model, to exclude the output from being affected by a nonlinearity, use the following syntax:

```
m = nlhw(Data,Orders,'saturation','unitgain')
```

In this case, the input has a saturation nonlinearity.

If nonlinearities are absent in input or output channels, you can replace `unitgain` with an empty matrix. For example, to specify a Wiener model with a sigmoid nonlinearity at the output and a unit gain at the input, use the following command:

```
m = nlhw(Data,Orders,[],'sigmoid');
```

## Tips

Use the `unitgain` object to exclude specific input and output channels from being affected by a nonlinearity in Hammerstein-Wiener models. `unitgain` is a linear function

$y = F(x)$ , where  $F(x)=x$ .

## See Also

[deadzone](#) | [nlhw](#) | [saturation](#) | [sigmoidnet](#)

**Introduced in R2007a**

## unscentedKalmanFilter

Create unscented Kalman filter object for online state estimation

### Syntax

```
obj = unscentedKalmanFilter(StateTransitionFcn,MeasurementFcn,  
InitialState)
```

```
obj = unscentedKalmanFilter(StateTransitionFcn,MeasurementFcn,  
InitialState,Name,Value)
```

```
obj = unscentedKalmanFilter(StateTransitionFcn,MeasurementFcn)
```

```
obj = unscentedKalmanFilter(StateTransitionFcn,MeasurementFcn,  
Name,Value)
```

```
obj = unscentedKalmanFilter(Name,Value)
```

### Description

`obj = unscentedKalmanFilter(StateTransitionFcn,MeasurementFcn,InitialState)` creates an unscented Kalman filter object for online state estimation of a discrete-time nonlinear system. `StateTransitionFcn` is a function that calculates the state of the system at time  $k$ , given the state vector at time  $k-1$ . `MeasurementFcn` is a function that calculates the output measurement of the system at time  $k$ , given the state at time  $k$ . `InitialState` specifies the initial value of the state estimates.

After creating the object, use the `correct` and `predict` commands to update state estimates and state estimation error covariance values using a discrete-time unscented Kalman filter algorithm and real-time data.

`obj = unscentedKalmanFilter(StateTransitionFcn,MeasurementFcn,InitialState,Name,Value)` specifies additional attributes of the unscented Kalman filter object using one or more `Name,Value` pair arguments.

`obj = unscentedKalmanFilter(StateTransitionFcn,MeasurementFcn)` creates an unscented Kalman filter object using the specified state transition and measurement functions. Before using the `predict` and `correct` commands, specify the initial state

values using dot notation. For example, for a two-state system with initial state values  $[1;0]$ , specify `obj.State = [1;0]`.

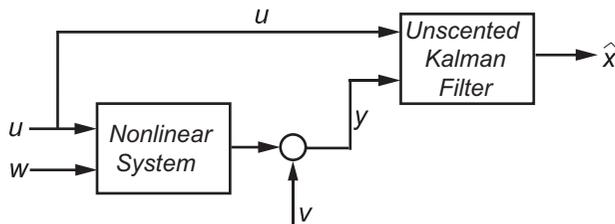
`obj = unscentedKalmanFilter(StateTransitionFcn,MeasurementFcn, Name, Value)` specifies additional attributes of the unscented Kalman filter object using one or more `Name, Value` pair arguments. Before using the `predict` and `correct` commands, specify the initial state values using `Name, Value` pair arguments or dot notation.

`obj = unscentedKalmanFilter(Name, Value)` creates an unscented Kalman filter object with properties specified using one or more `Name, Value` pair arguments. Before using the `predict` and `correct` commands, specify the state transition function, measurement function, and initial state values using `Name, Value` pair arguments or dot notation.

## Object Description

`unscentedKalmanFilter` creates an object for online state estimation of a discrete-time nonlinear system using the discrete-time unscented Kalman filter algorithm.

Consider a plant with states  $x$ , input  $u$ , output  $y$ , process noise  $w$ , and measurement noise  $v$ . Assume that you can represent the plant as a nonlinear system.



The algorithm computes the state estimates  $\hat{x}$  of the nonlinear system using state transition and measurement functions specified by you. The software lets you specify the noise in these functions as additive or nonadditive:

- **Additive Noise Terms** — The state transition and measurements equations have the following form:

$$x[k] = f(x[k-1], u_s[k-1]) + w[k-1]$$

$$y[k] = h(x[k], u_m[k]) + v[k]$$

Here  $f$  is a nonlinear state transition function that describes the evolution of states  $x$  from one time step to the next. The nonlinear measurement function  $h$  relates  $x$  to the measurements  $y$  at time step  $k$ .  $w$  and  $v$  are the zero-mean, uncorrelated process and measurement noises, respectively. These functions can also have additional input arguments that are denoted by  $u_s$  and  $u_m$  in the equations. For example, the additional arguments could be time step  $k$  or the inputs  $u$  to the nonlinear system. There can be multiple such arguments.

Note that the noise terms in both equations are additive. That is,  $x(k)$  is linearly related to the process noise  $w(k-1)$ , and  $y(k)$  is linearly related to the measurement noise  $v(k)$ .

- **Nonadditive Noise Terms** — The software also supports more complex state transition and measurement functions where the state  $x[k]$  and measurement  $y[k]$  are nonlinear functions of the process noise and measurement noise, respectively. When the noise terms are nonadditive, the state transition and measurements equation have the following form:

$$x[k] = f(x[k-1], w[k-1], u_s[k-1])$$

$$y[k] = h(x[k], v[k], u_m[k])$$

When you perform online state estimation, you first create the nonlinear state transition function  $f$  and measurement function  $h$ . You then construct the `unscentedKalmanFilter` object using these nonlinear functions, and specify whether the noise terms are additive or nonadditive. After you create the object, you use the `predict` command to predict state estimates at the next time step, and `correct` to correct state estimates using the unscented Kalman filter algorithm and real-time data. For information about the algorithm, see “Extended and Unscented Kalman Filter Algorithms for Online State Estimation”.

You can use the following commands with `unscentedKalmanFilter` objects:

Command	Description
<code>correct</code>	Correct the state and state estimation error covariance at time step $k$ using measured data at time step $k$ .
<code>predict</code>	Predict the state and state estimation error covariance at time the next time step.

Command	Description
clone	<p>Create another object with the same object property values.</p> <p>Do not create additional objects using syntax <code>obj2 = obj</code>. Any changes made to the properties of the new object created in this way (<code>obj2</code>) also change the properties of the original object (<code>obj</code>).</p>

For `unscentedKalmanFilter` object properties, see “Properties” on page 1-1859.

## Examples

### Create Unscented Kalman Filter Object for Online State Estimation

To define an unscented Kalman filter object for estimating the states of your system, you write and save the state transition function and measurement function for the system.

In this example, use the previously written and saved state transition and measurement functions, `vdpStateFcn.m` and `vdpMeasurementFcn.m`. These functions describe a discrete-approximation to van der Pol oscillator with nonlinearity parameter,  $\mu$ , equal to 1. The oscillator has two states.

Specify an initial guess for the two states. You specify the initial state guess as an  $M$ -element row or column vector, where  $M$  is the number of states.

```
initialStateGuess = [1;0];
```

Create the unscented Kalman filter object. Use function handles to provide the state transition and measurement functions to the object.

```
obj = unscentedKalmanFilter(@vdpStateFcn,@vdpMeasurementFcn,initialStateGuess);
```

The object has a default structure where the process and measurement noise are additive.

To estimate the states and state estimation error covariance from the constructed object, use the `correct` and `predict` commands and real-time data.

## Specify Process and Measurement Noise Covariances in Unscented Kalman Filter Object

Create an unscented Kalman filter object for a van der Pol oscillator with two states and one output. Use the previously written and saved state transition and measurement functions, `vdpStateFcn.m` and `vdpMeasurementFcn.m`. These functions are written for additive process and measurement noise terms. Specify the initial state values for the two states as `[2;0]`.

Since the system has two states and the process noise is additive, the process noise is a 2-element vector and the process noise covariance is a 2-by-2 matrix. Assume there is no cross-correlation between process noise terms, and both the terms have the same variance 0.01. You can specify the process noise covariance as a scalar. The software uses the scalar value to create a 2-by-2 diagonal matrix with 0.01 on the diagonals.

Specify the process noise covariance during object construction.

```
obj = unscentedKalmanFilter(@vdpStateFcn,@vdpMeasurementFcn,[2;0],...  
    'ProcessNoise',0.01);
```

Alternatively, you can specify noise covariances after object construction using dot notation. For example, specify the measurement noise covariance as 0.2.

```
obj.MeasurementNoise = 0.2;
```

Since the system has only one output, the measurement noise is a 1-element vector and the `MeasurementNoise` property denotes the variance of the measurement noise.

## Specify Nonadditive Measurement Noise in Unscented Kalman Filter Object

Create an unscented Kalman filter object for a van der Pol oscillator with two states and one output. Assume that the process noise terms in the state transition function are additive. That is, there is a linear relation between the state and process noise. Also assume that the measurement noise terms are nonadditive. That is, there is a nonlinear relation between the measurement and measurement noise.

```
obj = unscentedKalmanFilter('HasAdditiveMeasurementNoise',false);
```

Specify the state transition function and measurement functions. Use the previously written and saved functions, `vdpStateFcn.m` and `vdpMeasurementNonAdditiveNoiseFcn.m`.

The state transition function is written assuming the process noise is additive. The measurement function is written assuming the measurement noise is nonadditive.

```
obj.StateTransitionFcn = @vdpStateFcn;
obj.MeasurementFcn = @vdpMeasurementNonAdditiveNoiseFcn;
```

Specify the initial state values for the two states as `[2;0]`.

```
obj.State = [2;0];
```

You can now use the `correct` and `predict` commands to estimate the state and state estimation error covariance values from the constructed object.

### Specify Additional Inputs in State Transition and Measurement Functions

Consider a nonlinear system with input  $u$  whose state  $x$  and measurement  $y$  evolve according to the following state transition and measurement equations:

$$x[k] = \sqrt{x[k-1] + u[k-1]} + w[k-1]$$

$$y[k] = x[k] + 2 * u[k] + v[k]^2$$

The process noise  $w$  of the system is additive while the measurement noise  $v$  is nonadditive.

Create the state transition function and measurement function for the system. Specify the functions with an additional input  $u$ .

```
f = @(x,u)(sqrt(x+u));
h = @(x,v,u)(x+2*u+v^2);
```

`f` and `h` are function handles to the anonymous functions that store the state transition and measurement functions, respectively. In the measurement function, because the measurement noise is nonadditive,  $v$  is also specified as an input. Note that  $v$  is specified as an input before the additional input  $u$ .

Create an unscented Kalman filter object for estimating the state of the nonlinear system using the specified functions. Specify the initial value of the state as 1, and the measurement noise as nonadditive.

```
obj = unscentedKalmanFilter(f,h,1,'HasAdditiveMeasurementNoise',false);
```

Specify the measurement noise covariance.

```
obj.MeasurementNoise = 0.01;
```

You can now estimate the state of the system using the `predict` and `correct` commands. You pass the values of `u` to `predict` and `correct`, which in turn pass them to the state transition and measurement functions, respectively.

Correct the state estimate with measurement  $y[k]=0.8$  and input  $u[k]=0.2$  at time step  $k$ .

```
correct(obj,0.8,0.2)
```

Predict the state at next time step, given  $u[k]=0.2$ .

```
predict(obj,0.2)
```

## Input Arguments

### StateTransitionFcn — State transition function

function handle

State transition function  $f$ , specified as a function handle. The function calculates the  $N_s$ -element state vector of the system at time step  $k$ , given the state vector at time step  $k-1$ .  $N_s$  is the number of states of the nonlinear system.

You write and save the state transition function for your nonlinear system, and use it to construct the object. For example, if `vdpStateFcn.m` is the state transition function, specify `StateTransitionFcn` as `@vdpStateFcn`. You can also specify `StateTransitionFcn` as a function handle to an anonymous function.

The inputs to the function you write depend on whether you specify the process noise as additive or nonadditive in the `HasAdditiveProcessNoise` property of the object:

- `HasAdditiveProcessNoise` is true — The process noise  $w$  is additive, and the state transition function specifies how the states evolve as a function of state values at the previous time step:

$$x(k) = f(x(k-1), Us1, \dots, Usn)$$

Where  $x(k)$  is the estimated state at time  $k$ , and  $Us1, \dots, Usn$  are any additional input arguments required by your state transition function, such as system inputs or the sample time. During estimation, you pass these additional arguments to the `predict` command, which in turn passes them to the state transition function.

- `HasAdditiveProcessNoise` is false — The process noise is nonadditive, and the state transition function also specifies how the states evolve as a function of the process noise:

$$x(k) = f(x(k-1), w(k-1), Us1, \dots, Usn)$$

To see an example of a state transition function with additive process noise, type `edit vdpStateFcn` at the command line.

### MeasurementFcn — Measurement function

function handle

Measurement function  $h$ , specified as a function handle. The function calculates the  $N$ -element output measurement vector of the nonlinear system at time step  $k$ , given the state vector at time step  $k$ .  $N$  is the number of measurements of the system. You write and save the measurement function, and use it to construct the object. For example, if `vdpMeasurementFcn.m` is the measurement function, specify `MeasurementFcn` as `@vdpMeasurementFcn`. You can also specify `MeasurementFcn` as a function handle to an anonymous function.

The inputs to the function depend on whether you specify the measurement noise as additive or nonadditive in the `HasAdditiveMeasurementNoise` property of the object:

- `HasAdditiveMeasurementNoise` is true — The measurement noise  $v$  is additive, and the measurement function specifies how the measurements evolve as a function of state values:

$$y(k) = h(x(k), Um1, \dots, Umn)$$

Where  $y(k)$  and  $x(k)$  are the estimated output and estimated state at time  $k$ , and  $Um1, \dots, Umn$  are any optional input arguments required by your measurement function. For example, if you are using multiple sensors for tracking an object, an

additional input could be the sensor position. During estimation, you pass these additional arguments to the `correct` command, which in turn passes them to the measurement function.

- `HasAdditiveMeasurementNoise` is false — The measurement noise is nonadditive, and the measurement function also specifies how the output measurement evolves as a function of the measurement noise:

$$y(k) = h(x(k), v(k), Um1, \dots, Umn)$$

To see an example of a measurement function with additive process noise, type `edit vdpMeasurementFcn` at the command line. To see an example of a measurement function with nonadditive process noise, type `edit vdpMeasurementNonAdditiveNoiseFcn`.

### InitialState — Initial state estimates

vector

Initial state estimates, specified as an  $N_s$ -element vector, where  $N_s$  is the number of states in the system. Specify the initial state values based on your knowledge of the system.

The specified value is stored in the `State` property of the object. If you specify `InitialState` as a column vector then `State` is also a column vector, and `predict` and `correct` commands return state estimates as a column vector. Otherwise, a row vector is returned.

If you want a filter with single-precision floating-point variables, specify `InitialState` as a single-precision vector variable. For example, for a two-state system with state transition and measurement functions `vdpStateFcn.m` and `vdpMeasurementFcn.m`, create the unscented Kalman filter object with initial states `[1;2]` as follows:

```
obj = unscentedKalmanFilter(@vdpStateFcn,@vdpMeasurementFcn,single([1;2]))
```

Data Types: `double` | `single`

### Name-Value Pair Arguments

Specify optional comma-separated pairs of `Name`, `Value` arguments. `Name` is the argument name and `Value` is the corresponding value. `Name` must appear inside quotes. You can specify several name and value pair arguments in any order as `Name1, Value1, ..., NameN, ValueN`.

Use `Name`, `Value` arguments to specify properties on page 1-1859 of `unscentedKalmanFilter` object during object creation. For example, to create an unscented Kalman filter object and specify the process noise covariance as 0.01:

```
obj = unscentedKalmanFilter(StateTransitionFcn,MeasurementFcn,InitialState, 'ProcessNoise',0.01);
```

## Properties

`unscentedKalmanFilter` object properties are of three types:

- Tunable properties that you can specify multiple times, either during object construction using `Name`, `Value` arguments, or any time afterwards during state estimation. After object creation, use dot notation to modify the tunable properties.

```
obj = unscentedKalmanFilter(StateTransitionFcn,MeasurementFcn,InitialState);
obj.ProcessNoise = 0.01;
```

The tunable properties are `State`, `StateCovariance`, `ProcessNoise`, `MeasurementNoise`, `Alpha`, `Beta`, and `Kappa`.

- Nontunable properties that you can specify once, either during object construction or afterward using dot notation. Specify these properties before state estimation using `correct` and `predict`. The `StateTransitionFcn` and `MeasurementFcn` properties belong to this category.
- Nontunable properties that you must specify during object construction. The `HasAdditiveProcessNoise` and `HasAdditiveMeasurementNoise` properties belong to this category.

### Alpha — Spread of sigma points

1e-3 (default) | scalar value between 0 and 1

Spread of sigma points around mean state value, specified as a scalar value between 0 and 1 ( $0 < \text{Alpha} \leq 1$ ).

The unscented Kalman filter algorithm treats the state of the system as a random variable with mean value `State` and variance `StateCovariance`. To compute the state and its statistical properties at the next time step, the algorithm first generates a set of state values distributed around the mean `State` value by using the unscented transformation. These generated state values are called sigma points. The algorithm uses each of the sigma points as an input to the state transition and measurement functions to get a new set of transformed state points and measurements. The transformed points are used to compute the state and state estimation error covariance value at the next time step.

The spread of the sigma points around the mean state value is controlled by two parameters `Alpha` and `Kappa`. A third parameter, `Beta`, impacts the weights of the transformed points during state and measurement covariance calculations:

- `Alpha` — Determines the spread of the sigma points around the mean state value. It is usually a small positive value. The spread of sigma points is proportional to `Alpha`. Smaller values correspond to sigma points closer to the mean state.
- `Kappa` — A second scaling parameter that is usually set to 0. Smaller values correspond to sigma points closer to the mean state. The spread is proportional to the square-root of `Kappa`.
- `Beta` — Incorporates prior knowledge of the distribution of the state. For Gaussian distributions, `Beta = 2` is optimal.

If you know the distribution of state and state covariance, you can adjust these parameters to capture the transformation of higher-order moments of the distribution. The algorithm can track only a single peak in the probability distribution of the state. If there are multiple peaks in the state distribution of your system, you can adjust these parameters so that the sigma points stay around a single peak. For example, choose a small `Alpha` to generate sigma points close to the mean state value.

For more information, see “Unscented Kalman Filter Algorithm”.

`Alpha` is a tunable property. You can change it using dot notation.

### **Beta — Characterization of state distribution**

2 (default) | scalar value greater than or equal to 0

Characterization of the state distribution that is used to adjust weights of transformed sigma points, specified as a scalar value greater than or equal to 0. For Gaussian distributions, `Beta = 2` is an optimal choice.

For more information, see the `Alpha` property description.

`Beta` is a tunable property. You can change it using dot notation.

### **HasAdditiveMeasurementNoise — Measurement noise characteristics**

true (default) | false

Measurement noise characteristics, specified as one of the following values:

- `true` — Measurement noise  $v$  is additive. The measurement function  $h$  that is specified in `MeasurementFcn` has the following form:

$$y(k) = h(x(k), Um1, \dots, Umn)$$

Where  $y(k)$  and  $x(k)$  are the estimated output and estimated state at time  $k$ , and  $Um1, \dots, Umn$  are any optional input arguments required by your measurement function.

- `false` — Measurement noise is nonadditive. The measurement function specifies how the output measurement evolves as a function of the state *and* measurement noise:

$$y(k) = h(x(k), v(k), Um1, \dots, Umn)$$

`HasAdditiveMeasurementNoise` is a nontunable property, and you can specify it only during object construction. You cannot change it using dot notation.

### **HasAdditiveProcessNoise — Process noise characteristics**

`true` (default) | `false`

Process noise characteristics, specified as one of the following values:

- `true` — Process noise  $w$  is additive. The state transition function  $f$  specified in `StateTransitionFcn` has the following form:

$$x(k) = f(x(k-1), Us1, \dots, Usn)$$

Where  $x(k)$  is the estimated state at time  $k$ , and  $Us1, \dots, Usn$  are any additional input arguments required by your state transition function.

- `false` — Process noise is nonadditive. The state transition function specifies how the states evolve as a function of the state *and* process noise at the previous time step:

$$x(k) = f(x(k-1), w(k-1), Us1, \dots, Usn)$$

`HasAdditiveProcessNoise` is a nontunable property, and you can specify it only during object construction. You cannot change it using dot notation.

### **Kappa — Spread of sigma points**

`0` (default) | scalar value between 0 and 3

Spread of sigma points around mean state value, specified as a scalar value between 0 and 3 ( $0 \leq \text{Kappa} \leq 3$ ). `Kappa` is typically specified as `0`. Smaller values correspond to sigma points closer to the mean state. The spread is proportional to the square-root of `Kappa`. For more information, see the `Alpha` property description.

`Kappa` is a tunable property. You can change it using dot notation.

## MeasurementFcn — Measurement function

function handle

Measurement function  $h$ , specified as a function handle. The function calculates the  $N$ -element output measurement vector of the nonlinear system at time step  $k$ , given the state vector at time step  $k$ .  $N$  is the number of measurements of the system. You write and save the measurement function and use it to construct the object. For example, if `vdpMeasurementFcn.m` is the measurement function, specify `MeasurementFcn` as `@vdpMeasurementFcn`. You can also specify `MeasurementFcn` as a function handle to an anonymous function.

The inputs to the function depend on whether you specify the measurement noise as additive or nonadditive in the `HasAdditiveMeasurementNoise` property of the object:

- `HasAdditiveMeasurementNoise` is true — The measurement noise  $v$  is additive, and the measurement function specifies how the measurements evolve as a function of state values:

$$y(k) = h(x(k), Um1, \dots, Umn)$$

Where  $y(k)$  and  $x(k)$  are the estimated output and estimated state at time  $k$ , and  $Um1, \dots, Umn$  are any optional input arguments required by your measurement function. For example, if you are using multiple sensors for tracking an object, an additional input could be the sensor position. During estimation, you pass these additional arguments to the correct command which in turn passes them to the measurement function.

- `HasAdditiveMeasurementNoise` is false — The measurement noise is nonadditive, and the measurement function also specifies how the output measurement evolves as a function of the measurement noise:

$$y(k) = h(x(k), v(k), Um1, \dots, Umn)$$

To see an example of a measurement function with additive process noise, type `edit vdpMeasurementFcn` at the command line. To see an example of a measurement function with nonadditive process noise, type `edit vdpMeasurementNonAdditiveNoiseFcn`.

`MeasurementFcn` is a nontunable property. You can specify it once before using the correct command either during object construction or using dot notation after object construction. You cannot change it after using the correct command.

## MeasurementNoise — Measurement noise covariance

1 (default) | scalar | matrix

Measurement noise covariance, specified as a scalar or matrix depending on the value of the `HasAdditiveMeasurementNoise` property:

- `HasAdditiveMeasurementNoise` is true — Specify the covariance as a scalar or an  $N$ -by- $N$  matrix, where  $N$  is the number of measurements of the system. Specify a scalar if there is no cross-correlation between measurement noise terms and all the terms have the same variance. The software uses the scalar value to create an  $N$ -by- $N$  diagonal matrix.
- `HasAdditiveMeasurementNoise` is false — Specify the covariance as a  $V$ -by- $V$  matrix, where  $V$  is the number of measurement noise terms. `MeasurementNoise` must be specified before using `correct`. After you specify `MeasurementNoise` as a matrix for the first time, to then change `MeasurementNoise` you can also specify it as a scalar. Specify as a scalar if there is no cross-correlation between the measurement noise terms and all the terms have the same variance. The software extends the scalar to a  $V$ -by- $V$  diagonal matrix with the scalar on the diagonals.

`MeasurementNoise` is a tunable property. You can change it using dot notation.

## ProcessNoise — Process noise covariance

1 (default) | scalar | matrix

Process noise covariance, specified as a scalar or matrix depending on the value of the `HasAdditiveProcessNoise` property:

- `HasAdditiveProcessNoise` is true — Specify the covariance as a scalar or an  $N_s$ -by- $N_s$  matrix, where  $N_s$  is the number of states of the system. Specify a scalar if there is no cross-correlation between process noise terms, and all the terms have the same variance. The software uses the scalar value to create an  $N_s$ -by- $N_s$  diagonal matrix.
- `HasAdditiveProcessNoise` is false — Specify the covariance as a  $W$ -by- $W$  matrix, where  $W$  is the number of process noise terms. `ProcessNoise` must be specified before using `predict`. After you specify `ProcessNoise` as a matrix for the first time, to then change `ProcessNoise` you can also specify it as a scalar. Specify as a scalar if there is no cross-correlation between the process noise terms and all the terms have the same variance. The software extends the scalar to a  $W$ -by- $W$  diagonal matrix.

`ProcessNoise` is a tunable property. You can change it using dot notation.

## State — State of nonlinear system

[] (default) | vector

State of the nonlinear system, specified as a vector of size  $N_s$ , where  $N_s$  is the number of states of the system.

When you use the `predict` command, `State` is updated with the predicted value at time step  $k$  using the state value at time step  $k-1$ . When you use the `correct` command, `State` is updated with the estimated value at time step  $k$  using measured data at time step  $k$ .

The initial value of `State` is the value you specify in the `InitialState` input argument during object creation. If you specify `InitialState` as a column vector, then `State` is also a column vector, and the `predict` and `correct` commands return state estimates as a column vector. Otherwise, a row vector is returned. If you want a filter with single-precision floating-point variables, you must specify `State` as a single-precision variable during object construction using the `InitialState` input argument.

`State` is a tunable property. You can change it using dot notation.

### **StateCovariance — State estimation error covariance**

1 (default) | scalar | matrix

State estimation error covariance, specified as a scalar or an  $N_s$ -by- $N_s$  matrix, where  $N_s$  is the number of states of the system. If you specify a scalar, the software uses the scalar value to create an  $N_s$ -by- $N_s$  diagonal matrix.

Specify a high value for the covariance when you do not have confidence in the initial state values that you specify in the `InitialState` input argument.

When you use the `predict` command, `StateCovariance` is updated with the predicted value at time step  $k$  using the state value at time step  $k-1$ . When you use the `correct` command, `StateCovariance` is updated with the estimated value at time step  $k$  using measured data at time step  $k$ .

`StateCovariance` is a tunable property. You can change it using dot notation after using the `correct` or `predict` commands.

### **StateTransitionFcn — State transition function**

function handle

State transition function  $f$ , specified as a function handle. The function calculates the  $N_s$ -element state vector of the system at time step  $k$ , given the state vector at time step  $k-1$ .  $N_s$  is the number of states of the nonlinear system.

You write and save the state transition function for your nonlinear system and use it to construct the object. For example, if `vdpStateFcn.m` is the state transition function, specify `StateTransitionFcn` as `@vdpStateFcn`. You can also specify `StateTransitionFcn` as a function handle to an anonymous function.

The inputs to the function you write depend on whether you specify the process noise as additive or nonadditive in the `HasAdditiveProcessNoise` property of the object:

- `HasAdditiveProcessNoise` is true — The process noise  $w$  is additive, and the state transition function specifies how the states evolve as a function of state values at previous time step:

$$x(k) = f(x(k-1), Us1, \dots, Usn)$$

Where  $x(k)$  is the estimated state at time  $k$ , and  $Us1, \dots, Usn$  are any additional input arguments required by your state transition function, such as system inputs or the sample time. During estimation, you pass these additional arguments to the `predict` command, which in turn passes them to the state transition function.

- `HasAdditiveProcessNoise` is false — The process noise is nonadditive, and the state transition function also specifies how the states evolve as a function of the process noise:

$$x(k) = f(x(k-1), w(k-1), Us1, \dots, Usn)$$

To see an example of a state transition function with additive process noise, type `edit vdpStateFcn` at the command line.

`StateTransitionFcn` is a nontunable property. You can specify it once before using the `predict` command either during object construction or using dot notation after object construction. You cannot change it after using the `predict` command.

## Output Arguments

### **obj** — unscented Kalman filter object for online state estimation

`unscentedKalmanFilter` object

Unscented Kalman filter object for online state estimation, returned as an `unscentedKalmanFilter` object. This object is created using the specified properties on page 1-1859. Use the `correct` and `predict` commands to estimate the state and state estimation error covariance using the unscented Kalman filter algorithm.

When you use `predict`, `obj.State` and `obj.StateCovariance` are updated with the predicted value at time step  $k$  using the state value at time step  $k-1$ . When you use `correct`, `obj.State` and `obj.StateCovariance` are updated with the estimated values at time step  $k$  using measured data at time step  $k$ .

## Extended Capabilities

### C/C++ Code Generation

Generate C and C++ code using MATLAB® Coder™.

Usage notes and limitations:

For more information, see “Generate Code for Online State Estimation in MATLAB”.

Supports MATLAB Function block: No

## See Also

### Functions

`clone` | `correct` | `extendedKalmanFilter` | `predict`

### Blocks

Extended Kalman Filter | Kalman Filter | Unscented Kalman Filter

### Topics

“Nonlinear State Estimation Using Unscented Kalman Filter and Particle Filter”

“Generate Code for Online State Estimation in MATLAB”

“What Is Online Estimation?”

“Extended and Unscented Kalman Filter Algorithms for Online State Estimation”

“Validate Online State Estimation at the Command Line”

“Troubleshoot Online State Estimation”

### Introduced in R2016b

# wavenet

Create a wavelet network nonlinearity estimator object

## Syntax

NL = wavenet  
 NL = wavenet(Name, Value)

## Description

NL = wavenet creates a default wavelet network nonlinearity estimator object for estimating nonlinear ARX and Hammerstein-Wiener models. Use dot notation to customize the object properties, if needed.

NL = wavenet(Name, Value) creates a wavelet network nonlinearity estimator object with properties specified by one or more Name, Value pair arguments. The properties that you do not specify retain their default value.

## Object Description

wavenet is an object that stores the wavelet network nonlinearity estimator for estimating nonlinear ARX and Hammerstein-Wiener models.

Use wavenet to define a nonlinear function  $y = F(x, \theta)$ , where  $y$  is scalar,  $x$  is an  $m$ -dimensional row vector of regressors, and  $\theta$  represent the parameters in wavelet expansion. The wavelet network function is based on the following function expansion:

$$\begin{aligned}
 F(x, \theta) = & (x - r)PL + a_{s\_1}f(b_{s\_1}((x - r)Q - c_{s\_1})) + \dots \\
 & + a_{s\_ns}f(b_{s\_ns}((x - r)Q - c_{s\_ns})) \\
 & + a_{w\_1}g(b_{w\_1}((x - r)Q - c_{w\_1})) + \dots \\
 & + a_{w\_nw}g(b_{w\_nw}((x - r)Q - c_{w\_nw})) + d
 \end{aligned}$$

$$f(z) = e^{-0.5zz^T}$$

$$g(z) = (m - zz^T)e^{-0.5zz^T}$$

Here,

- $f(z)$  is a radial function called the scaling function, and  $z$  is the input to the scaling function.  $z$  is a 1-by- $q$  row vector.  $q$  is the number of components of  $x$  used in the scaling and wavelet functions.
- $g(z)$  is a radial function called the wavelet function, and  $z$  is the input to the wavelet function.
- $\theta$  represents the following parameters of the nonlinearity estimator:
  - $P$  and  $Q$  — Projection matrices of size  $m$ -by- $p$  and  $m$ -by- $q$ , respectively.

