# **NAG Fortran Library Routine Document**

# F11DXF

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

F11DXF computes the **approximate** solution of a complex, Hermitian or non-Hermitian, sparse system of linear equations applying a number of Jacobi iterations. It is expected that F11DXF will be used as a preconditioner for the iterative solution of complex sparse systems of equations.

# 2 Specification

```
SUBROUTINE F11DXF(STORE, TRANS, INIT, NITER, N, NNZ, A, IROW, ICOL,

CHECK, B, X, DIAG, WORK, IFAIL)

INTEGER

NITER, N, NNZ, IROW(NNZ), ICOL(NNZ), IFAIL

complex

A(NNZ), B(N), X(N), DIAG(N), WORK(N)

CHARACTER*1

STORE, TRANS, INIT, CHECK
```

# 3 Description

F11DXF computes the **approximate** solution of the complex sparse system of linear equations Ax = b using NITER iterations of the Jacobi algorithm (see also Golub and van Loan (1996), Young (1971)):

$$x_{k+1} = x_k + D^{-1}(b - Ax_k) (1)$$

where k = 1, ..., NITER and  $x_0 = 0$ .

F11DXF can be used both for non-Hermitian and Hermitian systems of equations. For Hermitian matrices, either all non-zero elements of the matrix A can be supplied using coordinate storage (CS), or only the non-zero elements of the lower triangle of A, using symmetric coordinate storage (SCS) (see the F11 Chapter Introduction).

It is expected that F11DXF will be used as a preconditioner for the iterative solution of complex sparse systems of equations, using either the suite comprising the routines F11GRF, F11GSF and F11GTF, for Hermitian systems, or the suite comprising the routines F11BRF, F11BSF and F11BTF, for non-Hermitian systems of equations.

#### 4 References

Golub G H and van Loan C F (1996) Matrix Computations (3rd Edition) Johns Hopkins University Press, Baltimore

Young D (1971) Iterative Solution of Large Linear Systems Academic Press, New York

# 5 Parameters

## 1: STORE – CHARACTER\*1

Input

On entry: specifies whether the matrix A is stored using symmetric coordinate storage (SCS) (applicable only to a Hermitian matrix A) or coordinate storage (CS) (applicable to both Hermitian and non-Hermitian matrices).

If STORE = N', then the complete matrix A is stored in CS format.

If STORE = 'S', then the lower triangle of the Hermitian matrix A is stored in SCS format.

Constraint: STORE = 'N' or 'S'.

## 2: TRANS - CHARACTER\*1

Input

On entry: if STORE = 'N', specifies whether the approximate solution of Ax = b or of  $A^Tx = b$  is required:

if TRANS = 'N', then the approximate solution of Ax = b is calculated;

if TRANS = 'T', then the approximate solution of  $A^Tx = b$  is calculated.

Constraint: TRANS = 'N' or 'T'.

Suggested value: if the matrix A is Hermitian and stored in CS format, it is recommended that TRANS = N' for reasons of efficiency.

#### 3: INIT – CHARACTER\*1

Input

On entry: on first entry, INIT should be set to 'I', unless the diagonal elements of A are already stored in the array DIAG. Otherwise, if DIAG already contains the diagonal of A, it can be set to 'N'.

If INIT = 'N', then DIAG must contain the diagonal of A.

If INIT = 'I', then DIAG will store the diagonal of A on exit.

Constraint: INIT = 'N' or 'I'.

Suggested value: INIT = 'I' on first entry; INIT = 'N', subsequently, unless DIAG has been overwritten.

#### 4: NITER – INTEGER

Input

On entry: the number of Jacobi iterations requested.

Constraint: NITER > 0.

#### 5: N – INTEGER

Input

On entry: n, the order of the matrix A.

Constraint: N > 1.

#### 6: NNZ – INTEGER

Input

On entry: if STORE = 'N', the number of non-zero elements in the matrix A. If STORE = 'S', the number of non-zero elements in the lower triangle of the matrix A.

