

nag_ode_ivp_rk_range (d02pcc)

1. Purpose

nag_ode_ivp_rk_range (d02pcc) is a function for solving the initial value problem for a first order system of ordinary differential equations using Runge-Kutta methods.

2. Specification

```
#include <nag.h>
#include <nagd02.h>

void nag_ode_ivp_rk_range(Integer neq,
    void (*f) (Integer neq, double t, double y[],
               double yp[], Nag_User *comm),
    double twant, double *tgot,
    double ygot[], double ypgot[], double ymax[],
    Nag_ODE_RK *opt, Nag_User *comm, NagError *fail)
```

3. Description

This function and its associated functions (**nag_ode_ivp_rk_setup (d02pvc)**, **nag_ode_ivp_rk_errass (d02pzc)**) solve the initial value problem for a first order system of ordinary differential equations. The functions, based on Runge-Kutta methods and derived from RKSUITE (Brankin *et al*, 1991) integrate

$$y' = f(t, y) \quad \text{given} \quad y(t_0) = y_0$$

where y is the vector of **neq** solution components and t is the independent variable.

This function is designed for the usual task, namely to compute an approximate solution at a sequence of points. You must first call **nag_ode_ivp_rk_setup (d02pvc)** to specify the problem and how it is to be solved. Thereafter you call **nag_ode_ivp_rk_range** repeatedly with successive values of **twant**, the points at which you require the solution, in the range from **tstart** to **tend** (as specified in **nag_ode_ivp_rk_setup (d02pvc)**). In this manner **nag_ode_ivp_rk_range** returns the point at which it has computed a solution **tgot** (usually **twant**), the solution there **ygot** and its derivative **ypgot**. If **nag_ode_ivp_rk_range** encounters some difficulty in taking a step toward **twant**, then it returns the point of difficulty **tgot** and the solution and derivative computed there **ygot** and **ypgot**.

In the call to **nag_ode_ivp_rk_setup (d02pvc)** you can specify the first step size for **nag_ode_ivp_rk_range** to attempt or that it compute automatically an appropriate value. Thereafter **nag_ode_ivp_rk_range** estimates an appropriate step size for its next step. This value and other details of the integration can be obtained after any call to **nag_ode_ivp_rk_range** by examining the contents of the structure **opt**, see Section 4. The local error is controlled at every step as specified in **nag_ode_ivp_rk_setup (d02pvc)**. If you wish to assess the true error, you must set **errass = Nag_ErrorAssess_on** in the call to **nag_ode_ivp_rk_setup (d02pvc)**. This assessment can be obtained after any call to **nag_ode_ivp_rk_range** by a call to the function **nag_ode_ivp_rk_errass (d02pzc)**.

For more complicated tasks, you are referred to functions **nag_ode_ivp_rk_onestep (d02pdc)**, **nag_ode_ivp_rk_interp (d02pxc)** and **nag_ode_ivp_rk_reset_tend (d02pwc)**.

4. Parameters

neq

Input: the number of ordinary differential equations in the system to be solved.

Constraint: **neq** ≥ 1.

f

This function must evaluate the first derivatives y'_i (that is the functions f_i) for given values of the arguments t, y_i .

```
void f (Integer neq, double t, double y[], double yp[], Nag_User *comm)
```

neq
Input: the number of differential equations.

t
Input: the current value of the independent variable, t .

y[neq]
Input: the current values of the dependent variables, y_i for $i = 1, 2, \dots, \mathbf{neq}$.

yp[neq]
Output: the values of f_i for $i = 1, 2, \dots, \mathbf{neq}$.

comm
Input/Output: pointer to a structure of type Nag_User with the following member:

p - Pointer
Input/Output: The pointer **comm**->**p** should be cast to the required type, e.g. `struct user *s = (struct user *)comm->p`, to obtain the original object's address with appropriate type. (See the argument **comm** below.)

twant

Input: the next value of the independent variable, t , where a solution is desired.

Constraints: **twant** must be closer to **tend** than the previous of **tgot** (or **tstart** on the first call to nag_ode_ivp_rk_range); see nag_ode_ivp_rk_setup (d02pvc) for a description of **tstart** and **tend**. **twant** must not lie beyond **tend** in the direction of integration.

tgot

Output: the value of the independent variable t at which a solution has been computed. On successful exit with fail.code = **NE_NOERROR**, **tgot** will equal **twant**. For non-trivial values of fail.code (i.e., those not related to an invalid call of nag_ode_ivp_rk_range) a solution has still been computed at the value of **tgot** but in general **tgot** will not equal **twant**.

ygot[neq]

Input: on the first call to nag_ode_ivp_rk_range, **ygot** need not be set. On all subsequent calls **ygot** must remain unchanged.

