

NAG Library Routine Document

F07JVF (ZPTRFS)

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

F07JVF (ZPTRFS) computes error bounds and refines the solution to a complex system of linear equations $AX = B$, where A is an n by n Hermitian positive definite tridiagonal matrix and X and B are n by r matrices, using the modified Cholesky factorization returned by F07JRF (ZPTTRF) and an initial solution returned by F07JSF (ZPTTRS). Iterative refinement is used to reduce the backward error as much as possible.

2 Specification

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SUBROUTINE F07JVF (UPLO, N, NRHS, D, E, DF, EF, B, LDB, X, LDX, FERR, BERR,      &
                  WORK, RWORK, INFO)
INTEGER          N, NRHS, LDB, LDX, INFO
REAL (KIND=nag_wp)  D(*), DF(*), FERR(NRHS), BERR(NRHS), RWORK(N)
COMPLEX (KIND=nag_wp) E(*), EF(*), B(LDB,*), X(LDX,*), WORK(N)
CHARACTER(1)      UPLO

```

The routine may be called by its LAPACK name *zptrfs*.

3 Description

F07JVF (ZPTRFS) should normally be preceded by calls to F07JRF (ZPTTRF) and F07JSF (ZPTTRS). F07JRF (ZPTTRF) computes a modified Cholesky factorization of the matrix A as

$$A = LDL^H,$$

where L is a unit lower bidiagonal matrix and D is a diagonal matrix, with positive diagonal elements. F07JSF (ZPTTRS) then utilizes the factorization to compute a solution, \hat{X} , to the required equations. Letting \hat{x} denote a column of \hat{X} , F07JVF (ZPTRFS) computes a *component-wise backward error*, β , the smallest relative perturbation in each element of A and b such that \hat{x} is the exact solution of a perturbed system

$$(A + E)\hat{x} = b + f, \quad \text{with} \quad |e_{ij}| \leq \beta |a_{ij}|, \quad \text{and} \quad |f_j| \leq \beta |b_j|.$$

The routine also estimates a bound for the *component-wise forward error* in the computed solution defined by $\max |x_i - \hat{x}_i| / \max |\hat{x}_i|$, where x is the corresponding column of the exact solution, X .

Note that the modified Cholesky factorization of A can also be expressed as

$$A = U^H D U,$$

where U is unit upper bidiagonal.

4 References

Anderson E, Bai Z, Bischof C, Blackford S, Demmel J, Dongarra J J, Du Croz J J, Greenbaum A, Hammarling S, McKenney A and Sorensen D (1999) *LAPACK Users' Guide* (3rd Edition) SIAM, Philadelphia <http://www.netlib.org/lapack/lug>

5 Parameters

- 1: UPLO – CHARACTER(1) *Input*
On entry: specifies the form of the factorization as follows:
 UPLO = 'U'
 $A = U^H DU$.
 UPLO = 'L'
 $A = LDL^H$.
Constraint: UPLO = 'U' or 'L'.
- 2: N – INTEGER *Input*
On entry: n , the order of the matrix A .
Constraint: $N \geq 0$.
- 3: NRHS – INTEGER *Input*
On entry: r , the number of right-hand sides, i.e., the number of columns of the matrix B .
Constraint: NRHS ≥ 0 .
- 4: D(*) – REAL (KIND=nag_wp) array *Input*
Note: the dimension of the array D must be at least $\max(1, N)$.
On entry: must contain the n diagonal elements of the matrix of A .
- 5: E(*) – COMPLEX (KIND=nag_wp) array *Input*
Note: the dimension of the array E must be at least $\max(1, N - 1)$.
On entry: if UPLO = 'U', E must contain the $(n - 1)$ superdiagonal elements of the matrix A .
 If UPLO = 'L', E must contain the $(n - 1)$ subdiagonal elements of the matrix A .
- 6: DF(*) – REAL (KIND=nag_wp) array *Input*
Note: the dimension of the array DF must be at least $\max(1, N)$.
On entry: must contain the n diagonal elements of the diagonal matrix D from the LDL^T factorization of A .
- 7: EF(*) – COMPLEX (KIND=nag_wp) array *Input*
Note: the dimension of the array EF must be at least $\max(1, N - 1)$.
On entry: if UPLO = 'U', EF must contain the $(n - 1)$ superdiagonal elements of the unit upper bidiagonal matrix U from the $U^H DU$ factorization of A .
