F08USF (CHBGST/ZHBGST) - 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.

1 Purpose

F08USF (CHBGST/ZHBGST) reduces a complex Hermitian-definite generalized eigenproblem $Az = \lambda Bz$ to the standard form $Cy = \lambda y$, where A and B are band matrices, A is a complex Hermitian matrix, and B has been factorized by F08UTF (CPBSTF/ZPBSTF).

2 Specification

```
SUBROUTINE FO8USF(VECT, UPLO, N, KA, KB, AB, LDAB, BB, LDBB, X,

LDX, WORK, RWORK, INFO)

ENTRY chbgst(VECT, UPLO, N, KA, KB, AB, LDAB, BB, LDBB, X,

LDX, WORK, RWORK, INFO)

INTEGER N, KA, KB, LDAB, LDBB, LDX, INFO

complex AB(LDAB,*), BB(LDBB,*), X(LDX,*), WORK(*)

real RWORK(*)

CHARACTER*1 VECT, UPLO
```

The ENTRY statement enables the routine to be called by its LAPACK name.

3 Description

To reduce the complex Hermitian-definite generalized eigenproblem $Az = \lambda Bz$ to the standard form $Cy = \lambda y$, where A, B and C are banded, this routine must be preceded by a call to F08UTF (CPBSTF/ZPBSTF) which computes the split Cholesky factorization of the positive-definite matrix B: $B = S^H S$. The split Cholesky factorization, compared with the ordinary Cholesky factorization, allows the work to be approximately halved.

This routine overwrites A with $C = X^H A X$, where $X = S^{-1}Q$ and Q is a unitary matrix chosen (implicitly) to preserve the bandwidth of A. The routine also has an option to allow the accumulation of X, and then, if Z is an eigenvector of C, XZ is an eigenvector of the original system.

4 References

- [1] Crawford C R (1973) Reduction of a band-symmetric generalized eigenvalue problem Comm. ACM 16 41–44
- [2] Kaufman L (1984) Banded eigenvalue solvers on vector machines ACM Trans. Math. Software 10 73–86

5 Parameters

1: VECT — CHARACTER*1

Input

On entry: indicates whether X is to be returned as follows:

```
if VECT = 'N', then X is not returned; if VECT = 'V', then X is returned.
```

Constraint: VECT = 'N' or 'V'.

2: UPLO — CHARACTER*1

Input

On entry: indicates whether the upper or lower triangular part of A is stored as follows:

if UPLO = 'U', then the upper triangular part of A is stored;

if UPLO = 'L', then the lower triangular part of A is stored.

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

3: N — INTEGER

On entry: n, the order of the matrices A and B.

Constraint: $N \geq 0$.

4: KA — INTEGER Input

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

Constraint: $KA \ge 0$.

5: KB — INTEGER

On entry: k_B , the number of super-diagonals of the matrix B if UPLO = 'U', or the number of sub-diagonals if UPLO = 'L'.

Constraint: $KA \ge KB \ge 0$.

6: AB(LDAB,*) - complex array

Input/Output

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

On entry: the n by n Hermitian band matrix A, stored in rows 1 to $k_A + 1$. More precisely, if UPLO = 'U', the elements of the upper triangle of A within the band must be stored with element a_{ij} in $AB(k_A + 1 + i - j, j)$ for $\max(1, j - k_A) \le i \le j$; if UPLO = 'L', the elements of the lower triangle of A within the band must be stored with element a_{ij} in AB(1 + i - j, j) for $j \le i \le \min(n, j + k_A)$.

On exit: the upper or lower triangle of A is overwritten by the corresponding upper or lower triangle of C as specified by UPLO.

7: LDAB — INTEGER Input

On entry: the first dimension of the array AB as declared in the (sub)program from which F08USF (CHBGST/ZHBGST) is called.

Constraint: LDAB \geq KA+1.

8: BB(LDBB,*) — complex array

Input

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

On entry: the banded split Cholesky factor of B as specified by UPLO, N and KB and returned by F08UTF (CPBSTF/ZPBSTF).

9: LDBB — INTEGER Input

On entry: the first dimension of the array BB as declared in the (sub)program from which F08USF (CHBGST/ZHBGST) is called.

Constraint: LDBB \geq KB+1.

10: X(LDX,*) — complex array

Output

Note: the second dimension of the array X must be at least max(1,N) if VECT = 'V', and at least 1 if VECT = 'N'.

On exit: the n by n matrix $X = S^{-1}Q$, if VECT = 'V'.

X is not referenced if VECT = 'N'.

11: LDX — INTEGER Input

On entry: the first dimension of the array X as declared in the (sub)program from which F08USF (CHBGST/ZHBGST) is called.

Constraints:

LDX
$$\geq \max(1,N)$$
 if VECT = 'V',
LDX ≥ 1 if VECT = 'N'.

12: WORK(*) — complex array

Workspace

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

13: RWORK(*) — real array

Workspace

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

14: INFO — INTEGER Output

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

6 Error Indicators and Warnings

INFO < 0

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

7 Accuracy

Forming the reduced matrix C is a stable procedure. However it involves implicit multiplication by B^{-1} . When the routine is used as a step in the computation of eigenvalues and eigenvectors of the original problem, there may be a significant loss of accuracy if B is ill-conditioned with respect to inversion.

8 Further Comments

The total number of real floating-point operations is approximately $20n^2k_B$, when VECT = 'N', assuming $n \gg k_A, k_B$; there are an additional $5n^3(k_B/k_A)$ operations when VECT = 'V'.