$P$  and  $Q$  are determined by principal component analysis of estimation data. Usually,  $p = m$ . If the components of  $x$  in the estimation data are linearly dependent, then  $p < m$ . The number of columns of  $Q$  is  $q$ .  $q$  is the number of components of  $x$  used in the scaling and wavelet function.

When used in a nonlinear ARX model,  $q$  is equal to the size of the `NonlinearRegressors` property of the `idnlarx` object. When used in a Hammerstein-Wiener model,  $m=q=1$  and  $Q$  is a scalar.

- $r$  - Mean value of the regressor vector computed from estimation data, specified as a 1-by- $m$  vector.
- $a_s, b_s, a_w,$  and  $b_w$  — Scaling and wavelet parameters, specified as scalars. Parameters with the  $s$  subscript are scaling parameters, and parameters with the  $w$  subscript are wavelet parameters.
- $L$  — Specified as a  $p$ -by-1 vector.
- $c_s$  and  $c_w$  - Specified as a 1-by- $q$  vectors.
- $d$  — Output offset, specified as a scalar.

The value  $F(x)$  is computed by `evaluate(NL, x)`, where `NL` is the `wavenet` object.

For `wavenet` object properties, see “Properties” on page 1-1872.

## Examples

### Create a Default Wavelet Nonlinearity Estimator

```
NL = wavenet;
```

Exclude the linear term from the wavelet expansion.

```
NL.LinearTerm = 'off';
```

### Estimate Nonlinear ARX Model with Specific Nonlinearity

Load the estimation data.

```
load twotankdata;
```

Create an `iddata` object from the estimation data.

```
z = iddata(y,u,0.2);
```

Create a wavelet network nonlinearity estimator with 5 units.

```
NL = wavenet('NumberOfUnits',5);
```

Estimate the nonlinear ARX model.

```
sys = nlarx(z,[4 4 1],NL);
```

### Estimate MIMO Nonlinear ARX Model

Load the estimation data.

```
load motorizedcamera;
```

Create an `iddata` object.

```
z = iddata(y,u,0.02,'Name','Motorized Camera','TimeUnit','s');
```

`z` is an `iddata` object with 6 inputs and 2 outputs.

Specify the model orders.

```
Orders = [ones(2,2),2*ones(2,6),ones(2,6)];
```

Specify different nonlinearity estimators for each output channel.

```
NL = [wavenet('NumberOfUnits',2),linear];
```

Estimate the nonlinear ARX model.

```
sys = nlarx(z,Orders,NL);
```

### **Estimate MIMO Hammerstein-Wiener Model**

Load the estimation data.

```
load motorizedcamera;
```

Create an `iddata` object.

```
z = iddata(y,u,0.02,'Name','Motorized Camera','TimeUnit','s');
```

`z` is an `iddata` object with 6 inputs and 2 outputs.

Specify the model orders and delays.

```
Orders = [ones(2,6),ones(2,6),ones(2,6)];
```

Specify the same nonlinearity estimator for each input channel.

```
InputNL = saturation;
```

Specify different nonlinearity estimators for each output channel.

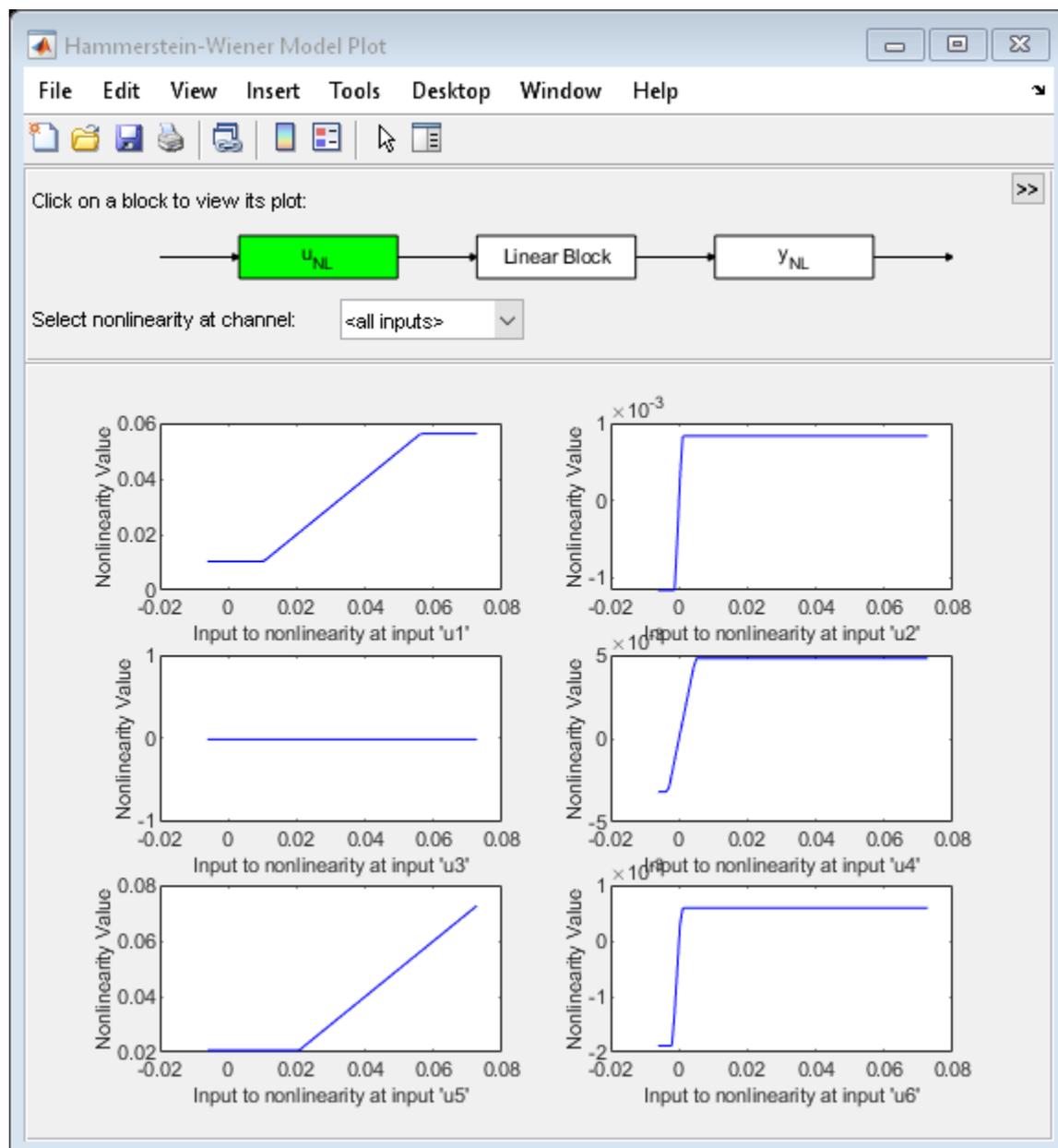
```
OutputNL = [deadzone,wavenet];
```

Estimate the Hammerstein-Wiener model.

```
sys = nlhw(z,Orders,InputNL,OutputNL);
```

To see the shape of the estimated input and output nonlinearities, plot the nonlinearities.

```
plot(sys)
```



Click on the input and output nonlinearity blocks on the top of the plot to see the nonlinearities.

## Input Arguments

### Name-Value Pair Arguments

Specify optional comma-separated pairs of `Name`, `Value` arguments. `Name` is the argument name and `Value` is the corresponding value. `Name` must appear inside quotes. You can specify several name and value pair arguments in any order as `Name1, Value1, ..., NameN, ValueN`.

Use `Name`, `Value` arguments to specify additional properties on page 1-662 of `wavenet` nonlinearity. For example, `NL= wavenet('NumberOfUnits',5)` creates a wavelet nonlinearity estimator object with five nonlinearity units in wavelet expansion.

## Properties

### NumberOfUnits

Number of nonlinearity units in wavelet expansion, specified as a positive integer or one of the following values:

- `'Auto'` — The number of units are automatically determined from estimation data.
- `'Interactive'` — The number of units are determined during model estimation.

**Default:** `'Auto'`

### LinearTerm

Inclusion of linear term, specified as one of the following values:

- `'on'` — The linear term,  $(x-r)PL$ , is included in the wavelet expansion.
- `'off'` — The linear term is not included in the wavelet expansion.

**Default:** `'on'`

## Parameters

Parameters in wavelet expansion, specified as a structure with the following fields:

Field Name	Description	Default
RegressorMean	Means of the regressors, $r$ , specified as a 1-by- $m$ vector. $m$ is the number of regressors, $x$ .  For Hammerstein-Wiener models, $m = 1$ .	[]
NonLinearSubspace	Projection matrix, $Q$ , specified as an $m$ -by- $q$ matrix.	[]
LinearSubspace	Projection matrix, $P$ , specified as an $m$ -by- $p$ matrix.	[]
LinearCoef	Linear coefficients, $L$ , specified as a $p$ -by-1 vector.	[]
ScalingDilation	Scaling function dilation, $b_{s\_ns}$ , specified as an $ns$ -by-1 matrix.	[]
WaveletDilation	Wavelet function dilation, $b_{w\_nw}$ , specified as an $nw$ -by-1 matrix.	[]
ScalingTranslation	Scaling function translation, $c_{s\_ns}$ , specified as an $ns$ -by- $q$ matrix.	[]
WaveletTranslation	Wavelet function translation, $c_{w\_nw}$ , specified as an $nw$ -by- $q$ matrix.	[]
ScalingCoef	Scaling function coefficients, $a_{s\_ns}$ , specified as an $ns$ -by-1 vector.	[]
WaveletCoef	Wavelet function coefficients, $a_{w\_nw}$ , specified as an $nw$ -by-1 vector.	[]
OutputOffset	Output offset, $d$ , specified as a scalar.	[]

The parameters are typically not assigned directly. They are estimated by the identification algorithm (`nlarx` or `nllhw`) when `wavenet` is used in a Nonlinear ARX (`idnlarx`) or Hammerstein-Wiener (`idnllhw`) model.

## Options

Options specifying the initial wavelet nonlinearity structure, specified as a structure with the following fields:

<b>Field Name</b>	<b>Description</b>	<b>Default</b>
FinestCell	Minimum number of data points in the smallest cell, specified as an integer or character vector. A cell is the area covered by the significantly nonzero portion of a wavelet.	'auto' — Compute the value from the data.
MinCells	Minimum number of cells in the partition, specified as an integer.	16
MaxCells	Maximum number of cells in the partition, specified as an integer.	128
MaxLevels	Maximum number of wavelet levels, specified as an integer.	10
DilationStep	Dilation step size, specified as real scalar.	2
TranslationStep	Translation step size, specified as real scalar.	1

## Output Arguments

### **NL — Wavelet nonlinearity estimator object**

wavenet object

Wavelet nonlinearity estimator object, returned as a wavenet object.

## Algorithms

wavenet can be used in both Nonlinear ARX and Hammerstein-Wiener models.

- When used in a Nonlinear ARX model:
  - If the Focus estimation option (see, nlarxOptions) is 'prediction', wavenet uses a fast, noniterative technique for estimating parameters [1]. Successive refinements after the first estimation use an iterative algorithm.
  - If Focus is 'simulation', wavenet uses an iterative technique for estimating parameters.

To always use a noniterative or iterative algorithm, specify the `IterativeWavenet` option of `nlarxOptions`.

- When used in a Hammerstein-Wiener model, `wavenet` parameters are determined by iterative minimization.

## References

- [1] Zhang, Q. "Using wavelet network in nonparametric estimation." *IEEE Trans. on Neural Networks*, Vol. 8, Number 2, March 1997, pp. 227-236.

## See Also

`nlarx` | `nllhw`

**Introduced in R2007a**

## xperm

Reorder states in state-space models

## Syntax

```
sys = xperm(sys,P)
```

## Description

`sys = xperm(sys,P)` reorders the states of the state-space model `sys` according to the permutation `P`. The vector `P` is a permutation of `1:NX`, where `NX` is the number of states in `sys`. For information about creating state-space models, see `ss` and `dss`.

## Examples

### Alphabetically Order States of State-Space Model

Load a previously saved state space model `ssF8` with four states.

```
load ltiexamples  
ssF8
```

```
ssF8 =
```

```
A =  
      PitchRate  Velocity  AOA  PitchAngle  
PitchRate      -0.7    -0.0458  -12.2      0  
Velocity         0     -0.014   -0.2904  -0.562  
AOA              1     -0.0057   -1.4      0  
PitchAngle       1         0         0      0
```

```
B =  
      Elevator  Flaperon  
PitchRate     -19.1   -3.1  
Velocity      -0.0119 -0.0096  
AOA           -0.14   -0.72
```

```

PitchAngle      0      0

C =
      PitchRate  Velocity      AOA  PitchAngle
FlightPath      0      0      -1      1
Acceleration    0      0      0.733    0

D =
      Elevator  Flaperon
FlightPath      0      0
Acceleration    0.0768  0.1134

```

Continuous-time state-space model.

Order the states in alphabetical order.

```

[y,P] = sort(ssF8.StateName);
sys = xperm(ssF8,P)

```

```

sys =

A =
      AOA  PitchAngle  PitchRate  Velocity
AOA      -1.4      0      1      -0.0057
PitchAngle  0      0      1      0
PitchRate  -12.2      0      -0.7      -0.0458
Velocity   -0.2904      -0.562      0      -0.014

B =
      Elevator  Flaperon
AOA      -0.14      -0.72
PitchAngle  0      0
PitchRate  -19.1      -3.1
Velocity   -0.0119      -0.0096

C =
      AOA  PitchAngle  PitchRate  Velocity
FlightPath  -1      1      0      0
Acceleration  0.733      0      0      0

D =
      Elevator  Flaperon
FlightPath      0      0
Acceleration    0.0768  0.1134

```

Continuous-time state-space model.

The states in `ssF8` now appear in alphabetical order.

### **See Also**

`dss` | `ss`

**Introduced in R2008b**

## zero

Zeros and gain of SISO dynamic system

### Syntax

```
Z = zero(sys)
[Z,gain] = zero(sys)
[Z,gain] = zero(sys,J1,...,JN)
```

### Description

`Z = zero(sys)` returns the zeros of the single-input, single-output (SISO) dynamic system model, `sys`. The output is expressed as the reciprocal of the time units specified in `sys.TimeUnit`.

`[Z,gain] = zero(sys)` also returns the zero-pole-gain of `sys`.

`[Z,gain] = zero(sys,J1,...,JN)` returns the zeros and gain of the entries in the model array `sys` with subscripts `J1,...,JN`.

### Examples

#### Zeros of Transfer Function

Compute the zeros of the following transfer function:

$$\text{sys}(s) = \frac{4.2s^2 + 0.25s - 0.004}{s^2 + 9.6s + 17}$$

```
sys = tf([4.2,0.25,-0.004],[1,9.6,17]);
Z = zero(sys)
```

```
Z = 2×1
```

```
-0.0726  
0.0131
```

## Zeros and Gain of Transfer Function

Calculate the zero locations and zero-pole gain of the following transfer function:

$$\text{sys}(s) = \frac{4.2s^2 + 0.25s - 0.004}{s^2 + 9.6s + 17}$$

```
sys = tf([4.2,0.25,-0.004],[1,9.6,17]);  
[z,gain] = zero(sys)
```

```
z = 2×1
```

```
-0.0726  
0.0131
```

```
gain = 4.2000
```

The zero locations are expressed in  $\text{second}^{-1}$ , because the time unit of the transfer function (`H.TimeUnit`) is seconds.

## Zeros and Gain of Models in an Array

For this example, load a 3-by-1 array of transfer function models.

```
load('tfArray.mat','sys');  
size(sys)
```

```
3×1 array of transfer functions.  
Each model has 1 outputs and 1 inputs.
```

Find the zeros and gain values of the models in the array.

```
[Z, gain] = zero(sys);  
Z(:, :, 1, 1)
```

```
ans = 2×1
    -0.0726
     0.0131
```

```
gain(:, :, 1, 1)
ans = 4.2000
```

`zero` returns an array each for the zeros and the gain values respectively. Here, `Z(:, :, 1, 1)` and `gain(:, :, 1, 1)` corresponds to the zero and the gain value of the first model in the array, that is, `sys(:, :, 1, 1)`.

## Input Arguments

### **sys** — Dynamic system

dynamic system model | model array

Dynamic system, specified as a SISO dynamic system model, or an array of SISO dynamic system models. Dynamic systems that you can use include continuous-time or discrete-time numeric LTI models such as `tf`, `zpk`, or `ss` models.

If `sys` is a generalized state-space model `genss` or an uncertain state-space model `uss`, `zero` returns the zeros of the current or nominal value of `sys`. If `sys` is an array of models, `zero` returns the zeros of the model corresponding to its subscript `J1, . . . , JN` in `sys`. For more information on model arrays, see “Model Arrays” (Control System Toolbox).

### **J1, . . . , JN** — Indices of models in array whose zeros you want to extract

positive integer

Indices of models in array whose zeros you want to extract, specified as a positive integer. You can provide as many indices as there are array dimensions in `sys`. For example, if `sys` is a 4-by-5 array of dynamic system models, the following command extracts the zeros for entry (2,3) in the array.

```
Z = zero(sys,2,3);
```

## Output Arguments

### **Z — Zeros of the dynamic system**

column vector | array

Zeros of the dynamic system, returned as a column vector or an array. If `sys` is:

- A single model, then `Z` is a column vector of zeros of the dynamic system model `sys`
- A model array, then `Z` is an array containing the zeros of each model in `sys`

`Z` is expressed as the reciprocal of the time units specified in `sys.TimeUnit`. For example, zero is expressed in 1/minute if `sys.TimeUnit = 'minutes'`.

### **gain — Zero-pole-gain of the dynamic system**

scalar

Zero-pole-gain of the dynamic system, returned as a scalar. In other words, `gain` is the value of `K` when the model is written in `zpk` form.

## Tips

- If `sys` has internal delays, `zero` sets all internal delays to zero, creating a zero-order Padé approximation. This approximation ensures that the system has a finite number of zeros. `zero` returns an error if setting internal delays to zero creates singular algebraic loops. To assess the stability of models with internal delays, use `step` or `impulse`.
- To calculate the transmission zeros of a multi-input, multi-output (MIMO) system, use `tzero`.

## See Also

`impulse` | `pole` | `pzmap` | `pzplot` | `step` | `tzero`

## Topics

“Pole and Zero Locations” (Control System Toolbox)

**Introduced in R2012a**

## zgrid

Generate z-plane grid of constant damping factors and natural frequencies

### Syntax

```
zgrid  
zgrid(z,wn)  
zgrid([],[])
```

### Description

`zgrid` generates, for root locus and pole-zero maps, a grid of constant damping factors from zero to one in steps of 0.1 and natural frequencies from zero to  $\pi$  in steps of  $\pi/10$ , and plots the grid over the current axis. If the current axis contains a discrete z-plane root locus diagram or pole-zero map, `zgrid` draws the grid over the plot without altering the current axis limits.

`zgrid(z,wn)` plots a grid of constant damping factor and natural frequency lines for the damping factors and normalized natural frequencies in the vectors `z` and `wn`, respectively. If the current axis contains a discrete z-plane root locus diagram or pole-zero map, `zgrid(z,wn)` draws the grid over the plot. The frequency lines for unnormalized (true) frequencies can be plotted using

```
zgrid(z,wn/Ts)
```

where `Ts` is the sample time.

`zgrid([],[])` draws the unit circle.

Alternatively, you can select **Grid** from the right-click menu to generate the same z-plane grid.

### Examples

## Plot z-plane grid lines on the root locus

To see the z-plane grid on the root locus plot, type

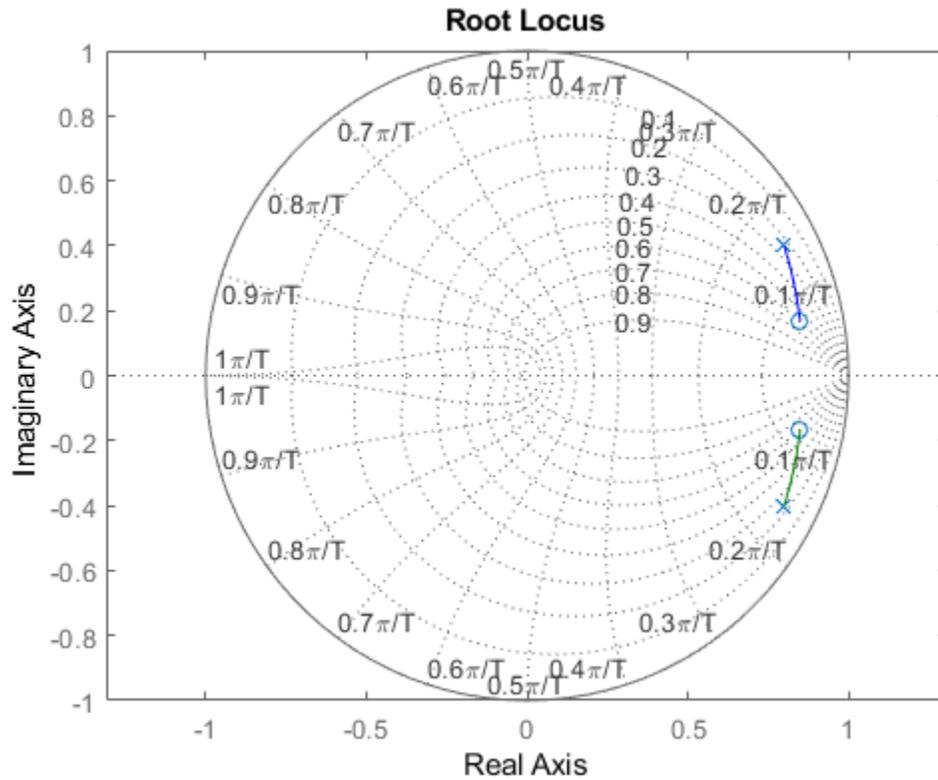
```
H = tf([2 -3.4 1.5],[1 -1.6 0.8],-1)
```

```
H =
```

$$\frac{2 z^2 - 3.4 z + 1.5}{z^2 - 1.6 z + 0.8}$$

```
Sample time: unspecified  
Discrete-time transfer function.
```

```
rlocus(H)  
zgrid  
axis equal
```



## See Also

[pzmap](#) | [rlocus](#) | [sgrid](#)

Introduced in R2012a

## zpkdata

Access zero-pole-gain data

### Syntax

```
[z,p,k] = zpkdata(sys)
[z,p,k,Ts] = zpkdata(sys)
[z,p,k,Ts,covz,covp,covk] = zpkdata(sys)
```

### Description

`[z,p,k] = zpkdata(sys)` returns the zeros  $z$ , poles  $p$ , and gain(s)  $k$  of the zero-pole-gain model `sys`. The outputs  $z$  and  $p$  are cell arrays with the following characteristics:

- $z$  and  $p$  have as many rows as outputs and as many columns as inputs.
- The  $(i,j)$  entries  $z\{i,j\}$  and  $p\{i,j\}$  are the (column) vectors of zeros and poles of the transfer function from input  $j$  to output  $i$ .

The output  $k$  is a matrix with as many rows as outputs and as many columns as inputs such that  $k(i,j)$  is the gain of the transfer function from input  $j$  to output  $i$ . If `sys` is a transfer function or state-space model, it is first converted to zero-pole-gain form using `zpk`.

For SISO zero-pole-gain models, the syntax

```
[z,p,k] = zpkdata(sys,'v')
```

forces `zpkdata` to return the zeros and poles directly as column vectors rather than as cell arrays (see example below).

`[z,p,k,Ts] = zpkdata(sys)` also returns the sample time  $T_s$ .

`[z,p,k,Ts,covz,covp,covk] = zpkdata(sys)` also returns the covariances of the zeros, poles and gain of the identified model `sys`. `covz` is a cell array such that `covz{ky,ku}` contains the covariance information about the zeros in the vector  $z\{ky,ku\}$ . `covz{ky,ku}` is a 3-D array of dimension 2-by-2-by- $N_z$ , where  $N_z$  is the

length of  $z\{k_y, k_u\}$ , so that the (1,1) element is the variance of the real part, the (2,2) element is the variance of the imaginary part, and the (1,2) and (2,1) elements contain the covariance between the real and imaginary parts.  $\text{covp}$  has a similar relationship to  $\text{p}$ .  $\text{covk}$  is a matrix containing the variances of the elements of  $k$ .

You can access the remaining LTI properties of `sys` with `get` or by direct referencing, for example,

```
sys.Ts
sys.inputname
```

## Examples

### Example 1

Given a zero-pole-gain model with two outputs and one input

```
H = zpk([0];[-0.5]},{[0.3];[0.1+i 0.1-i]},{[1;2],-1)
Zero/pole/gain from input to output...
```

```

      z
#1:  -----
      (z-0.3)

      2 (z+0.5)
#2:  -----
      (z^2 - 0.2z + 1.01)
```

Sample time: unspecified

you can extract the zero/pole/gain data embedded in `H` with

```
[z,p,k] = zpkdata(H)
z =
     [      0]
     [-0.5000]
p =
     [  0.3000]
     [2x1 double]
k =
     1
     2
```

To access the zeros and poles of the second output channel of H, get the content of the second cell in z and p by typing

```
z{2,1}
ans =
    -0.5000
p{2,1}
ans =
    0.1000+ 1.0000i
    0.1000- 1.0000i
```

## Example 2

Extract the ZPK matrices and their standard deviations for a 2-input, 1 output identified transfer function.

```
load iddata7
```

```
transfer function model
```

```
sys1 = tfest(z7, 2, 1, 'InputDelay',[1 0]);
```

```
an equivalent process model
```

```
sys2 = procest(z7, {'P2UZ', 'P2UZ'}, 'InputDelay',[1 0]);
```

```
l, p1, k1, ~, dz1, dp1, dk1] = zpkdata(sys1);
[z2, p2, k2, ~, dz2, dp2, dk2] = zpkdata(sys2);
```

Use `iopzplot` to visualize the pole-zero locations and their covariances

```
h = iopzplot(sys1, sys2);
showConfidence(h)
```

## See Also

[get](#) | [ssdata](#) | [tfdata](#) | [zpk](#)

**Introduced before R2006a**

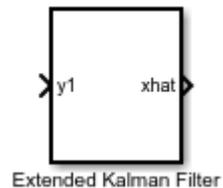
# Blocks — Alphabetical List

---

## Extended Kalman Filter

Estimate states of discrete-time nonlinear system using extended Kalman filter

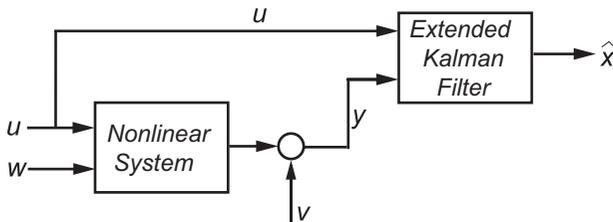
**Library:** Control System Toolbox / State Estimation  
System Identification Toolbox / Estimators



### Description

The Extended Kalman Filter block estimates the states of a discrete-time nonlinear system using the first-order discrete-time extended Kalman filter algorithm.

Consider a plant with states  $x$ , input  $u$ , output  $y$ , process noise  $w$ , and measurement noise  $v$ . Assume that you can represent the plant as a nonlinear system.



Using the state transition and measurement functions of the system and the extended Kalman filter algorithm, the block produces state estimates  $\hat{x}$  for the current time step. For information about the algorithm, see “Extended and Unscented Kalman Filter Algorithms for Online State Estimation”.

You create the nonlinear state transition function and measurement functions for the system and specify these functions in the block. The block supports state estimation of a system with multiple sensors that are operating at different sampling rates. You can specify up to five measurement functions, each corresponding to a sensor in the system. You can also specify the Jacobians of the state transition and measurement functions. If you do not specify them, the software numerically computes the Jacobians. For more information, see “State Transition and Measurement Functions” on page 2-19.

## Ports

### Input

#### **y1, y2, y3, y4, y5 — Measured system outputs**

vector

Measured system outputs corresponding to each measurement function that you specify in the block. The number of ports equals the number of measurement functions in your system. You can specify up to five measurement functions. For example, if your system has two sensors, you specify two measurement functions in the block. The first port **y1** is available by default. When you click **Apply**, the software generates port **y2** corresponding to the second measurement function.

Specify the ports as  $N$ -dimensional vectors, where  $N$  is the number of quantities measured by the corresponding sensor. For example, if your system has one sensor that measures the position and velocity of an object, then there is only one port **y1**. The port is specified as a 2-dimensional vector with values corresponding to position and velocity.

#### **Dependencies**

The first port **y1** is available by default. Ports **y2** to **y5** are generated when you click **Add Measurement**, and click **Apply**.

Data Types: `single` | `double`

#### **StateTransitionFcnInputs — Additional optional input argument to state transition function**

scalar | vector | matrix

Additional optional input argument to the state transition function  $f$  other than the state  $x$  and process noise  $w$ . For information about state transition functions see, “State Transition and Measurement Functions” on page 2-19.

Suppose that your system has nonadditive process noise, and the state transition function  $f$  has the following form:

$$x(k+1) = f(x(k), w(k), \text{StateTransitionFcnInputs})$$

Here  $k$  is the time step, and `StateTransitionFcnInputs` is an additional input argument other than  $x$  and  $w$ .

If you create  $f$  using a MATLAB function (.m file), the software generates the port **StateTransitionFcnInputs** when you click **Apply**. You can specify the inputs to this port as a scalar, vector, or matrix.

If your state transition function has more than one additional input, use a Simulink Function block to specify the function. When you use a Simulink Function block, you provide the additional inputs directly to the Simulink Function block using Inport blocks. No input ports are generated for the additional inputs in the Extended Kalman Filter block.

### Dependencies

This port is generated only if both of the following conditions are satisfied:

- You specify  $f$  in **Function** using a MATLAB function, and  $f$  is on the MATLAB path.
- $f$  requires only one additional input argument apart from  $x$  and  $w$ .

Data Types: `single` | `double`

### **MeasurementFcn1Inputs, MeasurementFcn2Inputs, MeasurementFcn3Inputs, MeasurementFcn4Inputs, MeasurementFcn5Inputs** — Additional optional input argument to each measurement function

`scalar` | `vector` | `matrix`

Additional optional inputs to the measurement functions other than the state  $x$  and measurement noise  $v$ . For information about measurement functions see, “State Transition and Measurement Functions” on page 2-19.

