

NAG Toolbox for MATLAB

d02cj

1 Purpose

d02cj 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 you, if desired.

2 Syntax

```
[x, y, ifail] = d02cj(x, xend, y, fcn, tol, relabs, output, g, 'n', n)
```

3 Description

d02cj advances the solution of a system of ordinary differential equations

$$y'_i = f_i(x, y_1, y_2, \dots, y_n), \quad i = 1, 2, \dots, n,$$

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

The solution is returned via the (sub)program **output** at points specified by you, 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 by C^1 interpolation to the solution. 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 **relabs**.

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

4 References

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

5 Parameters

5.1 Compulsory Input Parameters

1: **x – double scalar**

The initial value of the independent variable x .

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

2: **xend – double scalar**

The final value of the independent variable. If $\mathbf{xend} < \mathbf{x}$, integration will proceed in the negative direction.

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

3: **y(n) – double array**

The initial values of the solution $y_1, y_2, \dots, y_n$ at $x = \mathbf{x}$.

4: **fcn – string containing name of m-file**

fcn must evaluate the functions f_i (i.e., the derivatives y_i') for given values of its arguments $x, y_1, \dots, y_n$.

Its specification is:

$[f] = \text{fcn}(x, y)$	
Input Parameters	
1:	x – double scalar The value of the independent variable x .
2:	y(n) – double array The value of the variable y_i , for $i = 1, 2, \dots, n$.
Output Parameters	
1:	f(n) – double array The value of f_i , for $i = 1, 2, \dots, n$.

5: **tol – double scalar**

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.

d02cj 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. You are strongly recommended to call d02cj with more than one value for **tol** and to compare the results obtained to estimate their accuracy. In the absence of any prior knowledge, you might compare the results obtained by calling d02cj 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.

6: **relabs – string**

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:

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

where τ_r and τ_a are defined by

relabs	τ_r	τ_a
'M'	tol	tol
'A'	0.0	tol
'R'	tol	ϵ
'D'	tol	tol

where ϵ 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 you wish to measure the error in the computed solution in terms of the number of correct decimal places, then **relabs** should be set to 'A' on entry, whereas if the error requirement is in terms of the number of correct significant digits, then **relabs** should be set to 'R'. If you prefer a mixed error test, then **relabs** should be set to 'M', otherwise if you have no preference, **relabs** should be set to the default 'D'. Note that in this case 'D' is taken to be 'M'.

Constraint: **relabs** = 'M', 'A', 'R' or 'D'.

7: **output** – string containing name of m-file

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 d02cj with **xsol** = **x** (the initial value of x). You must reset **xsol** to the next point (between the current **xsol** and **xend**) where **output** is to be called, and so on at each call to **output**. If, after a call to **output**, the reset point **xsol** is beyond **xend**, d02cj will integrate to **xend** with no further calls to **output**; if a call to **output** is required at the point **xsol** = **xend**, then **xsol** must be given precisely the value **xend**.

Its specification is:

```
[xsol] = output(xsol, y)
```

Input Parameters1: **xsol** – double scalar

The output value of the independent variable x .

You must set **xsol** to the next value of x at which **output** is to be called.

2: **y(n)** – double array

The computed solution at the point **xsol**.

Output Parameters1: **xsol** – double scalar

The output value of the independent variable x .

You must set **xsol** to the next value of x at which **output** is to be called.

If you do not wish to access intermediate output, the actual argument **output** must be the string 'd02cjx'. **d02cjx** is included in the NAG Fortran Library.

8: **g** – string containing name of m-file

g must evaluate the function $g(x,y)$ 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.

Its specification is:

```
[result] = g(x, y)
```

Input Parameters1: **x** – double scalar

The value of the independent variable x .

2: **y(n)** – double array

The value of the variable y_i , for $i = 1, 2, \dots, n$.

Output Parameters1: **result** – double scalar

The result of the function.

If you do not require the root-finding option, the actual parameter **g** must be the string 'd02cjw'. **d02cjw** is included in the NAG Fortran Library.

