

NAG Toolbox

nag_ode_bvp_coll_nlin_solve (d02tl)

1 Purpose

nag_ode_bvp_coll_nlin_solve (d02tl) solves a general two-point boundary value problem for a nonlinear mixed order system of ordinary differential equations.

2 Syntax

```
[rcomm, icomm, user, ifail] = nag_ode_bvp_coll_nlin_solve(ffun, fjac, gafun,
gbfun, gajac, gbjac, guess, rcomm, icomm, 'user', user)

[rcomm, icomm, user, ifail] = d02tl(ffun, fjac, gafun, gbfun, gajac, gbjac,
guess, rcomm, icomm, 'user', user)
```

3 Description

nag_ode_bvp_coll_nlin_solve (d02tl) and its associated functions (nag_ode_bvp_coll_nlin_setup (d02tv), nag_ode_bvp_coll_nlin_contin (d02tx), nag_ode_bvp_coll_nlin_interp (d02ty) and nag_ode_bvp_coll_nlin_diag (d02tz)) solve the two-point boundary value problem for a nonlinear mixed order system of ordinary differential equations

$$\begin{aligned} y_1^{(m_1)}(x) &= f_1\left(x, y_1, y_1^{(1)}, \dots, y_1^{(m_1-1)}, y_2, \dots, y_n^{(m_n-1)}\right) \\ y_2^{(m_2)}(x) &= f_2\left(x, y_1, y_1^{(1)}, \dots, y_1^{(m_1-1)}, y_2, \dots, y_n^{(m_n-1)}\right) \\ &\vdots \\ y_n^{(m_n)}(x) &= f_n\left(x, y_1, y_1^{(1)}, \dots, y_1^{(m_1-1)}, y_2, \dots, y_n^{(m_n-1)}\right) \end{aligned}$$

over an interval $[a, b]$ subject to p (> 0) nonlinear boundary conditions at a and q (> 0) nonlinear boundary conditions at b , where $p + q = \sum_{i=1}^n m_i$. Note that $y_i^{(m)}(x)$ is the m th derivative of the i th solution component. Hence $y_i^{(0)}(x) = y_i(x)$. The left boundary conditions at a are defined as

$$g_i(z(y(a))) = 0, \quad i = 1, 2, \dots, p,$$

and the right boundary conditions at b as

$$\bar{g}_j(z(y(b))) = 0, \quad j = 1, 2, \dots, q,$$

where $y = (y_1, y_2, \dots, y_n)$ and

$$z(y(x)) = \left(y_1(x), y_1^{(1)}(x), \dots, y_1^{(m_1-1)}(x), y_2(x), \dots, y_n^{(m_n-1)}(x)\right).$$

First, nag_ode_bvp_coll_nlin_setup (d02tv) must be called to specify the initial mesh, error requirements and other details. Note that the error requirements apply only to the solution components y_1, y_2, \dots, y_n and that no error control is applied to derivatives of solution components. (If error control is required on derivatives then the system must be reduced in order by introducing the derivatives whose error is to be controlled as new variables. See Section 9 in nag_ode_bvp_coll_nlin_setup (d02tv).) Then, nag_ode_bvp_coll_nlin_solve (d02tl) can be used to solve the boundary value problem. After successful computation, nag_ode_bvp_coll_nlin_diag (d02tz) can be used to ascertain details about the final mesh and other details of the solution procedure, and nag_ode_bvp_coll_nlin_interp (d02ty) can be used to compute the approximate solution anywhere on the interval $[a, b]$.

A description of the numerical technique used in nag_ode_bvp_coll_nlin_solve (d02tl) is given in Section 3 in nag_ode_bvp_coll_nlin_setup (d02tv).

`nag_ode_bvp_coll_nlin_solve` (d02tl) can also be used in the solution of a series of problems, for example in performing continuation, when the mesh used to compute the solution of one problem is to be used as the initial mesh for the solution of the next related problem. `nag_ode_bvp_coll_nlin_contin` (d02tx) should be used in between calls to `nag_ode_bvp_coll_nlin_solve` (d02tl) in this context.

See Section 9 in `nag_ode_bvp_coll_nlin_setup` (d02tv) for details of how to solve boundary value problems of a more general nature.

The functions are based on modified versions of the codes COLSYS and COLNEW (see Ascher *et al.* (1979) and Ascher and Bader (1987)). A comprehensive treatment of the numerical solution of boundary value problems can be found in Ascher *et al.* (1988) and Keller (1992).