**MeasurementFcn1Inputs** corresponds to the first measurement function that you specify, and so on. For example, suppose that your system has three sensors and nonadditive measurement noise, and the three measurement functions  $h_1$ ,  $h_2$ , and  $h_3$  have the following form:

$$y_1[k] = h_1(x[k], v_1[k], \text{MeasurementFcn1Inputs})$$

$$y_2[k] = h_2(x[k], v_2[k], \text{MeasurementFcn2Inputs})$$

$$y_3[k] = h_3(x[k], v_3[k])$$

Here  $k$  is the time step, and **MeasurementFcn1Inputs** and **MeasurementFcn2Inputs** are the additional input arguments to  $h_1$  and  $h_2$ .

If you specify  $h_1$ ,  $h_2$ , and  $h_3$  using MATLAB functions (.m files) in **Function**, the software generates ports **MeasurementFcn1Inputs** and **MeasurementFcn2Inputs** when you click **Apply**. You can specify the inputs to these ports as scalars, vectors, or matrices.

If your measurement functions have more than one additional input, use Simulink Function blocks to specify the functions. When you use a Simulink Function block, you provide the additional inputs directly to the Simulink Function block using Inport blocks. No input ports are generated for the additional inputs in the Extended Kalman Filter block.

### Dependencies

A port corresponding to a measurement function  $h$  is generated only if both of the following conditions are satisfied:

- You specify  $h$  in **Function** using a MATLAB function, and  $h$  is on the MATLAB path.
- $h$  requires only one additional input argument apart from  $x$  and  $v$ .

Data Types: `single` | `double`

### Q — Time-varying process noise covariance

scalar | vector | matrix

Time-varying process noise covariance, specified as a scalar, vector, or matrix depending on the value of the **Process noise** parameter:

- **Process noise** is **Additive** — Specify the covariance as a scalar, an  $N_s$ -element vector, or an  $N_s$ -by- $N_s$  matrix, where  $N_s$  is the number of states of the system. Specify a scalar if there is no cross-correlation between process noise terms, and all the terms have the same variance. Specify a vector of length  $N_s$ , if there is no cross-correlation between process noise terms, but all the terms have different variances.
- **Process noise** is **Nonadditive** — Specify the covariance as a  $W$ -by- $W$  matrix, where  $W$  is the number of process noise terms in the state transition function.

### Dependencies

This port is generated if you specify the process noise covariance as **Time-Varying**. The port appears when you click **Apply**.

Data Types: `single` | `double`

### R1, R2, R3, R4, R5 — Time-varying measurement noise covariance

matrix

Time-varying measurement noise covariances for up to five measurement functions of the system, specified as matrices. The sizes of the matrices depend on the value of the **Measurement noise** parameter for the corresponding measurement function:

- **Measurement noise** is **Additive** — Specify the covariance as an  $N$ -by- $N$  matrix, where  $N$  is the number of measurements of the system.
- **Measurement noise** is **Nonadditive** — Specify the covariance as a  $V$ -by- $V$  matrix, where  $V$  is the number of measurement noise terms in the corresponding measurement function.

### Dependencies

A port is generated if you specify the measurement noise covariance as **Time-Varying** for the corresponding measurement function. The port appears when you click **Apply**.

Data Types: `single` | `double`

### **Enable1, Enable2, Enable3, Enable4, Enable5** — Enable correction of estimated states when measured data is available

scalar

Suppose that measured output data is not available at all time points at the port **y1** that corresponds to the first measurement function. Use a signal value other than  $\emptyset$  at the **Enable1** port to enable the correction of estimated states when measured data is available. Specify the port value as  $\emptyset$  when measured data is not available. Similarly, if measured output data is not available at all time points at the port **y $i$**  for the  $i^{\text{th}}$  measurement function, specify the corresponding port **Enable $i$**  as a value other than  $\emptyset$ .

### Dependencies

A port corresponding to a measurement function is generated if you select **Add Enable port** for that measurement function. The port appears when you click **Apply**.

Data Types: `single` | `double` | `Boolean`

## Output

### **xhat** — Estimated states

vector

Estimated states, returned as a vector of size  $N_s$ , where  $N_s$  is the number of states of the system. To access the individual states, use the Selector block.

When the **Use the current measurements to improve state estimates** parameter is selected, the block outputs the corrected state estimate  $\hat{x}[k|k]$  at time step  $k$ , estimated using measured outputs until time  $k$ . If you clear this parameter, the block returns the predicted state estimate  $\hat{x}[k|k-1]$  for time  $k$ , estimated using measured output until a previous time  $k-1$ . Clear this parameter if your filter is in a feedback loop and there is an algebraic loop in your Simulink model.

Data Types: `single` | `double`

### **P — State estimation error covariance**

matrix

State estimation error covariance, returned as an  $N_s$ -by- $N_s$  matrix, where  $N_s$  is the number of states of the system. To access the individual covariances, use the Selector block.

#### **Dependencies**

This port is generated if you select **Output state estimation covariance** in the **System Model** tab, and click **Apply**.

Data Types: `single` | `double`

## **Parameters**

### **System Model Tab**

#### **State Transition**

##### **Function — State transition function name**

`myStateTransitionFcn` (default) | function name

The state transition function calculates the  $N_s$ -element state vector of the system at time step  $k+1$ , given the state vector at time step  $k$ .  $N_s$  is the number of states of the nonlinear system. You create the state transition function and specify the function name in **Function**. For example, if `vdpStateFcn.m` is the state transition function that you created and saved, specify **Function** as `vdpStateFcn`.

The inputs to the function you create depend on whether you specify the process noise as additive or nonadditive in **Process noise**.

- **Process noise is Additive** — The state transition function  $f$  specifies how the states evolve as a function of state values at previous time step:

$$x(k+1) = f(x(k), Us1(k), \dots, Usn(k)),$$

where  $x(k)$  is the estimated state at time  $k$ , and  $Us1, \dots, Usn$  are any additional input arguments required by your state transition function, such as system inputs or the sample time. To see an example of a state transition function with additive process noise, type `edit vdpStateFcn` at the command line.

- **Process noise is Nonadditive** — The state transition function also specifies how the states evolve as a function of the process noise  $w$ :

$$x(k+1) = f(x(k), w(k), Us1(k), \dots, Usn(k)).$$

For more information, see “State Transition and Measurement Functions” on page 2-19.

You can create  $f$  using a Simulink Function block or as a MATLAB function (.m file).

- You can use a MATLAB function only if  $f$  has one additional input argument  $Us1$  other than  $x$  and  $w$ .

$$x(k+1) = f(x(k), w(k), Us1(k))$$

The software generates an additional input port **StateTransitionFcnInputs** to specify this argument.

- If you are using a Simulink Function block, specify  $x$  and  $w$  using Argument Inport blocks and the additional inputs  $Us1, \dots, Usn$  using Inport blocks in the Simulink Function block. You do not provide  $Us1, \dots, Usn$  to the Extended Kalman Filter block.

### Programmatic Use

**Block Parameter:** StateTransitionFcn

**Type:** character vector, string

**Default:** 'myStateTransitionFcn'

### Jacobian — Jacobian of state transition function

off (default) | on

Jacobian of state transition function  $f$ , specified as one of the following:

- **off** — The software computes the Jacobian numerically. This computation may increase processing time and numerical inaccuracy of the state estimation.

- **on** — You create a function to compute the Jacobian, and specify the name of the function in **Jacobian**. For example, if `vdpStateJacobianFcn.m` is the Jacobian function, specify **Jacobian** as `vdpStateJacobianFcn`. If you create the state transition function  $f$  using a Simulink Function block, then create the Jacobian using a Simulink Function block. If you create  $f$  using a MATLAB function (`.m` file), then create the Jacobian using a MATLAB function.

The function calculates the partial derivatives of the state transition function with respect to the states and process noise. The number of inputs to the Jacobian function must equal the number of inputs of the state transition function and must be specified in the same order in both functions. The number of outputs of the Jacobian function depends on the **Process noise** parameter:

- **Process noise** is **Additive** — The function calculates the partial derivative of the state transition function  $f$  with respect to the states ( $\partial f / \partial x$ ). The output is an  $N_s$ -by- $N_s$  Jacobian matrix, where  $N_s$  is the number of states.

To see an example of a Jacobian function for additive process noise, type `edit vdpStateJacobianFcn` at the command line.

- **Process noise** is **Nonadditive** — The function must also return a second output that is the partial derivative of the state transition function  $f$  with respect to the process noise terms ( $\partial f / \partial w$ ). The second output is returned as an  $N_s$ -by- $W$  matrix, where  $W$  is the number of process noise terms in the state transition function.

### Programmatic Use

**Block Parameter:** `HasStateTransitionJacobianFcn`

**Type:** character vector

**Values:** `'off'`, `'on'`

**Default:** `'off'`

**Block Parameter:** `StateTransitionJacobianFcn`

**Type:** character vector, string

**Default:** `''`

### Process noise — Process noise characteristics

Additive (default) | Nonadditive

Process noise characteristics, specified as one of the following values:

- **Additive** — Process noise  $w$  is additive, and the state transition function  $f$  that you specify in **Function** has the following form:

$$x(k+1) = f(x(k), Us1(k), \dots, Usn(k)),$$

where  $x(k)$  is the estimated state at time  $k$ , and  $Us1, \dots, Usn$  are any additional input arguments required by your state transition function.

- **Nonadditive** — Process noise is nonadditive, and the state transition function specifies how the states evolve as a function of the state *and* process noise at the previous time step:

$$x(k+1) = f(x(k), w(k), Us1(k), \dots, Usn(k)).$$

### Programmatic Use

**Block Parameter:** HasAdditiveProcessNoise

**Type:** character vector

**Values:** 'Additive', 'Nonadditive'

**Default:** 'Additive'

### Covariance — Time-invariant process noise covariance

1 (default) | scalar | vector | matrix

Time-invariant process noise covariance, specified as a scalar, vector, or matrix depending on the value of the **Process noise** parameter:

- **Process noise** is **Additive** — Specify the covariance as a scalar, an  $Ns$ -element vector, or an  $Ns$ -by- $Ns$  matrix, where  $Ns$  is the number of states of the system. Specify a scalar if there is no cross-correlation between process noise terms and all the terms have the same variance. Specify a vector of length  $Ns$ , if there is no cross-correlation between process noise terms but all the terms have different variances.
- **Process noise** is **Nonadditive** — Specify the covariance as a  $W$ -by- $W$  matrix, where  $W$  is the number of process noise terms.

If the process noise covariance is time-varying, select **Time-varying**. The block generates input port **Q** to specify the time-varying covariance.

### Dependencies

This parameter is enabled if you do not specify the process noise as **Time-Varying**.

### Programmatic Use

**Block Parameter:** ProcessNoise

**Type:** character vector, string

**Default:** '1'

**Time-varying — Time-varying process noise covariance**

'off' (default) | 'on'

If you select this parameter, the block includes an additional input port **Q** to specify the time-varying process noise covariance.

**Programmatic Use****Block Parameter:** HasTimeVaryingProcessNoise**Type:** character vector**Values:** 'off', 'on'**Default:** 'off'**Initialization****Initial state — Initial state estimate** $\theta$  (default) | vector

Initial state estimate value, specified as an  $N_s$ -element vector, where  $N_s$  is the number of states in the system. Specify the initial state values based on your knowledge of the system.

**Programmatic Use****Block Parameter:** InitialState**Type:** character vector, string**Default:** ' $\theta$ '**Initial covariance — State estimation error covariance**

1 (default) | scalar | vector | matrix

State estimation error covariance, specified as a scalar, an  $N_s$ -element vector, or an  $N_s$ -by- $N_s$  matrix, where  $N_s$  is the number of states of the system. If you specify a scalar or vector, the software creates an  $N_s$ -by- $N_s$  diagonal matrix with the scalar or vector elements on the diagonal.

Specify a high value for the covariance when you do not have confidence in the initial state values that you specify in **Initial state**.

**Programmatic Use****Block Parameter:** InitialStateCovariance**Type:** character vector, string**Default:** '1'

## Measurement

### Function — Measurement function name

myMeasurementFcn (default) | function name

The measurement function calculates the  $N$ -element output measurement vector of the nonlinear system at time step  $k$ , given the state vector at time step  $k$ . You create the measurement function and specify the function name in **Function**. For example, if `vdpMeasurementFcn.m` is the measurement function that you created and saved, specify **Function** as `vdpMeasurementFcn`.

The inputs to the function you create depend on whether you specify the measurement noise as additive or nonadditive in **Measurement noise**.

- **Measurement noise** is Additive — The measurement function  $h$  specifies how the measurements evolve as a function of state Values:

$$y(k) = h(x(k), Um1(k), \dots, Umn(k)),$$

where  $y(k)$  and  $x(k)$  are the estimated output and estimated state at time  $k$ , and  $Um1, \dots, Umn$  are any optional input arguments required by your measurement function. For example, if you are using a sensor for tracking an object, an additional input could be the sensor position.

To see an example of a measurement function with additive process noise, type `edit vdpMeasurementFcn` at the command line.

- **Measurement noise** is Nonadditive— The measurement function also specifies how the output measurement evolves as a function of the measurement noise  $v$ :

$$y(k) = h(x(k), v(k), Um1(k), \dots, Umn(k)).$$

To see an example of a measurement function with nonadditive process noise, type `edit vdpMeasurementNonAdditiveNoiseFcn`.

For more information, see “State Transition and Measurement Functions” on page 2-19.

You can create  $h$  using a Simulink Function block or as a MATLAB function (.m file).

- You can use a MATLAB function only if  $h$  has one additional input argument  $Um1$  other than  $x$  and  $v$ .

$$y[k] = h(x[k], v[k], Um1(k))$$

The software generates an additional input port **MeasurementFcnInputs** to specify this argument for the  $i$ th measurement function.

- If you are using a Simulink Function block, specify  $x$  and  $v$  using Argument Inport blocks and the additional inputs  $Um_1, \dots, Um_n$  using Inport blocks in the Simulink Function block. You do not provide  $Um_1, \dots, Um_n$  to the Extended Kalman Filter block.

If you have multiple sensors in your system, you can specify multiple measurement functions. You can specify up to five measurement functions using the **Add Measurement** button. To remove measurement functions, use **Remove Measurement**.

#### Programmatic Use

**Block Parameter:** MeasurementFcn1, MeasurementFcn2, MeasurementFcn3, MeasurementFcn4, MeasurementFcn5

**Type:** character vector, string

**Default:** 'myMeasurementFcn'

#### Jacobian — Jacobian of measurement function

off (default) | on

Jacobian of measurement function  $h$ , specified as one of the following:

- **off** — The software computes the Jacobian numerically. This computation may increase processing time and numerical inaccuracy of the state estimation.
- **on** — You create a function to compute the Jacobian of the measurement function  $h$ , and specify the name of the function in **Jacobian**. For example, if `vdpMeasurementJacobianFcn.m` is the Jacobian function, specify `MeasurementJacobianFcn` as `vdpMeasurementJacobianFcn`. If you create  $h$  using a Simulink Function block, then create the Jacobian using a Simulink Function block. If you create  $h$  using a MATLAB function (.m file), then create the Jacobian using a MATLAB function.

The function calculates the partial derivatives of the measurement function  $h$  with respect to the states and measurement noise. The number of inputs to the Jacobian function must equal the number of inputs to the measurement function and must be specified in the same order in both functions. The number of outputs of the Jacobian function depends on the **Measurement noise** parameter:

- **Measurement noise** is **Additive** — The function calculates the partial derivatives of the measurement function with respect to the states ( $\partial h / \partial x$ ). The

output is as an  $N$ -by- $N_s$  Jacobian matrix, where  $N$  is the number of measurements of the system and  $N_s$  is the number of states.

To see an example of a Jacobian function for additive measurement noise, type `edit vdpMeasurementJacobianFcn` at the command line.

- **Measurement noise** is **Nonadditive** — The function also returns a second output that is the partial derivative of the measurement function with respect to the measurement noise terms ( $\partial h / \partial v$ ). The second output is returned as an  $N$ -by- $V$  Jacobian matrix, where  $V$  is the number of measurement noise terms.

**Programmatic Use**

**Block Parameter:** HasMeasurementJacobianFcn1, HasMeasurementJacobianFcn2, HasMeasurementJacobianFcn3, HasMeasurementJacobianFcn4, HasMeasurementJacobianFcn5

**Type:** character vector

**Values:** 'off', 'on'

**Default:** 'off'

**Block Parameter:** MeasurementJacobianFcn1, MeasurementJacobianFcn2, MeasurementJacobianFcn3, MeasurementJacobianFcn4, MeasurementJacobianFcn5

**Type:** character vector

**Default:** ''

**Measurement noise — Measurement noise characteristics**

Additive (default) | Nonadditive

Measurement noise characteristics, specified as one of the following values:

- **Additive** — Measurement noise  $v$  is additive, and the measurement function  $h$  that you specify in **Function** has the following form:

$$y(k) = h(x(k), Um1(k), \dots, Umn(k)),$$

where  $y(k)$  and  $x(k)$  are the estimated output and estimated state at time  $k$ , and  $Um1, \dots, Umn$  are any optional input arguments required by your measurement function.

- **Nonadditive** — Measurement noise is nonadditive, and the measurement function specifies how the output measurement evolves as a function of the state *and* measurement noise:

$$y(k) = h(x(k), v(k), Um1(k), \dots, Umn(k)).$$

**Programmatic Use**

**Block Parameter:** HasAdditiveMeasurementNoise1, HasAdditiveMeasurementNoise2, HasAdditiveMeasurementNoise3, HasAdditiveMeasurementNoise4, HasAdditiveMeasurementNoise5

**Type:** character vector

**Values:** 'Additive', 'Nonadditive'

**Default:** 'Additive'

**Covariance — Time-invariant measurement noise covariance**

1 (default) | matrix

Time-invariant measurement noise covariance, specified as a matrix. The size of the matrix depends on the value of the **Measurement noise** parameter:

- **Measurement noise** is **Additive** — Specify the covariance as an  $N$ -by- $N$  matrix, where  $N$  is the number of measurements of the system.
- **Measurement noise** is **Nonadditive** — Specify the covariance as a  $V$ -by- $V$  matrix, where  $V$  is the number of measurement noise terms.

If the measurement noise covariance is time-varying, select **Time-varying**. The block generates input port **R $i$**  to specify the time-varying covariance for the  $i^{\text{th}}$  measurement function.

**Dependencies**

This parameter is enabled if you do not specify the process noise as **Time-Varying**.

**Programmatic Use**

**Block Parameter:** MeasurementNoise1, MeasurementNoise2, MeasurementNoise3, MeasurementNoise4, MeasurementNoise5

**Type:** character vector, string

**Default:** '1'

**Time-varying — Time-varying measurement noise covariance**

off (default) | on

If you select this parameter for the measurement noise covariance of the first measurement function, the block includes an additional input port **R1**. You specify the time-varying measurement noise covariance in **R1**. Similarly, if you select **Time-varying** for the  $i^{\text{th}}$  measurement function, the block includes an additional input port **R $i$**  to specify the time-varying measurement noise covariance for that function.

### Programmatic Use

**Block Parameter:** HasTimeVaryingMeasurementNoise1, HasTimeVaryingMeasurementNoise2, HasTimeVaryingMeasurementNoise3, HasTimeVaryingMeasurementNoise4, HasTimeVaryingMeasurementNoise5

**Type:** character vector

**Values:** 'off', 'on'

**Default:** 'off'

### Add Enable Port — Enable correction of estimated states only when measured data is available

off (default) | on

Suppose that measured output data is not available at all time points at the port **y1** that corresponds to the first measurement function. Select **Add Enable port** to generate an input port **Enable1**. Use a signal at this port to enable the correction of estimated states only when measured data is available. Similarly, if measured output data is not available at all time points at the port **y $i$**  for the  $i^{th}$  measurement function, select the corresponding **Add Enable port**.

### Programmatic Use

**Block Parameter:** HasMeasurementEnablePort1, HasMeasurementEnablePort2, HasMeasurementEnablePort3, HasMeasurementEnablePort4, HasMeasurementEnablePort5

**Type:** character vector

**Values:** 'off', 'on'

**Default:** 'off'

### Settings

#### Use the current measurements to improve state estimates — Choose between corrected or predicted state estimate

on (default) | off

When this parameter is selected, the block outputs the corrected state estimate  $\hat{x}[k|k]$  at time step  $k$ , estimated using measured outputs until time  $k$ . If you clear this parameter,

the block returns the predicted state estimate  $\hat{x}[k|k-1]$  for time  $k$ , estimated using measured output until a previous time  $k-1$ . Clear this parameter if your filter is in a feedback loop and there is an algebraic loop in your Simulink model.

### Programmatic Use

**Block Parameter:** UseCurrentEstimator

**Type:** character vector

**Values:** 'off', 'on'

**Default:** 'on'

### Output state estimation covariance — Output state estimation error covariance

off (default) | on

If you select this parameter, a state estimation error covariance output port **P** is generated in the block.

#### Programmatic Use

**Block Parameter:** OutputStateCovariance

**Type:** character vector

**Values:** 'off', 'on'

**Default:** 'off'

### Data type — Data type for block parameters

double (default) | single

Use this parameter to specify the data type for all block parameters.

#### Programmatic Use

**Block Parameter:** DataType

**Type:** character vector

**Values:** 'single', 'double'

**Default:** 'double'

### Sample time — Block sample time

1 (default) | positive scalar

Block sample time, specified as a positive scalar. If the sample times of your state transition and measurement functions are different, select **Enable multirate operation** in the **Multirate** tab, and specify the sample times in the **Multirate** tab instead.

#### Dependencies

This parameter is available if in the **Multirate** tab, the **Enable multirate operation** parameter is off.

#### Programmatic Use

**Block Parameter:** SampleTime

**Type:** character vector, string

**Default:** '1'

## Multirate Tab

### Enable multirate operation — Enable specification of different sample times for state transition and measurement functions

off (default) | on

Select this parameter if the sample times of the state transition and measurement functions are different. You specify the sample times in the **Multirate** tab, in **Sample time**.

#### Programmatic Use

**Block Parameter:** EnableMultirate

**Type:** character vector

**Values:** 'off', 'on'

**Default:** 'off'

### Sample time — State transition and measurement function sample times

positive scalar

If the sample times for state transition and measurement functions are different, specify **Sample time**. Specify the sample times for the measurement functions as positive integer multiples of the state transition sample time. The sample times you specify correspond to the following input ports:

- Ports corresponding to state transition function — Additional input to state transition function **StateTransitionFcnInputs** and time-varying process noise covariance **Q**. The sample times of these ports must always equal the state transition function sample time, but can differ from the sample time of the measurement functions.
- Ports corresponding to  $i^{\text{th}}$  measurement function — Measured output  **$y_i$** , additional input to measurement function **MeasurementFcnInputs**, enable signal at port **Enable $i$** , and time-varying measurement noise covariance  **$R_i$** . The sample times of these ports for the same measurement function must always be the same, but can differ from the sample time for the state transition function and other measurement functions.

#### Dependencies

This parameter is available if in the **Multirate** tab, the **Enable multirate operation** parameter is on.

**Programmatic Use**

**Block Parameter:** StateTransitionFcnSampleTime,  
MeasurementFcn1SampleTime1, MeasurementFcn1SampleTime2,  
MeasurementFcn1SampleTime3, MeasurementFcn1SampleTime4,  
MeasurementFcn1SampleTime5

**Type:** character vector; string

**Default:** '1'

## Definitions

### State Transition and Measurement Functions

The algorithm computes the state estimates  $\hat{x}$  of the nonlinear system using state transition and measurement functions specified by you. You can specify up to five measurement functions, each corresponding to a sensor in the system. The software lets you specify the noise in these functions as additive or nonadditive.

- **Additive Noise Terms** — The state transition and measurements equations have the following form:

$$x[k+1] = f(x[k], u_s[k]) + w[k]$$

$$y[k] = h(x[k], u_m[k]) + v[k]$$

Here  $f$  is a nonlinear state transition function that describes the evolution of states  $x$  from one time step to the next. The nonlinear measurement function  $h$  relates  $x$  to the measurements  $y$  at time step  $k$ .  $w$  and  $v$  are the zero-mean, uncorrelated process and measurement noises, respectively. These functions can also have additional optional input arguments that are denoted by  $u_s$  and  $u_m$  in the equations. For example, the additional arguments could be time step  $k$  or the inputs  $u$  to the nonlinear system. There can be multiple such arguments.

Note that the noise terms in both equations are additive. That is,  $x(k+1)$  is linearly related to the process noise  $w(k)$ , and  $y(k)$  is linearly related to the measurement noise  $v(k)$ . For additive noise terms, you do not need to specify the noise terms in the state transition and measurement functions. The software adds the terms to the output of the functions.

- **Nonadditive Noise Terms** — The software also supports more complex state transition and measurement functions where the state  $x[k]$  and measurement  $y[k]$  are

nonlinear functions of the process noise and measurement noise, respectively. When the noise terms are nonadditive, the state transition and measurements equation have the following form:

$$\begin{aligned}x[k+1] &= f(x[k], w[k], u_s[k]) \\ y[k] &= h(x[k], v[k], u_m[k])\end{aligned}$$

## Extended Capabilities

### C/C++ Code Generation

Generate C and C++ code using Simulink® Coder™.

The state transition, measurement, and Jacobian functions that you specify must use only the MATLAB commands and Simulink blocks that support code generation. For a list of blocks that support code generation, see “Simulink Built-In Blocks That Support Code Generation” (Simulink Coder). For a list of commands that support code generation, see “Functions and Objects Supported for C/C++ Code Generation — Category List” (MATLAB Coder).

### See Also

#### Blocks

Kalman Filter | Particle Filter | Unscented Kalman Filter

#### Functions

extendedKalmanFilter | particleFilter | unscentedKalmanFilter

#### Topics

“What Is Online Estimation?”

“Extended and Unscented Kalman Filter Algorithms for Online State Estimation”

“Validate Online State Estimation in Simulink”

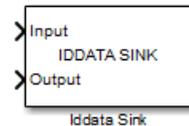
“Troubleshoot Online State Estimation”

**Introduced in R2017a**

## Iddata Sink

Export simulation data as `iddata` object to MATLAB workspace

**Library:** System Identification Toolbox



## Description

The Iddata Sink block exports simulation data as an `iddata` object to the MATLAB workspace. The object stores the input and simulated output signals, sampled at the sample time that you specify. If you simulate your model from the model window, the block exports the object to the MATLAB base workspace. If you simulate the model programmatically, the object is exported to the MATLAB caller workspace. The caller workspace is the workspace of the function that called the currently running function.

## Ports

### Input

#### Input — Input of `iddata` object

scalar | vector

Input of `iddata` object, specified as a scalar for single-input data. For multichannel data with  $N_u$  inputs, specify `Input` as a vector of length  $N_u$ .

Data Types: `double`

#### Output — Output of `iddata` object

scalar | vector

Output of `iddata` object, specified as a scalar for single-output data. For multichannel data with  $N_y$  outputs, specify `Output` as a vector of length  $N_y$ .

Data Types: `double`

## Parameters

### **IDDATA Name — Name of iddata object**

data (default) | variable name

Name of `iddata` object, specified as a MATLAB variable name. The object is exported with this name to the MATLAB workspace.

### **Sample Time — Sample time in seconds**

0.1 (default) | finite positive number

Sample time in seconds, specified as a finite positive number. The `iddata` object stores the input and output signals, sampled at the sample time that you specify.

## See Also

IDDATA Source

## Topics

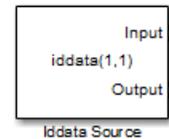
“Simulate Identified Model in Simulink”

**Introduced in R2008a**

## Iddata Source

Import time-domain data stored in `iddata` object in MATLAB workspace

**Library:** System Identification Toolbox



## Description

The Iddata Source block imports the input-output time-domain data stored in an `iddata` object in the MATLAB workspace. You can use this block to import data for simulating a model in Simulink.

## Ports

### Output

#### Input — Input data stored in `iddata` object

scalar | vector

Input data stored in `iddata` object, returned as a scalar for single-input data and a vector of length  $N_u$  for multichannel data with  $N_u$  inputs. If  $z$  is the `iddata` object, the output at this port at simulation time  $t$  is  $z.InputData$  at time  $t$ . If  $t$  is greater than  $z.SamplingInstants(end)$ , the maximum time in  $z$ , the input data is returned as  $\emptyset$  for each input channel.

Data Types: `double`

#### Output — Output data stored in `iddata` object

scalar | vector

Output data stored in `iddata` object, returned as a scalar for single-output data and as a vector of length  $N_y$  for multichannel data with  $N_y$  outputs. If  $z$  is the `iddata` object, the output at this port at simulation time  $t$  is  $z.OutputData$  at time  $t$ . If  $t$  is greater than

`z.SamplingInstants(end)`, the maximum time in `z`, the output data is returned as 0 for each output channel.

Data Types: `double`

## Parameters

**IDDATA object – Time-domain data to be imported**

`iddata(1,1)` (default) | `iddata` object

Time-domain data that is to be imported, specified as an `iddata` object that you have created in the MATLAB workspace. The `iddata` object must contain only one experiment. For a multiple-experiment `iddata` object `z`, to specify an `iddata` object for experiment number `kexp`, specify **IDDATA object** as `getexp(z, kexp)`.

## Extended Capabilities

### C/C++ Code Generation

Generate C and C++ code using Simulink® Coder™.

### See Also

Iddata Sink

### Topics

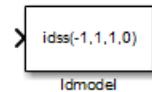
“Simulate Identified Model in Simulink”

**Introduced in R2008a**

## Idmodel

Simulate identified linear model in Simulink software

**Library:** System Identification Toolbox / Models



## Description

The Idmodel block simulates the output of an identified linear model using time-domain input data. The model is a state-space (`idss`), linear grey-box (`idgrey`), polynomial (`idpoly`), transfer function (`idtf`), or process (`idproc`) model that you previously estimated or created. For the simulation of state-space and linear grey-box models, you can specify the initial state values. For other linear models, initial conditions are set to zero. You can also add noise to the simulated output.

## Ports

### Input

#### Port\_1(In1) — Simulation input data

scalar | vector

Simulation input data, specified as a scalar for a single-input model. The data must be time-domain data. For multi-input models, specify the input as an  $Nu$ -element vector, where  $Nu$  is the number of inputs. For example, you can use a Vector Concatenate block to concatenate scalar signals into a vector signal.

---

**Note** Do not use a Bus Creator or Mux block to produce the vector signal.

---

Data Types: double

## Output

### Port\_1(Out1) — Simulated output

scalar | vector

Simulated output from linear model, returned as a scalar for a single-output model and an  $N_y$ -element vector for a model with  $N_y$  outputs.

Data Types: double

## Parameters

### Identified model — Linear model to be simulated

`idss(-1,1,1,0,'Ts',1)` (default) | `idss` object | `idgrey` object | `idpoly` object | `idtf` object | `idproc` object

Linear model to be simulated, specified as an `idss`, `idgrey`, `idpoly`, `idtf`, or `idproc` object. The model can be continuous-time or discrete-time, with or without input-output delays. You previously estimate or construct the linear model in the MATLAB workspace.

