

NAG Fortran Library Routine Document

D03PPF/D03PPA

Note: before using this routine, please read the Users' Note for your implementation to check the interpretation of *bold italicised* terms and other implementation-dependent details.

1 Purpose

D03PPF/D03PPA integrates a system of linear or nonlinear parabolic partial differential equations (PDEs) in one space variable, with scope for coupled ordinary differential equations (ODEs), and automatic adaptive spatial remeshing. The spatial discretization is performed using finite differences, and the method of lines is employed to reduce the PDEs to a system of ODEs. The resulting system is solved using a Backward Differentiation Formula (BDF) method or a Theta method (switching between Newton's method and functional iteration).

D03PPA is a version of D03PPF that has additional parameters in order to make it safe for use in multithreaded applications (see Section 5).

2 Specification

2.1 Specification for D03PPF

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SUBROUTINE D03PPF (NPDE, M, TS, TOUT, PDEDEF, BNDARY, UVINIT, U, NPTS,
1 X, NCODE, ODEDEF, NXI, XI, NEQN, RTOL, ATOL, ITOL,
2 NORM, LAOPT, ALGOPT, REMESH, NXFIX, XFIX, NRMESH,
3 DXMESH, TRMESH, IPMINF, XRATIO, CON, MONITF, RSAVE,
4 LRSAVE, ISAVE, LISAVE, ITASK, ITRACE, IND, IFAIL)

INTEGER NPDE, M, NPTS, NCODE, NXI, NEQN, ITOL, NXFIX, NRMESH,
1 IPMINF, LRSAVE, ISAVE(LISAVE), LISAVE, ITASK, ITRACE,
2 IND, IFAIL
double precision TS, TOUT, U(NEQN), X(NPTS), XI(*), RTOL(*), ATOL(*),
1 ALGOPT(30), XFIX(*), DXMESH, TRMESH, XRATIO, CON,
2 RSAVE(LRSAVE)
LOGICAL REMESH
CHARACTER*1 NORM, LAOPT
EXTERNAL PDEDEF, BNDARY, UVINIT, ODEDEF, MONITF

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2.2 Specification for D03PPA

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SUBROUTINE D03PPA (NPDE, M, TS, TOUT, PDEDEF, BNDARY, UVINIT, U, NPTS,
1 X, NCODE, ODEDEF, NXI, XI, NEQN, RTOL, ATOL, ITOL,
2 NORM, LAOPT, ALGOPT, REMESH, NXFIX, XFIX, NRMESH,
3 DXMESH, TRMESH, IPMINF, XRATIO, CON, MONITF, RSAVE,
4 LRSAVE, ISAVE, LISAVE, ITASK, ITRACE, IND, IUSER,
5 RUSER, CWSAV, LWSAV, IWSAV, RWSAV, IFAIL)

INTEGER NPDE, M, NPTS, NCODE, NXI, NEQN, ITOL, NXFIX, NRMESH,
1 IPMINF, LRSAVE, ISAVE(LISAVE), LISAVE, ITASK, ITRACE,
2 IND, IUSER(*), IWSAV(505), IFAIL
double precision TS, TOUT, U(NEQN), X(NPTS), XI(*), RTOL(*), ATOL(*),
1 ALGOPT(30), XFIX(*), DXMESH, TRMESH, XRATIO, CON,
2 RSAVE(LRSAVE), RUSER(*), RWSAV(1100)
LOGICAL REMESH, LWSAV(100)
CHARACTER*1 NORM, LAOPT
CHARACTER*80 CWSAV(10)
EXTERNAL PDEDEF, BNDARY, UVINIT, ODEDEF, MONITF

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3 Description

D03PPF/D03PPA integrates the system of parabolic-elliptic equations and coupled ODEs

$$\sum_{j=1}^{\text{NPDE}} P_{ij} \frac{\partial U_j}{\partial t} + Q_i = x^{-m} \frac{\partial}{\partial x} (x^m R_i), \quad i = 1, 2, \dots, \text{NPDE}, \quad a \leq x \leq b, t \geq t_0, \quad (1)$$

$$F_i(t, V, \dot{V}, \xi, U^*, U_x^*, R^*, U_t^*, U_{xt}^*) = 0, \quad i = 1, 2, \dots, \text{NCODE}, \quad (2)$$

where (1) defines the PDE part and (2) generalizes the coupled ODE part of the problem.

In (1), P_{ij} and R_i depend on x , t , U , U_x , and V ; Q_i depends on x , t , U , U_x , V and **linearly** on \dot{V} . The vector U is the set of PDE solution values

$$U(x, t) = \left[U_1(x, t), \dots, U_{\text{NPDE}}(x, t) \right]^T,$$

and the vector U_x is the partial derivative with respect to x . The vector V is the set of ODE solution values

$$V(t) = \left[V_1(t), \dots, V_{\text{NCODE}}(t) \right]^T,$$

and \dot{V} denotes its derivative with respect to time.

In (2), ξ represents a vector of n_ξ spatial coupling points at which the ODEs are coupled to the PDEs. These points may or may not be equal to some of the PDE spatial mesh points. U^* , U_x^* , R^* , U_t^* and U_{xt}^* are the functions U , U_x , R , U_t and U_{xt} evaluated at these coupling points. Each F_i may only depend linearly on time derivatives. Hence the equation (2) may be written more precisely as

$$F = G - A\dot{V} - B \begin{pmatrix} U_t^* \\ U_{xt}^* \end{pmatrix}, \quad (3)$$

where $F = \left[F_1, \dots, F_{\text{NCODE}} \right]^T$, G is a vector of length NCODE, A is an NCODE by NCODE matrix, B is an NCODE by $(n_\xi \times \text{NPDE})$ matrix and the entries in G , A and B may depend on t , ξ , U^* , U_x^* and V . In practice you only need to supply a vector of information to define the ODEs and not the matrices A and B . (See Section 5 for the specification of the user-supplied (sub)program ODEDEF.)

The integration in time is from t_0 to t_{out} , over the space interval $a \leq x \leq b$, where $a = x_1$ and $b = x_{\text{NPTS}}$ are the leftmost and rightmost points of a mesh $x_1, x_2, \dots, x_{\text{NPTS}}$ defined initially by you and (possibly) adapted automatically during the integration according to user-specified criteria. The co-ordinate system in space is defined by the following values of m ; $m = 0$ for Cartesian co-ordinates, $m = 1$ for cylindrical polar co-ordinates and $m = 2$ for spherical polar co-ordinates.

The PDE system which is defined by the functions P_{ij} , Q_i and R_i must be specified in the user-supplied (sub)program PDEDEF.

The initial ($t = t_0$) values of the functions $U(x, t)$ and $V(t)$ must be specified in a (sub)program UVINIT supplied by you. Note that UVINIT will be called again following any initial remeshing, and so $U(x, t_0)$ should be specified for **all** values of x in the interval $a \leq x \leq b$, and not just the initial mesh points.

The functions R_i which may be thought of as fluxes, are also used in the definition of the boundary conditions. The boundary conditions must have the form

$$\beta_i(x, t) R_i(x, t, U, U_x, V) = \gamma_i(x, t, U, U_x, V, \dot{V}), \quad i = 1, 2, \dots, \text{NPDE}, \quad (4)$$

where $x = a$ or $x = b$.

The boundary conditions must be specified in a (sub)program BNDARY provided by you. The function γ_i may depend **linearly** on \dot{V} .

The problem is subject to the following restrictions:

- (i) In (1), $\dot{V}_j(t)$, for $j = 1, 2, \dots, \text{NCODE}$, may only appear **linearly** in the functions Q_i , for $i = 1, 2, \dots, \text{NPDE}$, with a similar restriction for γ ;

- (ii) $P_{i,j}$ and the flux R_i must not depend on any time derivatives;
- (iii) $t_0 < t_{\text{out}}$, so that integration is in the forward direction;
- (iv) The evaluation of the terms $P_{i,j}$, Q_i and R_i is done approximately at the mid-points of the mesh $X(i)$, for $i = 1, 2, \dots, \text{NPTS}$, by calling the user-supplied (sub)program PDEDEF for each mid-point in turn. Any discontinuities in these functions **must** therefore be at one or more of the fixed mesh points specified by XFIX;
- (v) At least one of the functions $P_{i,j}$ must be non-zero so that there is a time derivative present in the PDE problem;
- (vi) If $m > 0$ and $x_1 = 0.0$, which is the left boundary point, then it must be ensured that the PDE solution is bounded at this point. This can be done by either specifying the solution at $x = 0.0$ or by specifying a zero flux there, that is $\beta_i = 1.0$ and $\gamma_i = 0.0$. See also Section 8.

The algebraic-differential equation system which is defined by the functions F_i must be specified in the user-supplied (sub)program ODEDEF. You must also specify the coupling points ξ in the array XI.

The parabolic equations are approximated by a system of ODEs in time for the values of U_i at mesh points. For simple problems in Cartesian co-ordinates, this system is obtained by replacing the space derivatives by the usual central, three-point finite-difference formula. However, for polar and spherical problems, or problems with nonlinear coefficients, the space derivatives are replaced by a modified three-point formula which maintains second order accuracy. In total there are $\text{NPDE} \times \text{NPTS} + \text{NCODE}$ ODEs in time direction. This system is then integrated forwards in time using a Backward Differentiation Formula (BDF) or a Theta method.

The adaptive space remeshing can be used to generate meshes that automatically follow the changing time-dependent nature of the solution, generally resulting in a more efficient and accurate solution using fewer mesh points than may be necessary with a fixed uniform or non-uniform mesh. Problems with travelling wavefronts or variable-width boundary layers for example will benefit from using a moving adaptive mesh. The discrete time-step method used here (developed by Furzeland (1984)) automatically creates a new mesh based on the current solution profile at certain time-steps, and the solution is then interpolated onto the new mesh and the integration continues.

