## nag\_ode\_ivp\_adams\_gen (d02cjc)

## 1. Purpose

**nag\_ode\_ivp\_adams\_gen (d02cjc)** integrates a system of first order ordinary differential equations over a range with suitable initial conditions, using a variable-order, variable-step Adams method until a user-specified function, if supplied, of the solution is zero, and returns the solution at points specified by the user, if desired.

## 2. Specification

## 3. Description

The function advances the solution of a system of ordinary differential equations

 $y'_i = f_i(x, y_1, y_2, \dots, y_{neq}), \qquad i = 1, 2, \dots, neq,$ 

from  $x = \mathbf{x}$  to  $x = \mathbf{xend}$  using a variable-order, variable-step Adams method. The system is defined by a function **fcn** supplied by the user, which evaluates  $f_i$  in terms of x and  $y_1, y_2, \ldots, y_{neq}$ . The initial values of  $y_1, y_2, \ldots, y_{neq}$  must be given at  $x = \mathbf{x}$ .

The solution is returned via the user-supplied function **output** at points specified by the user, if desired: this solution is obtained by  $C^1$  interpolation on solution values produced by the method. As the integration proceeds a check can be made on the user-specified function g(x, y) to determine an interval where it changes sign. The position of this sign change is then determined accurately. It is assumed that g(x, y) is a continuous function of the variables, so that a solution of g(x, y) = 0.0 can be determined by searching for a change in sign in g(x, y). The accuracy of the integration, the interpolation and, indirectly, of the determination of the position where g(x, y) = 0.0, is controlled by the parameters **tol** and **err\_c**.

For a description of Adams methods and their practical implementation see Hall and Watt (1976).

## 4. Parameters

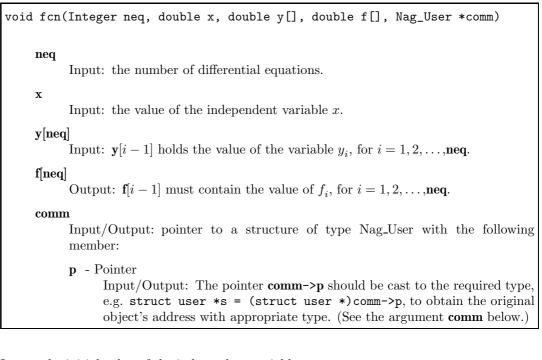
#### neq

Input: the number of differential equations. Constraint:  $\mathbf{neq} \ge 1$ .

#### fcn

The function **fcn**, supplied by the user, must evaluate the first derivatives  $y'_i$  (i.e., the functions  $f_i$ ) for given values of their arguments  $x, y_1, y_2, \ldots, y_{neq}$ .

The specification of **fcn** is:



x

Input: the initial value of the independent variable x.

Constraint:  $\mathbf{x} \neq \mathbf{xend}$ .

Output: if g is supplied by the user, **x** contains the point where g(x, y) = 0.0, unless  $g(x, y) \neq 0.0$  anywhere on the range **x** to **xend**, in which case, **x** will contain **xend**. If g is not supplied by the user **x** contains **xend**, unless an error has occurred, when it contains the value of x at the error.

#### y[neq]

Input: the initial values of the solution  $y_1, y_2, \ldots, y_{neq}$  at  $x = \mathbf{x}$ . Output: the computed values of the solution at the final point  $x = \mathbf{x}$ .

#### xend

Input: the final value of the independent variable. If xend < x, integration proceeds in the negative direction.

Constraint: **xend**  $\neq$  **x**.

tol

Input: a positive tolerance for controlling the error in the integration. Hence **tol** affects the determination of the position where g(x, y) = 0.0, if g is supplied.

nag\_ode\_ivp\_adams\_gen has been designed so that, for most problems, a reduction in **tol** leads to an approximately proportional reduction in the error in the solution. However, the actual relation between **tol** and the accuracy achieved cannot be guaranteed. The user is strongly recommended to call nag\_ode\_ivp\_adams\_gen with more than one value for **tol** and to compare the results obtained to estimate their accuracy. In the absence of any prior knowledge, the user might compare the results obtained by calling nag\_ode\_ivp\_adams\_gen with **tol** =  $10.0^{-p}$ and **tol** =  $10.0^{-p-1}$  where p correct decimal digits are required in the solution. Constraint: **tol** > 0.0.