4 References

Ascher U M and Bader G (1987) A new basis implementation for a mixed order boundary value ODE solver *SIAM J. Sci. Stat. Comput.* **8** 483–500

Ascher U M, Christiansen J and Russell R D (1979) A collocation solver for mixed order systems of boundary value problems *Math. Comput.* **33** 659–679

Ascher U M, Mattheij R M M and Russell R D (1988) *Numerical Solution of Boundary Value Problems for Ordinary Differential Equations* Prentice–Hall

Keller H B (1992) *Numerical Methods for Two-point Boundary-value Problems* Dover, New York

5 Parameters

5.1 Compulsory Input Parameters

1: **ffun** – SUBROUTINE, supplied by the user.

ffun must evaluate the functions f_i for given values $x, z(y(x))$.

```
[f, user] = ffun(x, y, neq, m, user)
```

Input Parameters

1: **x** – REAL (KIND=nag_wp)

x , the independent variable.

2: **y(neq, 0 : *)** – REAL (KIND=nag_wp) array

$y(i, j)$ contains $y_i^{(j)}(x)$, for $i = 1, 2, \dots, \mathbf{neq}$ and $j = 0, 1, \dots, \mathbf{m}(i) - 1$.

Note: $y_i^{(0)}(x) = y_i(x)$.

3: **neq** – INTEGER

The number of differential equations.

4: **m(neq)** – INTEGER array

$\mathbf{m}(i)$ contains m_i , the order of the i th differential equation, for $i = 1, 2, \dots, \mathbf{neq}$.

5: **user** – INTEGER array

ffun is called from `nag_ode_bvp_coll_nlin_solve` (d02tl) with the object supplied to `nag_ode_bvp_coll_nlin_solve` (d02tl).

Output Parameters

- 1: **f(neq)** – REAL (KIND=nag_wp) array
f(i) must contain f_i , for $i = 1, 2, \dots, \mathbf{neq}$.
- 2: **user** – INTEGER array

- 2: **fjac** – SUBROUTINE, supplied by the user.

fjac must evaluate the partial derivatives of f_i with respect to the elements of $z(y(x)) = (y_1(x), y_1^1(x), \dots, y_1^{(m_1-1)}(x), y_2(x), \dots, y_n^{(m_n-1)}(x))$.

```
[dfdy, user] = fjac(x, y, neq, m, dfdy, user)
```

Input Parameters

- 1: **x** – REAL (KIND=nag_wp)
 x , the independent variable.
- 2: **y(neq, 0 : *)** – REAL (KIND=nag_wp) array
y(i, j) contains $y_i^{(j)}(x)$, for $i = 1, 2, \dots, \mathbf{neq}$ and $j = 0, 1, \dots, \mathbf{m}(i) - 1$.
Note: $y_i^{(0)}(x) = y_i(x)$.
- 3: **neq** – INTEGER
The number of differential equations.
- 4: **m(neq)** – INTEGER array
m(i) contains m_i , the order of the i th differential equation, for $i = 1, 2, \dots, \mathbf{neq}$.
- 5: **dfdy(neq, neq, 0 : *)** – REAL (KIND=nag_wp) array
Set to zero.
- 6: **user** – INTEGER array
fjac is called from nag_ode_bvp_coll_nlin_solve (d02tl) with the object supplied to nag_ode_bvp_coll_nlin_solve (d02tl).

Output Parameters

- 1: **dfdy(neq, neq, 0 : *)** – REAL (KIND=nag_wp) array
dfdy(i, j, k) must contain the partial derivative of f_i with respect to $y_j^{(k)}$, for $i = 1, 2, \dots, \mathbf{neq}$, $j = 1, 2, \dots, \mathbf{neq}$ and $k = 0, 1, \dots, \mathbf{m}(j) - 1$. Only nonzero partial derivatives need be set.
- 2: **user** – INTEGER array

- 3: **gafun** – SUBROUTINE, supplied by the user.

gafun must evaluate the boundary conditions at the left-hand end of the range, that is functions $g_i(z(y(a)))$ for given values of $z(y(a))$.

```
[ga, user] = gafun(ya, neq, m, nlbc, user)
```

Input Parameters

- 1: **ya**(**neq**, 0 : *) – REAL (KIND=nag_wp) array
ya(*i*, *j*) contains $y_i^{(j)}(a)$, for $i = 1, 2, \dots, \mathbf{neq}$ and $j = 0, 1, \dots, \mathbf{m}(i) - 1$.
Note: $y_i^{(0)}(a) = y_i(a)$.
- 2: **neq** – INTEGER
The number of differential equations.
- 3: **m**(**neq**) – INTEGER array
m(*i*) contains m_i , the order of the *i*th differential equation, for $i = 1, 2, \dots, \mathbf{neq}$.
- 4: **nlbc** – INTEGER
The number of boundary conditions at *a*.
- 5: **user** – INTEGER array
gafun is called from nag_ode_bvp_coll_nlin_solve (d02tl) with the object supplied to nag_ode_bvp_coll_nlin_solve (d02tl).