Output: an approximation to the true solution at the value of **tgot**. At each step of the integration to **tgot**, the local error has been controlled as specified in nag_ode_ivp_rk_setup (d02pvc). The local error has still been controlled even when **tgot** \neq **twant**, that is after a return with a non-trivial error.

ypgot[neq]

Output: an approximation to the first derivative of the true solution at **tgot**.

ymax[neq]

Input: on the first call to nag_ode_ivp_rk_range, **ymax** need not be set. On all subsequent calls **ymax** must remain unchanged.

Output: **ymax**[$i-1$] contains the largest value of $|y_i|$ computed at any step in the integration so far.

opt

Input: pointer to a structure of type Nag_ODE_RK as initialised by the setup function nag_ode_ivp_rk_setup (d02pvc).

Output: the following structure members hold information as follows:

totfcn - Integer

The total number of evaluations of f used in the primary integration so far; this does not include evaluations of f for the secondary integration specified by a prior call to nag_ode_ivp_rk_setup (d02pvc) with errass = **Nag_ErrorAssess_on**.

stpcst - Integer

The cost in terms of number of evaluations of f of a typical step with the method being

used for the integration. The method is specified by the parameter **method** in a prior call to `nag_ode_ivp_rk_setup` (d02pvc).

waste - double

The number of attempted steps that failed to meet the local error requirement divided by the total number of steps attempted so far in the integration. A “large” fraction indicates that the integrator is having trouble with the problem being solved. This can happen when the problem is “stiff” and also when the solution has discontinuities in a low order derivative.

stpsok - Integer

The number of accepted steps.

hnext - double

The step size the integrator plans to use for the next step.

comm

Input/Output: pointer to a structure of type `Nag_User` with the following member:

p - Pointer

Input/Output: The pointer **p**, of type `Pointer`, allows the user to communicate information to and from the user-defined function `f()`. An object of the required type should be declared by the user, e.g. a structure, and its address assigned to the pointer **p** by means of a cast to `Pointer` in the calling program, e.g. `comm.p = (Pointer)&s`. The type pointer will be `void *` with a C compiler that defines `void *` and `char *` otherwise.

fail

The NAG error parameter, see the Essential Introduction to the NAG C Library.

5. Error Indications and Warnings

NE_PREV_CALL

The previous call to a function had resulted in a severe error. You must call `nag_ode_ivp_rk_setup` (d02pvc) to start another problem.

NE_NO_SETUP

The setup function `nag_ode_ivp_rk_setup` (d02pvc) has not been called.

NE_RK_INVALID_CALL

The function to be called as specified in the setup routine `nag_ode_ivp_rk_setup` (d02pvc) was `nag_ode_ivp_rk_onestep` (d02pdc). However the actual call was made to `nag_ode_ivp_rk_range`. This is not permitted.

NE_PREV_CALL_INI

The previous call to the function `nag_ode_ivp_rk_range` had resulted in a severe error. You must call `nag_ode_ivp_rk_setup` (d02pvc) to start another problem.

NE_NEQ

The value of **neq** supplied is not the same as that given to the setup function `nag_ode_ivp_rk_setup` (d02pvc).

neq = *<value>* but the value given to `nag_ode_ivp_rk_setup` (d02pvc) was *<value>*.

NE_RK_TGOT_EQ_TEND

The call to `nag_ode_ivp_rk_range` has been made after reaching **tend**. The previous call to `nag_ode_ivp_rk_range` resulted in **tgot** (**tstart** on the first call) = **tend**. You must call `nag_ode_ivp_rk_setup` (d02pvc) to start another problem.

NE_RK_TGOT_RANGE_TEND

The call to `nag_ode_ivp_rk_range` has been made with a **twant** that does not lie between the previous value of **tgot** (**tstart** on the first call) and **tend**. This is not permitted.

NE_RK_TGOT_RANGE_TEND_CLOSE

The call to `nag_ode_ivp_rk_range` has been made with a **twant** that does not lie between the previous value of **tgot** (**tstart** on the first call) and **tend**. This is not permitted. However **twant** is very close to **tend**, so you may have meant it to be **tend** exactly. Check your program.

NE_RK_TWANT_CLOSE_TGOT

The call to nag_ode_ivp_rk_range has been made with a **twant** that is not sufficiently different from the last value of **tgot** (**tstart** on the first call). When using **method = Nag_RK_7_8**, it must differ by at least $\langle value \rangle$.