Constraints:

if STORE = 'N', 
$$1 \le NNZ \le N^2$$
,  
If STORE = 'S',  $1 \le NNZ \le [N(N+1)]/2$ .

# 7: A(NNZ) - complex array

Input

On entry: if STORE = 'N', the non-zero elements in the matrix A (CS format). If STORE = 'S', the non-zero elements in the lower triangle of the matrix A (SCS format). In both cases, the elements of either A or of its lower triangle must be ordered by increasing row index and by increasing column index within each row. Multiple entries for the same row and columns indices are not permitted. The routine F11ZNF or F11ZPF may be used to reorder the elements in this way for CS and SCS storage, respectively.

# 8: IROW(NNZ) – INTEGER array

Input

9: ICOL(NNZ) – INTEGER array

Input

On entry: if STORE = 'N', the row and column indices of the non-zero elements supplied in A. If STORE = 'S', the row and column indices of the non-zero elements of the lower triangle of the matrix A supplied in A.

F11DXF.2 [NP3546/20A]

Constraints:

```
1 \leq IROW(i) \leq N, for i = 1, 2, ..., NNZ; 1 \leq ICOL(i) \leq N, for i = 1, 2, ..., NNZ, when STORE = 'N', or 1 \leq ICOL(i) \leq IROW(i), for i = 1, 2, ..., NNZ, when STORE = 'S', IROW(i-1) < IROW(i) or IROW(i-1) = IROW(i) and ICOL(i-1) < ICOL(i), for i = 2, 3, ..., NNZ.
```

#### 10: CHECK - CHARACTER\*1

Input

On entry: specifies whether or not the CS representation of the matrix A should be checked:

if CHECK = 'C', checks are carried out on the values of N, NNZ, IROW, ICOL; if INIT = 'N', DIAG is also checked;

if CHECK = 'N', none of these checks are carried out.

See also Section 8.2.

Constraint: CHECK = 'C' or 'N'.

# 11: B(N) - complex array

Input

On entry: the right-hand side vector b.

12: X(N) - complex array

Output

On exit: the approximate solution vector  $x_{\text{NITER}}$ .

13: DIAG(N) - complex array

Input/Output

On entry: if INIT = 'N', the diagonal elements of A.

On exit: if INIT = 'N', unchanged on exit. If INIT = 'I', the diagonal elements of A.

14: WORK(N) - complex array

Workspace

15: IFAIL – INTEGER

Input/Output

On entry: IFAIL must be set to 0, -1 or 1. Users who are unfamiliar with this parameter should refer to Chapter P01 for details.

On exit: IFAIL = 0 unless the routine detects an error (see Section 6).

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, for users not familiar with this parameter the recommended value is 0. When the value -1 or 1 is used it is essential to test the value of IFAIL on exit.

# 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, STORE \neq 'N' or 'S', or TRANS \neq 'N' or 'T', or INIT \neq 'N' or 'I', or CHECK \neq 'C' or 'N', or NITER \leq 0.
```

```
\begin{split} \text{IFAIL} &= 2 \\ &\quad \text{On entry, } N < 1, \\ &\quad \text{or} \qquad NNZ < 1, \\ &\quad \text{or} \qquad NNZ > N^2, \text{ if STORE} = \text{'N',} \\ &\quad \text{or} \qquad 1 \leq NNZ \leq [N(N+1)]/2, \text{ if STORE} = \text{'S'.} \end{split}
```

IFAIL = 3

On entry, the arrays IROW and ICOL fail to satisfy the following constraints:

```
1 \leq \mathrm{IROW}(i) \leq \mathrm{N} and  \mathrm{if} \ \mathrm{STORE} = \mathrm{'N'} \ \mathrm{then} \ 1 \leq \mathrm{ICOL}(i) \leq \mathrm{N}, \ \mathrm{or}   \mathrm{if} \ \mathrm{STORE} = \mathrm{'S'} \ \mathrm{then} \ 1 \leq \mathrm{ICOL}(i) \leq \mathrm{IROW}(i), \ \mathrm{for} \ i = 1, 2, \ldots, \mathrm{NNZ}.   \mathrm{IROW}(i-1) < \mathrm{IROW}(i) \ \ \mathrm{or} \ \ \mathrm{IROW}(i-1) = \mathrm{IROW}(i) \ \ \mathrm{and} \ \ \mathrm{ICOL}(i-1) < \mathrm{ICOL}(i), \ \ \mathrm{for}   i = 2, 3, \ldots, \mathrm{NNZ}.
```

Therefore a non-zero element has been supplied which does not lie within the matrix A, is out of order, or has duplicate row and column indices. Call either F11ZAF or F11ZBF to reorder and sum or remove duplicates when STORE = 'N' or STORE = 'S', respectively.

IFAIL = 4

On entry, INIT = 'N' and some diagonal elements of A stored in DIAG are zero.

IFAIL = 5

On entry, INIT = I' and some diagonal elements of A are zero.

# 7 Accuracy

In general, the Jacobi method cannot be used on its own to solve systems of linear equations. The rate of convergence is bound by its spectral properties (see, for example, Golub and van Loan (1996)) and as a solver, the Jacobi method can only be applied to a limited set of matrices. One condition that guarantees convergence is strict diagonal dominance.

However, the Jacobi method can be used successfully as a preconditioner to a wider class of systems of equations. The Jacobi method has good vector/parallel properties, hence it can be applied very efficiently. Unfortunately, it is not possible to provide criteria which define the applicability of the Jacobi method as a preconditioner, and its usefulness must be judged for each case.

## **8** Further Comments

# 8.1 Timing

The time taken for a call to F11DXF is proportional to NITER  $\times$  NNZ.

#### 8.2 Use of CHECK

It is expected that a common use of F11DXF will be as preconditioner for the iterative solution of complex, Hermitian or non-Hermitian, linear systems. In this situation, F11DXF is likely to be called many times. In the interests of both reliability and efficiency, you are recommended to set CHECK to 'C' for the first of such calls, and to 'N' for all subsequent calls.

# 9 Example

This example solves the complex sparse non-Hermitian system of equations Ax = b iteratively using F11DXF as a preconditioner.

F11DXF.4 [NP3546/20A]

## 9.1 Program Text

**Note:** the listing of the example program presented below uses **bold italicised** terms to denote precision-dependent details. Please read the Users' Note for your implementation to check the interpretation of these terms. As explained in the Essential Introduction to this manual, the results produced may not be identical for all implementations.