Output Parameters

- 1: **ga**(**nlbc**) – REAL (KIND=nag_wp) array
ga(*i*) must contain $g_i(z(y(a)))$, for $i = 1, 2, \dots, \mathbf{nlbc}$.
- 2: **user** – INTEGER array

- 4: **gbfun** – SUBROUTINE, supplied by the user.

gbfun must evaluate the boundary conditions at the right-hand end of the range, that is functions $\bar{g}_i(z(y(b)))$ for given values of $z(y(b))$.

```
[gb, user] = gbfun(yb, neq, m, nrbc, user)
```

Input Parameters

- 1: **yb**(**neq**, 0 : *) – REAL (KIND=nag_wp) array
yb(*i*, *j*) contains $y_i^{(j)}(b)$, for $i = 1, 2, \dots, \mathbf{neq}$ and $j = 0, 1, \dots, \mathbf{m}(i) - 1$.
Note: $y_i^{(0)}(b) = y_i(b)$.
- 2: **neq** – INTEGER
The number of differential equations.
- 3: **m**(**neq**) – INTEGER array
m(*i*) contains m_i , the order of the *i*th differential equation, for $i = 1, 2, \dots, \mathbf{neq}$.
- 4: **nrbc** – INTEGER
The number of boundary conditions at *b*.

5: **user** – INTEGER array
gbfun is called from `nag_ode_bvp_coll_nlin_solve (d02tl)` with the object supplied to `nag_ode_bvp_coll_nlin_solve (d02tl)`.

Output Parameters

1: **gb(nrbc)** – REAL (KIND=`nag_wp`) array
gb(*i*) must contain $\bar{g}_i(z(y(b)))$, for $i = 1, 2, \dots, \mathbf{nrbc}$.

2: **user** – INTEGER array

5: **gajac** – SUBROUTINE, supplied by the user.

gajac must evaluate the partial derivatives of $g_i(z(y(a)))$ with respect to the elements of $z(y(a)) = (y_1(a), y_1^1(a), \dots, y_1^{(m_1-1)}(a), y_2(a), \dots, y_n^{(m_n-1)}(a))$.

```
[dgady, user] = gajac(ya, neq, m, nlbc, dgady, user)
```

Input Parameters

1: **ya(neq, 0 : *)** – REAL (KIND=`nag_wp`) array
ya(*i, j*) contains $y_i^{(j)}(a)$, for $i = 1, 2, \dots, \mathbf{neq}$ and $j = 0, 1, \dots, \mathbf{m}(i) - 1$.
Note: $y_i^{(0)}(a) = y_i(a)$.

2: **neq** – INTEGER
The number of differential equations.

3: **m(neq)** – INTEGER array
m(*i*) contains m_i , the order of the *i*th differential equation, for $i = 1, 2, \dots, \mathbf{neq}$.

4: **nlbc** – INTEGER
The number of boundary conditions at *a*.

5: **dgady(nlbc, neq, 0 : *)** – REAL (KIND=`nag_wp`) array
Set to zero.

6: **user** – INTEGER array
gajac is called from `nag_ode_bvp_coll_nlin_solve (d02tl)` with the object supplied to `nag_ode_bvp_coll_nlin_solve (d02tl)`.

Output Parameters

1: **dgady(nlbc, neq, 0 : *)** – REAL (KIND=`nag_wp`) array
dgady(*i, j, k*) must contain the partial derivative of $g_i(z(y(a)))$ with respect to $y_j^{(k)}(a)$, for $i = 1, 2, \dots, \mathbf{nlbc}$, $j = 1, 2, \dots, \mathbf{neq}$ and $k = 0, 1, \dots, \mathbf{m}(j) - 1$. Only nonzero partial derivatives need be set.

2: **user** – INTEGER array

6: **gbjac** – SUBROUTINE, supplied by the user.

gbjac must evaluate the partial derivatives of $\bar{g}_i(z(y(b)))$ with respect to the elements of $z(y(b)) = (y_1(b), y_1^1(b), \dots, y_1^{(m_1-1)}(b), y_2(b), \dots, y_n^{(m_n-1)}(b))$.

```
[dgbdy, user] = gbjac(yb, neq, m, nrbc, dgbdy, user)
```

Input Parameters

1: **yb**(**neq**, 0 : *) – REAL (KIND=nag_wp) array

yb(*i*, *j*) contains $y_i^{(j)}(b)$, for $i = 1, 2, \dots, \mathbf{neq}$ and $j = 0, 1, \dots, \mathbf{m}(i) - 1$.

Note: $y_i^{(0)}(b) = y_i(b)$.

2: **neq** – INTEGER

The number of differential equations.

3: **m**(**neq**) – INTEGER array

m(*i*) contains m_i , the order of the *i*th differential equation, for $i = 1, 2, \dots, \mathbf{neq}$.

4: **nrbc** – INTEGER

The number of boundary conditions at *b*.