NE_RK_PDC_STEP

In order to satisfy the error requirements nag_ode_ivp_rk_range would have to use a step size of $\langle value \rangle$ at current **t** = $\langle value \rangle$. This is too small for the machine precision.

NE_RK_PDC_GLOBAL_ERROR_T

The global error assessment may not be reliable for t past **tgot**. **tgot** = $\langle value \rangle$.

NE_RK_PDC_GLOBAL_ERROR_S

The global error assessment algorithm failed at the start of the integration.

NE_STIFF_PROBLEM

The problem appears to be stiff.

NW_RK_TOO_MANY

Approximately $\langle value \rangle$ function evaluations have been used to compute the solution since the integration started or since this message was last printed.

NE_RK_PCC_METHOD

The efficiency of the integration has been degraded. Consider calling the set up function nag_ode_ivp_rk_setup (d02pvc) to re-initialize the integration at the current point with the method changed to **NE_RK_4_5**. Alternatively nag_ode_ivp_rk_range (d02pcc) can be called again to resume at the current point.

NE_INTERNAL_ERROR

An internal error has occurred in this function. Check the function call and any array sizes. If the call is correct then please consult NAG for assistance.

NE_MEMORY_FREED

Internally allocated memory has been freed by a call to nag_ode_ivp_rk_free (d02ppc) without a subsequent call to the set up function nag_ode_ivp_rk_setup (d02pvc).

6. Further Comments

If nag_ode_ivp_rk_range returns with fail.code = **NE_RK_PDC_STEP** and the accuracy specified by **tol** and **thres** is really required then you should consider whether there is a more fundamental difficulty. For example, the solution may contain a singularity. In such a region the solution components will usually be of a large magnitude. Successive output values of **ygot** and **ymax** should be monitored (or the routine nag_ode_ivp_rk_onestep (d02pdc) should be used since this takes one integration step at a time) with the aim of trapping the solution before the singularity. In any case numerical solution cannot be continued through a singularity, and analytical treatment may be necessary.

Performance statistics are available after any return from nag_ode_ivp_rk_range by examining the structure **opt** see Section 4. If **errass** was set to **Nag_ErrorAssess_on** in the call to nag_ode_ivp_rk_setup (d02pvc), global error assessment is available after any return from nag_ode_ivp_rk_range (except when the error is due to incorrect input arguments or incorrect setup) by a call to the routine nag_ode_ivp_rk_errass (d02pzc). The approximate extra number of evaluations of f used is given by $2 \times \mathbf{stpsok} \times \mathbf{stpcst}$ for method **NAG_RK_4_5** or **NAG_RK_7_8** and $3 \times \mathbf{stpsok} \times \mathbf{stpcst}$ for method = **NAG_RK_2_3**.

After a failure with fail.code = **NE_RK_PDC_STEP**, **NE_RK_PDC_GLOBAL_ERROR_T** or **NE_RK_PDC_GLOBAL_ERROR_S** the diagnostic routine nag_ode_ivp_rk_errass (d02pzc) may be called only once.

If nag_ode_ivp_rk_range returns with fail.code = **NE_STIFF_PROBLEM** then it is advisable to change to another code more suited to the solution of stiff problems. nag_ode_ivp_rk_range will not return with fail.code = **NE_STIFF_PROBLEM** if the problem is actually stiff but it is estimated that integration can be completed using less function evaluations than already computed.

6.1. Accuracy

The accuracy of integration is determined by the parameters **tol** and **thres** in a prior call to `nag_ode_ivp_rk_setup` (d02pvc). Note that only the local error at each step is controlled by these parameters. The error estimates obtained are not strict bounds but are usually reliable over one step. Over a number of steps the overall error may accumulate in various ways, depending on the properties of the differential system.

6.2. References

Brankin R W, Gladwell I and Shampine L F (1991) *RKSUITE: a suite of Runge-Kutta codes for the initial value problem for ODEs* SoftReport 91-S1, Department of Mathematics, Southern Methodist University, Dallas, TX 75275, U.S.A.

