

NAG Library Routine Document

D03FAF

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

D03FAF solves the Helmholtz equation in Cartesian coordinates in three dimensions using the standard seven-point finite difference approximation. This routine is designed to be particularly efficient on vector processors.

2 Specification

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SUBROUTINE D03FAF (XS, XF, L, LBDCND, BDXS, BDXF, YS, YF, M, MBDCND,      &
                  BDYS, BDYF, ZS, ZF, N, NBDCND, BDZS, BDZF, LAMBDA,    &
                  LDF, LDF2, F, PERTRB, W, LWRK, IFAIL)
INTEGER           L, LBDCND, M, MBDCND, N, NBDCND, LDF, LDF2, LWRK,      &
                  IFAIL
REAL (KIND=nag_wp) XS, XF, BDXS(LDF2,N+1), BDXF(LDF2,N+1), YS, YF,      &
                  BDYS(LDF,N+1), BDYF(LDF,N+1), ZS, ZF,                &
                  BDZS(LDF,M+1), BDZF(LDF,M+1), LAMBDA,                &
                  F(LDF,LDF2,N+1), PERTRB, W(LWRK)

```

3 Description

D03FAF solves the three-dimensional Helmholtz equation in Cartesian coordinates:

$$\frac{\partial^2 u}{\partial x^2} + \frac{\partial^2 u}{\partial y^2} + \frac{\partial^2 u}{\partial z^2} + \lambda u = f(x, y, z).$$

This subroutine forms the system of linear equations resulting from the standard seven-point finite difference equations, and then solves the system using a method based on the fast Fourier transform (FFT) described by Swarztrauber (1984). This subroutine is based on the routine HW3CRT from FISHPACK (see Swarztrauber and Sweet (1979)).

More precisely, the routine replaces all the second derivatives by second-order central difference approximations, resulting in a block tridiagonal system of linear equations. The equations are modified to allow for the prescribed boundary conditions. Either the solution or the derivative of the solution may be specified on any of the boundaries, or the solution may be specified to be periodic in any of the three dimensions. By taking the discrete Fourier transform in the x - and y -directions, the equations are reduced to sets of tridiagonal systems of equations. The Fourier transforms required are computed using the multiple FFT routines found in Chapter C06.

4 References

Swarztrauber P N (1984) Fast Poisson solvers *Studies in Numerical Analysis* (ed G H Golub) Mathematical Association of America

Swarztrauber P N and Sweet R A (1979) Efficient Fortran subprograms for the solution of separable elliptic partial differential equations *ACM Trans. Math. Software* **5** 352–364

5 Arguments

1: XS – REAL (KIND=nag_wp) *Input*
On entry: the lower bound of the range of x , i.e., $XS \leq x \leq XF$.
Constraint: $XS < XF$.

- 2: XF – REAL (KIND=nag_wp) Input
On entry: the upper bound of the range of x , i.e., $XS \leq x \leq XF$.
Constraint: $XS < XF$.
- 3: L – INTEGER Input
On entry: the number of panels into which the interval (XS,XF) is subdivided. Hence, there will be $L + 1$ grid points in the x -direction given by $x_i = XS + (i - 1) \times \delta x$, for $i = 1, 2, \dots, L + 1$, where $\delta x = (XF - XS)/L$ is the panel width.
Constraint: $L \geq 5$.
- 4: LBDCND – INTEGER Input
On entry: indicates the type of boundary conditions at $x = XS$ and $x = XF$.
 LBDCND = 0
 If the solution is periodic in x , i.e., $u(XS, y, z) = u(XF, y, z)$.
 LBDCND = 1
 If the solution is specified at $x = XS$ and $x = XF$.
 LBDCND = 2
 If the solution is specified at $x = XS$ and the derivative of the solution with respect to x is specified at $x = XF$.
 LBDCND = 3
 If the derivative of the solution with respect to x is specified at $x = XS$ and $x = XF$.
 LBDCND = 4
 If the derivative of the solution with respect to x is specified at $x = XS$ and the solution is specified at $x = XF$.
Constraint: $0 \leq LBDCND \leq 4$.
- 5: BDXS(LDF2, N + 1) – REAL (KIND=nag_wp) array Input
On entry: the values of the derivative of the solution with respect to x at $x = XS$. When $LBDCND = 3$ or 4 , $BDXS(j, k) = u_x(XS, y_j, z_k)$, for $j = 1, 2, \dots, M + 1$ and $k = 1, 2, \dots, N + 1$.