The method requires you to supply a (sub)program MONITF which specifies in an analytical or numerical form the particular aspect of the solution behaviour you wish to track. This so-called monitor function is used to choose a mesh which equally distributes the integral of the monitor function over the domain. A typical choice of monitor function is the second space derivative of the solution value at each point (or some combination of the second space derivatives if there is more than one solution component), which results in refinement in regions where the solution gradient is changing most rapidly.

You must specify the frequency of mesh updates together with certain other criteria such as adjacent mesh ratios. Remeshing can be expensive and you are encouraged to experiment with the different options in order to achieve an efficient solution which adequately tracks the desired features of the solution.

Note that unless the monitor function for the initial solution values is zero at all user-specified initial mesh points, a new initial mesh is calculated and adopted according to the user-specified remeshing criteria. The user-supplied (sub)program UVINIT will then be called again to determine the initial solution values at the new mesh points (there is no interpolation at this stage) and the integration proceeds.

4 References

- Berzins M (1990) Developments in the NAG Library software for parabolic equations *Scientific Software Systems* (ed J C Mason and M G Cox) 59–72 Chapman and Hall
- Berzins M, Dew P M and Furzeland R M (1989) Developing software for time-dependent problems using the method of lines and differential-algebraic integrators *Appl. Numer. Math.* **5** 375–397
- Berzins M and Furzeland R M (1992) An adaptive theta method for the solution of stiff and nonstiff differential equations *Appl. Numer. Math.* **9** 1–19
- Furzeland R M (1984) The construction of adaptive space meshes *TNER.85.022* Thornton Research Centre, Chester

Skeel R D and Berzins M (1990) A method for the spatial discretization of parabolic equations in one space variable *SIAM J. Sci. Statist. Comput.* **11 (1)** 1–32

5 Parameters

- 1: NPDE – INTEGER *Input*
On entry: the number of PDEs to be solved.
Constraint: NPDE \geq 1.
- 2: M – INTEGER *Input*
On entry: the co-ordinate system used:
M = 0
Indicates Cartesian co-ordinates.
M = 1
Indicates cylindrical polar co-ordinates.
M = 2
Indicates spherical polar co-ordinates.
Constraint: 0 \leq M \leq 2.
- 3: TS – *double precision* *Input/Output*
On entry: the initial value of the independent variable t .
On exit: the value of t corresponding to the solution values in U. Normally TS = TOUT.
Constraint: TS < TOUT.
- 4: TOUT – *double precision* *Input*
On entry: the final value of t to which the integration is to be carried out.
- 5: PDEDEF – SUBROUTINE, supplied by the user. *External Procedure*
PDEDEF must evaluate the functions $P_{i,j}$, Q_i and R_i which define the system of PDEs. The functions may depend on x , t , U , U_x and V . Q_i may depend linearly on \dot{V} . PDEDEF is called approximately midway between each pair of mesh points in turn by D03PPF/D03PPA.

The specification of PDEDEF for D03PPF is:

```

SUBROUTINE PDEDEF (NPDE, T, X, U, UX, NCODE, V, VDOT, P, Q, R, IRES)
  INTEGER          NPDE, NCODE, IRES
  double precision T, X, U(NPDE), UX(NPDE), V(*), VDOT(*),
1                 P(NPDE,NPDE), Q(NPDE), R(NPDE)

```

The specification of PDEDEF for D03PPA is:

```

SUBROUTINE PDEDEF (NPDE, T, X, U, UX, NCODE, V, VDOT, P, Q, R, IRES,
1                 IUSER, RUSER)
  INTEGER          NPDE, NCODE, IRES, IUSER(*)
  double precision T, X, U(NPDE), UX(NPDE), V(*), VDOT(*),
1                 P(NPDE,NPDE), Q(NPDE), R(NPDE), RUSER(*)

```

- 1: NPDE – INTEGER *Input*
On entry: the number of PDEs in the system.

2:	T – double precision <i>On entry:</i> the current value of the independent variable t .	<i>Input</i>
3:	X – double precision <i>On entry:</i> the current value of the space variable x .	<i>Input</i>
4:	U(NPDE) – double precision array <i>On entry:</i> U(i) contains the value of the component $U_i(x, t)$, for $i = 1, 2, \dots, \text{NPDE}$.	<i>Input</i>
5:	UX(NPDE) – double precision array <i>On entry:</i> UX(i) contains the value of the component $\frac{\partial U_i(x, t)}{\partial x}$, for $i = 1, 2, \dots, \text{NPDE}$.	<i>Input</i>
6:	NCODE – INTEGER <i>On entry:</i> the number of coupled ODEs in the system.	<i>Input</i>
7:	V(*) – double precision array Note: the dimension of the array V must be at least NCODE. <i>On entry:</i> V(i) contains the value of component $V_i(t)$, for $i = 1, 2, \dots, \text{NCODE}$.	<i>Input</i>
8:	VDOT(*) – double precision array Note: the dimension of the array VDOT must be at least NCODE. <i>On entry:</i> VDOT(i) contains the value of component $\dot{V}_i(t)$, for $i = 1, 2, \dots, \text{NCODE}$. Note: $\dot{V}_i(t)$, for $i = 1, 2, \dots, \text{NCODE}$, may only appear linearly in Q_j , for $j = 1, 2, \dots, \text{NPDE}$.	<i>Input</i>
9:	P(NPDE,NPDE) – double precision array <i>On exit:</i> P(i, j) must be set to the value of $P_{i,j}(x, t, U, U_x, V)$, for $i, j = 1, 2, \dots, \text{NPDE}$.	<i>Output</i>
10:	Q(NPDE) – double precision array <i>On exit:</i> Q(i) must be set to the value of $Q_i(x, t, U, U_x, V, \dot{V})$, for $i = 1, 2, \dots, \text{NPDE}$.	<i>Output</i>
11:	R(NPDE) – double precision array <i>On exit:</i> R(i) must be set to the value of $R_i(x, t, U, U_x, V)$, for $i = 1, 2, \dots, \text{NPDE}$.	<i>Output</i>
12:	IRES – INTEGER <i>On entry:</i> set to -1 or 1 . <i>On exit:</i> should usually remain unchanged. However, you may set IRES to force the integration routine to take certain actions as described below: IRES = 2 Indicates to the integrator that control should be passed back immediately to the calling (sub)program with the error indicator set to IFAIL = 6. IRES = 3 Indicates to the integrator that the current time step should be abandoned and a smaller time step used instead. You may wish to set IRES = 3 when a physically meaningless input or output value has been generated. If you consecutively set IRES = 3, then D03PPF/D03PPA returns to the calling (sub)program with the error indicator set to IFAIL = 4.	<i>Input/Output</i>

Note: the following are additional parameters for specific use with D03PPA. Users of D03PPF therefore need not read the remainder of this description.

- 13: IUSER(*) – INTEGER array User Workspace
 14: RUSER(*) – **double precision** array User Workspace

PDEDEF is called from D03PPA with the parameters IUSER and RUSER as supplied to D03PPA. You are free to use the arrays IUSER and RUSER to supply information to PDEDEF.

PDEDEF must be declared as EXTERNAL in the (sub)program from which D03PPF/D03PPA is called. Parameters denoted as *Input* must **not** be changed by this procedure.

- 6: BNDARY – SUBROUTINE, supplied by the user. External Procedure

BNDARY must evaluate the functions β_i and γ_i which describe the boundary conditions, as given in (4).

The specification of BNDARY for D03PPF is:

```

SUBROUTINE BNDARY (NPDE, T, U, UX, NCODE, V, VDOT, IBND, BETA,
1                GAMMA, IRES)
  INTEGER
  double precision      NPDE, NCODE, IBND, IRES
                        T, U(NPDE), UX(NPDE), V(*), VDOT(*), BETA(NPDE),
1                GAMMA(NPDE)

```

The specification of BNDARY for D03PPA is:

```

SUBROUTINE BNDARY (NPDE, T, U, UX, NCODE, V, VDOT, IBND, BETA,
1                GAMMA, IRES, IUSER, RUSER)
  INTEGER
  double precision      NPDE, NCODE, IBND, IRES, IUSER(*)
                        T, U(NPDE), UX(NPDE), V(*), VDOT(*), BETA(NPDE),
1                GAMMA(NPDE), RUSER(*)

```

- 1: NPDE – INTEGER Input

On entry: the number of PDEs in the system.

- 2: T – **double precision** Input

On entry: the current value of the independent variable t .

- 3: U(NPDE) – **double precision** array Input

On entry: U(i) contains the value of the component $U_i(x, t)$ at the boundary specified by IBND, for $i = 1, 2, \dots, \text{NPDE}$.

- 4: UX(NPDE) – **double precision** array Input

On entry: UX(i) contains the value of the component $\frac{\partial U_i(x, t)}{\partial x}$ at the boundary specified by IBND, for $i = 1, 2, \dots, \text{NPDE}$.

- 5: NCODE – INTEGER Input

On entry: the number of coupled ODEs in the system.

- 6: V(*) – **double precision** array Input

Note: the dimension of the array V must be at least NCODE.

On entry: V(i) contains the value of component $V_i(t)$, for $i = 1, 2, \dots, \text{NCODE}$.