err\_c

Input: the type of error control. At each step in the numerical solution an estimate of the local error, *est*, is made. For the current step to be accepted the following condition must be satisfied:

$$est = \sqrt{\sum_{i=1}^{\mathbf{neq}} (e_i/(\tau_r \times |y_i| + \tau_a))^2} \leq 1.0$$

where  $\tau_r$  and  $\tau_a$  are defined by

err_c	$ au_r$	$ au_a$
Nag_Relative	tol	ε
Nag_Absolute	0.0	tol
Nag_Mixed	tol	$\mathbf{tol}$

where  $\varepsilon$  is a small machine-dependent number and  $e_i$  is an estimate of the local error at  $y_i$ , computed internally. If the appropriate condition is not satisfied, the step size is reduced and the solution is recomputed on the current step. If the user wishes to measure the error in the computed solution in terms of the number of correct decimal places, then **err\_c** should be set to **Nag\_Absolute** on entry, whereas if the error requirement is in terms of the number of correct significant digits, then **err\_c** should be set to **Nag\_Relative**. If the user prefers a mixed error test, then **err\_c** should be set to **Nag\_Mixed**. The recommended value for **err\_c** is **Nag\_Mixed** and this should be chosen unless there are good reasons for a different choice. Constraint: **err\_c = Nag\_Relative**, **Nag\_Absolute** or **Nag\_Mixed**.

#### output

The function **output** permits access to intermediate values of the computed solution (for example to print or plot them), at successive user-specified points. It is initially called by nag\_ode\_ivp\_adams\_gen with  $\mathbf{xsol} = \mathbf{x}$  (the initial value of x). The user must reset  $\mathbf{xsol}$  to the next point (between the current  $\mathbf{xsol}$  and  $\mathbf{xend}$ ) where **output** is to be called, and so on at each call to **output**. If, after a call to **output**, the reset point  $\mathbf{xsol}$  is beyond  $\mathbf{xend}$ , nag\_ode\_ivp\_adams\_gen will integrate to  $\mathbf{xend}$  with no further calls to **output**; if a call to **output** is required at the point  $\mathbf{xsol} = \mathbf{xend}$ , then  $\mathbf{xsol}$  must be given precisely the value  $\mathbf{xend}$ . The specification of **output** is:

void output(Integer neq, double \*xsol, double y[], Nag\_User \*comm) neg Input: the number of differential equations. xsol Input: the value of the independent variable x. Output: the user must set **xsol** to the next value of x at which **output** is to be called. y[neq] Input: the computed solution at the point **xsol**. 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.)

If the user does not wish to access intermediate output, the actual argument **output** must be the NAG defined null function pointer NULLFN.

g

The function **g** must evaluate g(x, y) for specified values x, y. It specifies the function g for which the first position x where g(x, y) = 0 is to be found. The specification of **g** is: double g(Integer neq, double x, double y[], Nag\_User \*comm)
neq
Input: the number of differential equations.
x
Input: the value of the independent variable x.
y[neq]
Input: y[i - 1] holds the value of the variable y<sub>i</sub>, for i = 1, 2, ...,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.)

If the user does not require the root finding option, the actual argument  ${\bf g}$  must be the Nag defined null double function pointer NULLDFN.

### comm

Input/Output: pointer to a structure of type Nag\_User with the following member:

 ${\bf p}\,$  - Pointer

Input/Output: The pointer **p**, of type Pointer, allows the user to communicate information to and from the user-defined functions **fcn()**, **output()** and **g()**. 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\_INT\_ARG\_LT

On entry, **neq** must not be less than 1:  $\mathbf{neq} = \langle value \rangle$ .

## NE\_REAL\_ARG\_LE

On entry, **tol** must not be less than or equal to 0.0:  $\mathbf{tol} = \langle value \rangle$ .

## NE\_2\_REAL\_ARG\_EQ

On entry,  $\mathbf{x} = \langle value \rangle$  while **xend** =  $\langle value \rangle$ . These parameters must satisfy  $\mathbf{x} \neq \mathbf{xend}$ .

# NE\_BAD\_PARAM

On entry parameter **err\_c** had an illegal value.

#### NE\_TOL\_TOO\_SMALL

The value of **tol**,  $\langle value \rangle$ , is too small for the function to take an initial step.