The real analogue of this routine is F08UEF (SSBGST/DSBGST).

9 Example

To compute all the eigenvalues of $Az = \lambda Bz$, where

$$A = \begin{pmatrix} -1.13 + 0.00i & 1.94 - 2.10i & -1.40 + 0.25i & 0.00 + 0.00i \\ 1.94 + 2.10i & -1.91 + 0.00i & -0.82 - 0.89i & -0.67 + 0.34i \\ -1.40 - 0.25i & -0.82 + 0.89i & -1.87 + 0.00i & -1.10 - 0.16i \\ 0.00 + 0.00i & -0.67 - 0.34i & -1.10 + 0.16i & 0.50 + 0.00i \end{pmatrix}$$

and

$$B = \begin{pmatrix} 9.89 + 0.00i & 1.08 - 1.73i & 0.00 + 0.00i & 0.00 + 0.00i \\ 1.08 + 1.73i & 1.69 + 0.00i & -0.04 + 0.29i & 0.00 + 0.00i \\ 0.00 + 0.00i & -0.04 - 0.29i & 2.65 + 0.00i & -0.33 + 2.24i \\ 0.00 + 0.00i & 0.00 + 0.00i & -0.33 - 2.24i & 2.17 + 0.00i \end{pmatrix}$$

Here A is Hermitian, B is Hermitian positive-definite, and A and B are treated as band matrices. B must first be factorized by F08UTF (CPBSTF/ZPBSTF). The program calls F08USF (CHBGST/ZHBGST) to reduce the problem to the standard form $Cy = \lambda y$, then F08HSF (CHBTRD/ZHBTRD) to reduce C to tridiagonal form, and F08JFF (SSTERF/DSTERF) to compute the eigenvalues.

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.

```
FO8USF Example Program Text.
   Mark 19 Release. NAG Copyright 1999.
   .. Parameters ..
   INTEGER
                    NIN, NOUT
   PARAMETER
                    (NIN=5,NOUT=6)
   INTEGER
                    NMAX, KMAX, LDAB, LDBB, LDX
   PARAMETER
                    (NMAX=8,KMAX=8,LDAB=KMAX-1,LDBB=KMAX-1,LDX=NMAX)
   .. Local Scalars ..
   INTEGER I, INFO, J, KA, KB, N
   CHARACTER
                   UPLO
   .. Local Arrays ..
                    AB(LDAB, NMAX), BB(LDBB, NMAX), WORK(NMAX),
   complex
                    X(LDX,NMAX)
                    D(NMAX), E(NMAX-1), RWORK(NMAX)
   real
   .. External Subroutines ..
   EXTERNAL
                    ssterf, chbgst, chbtrd, cpbstf
   .. Intrinsic Functions ..
   INTRINSIC
                    MAX, MIN
   .. Executable Statements ..
   WRITE (NOUT,*) 'FO8USF Example Program Results'
   Skip heading in data file
   READ (NIN,*)
   READ (NIN,*) N, KA, KB
   IF (N.LE.NMAX .AND. KA.LE.KMAX .AND. KB.LE.KA) THEN
      Read A and B from data file
      READ (NIN,*) UPLO
      IF (UPLO.EQ.'U') THEN
         DO 20 I = 1, N
            READ (NIN,*) (AB(KA+1+I-J,J),J=I,MIN(N,I+KA))
20
         CONTINUE
         DO 40 I = 1, N
            READ (NIN,*) (BB(KB+1+I-J,J),J=I,MIN(N,I+KB))
40
         CONTINUE
      ELSE IF (UPLO.EQ.'L') THEN
         DO 60 I = 1, N
            READ (NIN,*) (AB(1+I-J,J),J=MAX(1,<math>I-KA),I)
         CONTINUE
60
         DO 80 I = 1, N
            READ (NIN,*) (BB(1+I-J,J),J=MAX(1,I-KB),I)
80
         CONTINUE
      END IF
      Compute the split Cholesky factorization of B
      CALL cpbstf(	ext{UPLO}, 	ext{N}, 	ext{KB}, 	ext{BB}, 	ext{LDBB}, 	ext{INFO})
      WRITE (NOUT, *)
      IF (INFO.GT.O) THEN
         WRITE (NOUT,*) 'B is not positive-definite.'
      ELSE
         Reduce the problem to standard form C*y = lambda*y, storing
```

```
the result in A
            CALL chbgst('N', UPLO, N, KA, KB, AB, LDAB, BB, LDBB, X, LDX, WORK,
                         RWORK, INFO)
            Reduce C to tridiagonal form T = (Q**H)*C*Q
            CALL chbtrd('N', UPLO, N, KA, AB, LDAB, D, E, X, LDX, WORK, INFO)
            Calclate the eigenvalues of T (same as C)
            CALL ssterf(N,D,E,INFO)
            IF (INFO.GT.O) THEN
               WRITE (NOUT,*) 'Failure to converge.'
            ELSE
               Print eigenvalues
               WRITE (NOUT,*) 'Eigenvalues'
               WRITE (NOUT, 99999) (D(I), I=1, N)
            END IF
         END IF
      END IF
      STOP
99999 FORMAT (3X,(8F8.4))
      END
```

9.2 Program Data

9.3 Program Results

```
F08USF Example Program Results

Eigenvalues
-6.6089 -2.0416 0.1603 1.7712
```