 If UPLO = 'L', EF must contain the $(n - 1)$ subdiagonal elements of the unit lower bidiagonal matrix L from the LDL^H factorization of A .
- 8: B(LDB,*) – COMPLEX (KIND=nag_wp) array *Input*
Note: the second dimension of the array B must be at least $\max(1, NRHS)$.
On entry: the n by r matrix of right-hand sides B .
- 9: LDB – INTEGER *Input*
On entry: the first dimension of the array B as declared in the (sub)program from which F07JVF (ZPTRFS) is called.
Constraint: LDB $\geq \max(1, N)$.

- 10: X(LDX,*) – COMPLEX (KIND=nag_wp) array Input/Output
Note: the second dimension of the array X must be at least $\max(1, \text{NRHS})$.
On entry: the n by r initial solution matrix X .
On exit: the n by r refined solution matrix X .
- 11: LDX – INTEGER Input
On entry: the first dimension of the array X as declared in the (sub)program from which F07JVF (ZPTRFS) is called.
Constraint: $\text{LDX} \geq \max(1, N)$.
- 12: FERR(NRHS) – REAL (KIND=nag_wp) array Output
On exit: estimate of the forward error bound for each computed solution vector, such that $\|\hat{x}_j - x_j\|_\infty / \|\hat{x}_j\|_\infty \leq \text{FERR}(j)$, where \hat{x}_j is the j th column of the computed solution returned in the array X and x_j is the corresponding column of the exact solution X . The estimate is almost always a slight overestimate of the true error.
- 13: BERR(NRHS) – REAL (KIND=nag_wp) array Output
On exit: estimate of the component-wise relative backward error of each computed solution vector \hat{x}_j (i.e., the smallest relative change in any element of A or B that makes \hat{x}_j an exact solution).
- 14: WORK(N) – COMPLEX (KIND=nag_wp) array Workspace
- 15: RWORK(N) – REAL (KIND=nag_wp) array Workspace
- 16: INFO – INTEGER Output
On exit: INFO = 0 unless the routine detects an error (see Section 6).

6 Error Indicators and Warnings

Errors or warnings detected by the routine:

INFO < 0

If INFO = $-i$, the i th argument had an illegal value. An explanatory message is output, and execution of the program is terminated.

7 Accuracy

The computed solution for a single right-hand side, \hat{x} , satisfies an equation of the form

$$(A + E)\hat{x} = b,$$

where

$$\|E\|_\infty = O(\epsilon)\|A\|_\infty$$

and ϵ is the *machine precision*. An approximate error bound for the computed solution is given by

$$\frac{\|\hat{x} - x\|_\infty}{\|x\|_\infty} \leq \kappa(A) \frac{\|E\|_\infty}{\|A\|_\infty},$$

where $\kappa(A) = \|A^{-1}\|_\infty \|A\|_\infty$, the condition number of A with respect to the solution of the linear equations. See Section 4.4 of Anderson *et al.* (1999) for further details.

Routine F07JUF (ZPTCON) can be used to compute the condition number of A .

8 Further Comments

The total number of floating point operations required to solve the equations $AX = B$ is proportional to nr . At most five steps of iterative refinement are performed, but usually only one or two steps are required.

The real analogue of this routine is F07JHF (DPTRFS).

9 Example

This example solves the equations

$$AX = B,$$

where A is the Hermitian positive definite tridiagonal matrix

$$A = \begin{pmatrix} 16.0 & 16.0 - 16.0i & 0 & 0 \\ 16.0 + 16.0i & 41.0 & 18.0 + 9.0i & 0 \\ 0 & 18.0 - 9.0i & 46.0 & 1.0 + 4.0i \\ 0 & 0 & 1.0 - 4.0i & 21.0 \end{pmatrix}$$

and

$$B = \begin{pmatrix} 64.0 + 16.0i & -16.0 - 32.0i \\ 93.0 + 62.0i & 61.0 - 66.0i \\ 78.0 - 80.0i & 71.0 - 74.0i \\ 14.0 - 27.0i & 35.0 + 15.0i \end{pmatrix}.$$

Estimates for the backward errors and forward errors are also output.

9.1 Program Text

```

Program f07jvfe

!      F07JVF Example Program Text

!      Mark 24 Release. NAG Copyright 2012.