5: **dgbdy**(**nrbc**, **neq**, 0 : *) – REAL (KIND=nag_wp) array

Set to zero.

6: **user** – INTEGER array

gbjac is called from `nag_ode_bvp_coll_nlin_solve (d02tl)` with the object supplied to `nag_ode_bvp_coll_nlin_solve (d02tl)`.

Output Parameters

1: **dgbdy**(**nrbc**, **neq**, 0 : *) – REAL (KIND=nag_wp) array

dgbdy(*i*, *j*, *k* + 1) must contain the partial derivative of $\bar{g}_i(z(y(b)))$ with respect to $y_j^{(k)}(b)$, for $i = 1, 2, \dots, \mathbf{nrbc}$, $j = 1, 2, \dots, \mathbf{neq}$ and $k = 0, 1, \dots, \mathbf{m}(j) - 1$. Only nonzero partial derivatives need be set.

2: **user** – INTEGER array

7: **guess** – SUBROUTINE, supplied by the user.

guess must return initial approximations for the solution components $y_i^{(j)}$ and the derivatives $y_i^{(m_i)}$, for $i = 1, 2, \dots, \mathbf{neq}$ and $j = 0, 1, \dots, \mathbf{m}(i) - 1$. Try to compute each derivative $y_i^{(m_i)}$ such that it corresponds to your approximations to $y_i^{(j)}$, for $j = 0, 1, \dots, \mathbf{m}(i) - 1$. You should **not** call **ffun** to compute $y_i^{(m_i)}$.

If `nag_ode_bvp_coll_nlin_solve (d02tl)` is being used in conjunction with `nag_ode_bvp_coll_nlin_contin (d02tx)` as part of a continuation process, then **guess** is not called by `nag_ode_bvp_coll_nlin_solve (d02tl)` after the call to `nag_ode_bvp_coll_nlin_contin (d02tx)`.

```
[y, dym, user] = guess(x, neq, m, user)
```

Input Parameters

- 1: **x** – REAL (KIND=nag_wp)
 x , the independent variable; $x \in [a, b]$.
- 2: **neq** – INTEGER
The number of differential equations.
- 3: **m(neq)** – INTEGER array
m(i) contains m_i , the order of the i th differential equation, for $i = 1, 2, \dots, \mathbf{neq}$.
- 4: **user** – INTEGER array
guess is called from `nag_ode_bvp_coll_nlin_solve (d02tl)` with the object supplied to `nag_ode_bvp_coll_nlin_solve (d02tl)`.

Output Parameters

- 1: **y(neq, 0 : *)** – REAL (KIND=nag_wp) array
y(i, j) must contain $y_i^{(j)}(x)$, for $i = 1, 2, \dots, \mathbf{neq}$ and $j = 0, 1, \dots, \mathbf{m}(i) - 1$.
Note: $y_i^{(0)}(x) = y_i(x)$.
- 2: **dym(neq)** – REAL (KIND=nag_wp) array
dym(i) must contain $y_i^{(m_i)}(x)$, for $i = 1, 2, \dots, \mathbf{neq}$.
- 3: **user** – INTEGER array

- 8: **rcomm(*)** – REAL (KIND=nag_wp) array

Note: the dimension of this array is dictated by the requirements of associated functions that must have been previously called. This array **must** be the same array passed as argument **rcomm** in the previous call to `nag_ode_bvp_coll_nlin_setup (d02tv)`.

This must be the same array as supplied to `nag_ode_bvp_coll_nlin_setup (d02tv)` and **must** remain unchanged between calls.

- 9: **icomm(*)** – INTEGER array

Note: the dimension of this array is dictated by the requirements of associated functions that must have been previously called. This array **must** be the same array passed as argument **icomm** in the previous call to `nag_ode_bvp_coll_nlin_setup (d02tv)`.

This must be the same array as supplied to `nag_ode_bvp_coll_nlin_setup (d02tv)` and **must** remain unchanged between calls.

5.2 Optional Input Parameters

- 1: **user** – INTEGER array

user is not used by `nag_ode_bvp_coll_nlin_solve (d02tl)`, but is passed to **ffun**, **fjac**, **gafun**, **gbfun**, **gajac**, **gbjac** and **guess**. Note that for large objects it may be more efficient to use a global variable which is accessible from the m-files than to use **user**.

5.3 Output Parameters

1: **rcomm**(*) – REAL (KIND=nag_wp) array

Note: the dimension of this array is dictated by the requirements of associated functions that must have been previously called. This array **must** be the same array passed as argument **rcomm** in the previous call to `nag_ode_bvp_coll_nlin_setup` (d02tv).

Contains information about the solution for use on subsequent calls to associated functions.

2: **icomm**(*) – INTEGER array

Note: the dimension of this array is dictated by the requirements of associated functions that must have been previously called. This array **must** be the same array passed as argument **icomm** in the previous call to `nag_ode_bvp_coll_nlin_setup` (d02tv).

