# NAG Fortran Library Routine Document F07JPF (ZPTSVX)

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

F07JPF (ZPTSVX) uses the factorization

$$A = LDL^H$$

to compute 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. Error bounds on the solution and a condition estimate are also provided.

### 2 Specification

The routine may be called by its LAPACK name zptsvx.

### 3 Description

The following steps are performed:

- 1. If FACT = 'N', the matrix A is factorized as  $A = LDL^H$ , where L is a unit lower bidiagonal matrix and D is diagonal. The factorization can also be regarded as having the form  $A = U^H DU$ .
- 2. If the leading i by i principal minor is not positive-definite, then the routine returns with INFO = i. Otherwise, the factored form of A is used to estimate the condition number of the matrix A. If the reciprocal of the condition number is less than **machine precision**, INFO = N + 1 is returned as a warning, but the routine still goes on to solve for X and compute error bounds as described below.
- 3. The system of equations is solved for X using the factored form of A.
- 4. Iterative refinement is applied to improve the computed solution matrix and to calculate error bounds and backward error estimates for it.

### 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 URL: http://www.netlib.org/lapack/lug

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

Higham N J (2002) Accuracy and Stability of Numerical Algorithms (2nd Edition) SIAM, Philadelphia

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#### 5 Parameters

### 1: FACT – CHARACTER\*1

Input

On entry: specifies whether or not the factored form of the matrix A is supplied on entry:

FACT = 'F' on entry, DF and EF contain the factored form of A. D, E, DF and EF will not be modified;

if FACT = 'N', the matrix A will be copied to DF and EF and factored.

Constraint: FACT = 'F' or 'N'.

#### 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  $\geq 0$ .

#### 4: D(\*) – *double precision* array

Input

**Note**: the dimension of the array D must be at least max(1, N).

On entry: the n diagonal elements of the tridiagonal matrix A.

#### 5: E(\*) - complex\*16 array

Input

**Note**: the dimension of the array E must be at least max(1, N - 1).

On entry: the (n-1) sub-diagonal elements of the tridiagonal matrix A.

### 6: DF(\*) – *double precision* array

Input/Output

**Note**: the dimension of the array DF must be at least max(1, N).

On entry: if FACT = 'F', DF contains the n diagonal elements of the diagonal matrix D from the  $LDL^H$  factorization of A.

On exit: if FACT = 'N', DF contains the n diagonal elements of the diagonal matrix D from the  $LDL^H$  factorization of A.

### 7: EF(\*) - complex\*16 array

Input/Output

**Note**: the dimension of the array EF must be at least max(1, N - 1).

On entry: if FACT = 'F', EF contains the (n-1) sub-diagonal elements of the unit bidiagonal factor L from the  $LDL^H$  factorization of A.

On exit: if FACT = 'N', EF contains the (n-1) sub-diagonal elements of the unit bidiagonal factor L from the  $LDL^H$  factorization of A.

### 8: B(LDB,\*) - complex\*16 array

Input

**Note**: the second dimension of the array B must be at least max(1, NRHS).

On entry: the n by r right-hand side matrix B.

### 9: LDB – INTEGER

Input

On entry: the first dimension of the array B as declared in the (sub)program from which F07JPF (ZPTSVX) is called.

Constraint: LDB  $\geq \max(1, N)$ .

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### 10: X(LDX,\*) - complex\*16 array

Output

**Note**: the second dimension of the array X must be at least max(1, NRHS).

On exit: if INFO = 0 or INFO = N + 1, the n by r solution matrix X.

#### 11: LDX – INTEGER

Input

On entry: the first dimension of the array X as declared in the (sub)program from which F07JPF (ZPTSVX) is called.

*Constraint*: LDX  $\geq$  max(1, N).

### 12: RCOND – double precision

Output

On exit: the reciprocal condition number of the matrix A. If RCOND is less than the **machine precision** (in particular, if RCOND = 0), the matrix is singular to working precision. This condition is indicated by a return code of INFO > 0.

#### 13: FERR(\*) – *double precision* array

Output

**Note**: the dimension of the array FERR must be at least max(1, NRHS).

On exit: the forward error bound for each solution vector X(j) (the jth column of the solution matrix X). If  $x_j$  is the true solution corresponding to X(j), FERR(j) is an estimated upper bound for the magnitude of the largest element in  $(X(j) - x_j)$  divided by the magnitude of the largest element in X(j).

### 14: BERR(\*) – *double precision* array

Output

**Note**: the dimension of the array BERR must be at least max(1, NRHS).

On exit: the componentwise relative backward error of each solution vector X(j) (i.e., the smallest relative change in any element of A or B that makes X(j) an exact solution).

### 15: WORK(\*) - complex\*16 array

Workspace

**Note**: the dimension of the array WORK must be at least max(1, N).

#### 16: RWORK(\*) – *double precision* array

Workspace

**Note**: the dimension of the array RWORK must be at least max(1, N).

### 17: 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.

INFO > 0

If INFO = i and  $i \le N$ , the leading minor of order i of A is not positive-definite, so the factorization could not be completed, and the solution has not been computed. RCOND = 0 is returned.