### Initial states (state space only: `idss`, `idgrey`) — Initial state values

`0` (default) | vector

Initial state values of state-space (`idss`) and linear grey-box (`idgrey`) models, specified as an  $N_x$ -element vector, where  $N_x$  is the number of states of the model. To estimate the initial states that provide a best fit between measured data and the simulated response of the model for the same input, use the `findstates` command.

For example, to compute initial states such that the response of the model `M` matches the simulated output data in the data set `z`, specify `X0`, such that:

```
X0 = findstates(M,z)
```

For linear models other than `idss` or `idgrey`, the block assumes that initial conditions are zero.

If you want to reproduce the simulation results that you get in the Model Output plot window in the System Identification app, or from the `compare` command:

- 1 If the identified model `m` is not a state-space or grey-box model, convert the model into state-space form (`idss` model), and specify the state-space model `mss` in the block.

```
mss = idss(m);
```

- 2 Compute the initial state values that produce the best fit between the model output and the measured output signal using `findstates`. Specify the prediction horizon as `Inf`, that is, minimize the simulation error.

```
X0 = findstates(mss,z,Inf);
```

- 3 Use the model `mss` and initial states `X0` in the `Idmodel` block to perform the simulation. Specify the same input signal `z` for simulation that you used as validation data in the `app` or `compare`.

### Add noise — Add noise to simulated output

`on` (default) | `off`

When you select this parameter, the block derives the noise amplitude from the linear model property `model.NoiseVariance`. The software filters random Gaussian white noise with the noise transfer function of the model and adds the resulting noise to the simulated model response. If you want to add the same noise every time you run the Simulink model, specify the `Noise seed(s)` property.

For continuous-time models, the ideal variance of the noise term is infinite. In reality, you see a band-limited noise that accounts for the time constants of the system. You can interpret the resulting simulated output as filtered using a lowpass filter with a passband that does not distort the dynamics from the input.

### Noise seed(s) — Add same noise to output for multiple simulations

`[]` (default) | nonnegative integer | vector

The `Noise seed(s)` property seeds the random number generator such that the block adds the same noise to the simulated output every time you run the Simulink model. For information about using seeds, see `rng`.

For multi-output models, you can use independent noise realizations that generate the outputs with additive noise. Enter a vector of  $N_y$  nonnegative integer entries, where  $N_y$  is the number of output channels.

For random restarts that vary from one simulation to another, specify `Noise seed(s)` as `[]`.

### Dependency

To enable this parameter, select `Add noise`.

## Extended Capabilities

### C/C++ Code Generation

Generate C and C++ code using Simulink® Coder™.

### See Also

#### Functions

`findstates` | `idgrey` | `idpoly` | `idproc` | `idss` | `idtf` | `sim`

#### Blocks

Iddata Sink | Iddata Source

### Topics

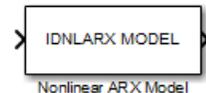
“Simulate Identified Model in Simulink”

### Introduced in R2008a

## Nonlinear ARX Model

Simulate nonlinear ARX model in Simulink software

**Library:** System Identification Toolbox / Models



### Description

The Nonlinear ARX Model block simulates the output of a nonlinear ARX model using time-domain input data. The model is an `idnlarx` model that you previously estimated or constructed in the MATLAB workspace. You specify initial conditions for the simulation as either steady-state input and output signal levels or as an initial state vector.

### Ports

#### Input

**Port\_1(In1) — Simulation input data**

scalar | vector

Simulation input data, specified as a scalar for a single-input model. The data must be time-domain data. For multi-input models, specify the input as an  $Nu$ -element vector, where  $Nu$  is the number of inputs. For example, you can use a Vector Concatenate block to concatenate scalar signals into a vector signal.

---

**Note** Do not use a Bus Creator or Mux block to produce the vector signal.

---

Data Types: `double`

## Output

### Port\_1(Out1) — Simulated output

scalar | vector

Simulated output from nonlinear ARX model, returned as a scalar for a single-output model and an  $N_y$ -element vector for a model with  $N_y$  outputs.

Data Types: double

## Parameters

### Model — Nonlinear ARX model to be simulated

`idnlarx` object

Nonlinear ARX model to be simulated, specified as an `idnlarx` object. You previously estimate or construct the `idnlarx` model in the MATLAB workspace.

### Initial conditions — Initial condition specification for simulation

Input and output values (default) | State values

The states of a nonlinear ARX model correspond to the dynamic elements of the nonlinear ARX model structure. The dynamic elements are the model regressors. Regressors can be the delayed input or output variables (standard regressors) or user-defined transformations of delayed input-output variables (custom regressors). For more information about the states of a nonlinear ARX model, see the `idnlarx` reference page.

For simulating nonlinear ARX models, you can specify the initial conditions one of the following:

- **Input and output values** — Specify steady-state input and output signal levels in `Input level` and `Output level`, respectively.
- **State values** — Specify a vector of length equal to the number of states in the model in `Specify initial states as a vector`.

### Input level — Steady-state input signal level

0 (default) | scalar

Steady-state input signal level before simulation, specified as a scalar.

### Dependency

To enable this parameter, specify `Initial` conditions as Input and output values.

### Output level — Steady-state output signal level

0 (default) | scalar

Steady-state output signal level before simulation, specified as a scalar.

### Dependency

To enable this parameter, specify `Initial` conditions as Input and output values.

### Specify initial states as a vector — Initial state values

0 (default) | vector

Initial state values of the model, specified as an  $N_x$ -element vector, where  $N_x$  is the number of states of the model. This parameter is named **Vector of state values** until you specify `Model`.

If you do not know the initial states, you can estimate these states as follows:

- To simulate the model around a given input level when you do not know the corresponding output level, estimate the equilibrium state values using the `findop` command. For example, to simulate a model `M` about a steady-state point where the input is 1 and the output is unknown, specify the initial state values as `X0`, where

```
X0 = findop(M, 'steady', 1, NaN)
```
- To estimate the initial states that provide a best fit between measured data and the simulated response of the model for the same input, use the `findstates` command. For example, to compute initial states such that the response of the model `M` matches the output data in the data set `z`, specify `X0`, such that:

```
X0 = findstates(M, z, Inf)
```
- To continue a simulation from a previous simulation run, use the simulated input-output values from the previous simulation to compute the initial states `X0` for the current simulation. Use the `data2state` command to compute `X0`. For example, suppose that `firstSimData` is a variable that stores the input and output values from a previous simulation. For a model `M`, you can specify `X0`, such that:

```
X0 = data2state(M, firstSimData)
```

**Dependency**

To enable this parameter, specify `Initial` conditions as `State` values.

## Extended Capabilities

### C/C++ Code Generation

Generate C and C++ code using Simulink® Coder™.

## See Also

**Functions**

`findstates` | `idnlarx` | `idnlarx/findop` | `sim`

**Blocks**

`Iddata Sink` | `Iddata Source`

**Topics**

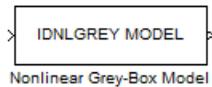
“Identifying Nonlinear ARX Models”

“Simulate Identified Model in Simulink”

**Introduced in R2008a**

## Nonlinear Grey-Box Model

Simulate nonlinear grey-box model in Simulink software



### Library

System Identification Toolbox

### Description

Simulates systems of nonlinear grey-box (`idnlgrey`) models.

### Input

Input signal to the model.

### Output

Output signal from the model.

### Parameters

#### **IDNLGREY model**

Name of `idnlgrey` variable in the MATLAB workspace.

#### **Initial state**

Specify the initial states as one of the following:

- 'z': Specifies zero, which corresponds to a system starting from rest.

- 'm': Specifies the internal initial states of the model.
- Vector of size equal to the number of states in the `idnlgrey` object.
- An initial state structure array. For information about creating this structure, type `help idnlgrey/sim` in the MATLAB Command Window.

## See Also

### Functions

`idnlgrey`

### Blocks

Iddata Sink | Iddata Source

### Topics

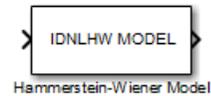
“Estimate Nonlinear Grey-Box Models”

**Introduced in R2008a**

## Hammerstein-Wiener Model

Simulate Hammerstein-Wiener model in Simulink software

**Library:** System Identification Toolbox / Models



### Description

The Hammerstein-Wiener Model block simulates the output of a Hammerstein-Wiener model using time-domain input data. The model is an `idnlhw` model that you previously estimated or constructed in the MATLAB workspace. You specify initial conditions for the simulation as one of the following:

- Zero for all states
- Initial state vector representing the initial states of the linear block

For information about the structure of a Hammerstein-Wiener model, see “What are Hammerstein-Wiener Models?”.

### Ports

#### Input

##### Port\_1(In1) — Simulation input data

scalar | vector

Simulation input data, specified as a scalar for a single-input model. The data must be time-domain data. For multi-input models, specify the input as an  $Nu$ -element vector, where  $Nu$  is the number of inputs. For example, you can use a Vector Concatenate block to concatenate scalar signals into a vector signal.

---

**Note** Do not use a Bus Creator or Mux block to produce the vector signal.

---

Data Types: double

## Output

### Port\_1(Out1) — Simulated output

scalar | vector

Simulated output from Hammerstein-Wiener model, returned as a scalar for a single-output model and as an  $N_y$ -element vector for a model with  $N_y$  outputs.

Data Types: double

## Parameters

### Model — Hammerstein-Wiener model to be simulated

idnlhw object

Hammerstein-Wiener model to be simulated, specified as an `idnlhw` object. You previously estimate or construct the `idnlhw` model in the MATLAB workspace.

### Initial conditions — Initial condition specification for simulation

Zero (default) | State values

The states of a Hammerstein-Wiener model correspond to the states of the embedded linear `idpoly` or `idss` model. For more information about the states, see the `idnlhw` reference page. You specify `Initial conditions` as one of the following:

- **Zero** — Specifies zero initial state values, which correspond to a simulation starting from a state of rest.
- **State values** — You specify the state values in **Specify a vector of state values**. Specify the states as a vector of length equal to the number of states in the model.

If you do not know the initial states, you can estimate these states as follows:

- To simulate the model around a given input level when you do not know the corresponding output level, you can estimate the equilibrium state values using the `findop` command.

For example, to simulate a model `M` about a steady-state point where the input is 1 and the output is unknown, you can specify the initial state values as `X0`, where:

```
X0 = findop(M, 'steady', 1, NaN)
```

- To estimate the initial states that provide a best fit between measured data and the simulated response of the model for the same input, use the `findstates` command.

For example, to compute initial states such that the response of the model `M` matches the simulated output data in the data set `z`, specify `X0`, such that:

```
X0 = findstates(M, z)
```

## Extended Capabilities

### C/C++ Code Generation

Generate C and C++ code using Simulink® Coder™.

## See Also

### Functions

`findstates` | `idnlhw` | `idnlhw/findop` | `sim`

### Blocks

`Iddata Sink` | `Iddata Source`

### Topics

“Identifying Hammerstein-Wiener Models”

“Simulate Identified Model in Simulink”

### Introduced in R2008a

# Kalman Filter

Estimate states of discrete-time or continuous-time linear system



## Library

Estimators

## Description

Use the Kalman Filter block to estimate states of a state-space plant model given process and measurement noise covariance data. The state-space model can be time-varying. A steady-state Kalman filter implementation is used if the state-space model and the noise covariance matrices are all time-invariant. A time-varying Kalman filter is used otherwise.

Kalman filter provides the optimal solution to the following continuous or discrete estimation problems:

### Continuous-Time Estimation

Given the continuous plant

$$\dot{x}(t) = A(t)x(t) + B(t)u(t) + G(t)w(t) \quad (\text{state equation})$$

$$y(t) = C(t)x(t) + D(t)u(t) + H(t)w(t) + v(t) \quad (\text{measurement equation})$$

with known inputs  $u$ , white process noise  $w$ , and white measurement noise  $v$  satisfying:

$$E[w(t)] = E[v(t)] = 0$$

$$E[w(t)w^T(t)] = Q(t)$$

$$E[w(t)v^T(t)] = N(t)$$

$$E[v(t)v^T(t)] = R(t)$$

construct a state estimate  $\hat{x}$  that minimizes the state estimation error covariance

$$P(t) = E[(x - \hat{x})(x - \hat{x})^T].$$

The optimal solution is the Kalman filter with equations

$$L(t) = (P(t)C^T(t) + \bar{N}),$$

$$\dot{P}(t) = A(t)P(t) + P(t)A^T(t) + \bar{Q}(t) - L(t)\bar{R}(t)L^T(t),$$

$$\dot{\hat{x}}(t) = A(t)\hat{x}(t) + B(t)u(t) + L(t)(y(t) - C(t)\hat{x}(t) - D(t)u(t)),$$

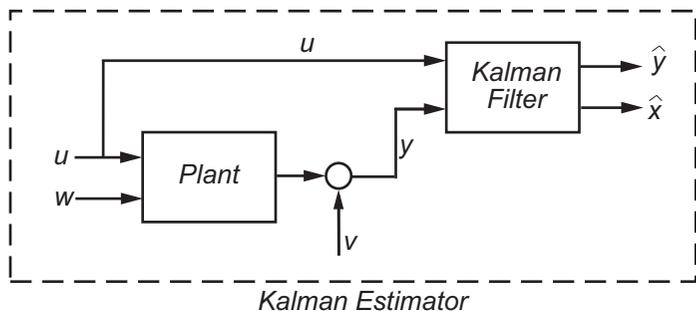
where

$$\bar{Q}(t) = G(t)Q(t)G^T(t),$$

$$\bar{R}(t) = R(t) + H(t)N(t) + N^T(t)H^T(t) + H(t)Q(t)H^T(t),$$

$$\bar{N}(t) = G(t)(Q(t)H^T(t) + N(t)).$$

The Kalman filter uses the known inputs  $u$  and the measurements  $y$  to generate the state estimates  $\hat{x}$ . If you want, the block can also output the estimates of the true plant output  $\hat{y}$ .



The block implements the steady-state Kalman filter when the system matrices ( $A(t)$ ,  $B(t)$ ,  $C(t)$ ,  $D(t)$ ,  $G(t)$ ,  $H(t)$ ) and noise covariance matrices ( $Q(t)$ ,  $R(t)$ ,  $N(t)$ ) are constant (specified in the Block Parameters dialog box). The steady-state Kalman filter uses a constant matrix  $P$  that minimizes the steady-state estimation error covariance and solves the associated continuous-time algebraic Riccati equation:

$$P = \lim_{t \rightarrow \infty} E[(x - \hat{x})(x - \hat{x})^T].$$

### Discrete-Time Estimation

Given the discrete plant

$$\begin{aligned} x[n+1] &= A[n] x[n] + B[n] u[n] + G[n] w[n], \\ y[n] &= C[n] x[n] + D[n] u[n] + H[n] w[n] + v[n], \end{aligned}$$

with known inputs  $u$ , white process noise  $w$  and white measurement noise  $v$  satisfying

$$\begin{aligned} E[u[n]] &= E[v[n]] = 0, \\ E[u[n]w^T[n]] &= Q[n], \\ E[v[n]w^T[n]] &= R[n], \\ E[u[n]v^T[n]] &= N[n]. \end{aligned}$$

The estimator has the following state equation

$$\hat{x}[n+1 | n] = A[n] \hat{x}[n | n-1] + B[n] u[n] + L[n](y[n] - C[n] \hat{x}[n | n-1] - D[n] u[n]),$$

where the gain  $L[n]$  is calculated through the discrete Riccati equation:

$$\begin{aligned} L[n] &= (A[n]P[n]C^T[n] + \bar{N}[n])(C[n]P[n]C^T[n] + \bar{R}[n])^{-1}, \\ M[n] &= P[n]C^T[n](C[n]P[n]C^T[n] + \bar{R}[n])^{-1}, \\ Z[n] &= (I - M[n]C[n])P[n](I - M[n]C[n])^T + M[n]\bar{R}[n]M^T[n], \\ P[n+1] &= (A[n] - \bar{N}[n]\bar{R}^{-1}[n]C[n])Z[A[n] - \bar{N}[n]\bar{R}^{-1}[n]C[n]]^T + \bar{Q}[n] - N[n]\bar{R}^{-1}[n]N^T[n], \end{aligned}$$

where  $I$  is the identity matrix of appropriate size and

$$\begin{aligned}\bar{Q}[n] &= G[n]Q[n]G^T[n], \\ \bar{R}[n] &= R[n] + H[n]N[n] + N^T[n]H^T[n] + H[n]Q[n]H^T[n], \\ \bar{N}[n] &= G[n](Q[n]H^T[n] + N[n]), \\ \text{and} \\ P[n] &= \mathbf{E}[(x - \hat{x}[n | n - 1])(x - \hat{x}[n | n - 1])^T], \\ Z[n] &= \mathbf{E}[(x - \hat{x}[n | n])(x - \hat{x}[n | n])^T],\end{aligned}$$

The steady-state Kalman filter uses a constant matrix  $P$  that minimizes the steady-state estimation error covariance and solves the associated discrete-time algebraic Riccati equation.

There are two variants of discrete-time Kalman filters:

- The current estimator generates the state estimates  $\hat{x}[n | n]$  using all measurement available, including  $y[n]$ . The filter updates  $\hat{x}[n | n - 1]$  with  $y[n]$  and outputs:

$$\begin{aligned}\hat{x}[n | n] &= \hat{x}[n | n - 1] + M[n](y[n] - C[n]\hat{x}[n | n - 1] - D[n]u[n]), \\ \hat{y}[n | n] &= C[n]\hat{x}[n | n] + D[n]u[n].\end{aligned}$$

- The delayed estimator generates the state estimates  $\hat{x}[n | n - 1]$  using measurements up to  $y[n - 1]$ . The filter outputs  $\hat{x}[n | n - 1]$  as defined previously, along with the optional output  $\hat{y}[n | n - 1]$

$$\hat{y}[n | n - 1] = C[n]\hat{x}[n | n - 1] + D[n]u[n]$$

The current estimator has better estimation accuracy compared to the delayed estimator, which is important for slow sample times. However, it has higher computational cost, making it harder to implement inside control loops. More specifically, it has direct feedthrough. This leads to an algebraic loop if the Kalman filter is used in a feedback loop that does not contain any delays (the feedback loop itself also has direct feedthrough). The algebraic loop can impact the speed of simulation. You cannot generate code if your model contains algebraic loops.

The Kalman Filter block differs from the `kalman` command in the following ways:

- When calling `kalman(sys, ...)`, `sys` includes the `G` and `H` matrices. Specifically, `sys.B` has `[B G]` and `sys.D` has `[D H]`. When you provide a LTI variable to the Kalman Filter block, it does not assume that the LTI variable provided contains `G` and `H`. They are optional and separate.
- The `kalman` command outputs `[yhat;xhat]` by default. The block only outputs `xhat` by default.

## Parameters

The following table summarizes the Kalman Filter block parameters, accessible via the Block Parameter dialog box.

Task	Parameters
Specify filter settings	<ul style="list-style-type: none"> <li>• <b>Time domain</b> on page 2-44</li> <li>• <b>Use the current measurement <math>y[n]</math> to improve <math>\hat{x}[n]</math></b> on page 2-44</li> </ul>
Specify the system model	<b>Model source</b> on page 2-44 in <b>Model Parameters</b> tab
Specify initial state estimates	<b>Source</b> on page 2-46 in <b>Model Parameters</b> tab
Specify noise characteristics	In <b>Model Parameters</b> tab: <ul style="list-style-type: none"> <li>• <b>Use <math>G</math> and <math>H</math> matrices (default <math>G=I</math> and <math>H=0</math>)</b> on page 2-47</li> <li>• <b><math>Q</math></b> on page 2-47, <b>Time-invariant <math>Q</math></b> on page 2-47</li> <li>• <b><math>R</math></b> on page 2-47, <b>Time-invariant <math>R</math></b> on page 2-48</li> <li>• <b><math>N</math></b> on page 2-48, <b>Time-invariant <math>N</math></b> on page 2-48</li> </ul>

Task	Parameters
Specify additional inports	In <b>Options</b> tab: <ul style="list-style-type: none"> <li>• <b>Add input port u</b> on page 2-48</li> <li>• <b>Add input port Enable to control measurement updates</b> on page 2-48</li> <li>• <b>External reset</b> on page 2-49</li> </ul>
Specify additional outports	In <b>Options</b> tab: <ul style="list-style-type: none"> <li>• <b>Output estimated model output y</b> on page 2-50</li> <li>• <b>Output state estimation error covariance Z</b> on page 2-50</li> </ul>

## Time domain

Specify whether to estimate continuous-time or discrete-time states:

- **Discrete-Time (Default)** — Block estimates discrete-time states
- **Continuous-Time** — Block estimates continuous-time states

When the Kalman Filter block is in a model with synchronous state control (see the State Control block), you cannot select **Continuous-time**.

## Use the current measurement $y[n]$ to improve $\hat{x}[n]$

Use the current estimator variant of the discrete-time Kalman filter. When not selected, the delayed estimator (variant) is used.

This option is available only when **Time Domain** is **Discrete-Time**.

## Model source

Specify how the A, B, C, D matrices are provided to the block. Must be one of the following:

- **Dialog: LTI State-Space Variable** — Use the values specified in the LTI state-space variable. You must also specify the variable name in **Variable**. The sample time

of the model must match the setting in the **Time domain** option, i.e. the model must be discrete-time if the **Time domain** is discrete-time.

- **Dialog: Individual A, B, C, D matrices** — Specify values in the following block parameters:
  - **A** — Specify the A matrix. It must be real and square.
  - **B** — Specify the B matrix. It must be real and have as many rows as the A matrix. This option is available only when **Add input port u** is selected in the **Options** tab.
  - **C** — Specify the C matrix. It must be real and have as many columns as the A matrix.
  - **D** — Specify the D matrix. It must be real. It must have as many rows as the C matrix and as many columns as the B matrix. This option is available only when **Add input port u** is selected in the **Options** tab.
- **External** — Specify the A, B, C, D matrices as input signals to the Kalman Filter block. If you select this option, the block includes additional input ports A, B, C and D. You must also specify the following in the block parameters:
  - **Number of states** — Number of states to be estimated, specified as a positive integer. The default value is 2.
  - **Number of inputs** — Number of known inputs in the model, specified as a positive integer. The default value is 2. This option is only available when **Add input port u** is selected.
  - **Number of outputs** — Number of measured outputs in the model, specified as a positive integer. The default value is 2.

## Sample Time

Block sample time, specified as -1 or a positive scalar.

This option is available only when **Time Domain** is **Discrete Time** and **Model Source** is **Dialog: Individual A, B, C, D matrices** or **External**. The sample time is obtained from the LTI state-space variable if the Model Source is **Dialog: LTI State-Space Variable**.

The default value is -1, which implies that the block inherits its sample time based on the context of the block within the model. All block input ports must have the same sample time.

## Source

Specify how to enter the initial state estimates and initial state estimation error covariance:

- **Dialog** — Specify the values directly in the dialog box. You must also specify the following parameters:
  - **Initial states  $\mathbf{x}[0]$**  — Specify the initial state estimate as a real scalar or vector. If you specify a scalar, all initial state estimates are set to this scalar. If you specify a vector, the length of the vector must match with the number of states in the model.
  - **State estimation error covariance  $\mathbf{P}[0]$**  (only when time-varying Kalman filter is used) — Specify the initial state estimation error covariance  $\mathbf{P}[0]$  for discrete-time Kalman filter or  $\mathbf{P}(0)$  for continuous-time Kalman filter. Must be specified as one of the following:
    - Real nonnegative scalar.  $\mathbf{P}$  is an  $N_s$ -by- $N_s$  diagonal matrix with the scalar on the diagonals.  $N_s$  is the number of states in the model.
    - Vector of real nonnegative scalars.  $\mathbf{P}$  is an  $N_s$ -by- $N_s$  diagonal matrix with the elements of the vector on the diagonals of  $\mathbf{P}$ .
    - $N_s$ -by- $N_s$  positive semi-definite matrix.
- **External** — Inherit the values from input ports. The block includes an additional input port  $X0$ . A second additional input port  $P0$  is added when time-varying Kalman filter is used.  $X0$  and  $P0$  must satisfy the same conditions described previously when you specify them in the dialog box.

## Use the Kalman Gain $\mathbf{K}$ from the model variable

Specify whether to use the pre-identified Kalman Gain contained in the state-space plant model. This option is available only when:

- **Model Source** is **Dialog: LTI State-Space Variable** and **Variable** is an identified state-space model (`ids`) with a nonzero  $\mathbf{K}$  matrix.
- **Time Invariant Q**, **Time Invariant R** and **Time Invariant N** options are selected.

If the **Use G and H matrices (default  $\mathbf{G}=\mathbf{I}$  and  $\mathbf{H}=\mathbf{0}$ )** option is selected, **Time Invariant G** and **Time Invariant H** options must also be selected.

## Use G and H matrices (default $G=I$ and $H=0$ )

Specify whether to use non-default values for the G and H matrices. If you select this option, you must specify:

- **G** — Specify the G matrix. It must be a real matrix with as many rows as the A matrix. The default value is 1.
- **Time-invariant G** — Specify if the G matrix is time invariant. If you unselect this option, the block includes an additional input port G.
- **H** — Specify the H matrix. It must be a real matrix with as many rows as the C matrix and as many columns as the G matrix. The default value is 0.
- **Time-invariant H** — Specify if the H matrix is time invariant. If you unselect this option, the block includes an additional input port G.
- **Number of process noise inputs** — Specify the number of process noise inputs in the model. The default value is 1.

This option is available only when **Time-invariant G** and **Time-invariant H** are cleared. Otherwise, this information is inferred from the G or H matrix.

## Q

Process noise covariance matrix, specified as one of the following:

- Real nonnegative scalar. Q is an  $N_w$ -by- $N_w$  diagonal matrix with the scalar on the diagonals.  $N_w$  is the number of process noise inputs in the model.
- Vector of real nonnegative scalars. Q is an  $N_w$ -by- $N_w$  diagonal matrix with the elements of the vector on the diagonals of Q.
- $N_w$ -by- $N_w$  positive semi-definite matrix.

## Time Invariant Q

Specify if the Q matrix is time invariant. If you unselect this option, the block includes an additional input port Q.

## R

Measurement noise covariance matrix, specified as one of the following:

- Real positive scalar.  $R$  is an  $N_y$ -by- $N_y$  diagonal matrix with the scalar on the diagonals.  $N_y$  is the number of measured outputs in the model.
- Vector of real positive scalars.  $R$  is an  $N_y$ -by- $N_y$  diagonal matrix with the elements of the vector on the diagonals of  $R$ .
- $N_y$ -by- $N_y$  positive-definite matrix.

### Time Invariant R

Specify if the  $R$  matrix is time invariant. If you unselect this option, the block includes an additional input port  $R$ .

### N

Process and measurement noise cross-covariance matrix. Specify it as a  $N_w$ -by- $N_y$  matrix. The matrix  $[Q \ N; N^T \ R]$  must be positive definite.

### Time Invariant N

Specify if the  $N$  matrix is time invariant. If you unselect this option, the block includes an additional input port  $N$ .

### Add input port $u$

Select this option if your model contains known inputs  $u(t)$  or  $u[k]$ . The option is selected by default. Unselecting this option removes the input port  $u$  from the block and removes the **B**, **D** and **Number of inputs** parameters from the block dialog box.

### Add input port Enable to control measurement updates

Select this option if you want to control the measurement updates. The block includes an additional input port `Enable`. The `Enable` input port takes a scalar signal. This option is cleared by default.

By default the block does measurement updates at each time step to improve the state and output estimates  $\hat{x}$  and  $\hat{y}$  based on measured outputs. The measurement update is skipped for the current sample time when the signal in the `Enable` port is 0. Concretely,

the equation for state estimates become  $\dot{\hat{x}}(t) = A(t)\hat{x}(t) + B(t)u(t)$  for continuous-time Kalman filter and  $\hat{x}[n+1|n] = A[n]\hat{x}[n|n-1] + B[n]u[n]$  for discrete-time.

## External Reset

Option to reset estimated states and parameter covariance matrix using specified initial values.

Suppose you reset the block at a time step,  $t$ . If the block is enabled at  $t$ , the software uses the initial parameter values specified either in the block dialog or the input ports P0 and X0 to estimate the states. In other words, at  $t$ , the block performs a time update and if it is enabled, a measurement update after the reset. The block outputs these updated estimates.

Specify one of the following:

- **None (Default)** — Estimated states  $\hat{x}$  and state estimation error covariance matrix P values are not reset.
- **Rising** — Triggers a reset when the control signal rises from a negative or zero value to a positive value. If the initial value is negative, rising to zero triggers a reset.
- **Falling** — Triggers a reset when the control signal falls from a positive or a zero value to a negative value. If the initial value is positive, falling to zero triggers a reset.
- **Either** — Triggers a reset when the control signal is either rising or falling.
- **Level** — Triggers a reset in either of these cases:
  - The control signal is nonzero at the current time step.
  - The control signal changes from nonzero at the previous time step to zero at the current time step.
- **Level hold** — Triggers reset when the control signal is nonzero at the current time step.

When you choose an option other than **None**, a **Reset** input port is added to the block to provide the reset control input signal.

## Output estimated model output $\hat{y}$

Add  $\hat{y}$  output port to the block to output the estimated model outputs. The option is cleared by default.

## Output state estimation error covariance $P$ or $Z$

Add  $P$  output port or  $Z$  output port to the block. The  $Z$  matrix is provided only when **Time Domain** is **Discrete Time** and the **Use the current measurement  $y[n]$  to improve  $\hat{x}[n]$**  is selected. Otherwise, the  $P$  matrix, as described in the “Description” on page 2-39 section previously, is provided.

The option is cleared by default.

## Ports

Port Name	Port Type (In/ Out)	Description
u (Optional)	In	Known inputs, specified as a real scalar or vector.
y	In	Measured outputs, specified as a real scalar or vector.
xhat	Out	Estimated states, returned as a real scalar or vector.
yhat (Optional)	Out	Estimated outputs, returned as a real scalar or vector.
P or Z (Optional)	Out	State estimation error covariance, returned as a matrix.
A (Optional)	In	A matrix, specified as a real matrix.
B (Optional)	In	B matrix, specified as a real matrix.
C (Optional)	In	C matrix, specified as a real matrix.
D (Optional)	In	D matrix, specified as a real matrix.
G (Optional)	In	G matrix, specified as a real matrix.
H (Optional)	In	H matrix, specified as a real matrix.

Port Name	Port Type (In/ Out)	Description
Q (Optional)	In	Q matrix, specified as a real scalar, vector or matrix.
R (Optional)	In	R matrix, specified as a real scalar, vector or matrix.
N (Optional)	In	N matrix, specified as a real matrix.
P0 (Optional)	In	P matrix at initial time, specified as a real scalar, vector, or matrix.
X0 (Optional)	In	Initial state estimates, specified as a real scalar or vector.
Enable (Optional)	In	Control signal to enable measurement updates, specified as a real scalar.
Reset (Optional)	In	Control signal to reset state estimates, specified as a real scalar.

