

F11DDF – NAG Fortran Library Routine Document

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.

Note. This routine was introduced into the NAG Fortran Library at Mark 19 and may therefore not be available to all users of the NAG Fortran SMP Library.

1 Purpose

F11DDF solves a system of linear equations involving the preconditioning matrix corresponding to SSOR applied to a real sparse nonsymmetric matrix, represented in coordinate storage format.

2 Specification

```
SUBROUTINE F11DDF(TRANS, N, NNZ, A, IROW, ICOL, RDIAG, OMEGA,
1              CHECK, Y, X, IWORK, IFAIL)
  INTEGER      N, NNZ, IROW(NNZ), ICOL(NNZ), IWORK(2*N+1), IFAIL
  real         A(NNZ), RDIAG(N), OMEGA, Y(N), X(N)
  CHARACTER*1  TRANS, CHECK
```

3 Description

This routine solves a system of linear equations

$$Mx = y, \quad \text{or} \quad M^T x = y,$$

according to the value of the argument TRANS, where the matrix

$$M = \frac{1}{\omega(2 - \omega)}(D + \omega L)D^{-1}(D + \omega U)$$

corresponds to symmetric successive-over-relaxation (SSOR) [1] applied to a linear system $Ax = b$, where A is a real sparse nonsymmetric matrix stored in coordinate storage (CS) format (see Section 2.1.1 of the Chapter Introduction).

In the definition of M given above D is the diagonal part of A , L is the strictly lower triangular part of A , U is the strictly upper triangular part of A , and ω is a user-defined relaxation parameter.

It is envisaged that a common use of F11DDF will be to carry out the preconditioning step required in the application of F11BBF to sparse linear systems. For an illustration of this use of F11DDF see the example program given in Section 9. F11DDF is also used for this purpose by the black-box routine F11DEF.

4 References

- [1] Young D (1971) *Iterative Solution of Large Linear Systems* Academic Press, New York

5 Parameters

- 1:** TRANS — CHARACTER*1 *Input*
On entry: specifies whether or not the matrix M is transposed:
 if TRANS = 'N', then $Mx = y$ is solved;
 if TRANS = 'T', then $M^T x = y$ is solved.
Constraint: TRANS = 'N' or 'T'.
- 2:** N — INTEGER *Input*
On entry: n , the order of the matrix A .
Constraint: $N \geq 1$.

- 3:** NNZ — INTEGER *Input*
On entry: the number of non-zero elements in the matrix A .
Constraint: $1 \leq \text{NNZ} \leq N^2$.
- 4:** A(NNZ) — *real* array *Input*
On entry: the non-zero elements in the matrix A , ordered by increasing row index, and by increasing column index within each row. Multiple entries for the same row and column indices are not permitted. The routine F11ZAF may be used to order the elements in this way.
- 5:** IROW(NNZ) — INTEGER array *Input*
6: ICOL(NNZ) — INTEGER array *Input*
On entry: the row and column indices of the non-zero elements supplied in A .
Constraints: IROW and ICOL must satisfy the following constraints (which may be imposed by a call to F11ZAF):
- $$1 \leq \text{IROW}(i) \leq N, 1 \leq \text{ICOL}(i) \leq N, \text{ for } i = 1, 2, \dots, \text{NNZ}.$$
- $$\text{IROW}(i-1) < \text{IROW}(i), \text{ or}$$
- $$\text{IROW}(i-1) = \text{IROW}(i) \text{ and } \text{ICOL}(i-1) < \text{ICOL}(i), \text{ for } i = 2, 3, \dots, \text{NNZ}.$$
- 7:** RDIAG(N) — *real* array *Input*
On entry: the elements of the diagonal matrix D^{-1} , where D is the diagonal part of A .
- 8:** OMEGA — *real* *Input*
On entry: the relaxation parameter ω .
Constraint: $0.0 < \text{OMEGA} < 2.0$.
- 9:** CHECK — CHARACTER*1 *Input*
On entry: specifies whether or not the CS representation of the matrix M should be checked:
 if CHECK = 'C', checks are carried on the values of N, NNZ, IROW, ICOL and OMEGA;
 if CHECK = 'N', none of these checks are carried out.
 See also Section 8.2.
Constraint: CHECK = 'C' or 'N'.
- 10:** Y(N) — *real* array *Input*
On entry: the right-hand side vector y .
- 11:** X(N) — *real* array *Output*
On exit: the solution vector x .
- 12:** IWORK(2*N+1) — INTEGER array *Workspace*
- 13:** IFAIL — INTEGER *Input/Output*
On entry: IFAIL must be set to 0, -1 or 1. For users not familiar with this parameter (described in Chapter P01) the recommended value is 0.
On exit: IFAIL = 0 unless the routine detects an error (see Section 6).

6 Errors 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 detected by the routine:

`IFAIL = 1`

On entry, `TRANS` \neq 'N' or 'T',
or `CHECK` \neq 'C' or 'N'.