Contains information about the solution for use on subsequent calls to associated functions.

3: **user** – INTEGER array

4: **ifail** – INTEGER

ifail = 0 unless the function detects an error (see Section 5).

6 Error Indicators and Warnings

Note: `nag_ode_bvp_coll_nlin_solve` (d02tl) may return useful information for one or more of the following detected errors or warnings.

Errors or warnings detected by the function:

ifail = 1

Either the setup function has not been called or the communication arrays have become corrupted. No solution will be computed.

ifail = 2

Numerical singularity has been detected in the Jacobian used in the Newton iteration. No results have been generated. Check the coding of the functions for calculating the Jacobians of system and boundary conditions.

ifail = 3

All Newton iterations that have been attempted have failed to converge. No results have been generated. Check the coding of the functions for calculating the Jacobians of system and boundary conditions. Try to provide a better initial solution approximation.

ifail = 4

A Newton iteration has failed to converge. The computation has not succeeded but results have been returned for an intermediate mesh on which convergence was achieved. These results should be treated with extreme caution.

ifail = 5

The expected number of sub-intervals required to continue the computation exceeds the maximum specified.

ifail = -99

An unexpected error has been triggered by this routine. Please contact NAG.

ifail = –399

Your licence key may have expired or may not have been installed correctly.

ifail = –999

Dynamic memory allocation failed.

7 Accuracy

The accuracy of the solution is determined by the argument **tols** in the prior call to `nag_ode_bvp_coll_nlin_setup` (d02tv) (see Section 3 and Section 9 in `nag_ode_bvp_coll_nlin_setup` (d02tv) for details and advice). Note that error control is applied only to solution components (variables) and not to any derivatives of the solution. An estimate of the maximum error in the computed solution is available by calling `nag_ode_bvp_coll_nlin_diag` (d02tz).

8 Further Comments

If `nag_ode_bvp_coll_nlin_solve` (d02tl) returns with **ifail** = 2, 3, 4 or 5 and the call to `nag_ode_bvp_coll_nlin_solve` (d02tl) was a part of some continuation procedure for which successful calls to `nag_ode_bvp_coll_nlin_solve` (d02tl) have already been made, then it is possible that the adjustment(s) to the continuation parameter(s) between calls to `nag_ode_bvp_coll_nlin_solve` (d02tl) is (are) too large for the problem under consideration. More conservative adjustment(s) to the continuation parameter(s) might be appropriate.

9 Example

The following example is used to illustrate the treatment of a high-order system, control of the error in a derivative of a component of the original system, and the use of continuation. See also `nag_ode_bvp_coll_nlin_setup` (d02tv), `nag_ode_bvp_coll_nlin_contin` (d02tx), `nag_ode_bvp_coll_nlin_interp` (d02ty) and `nag_ode_bvp_coll_nlin_diag` (d02tz), for the illustration of other facilities.

Consider the steady flow of an incompressible viscous fluid between two infinite coaxial rotating discs. See Ascher *et al.* (1979) and the references therein. The governing equations are

$$\begin{aligned}\frac{1}{\sqrt{R}}f''' + ff''' + gg' &= 0 \\ \frac{1}{\sqrt{R}}g'' + fg' - f'g &= 0\end{aligned}$$

subject to the boundary conditions

$$f(0) = f'(0) = 0, \quad g(0) = \Omega_0, \quad f(1) = f'(1) = 0, \quad g(1) = \Omega_1,$$

where R is the Reynolds number and Ω_0, Ω_1 are the angular velocities of the disks.

We consider the case of counter-rotation and a symmetric solution, that is $\Omega_0 = 1, \Omega_1 = -1$. This problem is more difficult to solve, the larger the value of R . For illustration, we use simple continuation to compute the solution for three different values of R ($= 10^6, 10^8, 10^{10}$). However, this problem can be addressed directly for the largest value of R considered here. Instead of the values suggested in Section 5 in `nag_ode_bvp_coll_nlin_contin` (d02tx) for **nmesh**, **ipmesh** and **mesh** in the call to `nag_ode_bvp_coll_nlin_contin` (d02tx) prior to a continuation call, we use every point of the final mesh for the solution of the first value of R , that is we must modify the contents of **ipmesh**. For illustrative purposes we wish to control the computed error in f' and so recast the equations as

$$\begin{aligned}y_1' &= y_2 \\ y_2'' &= -\sqrt{R}(y_1y_2'' + y_3y_3') \\ y_3'' &= \sqrt{R}(y_2y_3 - y_1y_3')\end{aligned}$$

subject to the boundary conditions

$$y_1(0) = y_2(0) = 0, \quad y_3(0) = \Omega, \quad y_1(1) = y_2(1) = 0, \quad y_3(1) = -\Omega, \quad \Omega = 1.$$

