NAG Fortran Library Routine Document

F07HBF (DPBSVX)

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

F07HBF (DPBSVX) uses the Cholesky factorization

$$A = U^T U$$
 or $A = L L^T$

to compute the solution to a real system of linear equations

AX = B,

where A is an n by n symmetric positive-definite band matrix of bandwidth $(2k_d + 1)$ and X and B are n by r matrices. Error bounds on the solution and a condition estimate are also provided.

2 Specification

```
SUBROUTINE F07HBF(FACT, UPLO, N, KD, NRHS, AB, LDAB, AFB, LDAFB, EQUED,<br/>S, B, LDB, X, LDX, RCOND, FERR, BERR, WORK, IWORK,<br/>INFO1N, KD, NRHS, LDAB, LDAFB, LDB, LDX, IWORK(*), INFO<br/>AB(LDAB,*), AFB(LDAFB,*), S(*), B(LDB,*), X(LDX,*),<br/>RCOND, FERR(*), BERR(*), WORK(*)<br/>FACT, UPLO, EQUED
```

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

3 Description

The following steps are performed:

1. If FACT = 'E', real diagonal scaling factors, D_S , are computed to equilibrate the system:

$$(D_S A D_S) \left(D_S^{-1} X \right) = D_S B A$$

Whether or not the system will be equilibrated depends on the scaling of the matrix A, but if equilibration is used, A is overwritten by D_SAD_S and B by D_SB .

- 2. If FACT = 'N' or 'E', the Cholesky decomposition is used to factor the matrix A (after equilibration if FACT = 'E') as $A = U^T U$, if UPLO = 'U', or $A = LL^T$, if UPLO = 'L', where U is an upper triangular matrix and L is a lower triangular matrix.
- 3. 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.
- 4. The system of equations is solved for X using the factored form of A.
- 5. Iterative refinement is applied to improve the computed solution matrix and to calculate error bounds and backward error estimates for it.
- 6. If equilibration was used, the matrix X is premultiplied by D_S so that it solves the original system before equilibration.

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

5 **Parameters**

FACT - CHARACTER*1 1:

> On entry: specifies whether or not the factored form of the matrix A is supplied on entry, and if not, whether the matrix A should be equilibrated before it is factored:

if FACT = 'F' on entry, AFB contains the factored form of A. If EQUED = 'Y', the matrix A has been equilibrated with scaling factors given by S. AB and AFB will not be modified; if FACT = 'N', the matrix A will be copied to AFB and factored;

if FACT = 'E', the matrix A will be equilibrated if necessary, then copied to AFB and factored.

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

2: UPLO - CHARACTER*1

On entry: if UPLO = 'U', the upper triangle of A is stored.

If UPLO = 'L', the lower triangle of A is stored.

Constraint: UPLO = 'U' or 'L'.

N - INTEGER 3:

On entry: n, the number of linear equations, i.e., the order of the matrix A.

Constraint: N > 0.

KD – **INTEGER** 4:

> On entry: k_d , the number of super-diagonals of the matrix A if UPLO = 'U', or the number of subdiagonals if UPLO = 'L'.

Constraint: $KD \ge 0$.

NRHS - INTEGER 5:

On entry: r, the number of right-hand sides, i.e., the number of columns of the matrix B.

Constraint: NRHS \geq 0.

AB(LDAB,*) – *double precision* array 6:

> Note: the second dimension of the array AB must be at least $\max(1, N)$. To solve the equations Ax = b, where b is a single right-hand side, B may be supplied as a one-dimensional array with length LDB = max(1, N).

> On entry: the upper or lower triangle of the symmetric band matrix A, stored in the first KD + 1rows of the array, except if FACT = 'F' and EQUED = 'Y', AB must contain the equilibrated matrix D_SAD_S . The *j*th column of A is stored in the *j*th column of the array AB as follows:

if UPLO = 'U', $AB(k_d + 1 + i - j, j) = a_{ij}$ for $max(1, j - k_d) \le i \le j$; if UPLO = 'L', $AB(1 + i - j, j) = a_{ij}$ for $j \le i \le \min(n, j + k_d)$.

On exit: if FACT = 'E' and EQUED = 'Y', AB is overwritten by D_SAD_S .

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Input

Input/Output

Input

Input

Input

Input

7: LDAB – INTEGER

On entry: the first dimension of the array AB as declared in the (sub)program from which F07HBF (DPBSVX) is called.

Constraint: $LDAB \ge KD + 1$.

8: AFB(LDAFB,*) – *double precision* array

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

On entry: if FACT = 'F', AFB contains the triangular factor U or L from the Cholesky factorization $A = U^T U$ or $A = LL^T$ of the band matrix A, in the same storage format as A). If EQUED = 'Y', AFB is the factored form of the equilibrated matrix A.