#### NE\_XSOL\_NOT\_RESET

On call  $\langle value \rangle$  to the supplied print function **xsol** was not reset.

#### NE\_XSOL\_SET\_WRONG

**xsol** was set to a value behind  $\mathbf{x}$  in the direction of integration by the first call to the supplied print function.

The integration range is  $[\langle value1 \rangle, \langle value2 \rangle], \mathbf{xsol} = \langle value \rangle.$ 

#### NE\_XSOL\_INCONSIST

On call  $\langle value \rangle$  to the supplied print function **xsol** was set to a value behind the previous value of **xsol** in the direction of integration.

Previous  $\mathbf{xsol} = \langle value \rangle$ ,  $\mathbf{xend} = \langle value \rangle$ , new  $\mathbf{xsol} = \langle value \rangle$ .

## NE\_NO\_SIGN\_CHANGE

No change in sign of the function g(x, y) was detected in the integration range.

## NE\_TOL\_PROGRESS

The value of **tol**,  $\langle value \rangle$ , is too small for the function to make any further progress across the integration range. Current value of  $\mathbf{x} = \langle value \rangle$ .

## NE\_ALLOC\_FAIL

Memory allocation failed.

## 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.

## 6. Further Comments

If more than one root is required then nag\_ode\_ivp\_adams\_roots (d02qfc) should be used.

If the function fails with error exit of **fail.code** = **NE\_TOL\_TOO\_SMALL**, then it can be called again with a larger value of **tol** if this has not already been tried. If the accuracy requested is really needed and cannot be obtained with this function, the system may be very stiff (see below) or so badly scaled that it cannot be solved to the required accuracy.

If the function fails with error exit of **fail.code** = **NE\_TOL\_PROGRESS**, it is probable that it has been called with a value of **tol** which is so small that a solution cannot be obtained on the range  $\mathbf{x}$  to **xend**. This can happen for well-behaved systems and very small values of **tol**. The user should, however, consider whether there is a more fundamental difficulty. For example:

- (a) in the region of a singularity (infinite value) of the solution, the function will usually stop with error exit of **fail.code** = **NE\_TOL\_PROGRESS**, unless overflow occurs first. Numerical integration cannot be continued through a singularity, and analytic treatment should be considered;
- (b) for 'stiff' equations where the solution contains rapidly decaying components, the function will use very small steps in x (internally to nag\_ode\_ivp\_adams\_gen) to preserve stability. This will exhibit itself by making the computing time excessively long, or occasionally by an exit with **fail.code** = **NE\_TOL\_PROGRESS**. Adams methods are not efficient in such cases.

## 6.1. Accuracy

The accuracy of the computation of the solution vector  $\mathbf{y}$  may be controlled by varying the local error tolerance **tol**. In general, a decrease in local error tolerance should lead to an increase in accuracy. Users are advised to choose **err\_c** = **Nag\_Mixed** unless they have a good reason for a different choice.

If the problem is a root-finding one, then the accuracy of the root determined will depend on the properties of g(x, y). The user should try to code **g** without introducing any unnecessary cancellation errors.

## 6.2. References

Hall G and Watt J M (ed) (1976) Modern Numerical Methods for Ordinary Differential Equations Clarendon Press, Oxford.

## 7. See Also

nag\_ode\_ivp\_bdf\_gen (d02ejc) nag\_ode\_ivp\_rk\_range (d02pcc) nag\_ode\_ivp\_adams\_roots (d02qfc)

## 8. Example

We illustrate the solution of four different problems. In each case the differential system (for a projectile) is

$$y' = \tan \phi$$
$$v' = \frac{-0.032 \tan \phi}{v} - \frac{0.02v}{\cos \phi}$$
$$\phi' = \frac{-0.032}{v^2}$$

over an interval  $\mathbf{x} = 0.0$  to  $\mathbf{xend} = 10.0$  starting with values y = 0.5, v = 0.5 and  $\phi = \pi/5$ . We solve each of the following problems with local error tolerances 1.0e-4 and 1.0e-5.

- (i) To integrate to x = 10.0 producing output at intervals of 2.0 until a point is encountered where y = 0.0.
- (ii) As (i) but with no intermediate output.
- (iii) As (i) but with no termination on a root-finding condition.
- (iv) As (i) but with no intermediate output and no root-finding termination condition.