```
F11DXF Example Program Text.
  Mark 20 Release. NAG Copyright 2001.
   .. Parameters ..
                    NIN, NOUT
   INTEGER
  PARAMETER
                    (NIN=5, NOUT=6)
   INTEGER
                    NMAX, LA, LWORK
                    (NMAX=1000, LA=10000, LWORK=10*NMAX)
  PARAMETER
   .. Local Scalars ..
  real
                    ANORM, SIGMAX, STPLHS, STPRHS, TOL
  INTEGER
                    I, IFAIL, IFAIL1, IFAILX, IREVCM, ITERM, ITN,
                    LWREQ, M, MAXITN, MONIT, N, NITER, NNZ
  LOGICAL
                    LOOP
                    INIT, NORM, PRECON, WEIGHT
   CHARACTER
   CHARACTER*8
                    METHOD
   .. Local Arrays ..
  complex
                    A(LA), B(NMAX), DIAG(NMAX), WORK(LWORK), X(NMAX)
   real
                    WGT (NMAX)
   INTEGER
                    ICOL(LA), IROW(LA)
   .. External Subroutines ..
                   F11BRF, F11BSF, F11BTF, F11DXF, F11XNF
   EXTERNAL
   .. Executable Statements ..
   WRITE (NOUT,*) 'F11DXF Example Program Results'
   Skip heading in data file
   READ (NIN, *)
  READ (NIN,*) N
   IF (N.LE.NMAX) THEN
      Read or initialize the parameters for the iterative solver
      READ (NIN, *) METHOD
      READ (NIN,*) PRECON, NORM, WEIGHT, ITERM
      READ (NIN,*) M, TOL, MAXITN
      READ (NIN,*) MONIT
      ANORM = 0.0e0
      SIGMAX = 0.0e0
      Read the parameters for the preconditioner
      READ (NIN, *) NITER
      Read the number of non-zero elements of the matrix A, then read
      the non-zero elements
      READ (NIN,*) NNZ
      DO 20 I = 1, NNZ
         READ (NIN,*) A(I), IROW(I), ICOL(I)
20
      CONTINUE
      Read right-hand side vector b and initial approximate solution
      READ (NIN, \star) (B(I), I=1, N)
      READ (NIN, *) (X(I), I=1, N)
      Call F11BDF to initialize the solver
      IFAIL = 0
      CALL F11BRF(METHOD, PRECON, NORM, WEIGHT, ITERM, N, M, TOL, MAXITN,
                  ANORM, SIGMAX, MONIT, LWREQ, WORK, LWORK, IFAIL)
      Call repeatedly F11BSF to solve the equations
      Note that the arrays B and X are overwritten
      On final exit, X will contain the solution and B the residual
```

```
vector
         IFAIL = 0
         IREVCM = 0
         LOOP = .TRUE.
         INIT = 'I'
   40
         CONTINUE
         CALL F11BSF(IREVCM, X, B, WGT, WORK, LWREQ, IFAIL)
         IF (IREVCM.EQ.-1) THEN
            IFAILX = -1
            CALL F11XNF('Transpose', N, NNZ, A, IROW, ICOL, 'No checking', X, B,
                          IFAILX)
         ELSE IF (IREVCM.EQ.1) THEN
             IFAILX = -1
             CALL F11XNF('No transpose', N, NNZ, A, IROW, ICOL, 'No checking',
                          X,B,IFAILX)
         ELSE IF (IREVCM.EQ.2) THEN
             IFAIL1 = -1
             CALL F11DXF('Non Hermitian','N', INIT, NITER, N, NNZ, A, IROW,
                          ICOL, 'Check', X, B, DIAG, WORK(LWREQ+1), IFAIL1)
             INIT = 'N'
             IF (IFAIL1.NE.O) IREVCM = 6
         ELSE IF (IREVCM.EQ.3) THEN
             IFAIL1 = 0
             CALL F11BTF (ITN, STPLHS, STPRHS, ANORM, SIGMAX, WORK, LWREQ,
                          IFAIL1)
             WRITE (NOUT, 99999) ITN, STPLHS
         ELSE IF (IREVCM.EQ.4) THEN
            LOOP = .FALSE.
         END IF
         IF (LOOP) GO TO 40
         Obtain information about the computation
         CALL F11BTF(ITN, STPLHS, STPRHS, ANORM, SIGMAX, WORK, LWREQ, IFAIL1)
         Print the output data
         WRITE (NOUT, 99998)
         WRITE (NOUT, 99997) 'Number of iterations for convergence:
           ITN
     +
         WRITE (NOUT, 99996) 'Residual norm:
           STPLHS
         WRITE (NOUT, 99996) 'Right-hand side of termination criterion:',
     +
           STPRHS
         WRITE (NOUT, 99996) '1-norm of matrix A:
           ANORM
         Output x
         WRITE (NOUT, 99995)
         WRITE (NOUT, 99993) (X(I), I=1, N)
         WRITE (NOUT, 99994)
         WRITE (NOUT, 99993) (B(I), I=1, N)
      END IF
      STOP
99999 FORMAT (/1X,'Monitoring at iteration no.', I4,/1X,1P,'residual no',
              'rm: ',e14.4)
99998 FORMAT (/1X, 'Final Results')
99997 FORMAT (1X,A,I4)
99996 FORMAT (1X,A,1P,e14.4)
99995 FORMAT (/2X,' Solution vector') 99994 FORMAT (/2X,' Residual vector')
99993 FORMAT (1X,'(',1P,e16.4,',',1P,e16.4,')')
      END
```

F11DXF.6 [NP3546/20A]

## 9.2 Program Data