## Supported Data Types

- Double-precision floating point
- Single-precision floating point (for discrete-time Kalman filter only)

---

### Note

- All input ports except **Enable** and **Reset** must have the same data type (single or double).
  - **Enable** and **Reset** ports support `single`, `double`, `int8`, `uint8`, `int16`, `uint16`, `int32`, `uint32`, and `boolean` data types.
- 

## Limitations

- The plant and noise data must satisfy:
  - $(C,A)$  detectable

- $\bar{R} > 0$  and  $\bar{Q} - \bar{N}\bar{R}^{-1}\bar{N}^T \geq 0$
- $(A - \bar{N}\bar{R}^{-1}C, \bar{Q} - \bar{N}\bar{R}^{-1}\bar{N}^T)$  has no uncontrollable mode on the imaginary axis (or unit circle in discrete time) with the notation

$$\bar{Q} = GQG^T$$

$$\bar{R} = R + HN + N^T H^T + HQH^T$$

$$\bar{N} = G(QH^T + N)$$

- The continuous-time Kalman filter cannot be used in Function-Call Subsystems or Triggered Subsystems.

## References

- [1] Franklin, G.F., J.D. Powell, and M.L. Workman, *Digital Control of Dynamic Systems*, Second Edition, Addison-Wesley, 1990.
- [2] Lewis, F., *Optimal Estimation*, John Wiley & Sons, Inc, 1986.

## Extended Capabilities

### C/C++ Code Generation

Generate C and C++ code using Simulink® Coder™.

### PLC Code Generation

Generate Structured Text code using Simulink® PLC Coder™.

## See Also

### Functions

`extendedKalmanFilter` | `kalman` | `particleFilter` | `unscentedKalmanFilter`

### **Blocks**

Extended Kalman Filter | Particle Filter | Unscented Kalman Filter

### **Topics**

“State Estimation Using Time-Varying Kalman Filter”

“What Is Online Estimation?”

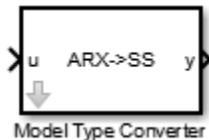
“Validate Online State Estimation in Simulink”

“Troubleshoot Online State Estimation”

### **Introduced in R2014b**

## Model Type Converter

Convert polynomial model coefficients to state-space model matrices



## Library

Estimators

## Description

Use the Model Type Converter block to convert the ARX, ARMAX, OE, or BJ model coefficients into state-space model matrices.

The block inport,  $u$ , requires a bus. The number of elements depends on the input polynomial model type:

- ARX — A, B
- ARMAX — A, B, C
- OE — B, F
- BJ — B,C, D, F

These bus elements must contain row vectors of the estimated coefficient values as outputted by the Recursive Polynomial Model Estimator block. For MISO data, specify  $B$  polynomial coefficients as a matrix where the  $i$ -th row parameters correspond to the  $i$ -th input. The coefficient values can vary with time. The Model Type Converter block converts these coefficients into the A, B, C, and D matrices of a discrete-time state-space model. The Model Type Converter block output,  $y$ , returns a bus with elements that correspond to the A, B, C, and D matrices of the state-space model. If the signals in  $u$  are time-varying, then the state-space matrices are time-varying too.

You can also estimate a state-space model online by using the Recursive Polynomial Model Estimator and Model Type Converter blocks together. Connect the output of the

Recursive Polynomial Model Estimator block to the inport of the Model Type Converter block to obtain online values of the state-space matrices. The conversion ignores the noise component of the models. In other words, the state-space matrices only capture the  $y(t)/u(t)$  relationship.

## Parameters

### Input model type

Specify the model type coefficients to convert to state-space model matrices. Specify one of the following model types:

- ARX
- ARMAX
- OE
- BJ

## Ports

Port	Port Type (In/ Out)	Description
u	In	Estimated A, B, C, D and F polynomial coefficients, specified as a bus with elements: A, B, C, D and F.
y	Out	State-space model, returned as a bus with elements that correspond to the A, B, C, and D matrices.

## Supported Data Types

- Double-precision floating point
- Single-precision floating point

## **Extended Capabilities**

### **C/C++ Code Generation**

Generate C and C++ code using Simulink® Coder™.

### **PLC Code Generation**

Generate Structured Text code using Simulink® PLC Coder™.

### **See Also**

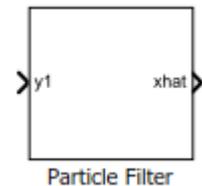
Recursive Polynomial Model Estimator

**Introduced in R2014a**

## Particle Filter

Estimate states of discrete-time nonlinear system using particle filter

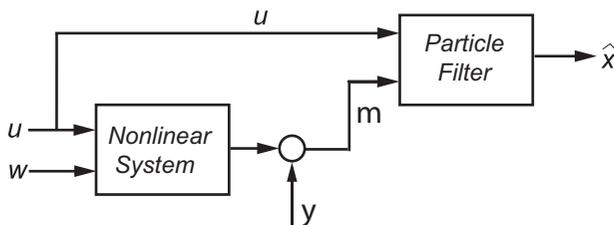
**Library:** Control System Toolbox / State Estimation  
System Identification Toolbox / Estimators



## Description

The Particle Filter block estimates the states of a discrete-time nonlinear system using the discrete-time particle filter algorithm.

Consider a plant with states  $x$ , input  $u$ , output  $m$ , process noise  $w$ , and measurement  $y$ . Assume that you can represent the plant as a nonlinear system.



The algorithm computes the state estimates  $\hat{x}$  of the nonlinear system using the state transition and measurement likelihood functions you specify.

You create the nonlinear state transition function and measurement likelihood functions for the system and specify these functions in the block. The block supports state estimation of a system with multiple sensors that are operating at different sampling rates. You can specify up to five measurement likelihood functions, each corresponding to a sensor in the system.

## Ports

### Input

#### **y1, y2, y3, y4, y5 — Measured system outputs**

vector

Measured system outputs corresponding to each measurement likelihood function that you specify in the block. The number of ports equals the number of measurement likelihood functions in your system. You can specify up to five measurement likelihood functions. For example, if your system has two sensors, you specify two measurement likelihood functions in the block. The first port **y1** is available by default. Click **Add Measurement**, to generate port **y2** corresponding to the second measurement likelihood function.

Specify the ports as  $N$ -dimensional vectors, where  $N$  is the number of quantities measured by the corresponding sensor. For example, if your system has one sensor that measures the position and velocity of an object, then there is only one port **y1**. The port is specified as a two-dimensional vector with values corresponding to position and velocity.

#### **Dependencies**

The first port **y1** is available by default. Ports **y2** to **y5** are generated when you click **Add Measurement**.

#### **StateTransitionFcnInputs — Optional input argument to state transition function**

scalar | vector | matrix

Optional input argument to the state transition function  $f$  other than the state  $x$ .

If you create  $f$  using a MATLAB function (.m file), the software generates the port **StateTransitionFcnInputs** when you enter the name of your function, and click **Apply**.

If your state transition function has more than one additional input, use a Simulink Function block to specify the function. When you use a Simulink Function block, you provide the additional inputs directly to the Simulink Function block using Inport blocks. No input ports are generated for the additional inputs in the Particle Filter block.

#### **Dependencies**

This port is generated only if both of the following conditions are satisfied:

- You specify  $f$  in **Function** using a MATLAB function, and  $f$  is on the MATLAB path.
- $f$  requires only one additional input argument apart from particles.

### **MeasurementLikelihoodFcn1Inputs, . . . , MeasurementLikelihoodFcn5Inputs** — **Optional input argument to each measurement likelihood function**

scalar | vector | matrix

Optional inputs to the measurement likelihood functions other than the state  $x$  and measurement  $y$ .

**MeasurementLikelihoodFcn1Inputs** corresponds to the first measurement likelihood function that you specify, and so on.

If you specify two measurement inputs using MATLAB functions (.m files) in **Function**, the software generates ports **MeasurementLikelihoodFcn1Inputs** and **MeasurementLikelihoodFcn2Inputs** when you click **Apply**. You can specify the inputs to these ports as scalars, vectors, or matrices.

If your measurement likelihood functions have more than one additional input, use Simulink Function blocks to specify the functions. When you use a Simulink Function block, you provide the additional inputs directly to the Simulink Function block using Inport blocks. No input ports are generated for the additional inputs in the Particle Filter block.

### **Dependencies**

A port corresponding to a measurement likelihood function  $h$  is generated only if both of the following conditions are satisfied:

- You specify measurement input  $h$  in **Function** using a MATLAB function, and  $h$  is on the MATLAB path.
- $h$  requires only one additional input argument apart from particles and measurement.

### **Enable1, Enable2, Enable3, Enable4, Enable5** — **Enable correction of estimated states when measured data is available**

scalar

Enable correction of estimated states when measured data is available.

For example, consider that measured output data is not available at all time points at the port **y1** that corresponds to the first measurement likelihood function. Then, use a signal value other than 0 at the **Enable1** port to enable the correction of estimated states when

measured data is available. Specify the port value as  $\emptyset$  when measured data is not available. Similarly, if measured output data is not available at all time points at the port  $\mathbf{y}_i$  for the  $i^{\text{th}}$  measurement likelihood function, specify the corresponding port **Enable $i$**  as a value other than  $\emptyset$ .

### Dependencies

If you select **Add Enable port** for a measurement likelihood function, a port corresponding to that measurement likelihood function is generated. The port appears when you click **Apply**.

## Output

### **xhat** — Estimated states

vector

Estimated states, returned as a vector of size  $N_s$ , where  $N_s$  is the number of states of the system. To access the individual states, use the Selector block.

When the **Use the current measurements to improve state estimates** parameter is selected, the block outputs the corrected state estimate  $\hat{x}[k|k]$  at time step  $k$ , estimated using measured outputs until time  $k$ . If you clear this parameter, the block returns the predicted state estimate  $\hat{x}[k|k-1]$  for time  $k$ , estimated using measured output until a previous time  $k-1$ . Clear this parameter if your filter is in a feedback loop and there is an algebraic loop in your Simulink model.

### **P** — State estimation error covariance

matrix

State estimation error covariance, returned as an  $N_s$ -by- $N_s$  matrix, where  $N_s$  is the number of states of the system. To access the individual covariances, use the Selector block.

You can output the error covariance only if you select **Output state estimation error covariance** in the **Block outputs, Multirate** tab, and click **Apply**.

### Dependencies

This parameter is available if in the **Block outputs, Multirate** tab, the **State estimation method** parameter is set to 'Mean'.

## Particles — Particle values used for state estimation

array

Particle values used for state estimation, returned as an  $N_s$ -by- $N_p$  or  $N_p$ -by- $N_s$  array.  $N_s$  is the number of states of the system, and  $N_p$  is the number of particles.

- If the `StateOrientation` parameter is specified as 'column', then **Particles** is returned as an  $N_s$ -by- $N_p$  array.
- If the `StateOrientation` parameter is specified as 'row', then **Particles** is returned as an  $N_p$ -by- $N_s$  array.

### Dependencies

This port is generated if you select **Output all particles** in the **Block outputs**, **Multirate** tab, and click **Apply**.

## Weights — Particle weights used for state estimation

vector

Particle weights used for state estimation, returned as a 1-by- $N_p$  or  $N_p$ -by-1 vector, where  $N_p$  is the number of particles used for state estimation.

- If the `StateOrientation` parameter is specified as 'column', then **Weights** is returned as a 1-by- $N_p$  vector, where each weight is associated with the particle in the same column in the **Particles** array.
- If the `StateOrientation` parameter is specified as 'row', then **Weights** is returned as a  $N_p$ -by-1 vector, where each weight is associated with the particle in the same row in the **Particles** array.

### Dependencies

This port is generated if you select **Output weights** in the **Block outputs**, **Multirate** tab, and click **Apply**.

## Parameters

### System Model Tab

#### State Transition

##### Function — State transition function name

'vdpParticleFilterStateFcn' (default) | function name

The particle filter state transition function calculates the particles at time step  $k+1$ , given particles at time step  $k$  per the dynamics of your system and process noise. This function has the syntax:

```
particlesNext = f(particles, param1, param2, ...)
```

where, *particles* and *particlesNext* have dimensions  $N_s$ -by- $N_p$  if **State Orientation** is specified as 'column', or  $N_p$ -by- $N_s$  if **State Orientation** is specified as 'row'. Also, *param\_i* represents optional input arguments you may specify. For more information on optional input arguments, see “StateTransitionFcnInputs” on page 2-0 .

You create the state transition function and specify the function name in **Function**. For example, if `vdpParticleFilterStateFcn.m` is the state transition function that you created and saved, specify **Function** as 'vdpParticleFilterStateFcn'.

You can create **Function** using a Simulink Function block or as a MATLAB function (.m file).

#### Programmatic Use

**Block Parameter:** StateTransitionFcn

**Type:** character vector, string

**Default:** 'vdpParticleFilterStateFcn'

#### Initialization

##### Number of Particles — Number of particles used in the filter

1000 (default) | positive scalar integer

Number of particles used in the filter, specified as a positive scalar integer. Each particle represents a state hypothesis in the system. A higher number of particles increases the state estimation accuracy, but also increases the computational effort required to run the filter.

**Programmatic Use****Block Parameter:** NumberOfParticles**Type:** positive scalar integer**Default:** 1000**Distribution – Initial distribution of particles**

'Gaussian' (default) | 'Uniform' | 'Custom'

Initial distribution of particles, specified as 'Gaussian', 'Uniform', or 'Custom'.

If you choose 'Gaussian', the initial set of particles or state hypotheses are distributed per the multivariate Gaussian distribution, where you specify the **Mean** and **Covariance**. The initial weight of all particles is assumed to be equal.

If you choose 'Uniform', the initial set of particles are distributed per the uniform distribution, where you specify the upper and lower **State bounds**. The initial weight of all particles is assumed to be equal.

'Custom' allows you to specify your own set of initial particles and their weights. You can use arbitrary probability distributions for **Particles** and **Weights** to initialize the filter.

**Programmatic Use****Block Parameter:** InitialDistribution**Type:** character vector**Values:** 'Gaussian', 'Uniform', 'Custom'**Default:** 'Gaussian'**Mean – Initial mean value of particles**

[0,0] (default) | vector

Initial mean value of particles, specified as a vector. The number of states to be estimated defines the length of the vector.

**Dependencies**

This parameter is available if in the **System model** tab, the **Distribution** parameter is set to **Gaussian**.

**Programmatic Use****Block Parameter:** InitialMean**Type:** array**Default:** [0,0]

### **Covariance — Initial covariance of particles**

1 (default) | scalar | vector | matrix

Initial covariance of particles, specified as a scalar, vector, or matrix.

If **Covariance** is specified as:

- A scalar, then it must be positive. The covariance is assumed to be a [ $N_s$   $N_s$ ] matrix with this scalar on the diagonals. Here,  $N_s$  is the number of states.
- A vector, then each element must be positive. The covariance is assumed to be a [ $N_s$   $N_s$ ] matrix with the elements of the vector on the diagonals.
- A matrix, then it must be positive semidefinite.

#### **Dependencies**

This parameter is available if in the **System model** tab, the **Distribution** parameter is set to Gaussian.

#### **Programmatic Use**

**Block Parameter:** InitialCovariance

**Type:** scalar, vector, or matrix

**Default:** 1

### **Circular Variables — Circular variables used for state estimation**

0 (default) | scalar | vector

Circular variables used for state estimation, specified as a scalar, or  $N_s$ -element vector, where  $N_s$  is the number of states.

If **Circular Variables** is specified as a scalar, the software extends it to a vector where each element is equal to this scalar. Circular (or angular) distributions use a probability density function with a range of  $[-\pi \pi]$ . Use circular variables if some of the states in your system represent angular quantities like the orientation of an object.

#### **Programmatic Use**

**Block Parameter:** CircularVariables

**Type:** scalar, vector

**Default:** 0

### **State Orientation — Orientation of input system states**

'column' (default) | 'row'

Orientation of system states, specified as 'column' or 'row'.

If **State Orientation** is specified as:

- 'column', then the first input argument to the state transition and measurement likelihood function is  $[Ns \ Np]$ . In this case,  $i^{th}$  column of this matrix is the  $i^{th}$  particle (state hypothesis). Also, the states estimates **xhat** is output as a  $[Ns \ 1]$  vector. Here,  $Ns$  is the number of states, and  $Np$  is the number of particles.
- 'row', then the first input argument to the state transition and measurement likelihood function is  $[Np \ Ns]$ , and each row of this matrix contains a particle. Also, the states estimates **xhat** is output as a  $[1 \ Ns]$  vector.

#### Programmatic Use

**Block Parameter:** StateOrientation

**Type:** character vector

**Values:** 'column', 'row'

**Default:** 'column'

#### State bounds — Initial bounds on system states

$[-3 \ 3; -3 \ 3]$  (default) | array

Initial bounds on system states, specified as an  $Ns$ -by-2 array, where  $Ns$  is the number of states.

The  $i^{th}$  row lists the lower and upper bound of the uniform distribution for the initial distribution of particles of the  $i^{th}$  state.

#### Dependencies

This parameter is available if in the **System model** tab, the **Distribution** parameter is set to Uniform.

#### Programmatic Use

**Block Parameter:** InitialStateBounds

**Type:** array

**Default:**  $[-3 \ 3; -3 \ 3]$

#### Particles — Custom particle distribution for state estimation

$[]$  (default) | array

Custom particle distribution for state estimation, specified as an  $Ns$ -by- $Np$  or  $Np$ -by- $Ns$  array.  $Ns$  is the number of states of the system, and  $Np$  is the number of particles.

- If the StateOrientation parameter is specified as 'column', then **Particles** is an  $Ns$ -by- $Np$  array.

- If the `StateOrientation` parameter is specified as 'row', then **Particles** is an  $Np$ -by- $Ns$  array.

### Dependencies

This parameter is available if in the **System model** tab, the **Distribution** parameter is set to Custom.

### Programmatic Use

**Block Parameter:** `InitialParticles`

**Type:** array

**Default:** `[]`

### Weights — Custom particle weight values for state estimation

`[]` (default) | positive vector

Custom particle weight values for state estimation, specified as a 1-by- $Np$  or  $Np$ -by-1 positive vector, where  $Np$  is the number of particles used for state estimation.

- If the `StateOrientation` parameter is specified as 'column', then **Weights** is a 1-by- $Np$  vector. Each weight in the vector is associated with the particle in the same column in the `Particles` array.
- If the `StateOrientation` parameter is specified as 'row', then **Weights** is a  $Np$ -by-1 vector. Each weight in the vector is associated with the particle in the same row in the `Particles` array.

### Dependencies

This parameter is available if in the **System model** tab, the **Distribution** parameter is set to Custom.

### Programmatic Use

**Block Parameter:** `InitialWeights`

**Type:** positive vector

**Default:** `[]`

### Measurement

#### Function — Measurement likelihood function name

'vdpMeasurementLikelihoodFcn' (default) | function name

The measurement likelihood function calculates the likelihood of particles (state hypotheses) using the sensor measurements. For each state hypothesis (particle), the

function first calculates an  $Nm$ -element measurement hypothesis vector. Then the likelihood of each measurement hypothesis is calculated based on the sensor measurement and the measurement noise probability distribution. this function has the syntax:

```
likelihood = h(particles, measurement, param1, param2, ...)
```

where, *likelihood* is an  $Np$ -element vector, where  $Np$  is the number of particles. *particles* have dimensions  $Ns$ -by- $Np$  if **State Orientation** is specified as 'column', or  $Np$ -by- $Ns$  if **State Orientation** is specified as 'row'. *measurement* is an  $Nm$ -element vector where,  $Nm$  is the number of measurements your sensor provides. *param<sub>i</sub>* represents optional input arguments you may specify. For more information on optional input arguments, see "MeasurementLikelihoodFcn1Inputs,...,MeasurementLikelihoodFcn5Inputs" on page 2-0 .

You create the measurement likelihood function and specify the function name in **Function**. For example, if `vdpMeasurementLikelihoodFcn.m` is the measurement likelihood function that you created and saved, specify **Function** as 'vdpMeasurementLikelihoodFcn'.

You can create **Function** using a Simulink Function block or as a MATLAB function (.m file).

- You can use a MATLAB function only if *h* has zero or one additional input argument *param<sub>i</sub>* other than **Particles** and **Measurement**.

The software generates an additional input port **MeasurementLikelihoodFcn<sub>i</sub>Inputs** to specify this argument for the  $i^{th}$  measurement likelihood function, and click **Apply**.

- If you are using a Simulink Function block, specify *x* and *y* using Argument Inport blocks and the additional inputs *param<sub>i</sub>* using Inport blocks in the Simulink Function block. You do not provide *param<sub>i</sub>* to the Particle Filter block.

If you have multiple sensors in your system, you can specify multiple measurement likelihood functions. You can specify up to five measurement likelihood functions using the **Add Measurement** button. To remove measurement likelihood functions, use **Remove Measurement**.

### Programmatic Use

**Block Parameter:** MeasurementLikelihoodFcn1, MeasurementLikelihoodFcn2, MeasurementLikelihoodFcn3, MeasurementLikelihoodFcn4, MeasurementLikelihoodFcn5

**Type:** character vector, string

**Default:** 'vdpMeasurementLikelihoodFcn'

### Add Enable Port — Enable correction of estimated states only when measured data is available

off (default) | on

Suppose that measured output data is not available at all time points at the port **y1** that corresponds to the first measurement likelihood function. To generate an input port **Enable1**, select **Add Enable port**. Use a signal at this port to enable the correction of estimated states only when measured data is available. Similarly, if measured output data is not available at all time points at the port **yi** for the  $i^{th}$  measurement likelihood function, select the corresponding **Add Enable port**.

### Programmatic Use

**Block Parameter:** HasMeasurementEnablePort1, HasMeasurementEnablePort2, HasMeasurementEnablePort3, HasMeasurementEnablePort4, HasMeasurementEnablePort5

**Type:** character vector

**Values:** 'off', 'on'

**Default:** 'off'

### Resampling

#### Resampling method — Method used for particle resampling

'Multinomial' (default) | 'Systemic' | 'Stratified'

Method used for particle resampling, specified as one of the following:

- 'Multinomial'
- 'Systematic'
- 'Stratified'

### Programmatic Use

**Block Parameter:** ResamplingMethod

**Type:** character vector

**Values:** 'Multinomial', 'Systemic', 'Stratified'

**Default:** 'Multinomial'

**Trigger method — Method to determine when resampling occurs**

'Ratio' (default) | 'Interval'

Method to determine when resampling occurs, specified as either 'Ratio' or 'Interval'. The 'Ratio' value triggers resampling based on the ratio of effective total particles. The 'Interval' value triggers resampling at regular time steps of the particle filter operation.

**Programmatic Use**

**Block Parameter:** TriggerMethod

**Type:** character vector

**Values:** 'Ratio', 'Interval'

**Default:** 'Ratio'

**Minimum effective particle ratio — Minimum desired ratio of the effective number of particles to the total number of particles**

0.5 (default) | positive scalar

Minimum desired ratio of the effective number of particles to the total number of particles, specified as a positive scalar. The effective number of particles is a measure of how well the current set of particles approximates the posterior distribution. A lower effective particle ratio implies that a lower number of particles are contributing to the estimation and resampling is required.

If the ratio of the effective number of particles to the total number of particles falls below the minimum effective particle ratio, a resampling step is triggered.

Specify minimum effective particle ratio as any value from 0 through 1.

**Dependencies**

This parameter is available if in the **System model** tab, the **Trigger method** parameter is set to Ratio.

**Programmatic Use**

**Block Parameter:** MinEffectiveParticleRatio

**Type:** scalar

**Values:** Range [0, 1]

**Default:** 0.5

**Sampling Interval — Fixed interval between resampling**

1 (default) | positive scalar integer

Fixed interval between resampling, specified as a positive scalar integer. The sampling interval determines during which correction steps the resampling is executed. For example, a value of two means the resampling is executed every second correction step. A value of `inf` means that resampling is never executed.

### Dependencies

This parameter is available if in the **System model** tab, the **Trigger method** parameter is set to `Interval`.

### Programmatic Use

**Block Parameter:** `SamplingInterval`

**Type:** positive scalar integer

**Default:** 1

### Random Number Generator Options

#### Randomness — Whether the random numbers are repeatable

`'Repeatable'` (default) | `'Not repeatable'`

Whether the random numbers are repeatable, specified as either `'Repeatable'` or `'Not repeatable'`. If you want to be able to produce the same result more than once, set **Randomness** to `'Repeatable'`, and specify the same random number generator seed value in **Seed**.

### Programmatic Use

**Block Parameter:** `Randomness`

**Type:** character vector

**Values:** `'Repeatable'`, `'Not repeatable'`

**Default:** `'Repeatable'`

#### Seed — Seed value for repeatable random numbers

0 (default) | scalar

Seed value for repeatable random numbers, specified as a scalar.

### Dependencies

This parameter is available if in the **System model** tab, the **Randomness** parameter is set to `'Repeatable'`.

### Programmatic Use

**Block Parameter:** `Seed`

**Type:** scalar

**Default:** 0

### Settings

#### Sample time — Block sample time

1 (default) | positive scalar

Block sample time, specified as a positive scalar.

Use the **Sample time** parameter if your state transition and all measurement likelihood functions have the same sample time. Otherwise, select the **Enable multirate operation** option in the **Multirate** tab, and specify sample times in the same tab.

### Dependencies

This parameter is available if in the **Block output, Multirate** tab, the **Enable multirate operation** parameter is off.

### Programmatic Use

**Block Parameter:** SampleTime

**Type:** character vector, string

**Default:** '1'

## Block Outputs, Multirate Tab

### Outputs

#### State Estimation Method — Method used for extracting a state estimate from particles

'Mean' (default) | 'MaxWeight' | 'None'

Method used for extracting a state estimate from particles, specified as one of the following:

- 'Mean' — The Particle Filter block outputs the weighted mean of the particles, depending on the parameters **Weights** and **Particles**, as the state estimate.
- 'Maxweight' — The Particle Filter block outputs the particle with the highest weight as the state estimate.
- 'None' — Use this option to implement a custom state estimation method by accessing all particles using the **Output all particles** parameter from the **Block outputs, Multirate** tab.

### Programmatic Use

**Block Parameter:** StateEstimationMethod

**Type:** character vector, string

**Values:** 'Mean', 'MaxWeight', 'None'

**Default:** 'Mean'

### Output all particles — Output all particles

'off' (default) | 'on'

If you select this parameter, an output port for particles used in the estimation, **Particles** is generated in the block.

- If the StateOrientation parameter is specified as 'column', then the particles are output as an  $N_s$ -by- $N_p$  array.  $N_s$  is the number of states of the system, and  $N_p$  is the number of particles.
- If the StateOrientation parameter is specified as 'row', then the particles are output as an  $N_p$ -by- $N_s$  array.

### Programmatic Use

**Block Parameter:** OutputParticles

**Type:** character vector

**Values:** 'off', 'on'

**Default:** 'off'

### Output weights — Output particle weights

'off' (default) | 'on'

If you select this parameter, an output port for particle weights used in the estimation, **Weights** is generated in the block.

- If the StateOrientation parameter is specified as 'column', then the particle weights are output as a 1-by- $N_p$  vector. Here, where each weight is associated with the particle in the same column in the **Particles** array.  $N_p$  is the number of particles used for state estimation.
- If the StateOrientation parameter is specified as 'row', then the particle weights are output as a  $N_p$ -by-1 vector.

### Programmatic Use

**Block Parameter:** OutputWeights

**Type:** character vector

**Values:** 'off', 'on'

**Default:** 'off'

### Output state estimation error covariance — Output state estimation error covariance

'off' (default) | 'on'

If you select this parameter, a state estimation error covariance output port, **P** is generated in the block.

#### Dependencies

This parameter is available if in the **Block outputs, Multirate** tab, the **State estimation method** parameter is set to 'Mean'.

#### Programmatic Use

**Block Parameter:** OutputStateCovariance

**Type:** character vector

**Values:** 'off', 'on'

**Default:** 'off'

### Use the current measurements to improve state estimates — Option to use current measurements for state estimation

'on' (default) | 'off'

When this parameter is selected, the block outputs the corrected state estimate  $\hat{x}[k|k]$  at time step  $k$ , estimated using measured outputs until time  $k$ . If you clear this parameter,

the block returns the predicted state estimate  $\hat{x}[k|k-1]$  for time  $k$ , estimated using measured output until a previous time  $k-1$ . Clear this parameter if your filter is in a feedback loop and there is an algebraic loop in your Simulink model.

#### Programmatic Use

**Block Parameter:** UseCurrentEstimator

**Type:** character vector

**Values:** 'on', 'off'

**Default:** 'on'

#### Multirate

### Enable multirate operation — Enable specification of different sample times for state transition and measurement likelihood functions

'off' (default) | 'on'

Select this parameter if the sample times of the state transition or any of the measurement likelihood functions differ from the rest. You specify the sample times in the **Multirate** tab, in **Sample time**.

### Programmatic Use

**Block Parameter:** EnableMultirate

**Type:** character vector

**Values:** 'off', 'on'

**Default:** 'off'

### Sample time — State transition and measurement likelihood function sample times

positive scalar

If the sample times for state transition and measurement likelihood functions are different, specify **Sample time**. Specify the sample times for the measurement functions as positive integer multiples of the state transition sample time. The sample times you specify correspond to the following input ports:

- Ports corresponding to state transition function — Additional input to state transition function **StateTransitionFcnInputs**. The sample times of these ports must always equal the state transition function sample time, but can differ from the sample time of the measurement likelihood functions.
- Ports corresponding to  $i^{\text{th}}$  measurement likelihood function — Measured output **y $i$** , additional input to measurement likelihood function **MeasurementLikelihoodFcn $i$ Inputs**, enable signal at port **Enable $i$** . The sample times of these ports for the same measurement likelihood function must always be the same, but can differ from the sample time for the state transition function and other measurement likelihood functions.

### Dependencies

This parameter is available if in the **Block outputs**, **Multirate** tab, the **Enable multirate operation** parameter is on.

### Programmatic Use

**Block Parameter:** StateTransitionFcnSampleTime,  
MeasurementLikelihoodFcn1SampleTime1,  
MeasurementLikelihoodFcn1SampleTime2,  
MeasurementLikelihoodFcn1SampleTime3,

MeasurementLikelihoodFcn1SampleTime4,  
MeasurementLikelihoodFcn1SampleTime5

**Type:** character vector, string

**Default:** '1'

## Extended Capabilities

### C/C++ Code Generation

Generate C and C++ code using Simulink® Coder™.

The state transition and measurement likelihood functions that you specify must use only the MATLAB commands and Simulink blocks that support code generation. For a list of blocks that support code generation, see “Simulink Built-In Blocks That Support Code Generation” (Simulink Coder). For a list of commands that support code generation, see “Functions and Objects Supported for C/C++ Code Generation — Category List” (MATLAB Coder).