## 8.1. Program Text

```
/* nag_ode_ivp_adams_gen(d02cjc) Example Program
```

```
* Copyright 1991 Numerical Algorithms Group.
 * Mark 2, 1991.
 * Mark 3 revised, 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 out(Integer neq, double *xsol, double y[], Nag_User *comm);
#else
static void out();
#endif
#ifdef NAG_PROTO
static void fcn(Integer neq, double x, double y[], double f[], Nag_User *comm);
#else
static void fcn();
#endif
#ifdef NAG_PROTO
static double g(Integer neq, double x, double y[], Nag_User *comm);
#else
static double g();
#endif
struct user
{
  double xend, h;
  Integer k;
};
main()
{
```

```
Integer i, j, neq;
double x, pi, tol;
double y[3];
Nag_User comm;
struct user s;
Vprintf("d02cjc Example Program Results\n");
/* For communication with function out()
* assign address of user defined structure
* to Nag pointer.
*/
comm.p = (Pointer)&s;
neq = 3;
s.xend = 10.0;
pi = XO1AAC;
Vprintf("\nCase 1: intermediate output, root-finding\n");
for (j = 4; j <= 5; ++j)
  {
    tol = pow(10.0, (double)(-j));
    Vprintf("\n Calculation with tol = %8.1e\n", tol);
    x = 0.0;
    y[0] = 0.5;
    y[1] = 0.5;
    y[2] = pi / 5.0;
    s.k = 4;
    s.k = 4;
s.h = (s.xend - x) / (double)(s.k + 1);
v(1) Y(2)
    Vprintf("\n
                                                           Y(3) n");
    d02cjc(neq, fcn, &x, y, s.xend, tol, Nag_Mixed, out, g, &comm,
           NAGERR_DEFAULT);
    Vprintf("\n Root of Y(1) = 0.0 at %7.3f\n", x);
    Vprintf("\n Solution is");
    for (i = 0; i < 3; ++i)
      Vprintf("%10.5f", y[i]);
    Vprintf("\n");
  }
Vprintf("\n\nCase 2: no intermediate output, root-finding\n");
for (j = 4; j <= 5; ++j)
  {
    tol = pow(10.0, (double)(-j));
    Vprintf("\n Calculation with tol = %8.1e\n", tol);
    x = 0.0;
    y[0] = 0.5;
    y[1] = 0.5;
    y[2] = pi / 5.0;
    d02cjc(neq, fcn, &x, y, s.xend, tol, Nag_Mixed, NULLFN, g, &comm,
NAGERR_DEFAULT);
    Vprintf("\n Root of Y(1) = 0.0 at \%7.3f\n", x);
    Vprintf("\n Solution is");
    for (i = 0; i < 3; ++i)
    Vprintf("%10.5f", y[i]);</pre>
    Vprintf("\n");
  }
Vprintf("\n\nCase 3: intermediate output, no root-finding\n");
for (j = 4; j <= 5; ++j)
  ſ
    tol = pow(10.0, (double)(-j));
    Vprintf("\n Calculation with tol = %8.1e\n", tol);
    x = 0.0;
    y[0] = 0.5;
    y[1] = 0.5;
    y[2] = pi / 5.0;
    s.k = 4;
    s.h = (s.xend - x) / (double)(s.k + 1);
                   Х
    Vprintf("\n
                                            Y(2)
                               Y(1)
                                                          Y(3)\n");
```

```
d02cjc(neq, fcn, &x, y, s.xend, tol, Nag_Mixed, out, NULLDFN, &comm,
             NAGERR_DEFAULT);
    }
  Vprintf("\n\nCase 4: no intermediate output, no root-finding");
  Vprintf(" ( integrate to xend)\n");
  for (j = 4; j <= 5; ++j)
    {
      tol = pow(10.0, (double)(-j));
      Vprintf("\n Calculation with tol = %8.1e\n", tol);
      x = 0.0;
      y[0] = 0.5;
      y[1] = 0.5;
      y[2] = pi / 5.0;
Vprintf("\n
                      Х
                                  Y(1)
                                              Y(2)
                                                             Y(3)\n");
      Vprintf("%8.2f", x);
for (i = 0; i < 3; ++i)
       Vprintf("%13.5f", y[i]);
      Vprintf("\n");
      d02cjc(neq, fcn, &x, y, s.xend, tol, Nag_Mixed, NULLFN,
             NULLDFN, &comm, NAGERR_DEFAULT);
      Vprintf("%8.2f", x);
for (i = 0; i < 3; ++i)</pre>
        Vprintf("%13.5f", y[i]);
      Vprintf("\n");
    }
  exit(EXIT_SUCCESS);
}
                                  /* main */
#ifdef NAG_PROTO
static void out(Integer neq, double *xsol, double y[], Nag_User *comm)
#else
     static void out(neq, xsol, y, comm)
     Integer neq;
     double *xsol;
     double y[];
     Nag_User *comm;
#endif
{
  Integer i;
  struct user *s = (struct user *)comm->p;
  Vprintf("%8.2f", *xsol);
  for (i = 0; i < 3; ++i)
    {
      Vprintf("%13.5f", y[i]);
    }
  Vprintf("\n");
  *xsol = s->xend - (double)s->k * s->h;
  s->k--;
}
                                  /* out */
#ifdef NAG_PROTO
static void fcn(Integer neq, double x, double y[], double f[], Nag_User *comm)
#else
     static void fcn(neq, x, y, f, comm)
     Integer neq;
     double x;
     double y[], f[];
     Nag_User *comm;
#endif
{
  double pwr;
```

```
f[0] = tan(y[2]);
  f[1] = -0.032*tan(y[2])/y[1] - 0.02*y[1]/cos(y[2]);
  pwr = y[1];
f[2] = -0.032/(pwr*pwr);
}
                                   /* fcn */
#ifdef NAG_PROTO
static double g(Integer neq, double x, double y[], Nag_User *comm)
#else
     static double g(neq, x, y, comm)
     Integer neq;
     double x;
     double y[];
Nag_User *comm;
#endif
{
 return y[0];
}
                                   /* g */
```