 When $LBDCND$ has any other value, BDXS is not referenced.
- 6: BDXF(LDF2, N + 1) – REAL (KIND=nag_wp) array Input
On entry: the values of the derivative of the solution with respect to x at $x = XF$. When $LBDCND = 2$ or 3 , $BDXF(j, k) = u_x(XF, y_j, z_k)$, for $j = 1, 2, \dots, M + 1$ and $k = 1, 2, \dots, N + 1$.
 When $LBDCND$ has any other value, BDXF is not referenced.
- 7: YS – REAL (KIND=nag_wp) Input
On entry: the lower bound of the range of y , i.e., $YS \leq y \leq YF$.
Constraint: $YS < YF$.
- 8: YF – REAL (KIND=nag_wp) Input
On entry: the upper bound of the range of y , i.e., $YS \leq y \leq YF$.
Constraint: $YS < YF$.

- 9: M – INTEGER *Input*
On entry: the number of panels into which the interval (YS,YF) is subdivided. Hence, there will be $M + 1$ grid points in the y -direction given by $y_j = YS + (j - 1) \times \delta y$, for $j = 1, 2, \dots, M + 1$, where $\delta y = (YF - YS)/M$ is the panel width.
Constraint: $M \geq 5$.
- 10: MBDCND – INTEGER *Input*
On entry: indicates the type of boundary conditions at $y = YS$ and $y = YF$.
 MBDCND = 0
 If the solution is periodic in y , i.e., $u(x, YF, z) = u(x, YS, z)$.
 MBDCND = 1
 If the solution is specified at $y = YS$ and $y = YF$.
 MBDCND = 2
 If the solution is specified at $y = YS$ and the derivative of the solution with respect to y is specified at $y = YF$.
 MBDCND = 3
 If the derivative of the solution with respect to y is specified at $y = YS$ and $y = YF$.
 MBDCND = 4
 If the derivative of the solution with respect to y is specified at $y = YS$ and the solution is specified at $y = YF$.
Constraint: $0 \leq \text{MBDCND} \leq 4$.
- 11: BDYS(LDF, N + 1) – REAL (KIND=nag_wp) array *Input*
On entry: the values of the derivative of the solution with respect to y at $y = YS$. When MBDCND = 3 or 4, $\text{BDYS}(i, k) = u_y(x_i, YS, z_k)$, for $i = 1, 2, \dots, L + 1$ and $k = 1, 2, \dots, N + 1$.
 When MBDCND has any other value, BDYS is not referenced.
- 12: BDYF(LDF, N + 1) – REAL (KIND=nag_wp) array *Input*
On entry: the values of the derivative of the solution with respect to y at $y = YF$. When MBDCND = 2 or 3, $\text{BDYF}(i, k) = u_y(x_i, YF, z_k)$, for $i = 1, 2, \dots, L + 1$ and $k = 1, 2, \dots, N + 1$.
 When MBDCND has any other value, BDYF is not referenced.
- 13: ZS – REAL (KIND=nag_wp) *Input*
On entry: the lower bound of the range of z , i.e., $ZS \leq z \leq ZF$.
Constraint: $ZS < ZF$.
- 14: ZF – REAL (KIND=nag_wp) *Input*
On entry: the upper bound of the range of z , i.e., $ZS \leq z \leq ZF$.
Constraint: $ZS < ZF$.
- 15: N – INTEGER *Input*
On entry: the number of panels into which the interval (ZS,ZF) is subdivided. Hence, there will be $N + 1$ grid points in the z -direction given by $z_k = ZS + (k - 1) \times \delta z$, for $k = 1, 2, \dots, N + 1$, where $\delta z = (ZF - ZS)/N$ is the panel width.
Constraint: $N \geq 5$.

- 16: NBDCND – INTEGER *Input*
- On entry:* specifies the type of boundary conditions at $z = ZS$ and $z = ZF$.
- NBDCND = 0
if the solution is periodic in z , i.e., $u(x, y, ZF) = u(x, y, ZS)$.
- NBDCND = 1
if the solution is specified at $z = ZS$ and $z = ZF$.
- NBDCND = 2
if the solution is specified at $z = ZS$ and the derivative of the solution with respect to z is specified at $z = ZF$.
- NBDCND = 3
if the derivative of the solution with respect to z is specified at $z = ZS$ and $z = ZF$.
- NBDCND = 4
if the derivative of the solution with respect to z is specified at $z = ZS$ and the solution is specified at $z = ZF$.
- Constraint:* $0 \leq \text{NBDCND} \leq 4$.
- 17: BDZS(LDF, M + 1) – REAL (KIND=nag_wp) array *Input*
- On entry:* the values of the derivative of the solution with respect to z at $z = ZS$. When NBDCND = 3 or 4, $\text{BDZS}(i, j) = u_z(x_i, y_j, ZS)$, for $i = 1, 2, \dots, L + 1$ and $j = 1, 2, \dots, M + 1$. When NBDCND has any other value, BDZS is not referenced.
- 18: BDZF(LDF, M + 1) – REAL (KIND=nag_wp) array *Input*