!      .. Use Statements ..
      Use nag_library, Only: nag_wp, x04dbf, zptrfs, zptrfs, zptrfs
!      .. Implicit None Statement ..
      Implicit None
!      .. Parameters ..
      Integer, Parameter          :: nin = 5, nout = 6
!      .. Local Scalars ..
      Integer                     :: i, ifail, info, ldb, ldx, n, nrhs
!      .. Local Arrays ..
      Complex (Kind=nag_wp), Allocatable :: b(:,,:), e(:,), ef(:,), work(:,), x(:,,:)
      Real (Kind=nag_wp), Allocatable   :: berr(:,), d(:,), df(:,), ferr(:,),      &
                                         rwork(:)
      Character (1)                :: clabs(1), rlabs(1)
!      .. Executable Statements ..
      Write (nout,*) 'F07JVF Example Program Results'
      Write (nout,*)
      Flush (nout)
!      Skip heading in data file
      Read (nin,*)
      Read (nin,*) n, nrhs
      ldb = n
      ldx = n
      Allocate (b(ldb,nrhs),e(n-1),ef(n-1),work(n),x(ldx,nrhs),berr(nrhs), &
               d(n),df(n),ferr(nrhs),rwork(n))

!      Read the lower bidiagonal part of the tridiagonal matrix A from
!      data file

      Read (nin,*) d(1:n)
      Read (nin,*) e(1:n-1)

```

```

!      Read the right hand matrix B
      Read (nin,*)(b(i,1:nrhs),i=1,n)

!      Copy A into DF and EF, and copy B into X
      df(1:n) = d(1:n)
      ef(1:n-1) = e(1:n-1)
      x(1:n,1:nrhs) = b(1:n,1:nrhs)

!      Factorize the copy of the tridiagonal matrix A
!      The NAG name equivalent of zptrrf is f07jrf
      Call zpttrf(n,df,ef,info)

      If (info==0) Then

!      Solve the equations AX = B
!      The NAG name equivalent of zptrrs is f07jsf
      Call zptrrs('Lower',n,nrhs,df,ef,x,ldx,info)

!      Improve the solution and compute error estimates
!      The NAG name equivalent of zptrfs is f07jvf
      Call zptrfs('Lower',n,nrhs,d,e,df,ef,b,ldb,x,ldx,ferr,berr,work,rwork, &
        info)

!      Print the solution and the forward and backward error estimates

!      ifail: behaviour on error exit
!      =0 for hard exit, =1 for quiet-soft, =-1 for noisy-soft
      ifail = 0
      Call x04dbf('General',' ',n,nrhs,x,ldx,'Bracketed','F7.4', &
        'Solution(s)','Integer',rlabs,'Integer',clabs,80,0,ifail)

      Write (nout,*)
      Write (nout,*) 'Backward errors (machine-dependent)'
      Write (nout,99999) berr(1:nrhs)
      Write (nout,*)
      Write (nout,*) 'Estimated forward error bounds (machine-dependent)'
      Write (nout,99999) ferr(1:nrhs)
    Else
      Write (nout,99998) 'The leading minor of order ', info, &
        ' is not positive definite'
    End If

99999 Format ((3X,1P,7E11.1))
99998 Format (1X,A,I3,A)
      End Program f07jvfe

```

9.2 Program Data

F07JVF Example Program Data

```

      4          2                               :Values of N and NRHS
      16.0      41.0      46.0      21.0       :End of diagonal D
( 16.0, 16.0) ( 18.0, -9.0) (  1.0, -4.0)     :End of sub-diagonal E
( 64.0, 16.0) (-16.0,-32.0)
( 93.0, 62.0) ( 61.0,-66.0)
( 78.0,-80.0) ( 71.0,-74.0)
( 14.0,-27.0) ( 35.0, 15.0)                   :End of matrix B

```

9.3 Program Results

F07JVF Example Program Results

Solution(s)

```

      1          2
1 ( 2.0000, 1.0000) (-3.0000,-2.0000)
2 ( 1.0000, 1.0000) ( 1.0000, 1.0000)
3 ( 1.0000,-2.0000) ( 1.0000,-2.0000)
4 ( 1.0000,-1.0000) ( 2.0000, 1.0000)

```

Backward errors (machine-dependent)

0.0E+00 0.0E+00

Estimated forward error bounds (machine-dependent)

9.0E-12 6.1E-12