`IFAIL = 2`

On entry, `N` < 1 ,
or `NNZ` < 1 ,
or `NNZ` $> N^2$,
or `OMEGA` lies outside the interval (0.0,2.0),

`IFAIL = 3`

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

$1 \leq \text{IROW}(i) \leq N$ and $1 \leq \text{ICOL}(i) \leq N$, for $i = 1, 2, \dots, \text{NNZ}$.
`IROW`($i - 1$) $<$ `IROW`(i), or
`IROW`($i - 1$) = `IROW`(i) and `ICOL`($i - 1$) $<$ `ICOL`(i), for $i = 2, 3, \dots, \text{NNZ}$.

Therefore a non-zero element has been supplied which does not lie in the matrix A , is out of order, or has duplicate row and column indices. Call `F11ZAF` to reorder and sum or remove duplicates.

`IFAIL = 4`

On entry, the matrix A has a zero diagonal element. The SSOR preconditioner is not appropriate for this problem.

7 Accuracy

If `TRANS = 'N'` the computed solution x is the exact solution of a perturbed system of equations $(M + \delta M)x = y$, where

$$|\delta M| \leq c(n)\epsilon|D + \omega L||D^{-1}||D + \omega U|,$$

$c(n)$ is a modest linear function of n , and ϵ is the *machine precision*. An equivalent result holds when `TRANS = 'T'`.

8 Further Comments

8.1 Timing

The time taken for a call to `F11DDF` is proportional to `NNZ`.

8.2 Use of CHECK

It is expected that a common use of `F11DDF` will be to carry out the preconditioning step required in the application of `F11BBF` to sparse linear systems. In this situation `F11DDF` is likely to be called many times with the same matrix M . 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 program solves a sparse linear system of equations:

$$Ax = b,$$

using RGMRES with SSOR preconditioning.

The RGMRES algorithm itself is implemented by the reverse communication routine F11BBF, which returns repeatedly to the calling program with various values of the parameter IREVCM. This parameter indicates the action to be taken by the calling program.

If IREVCM = 1 a matrix-vector product $v = Au$ is required. This is implemented by a call to F11XAF.

If IREVCM = -1 a transposed matrix-vector product $v = A^T u$ is required in the estimation of the norm of A . This is implemented by a call to F11XAF.

If IREVCM = 2 a solution of the preconditioning equation $Mv = u$ is required. This is achieved by a call to F11DDF.

If IREVCM = 4 F11BBF has completed its tasks. Either the iteration has terminated, or an error condition has arisen.

For further details see the routine document for F11BBF.

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.

```
*      F11DDF Example Program Text
*      Mark 19 Revised. NAG Copyright 1999.
*      .. Parameters ..
      INTEGER          NIN, NOUT
      PARAMETER        (NIN=5,NOUT=6)
      INTEGER          NMAX, LA, LIWORK, LWORK
      PARAMETER        (NMAX=1000,LA=10000,LIWORK=2*NMAX+1,LWORK=10000)
*      .. Local Scalars ..
      real             ANORM, OMEGA, SIGMAX, STPLHS, STPRHS, TOL
      INTEGER          I, IFAIL, IREVCM, ITERM, ITN, LWNEED, LWREQ, M,
+                     MAXITN, MONIT, N, NNZ
      CHARACTER        CKDDF, CKXAF, NORM, PRECON, TRANS, WEIGHT
      CHARACTER*8       METHOD
*      .. Local Arrays ..
      real             A(LA), B(NMAX), RDIAG(NMAX), WGT(NMAX),
+                     WORK(LWORK), X(NMAX)
      INTEGER          ICOL(LA), IROW(LA), IWORK(LIWORK)
*      .. External Subroutines ..
      EXTERNAL         F11BDF, F11BEF, F11BFF, F11DDF, F11XAF
*      .. Intrinsic Functions ..
      INTRINSIC        MAX
*      .. Executable Statements ..
      WRITE (NOUT,*) 'F11DDF Example Program Results'
      WRITE (NOUT,*)
*      Skip heading in data file
      READ (NIN,*)
*
*      Read algorithmic parameters
*
      READ (NIN,*) N
      IF (N.LE.NMAX) THEN
```

```

      READ (NIN,*) NNZ
      READ (NIN,*) METHOD
      READ (NIN,*) PRECON, NORM, ITERM
      READ (NIN,*) M, TOL, MAXITN
      READ (NIN,*) ANORM, SIGMAX
      READ (NIN,*) OMEGA

*
*      Check size of workspace
*
      LWREQ = MAX(N*(M+3)+M*(M+5)+101,7*N+100,(2*N+M)*(M+2)+N+100,
+      10*N+100)
      IF (LWORK.LT.LWREQ) THEN
        WRITE (NOUT,*) 'LWORK must be at least', LWREQ
        STOP
      END IF

*
*      Read the matrix A
*
      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 x
*
      READ (NIN,*) (B(I),I=1,N)
      READ (NIN,*) (X(I),I=1,N)

*
*      Call F11BDF to initialize solver
*
      WEIGHT = 'N'
      MONIT = 0
      IFAIL = 0
      CALL F11BDF(METHOD,PRECON,NORM,WEIGHT,ITERM,N,M,TOL,MAXITN,
+      ANORM,SIGMAX,MONIT,LWNEED,WORK,LWORK,IFAIL)

*
*      Calculate reciprocal diagonal matrix elements if necessary
*
      IF (PRECON.EQ.'P' .OR. PRECON.EQ.'p') THEN

*
        DO 40 I = 1, N
          IWORK(I) = 0
40    CONTINUE

*
        DO 60 I = 1, NNZ
          IF (IROW(I).EQ.ICOL(I)) THEN
            IWORK(IROW(I)) = IWORK(IROW(I)) + 1
            IF (A(I).NE.0.e0) THEN
              RDIAG(IROW(I)) = 1.e0/A(I)
            ELSE
              WRITE (NOUT,*) 'Matrix has a zero diagonal element'
              GO TO 140
            END IF
          END IF
60    CONTINUE

*
        DO 80 I = 1, N
          IF (IWORK(I).EQ.0) THEN
            WRITE (NOUT,*) 'Matrix has a missing diagonal element'