7:	VDOT(*) – double precision array	<i>Input</i>
	Note: the dimension of the array VDOT must be at least NCODE.	
	<i>On entry:</i> VDOT(<i>i</i>) contains the value of component $\dot{V}_i(t)$, for $i = 1, 2, \dots, \text{NCODE}$.	
	Note: $\dot{V}_i(t)$, for $i = 1, 2, \dots, \text{NCODE}$, may only appear linearly in γ_j , for $j = 1, 2, \dots, \text{NPDE}$.	
8:	IBND – INTEGER	<i>Input</i>
	<i>On entry:</i> specifies which boundary conditions are to be evaluated.	
	IBND = 0	
	BNDARY must set up the coefficients of the left-hand boundary, $x = a$.	
	IBND \neq 0	
	BNDARY must set up the coefficients of the right-hand boundary, $x = b$.	
9:	BETA(NPDE) – double precision array	<i>Output</i>
	<i>On exit:</i> BETA(<i>i</i>) must be set to the value of $\beta_i(x, t)$ at the boundary specified by IBND, for $i = 1, 2, \dots, \text{NPDE}$.	
10:	GAMMA(NPDE) – double precision array	<i>Output</i>
	<i>On exit:</i> GAMMA(<i>i</i>) must be set to the value of $\gamma_i(x, t, U, U_x, V, \dot{V})$ at the boundary specified by IBND, for $i = 1, 2, \dots, \text{NPDE}$.	
11:	IRES – INTEGER	<i>Input/Output</i>
	<i>On entry:</i> set to -1 or 1 .	
	<i>On exit:</i> should usually remain unchanged. However, you may set IRES to force the integration routine to take certain actions as described below:	
	IRES = 2	
	Indicates to the integrator that control should be passed back immediately to the calling (sub)program with the error indicator set to IFAIL = 6.	
	IRES = 3	
	Indicates to the integrator that the current time step should be abandoned and a smaller time step used instead. You may wish to set IRES = 3 when a physically meaningless input or output value has been generated. If you consecutively set IRES = 3, then D03PPF/D03PPA returns to the calling (sub)program with the error indicator set to IFAIL = 4.	
	Note: the following are additional parameters for specific use with D03PPA. Users of D03PPF therefore need not read the remainder of this description.	
12:	IUSER(*) – INTEGER array	<i>User Workspace</i>
13:	RUSER(*) – double precision array	<i>User Workspace</i>
	BNDARY is called from D03PPA with the parameters IUSER and RUSER as supplied to D03PPA. You are free to use the arrays IUSER and RUSER to supply information to BNDARY.	

BNDARY must be declared as EXTERNAL in the (sub)program from which D03PPF/D03PPA is called. Parameters denoted as *Input* must **not** be changed by this procedure.

7: UVINIT – SUBROUTINE, supplied by the user.

External Procedure

UVINIT must supply the initial ($t = t_0$) values of $U(x, t)$ and $V(t)$ for all values of x in the interval $a \leq x \leq b$.

The specification of UVINIT for D03PPF is:

```

SUBROUTINE UVINIT (NPDE, NPTS, NXI, X, XI, U, NCODE, V)
  INTEGER          NPDE, NPTS, NXI, NCODE
  double precision X(NPTS), XI(*), U(NPDE,NPTS), V(*)

```

The specification of UVINIT for D03PPA is:

```

SUBROUTINE UVINIT (NPDE, NPTS, NXI, X, XI, U, NCODE, V, IUSER,
1                 RUSER)
  INTEGER          NPDE, NPTS, NXI, NCODE, IUSER(*),
  double precision X(NPTS), XI(*), U(NPDE,NPTS), V(*), RUSER(*)

```

- 1: NPDE – INTEGER *Input*
On entry: the number of PDEs in the system.
- 2: NPTS – INTEGER *Input*
On entry: the number of mesh points in the interval $[a, b]$.
- 3: NXI – INTEGER *Input*
On entry: the number of ODE/PDE coupling points.
- 4: X(NPTS) – **double precision** array *Input*
On entry: the current mesh. $X(i)$ contains the value of x_i , for $i = 1, 2, \dots, NPTS$.
- 5: XI(*) – **double precision** array *Input*
Note: the dimension of the array XI must be at least NXI.
On entry: XI(i) contains the value of the ODE/PDE coupling point, ξ_i , for $i = 1, 2, \dots, NXI$.
- 6: U(NPDE,NPTS) – **double precision** array *Output*
On exit: U(i, j) contains the value of the component $U_i(x_j, t_0)$, for $i = 1, 2, \dots, NPDE$; $j = 1, 2, \dots, NPTS$.
- 7: NCODE – INTEGER *Input*
On entry: the number of coupled ODEs in the system.
- 8: V(*) – **double precision** array *Output*
Note: the dimension of the array V must be at least NCODE.
On exit: V(i) contains the value of component $V_i(t_0)$, for $i = 1, 2, \dots, NCODE$.
- Note:** the following are additional parameters for specific use with D03PPA. Users of D03PPF therefore need not read the remainder of this description.
- 9: IUSER(*) – INTEGER array *User Workspace*
- 10: RUSER(*) – **double precision** array *User Workspace*

UVINIT is called from D03PPA with the parameters IUSER and RUSER as supplied to D03PPA. You are free to use the arrays IUSER and RUSER to supply information to UVINIT.

UVINIT must be declared as EXTERNAL in the (sub)program from which D03PPF/D03PPA is called. Parameters denoted as *Input* must **not** be changed by this procedure.

- 8: U(NEQN) – **double precision** array *Input/Output*
On entry: if IND = 1, the value of U must be unchanged from the previous call.
On exit: U(NPDE × (j – 1) + i) contains the computed solution $U_i(x_j, t)$, for $i = 1, 2, \dots, \text{NPDE}$; $j = 1, 2, \dots, \text{NPTS}$, and U(NPTS × NPDE + k) contains $V_k(t)$, for $k = 1, 2, \dots, \text{NCODE}$, evaluated at $t = \text{TS}$.
- 9: NPTS – INTEGER *Input*
On entry: the number of mesh points in the interval $[a, b]$.
Constraint: NPTS ≥ 3.
- 10: X(NPTS) – **double precision** array *Input/Output*
On entry: the initial mesh points in the space direction. X(1) must specify the left-hand boundary, a and X(NPTS) must specify the right-hand boundary, b .
Constraint: X(1) < X(2) < ... < X(NPTS).
On exit: the final values of the mesh points.
- 11: NCODE – INTEGER *Input*
On entry: the number of coupled ODE in the system.
Constraint: NCODE ≥ 0.
- 12: ODEDEF – SUBROUTINE, supplied by the user. *External Procedure*
 ODEDEF must evaluate the functions F , which define the system of ODEs, as given in (3). If you wish to compute the solution of a system of PDEs only (NCODE = 0), ODEDEF must be the dummy routine D03PCK for D03PPF (or D53PCK for D03PPA). (D03PCK and D53PCK are included in the NAG Fortran Library; however, their names may be implementation-dependent: see the Users' Note for your implementation for details.)

The specification of ODEDEF for D03PPF is:

```

SUBROUTINE ODEDEF (NPDE, T, NCODE, V, VDOT, NXI, XI, UCP, UCPX, RCP,
1 UCPT, UCPTX, F, IRES)
  INTEGER
  double precision NPDE, NCODE, NXI, IRES
1 T, V(*), VDOT(*), XI(*), UCP(NPDE,*),
2 UCPX(NPDE,*), RCP(NPDE,*), UCPT(NPDE,*),
  UCPTX(NPDE,*), F(*)

```

The specification of ODEDEF for D03PPA is:

```

SUBROUTINE ODEDEF (NPDE, T, NCODE, V, VDOT, NXI, XI, UCP, UCPX, RCP,
1 UCPT, UCPTX, F, IRES, IUSER, RUSER)
  INTEGER
  double precision NPDE, NCODE, NXI, IRES, IUSER(*)
1 T, V(*), VDOT(*), XI(*), UCP(NPDE,*),
2 UCPX(NPDE,*), RCP(NPDE,*), UCPT(NPDE,*),
  UCPTX(NPDE,*), F(*), RUSER(*)

```

- 1: NPDE – INTEGER *Input*
On entry: the number of PDEs in the system.
- 2: T – **double precision** *Input*
On entry: the current value of the independent variable t .

3:	<p>NCODE – INTEGER</p> <p><i>On entry:</i> the number of coupled ODEs in the system.</p>	<i>Input</i>
4:	<p>V(*) – double precision array</p> <p>Note: the dimension of the array V must be at least NCODE.</p> <p><i>On entry:</i> V(i) contains the value of component $V_i(t)$, for $i = 1, 2, \dots, \text{NCODE}$.</p>	<i>Input</i>
5:	<p>VDOT(*) – double precision array</p> <p>Note: the dimension of the array VDOT must be at least NCODE.</p> <p><i>On entry:</i> VDOT(i) contains the value of component $\dot{V}_i(t)$, for $i = 1, 2, \dots, \text{NCODE}$.</p>	<i>Input</i>
6:	<p>NXI – INTEGER</p> <p><i>On entry:</i> the number of ODE/PDE coupling points.</p>	<i>Input</i>
7:	<p>XI(*) – double precision array</p> <p>Note: the dimension of the array XI must be at least NXI.</p> <p><i>On entry:</i> XI(i) contains the ODE/PDE coupling points, ξ_i, for $i = 1, 2, \dots, \text{NXI}$.</p>	<i>Input</i>
8:	<p>UCP(NPDE,*) – double precision array</p> <p>Note: the second dimension of the array UCP must be at least max(1, NXI).</p> <p><i>On entry:</i> UCP(i,j) contains the value of $U_i(x, t)$ at the coupling point $x = \xi_j$, for $i = 1, 2, \dots, \text{NPDE}$; $j = 1, 2, \dots, \text{NXI}$.</p>	<i>Input</i>
9:	<p>UCPX(NPDE,*) – double precision array</p> <p>Note: the second dimension of the array UCPX must be at least max(1, NXI).</p> <p><i>On entry:</i> UCPX(i,j) contains the value of $\frac{\partial U_i(x, t)}{\partial x}$ at the coupling point $x = \xi_j$, for $i = 1, 2, \dots, \text{NPDE}$; $j = 1, 2, \dots, \text{NXI}$.</p>	<i>Input</i>
10:	<p>RCP(NPDE,*) – double precision array</p> <p>Note: the second dimension of the array RCP must be at least max(1, NXI).</p> <p><i>On entry:</i> RCP(i,j) contains the value of the flux R_i at the coupling point $x = \xi_j$, for $i = 1, 2, \dots, \text{NPDE}$; $j = 1, 2, \dots, \text{NXI}$.</p>	<i>Input</i>
11:	<p>UCPT(NPDE,*) – double precision array</p> <p>Note: the second dimension of the array UCPT must be at least max(1, NXI).</p> <p><i>On entry:</i> UCPT(i,j) contains the value of $\frac{\partial U_i}{\partial t}$ at the coupling point $x = \xi_j$, for $i = 1, 2, \dots, \text{NPDE}$; $j = 1, 2, \dots, \text{NXI}$.</p>	<i>Input</i>
12:	<p>UCPTX(NPDE,*) – double precision array</p> <p>Note: the second dimension of the array UCPTX must be at least max(1, NXI).</p> <p><i>On entry:</i> UCPTX(i,j) contains the value of $\frac{\partial^2 U_i}{\partial x \partial t}$ at the coupling point $x = \xi_j$, for $i = 1, 2, \dots, \text{NPDE}$; $j = 1, 2, \dots, \text{NXI}$.</p>	<i>Input</i>

13: $F(*)$ – **double precision** array *Output*

Note: the dimension of the array F must be at least NCODE.