- On entry:* the values of the derivative of the solution with respect to z at $z = ZF$. When NBDCND = 2 or 3, $\text{BDZF}(i, j) = u_z(x_i, y_j, ZF)$, for $i = 1, 2, \dots, L + 1$ and $j = 1, 2, \dots, M + 1$. When NBDCND has any other value, BDZF is not referenced.
- 19: LAMBDA – REAL (KIND=nag_wp) *Input*
- On entry:* the constant λ in the Helmholtz equation. For certain positive values of λ a solution to the differential equation may not exist, and close to these values the solution of the discretized problem will be extremely ill-conditioned. If $\lambda > 0$, then D03FAF will set IFAIL = 3, but will still attempt to find a solution. However, since in general the values of λ for which no solution exists cannot be predicted *a priori*, you are advised to treat any results computed with $\lambda > 0$ with great caution.
- 20: LDF – INTEGER *Input*
- On entry:* the first dimension of the arrays F, BDYS, BDYF, BDZS and BDZF as declared in the (sub)program from which D03FAF is called.
- Constraint:* $\text{LDF} \geq L + 1$.
- 21: LDF2 – INTEGER *Input*
- On entry:* the second dimension of the array F and the first dimension of the arrays BDXS and BDXF as declared in the (sub)program from which D03FAF is called.
- Constraint:* $\text{LDF2} \geq M + 1$.
- 22: F(LDF, LDF2, N + 1) – REAL (KIND=nag_wp) array *Input/Output*
- On entry:* the values of the right-side of the Helmholtz equation and boundary values (if any).
- $$F(i, j, k) = f(x_i, y_j, z_k) \quad i = 2, 3, \dots, L, j = 2, 3, \dots, M \text{ and } k = 2, 3, \dots, N.$$
- On the boundaries F is defined by

LBDCND	$F(1, j, k)$	$F(L + 1, j, k)$	
0	$f(XS, y_j, z_k)$	$f(XS, y_j, z_k)$	
1	$u(XS, y_j, z_k)$	$u(XF, y_j, z_k)$	
2	$u(XS, y_j, z_k)$	$f(XF, y_j, z_k)$	$j = 1, 2, \dots, M + 1$
3	$f(XS, y_j, z_k)$	$f(XF, y_j, z_k)$	$k = 1, 2, \dots, N + 1$
4	$f(XS, y_j, z_k)$	$u(XF, y_j, z_k)$	
MBDCND	$F(i, 1, k)$	$F(i, M + 1, k)$	
0	$f(x_i, YS, z_k)$	$f(x_i, YS, z_k)$	
1	$u(x_i, YS, z_k)$	$u(x_i, YF, z_k)$	
2	$u(x_i, YS, z_k)$	$f(x_i, YF, z_k)$	$i = 1, 2, \dots, L + 1$
3	$f(x_i, YS, z_k)$	$f(x_i, YF, z_k)$	$k = 1, 2, \dots, N + 1$
4	$f(x_i, YS, z_k)$	$u(x_i, YF, z_k)$	
NBDCND	$F(i, j, 1)$	$F(i, j, N + 1)$	
0	$f(x_i, y_j, ZS)$	$f(x_i, y_j, ZS)$	
1	$u(x_i, y_j, ZS)$	$u(x_i, y_j, ZF)$	
2	$u(x_i, y_j, ZS)$	$f(x_i, y_j, ZF)$	$i = 1, 2, \dots, L + 1$
3	$f(x_i, y_j, ZS)$	$f(x_i, y_j, ZF)$	$j = 1, 2, \dots, M + 1$
4	$f(x_i, y_j, ZS)$	$u(x_i, y_j, ZF)$	

Note: if the table calls for both the solution u and the right-hand side f on a boundary, then the solution must be specified.

On exit: contains the solution $u(i, j, k)$ of the finite difference approximation for the grid point (x_i, y_j, z_k) , for $i = 1, 2, \dots, L + 1$, $j = 1, 2, \dots, M + 1$ and $k = 1, 2, \dots, N + 1$.

23: PERTRB – REAL (KIND=nag_wp) *Output*

On exit: PERTRB = 0, unless a solution to Poisson's equation ($\lambda = 0$) is required with a combination of periodic or derivative boundary conditions (LBDCND, MBDCND and NBDCND = 0 or 3). In this case a solution may not exist. PERTRB is a constant, calculated and subtracted from the array F, which ensures that a solution exists. D03FAF then computes this solution, which is a least squares solution to the original approximation. This solution is not unique and is unnormalized. The value of PERTRB should be small compared to the right-hand side F, otherwise a solution has been obtained to an essentially different problem. This comparison should always be made to ensure that a meaningful solution has been obtained.

24: W(LWRK) – REAL (KIND=nag_wp) array *Workspace*

25: LWRK – INTEGER *Input*

On entry: the dimension of the array W as declared in the (sub)program from which D03FAF is called. W is no longer used as workspace, LWRK can therefore be safely set to zero.

Constraint: LWRK \geq 0.

26: IFAIL – INTEGER *Input/Output*

On entry: IFAIL must be set to 0, -1 or 1. If you are unfamiliar with this argument you should refer to Section 3.4 in How to Use the NAG Library and its Documentation for details.

