NAG Library Routine Document

F08USF (ZHBGST)

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 (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 (ZPBSTF).

2 Specification

```
SUBROUTINE F08USF(VECT, UPLO, N, KA, KB, AB, LDAB, BB, LDBB, X, LDX,
1                 WORK, RWORK, INFO)
INTEGER           N, KA, KB, LDAB, LDBB, LDX, INFO
double precision  RWORK(N)
complex*16        AB(LDAB,*), BB(LDBB,*), X(LDX,*), WORK(N)
CHARACTER*1       VECT, UPLO
```

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

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, F08USF (ZHBGST) must be preceded by a call to F08UTF (ZPBSTF) which computes the split Cholesky factorization of the positive-definite matrix B: $B = S^{\mathrm{H}}S$. The split Cholesky factorization, compared with the ordinary Cholesky factorization, allows the work to be approximately halved.

This routine overwrites A with $C = X^{\mathrm{H}}AX$, 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

Crawford C R (1973) Reduction of a band-symmetric generalized eigenvalue problem *Comm. ACM* **16** 41–44

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.

VECT = 'N'
X is not returned.

VECT = 'V'
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.

UPLO = 'U'
The upper triangular part of A is stored.

UPLO = 'L'
The lower triangular part of A is stored.

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

3: N – INTEGER *Input*

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

Constraint: $\mathrm{N} \geq 0$.

4: KA – INTEGER *Input*

On entry: if UPLO = 'U', the number of superdiagonals, k_a, of the matrix A.

If UPLO = 'L', the number of subdiagonals, k_a, of the matrix A.

Constraint: $\mathrm{KA} \geq 0$.

5: KB – INTEGER *Input*

On entry: if UPLO = 'U', the number of superdiagonals, k_b, of the matrix B.

If UPLO = 'L', the number of subdiagonals, k_b, of the matrix B.

Constraint: $\mathrm{KA} \geq \mathrm{KB} \geq 0$.

6: AB(LDAB,$*$) – ***complex*16*** array *Input/Output*

Note: the second dimension of the array AB must be at least $\max(1, \mathrm{N})$.

On entry: the upper or lower triangle of the n by n Hermitian band matrix A.

The matrix is 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 $\mathrm{AB}(k_a + 1 + i - j, j)$ for $\max(1, j - k_a) \leq i \leq j$;

if UPLO = 'L', the elements of the lower triangle of A within the band must be stored with element A_{ij} in $\mathrm{AB}(1 + i - j, j)$ for $j \leq i \leq \min(n, j + k_a)$.

On exit: the upper or lower triangle of AB 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 (ZHBGST) is called.

Constraint: $\mathrm{LDAB} \geq \mathrm{KA} + 1$.

8: BB(LDBB,$*$) – ***complex*16*** array *Input*

Note: the second dimension of the array BB must be at least $\max(1, \mathrm{N})$.

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

9: LDBB – INTEGER *Input*

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

Constraint: $\mathrm{LDBB} \geq \mathrm{KB} + 1$.

10: X(LDX,$*$) – ***complex*16*** array *Output*

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

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

If VECT = 'N', X is not referenced.

11: LDX – INTEGER *Input*

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

Constraints:

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

12: WORK(N) – ***complex*16*** array *Workspace*

13: RWORK(N) – ***double precision*** array *Workspace*

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 $= -i$, argument i 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 F08USF (ZHBGST) 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 (DSBGST).

9 Example

This example computes 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 (ZPBSTF). The program calls F08USF (ZHBGST) to reduce the problem to the standard form $Cy = \lambda y$, then F08HSF (ZHBTRD) to reduce C to tridiagonal form, and F08JFF (DSTERF) to compute the eigenvalues.

9.1 Program Text

```
*     F08USF 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 ..
      COMPLEX *16      AB(LDAB,NMAX), BB(LDBB,NMAX), WORK(NMAX),
     +                 X(LDX,NMAX)
      DOUBLE PRECISION D(NMAX), E(NMAX-1), RWORK(NMAX)
*     .. External Subroutines ..
      EXTERNAL         DSTERF, ZHBGST, ZHBTRD, ZPBSTF
*     .. Intrinsic Functions ..
      INTRINSIC        MAX, MIN
*     .. Executable Statements ..
      WRITE (NOUT,*) 'F08USF 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,I-KA),I)
   60       CONTINUE
            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 ZPBSTF(UPLO,N,KB,BB,LDBB,INFO)
*
         WRITE (NOUT,*)
         IF (INFO.GT.0) 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 ZHBGST('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 ZHBTRD('N',UPLO,N,KA,AB,LDAB,D,E,X,LDX,WORK,INFO)
*
*           Calclate the eigenvalues of T (same as C)
*
            CALL DSTERF(N,D,E,INFO)
*
            IF (INFO.GT.0) 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
*
99999 FORMAT (3X,(8F8.4))
      END
```

9.2 Program Data

```
F08USF Example Program Data
  4  2  1                                                :Values of N, KA and KB
  'L'                                                    :Value of UPLO
 (-1.13, 0.00)
 ( 1.94, 2.10) (-1.91, 0.00)
 (-1.40,-0.25) (-0.82, 0.89) (-1.87, 0.00)
               (-0.67,-0.34) (-1.10, 0.16) ( 0.50, 0.00)    :End of matrix A
 ( 9.89, 0.00)
 ( 1.08, 1.73) ( 1.69, 0.00)
               (-0.04,-0.29) ( 2.65, 0.00)
                             (-0.33,-2.24) ( 2.17, 0.00)    :End of matrix B
```

9.3 Program Results

```
 F08USF Example Program Results

 Eigenvalues
    -6.6089 -2.0416  0.1603  1.7712
```