## See Also

### Blocks

Extended Kalman Filter | Kalman Filter | Unscented Kalman Filter

### Functions

extendedKalmanFilter | particleFilter | unscentedKalmanFilter

### Topics

“Parameter and State Estimation in Simulink Using Particle Filter Block”

“Validate Online State Estimation in Simulink”

“Troubleshoot Online State Estimation”

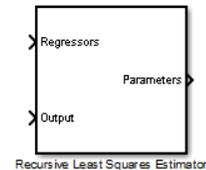
“Estimate States of Nonlinear System with Multiple, Multirate Sensors”

### Introduced in R2018a

## Recursive Least Squares Estimator

Estimate model coefficients using recursive least squares (RLS) algorithm

**Library:** System Identification Toolbox / Estimators



### Description

The Recursive Least Squares Estimator estimates the parameters of a system using a model that is linear in those parameters. Such a system has the following form:

$$y(t) = H(t)\theta(t).$$

$y$  and  $H$  are known quantities that you provide to the block to estimate  $\theta$ . The block can provide both infinite-history [1] and finite-history [2] (also known as sliding-window), estimates for  $\theta$ . For more information on these methods, see “Recursive Algorithms for Online Parameter Estimation”.

The block supports several estimation methods and data input formats. Configurable options in the block include:

- Sample-based or frame-based data format — See the **Input Processing** parameter.
- Infinite-history or finite- history estimation — See the **History** parameter.
- Multiple infinite-history estimation methods — See the **Estimation Method** parameter.
- Initial conditions, enable flag, and reset trigger — See the **Initial Estimate**, **Add enable port**, and **External Reset** parameters.

For a given time step  $t$ ,  $y(t)$  and  $H(t)$  correspond to the **Output** and **Regressors** inports of the Recursive Least Squares Estimator block, respectively.  $\theta(t)$  corresponds to the **Parameters** outport.

For example, suppose that you want to estimate a scalar gain,  $\theta$ , in the system  $y = h^2\theta$ . Here,  $y$  is linear with respect to  $\theta$ . You can use the Recursive Least Squares Estimator block to estimate  $\theta$ . Specify  $y$  and  $h^2$  as inputs to the **Output** and **Regressor** inports.

## Ports

### Input

#### Regressors — Regressors signal

vector | matrix

Regressors input signal  $H(t)$ . The **Input Processing** and **Number of Parameters** parameters define the dimensions of the signal:

- Sample-based input processing and  $N$  estimated parameters — 1-by- $N$  vector
- Frame-based input processing with  $M$  samples per frame and  $N$  estimated parameters —  $M$ -by- $N$  matrix

Data Types: single | double

#### Output — Measured output

scalar | vector

Measured output signal  $y(t)$ . The **Input Processing** parameter defines the dimensions of the signal:

- Sample-based input processing — Scalar
- Frame-based input processing with  $M$  samples per frame —  $M$ -by-1 vector

Data Types: single | double

#### Enable — Enable estimation updates

true (default) | false

External signal that allows you to enable and disable estimation updates. If the signal value is:

- `true` — Estimate and output the parameter values for the time step.
- `false` — Do not estimate the parameter values, and output the most recent previously estimated value.

### Dependencies

To enable this port, select the **Add enable port** parameter.

Data Types: `single` | `double` | `Boolean` | `int8` | `int16` | `int32` | `int64` | `uint8` | `uint16` | `uint32`

### Reset — Reset trigger

scalar

Reset parameter estimation to its initial conditions. The value of the **External reset** parameter determines the trigger type. The trigger type dictates whether the reset occurs on a signal that is rising, falling, either rising or falling, level, or on level hold.

### Dependencies

To enable this port, select any option other than **None** in the **External reset** dropdown.

Data Types: `single` | `double` | `Boolean` | `int8` | `int16` | `int32` | `uint8` | `uint16` | `uint32`

### InitialParameters — Initial parameter estimates

vector

Initial parameter estimates, supplied from a source external to the block. The block uses this inport at the beginning of the simulation or when you trigger an algorithm reset using the **Reset** signal.

The **Number of Parameters** parameter defines the dimensions of the signal. If there are  $N$  parameters, the signal is  $N$ -by-1.

### Dependencies

To enable this port, set **History** to **Infinite** and **Initial Estimate** to **External**.

Data Types: `single` | `double`

### InitialCovariance — Initial covariance of parameters

positive scalar (default) | vector of positive scalars | symmetric positive-definite matrix

Initial parameter covariances, supplied from a source external to the block. For details, see the **Parameter Covariance Matrix** parameter. The block uses this inport at the beginning of the simulation or when you trigger an algorithm reset using the **Reset** signal.

### Dependencies

To enable this port, set the following parameters:

- **History** to Infinite
- **Estimation Method** to Forgetting Factor or Kalman Filter
- **Initial Estimate** to External

Data Types: single | double

### InitialRegressors — Initial values of the regressors

matrix

Initial values of the regressors in the initial data window when using finite-history (sliding-window) estimation, supplied from an external source. The **Window length** parameter  $W$  and the **Number of Parameters** parameter  $N$  define the dimensions of this signal, which is  $W$ -by- $N$ .

The **InitialRegressors** signal controls the initial behavior of the algorithm. The block uses this inport at the beginning of the simulation or whenever the **Reset** signal triggers.

If the initial buffer is set to 0 or does not contain enough information, you see a warning message during the initial phase of your estimation. The warning should clear after a few cycles. The number of cycles it takes for sufficient information to be buffered depends upon the order of your polynomials and your input delays. If the warning persists, you should evaluate the content of your signals.

### Dependencies

To enable this port, set **History** to Finite and **Initial Estimate** to External.

Data Types: single | double

### InitialOutputs — Initial value of the measured output buffer

vector

Initial set of output measurements when using finite-history (sliding-window) estimation, supplied from an external source. The signal to this port must be a  $W$ -by-1 vector, where  $W$  is the window length.

The **InitialOutputs** signal controls the initial behavior of the algorithm. The block uses this inport at the beginning of the simulation or whenever the **Reset** signal triggers.

If the initial buffer is set to  $\emptyset$  or does not contain enough information, you see a warning message during the initial phase of your estimation. The warning should clear after a few cycles. The number of cycles it takes for sufficient information to be buffered depends upon the order of your polynomials and your input delays. If the warning persists, you should evaluate the content of your signals.

### Dependencies

To enable this port, set **History** to Finite, and **Initial Estimate** to External.

Data Types: single | double

## Output

### Parameters — Estimated parameters

vector

Estimated parameters  $\theta(t)$ , returned as an  $N$ -by-1 vector where  $N$  is the number of parameters.

Data Types: single | double

### Error — Estimation error

scalar | vector

Estimation error, returned as:

- Scalar — Sample-based input processing
- $M$ -by-1 vector — Frame-based input processing with  $M$  samples per frame

### Dependencies

To enable this port, select the **Output estimation error** parameter.

Data Types: single | double

### Covariance — Parameter estimation error covariance P

matrix

Parameter estimation error covariance  $P$ , returned as an  $N$ -by- $N$  matrix, where  $N$  is the number of parameters. For details, see the **Output Parameter Covariance Matrix** parameter.

## Dependencies

To enable this port:

- If **History** is Infinite, set **Estimation Method** to Forgetting Factor or Kalman Filter.
- Whether **History** is Infinite or Finite, select the **Output parameter covariance matrix** parameter.

Data Types: single | double

# Parameters

## Model Parameters

### Initial Estimate — Source of initial parameter estimates

None (default) | Internal | External

Specify how to provide initial parameter estimates to the block:

- None — Do not specify initial estimates.
  - If **History** is Infinite, the block uses 1 as the initial parameter estimate.
  - If **History** is Finite, the block calculates the initial parameter estimates from the initial **Regressors** and **Outputs** signals.

Specify **Number of Parameters**, and also, if **History** is Infinite, **Parameter Covariance Matrix**.

- Internal — Specify initial parameter estimates internally to the block
  - If **History** is Infinite, specify the **Initial Parameter Values** and **Parameter Covariance Matrix** parameters.
  - If **History** is Finite, specify the **Number of Parameters**, the **Initial Regressors**, and the **Initial Outputs** parameters.
- External — Specify initial parameter estimates as an input signal to the block.

Specify the **Number of Parameters** parameter. Your setting for the **History** parameter determines which additional signals to connect to the relevant ports:

- If **History** is Infinite — **InitialParameters** and **InitialCovariance**
- If **History** is Finite — **InitialRegressors** and **InitialOutputs**

### Programmatic Use

**Block Parameter:** InitialEstimateSource

**Type:** character vector, string

**Values:** 'None', 'Internal', 'External'

**Default:** 'None'

### Number of Parameters — Number of parameters to estimate

2 (default) | positive integer

Specify the number of parameters to estimate in the model, equal to the number of elements in the parameter  $\theta(t)$  vector.

### Dependencies

To enable this parameter, set either:

- **History** to Infinite and **Initial Estimate** to either None or External
- **History** to Finite

An alternative way to specify the number of parameters  $N$  to estimate is by using the **Initial Parameter Values** parameter, for which you define an initial estimate vector with  $N$  elements. This approach covers the one remaining combination, where **History** is Infinite and **Initial Estimate** is Internal. For more information, see **Initial Parameter Values**.

### Programmatic Use

**Block Parameter:** InitialParameterData

**Type:** positive integer

**Default:** 2

### Parameter Covariance Matrix — Initial parameter covariance

1e4 (default) | scalar | vector | matrix

Specify **Parameter Covariance Matrix** as a:

- Real positive scalar,  $\alpha$  — Covariance matrix is an  $N$ -by- $N$  diagonal matrix, with  $\alpha$  as the diagonal elements.
- Vector of real positive scalars,  $[\alpha_1, \dots, \alpha_N]$  — Covariance matrix is an  $N$ -by- $N$  diagonal matrix, with  $[\alpha_1, \dots, \alpha_N]$  as the diagonal elements.

- $N$ -by- $N$  symmetric positive-definite matrix.

Here,  $N$  is the number of parameters to be estimated.

#### Dependencies

To enable this parameter, set the following parameters:

- **History** to Infinite
- **Initial Estimate** to None or Internal
- **Estimation Method** to Forgetting Factor or Kalman Filter

#### Programmatic Use

**Block Parameter:**  $P_0$

**Type:** scalar, vector, or matrix

**Default:**  $1e4$

#### Initial Parameter Values — Initial values of the parameters to estimate

[1 1] (default) | vector

Specify initial parameter values as a vector of length  $N$ , where  $N$  is the number of parameters to estimate.

#### Dependencies

To enable this parameter, set **History** to Infinite and **Initial Estimate** to Internal.

#### Programmatic Use

**Block Parameter:** InitialParameterData

**Type:** real vector

**Default:** [1 1]

#### Initial Regressors — Initial values of the regressors buffer

$\emptyset$  (default) | matrix

Specify the initial values of the regressors buffer when using finite-history (sliding window) estimation. The **Window length** parameter  $W$  and the **Number of Parameters** parameter  $N$  define the dimensions of the regressors buffer, which is  $W$ -by- $N$ .

The **Initial Regressors** parameter controls the initial behavior of the algorithm. The block uses this parameter at the beginning of the simulation or whenever the **Reset** signal triggers.

When the initial value is set to  $\emptyset$ , the block populates the buffer with zeros.

If the initial buffer is set to  $\emptyset$  or does not contain enough information, you see a warning message during the initial phase of your estimation. The warning should clear after a few cycles. The number of cycles it takes for sufficient information to be buffered depends upon the order of your polynomials and your input delays. If the warning persists, you should evaluate the content of your signals.

### Dependencies

To enable this parameter, set **History** to Finite and **Initial Estimate** to Internal.

### Programmatic Use

**Block Parameter:** InitialRegressors

**Type:** real matrix

**Default:**  $\emptyset$

### Initial Outputs — Initial values of the measured outputs buffer

$\emptyset$  (default) | vector

Specify initial values of the measured outputs buffer when using finite-history (sliding-window) estimation. This parameter is a  $W$ -by-1 vector, where  $W$  is the window length.

When the initial value is set to  $\emptyset$ , the block populates the buffer with zeros.

If the initial buffer is set to  $\emptyset$  or does not contain enough information, you see a warning message during the initial phase of your estimation. The warning should clear after a few cycles. The number of cycles it takes for sufficient information to be buffered depends upon the order of your polynomials and your input delays. If the warning persists, you should evaluate the content of your signals.

The **Initial Outputs** parameter controls the initial behavior of the algorithm. The block uses this parameter at the beginning of the simulation or whenever the **Reset** signal triggers.

### Dependencies

To enable this parameter, set **History** to Finite and **Initial Estimate** to Internal.

### Programmatic Use

**Block Parameter:** InitialOutputs

**Type:** real vector

**Default:**  $\emptyset$

## Input Processing and Sample Time

### Input Processing — Choose sample-based or frame-based processing

Sample-based (default) | Frame-based

- Sample-based processing operates on signals streamed one sample at a time.
- Frame-based processing operates on signals containing samples from multiple time steps. Many machine sensor interfaces package multiple samples and transmit these samples together in frames. Frame-based processing allows you to input this data directly without having to first unpack it.

Specifying frame-based data adds an extra dimension of  $M$  to some of your data inports and outports, where  $M$  is the number of time steps in a frame. These ports are:

- **Regressors**
- **Output**
- **Error**

For more information, see the port descriptions in “Ports” on page 2-77.

#### Programmatic Use

**Block Parameter:** InputProcessing

**Type:** character vector, string

**Values:** 'Sample-based', 'Frame-based'

**Default:** 'Sample-based'

### Sample Time — Block sample time

-1 (default) | positive scalar

Specify the data sample time, whether by individual samples for sample-based processing ( $t_s$ ), or by frames for frame-based processing ( $t_f = Mt_s$ ), where  $M$  is the frame length. When you set **Sample Time** to its default value of -1, the block inherits its  $t_s$  or  $t_f$  based on the signal.

Specify **Sample Time** as a positive scalar to override the inheritance.

#### Programmatic Use

**Block Parameter:** Ts

**Type:** real scalar

**Default:** -1

## Algorithm and Block Options

### Algorithm Options

#### History — Choose infinite or finite data history

Infinite (default) | Finite

The **History** parameter determines what type of recursive algorithm you use:

- **Infinite** — Algorithms in this category aim to produce parameter estimates that explain all data since the start of the simulation. These algorithms retain the history in a data summary. The block maintains this summary within a fixed amount of memory that does not grow over time.

The block provides multiple algorithms of the **Infinite** type. Selecting this option enables the **Estimation Method** parameter with which you specify the algorithm.

- **Finite** — Algorithms in this category aim to produce parameter estimates that explain only a finite number of past data samples. The block uses all of the data within a finite window, and discards data once that data is no longer within the window bounds. This method is also called sliding-window estimation.

Selecting this option enables the **Window Length** parameter that sizes the sliding window.

For more information on recursive estimation methods, see “Recursive Algorithms for Online Parameter Estimation”

#### Programmatic Use

**Block Parameter:** History

**Type:** character vector, string

**Values:** 'Infinite', 'Finite'

**Default:** 'Infinite'

#### Window Length — Window size for finite sliding-window estimation

200 (default) | positive integer

The **Window Length** parameter determines the number of time samples to use for the sliding-window estimation method. Choose a window size that balances estimation performance with computational and memory burden. Sizing factors include the number and time variance of the parameters in your model. Always specify **Window Length** in samples, even if you are using frame-based input processing.

**Window Length** must be greater than or equal to the number of estimated parameters.

Suitable window length is independent of whether you are using sample-based or frame-based input processing. However, when using frame-based processing, **Window Length** must be greater than or equal to the number of samples (time steps) contained in the frame.

### Dependencies

To enable this parameter, set **History** to `Finite`.

### Programmatic Use

**Block Parameter:** `WindowLength`

**Type:** positive integer

**Default:** 200

### Estimation Method — Recursive estimation algorithm

`Forgetting Factor (default) | Kalman Filter | Normalized Gradient | Gradient`

Specify the estimation algorithm when performing infinite-history estimation. When you select any of these methods, the block enables additional related parameters.

Forgetting factor and Kalman filter algorithms are more computationally intensive than gradient and normalized gradient methods. However, these more intensive methods have better convergence properties than the gradient methods. For more information about these algorithms, see “Recursive Algorithms for Online Parameter Estimation”.

### Programmatic Use

**Block Parameter:** `EstimationMethod`

**Type:** character vector, string

**Values:** `'Forgetting Factor', 'Kalman Filter', 'Normalized Gradient', 'Gradient'`

**Default:** `'Forgetting Factor'`

### Forgetting Factor — Discount old data using forgetting factor

`1 (default) | positive scalar in (0 1] range`

The forgetting factor  $\lambda$  specifies if and how much old data is discounted in the estimation. Suppose that the system remains approximately constant over  $T_0$  samples. You can choose  $\lambda$  such that:

$$T_0 = \frac{1}{1 - \lambda}$$

- Setting  $\lambda = 1$  corresponds to “no forgetting” and estimating constant coefficients.
- Setting  $\lambda < 1$  implies that past measurements are less significant for parameter estimation and can be “forgotten.” Set  $\lambda < 1$  to estimate time-varying coefficients.

Typical choices of  $\lambda$  are in the [0.98 0.995] range.

### Dependencies

To enable this parameter, set **History** to Infinite and **Estimation Method** to Forgetting Factor.

### Programmatic Use

**Block Parameter:** AdaptationParameter

**Type:** scalar

**Values:** (0 1] range

**Default:** 1

### Process Noise Covariance — Process noise covariance for Kalman filter estimation method

1 (default) | nonnegative scalar | vector of nonnegative scalars | symmetric positive semidefinite matrix

**Process Noise Covariance** prescribes the elements and structure of the noise covariance matrix for the Kalman filter estimation. Using  $N$  as the number of parameters to estimate, specify the **Process Noise Covariance** as one of the following:

- Real nonnegative scalar,  $\alpha$  — Covariance matrix is an  $N$ -by- $N$  diagonal matrix, with  $\alpha$  as the diagonal elements.
- Vector of real nonnegative scalars,  $[\alpha_1, \dots, \alpha_N]$  — Covariance matrix is an  $N$ -by- $N$  diagonal matrix, with  $[\alpha_1, \dots, \alpha_N]$  as the diagonal elements.
- $N$ -by- $N$  symmetric positive semidefinite matrix.

The Kalman filter algorithm treats the parameters as states of a dynamic system and estimates these parameters using a Kalman filter. **Process Noise Covariance** is the covariance of the process noise acting on these parameters. Zero values in the noise covariance matrix correspond to constant coefficients, or parameters. Values larger than 0 correspond to time-varying parameters. Use large values for rapidly changing parameters. However, expect the larger values to result in noisier parameter estimates. The default value is 1.

**Dependencies**

To enable this parameter, set **History** to Infinite and **Estimation Method** to Kalman Filter.

**Programmatic Use**

**Block Parameter:** AdaptationParameter

**Type:** scalar, vector, matrix

**Default:** 1

**Adaptation Gain — Adaptation gain specification for gradient estimation methods**

1 (default) | positive scalar

The adaptation gain  $\gamma$  scales the influence of new measurement data on the estimation results for the gradient and normalized gradient methods. When your measurements are trustworthy, or in other words have a high signal-to-noise ratio, specify a larger value for  $\gamma$ . However, setting  $\gamma$  too high can cause the parameter estimates to diverge. This divergence is possible even if the measurements are noise free.

When **Estimation Method** is NormalizedGradient, **Adaptation Gain** should be less than 2. With either gradient method, if errors are growing in time (in other words, estimation is diverging), or parameter estimates are jumping around frequently, consider reducing **Adaptation Gain**.

**Dependencies**

To enable this parameter, set **History** to Infinite and **Estimation Method** to Normalized Gradient or to Gradient.

**Programmatic Use**

**Block Parameter:** AdaptationParameter

**Type:** scalar

**Default:** 1

**Normalization Bias — Bias for adaptation gain scaling for normalized gradient estimation method**

eps (default) | nonnegative scalar

The normalized gradient algorithm scales the adaptation gain at each step by the square of the two-norm of the gradient vector. If the gradient is close to zero, the near-zero denominator can cause jumps in the estimated parameters. **Normalization Bias** is the

term introduced to the denominator to prevent these jumps. Increase **Normalization Bias** if you observe jumps in estimated parameters.

### Dependencies

To enable this parameter, set **History** to Infinite and **Estimation Method** to Normalized Gradient.

### Programmatic Use

**Block Parameter:** NormalizationBias

**Type:** scalar

**Default:** eps

### Block Options

#### Output estimation error — Add Error output to block

off (default) | on

Use the **Error** output signal to validate the estimation. For a given time step  $t$ , the estimation error  $e(t)$  is calculated as:

$$e(t) = y(t) - y_{est}(t),$$

where  $y(t)$  is the measured output that you provide, and  $y_{est}(t)$  is the estimated output using the regressors  $H(t)$  and parameter estimates  $\theta(t-1)$ .

### Programmatic Use

**Block Parameter:** OutputError

**Type:** character vector, string

**Values:** 'off', 'on',

**Default:** 'off'

#### Output parameter covariance matrix — Add covariance output to block

off (default) | on

Use the **Covariance** output signal to examine parameter estimation uncertainty. The software computes parameter covariance  $P$  assuming that the residuals,  $e(t)$ , are white noise, and the variance of these residuals is 1.

The interpretation of  $P$  depends on the estimation approach you specify in **History** and **Estimation Method** as follows:

- If **History** is Infinite, then your **Estimation Method** selection results in:

- **Forgetting Factor** —  $(R_2/2)P$  is approximately equal to the covariance matrix of the estimated parameters, where  $R_2$  is the true variance of the residuals. The block outputs the residuals in the **Error** port.
- **Kalman Filter** —  $R_2P$  is the covariance matrix of the estimated parameters, and  $R_1/R_2$  is the covariance matrix of the parameter changes. Here,  $R_1$  is the covariance matrix that you specify in **Parameter Covariance Matrix**.
- **Normalized Gradient or Gradient** — Covariance  $P$  is not available.
- If **History** is **Finite** (sliding-window estimation) —  $R_2 P$  is the covariance of the estimated parameters. The sliding-window algorithm does not use this covariance in the parameter-estimation process. However, the algorithm does compute the covariance for output so that you can use it for statistical evaluation.

**Programmatic Use****Block Parameter:** OutputP**Type:** character vector, string**Values:** 'off', 'on'**Default:** 'off'**Add enable port — Add Enable inport to block**

off (default) | on

Use the **Enable** signal to provide a control signal that enables or disables parameter estimation. The block estimates the parameter values for each time step that parameter estimation is enabled. If you disable parameter estimation at a given step,  $t$ , then the software does not update the parameters for that time step. Instead, the block outputs the last estimated parameter values.

- You can use this option, for example, when or if:
  - Your regressors or output signal become too noisy, or do not contain information at some time steps
  - Your system enters a mode where the parameter values do not change in time

**Programmatic Use****Block Parameter:** AddEnablePort**Type:** character vector, string**Values:** 'off', 'on'**Default:** 'off'**External reset — Specify trigger for external reset**

None (default) | Rising | Falling | Either | Level | Level hold

Set the **External reset** parameter to both add a **Reset** inport and specify the inport signal condition that triggers a reset of algorithm states to their specified initial values. Reset the estimation, for example, if parameter covariance is becoming too large because of lack of either sufficient excitation or information in the measured signals.

Suppose that you reset the block at a time step,  $t$ . If the block is enabled at  $t$ , the software uses the initial parameter values specified in **Initial Estimate** to estimate the parameter values. In other words, at  $t$ , the block performs a parameter update using the initial estimate and the current values of the inports.

If the block is disabled at  $t$  and you reset the block, the block outputs the values specified in **Initial Estimate**.

Specify this option as one of the following:

- **None** — Algorithm states and estimated parameters are not reset.
- **Rising** — Trigger reset when the control signal rises from a negative or zero value to a positive value. If the initial value is negative, rising to zero triggers reset.
- **Falling** — Trigger reset when the control signal falls from a positive or a zero value to a negative value. If the initial value is positive, falling to zero triggers reset.
- **Either** — Trigger reset when the control signal is either rising or falling.
- **Level** — Trigger reset in either of these cases:
  - Control signal is nonzero at the current time step.
  - Control signal changes from nonzero at the previous time step to zero at the current time step.
- **Level hold** — Trigger reset when the control signal is nonzero at the current time step.

When you choose any option other than **None**, the software adds a **Reset** inport to the block. You provide the reset control input signal to this inport.

### **Programmatic Use**

**Block Parameter:** ExternalReset

**Type:** character vector, string

**Values:** 'None', 'Rising', 'Falling', 'Either', 'Level', 'Level hold'

**Default:** 'None'

## References

- [1] Ljung, L. *System Identification: Theory for the User*. Upper Saddle River, NJ: Prentice-Hall PTR, 1999, pp. 363-369.
- [2] Zhang, Q. "Some Implementation Aspects of Sliding Window Least Squares Algorithms." *IFAC Proceedings*. Vol. 33, Issue 15, 2000, pp. 763-768.

## Extended Capabilities

### C/C++ Code Generation

Generate C and C++ code using Simulink® Coder™.

### PLC Code Generation

Generate Structured Text code using Simulink® PLC Coder™.

## See Also

Kalman Filter | Recursive Polynomial Model Estimator

## Topics

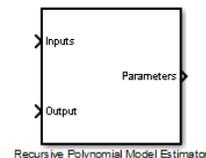
- "Estimate Parameters of System Using Simulink Recursive Estimator Block"
- "Online Recursive Least Squares Estimation"
- "Preprocess Online Parameter Estimation Data in Simulink"
- "Validate Online Parameter Estimation Results in Simulink"
- "Generate Online Parameter Estimation Code in Simulink"
- "Recursive Algorithms for Online Parameter Estimation"

**Introduced in R2014a**

## Recursive Polynomial Model Estimator

Estimate input-output and time-series polynomial model coefficients

**Library:** System Identification Toolbox / Estimators



### Description

#### Model Structures

Use the Recursive Polynomial Model Estimator block to estimate discrete-time input-output polynomial and time-series models.

These model structures are:

- AR —  $A(q)y(t) = e(t)$
- ARMA —  $A(q)y(t) = C(q)e(t)$
- ARX —  $A(q)y(t) = B(q)u(t-n_k) + e(t)$
- ARMAX —  $A(q)y(t) = B(q)u(t-n_k) + C(q)e(t)$
- 

$$\text{OE — } y(t) = \frac{B(q)}{F(q)}u(t-n_k) + e(t)$$

- BJ —  $y(t) = \frac{B(q)}{F(q)}u(t-n_k) + \frac{C(q)}{D(q)}e(t)$

$q$  is the time-shift operator and  $n_k$  is the input delay.  $u(t)$  is the input,  $y(t)$  is the output, and  $e(t)$  is the error. For MISO models, there are as many  $B(q)$  polynomials as the number of inputs.

The orders of these models correspond to the maximum number of time shifts, as represented by the exponent of  $q$ . For instance, the  $A(q)$  polynomial with order  $na$  can be represented as:

$$1+a_1q^{-1}+a_2q^{-2}+\dots+a_{na}q^{-na}.$$

An equivalent representation applies to the  $C(q)$ ,  $D(q)$ , and  $F(q)$  polynomials and their corresponding orders  $nc$ ,  $nd$ , and  $nf$ .

The  $B(q)$  polynomial is unique with respect to the others, because this polynomial operates on the input and contains the system zeros. For  $B(q)$ , the order  $nb$  is equal to the number of zeros plus 1:

$$b_1+b_2q^{-1}+b_3q^{-2}+\dots+b_{nb}q^{-(nb-1)}.$$

The orders  $na$ ,  $nb$ ,  $nc$ ,  $nd$ ,  $nf$ , and input delay  $nk$  are known ahead of time. Specify these values as block parameters. Provide  $u(t)$  and  $y(t)$  through the **Inputs** and **Outputs** inports, respectively. The block estimates the set of  $A(q)$ ,  $B(q)$ ,  $C(q)$ ,  $D(q)$ , and  $F(q)$  coefficients that the model structure uses and outputs them in the **Parameters** output. During the estimation, the block constrains the estimated  $C$ ,  $D$ , and  $F$  polynomials to a stable region with roots in the unit disk, while allowing the estimated  $A$  and  $B$  polynomials to be unstable. The **Parameters** output provides a bus signal with the following elements:

- A — Vector containing  $[1 \ a_1(t) \ \dots \ a_{na}(t)]$ .
- B — Vector containing  $[0_1 \ \dots \ 0_{nk}, \ b_1(t) \ \dots \ b_{nb}(t)]$ . For MISO data,  $B$  is a matrix where the  $i$ -th row parameters correspond to the  $i$ -th input.
- C — Vector containing  $[1 \ c_1(t) \ \dots \ c_{nc}(t)]$ .
- D — Vector containing  $[1 \ d_1(t) \ \dots \ d_{nd}(t)]$ .
- F — Vector containing  $[1 \ f_1(t) \ \dots \ f_{nf}(t)]$ .

For example, suppose that you want to estimate the coefficients for the following SISO ARMAX model:

$$y(t)+a_1y(t-1)+\dots+a_{na}y(t-na) = b_1u(t-nk)+\dots+b_{nb}u(t-nb-nk+1)+ e(t)+c_1e(t-1)+\dots+c_{nc}e(t-nc)$$

$y$ ,  $u$ ,  $na$ ,  $nb$ ,  $nc$ , and  $nk$  are known quantities that you provide to the block. For each time step,  $t$ , the block estimates the  $A$ ,  $B$ , and  $C$  parameter values, constraining only the  $C$  polynomial to a stable region. The block then outputs these estimated values using the **Parameters** output.

### Block Capabilities

The block supports several estimation methods and data input formats. The block can provide both infinite-history [1] and finite-history [2] (also known as sliding-window) estimates for  $\theta$ . Configurable options in the block include:

- Multiple inputs (ARX model structure only) — See the **Inputs** port.
- Sample-based or frame-based data format — See the **Input Processing** parameter.
- Multiple infinite-history estimation methods [1] — See the **Estimation Method** parameter.
- Infinite-history or finite-history (AR or ARX model structures only) — See the **History** parameter.
- Initial conditions, enable flag, and reset trigger — See the **Initial Estimate**, **Add enable port**, and **External Reset** parameters.

For more information on the estimation methods, see “Recursive Algorithms for Online Parameter Estimation”.

### Ports

#### Input

##### Inputs — Input signal

vector | matrix

Input signal  $u(t)$ . The **Input Processing** parameter and the number of inputs  $nu$  define the dimensions of the signal. Only the ARX model structure can be MISO, with  $nu$  greater than 1.

- Sample-based input processing and  $nu$  inputs —  $nu$ -by-1 vector
- Frame-based input processing with  $M$  samples per frame and  $nu$  inputs —  $M$ -by- $nu$  matrix

##### Dependencies

To enable this port, set the **Model Structure** parameter to ARX, ARMAX, BJ, or OE.

Data Types: single | double

**Output — Measured output signal**

scalar | vector

Measured output signal  $y(t)$ . The **Input Processing** parameter defines the dimensions of the signal:

- Sample-based input processing — Scalar
- Frame-based input processing with  $M$  samples per frame —  $M$ -by-1 vector

Data Types: single | double

**Enable — Enable estimation updates**

true (default) | false

External signal that allows you to enable and disable estimation updates. If the signal value is:

- `true` — Estimate and output the parameter values for the time step.
- `false` — Do not estimate the parameter values, and output the most recent previously estimated value.