For the symmetric boundary conditions considered, there exists an odd solution about $x = 0.5$. Hence,

to satisfy the boundary conditions, we use the following initial approximations to the solution in **guess**:

$$\begin{aligned}y_1(x) &= -x^2(x - \frac{1}{2})(x - 1)^2 \\y_2(x) &= -x(x - 1)(5x^2 - 5x + 1) \\y_3(x) &= -8\Omega(x - \frac{1}{2})^3.\end{aligned}$$

9.1 Program Text

```
function d02tl_example

fprintf('d02tl example results\n\n');

global omega sqrofr

% Initialize variables and arrays.
neq = nag_int(3);
nlbc = neq;
nrbc = neq;
ncol = nag_int(7);
mmax = nag_int(3);
m = nag_int([1; 3; 2]);
tols = [1; 1; 1]/10^4;

% Set values for problem-specific physical parameters.
omega = 1;
r = 10^6;

% Set up the mesh.
nmesh = nag_int(11);
mxmesh = nag_int(51);
ipmesh = zeros(mxmesh, 1, nag_int_name);
mesh = zeros(mxmesh, 1);

mesh(1:nmesh) = [0:0.1:1];
ipmesh(1) = 1;
ipmesh(2:nmesh-1) = 2;
ipmesh(nmesh) = 1;

% d02tv is a setup routine to be called prior to d02tl.
[rcomm, icomm, ifail] = d02tv(...
    m, nlbc, nrbc, ncol, tols, nmesh, mesh, ipmesh);

% Set number of continuation steps.
ncont = 3;

% We run through the calculation three times with different parameter sets.
for jcont = 1:ncont
    sqrofr = sqrt(r);
    fprintf('\n Tolerance = %8.1e R = %10.3e\n\n', tols(1), r);

    % Call d02tk to solve BVP for this set of parameters.
    [rcomm, icomm, ifail] = ...
        d02tk(...
            @ffun, @fjac, @gafun, @gbfun, @gajac, @gbjac, @guess, rcomm, icomm);

    % Call d02tz to extract mesh from solution.
    [nmesh, mesh, ipmesh, erm, ierm, ijerm, ifail] = ...
        d02tz( ...
            mxmesh, rcomm, icomm);

    % Output mesh results.
    fprintf(' Mesh size = %d\n', nmesh);
    fprintf(' Maximum error = %10.2e in interval %d for component %d\n\n', ...
        erm, ierm, ijerm);
    fprintf('Mesh points:\n');
    for imesh = 1:nmesh
        fprintf(' %4d(%d) %6.4f', imesh, ipmesh(imesh), mesh(imesh));
        if mod(imesh, 4) == 0
            fprintf('\n');
        end
    end
end
```

```

    end
end

% Output solution, and store it for plotting.
xarray = zeros(nmesh, 1);
yarray = zeros(nmesh, 3);
fprintf('\n\n    x          f          f''          g\n');
for imesh = 1:nmesh
    % Call d02ty to perform interpolation on the solution.
    [y, rcomm, ifail] = d02ty(...
        mesh(imesh), neq, mmax, rcomm, icomm);
    fprintf(' %6.3f %8.4f %8.4f %8.4f\n', mesh(imesh), y(:,1));
    xarray(imesh) = mesh(imesh);
    yarray(imesh, 1:3) = y(1:3,1);
end

% Plot results for this parameter set.
if jcont==1
    fig1 = figure;
else
    if jcont==2
        fig2 = figure;
    else
        fig3 = figure;
    end
end
display_plot(xarray, yarray, r);

% Select mesh for next calculation.
if jcont < ncont
    r = 100*r;
    ipmesh(2:nmesh-1) = 2;

    % d02tx allows the current solution to be used for continuation
    [rcomm, icomm, ifail] = d02tx(...
        nmesh, mesh, ipmesh, rcomm, icomm);
end
end

function [f, user] = ffun(x, y, neq, m, user)
% Evaluate derivative functions (rhs of system of ODEs).

global omega sqrofr; % For communication with main routine.
f = zeros(neq, 1);
f(1,1) = y(2,1);
f(2,1) = -(y(1,1)*y(2,3) + y(3,1)*y(3,2))*sqrofr;
f(3,1) = (y(2,1)*y(3,1) - y(1,1)*y(3,2))*sqrofr;

function [dfdy, user] = fjac(x, y, neq, m, dfdy, user)
% Evaluate Jacobians (partial derivatives of f).

global omega sqrofr
dfdy = zeros(neq, neq, 3);
dfdy(1,2,1) = 1.0;
dfdy(2,1,1) = -y(2,3)*sqrofr;
dfdy(2,2,3) = -y(1,1)*sqrofr;
dfdy(2,3,1) = -y(3,2)*sqrofr;
dfdy(2,3,2) = -y(3,1)*sqrofr;
dfdy(3,1,1) = -y(3,2)*sqrofr;
dfdy(3,2,1) = y(3,1)*sqrofr;
dfdy(3,3,1) = y(2,1)*sqrofr;
dfdy(3,3,2) = -y(1,1)*sqrofr;

function [ga, user] = gafun(ya, neq, m, nlbc, user)
% Evaluate boundary conditions at left-hand end of range.

global omega sqrofr
ga = zeros(nlbc, 1);
ga(1) = ya(1);
ga(2) = ya(2);
ga(3) = ya(3) - omega;