On exit: if FACT = 'N', AFB returns the triangular factor U or L from the Cholesky factorization $A = U^T U$ or $A = LL^T$.

If FACT = 'E', AFB returns the triangular factor U or L from the Cholesky factorization $A = U^T U$ or $A = LL^T$ of the equilibrated matrix A (see the description of AB for the form of the equilibrated matrix).

9: LDAFB – INTEGER

On entry: the first dimension of the array AFB as declared in the (sub)program from which F07HBF (DPBSVX) is called.

Constraint: $LDAFB \ge KD + 1$.

10: EQUED – CHARACTER*1

On entry: if FACT = 'N' or 'E', EQUED need not be set.

If FACT = 'F', EQUED must specify the form of the equilibration that was performed as follows:

if EQUED = 'N', no equilibration;

if EQUED = 'Y', equilibration was performed, i.e., A has been replaced by D_SAD_S .

On exit: if FACT = 'F', EQUED is unchanged from entry.

Otherwise, if $\mathrm{INFO}\geq0,$ EQUED specifies the form of the equilibration that was performed as specified above.

Constraint: if FACT = 'F', EQUED = 'N' or 'Y'.

11: S(*) - double precision array

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

On entry: if FACT = 'N' or 'E', S need not be set.

If FACT = 'F' and EQUED = 'Y', S must contain the scale factors, D_S , for A; each element of S must be positive.

On exit: if FACT = 'F', S is unchanged from entry.

Otherwise, if INFO ≥ 0 and EQUED = 'Y', S contains the scale factors, D_S , for A; each element of S is positive.

12: B(LDB,*) - double precision array

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.

On exit: if EQUED = 'N', B is not modified.

If EQUED = 'Y', B is overwritten by $D_S B$.

Input

Input/Output

Input

Input/Output

Input/Output

Input/Output

On entry: the first dimension of the array B as declared in the (sub)program from which F07HBF (DPBSVX) is called.

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

X(LDX,*) – *double precision* array 14:

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 to the original system of equations. Note that if EQUED = 'Y', A and B are modified on exit, and the solution to the equilibrated system is $D_S^{-1}X$.

15: LDX – INTEGER

On entry: the first dimension of the array X as declared in the (sub)program from which F07HBF (DPBSVX) is called.

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

RCOND - double precision 16:

On exit: if INFO ≥ 0 , an estimate of the reciprocal condition number of the matrix A (after equilibration if that is performed), computed as $\text{RCOND} = 1/(||A||_1 ||A^{-1}||_1)$.

FERR(*) – *double precision* array 17:

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

On exit: if INFO = 0 or INFO = N + 1, an estimate of the forward error bound for each computed solution vector, such that $\|\hat{x}_j - x_j\|_{\infty} / \|x_j\|_{\infty} \leq \text{FERR}(j)$ where \hat{x}_j is the *j*th column of the computed solution returned in the array X and x_i is the corresponding column of the exact solution The estimate is as reliable as the estimate for RCOND, and is almost always a slight X. overestimate of the true error.

BERR(*) – *double precision* array 18:

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

On exit: if INFO = 0 or INFO = N + 1, an estimate of the componentwise relative backward error of each computed solution vector \hat{x}_i (i.e., the smallest relative change in any element of A or B that makes \hat{x}_i an exact solution).

19: WORK(*) - double precision array Workspace

Note: the dimension of the array WORK must be at least $max(1, 3 \times N)$.

20: IWORK(*) - INTEGER array

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

INFO - INTEGER 21:

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.

Input

Input

Output

Output

Output

Output

Workspace

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 $\mathrm{INFO} > 0$

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.

7 Accuracy

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

$$|E| \le c(n)\epsilon |U^T| |U|,$$

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

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

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

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 *j*th 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**

When $n \gg k$, the factorization of A requires approximately $n(k+1)^2$ floating point operations, where k is the number of super-diagonals.

For each right-hand side, computation of the backward error involves a minimum of 8nk floating point operations. Each step of iterative refinement involves an additional 12nk operations. At most 5 steps of iterative refinement are performed, but usually only 1 or 2 steps are required. Estimating the forward error involves solving a number of systems of equations of the form Ax = b; the number is usually 4 or 5 and never more than 11. Each solution involves approximately 4nk operations.

The complex analogue of this routine is F07HPF (ZPBSVX).

9 Example

To solve the equations

$$AX = B$$

where A is the symmetric positive-definite band matrix

$$A = \begin{pmatrix} 5.49 & 2.68 & 0 & 0\\ 2.68 & 5.63 & -2.39 & 0\\ 0 & -2.39 & 2.60 & -2.22\\ 0 & 0 & -2.22 & 5.17 \end{pmatrix}$$

and

$$B = \begin{pmatrix} 22.09 & 5.10\\ 9.31 & 30.81\\ -5.24 & -25.82\\ 11.83 & 22.90 \end{pmatrix}.$$

Error estimates for the solutions, information on equilibration and an estimate of the reciprocal of the condition number of the scaled matrix A are also output.