If INFO = i and i = N + 1, U is nonsingular, but RCOND is less than *machine precision*, meaning that the matrix is singular to working precision. Nevertheless, the solution and error bounds are computed because there are a number of situations where the computed solution can be more accurate than the value of RCOND would suggest.

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### 7 Accuracy

For each right-hand side vector b, the computed solution  $\hat{x}$  is the exact solution of a perturbed system of equations  $(A+E)\hat{x}=b$ , where

$$|E| \le c(n)\epsilon |R^T||R|$$
, where  $R = D^{\frac{1}{2}}U$ ,

c(n) is a modest linear function of n, and  $\epsilon$  is the **machine precision**. See Section 10.1 of Higham (2002) for further details.

If x is the true solution, then the computed solution  $\hat{x}$  satisfies a forward error bound of the form

$$\frac{\|x - \hat{x}\|_{\infty}}{\|\hat{x}\|_{\infty}} \le w_c \operatorname{cond}(A, \hat{x}, b)$$

where  $\operatorname{cond}(A,\hat{x},b) = \||A^{-1}|(|A||\hat{x}|+|b|)\|_{\infty}/\|\hat{x}\|_{\infty} \leq \operatorname{cond}(A) = \||A^{-1}||A|\|_{\infty} \leq \kappa_{\infty}(A)$ . If  $\hat{x}$  is the jth column of X, then  $w_c$  is returned in  $\operatorname{BERR}(j)$  and a bound on  $\|x-\hat{x}\|_{\infty}/\|\hat{x}\|_{\infty}$  is returned in  $\operatorname{FERR}(j)$ . See Section 4.4 of Anderson et al. (1999) for further details.

### **8** Further Comments

The number of floating point operations required for the factorization, and for the estimation of the condition number of A is proportional to n. The number of floating point operations required for the solution of the equations, and for the estimation of the forward and backward error is proportional to  $n \times r$ , where r is the number of right-hand sides.

The condition estimation is based upon Equation (15.11) of Higham (2002). For further details of the error estimation, see Section 4.4 of Anderson *et al.* (1999).

The real analogue of this routine is F07JBF (DPTSVX).

# 9 Example

To solve 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

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

Error estimates for the solutions and an estimate of the reciprocal of the condition number of A are also output.

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### 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.

```
F07JPF Example Program Text
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.. Parameters ..
INTEGER
                 NIN, NOUT
PARAMETER
                 (NIN=5,NOUT=6)
INTEGER
                 NMAX
PARAMETER
                 (NMAX=8)
INTEGER
                 LDB, LDX, NRHSMX
                 (LDB=NMAX,LDX=NMAX,NRHSMX=NMAX)
PARAMETER
.. Local Scalars ..
DOUBLE PRECISION RCOND
INTEGER
                 I, IFAIL, INFO, J, N, NRHS
 . Local Arrays ..
COMPLEX *16
                 B(LDB, NRHSMX), E(NMAX-1), EF(NMAX-1), WORK(NMAX),
                 X(LDX,NRHSMX)
DOUBLE PRECISION BERR(NRHSMX), D(NMAX), DF(NMAX), FERR(NRHSMX),
                 RWORK (NMAX)
CHARACTER
                 CLABS(1), RLABS(1)
.. External Subroutines ..
EXTERNAL
                XO4DBF, ZPTSVX
.. Executable Statements ..
WRITE (NOUT,*) 'F07JPF Example Program Results'
WRITE (NOUT, *)
Skip heading in data file
READ (NIN, *)
READ (NIN,*) N, NRHS
IF (N.LE.NMAX .AND. NRHS.LE.NRHSMX) THEN
   Read the lower bidiagonal part of the tridiagonal matrix A and
   the right hand side b from data file
   READ (NIN, \star) (D(I), I=1, N)
   READ (NIN, *) (E(I), I=1, N-1)
   READ (NIN,*) ((B(I,J),J=1,NRHS),I=1,N)
   Solve the equations AX = B for X
   CALL ZPTSVX('Not factored',N,NRHS,D,E,DF,EF,B,LDB,X,LDX,RCOND,
               FERR, BERR, WORK, RWORK, INFO)
   IF ((INFO.EQ.O) .OR. (INFO.EQ.N+1)) THEN
      Print solution, error bounds and condition number
      IFAIL = 0
      CALL XO4DBF('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(J), J=1, NRHS)
      WRITE (NOUT, *)
      WRITE (NOUT, *)
        'Estimated forward error bounds (machine-dependent)'
      WRITE (NOUT, 99999) (FERR(J), J=1, NRHS)
      WRITE (NOUT, *)
      WRITE (NOUT,*) 'Estimate of reciprocal condition number'
      WRITE (NOUT, 99999) RCOND
      IF (INFO.EQ.N+1) THEN
         WRITE (NOUT, *)
         WRITE (NOUT, *)
           'The matrix A is singular to working precision'
      END IF
```

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```
ELSE

WRITE (NOUT,99998) 'The leading minor of order ', INFO,

' is not positive definite'

END IF

ELSE

WRITE (NOUT,*) 'NMAX and/or NRHSMX too small'

END IF

STOP

*
99999 FORMAT (1X,1P,7E11.1)

99998 FORMAT (1X,A,I3,A)

END
```

### 9.2 Program Data

```
F07JPF 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