7. See Also

`nag_ode_ivp_adams_gen` (d02cjc)
`nag_ode_ivp_adams_roots` (d02qfc)
`nag_ode_ivp_rk_setup` (d02pvc)
`nag_ode_ivp_rk_errass` (d02pzc)

8. Example

We solve the equation

$$y'' = -y, \quad y(0) = 0, y'(0) = 1$$

reposed as

$$y'_1 = y_2 \quad y'_2 = -y_1$$

over the range $[0, 2\pi]$ with initial conditions $y_1 = 0.0$ and $y_2 = 1.0$. We use relative error control with threshold values of $1.0e-8$ for each solution component and compute the solution at intervals of length $\pi/4$ across the range. We use a low order Runge-Kutta method (**method** = **Nag_RK_2.3**) with tolerances **tol** = $1.0e-3$ and **tol** = $1.0e-4$ in turn so that we may compare the solutions. The value of π is obtained by using X01AAC.

See also the example program for `nag_ode_ivp_rk_errass` (d02pzc).

8.1. Program Text

```
/* nag_ode_ivp_rk_range(d02pcc) Example Program
 *
 * Copyright 1994 Numerical Algorithms Group.
 *
 * Mark 3, 1994.
 *
 */

#include <nag.h>
#include <math.h>
#include <stdio.h>
#include <nag_stdlib.h>
#include <nagd02.h>
#include <nagx01.h>

#ifdef NAG_PROTO
static void f(Integer neq, double t1, double y[], double yp[], Nag_User *comm);
#else
static void f();
#endif

#define NEQ 2
#define ZERO 0.0
#define ONE 1.0
#define TWO 2.0
```

```

#define FOUR 4.0

main()
{
    Integer neq;
    Nag_RK_method method;
    double hstart, pi, tgot, tend, tinc;
    double tol, tstart, twant;
    Integer i, j, nout;
    double thres[NEQ], ygot[NEQ], ymax[NEQ], ypgot[NEQ], ystart[NEQ];
    Nag_ErrorAssess errass;
    Nag_ODE_RK opt;
    Nag_User comm;

    Vprintf("d02pcc Example Program Results\n");

    /* Set initial conditions and input for d02pvc */
    neq = NEQ;
    pi = X01AAC;
    tstart = ZERO;
    ystart[0] = ZERO;
    ystart[1] = ONE;
    tend = TWO*pi;
    for (i=0; i<neq; i++)
        thres[i] = 1.0e-8;
    errass = Nag_ErrorAssess_off;
    hstart = ZERO;
    method = Nag_RK_2_3;

    /*
     * Set control for output
     */
    nout = 8;
    tinc = (tend-tstart)/nout;

    for (i=1; i<=2; i++)
    {
        if (i==1) tol = 1.0e-3;
        if (i==2) tol = 1.0e-4;
        d02pvc(neq, tstart, ystart, tend, tol, thres, method,
            Nag_RK_range, errass, hstart, &opt, NAGERR_DEFAULT);

        Vprintf("\nCalculation with tol = %8.1e\n\n",tol);
        Vprintf("      t      y1      y2\n\n");
        Vprintf("%8.3f %8.3f %8.3f\n", tstart, ystart[0], ystart[1]);
        for (j=nout-1; j>=0; j--)
        {
            twant = tend - j*tinc;
            d02pcc(neq, f, twant, &tgot, ygot, ypgot, ymax, &opt, &comm,
                NAGERR_DEFAULT);
            Vprintf("%8.3f %8.3f %8.3f\n", tgot, ygot[0], ygot[1]);
        }
        Vprintf("\nCost of the integration in evaluations of f is %ld\n\n",
            opt.totfcn);
        d02ppc(&opt);
    }
    exit(EXIT_SUCCESS);
}
#ifdef NAG_PROTO
static void f(Integer neq, double t, double y[], double yp[], Nag_User *comm)
#else
static void f(neq, t, y, yp, comm)
    Integer neq;
    double t;
    double y[], yp[];
    Nag_User *comm;
#endif

{
    yp[0] = y[1];

```

```

    yp[1] = -y[0];
}

```

8.2. Program Data

None.

8.3. Program Results

d02pcc Example Program Results

Calculation with tol = 1.0e-03

t	y1	y2
0.000	0.000	1.000
0.785	0.707	0.707
1.571	0.999	0.000
2.356	0.706	-0.706
3.142	0.000	-0.999
3.927	-0.706	-0.706
4.712	-0.998	0.000
5.498	-0.705	0.706
6.283	0.001	0.997

Cost of the integration in evaluations of f is 124

Calculation with tol = 1.0e-04

t	y1	y2
0.000	0.000	1.000
0.785	0.707	0.707
1.571	1.000	0.000
2.356	0.707	-0.707
3.142	0.000	-1.000
3.927	-0.707	-0.707
4.712	-1.000	0.000
5.498	-0.707	0.707
6.283	0.000	1.000

Cost of the integration in evaluations of f is 235