```

```

function [dgady, user] = gajac(ya, neq, m, nlbc, dgady, user)
    % Evaluate Jacobians (partial derivatives of ga).

    dgady = zeros(nlbc, neq, 3);
    dgady(1,1,1) = 1;
    dgady(2,2,1) = 1;
    dgady(3,3,1) = 1;

function [gb, user] = gbfun(yb, neq, m, nrbc, user)
    % Evaluate boundary conditions at right-hand end of range.

    global omega sqrofr
    gb = zeros(nrbc, 1);
    gb(1) = yb(1);
    gb(2) = yb(2);
    gb(3) = yb(3) + omega;

function [dgbdy, user] = gbjac(yb, neq, m, nrbc, dgbdy, user)
    % Evaluate Jacobians (partial derivatives of gb).

    dgbdy = zeros(nrbc, neq, 3);
    dgbdy(1,1,1) = 1;
    dgbdy(2,2,1) = 1;
    dgbdy(3,3,1) = 1;

function [y, dym, user] = guess(x, neq, m, user)
    % Evaluate initial approximations to solution components and derivatives.

    global omega sqrofr
    y = zeros(neq, 3);
    dym = zeros(neq, 1);
    y(1,1) = -x^2*(x - 0.5)*(x - 1)^2;
    y(2,1) = -x*(x - 1)*(5*x^2 - 5*x + 1);
    y(3,1) = -8*omega*(x - 0.5)^3;
    y(2,2) = -(20*x^3 - 30*x^2 + 12*x - 1);
    y(2,3) = -(60*x^2 - 60*x + 12*x);
    y(3,2) = -24*omega*(x - 0.5)^2;

    dym(1) = y(2,1);
    dym(2) = -(120*x - 60);
    dym(3) = -56*omega*(x - 0.5);

function display_plot(x, y, r)
    % Plot two of the curves, then add the other one.
    [haxes, hline1, hline2] = plotyy(x, y(:,2), x, y(:,3));
    % We want the third curve to be plotted on the left-hand y-axis.
    hold(haxes(1), 'on');
    hline3 = plot(x, y(:,1));
    % Set the axis limits and the tick specifications to beautify the plot.
    set(haxes(1), 'YLim', [-0.1 0.4]);
    set(haxes(2), 'YLim', [-1 1]);
    set(haxes(1), 'XMinorTick', 'on', 'YMinorTick', 'on');
    set(haxes(2), 'YMinorTick', 'on');
    set(haxes(1), 'YTick', [-0.1:0.1:0.4]);
    set(haxes(2), 'YTick', [-1:0.5:1]);
    for iaxis = 1:2
        % These properties must be the same for both sets of axes.
        set(haxes(iaxis), 'XLim', [0 1]);
        set(haxes(iaxis), 'XTick', [0:0.2:1]);
    end
    set(gca, 'box', 'off');
    % Add title.
    ord = log10(r);
    title(['Flow between Discs, Re = 10^n, n = ', num2str(ord)]);
    % Label the axes.
    xlabel('x');
    ylabel(haxes(1), 'f and f''');
    ylabel(haxes(2), 'g');
    % Add a legend.
    legend('f''', 'f', 'g', 'Location', 'Best')

```

```
% Set some features of the three lines.
set(hline1, 'Linewidth', 0.25, 'Marker', '+', 'LineStyle', '-', ...
    'Color', 'Magenta');
set(hline2, 'Linewidth', 0.25, 'Marker', 'x', 'LineStyle', '--');
set(hline3, 'Linewidth', 0.25, 'Marker', '*', 'LineStyle', ':');
```

9.2 Program Results

d02tl example results

Tolerance = 1.0e-04 R = 1.000e+06

Mesh size = 21

Maximum error = 6.16e-10 in interval 20 for component 3

Mesh points:

1(1)	0.0000	2(3)	0.0500	3(2)	0.1000	4(3)	0.1500
5(2)	0.2000	6(3)	0.2500	7(2)	0.3000	8(3)	0.3500
9(2)	0.4000	10(3)	0.4500	11(2)	0.5000	12(3)	0.5500
13(2)	0.6000	14(3)	0.6500	15(2)	0.7000	16(3)	0.7500
17(2)	0.8000	18(3)	0.8500	19(2)	0.9000	20(3)	0.9500
21(1)	1.0000						

x	f	f'	g
0.000	0.0000	0.0000	1.0000
0.050	0.0070	0.1805	0.4416
0.100	0.0141	0.0977	0.1886
0.150	0.0171	0.0252	0.0952
0.200	0.0172	-0.0165	0.0595
0.250	0.0157	-0.0400	0.0427
0.300	0.0133	-0.0540	0.0322
0.350	0.0104	-0.0628	0.0236
0.400	0.0071	-0.0683	0.0156
0.450	0.0036	-0.0714	0.0078
0.500	0.0000	-0.0724	0.0000
0.550	-0.0036	-0.0714	-0.0078
0.600	-0.0071	-0.0683	-0.0156
0.650	-0.0104	-0.0628	-0.0236
0.700	-0.0133	-0.0540	-0.0322
0.750	-0.0157	-0.0400	-0.0427
0.800	-0.0172	-0.0165	-0.0595
0.850	-0.0171	0.0252	-0.0952
0.900	-0.0141	0.0977	-0.1886
0.950	-0.0070	0.1805	-0.4416
1.000	-0.0000	-0.0000	-1.0000

Tolerance = 1.0e-04 R = 1.000e+08

Mesh size = 21

Maximum error = 4.49e-09 in interval 6 for component 3

Mesh points:

1(1)	0.0000	2(3)	0.0176	3(2)	0.0351	4(3)	0.0520
5(2)	0.0689	6(3)	0.0859	7(2)	0.1030	8(3)	0.1351
9(2)	0.1672	10(3)	0.2306	11(2)	0.2939	12(3)	0.4713
13(2)	0.6486	14(3)	0.7455	15(2)	0.8423	16(3)	0.8824
17(2)	0.9225	18(3)	0.9449	19(2)	0.9673	20(3)	0.9836
21(1)	1.0000						

x	f	f'	g
0.000	0.0000	0.0000	1.0000
0.018	0.0025	0.1713	0.3923
0.035	0.0047	0.0824	0.1381
0.052	0.0056	0.0267	0.0521
0.069	0.0058	0.0025	0.0213
0.086	0.0057	-0.0073	0.0097
0.103	0.0056	-0.0113	0.0053
0.135	0.0052	-0.0135	0.0027
0.167	0.0047	-0.0140	0.0020
0.231	0.0038	-0.0142	0.0015

0.294	0.0029	-0.0142	0.0012
0.471	0.0004	-0.0143	0.0002
0.649	-0.0021	-0.0143	-0.0008
0.745	-0.0035	-0.0142	-0.0014
0.842	-0.0049	-0.0139	-0.0022
0.882	-0.0054	-0.0127	-0.0036
0.922	-0.0058	-0.0036	-0.0141
0.945	-0.0057	0.0205	-0.0439
0.967	-0.0045	0.0937	-0.1592
0.984	-0.0023	0.1753	-0.4208
1.000	-0.0000	0.0000	-1.0000

Tolerance = 1.0e-04 R = 1.000e+10

Mesh size = 21

Maximum error = 3.13e-06 in interval 7 for component 3

Mesh points:

1(1) 0.0000	2(3) 0.0063	3(2) 0.0125	4(3) 0.0185
5(2) 0.0245	6(3) 0.0308	7(2) 0.0370	8(3) 0.0500
9(2) 0.0629	10(3) 0.0942	11(2) 0.1256	12(3) 0.4190
13(2) 0.7125	14(3) 0.8246	15(2) 0.9368	16(3) 0.9544
17(2) 0.9719	18(3) 0.9803	19(2) 0.9886	20(3) 0.9943
21(1) 1.0000			

x	f	f'	g
0.000	0.0000	0.0000	1.0000
0.006	0.0009	0.1623	0.3422
0.013	0.0016	0.0665	0.1021
0.019	0.0018	0.0204	0.0318
0.025	0.0019	0.0041	0.0099
0.031	0.0019	-0.0014	0.0028
0.037	0.0019	-0.0031	0.0007
0.050	0.0019	-0.0038	-0.0002
0.063	0.0018	-0.0038	-0.0003
0.094	0.0017	-0.0039	-0.0003
0.126	0.0016	-0.0039	-0.0002
0.419	0.0004	-0.0041	-0.0001
0.712	-0.0008	-0.0042	0.0001
0.825	-0.0013	-0.0043	0.0002
0.937	-0.0018	-0.0043	0.0003
0.954	-0.0019	-0.0042	0.0001
0.972	-0.0019	-0.0003	-0.0049
0.980	-0.0019	0.0152	-0.0252
0.989	-0.0015	0.0809	-0.1279
0.994	-0.0008	0.1699	-0.3814
1.000	0.0000	0.0000	-1.0000