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 argument, the recommended value is 0. **When the value -1 or 1 is used it is essential to test the value of IFAIL on exit.**

On exit: IFAIL = 0 unless the routine detects an error or a warning has been flagged (see Section 6).

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, $XS \geq XF$,
 or $L < 5$,
 or $LBDCND < 0$,
 or $LBDCND > 4$,
 or $YS \geq YF$,
 or $M < 5$,
 or $MBDCND < 0$,
 or $MBDCND > 4$,
 or $ZS \geq ZF$,
 or $N < 5$,
 or $NBDCND < 0$,
 or $NBDCND > 4$,
 or $LDF < L + 1$,
 or $LDF2 < M + 1$.

$IFAIL = 2$

On entry, $LWRK < 0$.

$IFAIL = 3$

On entry, $\lambda > 0$.

$IFAIL = -99$

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

See Section 3.9 in *How to Use the NAG Library and its Documentation* for further information.

$IFAIL = -399$

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

See Section 3.8 in *How to Use the NAG Library and its Documentation* for further information.

$IFAIL = -999$

Dynamic memory allocation failed.

See Section 3.7 in *How to Use the NAG Library and its Documentation* for further information.

7 Accuracy

Not applicable.

8 Parallelism and Performance

D03FAF is threaded by NAG for parallel execution in multithreaded implementations of the NAG Library.

D03FAF makes calls to BLAS and/or LAPACK routines, which may be threaded within the vendor library used by this implementation. Consult the documentation for the vendor library for further information.

Please consult the X06 Chapter Introduction for information on how to control and interrogate the OpenMP environment used within this routine. Please also consult the Users' Note for your implementation for any additional implementation-specific information.

9 Further Comments

The execution time is roughly proportional to $L \times M \times N \times (\log_2 L + \log_2 M + 5)$, but also depends on input arguments LBDCND and MBDCND.

10 Example

This example solves the Helmholtz equation

$$\frac{\partial^2 u}{\partial x^2} + \frac{\partial^2 u}{\partial y^2} + \frac{\partial^2 u}{\partial z^2} + \lambda u = f(x, y, z)$$

for $(x, y, z) \in [0, 1] \times [0, 2\pi] \times [0, \frac{\pi}{2}]$, where $\lambda = -2$, and $f(x, y, z)$ is derived from the exact solution

$$u(x, y, z) = x^4 \sin y \cos z.$$

The equation is subject to the following boundary conditions, again derived from the exact solution given above.

$u(0, y, z)$ and $u(1, y, z)$ are prescribed (i.e., LBDCND = 1).

$u(x, 0, z) = u(x, 2\pi, z)$ (i.e., MBDCND = 0).

$u(x, y, 0)$ and $u_z(x, y, \frac{\pi}{2})$ are prescribed (i.e., NBDCND = 2).

10.1 Program Text

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Program d03faf

!      D03FAF Example Program Text

!      Mark 26 Release. NAG Copyright 2016.

!      .. Use Statements ..
Use nag_library, Only: d03faf, nag_wp
!      .. Implicit None Statement ..
Implicit None
!      .. Parameters ..
Integer, Parameter          :: nin = 5, nout = 6
!      .. Local Scalars ..
Real (Kind=nag_wp)         :: dx, dy, dz, error, lambda, pertrb, &
                             t, xf, xs, yf, ys, yy, zf, zs, zz
Integer                     :: i, ifail, j, k, l, lbdcnd, ldf, &
                             ldf2, lwrk, m, mbdcnd, n, nbdcnd
!      .. Local Arrays ..
Real (Kind=nag_wp), Allocatable :: bdxf(:, :), bdxs(:, :), bdyf(:, :), &
                                   bdys(:, :), bdzf(:, :), bdzs(:, :), &
                                   f(:, :, :), x(:), y(:), z(:)
Real (Kind=nag_wp)         :: w(0)
!      .. Intrinsic Procedures ..
Intrinsic                   :: abs, cos, real, sin
!      .. Executable Statements ..
Write (nout,*) 'D03FAF Example Program Results'
!      Skip heading in data file
Read (nin,*)
Read (nin,*) l, m, n
ldf = l + 1
ldf2 = m + 1
lwrk = 0
Allocate (bdxf(ldf2,n+1),bdxs(ldf2,n+1),bdyf(ldf,n+1),bdys(ldf,n+1), &
         bdzf(ldf,m+1),bdzs(ldf,m+1),f(ldf,ldf2,n+1),x(l+1),y(m+1),z(n+1))
Read (nin,*) lambda
Read (nin,*) xs, xf