```
F11DXF Example Program Data
  'TFQMR'
                              METHOD
  'P' '1' 'N' 1
                              PRECON, NORM, WEIGHT, ITERM
  2 1.0e-6 20
                              M, TOL, MAXITN
                              MONIT
                              NITER
 24
                              NNZ
 ( 2., 1.)
               1
                     1
 (-1., 1.)
(1.,-3.)
               1
                      4
                      8
 (4., 7.)
               2
                      1
 (-3., 0.)
                      2
 ( 2., 4.)
(-7.,-5.)
( 2., 1.)
               2
                      5
               3
                      6
 ( 3., 2.)
                      1
 (-4., 2.)
               4
                      3
 ( 0., 1.)
( 5.,-3.)
               4
                      4
                      7
               4
 (-1., 2.)
               5
                      2
 (8., 6.)
(-3.,-4.)
               5
                      5
               5
                      7
 (-6.,-2.)
               6
                      1
 (5.,-2.)
               6
                      3
 ( 2., 0.)
                      6
               6
 ( 0.,-5.)
               7
                      3
 (-1., 5.)
(6., 2.)
               7
                      5
 (-1., 4.)
                      2
               8
 ( 2., 0.)
               8
                      6
 ( 3., 3.)
               8
                      8
                              A(I), IROW(I), ICOL(I), I=1,...,NNZ
   7., 11.)
1., 24.)
 (-13.,-18.)
 (-10., 3.)
(23., 14.)
(17., -7.)
 ( 15., -3.)
 ( -3., 20.)
                             B(I), I=1,...,N
 ( 0., 0.)
( 0., 0.)
( 0., 0.)
 ( 0., 0.)
 ( 0., 0.)
 (0.,0.)
 (0., 0.)
                              X(I), I=1,...,N
```

## 9.3 Program Results

```
Monitoring at iteration no. residual norm: 1.5062E+02

Monitoring at iteration no. residual norm: 1.5704E+02

Monitoring at iteration no. residual norm: 1.4803E+02

Monitoring at iteration no. residual norm: 8.5215E+01

Monitoring at iteration no. residual norm: 8.5215E+01

Monitoring at iteration no. residual norm: 4.2951E+01
```

#### F11DXF

```
Monitoring at iteration no.
residual norm:
                 2.5055E+01
Monitoring at iteration no.
residual norm: 1.9090E-01
Final Results
Number of iterations for convergence:
                                             3.3292E-08
Residual norm:
Right-hand side of termination criterion:
                                            8.9100E-04
1-norm of matrix A:
                                             2.7000E+01
   Solution vector
                      1.0000E-00)
      1.0000E-00,
      2.0000E+00,
                     -1.0000E-00)
      3.0000E+00,
                      1.0000E-00)
      4.0000E+00,
                      -1.0000E-00)
      3.0000E+00,
                     -1.0000E-00)
      2.0000E+00,
                      1.0000E+00)
      1.0000E+00,
                     -1.0000E-00)
     -9.5326E-10,
                      3.0000E+00)
  Residual vector
     -1.2622E-09,
                      1.6809E-09)
      -2.7422E-09,
                      3.6071E-09)
                      -1.9902E-09)
3.7184E-09)
      6.6013E-10,
      -3.2883E-10,
      3.4339E-09,
                      2.1164E-09)
      7.377E-10,
                      -2.9778E-09)
       3.4214E-10,
                       1.2555E-09)
                       2.7728E-09)
      3.6653E-09,
```

F11DXF.8 (last) [NP3546/20A]