**Dependencies**

To enable this port, select the **Add enable port** parameter.

Data Types: single | double | Boolean | int8 | int16 | int32 | int64 | uint8 | uint16 | uint32

**Reset — Reset trigger**

scalar

Reset parameter estimation to its initial conditions. The value of the **External reset** parameter determines the trigger type. The trigger type dictates whether the reset occurs on a signal that is rising, falling, either rising or falling, level, or on level hold.

**Dependencies**

To enable this port, select any option other than **None** in the **External reset** dropdown.

Data Types: single | double | Boolean | int8 | int16 | int32 | uint8 | uint16 | uint32

**InitialParameters — Initial parameter estimates**

bus object

Initial parameter estimates, supplied from a source external to the block. The block uses this inport at the beginning of the simulation or when you trigger an algorithm reset using the **Reset** signal.

For information on the contents of the **InitialParameters** bus object, see the **Parameters** port description.

### Dependencies

To enable this port, set **History** to Infinite and **Initial Estimate** to External.

Data Types: single | double

### InitialCovariance — Initial covariance of parameters

positive scalar (default) | vector of positive scalars | symmetric positive-definite matrix

Initial parameter covariances, supplied from a source external to the block. For details, see the **Parameter Covariance Matrix** parameter. The block uses this inport at the beginning of the simulation or when you trigger an algorithm reset using the **Reset** signal.

### Dependencies

To enable this port, set the following parameters:

- **History** to Infinite
- **Estimation Method** to Forgetting Factor or Kalman Filter
- **Initial Estimate** to External

Data Types: single | double

### InitialInputs — Initial values of the inputs

matrix

Initial set of inputs when using finite-history (sliding-window) estimation, supplied from an external source. The signal to this port must be a  $(W-1+\max(nb)+\max(nk))$ -by- $nu$  matrix, where  $W$  is the window length and  $nu$  is the number of inputs.  $nb$  is the vector of  $B(q)$  polynomial orders and  $nk$  is the vector of input delays.

The block uses this inport at the beginning of the simulation or whenever the **Reset** signal triggers.

If the initial buffer is set to 0 or does not contain enough information, you see a warning message during the initial phase of your estimation. The warning should clear after a few

cycles. The number of cycles it takes for sufficient information to be buffered depends upon the order of your polynomials and your input delays. If the warning persists, you should evaluate the content of your signals.

### Dependencies

To enable this port, set:

- **History** to Finite
- **Model Structure** to ARX
- **Initial Estimate** to External

Data Types: `single` | `double`

### InitialOutputs — Initial values of the measured outputs buffer

vector

Initial set of output measurements when using finite-history (sliding-window) estimation, supplied from an external source. The signal to this port must be a  $(W+na)$ -by-1 vector, where  $W$  is the window length and  $na$  is the polynomial order of  $A(q)$ .

The block uses this inport at the beginning of the simulation or whenever the **Reset** signal triggers.

If the initial buffer is set to  $\theta$  or does not contain enough information, you see a warning message during the initial phase of your estimation. The warning should clear after a few cycles. The number of cycles it takes for sufficient information to be buffered depends upon the order of your polynomials and your input delays. If the warning persists, you should evaluate the content of your signals.

### Dependencies

To enable this port, set:

- **History** to Finite
- **Model Structure** to AR or ARX
- **Initial Estimate** to External

Data Types: `single` | `double`

### Output

#### Parameters — Estimated parameters

bus object

Estimated polynomial coefficients, returned as a bus. The bus contains an element for each of the  $A$ ,  $B$ ,  $C$ ,  $D$ , and  $F$  polynomials that correspond to the structure that you specify in **Model Structure** (see “Model Structures” on page 2-94 ).

Each bus element is a vector signal containing the associated polynomial coefficients. For example, the  $A$  element contains  $[1 \ a_1(t) \ \dots \ a_{nd}(t)]$ .

Estimated  $C$ ,  $D$ , and  $F$  values are constrained to be stable discrete-time polynomials. That is, these polynomials all have roots within the unit circle. Estimated  $A$  and  $B$  polynomials are allowed to be unstable.

Data Types: `single` | `double`

#### Error — Estimation error

scalar | vector

Estimation error, returned as:

- Scalar — Sample-based input processing
- $M$ -by-1 vector — Frame-based input processing with  $M$  samples per frame

#### Dependencies

To enable this port, select the **Output estimation error** parameter.

Data Types: `single` | `double`

#### Covariance — Parameter estimation error covariance $P$

matrix

Parameter estimation error covariance  $P$ , returned as an  $N$ -by- $N$  matrix, where  $N$  is the number of parameters. For details, see the **Output Parameter Covariance Matrix** parameter.

#### Dependencies

To enable this port:

- Select the **Output parameter covariance matrix** parameter.
- If **History** is Infinite, set **Estimation Method** to Forgetting Factor or Kalman Filter.

Data Types: single | double

## Parameters

### Model Structure

Estimated model structure, specified as one of the following:

- ARX — SISO or MISO ARX model
- ARMAX — SISO ARMAX model
- OE — SISO OE model
- BJ — SISO BJ model
- AR — Time-series AR model
- ARMA — Time-series ARMA model

### Model Parameters

#### Initial Estimate — Source of initial parameter estimates

None (default) | Internal | External

Specify how to provide initial parameter estimates to the block:

- None — Do not specify initial estimates.
  - If **History** is Infinite, the block uses  $\theta$  as the initial parameter estimate.
  - If **History** is Finite, the block calculates the initial parameter estimates from the initial **Inputs** and **Outputs** signals.

Specify the parameters that the block enables based on your choice of model structure and estimation method.

- Specify the set of **Number of Parameters ()** parameters that the block enables based on your **Model Structure**. For instance, if your setting for **Model Structure** is AR, specify the **Number of Parameters in A(q) (na)** parameter.

- Specify the **Input Delay (nk)** parameter that the block enables when your model structure uses a  $B(q)$  element.
- Specify the **Parameter Covariance Matrix** if **Estimation Method** is Forgetting Factor or Kalman Filter.
- **Internal** — Specify initial parameter estimates internally to the block.
  - Specify the initial parameter values **Initial ()** parameters that the block enables based on your **Model Structure** and **History**. For instance, if your setting for **Model Structure** is AR and **History** is Infinite, specify the **Initial A(q)** parameter.
  - Specify the **Input Delay (nk)** parameter that the block enables when your model structure uses a  $B(q)$  element.
  - Specify the **Parameter Covariance Matrix** parameter if **Estimation Method** is Forgetting Factor or Kalman Filter.
  - Specify the **Initial Inputs** parameter (ARX only) and the **Initial Outputs** parameter (ARX and AR) if **History** is Finite.
- **External** — Specify initial parameter estimates as an input signal to the block.

Specify the **Number of Parameters ()** parameters that the block enables based on your **Model Structure**. Your setting for **Model Structure** and for the **History** parameter determines which signals to connect to the relevant ports:

- If **History** is Infinite — **InitialParameters** and **InitialCovariance**
- If **History** is Finite — **InitialOutputs**, and, for the ARX model structure, also **InitialInputs**

### Programmatic Use

**Block Parameter:** InitialEstimateSource

**Type:** character vector, string

**Values:** 'None', 'Internal', 'External'

**Default:** 'None'

### Number of Parameters in $A(q)$ (na) — Number of estimated parameters in the $A(q)$ polynomial

1 (default) | non-negative integer

Specify the order  $na$  of the  $A(q)$  polynomial. This order is equivalent to the number of parameters in the polynomial that you are estimating.

**Dependencies**

To enable this parameter, either:

- Set **History** to Infinite, **Model Structure** to AR, ARX, ARMA, or ARMAX, and **Initial Estimate** to None or External.
- Set **History** to Finite and **Model Structure** to AR or ARX.

**Programmatic Use**

**Block Parameter:** A0

**Type:** non-negative integer

**Default:** 1

**Number of Parameters in B(q) (nb) – Number of estimated parameters in the B(q) polynomial**

1 (default) | vector of positive integers

Specify the order  $nb$  of the  $B(q)$  polynomial. This order is equivalent to the number of parameters minus one in the polynomial that you are estimating.

For MISO systems using an ARX model structure, specify  $nb$  as a vector with as many elements as there are inputs. Each element of this vector represents the order of the  $B(q)$  polynomial associated with the corresponding input. For example, suppose that you have

a two-input MISO system whose  $B(q)$  elements are: 
$$\begin{bmatrix} B_1 \\ B_2 \end{bmatrix} = \begin{bmatrix} 0 & b_{11} & 0 \\ 0 & b_{21} & b_{22} \end{bmatrix}$$
. The zero at the beginning of each polynomial represents a single input delay for each input (see the **Initial B(q)** parameter description). The order  $nb$  for each polynomial is equal to the number of elements following the initial zero, or 1 for input 1 and 2 for input 2. Specify **Number of Parameters in B(q) (nb)** as [1 2], and **Input Delay (nk)** as [1 1].

**Dependencies**

To enable this parameter, either:

- Set **History** to Infinite, **Model Structure** to ARX, ARMAX, BJ, or OE, and **Initial Estimate** to None or External.
- Set **History** to Finite with **Model Structure** of ARX.

**Programmatic Use**

**Block Parameter:** B0

**Type:** positive integer

**Default:** 1

**Number of Parameters in  $C(q)$  ( $nc$ ) — Number of estimated parameters in the  $C(q)$  polynomial**

1 (default) | positive integer

Specify the order  $nc$  of the  $C(q)$  polynomial. This order is equivalent to the number of parameters in the polynomial that you are estimating.

**Dependencies**

To enable this parameter, set **History** to Infinite, **Model Structure** to ARMA, ARMAX, or BJ, and **Initial Estimate** to None or External

**Programmatic Use**

**Block Parameter:** C0

**Type:** positive integer

**Default:** 1

**Number of Parameters in  $D(q)$  ( $nd$ ) — Number of estimated parameters in the  $D(q)$  polynomial**

1 (default) | positive integer

Specify the order  $nd$  of the  $D(q)$  polynomial. This order is equivalent to the number of parameters in the polynomial that you are estimating.

**Dependencies**

To enable this parameter, set **History** to Infinite, **Model Structure** to BJ, and **Initial Estimate** to None or External.

**Programmatic Use**

**Block Parameter:** D0

**Type:** positive integer

**Default:** 1

**Number of Parameters in  $F(q)$  ( $nf$ ) — Number of estimated parameters in the  $F(q)$  polynomial**

1 (default) | positive integer

Specify the order  $nf$  of the  $F(q)$  polynomial. This order is equivalent to the number of parameters in the polynomial that you are estimating.

**Dependencies**

To enable this parameter, set **History** to Infinite, **Model Structure** to OE or to BJ, and **Initial Estimate** to None or External.

**Programmatic Use**

**Block Parameter:** F0

**Type:** positive integer

**Default:** 1

**Input Delay (nk) — Input-output delay**

1 (default) | vector of non-negative integers

Specify the input-output delay as an integer representing the number of input samples that occur before the input affects the output. This delay is also called the *dead time* in the system. The block encodes the input delay as fixed leading zeros of the  $B(q)$  polynomial. For more information, see the  $B(q)$  parameter description.

For MISO systems with ARX model structure, specify  $nk$  as a vector with elements specifying the delay for each input. This vector is of length  $nu$ , where  $nu$  is the number of inputs.

For example, suppose that you have a two-input MISO system whose  $B(q)$  elements are:

$$\begin{bmatrix} B_1 \\ B_2 \end{bmatrix} = \begin{bmatrix} 0 & b_{11} & b_{12} \\ 0 & 0 & b_{21} \end{bmatrix}$$
. The zeros at the beginning of the polynomials represent a single-sample delay for the first input, and a two-sample delay for the second input (see the **Initial B(q)** parameter description). The order  $nb$  for each polynomial is equal to the number of elements following the initial zero, or 2 for input 1 and 1 for input 2. Specify **Input Delay (nk)** as [1 1], and **Number of Parameters in B(q) (nb)** as [21].

**Dependencies**

To enable this parameter, either:

- Set **History** to Infinite, **Model Structure** to ARX, ARMAX, OE, or BJ, and **Initial Estimate** to None or External.
- Set **History** to Finite, **Model Structure** to ARX, and **Initial Estimate** to None or External.

**Programmatic Use**

**Block Parameter:** nk

**Type:** non-negative integer vector

**Default:** 1

### Parameter Covariance Matrix — Initial parameter covariance

1e4 (default) | scalar | vector | matrix

- Real positive scalar,  $\alpha$  — Covariance matrix is an  $N$ -by- $N$  diagonal matrix, with  $\alpha$  as the diagonal elements.
- Vector of real positive scalars,  $[\alpha(a), \alpha(b), \alpha(c), \alpha(d), \alpha(f)]$  — Covariance matrix is an  $N$ -by- $N$  diagonal matrix, with  $[\alpha(a), \alpha(b), \alpha(c), \alpha(d), \alpha(f)]$  as the diagonal elements.  $\alpha(a)$  is a vector of the covariance for each coefficient of the  $A$  polynomial. Similarly,  $\alpha(b)$ ,  $\alpha(c)$ ,  $\alpha(d)$  and  $\alpha(f)$  are vectors containing the covariance of the coefficients of the  $B$ ,  $C$ ,  $D$  and  $F$  polynomials, respectively.
- $N$ -by- $N$  symmetric positive-definite matrix.

$N$  can be one of the following:

- AR —  $N = na$

- 

$$\text{ARX} \text{ — } N = na + \sum_{i=1}^N nb_i$$

- ARMA —  $N = na + nc$
- ARMAX —  $N = na + nb + nc$
- OE —  $N = nb + nf$
- BJ —  $N = nb + nc + nd + nf$

### Dependencies

To enable this parameter, set

- **History** to Infinite
- **Initial Estimate** to None or Internal
- **Estimation Method** to Forgetting Factor or Kalman Filter

The block uses this parameter at the beginning of the simulation or whenever the **Reset** signal triggers.

**Programmatic Use****Block Parameter:**  $P_0$ **Type:** scalar, vector, or matrix**Default:**  $1e4$ **Initial  $A(q)$  — Initial values of the  $A(q)$  polynomial** $[1 \text{ eps}]$  (default) | vector

Specify the initial estimate of the  $A(q)$  values as a row vector of length  $na+1$ .

The leading coefficient of  $A$  must be 1.

**Dependencies**

To enable this parameter, set:

- **History** to Infinite
- **Model Structure** to AR, ARX, ARMA, or ARMAX
- **Initial Estimate** to Internal

**Programmatic Use****Block Parameter:**  $A_0$ **Type:** real vector**Default:**  $[1 \text{ eps}]$ **Initial  $B(q)$  — Initial values of the  $B(q)$  polynomial** $[0 \text{ eps}]$  (default) | vector | matrix

Specify the initial estimate of the  $B(q)$  values as a row vector of length  $nb+nk$ . For multiple-input models, specify a matrix where the  $i$ th row corresponds to the  $i$ th input.

The block counts the leading zeros in  $B(q)$  and interprets them as input delay  $nk$ . Those zeros are fixed throughout the estimation.  $nb$  is the number of elements after the first nonzero element in  $B(q)$ . The block estimates the value of these  $nb$  elements.

For example:

- $[0 \text{ eps}]$  corresponds to  $nk=1$  and  $nb=1$ .
- $[0 \ 0 \ \text{eps}]$  corresponds to  $nk=2$  and  $nb=1$ .
- $[0 \ 0 \ \text{eps} \ 0 \ \text{eps}]$  corresponds to  $nk=2$  and  $nb=3$ .

The default value is  $[0 \ \text{eps}]$ .

### Dependencies

To enable this parameter, set:

- **History** to Infinite
- **Model Structure** to ARX, ARMAX, OE, or BJ
- **Initial Estimate** to Internal

### Programmatic Use

**Block Parameter:** B0

**Type:** real vector or matrix

**Default:** [0 eps]

### Initial $C(q)$ — Initial values of the $C(q)$ polynomial

[1 eps] (default) | vector

Specify the initial estimate of the  $C(q)$  values as a row vector of length  $nc+1$ .

The leading coefficient of  $C(q)$  must be 1.

The coefficients must define a stable discrete-time polynomial, that is, have all polynomial roots within the unit circle.

### Dependencies

To enable this parameter, set:

- **History** to Infinite
- **Model Structure** to ARMA, ARMAX, or BJ
- **Initial Estimate** to Internal

### Programmatic Use

**Block Parameter:** C0

**Type:** real vector

**Default:** [1 eps]

### Initial $D(q)$ — Initial values of the $D(q)$ polynomial

[1 eps] (default) | vector

Specify the initial estimate of the  $D(q)$  values as a row vector of length  $nd+1$ .

The leading coefficient of  $D(q)$  must be 1.

The coefficients must define a stable discrete-time polynomial, that is, have all polynomial roots within the unit circle.

### Dependencies

To enable this parameter, set:

- **History** to Infinite
- **Model Structure** to BJ
- **Initial Estimate** to Internal

### Programmatic Use

**Block Parameter:** D0

**Type:** real vector

**Default:** [1 eps]

### Initial $F(q)$ — Initial values of the $F(q)$ polynomial

[1 eps] (default) | vector

Specify the initial estimate of the  $F(q)$  values as a row vector of length  $nf+1$ .

The leading coefficient of  $F(q)$  must be 1.

The coefficients must define a stable discrete-time polynomial, that is, have all polynomial roots within the unit circle.

### Dependencies

To enable this parameter, set:

- **History** to Infinite
- **Model Structure** to OE or to BJ
- **Initial Estimate** to Internal

### Programmatic Use

**Block Parameter:** F0

**Type:** real vector

**Default:** [1 eps]

### Initial Inputs — Initial values of the inputs buffer

0 (default) | matrix

Specify initial values of the inputs buffer when using finite-history (sliding window) estimation. The buffer dimensions accommodate the specified window length, the regressors associated with polynomials within that window, the input delays, and the number of inputs. These elements drive a matrix size of  $(W-1+\max(nb)+\max(nk))$ -by- $nu$  matrix, where  $W$  is the window length and  $nu$  is the number of inputs.  $nb$  is the vector of  $B(q)$  polynomial orders and  $nk$  is the vector of input delays.

When the initial value is set to  $\mathbf{0}$ , the block populates the buffer with zeros.

If the initial buffer is set to  $\mathbf{0}$  or does not contain enough information, you see a warning message during the initial phase of your estimation. The warning should clear after a few cycles. The number of cycles it takes for sufficient information to be buffered depends upon the order of your polynomials and your input delays. If the warning persists, you should evaluate the content of your signals.

The block uses this parameter at the beginning of the simulation or whenever the **Reset** signal triggers.

### Dependencies

To enable this parameter, set

- **History** to Finite
- **Model Structure** to ARX
- **Initial Estimate** to Internal

.

### Programmatic Use

**Block Parameter:** InitialInputs

**Type:** real matrix

**Default:**  $\mathbf{0}$

### Initial Outputs — Initial values of the measured outputs buffer

$\mathbf{0}$  (default) | vector

Specify initial values of the measured outputs buffer when using finite-history (sliding-window) estimation. The buffer dimensions accommodate the specified window length and the regressors associated with the polynomials within that window. This parameter is a  $(W+na)$ -by-1 vector, where  $W$  is the window length and  $na$  is the polynomial order of  $A(q)$ .

When the initial value is set to  $\mathbf{0}$ , the block populates the buffer with zeros.

If the initial buffer is set to  $\emptyset$  or does not contain enough information, you see a warning message during the initial phase of your estimation. The warning should clear after a few cycles. The number of cycles it takes for sufficient information to be buffered depends upon the order of your polynomials and your input delays. If the warning persists, you should evaluate the content of your signals.

The block uses this parameter at the beginning of the simulation or whenever the **Reset** signal triggers.

### Dependencies

To enable this parameter, set **History** to Finite and **Initial Estimate** to Internal.

### Programmatic Use

**Block Parameter:** InitialOutputs

**Type:** real vector

**Default:**  $\emptyset$

### Input Processing and Sample Time

#### Input Processing — Choose sample-based or frame-based processing

Sample-based (default) | Frame-based

- Sample-based processing operates on signals streamed one sample at a time.
- Frame-based processing operates on signals containing samples from multiple time steps. Many machine sensor interfaces package multiple samples and transmit these samples together in frames. Frame-based processing allows you to input this data directly without having to first unpack it.

Specifying frame-based data adds an extra dimension of  $M$  to some of your data inports and outports, where  $M$  is the number of time steps in a frame. These ports are:

- **Inputs**
- **Output**
- **Error**

For more information, see the port descriptions in “Ports” on page 2-96.

### Programmatic Use

**Block Parameter:** InputProcessing

**Type:** character vector, string

**Values:** 'Sample-based', 'Frame-based'

**Default:** 'Sample-based'

### **Sample Time — Block sample time**

-1 (default) | positive scalar

Specify the data sample time, whether by individual samples for sample-based processing ( $t_s$ ), or by frames for frame-based processing ( $t_f = Mt_s$ ), where  $M$  is the frame length. When you set **Sample Time** to its default value of -1, the block inherits its  $t_s$  or  $t_f$  based on the signal.

Specify **Sample Time** as a positive scalar to override the inheritance.

#### **Programmatic Use**

**Block Parameter:** Ts

**Type:** real scalar

**Values:** -1, positive scalar

**Default:** -1

## **Algorithm and Block Options**

### **Algorithm Options**

#### **History — Choose infinite or finite data history**

Infinite (default) | Finite

The **History** parameter determines what type of recursive algorithm you use:

- **Infinite** — Algorithms in this category aim to produce parameter estimates that explain all data since the start of the simulation. These algorithms retain the history in a data summary. The block maintains this summary within a fixed amount of memory that does not grow over time.

The block provides multiple algorithms of the **Infinite** type. Selecting this option enables the **Estimation Method** parameter with which you specify the algorithm.

- **Finite** — Algorithms in this category aim to produce parameter estimates that explain only a finite number of past data samples. The block uses all of the data within a finite window, and discards data once that data is no longer within the window bounds. This method is also called sliding-window estimation.

The block provides one algorithm of the **Finite** type. You can use this option only with the AR and ARX model structures.

Selecting this option enables the **Window Length** parameter.

For more information on recursive estimation methods, see “Recursive Algorithms for Online Parameter Estimation”

**Programmatic Use**

**Block Parameter:** History

**Type:** character vector, string

**Values:** 'Infinite', 'Finite'

**Default:** 'Infinite'

**Window Length — Window size for finite sliding-window estimation**

200 (default) | positive integer

The **Window Length** parameter determines the number of time samples to use for the finite-history (sliding-window) estimation method. Choose a window size that balances estimation performance with computational and memory burden. Sizing factors include the number and time variance of the parameters in your model. Always specify **Window Length** in samples, even if you are using frame-based input processing.

**Window Length** must be greater than or equal to the number of estimated parameters.

Suitable window length is independent of whether you are using sample-based or frame-based input processing. However, when using frame-based processing, **Window Length** must be greater than or equal to the number of samples (time steps) contained in the frame.

**Dependencies**

To enable this parameter, set **History** to Finite.

**Programmatic Use**

**Block Parameter:** WindowLength

**Type:** positive integer

**Default:** 200

**Estimation Method — Recursive estimation algorithm**

Forgetting Factor (default) | Kalman Filter | Normalized Gradient | Gradient

Specify the estimation algorithm when performing infinite-history estimation. When you select any of these methods, the block enables additional related parameters.

Forgetting factor and Kalman filter algorithms are more computationally intensive than gradient and normalized gradient methods. However, these more intensive methods have

better convergence properties than the gradient methods. For more information about these algorithms, see “Recursive Algorithms for Online Parameter Estimation”.

### **Programmatic Use**

**Block Parameter:** EstimationMethod

**Type:** character vector, string

**Values:** 'Forgetting Factor', 'Kalman Filter', 'Normalized Gradient', 'Gradient'

**Default:** 'Forgetting Factor'

### **Forgetting Factor — Discount old data using forgetting factor**

1 (default) | positive scalar in (0 1] range

The forgetting factor  $\lambda$  specifies if and how much old data is discounted in the estimation. Suppose that the system remains approximately constant over  $T_0$  samples. You can choose  $\lambda$  such that:

$$T_0 = \frac{1}{1 - \lambda}$$

- Setting  $\lambda = 1$  corresponds to “no forgetting” and estimating constant coefficients.
- Setting  $\lambda < 1$  implies that past measurements are less significant for parameter estimation and can be “forgotten.” Set  $\lambda < 1$  to estimate time-varying coefficients.

Typical choices of  $\lambda$  are in the [0.98 0.995] range.

### **Dependencies**

To enable this parameter, set **History** to Infinite and **Estimation Method** to Forgetting Factor.

### **Programmatic Use**

**Block Parameter:** AdaptationParameter

**Type:** scalar

**Values:** (0 1] range

**Default:** 1

### **Process Noise Covariance — Process noise covariance for Kalman filter estimation method**

1 (default) | nonnegative scalar | vector of nonnegative scalars | symmetric positive semidefinite matrix

**Process Noise Covariance** prescribes the elements and structure of the noise covariance matrix for the Kalman filter estimation. Using  $N$  as the number of parameters to estimate, specify the **Process Noise Covariance** as one of the following:

- Real nonnegative scalar,  $\alpha$  — Covariance matrix is an  $N$ -by- $N$  diagonal matrix, with  $\alpha$  as the diagonal elements.
- Vector of real nonnegative scalars,  $[\alpha_1, \dots, \alpha_N]$  — Covariance matrix is an  $N$ -by- $N$  diagonal matrix, with  $[\alpha_1, \dots, \alpha_N]$  as the diagonal elements.
- $N$ -by- $N$  symmetric positive semidefinite matrix.

The Kalman filter algorithm treats the parameters as states of a dynamic system and estimates these parameters using a Kalman filter. **Process Noise Covariance** is the covariance of the process noise acting on these parameters. Zero values in the noise covariance matrix correspond to constant coefficients, or parameters. Values larger than 0 correspond to time-varying parameters. Use large values for rapidly changing parameters. However, expect the larger values to result in noisier parameter estimates. The default value is 1.

#### Dependencies

To enable this parameter, set **History** to Infinite and **Estimation Method** to Kalman Filter.

#### Programmatic Use

**Block Parameter:** AdaptationParameter

**Type:** scalar, vector, matrix

**Default:** 1

#### Adaptation Gain — Adaptation gain specification for gradient estimation methods

1 (default) | positive scalar

The adaptation gain  $\gamma$  scales the influence of new measurement data on the estimation results for the gradient and normalized gradient methods. When your measurements are trustworthy, or in other words have a high signal-to-noise ratio, specify a larger value for  $\gamma$ . However, setting  $\gamma$  too high can cause the parameter estimates to diverge. This divergence is possible even if the measurements are noise free.

When **Estimation Method** is NormalizedGradient, **Adaptation Gain** should be less than 2. With either gradient method, if errors are growing in time (in other words, estimation is diverging), or parameter estimates are jumping around frequently, consider reducing **Adaptation Gain**.

### Dependencies

To enable this parameter, set **History** to Infinite and **Estimation Method** to Normalized Gradient or to Gradient.

### Programmatic Use

**Block Parameter:** AdaptationParameter

**Type:** scalar

**Default:** 1

### Normalization Bias — Bias for adaptation gain scaling for normalized gradient estimation method

eps (default) | nonnegative scalar

The normalized gradient algorithm scales the adaptation gain at each step by the square of the two-norm of the gradient vector. If the gradient is close to zero, the near-zero denominator can cause jumps in the estimated parameters. **Normalization Bias** is the term introduced to the denominator to prevent these jumps. Increase **Normalization Bias** if you observe jumps in estimated parameters.

### Dependencies

To enable this parameter, set **History** to Infinite and **Estimation Method** to Normalized Gradient.

### Programmatic Use

**Block Parameter:** NormalizationBias

**Type:** scalar

**Default:** eps

### Block Options

#### Output estimation error — Add Error output to block

off (default) | on

Use the **Error** output signal to validate the estimation. For a given time step  $t$ , the estimation error  $e(t)$  is calculated as:

$$e(t) = y(t) - y_{est}(t),$$

where  $y(t)$  is the measured output that you provide, and  $y_{est}(t)$  is the estimated output using the regressors  $H(t)$  and parameter estimates  $\theta(t-1)$ .

**Programmatic Use****Block Parameter:** OutputError**Type:** character vector, string**Values:** 'off','on',**Default:** 'off'**Output parameter covariance matrix — Add covariance output to block**

off (default) | on

Use the **Covariance** output signal to examine parameter estimation uncertainty. The software computes parameter covariance  $P$  assuming that the residuals,  $e(t)$ , are white noise, and the variance of these residuals is 1.

The interpretation of  $P$  depends on the estimation approach you specify in **History** and **Estimation Method** as follows:

- If **History** is Infinite, then your **Estimation Method** selection results in:
  - Forgetting Factor —  $(R_2/2)P$  is approximately equal to the covariance matrix of the estimated parameters, where  $R_2$  is the true variance of the residuals. The block returns these residuals through the **Error** port.
  - Kalman Filter —  $R_2P$  is the covariance matrix of the estimated parameters, and  $R_1/R_2$  is the covariance matrix of the parameter changes. Here,  $R_1$  is the covariance matrix that you specify in **Parameter Covariance Matrix**.
  - Normalized Gradient or Gradient — Covariance  $P$  is not available.
- If **History** is Finite (sliding-window estimation) —  $R_2P$  is the covariance of the estimated parameters. The sliding-window algorithm does not use this covariance in the parameter-estimation process. However, the algorithm does compute the covariance for output so that you can use it for statistical evaluation.

**Programmatic Use****Block Parameter:** OutputP**Type:** character vector, string**Values:** 'off','on'**Default:** 'off'**Add enable port — Add Enable inport to block**

off (default) | on

Use the **Enable** signal to provide a control signal that enables or disables parameter estimation. The block estimates the parameter values for each time step that parameter

estimation is enabled. If you disable parameter estimation at a given step,  $t$ , then the software does not update the parameters for that time step. Instead, the block output contains the last estimated parameter values.

You can use this option, for example, when or if:

- Your regressors or output signal become too noisy, or do not contain information at some time steps
- Your system enters a mode where the parameter values do not change in time

### Programmatic Use

**Block Parameter:** AddEnablePort

**Type:** character vector, string

**Values:** 'off', 'on'

**Default:** 'off'

### External reset — Specify trigger for external reset

None (default) | Rising | Falling | Either | Level | Level hold

Set the **External reset** parameter to both add a **Reset** inport and specify the inport signal condition that triggers a reset of algorithm states to their specified initial values. Reset the estimation, for example, if parameter covariance is becoming too large because of lack of either sufficient excitation or information in the measured signals. The **External reset** parameter determines the timing for the reset.