On exit: $F(i)$ must contain the i th component of F , for $i = 1, 2, \dots, \text{NCODE}$, where F is defined as

$$F = G - A\dot{V} - B \begin{pmatrix} U_t^* \\ U_{xt}^* \end{pmatrix}, \quad (5)$$

or

$$F = -A\dot{V} - B \begin{pmatrix} U_t^* \\ U_{xt}^* \end{pmatrix}. \quad (6)$$

The definition of F is determined by the input value of IRES.

14: IRES – INTEGER *Input/Output*

On entry: the form of F that must be returned in the array F . If $\text{IRES} = 1$, then equation (5) must be used. If $\text{IRES} = -1$, then equation (6) must be used.

On exit: should usually remain unchanged. However, you may reset IRES to force the integration routine to take certain actions as described below:

IRES = 2

Indicates to the integrator that control should be passed back immediately to the calling (sub)program with the error indicator set to $\text{IFAIL} = 6$.

IRES = 3

Indicates to the integrator that the current time step should be abandoned and a smaller time step used instead. You may wish to set $\text{IRES} = 3$ when a physically meaningless input or output value has been generated. If you consecutively set $\text{IRES} = 3$, then D03PPF/D03PPA returns to the calling (sub)program with the error indicator set to $\text{IFAIL} = 4$.

Note: the following are additional parameters for specific use with D03PPA. Users of D03PPF therefore need not read the remainder of this description.

15: IUSER(*) – INTEGER array *User Workspace*

16: RUSER(*) – **double precision** array *User Workspace*

ODEDEF is called from D03PPA with the parameters IUSER and RUSER as supplied to D03PPA. You are free to use the arrays IUSER and RUSER to supply information to ODEDEF.

ODEDEF must be declared as EXTERNAL in the (sub)program from which D03PPF/D03PPA is called. Parameters denoted as *Input* must **not** be changed by this procedure.

13: NXI – INTEGER *Input*

On entry: the number of ODE/PDE coupling points.

Constraints:

if $\text{NCODE} = 0$, $\text{NXI} = 0$;
if $\text{NCODE} > 0$, $\text{NXI} \geq 0$.

14: XI(*) – **double precision** array *Input*

Note: the dimension of the array XI must be at least $\max(1, \text{NXI})$.

On entry: $\text{XI}(i)$, for $i = 1, 2, \dots, \text{NXI}$, must be set to the ODE/PDE coupling points.

Constraint: $\text{X}(1) \leq \text{XI}(1) < \text{XI}(2) < \dots < \text{XI}(\text{NXI}) \leq \text{X}(\text{NPTS})$.

15: NEQN – INTEGER *Input*

On entry: the number of ODEs in the time direction.

Constraint: NEQN = NPDE × NPTS + NCODE.

16: RTOL(*) – *double precision* array *Input*

Note: the dimension of the array RTOL must be at least 1 if ITOL = 1 or 2 and at least NEQN if ITOL = 3 or 4.

On entry: the relative local error tolerance.

Constraint: RTOL(*i*) ≥ 0 for all relevant *i*.

17: ATOL(*) – *double precision* array *Input*

Note: the dimension of the array ATOL must be at least 1 if ITOL = 1 or 3 and at least NEQN if ITOL = 2 or 4.

On entry: the absolute local error tolerance.

Constraints:

ATOL(*i*) ≥ 0 for all relevant *i*;

Corresponding elements of ATOL and RTOL cannot both be 0.0.

18: ITOL – INTEGER *Input*

On entry: a value to indicate the form of the local error test. ITOL indicates to D03PPF/D03PPA whether to interpret either or both of RTOL or ATOL as a vector or scalar. The error test to be satisfied is $\|e_i/w_i\| < 1.0$, where w_i is defined as follows:

ITOL	RTOL	ATOL	w_i
1	scalar	scalar	$\text{RTOL}(1) \times U_i + \text{ATOL}(1)$
2	scalar	vector	$\text{RTOL}(1) \times U_i + \text{ATOL}(i)$
3	vector	scalar	$\text{RTOL}(i) \times U_i + \text{ATOL}(1)$
4	vector	vector	$\text{RTOL}(i) \times U_i + \text{ATOL}(i)$

In the above, e_i denotes the estimated local error for the *i*th component of the coupled PDE/ODE system in time, $U(i)$, for $i = 1, 2, \dots, \text{NEQN}$.

The choice of norm used is defined by the parameter NORM.

Constraint: $1 \leq \text{ITOL} \leq 4$.

19: NORM – CHARACTER*1 *Input*

On entry: the type of norm to be used.

NORM = 'M'

Maximum norm.

NORM = 'A'

Averaged L_2 norm.

If U_{norm} denotes the norm of the vector U of length NEQN, then for the averaged L_2 norm

$$U_{\text{norm}} = \sqrt{\frac{1}{\text{NEQN}} \sum_{i=1}^{\text{NEQN}} (U(i)/w_i)^2},$$

while for the maximum norm

$$U_{\text{norm}} = \max_i |U(i)/w_i|.$$

See the description of ITOL for the formulation of the weight vector w .

Constraint: NORM = 'M' or 'A'.

20: LAOPT – CHARACTER*1

Input

On entry: the type of matrix algebra required.

LAOPT = 'F'

Full matrix methods to be used.

LAOPT = 'B'

Banded matrix methods to be used.

LAOPT = 'S'

Sparse matrix methods to be used.

Constraint: LAOPT = 'F', 'B' or 'S'.

Note: you are recommended to use the banded option when no coupled ODEs are present (i.e., NCODE = 0).

21: ALGOPT(30) – *double precision* array

Input

On entry: may be set to control various options available in the integrator. If you wish to employ all the default options, then ALGOPT(1) should be set to 0.0. Default values will also be used for any other elements of ALGOPT set to zero. The permissible values, default values, and meanings are as follows:

ALGOPT(1)

Selects the ODE integration method to be used. If ALGOPT(1) = 1.0, a BDF method is used and if ALGOPT(1) = 2.0, a Theta method is used. The default value is ALGOPT(1) = 1.0.

If ALGOPT(1) = 2.0, then ALGOPT(i), for $i = 2, 3, 4$ are not used.

ALGOPT(2)

Specifies the maximum order of the BDF integration formula to be used. ALGOPT(2) may be 1.0, 2.0, 3.0, 4.0 or 5.0. The default value is ALGOPT(2) = 5.0.

ALGOPT(3)

Specifies what method is to be used to solve the system of nonlinear equations arising on each step of the BDF method. If ALGOPT(3) = 1.0 a modified Newton iteration is used and if ALGOPT(3) = 2.0 a functional iteration method is used. If functional iteration is selected and the integrator encounters difficulty, then there is an automatic switch to the modified Newton iteration. The default value is ALGOPT(3) = 1.0.

ALGOPT(4)

Specifies whether or not the Petzold error test is to be employed. The Petzold error test results in extra overhead but is more suitable when algebraic equations are present, such as $P_{ij} = 0.0$, for $j = 1, 2, \dots, NPDE$ for some i or when there is no $\dot{V}_i(t)$ dependence in the coupled ODE system. If ALGOPT(4) = 1.0, then the Petzold test is used. If ALGOPT(4) = 2.0, then the Petzold test is not used. The default value is ALGOPT(4) = 1.0.

If ALGOPT(1) = 1.0, then ALGOPT(i), for $i = 5, 6, 7$ are not used.

ALGOPT(5)

Specifies the value of Theta to be used in the Theta integration method. $0.51 \leq \text{ALGOPT}(5) \leq 0.99$. The default value is $\text{ALGOPT}(5) = 0.55$.

ALGOPT(6)

Specifies what method is to be used to solve the system of nonlinear equations arising on each step of the Theta method. If $\text{ALGOPT}(6) = 1.0$, a modified Newton iteration is used and if $\text{ALGOPT}(6) = 2.0$, a functional iteration method is used. The default value is $\text{ALGOPT}(6) = 1.0$.

ALGOPT(7)

Specifies whether or not the integrator is allowed to switch automatically between modified Newton and functional iteration methods in order to be more efficient. If $\text{ALGOPT}(7) = 1.0$, then switching is allowed and if $\text{ALGOPT}(7) = 2.0$, then switching is not allowed. The default value is $\text{ALGOPT}(7) = 1.0$.

ALGOPT(11)

Specifies a point in the time direction, t_{crit} , beyond which integration must not be attempted. The use of t_{crit} is described under the parameter ITASK. If $\text{ALGOPT}(1) \neq 0.0$, a value of 0.0 for $\text{ALGOPT}(11)$, say, should be specified even if ITASK subsequently specifies that t_{crit} will not be used.

ALGOPT(12)

Specifies the minimum absolute step size to be allowed in the time integration. If this option is not required, $\text{ALGOPT}(12)$ should be set to 0.0.