5.2 Optional Input Parameters

1: **n** – **int32 scalar**

Default: The dimension of the array **y**.
the number of differential equations.

Constraint: $n \geq 1$.

5.3 Input Parameters Omitted from the MATLAB Interface

w

5.4 Output Parameters

1: **x** – **double scalar**

If *g* is supplied by you, it 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 you it contains **xend**, unless an error has occurred, when it contains the value of *x* at the error.

2: **y(n)** – **double array**

The computed values of the solution at the final point $x = \mathbf{x}$.

3: **ifail** – **int32 scalar**

0 unless the function detects an error (see Section 6).

6 Error Indicators and Warnings

Errors or warnings detected by the function:

ifail = 1

On entry, **tol** ≤ 0.0 ,
or **n** ≤ 0 ,
or **relabs** $\neq$ 'M', 'A', 'R' or 'D',
or **x** = **xend**.

ifail = 2

With the given value of **tol**, no further progress can be made across the integration range from the current point $x = \mathbf{x}$. (See Section 8 for a discussion of this error exit.) The components **y**(1), **y**(2), ..., **y**(**n**) contain the computed values of the solution at the current point $x = \mathbf{x}$. If you have supplied *g*, then no point at which $g(x,y)$ changes sign has been located up to the point $x = \mathbf{x}$.

ifail = 3

tol is too small for d02cj to take an initial step. **x** and **y**(1), **y**(2), ..., **y**(**n**) retain their initial values.

ifail = 4

xsol has not been reset or **xsol** lies behind **x** in the direction of integration, after the initial call to (sub)program **output**, if the **output** option was selected.

ifail = 5

A value of **xsol** returned by the (sub)program **output** has not been reset or lies behind the last value of **xsol** in the direction of integration, if the **output** option was selected.

ifail = 6

At no point in the range **x** to **xend** did the function $g(x,y)$ change sign, if g was supplied. It is assumed that $g(x,y) = 0$ has no solution.

ifail = 7

A serious error has occurred in an internal call. Check all (sub)program calls and array sizes. Seek expert help.

7 Accuracy

The accuracy of the computation of the solution vector **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. You are advised to choose **relabs** = 'M' unless you 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 $g(x,y)$. You should try to code user-supplied real function **g** without introducing any unnecessary cancellation errors.

8 Further Comments

If more than one root is required then d02qf should be used.

If d02cj fails with **ifail** = 3, 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 d02cj fails with **ifail** = 2, 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 **x** to **xend**. This can happen for well-behaved systems and very small values of **tol**. You 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 **ifail** = 2, 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 d02cj) to preserve stability. This will exhibit itself by making the computing time excessively long, or occasionally by an exit with **ifail** = 2. Adams methods are not efficient in such cases, and you should try d02ej.

9 Example

d02cj_fcn.m

```
function f = fcn(x,y)
    f = zeros(3,1);
    f(1) = tan(y(3));
    f(2) = -0.032*tan(y(3))/y(2) - 0.02*y(2)/cos(y(3));
    f(3) = -0.032/y(2)^2;
```

d02cj_g.m

```
function result = g(x,y)
    result = y(1);
```

d02cj_output.m

```
function xsolOut = output(xsol, y)
```

```
fprintf('%3.5f %3.5f %3.5f %3.5f\n', xsol, y(1), y(2), y(3));  
xsolOut = xsol + 2;
```

```
x = 0;  
xend = 10;  
y = [0.5;  
      0.5;  
      0.6283185307179586];  
tol = 0.0001;  
relabs = 'Default';  
[xOut, yOut, ifail] = ...  
    d02cj(x, xend, y, 'd02cj_fcn', tol, relabs, 'd02cj_output',  
    'd02cj_g')
```

```
0.00000 +0.50000 +0.50000 +0.62832  
2.00000 +1.54931 +0.40548 +0.30662  
4.00000 +1.74229 +0.37433 -0.12890  
6.00000 +1.00554 +0.41731 -0.55068  
xOut =  
      7.2883  
yOut =  
      -0.0000  
       0.4749  
      -0.7601  
ifail =  
       0
```