#### 8.2. Program Data

None.

#### 8.3. Program Results

d02cjc Example Program Results

Case 1: intermediate output, root-finding

Calculation with tol = 1.0e-04Х Y(1) Y(2) Y(3) 0.00 0.50000 0.50000 0.62832 2.00 1.54931 0.40548 0.30662 4.00 1.74229 0.37433 -0.12890 6.00 1.00554 0.41731 -0.55068 Root of Y(1) = 0.0 at 7.288 Solution is 0.00000 0.47486 -0.76011 Calculation with tol = 1.0e-05Х Y(1) Y(2) Y(3) 0.00 0.50000 0.50000 0.62832 2.00 1.54933 0.40548 0.30662 4.00 1.74232 0.37433 -0.12891 6.00 1.00552 0.41731 -0.55069 Root of Y(1) = 0.0 at 7.288 Solution is 0.00000 0.47486 -0.76010 Case 2: no intermediate output, root-finding Calculation with tol = 1.0e-04Root of Y(1) = 0.0 at 7.288

Solution is 0.00000 0.47486 -0.76011 Calculation with tol = 1.0e-05 Root of Y(1) = 0.0 at 7.288 Solution is 0.00000 0.47486 -0.76010 Case 3: intermediate output, no root-finding

Calculati	on with tol =	1.0e-04		
X 0.00 2.00 4.00 6.00 8.00 10.00	Y(1) 0.50000 1.54931 1.74229 1.00554 -0.74589 -3.62813	Y(2) 0.50000 0.40548 0.37433 0.41731 0.51299 0.63325	Y(3) 0.62832 0.30662 -0.12890 -0.55068 -0.85371 -1.05152	
Calculati	on with tol =	1.0e-05		
X 0.00 2.00 4.00 6.00 8.00 10.00	Y(1) 0.50000 1.54933 1.74232 1.00552 -0.74601 -3.62829	Y(2) 0.50000 0.40548 0.37433 0.41731 0.51299 0.63326	Y(3) 0.62832 0.30662 -0.12891 -0.55069 -0.85372 -1.05153	
		-	oot-finding (	integrate to xend)
Calculati	on with tol =	1.0e-04		
X 0.00 10.00	Y(1) 0.50000 -3.62813	Y(2) 0.50000 0.63325	Y(3) 0.62832 -1.05152	
Calculation with tol = $1.0e-05$				
X 0.00 10.00	Y(1) 0.50000 -3.62829	Y(2) 0.50000 0.63326	Y(3) 0.62832 -1.05153	