ALGOPT(13)

Specifies the maximum absolute step size to be allowed in the time integration. If this option is not required, $\text{ALGOPT}(13)$ should be set to 0.0.

ALGOPT(14)

Specifies the initial step size to be attempted by the integrator. If $\text{ALGOPT}(14) = 0.0$, then the initial step size is calculated internally.

ALGOPT(15)

Specifies the maximum number of steps to be attempted by the integrator in any one call. If $\text{ALGOPT}(15) = 0.0$, then no limit is imposed.

ALGOPT(23)

Specifies what method is to be used to solve the nonlinear equations at the initial point to initialize the values of U , U_t , V and \dot{V} . If $\text{ALGOPT}(23) = 1.0$, a modified Newton iteration is used and if $\text{ALGOPT}(23) = 2.0$, functional iteration is used. The default value is $\text{ALGOPT}(23) = 1.0$.

$\text{ALGOPT}(29)$ and $\text{ALGOPT}(30)$ are used only for the sparse matrix algebra option, $\text{LAOPT} = 'S'$.

ALGOPT(29)

Governs the choice of pivots during the decomposition of the first Jacobian matrix. It should lie in the range $0.0 < \text{ALGOPT}(29) < 1.0$, with smaller values biasing the algorithm towards maintaining sparsity at the expense of numerical stability. If $\text{ALGOPT}(29)$ lies outside this range then the default value is used. If the routines regard the Jacobian matrix as numerically singular then increasing $\text{ALGOPT}(29)$ towards 1.0 may help, but at the cost of increased fill-in. The default value is $\text{ALGOPT}(29) = 0.1$.

ALGOPT(30)

Is used as a relative pivot threshold during subsequent Jacobian decompositions (see $\text{ALGOPT}(29)$) below which an internal error is invoked. If $\text{ALGOPT}(30)$ is greater than 1.0 no check is made on the pivot size, and this may be a necessary option if the Jacobian is

found to be numerically singular (see ALGOPT(29)). The default value is ALGOPT(30) = 0.0001.

- 22: REMESH – LOGICAL *Input*
On entry: indicates whether or not spatial remeshing should be performed.
 REMESH = .TRUE.
 Indicates that spatial remeshing should be performed as specified.
 REMESH = .FALSE.
 Indicates that spatial remeshing should be suppressed.
Note: REMESH should **not** be changed between consecutive calls to D03PPF/D03PPA. Remeshing can be switched off or on at specified times by using appropriate values for the parameters NRMESH and TRMESH at each call.
- 23: NXFIX – INTEGER *Input*
On entry: the number of fixed mesh points.
Constraint: $0 \leq \text{NXFIX} \leq \text{NPTS} - 2$
Note: the end points X(1) and X(NPTS) are fixed automatically and hence should not be specified as fixed points.
- 24: XFIX(*) – *double precision* array *Input*
Note: the dimension of the array XFIX must be at least $\max(1, \text{NXFIX})$.
On entry: XFIX(*i*), for $i = 1, 2, \dots, \text{NXFIX}$, must contain the value of the *x* co-ordinate at the *i*th fixed mesh point.
Constraint: XFIX(*i*) < XFIX(*i* + 1), for $i = 1, 2, \dots, \text{NXFIX} - 1$, and each fixed mesh point must coincide with a user-supplied initial mesh point, that is XFIX(*i*) = X(*j*) for some j , $2 \leq j \leq \text{NPTS} - 1$.
Note: the positions of the fixed mesh points in the array X remain fixed during remeshing, and so the number of mesh points between adjacent fixed points (or between fixed points and end points) does not change. You should take this into account when choosing the initial mesh distribution.
- 25: NRMESH – INTEGER *Input*
On entry: specifies the spatial remeshing frequency and criteria for the calculation and adoption of a new mesh.
 NRMESH < 0
 Indicates that a new mesh is adopted according to the parameter DXMESH. The mesh is tested every |NRMESH| timesteps.
 NRMESH = 0
 Indicates that remeshing should take place just once at the end of the first time step reached when $t > \text{TRMESH}$.
 NRMESH > 0
 Indicates that remeshing will take place every NRMESH time steps, with no testing using DXMESH.
Note: NRMESH may be changed between consecutive calls to D03PPF/D03PPA to give greater flexibility over the times of remeshing.
- 26: DXMESH – *double precision* *Input*
On entry: determines whether a new mesh is adopted when NRMESH is set less than zero. A possible new mesh is calculated at the end of every |NRMESH| time steps, but is adopted only if

$$x_i^{(\text{new})} > x_i^{(\text{old})} + \text{DXMESH} \times (x_{i+1}^{(\text{old})} - x_i^{(\text{old})})$$

or

$$x_i^{(\text{new})} < x_i^{(\text{old})} - \text{DXMESH} \times (x_i^{(\text{old})} - x_{i-1}^{(\text{old})})$$

DXMESH thus imposes a lower limit on the difference between one mesh and the next.

Constraint: $\text{DXMESH} \geq 0.0$.

27: TRMESH – *double precision* *Input*

On entry: specifies when remeshing will take place when NRMESH is set to zero. Remeshing will occur just once at the end of the first time step reached when t is greater than TRMESH.

Note: TRMESH may be changed between consecutive calls to D03PPF/D03PPA to force remeshing at several specified times.

28: IPMINF – INTEGER *Input*

On entry: the level of trace information regarding the adaptive remeshing. Details are directed to the current advisory message unit (see X04ABF).

IPMINF = 0

No trace information.

IPMINF = 1

Brief summary of mesh characteristics.

IPMINF = 2

More detailed information, including old and new mesh points, mesh sizes and monitor function values.

Constraint: $0 \leq \text{IPMINF} \leq 2$.

29: XRATIO – *double precision* *Input*

On entry: an input bound on the adjacent mesh ratio (greater than 1.0 and typically in the range 1.5 to 3.0). The remeshing routines will attempt to ensure that

$$(x_i - x_{i-1})/\text{XRATIO} < x_{i+1} - x_i < \text{XRATIO} \times (x_i - x_{i-1}).$$

Suggested value: XRATIO = 1.5.

Constraint: XRATIO > 1.0.

30: CON – *double precision* *Input*

On entry: an input bound on the sub-integral of the monitor function $F^{\text{mon}}(x)$ over each space step. The remeshing routines will attempt to ensure that

$$\int_{x_i}^{x_{i+1}} F^{\text{mon}}(x) dx \leq \text{CON} \int_{x_1}^{x_{\text{NPTS}}} F^{\text{mon}}(x) dx,$$

(see Furzeland (1984)). CON gives you more control over the mesh distribution e.g., decreasing CON allows more clustering. A typical value is $2/(\text{NPTS} - 1)$, but you are encouraged to experiment with different values. Its value is not critical and the mesh should be qualitatively correct for all values in the range given below.

Suggested value: $\text{CON} = 2.0/(\text{NPTS} - 1)$.

Constraint: $0.1/(\text{NPTS} - 1) \leq \text{CON} \leq 10.0/(\text{NPTS} - 1)$.

31: MONITF – SUBROUTINE, supplied by the user.

External Procedure

MONITF must supply and evaluate a remesh monitor function to indicate the solution behaviour of interest.

If you specify REMESH = .FALSE., i.e., no remeshing, then MONITF will not be called and the dummy routine D03PCL for D03PPF (or D53PCL for D03PPA) may be used for MONITF. (D03PCL and D53PCL are included in the NAG Fortran Library; however, their names may be implementation-dependent: see the Users' Note for your implementation for details.)

The specification of MONITF for D03PPF is:

```
SUBROUTINE MONITF (T, NPTS, NPDE, X, U, R, FMON)
  INTEGER
  double precision
  1 NPTS, NPDE
  T, X(NPTS), U(NPDE,NPTS), R(NPDE,NPTS),
  FMON(NPTS)
```

The specification of MONITF for D03PPA is:

```
SUBROUTINE MONITF (T, NPTS, NPDE, X, U, R, FMON, IUSER, RUSER)
  INTEGER
  double precision
  1 NPTS, NPDE, IUSER(*)
  T, X(NPTS), U(NPDE,NPTS), R(NPDE,NPTS),
  FMON(NPTS), RUSER(*)
```

1: T – **double precision** *Input*

On entry: the current value of the independent variable t .

2: NPTS – INTEGER *Input*

On entry: the number of mesh points in the interval $[a, b]$.

3: NPDE – INTEGER *Input*

On entry: the number of PDEs in the system.

4: X(NPTS) – **double precision** array *Input*

On entry: the current mesh. $X(i)$ contains the value of x_i , for $i = 1, 2, \dots, NPTS$.

5: U(NPDE,NPTS) – **double precision** array *Input*

On entry: $U(i, j)$ contains the value of $U_i(x, t)$ at $x = X(j)$ and time t , for $i = 1, 2, \dots, NPDE$; $j = 1, 2, \dots, NPTS$.

6: R(NPDE,NPTS) – **double precision** array *Input*

On entry: $R(i, j)$ contains the value of $R_i(x, t, U, U_x, V)$ at $x = X(j)$ and time t , for $i = 1, 2, \dots, NPDE$; $j = 1, 2, \dots, NPTS$.

7: FMON(NPTS) – **double precision** array *Output*

On exit: $FMON(i)$ must contain the value of the monitor function $F^{\text{mon}}(x)$ at mesh point $x = X(i)$.

Note: the following are additional parameters for specific use with D03PPA. Users of D03PPF therefore need not read the remainder of this description.

8: IUSER(*) – INTEGER array *User Workspace*

9: RUSER(*) – **double precision** array *User Workspace*

MONITF is called from D03PPA with the parameters IUSER and RUSER as supplied to D03PPA. You are free to use the arrays IUSER and RUSER to supply information to MONITF.

MONITF must be declared as EXTERNAL in the (sub)program from which D03PPF/D03PPA is called. Parameters denoted as *Input* must **not** be changed by this procedure.