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.

```
FO7HBF Example Program Text
*
*
     Mark 21 Release. NAG Copyright 2004.
      .. Parameters ..
*
                       NIN, NOUT
      INTEGER
                       (NIN=5,NOUT=6)
     PARAMETER
                       KDMAX, NMAX
     INTEGER
      PARAMETER
                       (KDMAX=4,NMAX=8)
                       LDAB, LDAFB, LDB, LDX, NRHSMX
      INTEGER
                       (LDAB=KDMAX+1,LDAFB=KDMAX+1,LDB=NMAX,LDX=NMAX,
     PARAMETER
     +
                       NRHSMX=NMAX)
     CHARACTER
                       UPLO
                       (UPLO='U')
      PARAMETER
      .. Local Scalars ..
      DOUBLE PRECISION RCOND
      INTEGER
                       I, IFAIL, INFO, J, KD, N, NRHS
      CHARACTER
                       EQUED
      .. Local Arrays .
*
     DOUBLE PRECISION AB(LDAB,NMAX), AFB(LDAFB,NMAX), B(LDB,NRHSMX),
     +
                       BERR(NRHSMX), FERR(NRHSMX), S(NMAX),
     +
                       WORK(3*NMAX), X(LDX,NRHSMX)
      INTEGER
                       IWORK(NMAX)
      .. External Subroutines ..
      EXTERNAL
                 DPBSVX, XO4CAF
      .. Intrinsic Functions .
*
      INTRINSIC
                      MAX, MIN
      .. Executable Statements
*
      WRITE (NOUT, *) 'FO7HBF Example Program Results'
      WRITE (NOUT, *)
      Skip heading in data file
      READ (NIN, *)
     READ (NIN,*) N, KD, NRHS
      IF (N.LE.NMAX .AND. KD.LE.KDMAX .AND. NRHS.LE.NRHSMX) THEN
*
*
         Read the upper or lower triangular part of the band matrix A
*
         from data file
         IF (UPLO.EQ.'U') THEN
            READ (NIN, *) ((AB(KD+1+I-J,J),J=I,MIN(N,I+KD)),I=1,N)
         ELSE IF (UPLO.EQ.'L') THEN
            READ (NIN,*) ((AB(1+I-J,J),J=MAX(1,I-KD),I),I=1,N)
         END IF
*
         Read B from data file
*
*
         READ (NIN,*) ((B(I,J),J=1,NRHS),I=1,N)
*
*
         Solve the equations AX = B for X
         CALL DPBSVX('Equilibration', UPLO, N, KD, NRHS, AB, LDAB, AFB, LDAFB,
                     EQUED, S, B, LDB, X, LDX, RCOND, FERR, BERR, WORK, IWORK,
     +
     +
                     INFO)
*
         IF ((INFO.EQ.O) .OR. (INFO.EQ.N+1)) THEN
*
            Print solution, error bounds, condition number and the form
*
            of equilibration
            TFATL = 0
            CALL X04CAF('General',' ',N,NRHS,X,LDX,'Solution(s)',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
            WRITE (NOUT, *)
            IF (EQUED.EQ.'N') THEN
               WRITE (NOUT,*) 'A has not been equilibrated'
            ELSE IF (EQUED.EQ.'S') THEN
               WRITE (NOUT, *)
                 'A has been row and column scaled as diag(S)*A*diag(S)'
     +
            END IF
            IF (INFO.EQ.N+1) THEN
               WRITE (NOUT, *)
               WRITE (NOUT, *)
     +
                 'The matrix A is singular to working precision'
            END IF
         ELSE
            WRITE (NOUT, 99998) 'The leading minor of order ', INFO,
             ' is not positive definite'
     +
         END IF
      ELSE
         WRITE (NOUT, *)
           'One or more of NMAX, KDMAX and NRHSMX is too small'
     +
      END IF
      STOP
99999 FORMAT ((3X,1P,7E11.1))
99998 FORMAT (1X,A,I3,A)
      END
```

9.2 **Program Data**

*

*

FO7HBF Example Program Data

4 :Values of N, KD and NRHS 1 2 5.49 2.68 5.63 -2.39 2.60 -2.22 5.17 :End of matrix A 22.09 5.10 9.31 30.81 -5.24 -25.82 11.83 22.90 :End of matrix B

9.3 Program Results

FO7HBF Example Program Results

Solution(s) 1 2 1 5.0000 -2.0000 2 -2.0000 6.0000 3 -3.0000 -1.0000 4 1.0000 4.0000 Backward errors (machine-dependent) 6.4E-17 6.3E-17 Estimated forward error bounds (machine-dependent) 2.0E-14 2.9E-14 Estimate of reciprocal condition number 1.3E-02 A has not been equilibrated