```

```

      GO TO 140
    END IF
    IF (IWORK(I).GE.2) THEN
      WRITE (NOUT,*)
+      'Matrix has a multiple diagonal element'
      GO TO 140
    END IF
80    CONTINUE
*
    END IF
*
*    Call F11BEF to solve the linear system
*
    IREVCM = 0
    CKXAF = 'C'
    CKDDF = 'C'
*
100  CONTINUE
*
    CALL F11BEF(IREVCM,X,B,WGT,WORK,LWORK,IFAIL)
*
    IF (IREVCM.EQ.1) THEN
*
*      Compute matrix-vector product
*
      TRANS = 'Not transposed'
      CALL F11XAF(TRANS,N,NNZ,A,IROW,ICOL,CKXAF,X,B,IFAIL)
      CKXAF = 'N'
      GO TO 100
*
    ELSE IF (IREVCM.EQ.-1) THEN
*
*      Compute transposed matrix-vector product
*
      TRANS = 'Transposed'
      CALL F11XAF(TRANS,N,NNZ,A,IROW,ICOL,CKXAF,X,B,IFAIL)
      CKXAF = 'N'
      GO TO 100
*
    ELSE IF (IREVCM.EQ.2) THEN
*
*      SSOR preconditioning
*
      TRANS = 'Not transposed'
+      CALL F11DDF(TRANS,N,NNZ,A,IROW,ICOL,RDIAG,OMEGA,CKDDF,X,B,
+                  IWORK,IFAIL)
      CKDDF = 'N'
      GO TO 100
*
    ELSE IF (IREVCM.EQ.4) THEN
*
*      Termination
*
      CALL F11BFF(ITN,STPLHS,STPRHS,ANORM,SIGMAX,WORK,LWORK,IFAIL)
*
      WRITE (NOUT,'(A,I10,A)') ' Converged in', ITN, ' iterations'
      WRITE (NOUT,'(A,1P,D16.3)') ' Matrix norm =', ANORM
      WRITE (NOUT,'(A,1P,D16.3)') ' Final residual norm =', STPLHS

```

```

        WRITE (NOUT,*)
*
*      Output x
*
        WRITE (NOUT,*) '          X'
        DO 120 I = 1, N
            WRITE (NOUT,'(1X,1P,D16.4)') X(I)
120      CONTINUE
*
        END IF
*
140      CONTINUE
*
        END IF
        STOP
        END

```

9.2 Program Data

F11DDF Example Program Data

5	N
16	NNZ
'RGMRES'	METHOD
'P' 'I' 1	PRECON, NORM, ITERM
2 1.E-10 1000	M, TOL, MAXITN
0.E0 0.E0	ANORM, SIGMAX
1.1E0	OMEGA
2. 1 1	
1. 1 2	
-1. 1 4	
-3. 2 2	
-2. 2 3	
1. 2 5	
1. 3 1	
5. 3 3	
3. 3 4	
1. 3 5	
-2. 4 1	
-3. 4 4	
-1. 4 5	
4. 5 2	
-2. 5 3	
-6. 5 5	A(I), IROW(I), ICOL(I), I=1,...,NNZ
0. -7. 33.	
-19. -28.	B(I), I=1,...,N
0. 0. 0.	
0. 0.	X(I), I=1,...,N

9.3 Program Results

F11DDF Example Program Results

Converged in 12 iterations
Matrix norm = 1.200E+01
Final residual norm = 3.841E-09

 X
1.0000E+00
2.0000E+00
3.0000E+00
4.0000E+00
5.0000E+00