- 32: RSAVE(LRSAVE) – *double precision* array *Communication Array*

If IND = 0, RSAVE need not be set on entry.

If IND = 1, RSAVE must be unchanged from the previous call to the routine because it contains required information about the iteration.

- 33: LRSAVE – INTEGER *Input*

On entry: the dimension of the array RSAVE as declared in the (sub)program from which D03PPF/D03PPA is called. Its size depends on the type of matrix algebra selected.

If LAOPT = 'F', $LRSAVE \geq NEQN \times NEQN + NEQN + NWKRES + LENODE$.

If LAOPT = 'B', $LRSAVE \geq (3 \times MLU + 1) \times NEQN + NWKRES + LENODE$.

If LAOPT = 'S', $LRSAVE \geq 4 \times NEQN + 11 \times NEQN/2 + 1 + NWKRES + LENODE$.

Where

MLU = the lower or upper half bandwidths, and
 $MLU = 2 \times NPDE - 1$, for PDE problems only, and
 $MLU = NEQN - 1$, for coupled PDE/ODE problems.

$NWKRES = NPDE \times (3 \times NPDE + 6 \times NXI + NPTS + 15) + NXI + NCODE + 7 \times NPTS + NXFIX + 1$, when NCODE > 0 and NXI > 0, and
 $NWKRES = NPDE \times (3 \times NPDE + NPTS + 21) + NCODE + 7 \times NPTS + NXFIX + 2$, when NCODE > 0 and NXI = 0, and
 $NWKRES = NPDE \times (3 \times NPDE + NPTS + 21) + 7 \times NPTS + NXFIX + 3$, when NCODE = 0.

$LENODE = (6 + \text{int}(\text{ALGOPT}(2))) \times NEQN + 50$, when the BDF method is used, and
 $LENODE = 9 \times NEQN + 50$, when the Theta method is used.

Note: when using the sparse option, the value of LRSAVE may be too small when supplied to the integrator. An estimate of the minimum size of LRSAVE is printed on the current error message unit if ITRACE > 0 and the routine returns with IFAIL = 15.

- 34: ISAVE(LISAVE) – INTEGER array *Communication Array*

If IND = 0, ISAVE need not be set on entry.

If IND = 1, ISAVE must be unchanged from the previous call to the routine because it contains required information about the iteration required for subsequent calls. In particular:

ISAVE(1)

Contains the number of steps taken in time.

ISAVE(2)

Contains the number of residual evaluations of the resulting ODE system used. One such evaluation involves computing the PDE functions at all the mesh points, as well as one evaluation of the functions in the boundary conditions.

ISAVE(3)

Contains the number of Jacobian evaluations performed by the time integrator.

ISAVE(4)

Contains the order of the ODE method last used in the time integration.

ISAVE(5)

Contains the number of Newton iterations performed by the time integrator. Each iteration involves residual evaluation of the resulting ODE system followed by a back-substitution using the *LU* decomposition of the Jacobian matrix.

The rest of the array is used as workspace.

35: LISAVE – INTEGER *Input*

On entry: the dimension of the array ISAVE as declared in the (sub)program from which D03PPF/D03PPA is called.

Its size depends on the type of matrix algebra selected:

if LAOPT = 'B', LISAVE \geq NEQN + 25 + NXFIX;
 if LAOPT = 'F', LISAVE \geq 25 + NXFIX;
 if LAOPT = 'S', LISAVE \geq 25 \times NEQN + 25 + NXFIX.

Note: when using the sparse option, the value of LISAVE may be too small when supplied to the integrator. An estimate of the minimum size of LISAVE is printed on the current error message unit if ITRACE > 0 and the routine returns with IFAIL = 15.

36: ITASK – INTEGER *Input*

On entry: specifies the task to be performed by the ODE integrator.

ITASK = 1

Normal computation of output values U at $t = TOUT$.

ITASK = 2

One step and return.

ITASK = 3

Stop at first internal integration point at or beyond $t = TOUT$.

ITASK = 4

Normal computation of output values U at $t = TOUT$ but without overshooting $t = t_{crit}$ where t_{crit} is described under the parameter ALGOPT.

ITASK = 5

Take one step in the time direction and return, without passing t_{crit} , where t_{crit} is described under the parameter ALGOPT.

Constraint: $1 \leq ITASK \leq 5$.

37: ITRACE – INTEGER *Input*

On entry: the level of trace information required from D03PPF/D03PPA and the underlying ODE solver:

ITRACE ≤ -1

No output is generated.

ITRACE = 0

Only warning messages from the PDE solver are printed on the current error message unit (see X04AAF).

ITRACE = 1

Output from the underlying ODE solver is printed on the current advisory message unit (see X04ABF). This output contains details of Jacobian entries, the nonlinear iteration and the time integration during the computation of the ODE system.

ITRACE = 2

Output from the underlying ODE solver is similar to that produced when ITRACE = 1, except that the advisory messages are given in greater detail.

ITRACE \geq 3

Output from the underlying ODE solver is similar to that produced when ITRACE = 2, except that the advisory messages are given in greater detail.

You are advised to set ITRACE = 0, unless you are experienced with sub-chapter D02M/N.

38: IND – INTEGER *Input/Output*

On entry: must be set to 0 or 1.

IND = 0

Starts or restarts the integration in time.

IND = 1

Continues the integration after an earlier exit from the routine. In this case, only the parameters TOUT and IFAIL and the remeshing parameters NRMESH, DXMESH, TRMESH, XRATIO and CON may be reset between calls to D03PPF/D03PPA.

Constraint: $0 \leq \text{IND} \leq 1$.

On exit: IND = 1.

39: IFAIL – INTEGER *Input/Output*

Note: for D03PPA, IFAIL does not occur in this position in the parameter list. See the additional parameters described below.

On entry: IFAIL must be set to 0, -1 or 1. If you are unfamiliar with this parameter you should refer to Chapter P01 for details.

On exit: IFAIL = 0 unless the routine detects an error (see Section 6).

For environments where it might be inappropriate to halt program execution when an error is detected, the value -1 or 1 is recommended. If the output of error messages is undesirable, then the value 1 is recommended. Otherwise, if you are not familiar with this parameter the recommended value is 0. **When the value -1 or 1 is used it is essential to test the value of IFAIL on exit.**

Note: the following are additional parameters for specific use with D03PPA. Users of D03PPF therefore need not read the remainder of this description.

39: IUSER(*) – INTEGER array *User Workspace*

Note: the dimension of the array IUSER must be at least 1.

IUSER is not used by D03PPA, but is passed directly to the user-supplied (sub)programs PDEDEF, BNDARY, UVINIT, ODEDEF and MONITF and may be used to pass information to these routines.

40: RUSER(*) – *double precision* array *User Workspace*

Note: the dimension of the array RUSER must be at least 1.

RUSER is not used by D03PPA, but is passed directly to the user-supplied (sub)programs PDEDEF, BNDARY, UVINIT, ODEDEF and MONITF and may be used to pass information to these routines.

42:	CWSAV(10) – CHARACTER*80 array	Communication Array
43:	LWSAV(100) – LOGICAL array	Communication Array
44:	IWSAV(505) – INTEGER array	Communication Array
45:	RWSAV(1100) – <i>double precision</i> array	Communication Array
45:	IFAIL – INTEGER	Input/Output

Note: see the parameter description for IFAIL above.

6 Error Indicators and Warnings

If on entry IFAIL = 0 or -1, explanatory error messages are output on the current error message unit (as defined by X04AAF).

Errors or warnings detected by the routine:

IFAIL = 1

On entry, TOUT – TS is too small,
 or ITASK \neq 1, 2, 3, 4 or 5,
 or M \neq 0, 1 or 2,
 or at least one of the coupling points defined in array XI is outside the interval [X(1), X(NPTS)],
 or M > 0 and X(1) < 0.0,
 or NPTS < 3,
 or NPDE < 1,
 or NORM \neq 'A' or 'M',
 or LAOPT \neq 'F', 'B' or 'S',
 or ITOL \neq 1, 2, 3 or 4,
 or IND \neq 0 or 1,
 or mesh points X(i) badly ordered,
 or LRSAVE is too small,
 or LISAVE is too small,
 or NCODE and NXI are incorrectly defined,
 or an element of RTOL or ATOL < 0.0,
 or corresponding elements of RTOL and ATOL are both 0.0,
 or NEQN \neq NPDE \times NPTS + NCODE,
 or NXFIX not in the range 0 to NPTS – 2,
 or fixed mesh point(s) do not coincide with any of the user-supplied mesh points,
 or DXMESH < 0.0,
 or IPMINF \neq 0, 1 or 2,
 or XRATIO \leq 1.0,
 or CON not in the range 0.1/(NPTS – 1) to 10/(NPTS – 1).

IFAIL = 2

The underlying ODE solver cannot make any further progress, with the values of ATOL and RTOL, across the integration range from the current point $t = TS$. The components of U contain the computed values at the current point $t = TS$.

IFAIL = 3

In the underlying ODE solver, there were repeated error test failures on an attempted step, before completing the requested task, but the integration was successful as far as $t = TS$. The problem may have a singularity, or the error requirement may be inappropriate.

IFAIL = 4

In setting up the ODE system, the internal initialization routine was unable to initialize the derivative of the ODE system. This could be due to the fact that IRES was repeatedly set to 3 in at least one of the user-supplied (sub)programs PDEDEF, BNDARY or ODEDEF, when the residual in the underlying ODE solver was being evaluated.

IFAIL = 5

In solving the ODE system, a singular Jacobian has been encountered. You should check your problem formulation.

IFAIL = 6

When evaluating the residual in solving the ODE system, IRES was set to 2 in at least one of the user-supplied (sub)programs PDEDEF, BNDARY or ODEDEF. Integration was successful as far as $t = TS$.

IFAIL = 7

The values of ATOL and RTOL are so small that the routine is unable to start the integration in time.

IFAIL = 8

In one of the user-supplied (sub)programs, PDEDEF, BNDARY or ODEDEF, IRES was set to an invalid value.