```

```

Read (nin,*) ys, yf
Read (nin,*) zs, zf
Read (nin,*) lbdcnd, mbdcnd, nbdcnd

!   Define the grid points for later use.

dx = (xf-xs)/real(l,kind=nag_wp)
Do i = 1, l + 1
  x(i) = xs + real(i-1,kind=nag_wp)*dx
End Do
dy = (yf-ys)/real(m,kind=nag_wp)
Do j = 1, m + 1
  y(j) = ys + real(j-1,kind=nag_wp)*dy
End Do
dz = (zf-zs)/real(n,kind=nag_wp)
Do k = 1, n + 1
  z(k) = zs + real(k-1,kind=nag_wp)*dz
End Do

!   Define the array of derivative boundary values.

Do j = 1, m + 1
  yy = sin(y(j))
  bdzf(1:l+1,j) = -yy*x(1:l+1)**4
End Do

!   Note that for this example all other boundary arrays are
!   dummy variables.

!   We define the function boundary values in the F array.

f(1,1:m+1,1:n+1) = 0.0_nag_wp
Do k = 1, n + 1
  zz = cos(z(k))
  Do j = 1, m + 1
    f(1+1,j,k) = zz*sin(y(j))
  End Do
End Do
f(1:1+1,1:m+1,1) = -bdzf(1:1+1,1:m+1)

!   Define the values of the right hand side of the Helmholtz
!   equation.

Do k = 2, n + 1
  zz = 4.0_nag_wp*cos(z(k))
  Do j = 1, m + 1
    yy = sin(y(j))*zz
    Do i = 2, l
      f(i,j,k) = yy*x(i)**2*(3.0_nag_wp-x(i)**2)
    End Do
  End Do
End Do

!   Call D03FAF to generate and solve the finite difference equation.
!   ifail: behaviour on error exit
!           =0 for hard exit, =1 for quiet-soft, =-1 for noisy-soft
ifail = 0
Call d03faf(xs,xf,l,lbdcnd,bdxc,bdxf,ys,yf,m,mbdcnd,bdys,bdyf,zs,zf,n,
  nbdcnd,bdzc,bdzf,lambda,ldf,ldf2,f,pertrb,w,lwrk,ifail)

!   Compute discretization error. The exact solution to the
!   problem is

!   U(X,Y,Z) = X**4*SIN(Y)*COS(Z)

error = 0.0_nag_wp
Do k = 1, n + 1
  zz = cos(z(k))
  Do j = 1, m + 1
    yy = sin(y(j))*zz
    Do i = 1, l + 1

```



```

        t = abs(f(i,j,k)-yy*x(i)**4)
        If (t>error) Then
            error = t
        End If
    End Do
End Do
Write (nout,*)
Write (nout,99999) error

99999 Format (1X,'Maximum component of discretization error =',1P,E13.6)
End Program d03fafa

```

10.2 Program Data

```

D03FAF Example Program Data
 16 32 20           : l, m, n
 -2.0              : lambda
 0.0 1.0           : xs, xf
 0.0 6.28318530717958647692 : ys, yf
 0.0 1.57079632679489661923 : zs, zf
 1 0 2            : lbdcnd, mbdcnd, nbdcnd

```

10.3 Program Results

D03FAF Example Program Results

Maximum component of discretization error = 5.176553E-04