Suppose that you reset the block at a time step,  $t$ . If the block is enabled at  $t$ , the software uses the initial parameter values specified in **Initial Estimate** to estimate the parameter values. In other words, at  $t$ , the block performs a parameter update using the initial estimate and the current values of the inports.

If the block is disabled at  $t$  and you reset the block, the block output contains the values specified in **Initial Estimate**.

Specify this option as one of the following:

- **None** — Algorithm states and estimated parameters are not reset.
- **Rising** — Trigger reset when the control signal rises from a negative or zero value to a positive value. If the initial value is negative, rising to zero triggers reset.
- **Falling** — Trigger reset when the control signal falls from a positive or a zero value to a negative value. If the initial value is positive, falling to zero triggers reset.

- **Either** — Trigger reset when the control signal is either rising or falling.
- **Level** — Trigger reset in either of these cases:
  - Control signal is nonzero at the current time step.
  - Control signal changes from nonzero at the previous time step to zero at the current time step.
- **Level hold** — Trigger reset when the control signal is nonzero at the current time step.

When you choose any option other than **None**, the software adds a **Reset** inport to the block. You provide the reset control input signal to this inport.

**Programmatic Use****Block Parameter:** ExternalReset**Type:** character vector, string**Values:** 'None', 'Rising', 'Falling', 'Either', 'Level', 'Level hold'**Default:** 'None'**References**

- [1] Ljung, L. *System Identification: Theory for the User*. Upper Saddle River, NJ: Prentice-Hall PTR, 1999, pp. 363–369.
- [2] Zhang, Q. "Some Implementation Aspects of Sliding Window Least Squares Algorithms." *IFAC Proceedings*. Vol. 33, Issue 15, 2000, pp. 763-768.

**Extended Capabilities****C/C++ Code Generation**

Generate C and C++ code using Simulink® Coder™.

**PLC Code Generation**

Generate Structured Text code using Simulink® PLC Coder™.

## **See Also**

Kalman Filter | Recursive Polynomial Model Estimator

## **Topics**

*“Estimate Parameters of System Using Simulink Recursive Estimator Block”*

*“Online Recursive Least Squares Estimation”*

*“Preprocess Online Parameter Estimation Data in Simulink”*

*“Validate Online Parameter Estimation Results in Simulink”*

*“Generate Online Parameter Estimation Code in Simulink”*

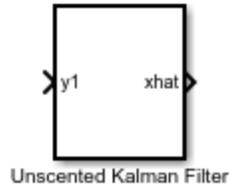
*“Recursive Algorithms for Online Parameter Estimation”*

**Introduced in R2014a**

# Unscented Kalman Filter

Estimate states of discrete-time nonlinear system using unscented Kalman filter

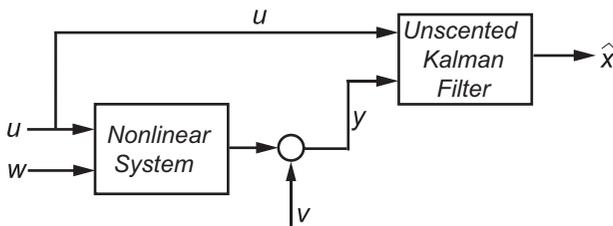
**Library:** Control System Toolbox / State Estimation  
System Identification Toolbox / Estimators



## Description

The Unscented Kalman Filter block estimates the states of a discrete-time nonlinear system using the discrete-time unscented Kalman filter algorithm.

Consider a plant with states  $x$ , input  $u$ , output  $y$ , process noise  $w$ , and measurement noise  $v$ . Assume that you can represent the plant as a nonlinear system.



Using the state transition and measurement functions of the system and the unscented Kalman filter algorithm, the block produces state estimates  $\hat{x}$  for the current time step. For information about the algorithm, see “Extended and Unscented Kalman Filter Algorithms for Online State Estimation”.

You create the nonlinear state transition function and measurement functions for the system and specify these functions in the block. The block supports state estimation of a system with multiple sensors that are operating at different sampling rates. You can specify up to five measurement functions, each corresponding to a sensor in the system. For more information, see “State Transition and Measurement Functions” on page 2-137.

## Ports

### Input

#### **y1, y2, y3, y4, y5 — Measured system outputs**

vector

Measured system outputs corresponding to each measurement function that you specify in the block. The number of ports equals the number of measurement functions in your system. You can specify up to five measurement functions. For example, if your system has two sensors, you specify two measurement functions in the block. The first port **y1** is available by default. When you click **Apply**, the software generates port **y2** corresponding to the second measurement function.

Specify the ports as  $N$ -dimensional vectors, where  $N$  is the number of quantities measured by the corresponding sensor. For example, if your system has one sensor that measures the position and velocity of an object, then there is only one port **y1**. The port is specified as a 2-dimensional vector with values corresponding to position and velocity.

#### **Dependencies**

The first port **y1** is available by default. Ports **y2** to **y5** are generated when you click **Add Measurement**, and click **Apply**.

Data Types: `single` | `double`

#### **StateTransitionFcnInputs — Additional optional input argument to state transition function**

scalar | vector | matrix

Additional optional input argument to the state transition function  $f$  other than the state  $x$  and process noise  $w$ . For information about state transition functions see, “State Transition and Measurement Functions” on page 2-137.

Suppose that your system has nonadditive process noise, and the state transition function  $f$  has the following form:

$$x(k+1) = f(x(k), w(k), \text{StateTransitionFcnInputs}).$$

Here  $k$  is the time step, and `StateTransitionFcnInputs` is an additional input argument other than  $x$  and  $w$ .

If you create  $f$  using a MATLAB function (.m file), the software generates the port **StateTransitionFcnInputs** when you click **Apply**. You can specify the inputs to this port as a scalar, vector, or matrix.

If your state transition function has more than one additional input, use a Simulink Function block to specify the function. When you use a Simulink Function block, you provide the additional inputs directly to the Simulink Function block using Inport blocks. No input ports are generated for the additional inputs in the Unscented Kalman Filter block.

### Dependencies

This port is generated only if both of the following conditions are satisfied:

- You specify  $f$  in **Function** using a MATLAB function, and  $f$  is on the MATLAB path.
- $f$  requires only one additional input argument apart from  $x$  and  $w$ .

Data Types: `single` | `double`

### **MeasurementFcn1Inputs, MeasurementFcn2Inputs, MeasurementFcn3Inputs, MeasurementFcn4Inputs, MeasurementFcn5Inputs** — Additional optional input argument to each measurement function

`scalar` | `vector` | `matrix`

Additional optional inputs to the measurement functions other than the state  $x$  and measurement noise  $v$ . For information about measurement functions see, “State Transition and Measurement Functions” on page 2-137.

**MeasurementFcn1Inputs** corresponds to the first measurement function that you specify, and so on. For example, suppose that your system has three sensors and nonadditive measurement noise, and the three measurement functions  $h_1$ ,  $h_2$ , and  $h_3$  have the following form:

$$y_1[k] = h_1(x[k], v[k], \text{MeasurementFcn1Inputs})$$

$$y_2[k] = h_2(x[k], v[k], \text{MeasurementFcn2Inputs})$$

$$y_3[k] = h_3(x[k], v[k])$$

Here  $k$  is the time step, and **MeasurementFcn1Inputs** and **MeasurementFcn2Inputs** are the additional input arguments to  $h_1$  and  $h_2$ .

If you specify  $h_1$ ,  $h_2$ , and  $h_3$  using MATLAB functions (.m files) in **Function**, the software generates ports **MeasurementFcn1Inputs** and **MeasurementFcn2Inputs** when you click **Apply**. You can specify the inputs to these ports as scalars, vectors, or matrices.

If your measurement functions have more than one additional input, use Simulink Function blocks to specify the functions. When you use a Simulink Function block, you provide the additional inputs directly to the Simulink Function block using Inport blocks. No input ports are generated for the additional inputs in the Unscented Kalman Filter block.

### Dependencies

A port corresponding to a measurement function  $h$  is generated only if both of the following conditions are satisfied:

- You specify  $h$  in **Function** using a MATLAB function, and  $h$  is on the MATLAB path.
- $h$  requires only one additional input argument apart from  $x$  and  $v$ .

Data Types: `single` | `double`

### Q — Time-varying process noise covariance

`scalar` | `vector` | `matrix`

Time-varying process noise covariance, specified as a scalar, vector, or matrix depending on the value of the **Process noise** parameter:

- **Process noise** is **Additive** — Specify the covariance as a scalar, an  $N_s$ -element vector, or an  $N_s$ -by- $N_s$  matrix, where  $N_s$  is the number of states of the system. Specify a scalar if there is no cross-correlation between process noise terms, and all the terms have the same variance. Specify a vector of length  $N_s$ , if there is no cross-correlation between process noise terms, but all the terms have different variances.
- **Process noise** is **Nonadditive** — Specify the covariance as a  $W$ -by- $W$  matrix, where  $W$  is the number of process noise terms in the state transition function.

### Dependencies

This port is generated if you specify the process noise covariance as **Time-Varying**. The port appears when you click **Apply**.

Data Types: `single` | `double`

### R1, R2, R3, R4, R5 — Time-varying measurement noise covariance

`matrix`

Time-varying measurement noise covariances for up to five measurement functions of the system, specified as matrices. The sizes of the matrices depend on the value of the **Measurement noise** parameter for the corresponding measurement function:

- **Measurement noise** is **Additive** — Specify the covariance as an  $N$ -by- $N$  matrix, where  $N$  is the number of measurements of the system.
- **Measurement noise** is **Nonadditive** — Specify the covariance as a  $V$ -by- $V$  matrix, where  $V$  is the number of measurement noise terms in the corresponding measurement function.

### Dependencies

A port is generated if you specify the measurement noise covariance as **Time-Varying** for the corresponding measurement function. The port appears when you click **Apply**.

Data Types: `single` | `double`

### **Enable1, Enable2, Enable3, Enable4, Enable5** — Enable correction of estimated states when measured data is available

scalar

Suppose that measured output data is not available at all time points at the port **y1** that corresponds to the first measurement function. Use a signal value other than  $\emptyset$  at the **Enable1** port to enable the correction of estimated states when measured data is available. Specify the port value as  $\emptyset$  when measured data is not available. Similarly, if measured output data is not available at all time points at the port **yi** for the  $i^{\text{th}}$  measurement function, specify the corresponding port **Enablei** as a value other than  $\emptyset$ .

### Dependencies

A port corresponding to a measurement function is generated if you select **Add Enable port** for that measurement function. The port appears when you click **Apply**.

Data Types: `single` | `double` | `Boolean`

## Output

### **xhat** — Estimated states

vector

Estimated states, returned as a vector of size  $N_s$ , where  $N_s$  is the number of states of the system. To access the individual states, use the Selector block.

When the **Use the current measurements to improve state estimates** parameter is selected, the block outputs the corrected state estimate  $\hat{x}[k|k]$  at time step  $k$ , estimated using measured outputs until time  $k$ . If you clear this parameter, the block returns the predicted state estimate  $\hat{x}[k|k-1]$  for time  $k$ , estimated using measured output until a previous time  $k-1$ . Clear this parameter if your filter is in a feedback loop and there is an algebraic loop in your Simulink model.

Data Types: `single` | `double`

### **P — State estimation error covariance**

matrix

State estimation error covariance, returned as an  $N_s$ -by- $N_s$  matrix, where  $N_s$  is the number of states of the system. To access the individual covariances, use the Selector block.

#### **Dependencies**

This port is generated if you select **Output state estimation covariance** in the **System Model** tab, and click **Apply**.

Data Types: `single` | `double`

## Parameters

### System Model Tab

#### **State Transition**

##### **Function — State transition function name**

`myStateTransitionFcn` (default) | function name

The state transition function calculates the  $N_s$ -element state vector of the system at time step  $k+1$ , given the state vector at time step  $k$ .  $N_s$  is the number of states of the nonlinear system. You create the state transition function and specify the function name in **Function**. For example, if `vdpStateFcn.m` is the state transition function that you created and saved, specify **Function** as `vdpStateFcn`.

The inputs to the function you create depend on whether you specify the process noise as additive or nonadditive in **Process noise**.

- **Process noise is Additive** — The state transition function  $f$  specifies how the states evolve as a function of state values at previous time step:

$$x(k+1) = f(x(k), Us1(k), \dots, Usn(k)),$$

where  $x(k)$  is the estimated state at time  $k$ , and  $Us1, \dots, Usn$  are any additional input arguments required by your state transition function, such as system inputs or the sample time. To see an example of a state transition function with additive process noise, type `edit vdpStateFcn` at the command line.

- **Process noise is Nonadditive** — The state transition function also specifies how the states evolve as a function of the process noise  $w$ :

$$x(k+1) = f(x(k), w(k), Us1(k), \dots, Usn(k)).$$

For more information, see “State Transition and Measurement Functions” on page 2-137.

You can create  $f$  using a Simulink Function block or as a MATLAB function (.m file).

- You can use a MATLAB function only if  $f$  has one additional input argument  $Us1$  other than  $x$  and  $w$ .

$$x(k+1) = f(x(k), w(k), Us1(k))$$

The software generates an additional input port **StateTransitionFcnInputs** to specify this argument.

- If you are using a Simulink Function block, specify  $x$  and  $w$  using Argument Inport blocks and the additional inputs  $Us1, \dots, Usn$  using Inport blocks in the Simulink Function block. You do not provide  $Us1, \dots, Usn$  to the Unscented Kalman Filter block.

### Programmatic Use

**Block Parameter:** StateTransitionFcn

**Type:** character vector, string

**Default:** 'myStateTransitionFcn'

### Process noise — Process noise characteristics

Additive (default) | Nonadditive

Process noise characteristics, specified as one of the following values:

- **Additive** — Process noise  $w$  is additive, and the state transition function  $f$  that you specify in **Function** has the following form:

$$x(k+1) = f(x(k), Us1(k), \dots, Usn(k)),$$

where  $x(k)$  is the estimated state at time  $k$ , and  $Us1, \dots, Usn$  are any additional input arguments required by your state transition function.

- **Nonadditive** — Process noise is nonadditive, and the state transition function specifies how the states evolve as a function of the state *and* process noise at the previous time step:

$$x(k+1) = f(x(k), w(k), Us1(k), \dots, Usn(k)).$$

### Programmatic Use

**Block Parameter:** HasAdditiveProcessNoise

**Type:** character vector

**Values:** 'Additive', 'Nonadditive'

**Default:** 'Additive'

### Covariance — Time-invariant process noise covariance

1 (default) | scalar | vector | matrix

Time-invariant process noise covariance, specified as a scalar, vector, or matrix depending on the value of the **Process noise** parameter:

- **Process noise** is **Additive** — Specify the covariance as a scalar, an  $Ns$ -element vector, or an  $Ns$ -by- $Ns$  matrix, where  $Ns$  is the number of states of the system. Specify a scalar if there is no cross-correlation between process noise terms and all the terms have the same variance. Specify a vector of length  $Ns$ , if there is no cross-correlation between process noise terms but all the terms have different variances.
- **Process noise** is **Nonadditive** — Specify the covariance as a  $W$ -by- $W$  matrix, where  $W$  is the number of process noise terms.

If the process noise covariance is time-varying, select **Time-varying**. The block generates input port **Q** to specify the time-varying covariance.

### Dependencies

This parameter is enabled if you do not specify the process noise as **Time-Varying**.

### Programmatic Use

**Block Parameter:** ProcessNoise

**Type:** character vector, string

**Default:** '1'

**Time-varying — Time-varying process noise covariance**

'off' (default) | 'on'

If you select this parameter, the block includes an additional input port **Q** to specify the time-varying process noise covariance.

**Programmatic Use****Block Parameter:** HasTimeVaryingProcessNoise**Type:** character vector**Values:** 'off', 'on'**Default:** 'off'**Initialization****Initial state — Initial state estimate** $\theta$  (default) | vector

Initial state estimate value, specified as an  $N_s$ -element vector, where  $N_s$  is the number of states in the system. Specify the initial state values based on your knowledge of the system.

**Programmatic Use****Block Parameter:** InitialState**Type:** character vector, string**Default:** '0'**Initial covariance — State estimation error covariance**

1 (default) | scalar | vector | matrix

State estimation error covariance, specified as a scalar, an  $N_s$ -element vector, or an  $N_s$ -by- $N_s$  matrix, where  $N_s$  is the number of states of the system. If you specify a scalar or vector, the software creates an  $N_s$ -by- $N_s$  diagonal matrix with the scalar or vector elements on the diagonal.

Specify a high value for the covariance when you do not have confidence in the initial state values that you specify in **Initial state**.

**Programmatic Use****Block Parameter:** InitialStateCovariance**Type:** character vector, string**Default:** '1'

### Unscented Transformation Parameters

#### Alpha — Spread of sigma points

1e-3 (default) | scalar value between 0 and 1

The unscented Kalman filter algorithm treats the state of the system as a random variable with a mean state value and variance. To compute the state and its statistical properties at the next time step, the algorithm first generates a set of state values distributed around the mean value by using the unscented transformation. These generated state values are called sigma points. The algorithm uses each of the sigma points as an input to the state transition and measurement functions to get a new set of transformed state points and measurements. The transformed points are used to compute the state and state estimation error covariance value at the next time step.

The spread of the sigma points around the mean state value is controlled by two parameters **Alpha** and **Kappa**. A third parameter, **Beta**, impacts the weights of the transformed points during state and measurement covariance calculations:

- **Alpha** — Determines the spread of the sigma points around the mean state value. Specify as a scalar value between 0 and 1 ( $0 < \mathbf{Alpha} \leq 1$ ). It is usually a small positive value. The spread of sigma points is proportional to **Alpha**. Smaller values correspond to sigma points closer to the mean state.
- **Kappa** — A second scaling parameter that is typically set to 0. Smaller values correspond to sigma points closer to the mean state. The spread is proportional to the square-root of Kappa.
- **Beta** — Incorporates prior knowledge of the distribution of the state. For Gaussian distributions, **Beta** = 2 is optimal.

If you know the distribution of state and state covariance, you can adjust these parameters to capture the transformation of higher-order moments of the distribution. The algorithm can track only a single peak in the probability distribution of the state. If there are multiple peaks in the state distribution of your system, you can adjust these parameters so that the sigma points stay around a single peak. For example, choose a small **Alpha** to generate sigma points close to the mean state value.

For more information, see “Unscented Kalman Filter Algorithm”.

#### Programmatic Use

**Block Parameter:** Alpha

**Type:** character vector, string

**Default:** '1e-3'

**Beta — Characterization of state distribution**

2 (default) | scalar value greater than or equal to 0

Characterization of the state distribution that is used to adjust weights of transformed sigma points, specified as a scalar value greater than or equal to 0. For Gaussian distributions,  $\text{Beta} = 2$  is the optimal choice.

For more information, see the description for **Alpha**.

**Programmatic Use**

**Block Parameter:** Beta

**Type:** character vector, string

**Default:** '2'

**Kappa — Spread of sigma points**

0 (default) | scalar value between 0 and 3

Spread of sigma points around mean state value, specified as a scalar value between 0 and 3 ( $0 \leq \text{Kappa} \leq 3$ ). **Kappa** is typically specified as 0. Smaller values correspond to sigma points closer to the mean state. The spread is proportional to the square root of **Kappa**. For more information, see the description for **Alpha**.

**Programmatic Use**

**Block Parameter:** Kappa

**Type:** character vector, string

**Default:** '0'

**Measurement****Function — Measurement function name**

myMeasurementFcn (default) | function name

The measurement function calculates the  $N$ -element output measurement vector of the nonlinear system at time step  $k$ , given the state vector at time step  $k$ . You create the measurement function and specify the function name in **Function**. For example, if `vdpMeasurementFcn.m` is the measurement function that you created and saved, specify **Function** as `vdpMeasurementFcn`.

The inputs to the function you create depend on whether you specify the measurement noise as additive or nonadditive in **Measurement noise**.

- **Measurement noise** is Additive — The measurement function  $h$  specifies how the measurements evolve as a function of state Values:

$$y(k) = h(x(k), Um1(k), \dots, Umn(k)),$$

where  $y(k)$  and  $x(k)$  are the estimated output and estimated state at time  $k$ , and  $Um1, \dots, Umn$  are any optional input arguments required by your measurement function. For example, if you are using a sensor for tracking an object, an additional input could be the sensor position.

To see an example of a measurement function with additive process noise, type `edit vdpMeasurementFcn` at the command line.

- **Measurement noise** is Nonadditive— The measurement function also specifies how the output measurement evolves as a function of the measurement noise  $v$ :

$$y(k) = h(x(k), v(k), Um1(k), \dots, Umn(k)).$$

To see an example of a measurement function with nonadditive process noise, type `edit vdpMeasurementNonAdditiveNoiseFcn`.

For more information, see “State Transition and Measurement Functions” on page 2-137.

You can create  $h$  using a Simulink Function block or as a MATLAB function (.m file).

- You can use a MATLAB function only if  $h$  has one additional input argument  $Um1$  other than  $x$  and  $v$ .

$$y[k] = h(x[k], v[k], Um1(k))$$

The software generates an additional input port **MeasurementFcnInput** to specify this argument.

- If you are using a Simulink Function block, specify  $x$  and  $v$  using Argument Inport blocks and the additional inputs  $Um1, \dots, Umn$  using Inport blocks in the Simulink Function block. You do not provide  $Um1, \dots, Umn$  to the Unscented Kalman Filter block.

If you have multiple sensors in your system, you can specify multiple measurement functions. You can specify up to five measurement functions using the **Add Measurement** button. To remove measurement functions, use **Remove Measurement**.

#### Programmatic Use

**Block Parameter:** MeasurementFcn1, MeasurementFcn2, MeasurementFcn3, MeasurementFcn4, MeasurementFcn5

**Type:** character vector, string

**Default:** 'myMeasurementFcn'

### Measurement noise — Measurement noise characteristics

Additive (default) | Nonadditive

Measurement noise characteristics, specified as one of the following values:

- **Additive** — Measurement noise  $v$  is additive, and the measurement function  $h$  that you specify in **Function** has the following form:

$$y(k) = h(x(k), Um1(k), \dots, Umn(k)),$$

where  $y(k)$  and  $x(k)$  are the estimated output and estimated state at time  $k$ , and  $Um1, \dots, Umn$  are any optional input arguments required by your measurement function.

- **Nonadditive** — Measurement noise is nonadditive, and the measurement function specifies how the output measurement evolves as a function of the state *and* measurement noise:

$$y(k) = h(x(k), v(k), Um1(k), \dots, Umn(k)).$$

#### Programmatic Use

**Block Parameter:** HasAdditiveMeasurementNoise1, HasAdditiveMeasurementNoise2, HasAdditiveMeasurementNoise3, HasAdditiveMeasurementNoise4, HasAdditiveMeasurementNoise5

**Type:** character vector

**Values:** 'Additive', 'Nonadditive'

**Default:** 'Additive'

### Covariance — Time-invariant measurement noise covariance

1 (default) | matrix

Time-invariant measurement noise covariance, specified as a matrix. The size of the matrix depends on the value of the **Measurement noise** parameter:

- **Measurement noise** is **Additive** — Specify the covariance as an  $N$ -by- $N$  matrix, where  $N$  is the number of measurements of the system.
- **Measurement noise** is **Nonadditive** — Specify the covariance as a  $V$ -by- $V$  matrix, where  $V$  is the number of measurement noise terms.

If the measurement noise covariance is time-varying, select **Time-varying**. The block generates input port **R $i$**  to specify the time-varying covariance for the  $i^{\text{th}}$  measurement function.

**Dependencies**

This parameter is enabled if you do not specify the process noise as **Time-Varying**.

**Programmatic Use**

**Block Parameter:** MeasurementNoise1, MeasurementNoise2, MeasurementNoise3, MeasurementNoise4, MeasurementNoise5

**Type:** character vector, string

**Default:** '1'

**Time-varying — Time-varying measurement noise covariance**

off (default) | on

If you select this parameter for the measurement noise covariance of the first measurement function, the block includes an additional input port **R1**. You specify the time-varying measurement noise covariance in **R1**. Similarly, if you select **Time-varying** for the  $i^{\text{th}}$  measurement function, the block includes an additional input port **R $i$**  to specify the time-varying measurement noise covariance for that function.

**Programmatic Use**

**Block Parameter:** HasTimeVaryingMeasurementNoise1, HasTimeVaryingMeasurementNoise2, HasTimeVaryingMeasurementNoise3, HasTimeVaryingMeasurementNoise4, HasTimeVaryingMeasurementNoise5

**Type:** character vector

**Values:** 'off', 'on'

**Default:** 'off'

**Add Enable Port — Enable correction of estimated states only when measured data is available**

off (default) | on

Suppose that measured output data is not available at all time points at the port **y1** that corresponds to the first measurement function. Select **Add Enable port** to generate an input port **Enable1**. Use a signal at this port to enable the correction of estimated states only when measured data is available. Similarly, if measured output data is not available at all time points at the port **y $i$**  for the  $i^{\text{th}}$  measurement function, select the corresponding **Add Enable port**.

**Programmatic Use**

**Block Parameter:** HasMeasurementEnablePort1, HasMeasurementEnablePort2, HasMeasurementEnablePort3, HasMeasurementEnablePort4, HasMeasurementEnablePort5

**Type:** character vector

**Values:** 'off', 'on'

**Default:** 'off'

**Settings****Use the current measurements to improve state estimates — Choose between corrected or predicted state estimate**

on (default) | off

When this parameter is selected, the block outputs the corrected state estimate  $\hat{x}[k|k]$  at time step  $k$ , estimated using measured outputs until time  $k$ . If you clear this parameter,

the block returns the predicted state estimate  $\hat{x}[k|k-1]$  for time  $k$ , estimated using measured output until a previous time  $k-1$ . Clear this parameter if your filter is in a feedback loop and there is an algebraic loop in your Simulink model.

**Programmatic Use**

**Block Parameter:** UseCurrentEstimator

**Type:** character vector

**Values:** 'off', 'on'

**Default:** 'on'

**Output state estimation covariance — Output state estimation error covariance**

off (default) | on

If you select this parameter, a state estimation error covariance output port **P** is generated in the block.

**Programmatic Use**

**Block Parameter:** OutputStateCovariance

**Type:** character vector

**Values:** 'off', 'on'

**Default:** 'off'

**Data type — Data type for block parameters**

double (default) | single

Use this parameter to specify the data type for all block parameters.

### **Programmatic Use**

**Block Parameter:** DataType

**Type:** character vector

**Values:** 'single', 'double'

**Default:** 'double'

### **Sample time — Block sample time**

1 (default) | positive scalar

Block sample time, specified as a positive scalar. If the sample times of your state transition and measurement functions are different, select **Enable multirate operation** in the **Multirate** tab, and specify the sample times in the **Multirate** tab instead.

### **Dependencies**

This parameter is available if in the **Multirate** tab, the **Enable multirate operation** parameter is off.

### **Programmatic Use**

**Block Parameter:** SampleTime

**Type:** character vector, string

**Default:** '1'

## **Multirate Tab**

### **Enable multirate operation — Enable specification of different sample times for state transition and measurement functions**

off (default) | on

Select this parameter if the sample times of the state transition and measurement functions are different. You specify the sample times in the **Multirate** tab, in **Sample time**.

### **Programmatic Use**

**Block Parameter:** EnableMultirate

**Type:** character vector

**Values:** 'off', 'on'

**Default:** 'off'

### **Sample time — State transition and measurement function sample times**

positive scalar

If the sample times for state transition and measurement functions are different, specify **Sample time**. Specify the sample times for the measurement functions as positive integer multiples of the state transition sample time. The sample times you specify correspond to the following input ports:

- Ports corresponding to state transition function — Additional input to state transition function **StateTransitionFcnInputs** and time-varying process noise covariance **Q**. The sample times of these ports must always equal the state transition function sample time, but can differ from the sample time of the measurement functions.
- Ports corresponding to  $i^{\text{th}}$  measurement function — Measured output **yi**, additional input to measurement function **MeasurementFcnInputs**, enable signal at port **Enablei**, and time-varying measurement noise covariance **Ri**. The sample times of these ports for the same measurement function must always be the same, but can differ from the sample time for the state transition function and other measurement functions.

### Dependencies

This parameter is available if in the **Multirate** tab, the **Enable multirate operation** parameter is on.

### Programmatic Use

**Block Parameter:** StateTransitionFcnSampleTime,  
MeasurementFcn1SampleTime1, MeasurementFcn1SampleTime2,  
MeasurementFcn1SampleTime3, MeasurementFcn1SampleTime4,  
MeasurementFcn1SampleTime5

**Type:** character vector, string

**Default:** '1'

## Definitions

### State Transition and Measurement Functions

The algorithm computes the state estimates  $\hat{x}$  of the nonlinear system using state transition and measurement functions specified by you. You can specify up to five measurement functions, each corresponding to a sensor in the system. The software lets you specify the noise in these functions as additive or nonadditive.

- **Additive Noise Terms** — The state transition and measurements equations have the following form:

$$x[k+1] = f(x[k], u_s[k]) + w[k]$$

$$y[k] = h(x[k], u_m[k]) + v[k]$$

Here  $f$  is a nonlinear state transition function that describes the evolution of states  $x$  from one time step to the next. The nonlinear measurement function  $h$  relates  $x$  to the measurements  $y$  at time step  $k$ .  $w$  and  $v$  are the zero-mean, uncorrelated process and measurement noises, respectively. These functions can also have additional optional input arguments that are denoted by  $u_s$  and  $u_m$  in the equations. For example, the additional arguments could be time step  $k$  or the inputs  $u$  to the nonlinear system. There can be multiple such arguments.

Note that the noise terms in both equations are additive. That is,  $x(k+1)$  is linearly related to the process noise  $w(k)$ , and  $y(k)$  is linearly related to the measurement noise  $v(k)$ . For additive noise terms, you do not need to specify the noise terms in the state transition and measurement functions. The software adds the terms to the output of the functions.

- **Nonadditive Noise Terms** — The software also supports more complex state transition and measurement functions where the state  $x[k]$  and measurement  $y[k]$  are nonlinear functions of the process noise and measurement noise, respectively. When the noise terms are nonadditive, the state transition and measurements equation have the following form:

$$x[k+1] = f(x[k], w[k], u_s[k])$$

$$y[k] = h(x[k], v[k], u_m[k])$$

## Extended Capabilities

### C/C++ Code Generation

Generate C and C++ code using Simulink® Coder™.

The state transition and measurement functions that you specify must use only the MATLAB commands and Simulink blocks that support code generation. For a list of blocks that support code generation, see “Simulink Built-In Blocks That Support Code Generation” (Simulink Coder). For a list of commands that support code generation, see “Functions and Objects Supported for C/C++ Code Generation — Category List” (MATLAB Coder).

## See Also

### Blocks

Extended Kalman Filter | Kalman Filter | Particle Filter

### Functions

extendedKalmanFilter | particleFilter | unscentedKalmanFilter

### Topics

“What Is Online Estimation?”

“Extended and Unscented Kalman Filter Algorithms for Online State Estimation”

“Validate Online State Estimation in Simulink”

“Troubleshoot Online State Estimation”

### Introduced in R2017a