IFAIL = 9 (D02NNF)

A serious error has occurred in an internal call to the specified routine. Check the problem specification and all parameters and array dimensions. Setting ITRACE = 1 may provide more information. If the problem persists, contact NAG.

IFAIL = 10

The required task has been completed, but it is estimated that a small change in ATOL and RTOL is unlikely to produce any change in the computed solution. (Only applies when you are not operating in one step mode, that is when ITASK \neq 2 or 5.)

IFAIL = 11

An error occurred during Jacobian formulation of the ODE system (a more detailed error description may be directed to the current error message unit). If using the sparse matrix algebra option, the values of ALGOPT(29) and ALGOPT(30) may be inappropriate.

IFAIL = 12

In solving the ODE system, the maximum number of steps specified in ALGOPT(15) have been taken.

IFAIL = 13

Some error weights w_i became zero during the time integration (see the description of ITOL). Pure relative error control ($ATOL(i) = 0.0$) was requested on a variable (the i th) which has become zero. The integration was successful as far as $t = TS$.

IFAIL = 14

The flux function R_i was detected as depending on time derivatives, which is not permissible.

IFAIL = 15

When using the sparse option, the value of LISAVE or LRSAVE was not sufficient (more detailed information may be directed to the current error message unit).

IFAIL = 16

REMESH has been changed between calls to D03PPF/D03PPA, which is not permissible.

IFAIL = 17

The remeshing process has produced zero or negative mesh spacing. You are advised to check the user-supplied function MONITF and to try adjusting the values of DXMESH, XRATIO and CON. Setting IPMINF = 1 may provide more information.

7 Accuracy

D03PPF/D03PPA controls the accuracy of the integration in the time direction but not the accuracy of the approximation in space. The spatial accuracy depends on both the number of mesh points and on their distribution in space. In the time integration only the local error over a single step is controlled and so the accuracy over a number of steps cannot be guaranteed. You should therefore test the effect of varying the accuracy parameters, ATOL and RTOL.

8 Further Comments

The parameter specification allows you to include equations with only first-order derivatives in the space direction but there is no guarantee that the method of integration will be satisfactory for such systems. The position and nature of the boundary conditions in particular are critical in defining a stable problem. It may be advisable in such cases to reduce the whole system to first-order and to use the Keller box scheme routine D03PRF.

The time taken depends on the complexity of the parabolic system, the accuracy requested, and the frequency of the mesh updates. For a given system with fixed accuracy and mesh-update frequency it is approximately proportional to NEQN.

9 Example

This example uses Burgers Equation, a common test problem for remeshing algorithms, given by

$$\frac{\partial U}{\partial t} = -U \frac{\partial U}{\partial x} + E \frac{\partial^2 U}{\partial x^2},$$

for $x \in [0, 1]$ and $t \in [0, 1]$, where E is a small constant.

The initial and boundary conditions are given by the exact solution

$$U(x, t) = \frac{0.1 \exp(-A) + 0.5 \exp(-B) + \exp(-C)}{\exp(-A) + \exp(-B) + \exp(-C)},$$

where

$$\begin{aligned} A &= \frac{50}{E}(x - 0.5 + 4.95t), \\ B &= \frac{250}{E}(x - 0.5 + 0.75t), \\ C &= \frac{500}{E}(x - 0.375). \end{aligned}$$

9.1 Program Text

Note: *the following program illustrates the use of D03PPF. An equivalent program illustrating the use of D03PPA is available with the supplied Library and is also available from the NAG web site.*

```

*      D03PPF Example Program Text
*      Mark 16 Release. NAG Copyright 1993.
*      .. Parameters ..
      INTEGER          NOUT
      PARAMETER       (NOUT=6)
      INTEGER          NPDE, NPTS, NCODE, M, NXI, NXFIX, NEQN, NIW,
+
      NWKRES, LENODE, NW, INTPTS, ITYPE
      PARAMETER       (NPDE=1,NPTS=61,NCODE=0,M=0,NXI=0,NXFIX=0,
+
      NEQN=NPDE*NPTS+NCODE,NIW=25+NXFIX,
+
      NWKRES=NPDE*(NPTS+3*NPDE+21)+7*NPTS+NXFIX+3,
+
      LENODE=11*NEQN+50,NW=NEQN*NEQN+NEQN+NWKRES+
+
      LENODE,INTPTS=5,ITYPE=1)
*      .. Scalars in Common ..
      DOUBLE PRECISION E
*      .. Local Scalars ..
      DOUBLE PRECISION CONST, DXMESH, TOUT, TRMESH, TS, XRATIO
      INTEGER          I, IFAIL, IND, IPMINF, IT, ITASK, ITOL, ITRACE,
+
      NRMESH
      LOGICAL          REMESH, THETA
      CHARACTER        LAOPT, NORM
*      .. Local Arrays ..
      DOUBLE PRECISION ALGOPT(30), ATOL(1), RTOL(1), U(NEQN),
+
      UE(INTPTS), UOUT(NPDE,INTPTS,ITYPE), W(NW),
+
      X(NPTS), XFIX(1), XI(1), XOUT(INTPTS)
      INTEGER          IW(NIW)
*      .. External Subroutines ..
      EXTERNAL         BNDARY, D03PCK, D03PPF, D03PZF, EXACT, MONITF,
+
      PDEDEF, UVINIT
*      .. Common blocks ..
      COMMON           /EPS/E
*      .. Executable Statements ..
      WRITE (NOUT,*) 'D03PPF Example Program Results'
      E = 0.005D0
      ITRACE = 0
      ITOL = 1
      ATOL(1) = 0.5D-4
      RTOL(1) = ATOL(1)
      WRITE (NOUT,99998) ATOL, NPTS
*
*      Initialise mesh ..
*
      DO 20 I = 1, NPTS
          X(I) = (I-1.0D0)/(NPTS-1.0D0)
20 CONTINUE
*
*      Set remesh parameters..
*
      REMESH = .TRUE.
      NRMESH = 3
      DXMESH = 0.5D0
      CONST = 2.0D0/(NPTS-1.0D0)
      XRATIO = 1.5D0
      IPMINF = 0
*
      WRITE (NOUT,99993) NRMESH
      WRITE (NOUT,99992) E
      WRITE (NOUT,*)
*
      XI(1) = 0.0D0
      NORM = 'A'
      LAOPT = 'F'
      IND = 0
      ITASK = 1
*
*      Set THETA to .TRUE. if the Theta integrator is required
*

```



```

      THETA = .FALSE.
      DO 40 I = 1, 30
        ALGOPT(I) = 0.0D0
40 CONTINUE
      IF (THETA) THEN
        ALGOPT(1) = 2.0D0
      ELSE
        ALGOPT(1) = 0.0D0
      END IF
*
*   Loop over output value of t
*
      TS = 0.0D0
      TOUT = 0.0D0
      DO 60 IT = 1, 5
        TOUT = 0.2D0*IT
        IFAIL = 0
*
        CALL D03PPF(NPDE,M,TS,TOUT,PDEDEF,BNDARY,UVINIT,U,NPTS,X,NCODE,
+           D03PCK,NXI,XI,NEQN,RTOL,ATOL,ITOL,NORM,LAOPT,
+           ALGOPT,REMESH,NXFIX,XFIX,NRMESH,DXMESH,TRMESH,
+           IPMINF,XRATIO,CONST,MONITF,W,NW,IW,NIW,ITASK,
+           ITRACE,IND,IFAIL)
*
*   Set output points ..
      IF (IT.EQ.1) THEN
        XOUT(1) = 0.3D0
        XOUT(2) = 0.4D0
        XOUT(3) = 0.5D0
        XOUT(4) = 0.6D0
        XOUT(5) = 0.7D0
      ELSE IF (IT.EQ.2) THEN
        XOUT(1) = 0.4D0
        XOUT(2) = 0.5D0
        XOUT(3) = 0.6D0
        XOUT(4) = 0.7D0
        XOUT(5) = 0.8D0
      ELSE IF (IT.EQ.3) THEN
        XOUT(1) = 0.6D0
        XOUT(2) = 0.65D0
        XOUT(3) = 0.7D0
        XOUT(4) = 0.75D0
        XOUT(5) = 0.8D0
      ELSE IF (IT.EQ.4) THEN
        XOUT(1) = 0.7D0
        XOUT(2) = 0.75D0
        XOUT(3) = 0.8D0
        XOUT(4) = 0.85D0
        XOUT(5) = 0.9D0
      ELSE IF (IT.EQ.5) THEN
        XOUT(1) = 0.8D0
        XOUT(2) = 0.85D0
        XOUT(3) = 0.9D0
        XOUT(4) = 0.95D0
        XOUT(5) = 1.0D0
      END IF
*
      WRITE (NOUT,99999) TS
      WRITE (NOUT,99996) (XOUT(I),I=1,INTPTS)
*   Interpolate at output points ..
      CALL D03PZF(NPDE,M,U,NPTS,X,XOUT,INTPTS,ITYPE,UOUT,IFAIL)
*
*   Check against exact solution ..
      CALL EXACT(TS,XOUT,INTPTS,UE)
*
      WRITE (NOUT,99995) (UOUT(1,I,1),I=1,INTPTS)
      WRITE (NOUT,99994) (UE(I),I=1,INTPTS)
*
60 CONTINUE
      WRITE (NOUT,99997) IW(1), IW(2), IW(3), IW(5)
      STOP

```

```

*
99999 FORMAT (' T = ',F6.3)
99998 FORMAT (/' Accuracy requirement =',E10.3,' Number of points = ',
+ I3,/)
99997 FORMAT (' Number of integration steps in time = ',I6,/' Number o',
+ 'f function evaluations = ',I6,/' Number of Jacobian eval',
+ 'uations = ',I6,/' Number of iterations = ',I6)
99996 FORMAT (1X,'X ',F9.4)
99995 FORMAT (1X,'Approx sol. ',F9.4)
99994 FORMAT (1X,'Exact sol. ',F9.4,/)
99993 FORMAT (2X,'Remeshing every',I3,' time steps',/)
99992 FORMAT (2X,'E =',F8.3)
END

*
SUBROUTINE UVINIT(NPDE,NPTS,NXI,X,XI,U,NCODE,V)
*
.. Scalar Arguments ..
INTEGER NCODE, NPDE, NPTS, NXI
*
.. Array Arguments ..
DOUBLE PRECISION U(NPDE,NPTS), V(*), X(NPTS), XI(*)
*
.. Scalars in Common ..
DOUBLE PRECISION E
*
.. Local Scalars ..
DOUBLE PRECISION A, B, C, T
INTEGER I
*
.. Intrinsic Functions ..
INTRINSIC EXP
*
.. Common blocks ..
COMMON /EPS/E
*
.. Executable Statements ..
T = 0.0D0
DO 20 I = 1, NPTS
  A = (X(I)-0.25D0-0.75D0*T)/(4.0D0*E)
  B = (0.9D0*X(I)-0.325D0-0.495D0*T)/(2.0D0*E)
  IF (A.GT.0.0D0 .AND. A.GT.B) THEN
    A = EXP(-A)
    C = (0.8D0*X(I)-0.4D0-0.24D0*T)/(4.0D0*E)
    C = EXP(C)
    U(1,I) = (0.5D0+0.1D0*C+A)/(1.0D0+C+A)
  ELSE IF (B.GT.0.0D0 .AND. B.GE.A) THEN
    B = EXP(-B)
    C = (-0.8D0*X(I)+0.4D0+0.24D0*T)/(4.0D0*E)
    C = EXP(C)
    U(1,I) = (0.1D0+0.5D0*C+B)/(1.0D0+C+B)
  ELSE
    A = EXP(A)
    B = EXP(B)
    U(1,I) = (1.0D0+0.5D0*A+0.1D0*B)/(1.0D0+A+B)
  END IF
20 CONTINUE
RETURN
END

*
SUBROUTINE PDEDEF(NPDE,T,X,U,UX,NCODE,V,VDOT,P,Q,R,IRES)
*
.. Scalar Arguments ..
DOUBLE PRECISION T, X
INTEGER IRES, NCODE, NPDE
*
.. Array Arguments ..
DOUBLE PRECISION P(NPDE,NPDE), Q(NPDE), R(NPDE), U(NPDE),
+ UX(NPDE), V(*), VDOT(*)
*
.. Scalars in Common ..
DOUBLE PRECISION E
*
.. Common blocks ..
COMMON /EPS/E
*
.. Executable Statements ..
P(1,1) = 1.0D0
R(1) = E*UX(1)
Q(1) = U(1)*UX(1)
RETURN
END

*
SUBROUTINE BNDARY(NPDE,T,U,UX,NCODE,V,VDOT,IBND,BETA,GAMMA,IRES)

```

```

*   .. Scalar Arguments ..
DOUBLE PRECISION T
INTEGER          IBND, IRES, NCODE, NPDE
*   .. Array Arguments ..
DOUBLE PRECISION BETA(NPDE), GAMMA(NPDE), U(NPDE), UX(NPDE),
+   V(*), VDOT(*)
*   .. Scalars in Common ..
DOUBLE PRECISION E
*   .. Local Scalars ..
DOUBLE PRECISION A, B, C, UE, X
*   .. Intrinsic Functions ..
INTRINSIC        EXP
*   .. Common blocks ..
COMMON          /EPS/E
*   .. Executable Statements ..
BETA(1) = 0.0D0
IF (IBND.EQ.0) THEN
  X = 0.0D0
  A = (X-0.25D0-0.75D0*T)/(4.0D0*E)
  B = (0.9D0*X-0.325D0-0.495D0*T)/(2.0D0*E)
  IF (A.GT.0.0D0 .AND. A.GT.B) THEN
    A = EXP(-A)
    C = (0.8D0*X-0.4D0-0.24D0*T)/(4.0D0*E)
    C = EXP(C)
    UE = (0.5D0+0.1D0*C+A)/(1.0D0+C+A)
  ELSE IF (B.GT.0.0D0 .AND. B.GE.A) THEN
    B = EXP(-B)
    C = (-0.8D0*X+0.4D0+0.24D0*T)/(4.0D0*E)
    C = EXP(C)
    UE = (0.1D0+0.5D0*C+B)/(1.0D0+C+B)
  ELSE
    A = EXP(A)
    B = EXP(B)
    UE = (1.0D0+0.5D0*A+0.1D0*B)/(1.0D0+A+B)
  END IF
ELSE
  X = 1.0D0
  A = (X-0.25D0-0.75D0*T)/(4.0D0*E)
  B = (0.9D0*X-0.325D0-0.495D0*T)/(2.0D0*E)
  IF (A.GT.0.0D0 .AND. A.GT.B) THEN
    A = EXP(-A)
    C = (0.8D0*X-0.4D0-0.24D0*T)/(4.0D0*E)
    C = EXP(C)
    UE = (0.5D0+0.1D0*C+A)/(1.0D0+C+A)
  ELSE IF (B.GT.0.0D0 .AND. B.GE.A) THEN
    B = EXP(-B)
    C = (-0.8D0*X+0.4D0+0.24D0*T)/(4.0D0*E)
    C = EXP(C)
    UE = (0.1D0+0.5D0*C+B)/(1.0D0+C+B)
  ELSE
    A = EXP(A)
    B = EXP(B)
    UE = (1.0D0+0.5D0*A+0.1D0*B)/(1.0D0+A+B)
  END IF
END IF
GAMMA(1) = U(1) - UE
RETURN
END

*
SUBROUTINE EXACT(T,X,NPTS,U)
*   Exact solution (for comparison purposes)
*   .. Scalar Arguments ..
DOUBLE PRECISION T
INTEGER          NPTS
*   .. Array Arguments ..
DOUBLE PRECISION U(NPTS), X(NPTS)
*   .. Scalars in Common ..
DOUBLE PRECISION E
*   .. Local Scalars ..
DOUBLE PRECISION A, B, C
INTEGER          I

```

```

* .. Intrinsic Functions ..
INTRINSIC      EXP
* .. Common blocks ..
COMMON        /EPS/E
* .. Executable Statements ..
DO 20 I = 1, NPTS
  A = (X(I)-0.25D0-0.75D0*T)/(4.0D0*E)
  B = (0.9D0*X(I)-0.325D0-0.495D0*T)/(2.0D0*E)
  IF (A.GT.0.0D0 .AND. A.GT.B) THEN
    A = EXP(-A)
    C = (0.8D0*X(I)-0.4D0-0.24D0*T)/(4.0D0*E)
    C = EXP(C)
    U(I) = (0.5D0+0.1D0*C+A)/(1.0D0+C+A)
  ELSE IF (B.GT.0.0D0 .AND. B.GE.A) THEN
    B = EXP(-B)
    C = (-0.8D0*X(I)+0.4D0+0.24D0*T)/(4.0D0*E)
    C = EXP(C)
    U(I) = (0.1D0+0.5D0*C+B)/(1.0D0+C+B)
  ELSE
    A = EXP(A)
    B = EXP(B)
    U(I) = (1.0D0+0.5D0*A+0.1D0*B)/(1.0D0+A+B)
  END IF
20 CONTINUE
RETURN
END

*
SUBROUTINE MONITF(T,NPTS,NPDE,X,U,R,FMON)
* .. Scalar Arguments ..
DOUBLE PRECISION T
INTEGER          NPDE, NPTS
* .. Array Arguments ..
DOUBLE PRECISION FMON(NPTS), R(NPDE,NPTS), U(NPDE,NPTS), X(NPTS)
* .. Local Scalars ..
DOUBLE PRECISION DRDX, H
INTEGER          I, K, L
* .. Intrinsic Functions ..
INTRINSIC        ABS, MAX, MIN
* .. Executable Statements ..
DO 20 I = 1, NPTS - 1
  K = MAX(1,I-1)
  L = MIN(NPTS,I+1)
  H = (X(L)-X(K))*0.5D0
* Second derivative ..
  DRDX = (R(1,I+1)-R(1,I))/H
  FMON(I) = ABS(DRDX)
20 CONTINUE
FMON(NPTS) = FMON(NPTS-1)
RETURN
END

```

9.2 Program Data

None.

9.3 Program Results

D03PPF Example Program Results

Accuracy requirement = 0.500E-04 Number of points = 61

Remeshing every 3 time steps

E = 0.005

T = 0.200

X	0.3000	0.4000	0.5000	0.6000	0.7000
Approx sol.	0.9968	0.7448	0.4700	0.1667	0.1018
Exact sol.	0.9967	0.7495	0.4700	0.1672	0.1015

T = 0.400					
X	0.4000	0.5000	0.6000	0.7000	0.8000
Approx sol.	1.0003	0.9601	0.4088	0.1154	0.1005
Exact sol.	0.9997	0.9615	0.4094	0.1157	0.1003
T = 0.600					
X	0.6000	0.6500	0.7000	0.7500	0.8000
Approx sol.	0.9966	0.9390	0.3978	0.1264	0.1037
Exact sol.	0.9964	0.9428	0.4077	0.1270	0.1033
T = 0.800					
X	0.7000	0.7500	0.8000	0.8500	0.9000
Approx sol.	1.0003	0.9872	0.5450	0.1151	0.1010
Exact sol.	0.9996	0.9878	0.5695	0.1156	0.1008
T = 1.000					
X	0.8000	0.8500	0.9000	0.9500	1.0000
Approx sol.	1.0001	0.9961	0.7324	0.1245	0.1004
Exact sol.	0.9999	0.9961	0.7567	0.1273	0.1004
Number of integration steps in time =	205				
Number of function evaluations =	4872				
Number of Jacobian evaluations =	71				
Number of iterations =	518				
