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 FO8USF (VECT, UPLO, N, KA, KB, AB, LDAB, BB, LDBB, X, LDX, WORK, RWORK, INFO)

INTEGER

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

REAL (KIND=nag_wp)

RWORK(N)

COMPLEX (KIND=nag_wp)

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: $B = S^{\rm H}S$. 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

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'.

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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: $N \ge 0$.

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: $KA \geq 0$.

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: $KA \ge KB \ge 0$.

6: AB(LDAB, *) - COMPLEX (KIND=nag_wp) array

Input/Output

Note: the second dimension of the array AB must be at least max(1, 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 $AB(k_a+1+i-j,j)$ 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 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: LDAB \geq KA + 1.

8: BB(LDBB,*) - COMPLEX (KIND=nag wp) 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 (ZPBSTF).

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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: LDBB \geq KB + 1.

10: X(LDX,*) - COMPLEX (KIND=nag wp) 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'.

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:

$$\begin{array}{l} \text{if VECT} = \text{'V', LDX} \geq max(1,N);\\ \text{if VECT} = \text{'N', LDX} \geq 1. \end{array}$$

12: WORK(N) – COMPLEX (KIND=nag wp) array

Workspace

13: RWORK(N) – REAL (KIND=nag wp) 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 Parallelism and Performance

F08USF (ZHBGST) is not threaded by NAG in any implementation.

F08USF (ZHBGST) makes calls to BLAS and/or LAPACK routines, which may be threaded within the vendor library used by this implementation. Consult the documentation for the vendor library for further information.

Please consult the X06 Chapter Introduction for information on how to control and interrogate the OpenMP environment used within this routine. Please also consult the Users' Note for your implementation for any additional implementation-specific information.

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9 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).

10 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.

10.1 Program Text

```
Program f08usfe
1
     FO8USF Example Program Text
!
     Mark 25 Release. NAG Copyright 2014.
      .. Use Statements ..
!
     Use nag_library, Only: dsterf, nag_wp, zhbgst, zhbtrd, zpbstf
!
      .. Implicit None Statement ..
     Implicit None
!
      .. Parameters ..
                                       :: nin = 5, nout = 6
     Integer, Parameter
1
      .. Local Scalars ..
                                        :: i, info, j, ka, kb, ldab, ldbb, ldx, n
     Integer
     Character (1)
                                        :: uplo
      .. Local Arrays ..
!
      Complex (Kind=nag_wp), Allocatable :: ab(:,:), bb(:,:), work(:), x(:,:)
     Real (Kind=nag_wp), Allocatable :: d(:), e(:), rwork(:)
     .. Intrinsic Procedures ..
     Intrinsic
                                        :: max, min
!
      .. Executable Statements ..
     Write (nout,*) 'FO8USF Example Program Results'
     Skip heading in data file
     Read (nin,*)
     Read (nin,*) n, ka, kb
      ldab = ka + 1
      ldbb = kb + 1
      ldx = n
     Allocate (ab(ldab,n),bb(ldbb,n),work(n),x(ldx,n),d(n),e(n-1),rwork(n))
!
     Read A and B from data file
     Read (nin,*) uplo
      If (uplo=='U') Then
        Do i = 1, n
         Read (nin,*)(ab(ka+1+i-j,j),j=i,min(n,i+ka))
```

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Do i = 1, n
          Read (nin,*)(bb(kb+1+i-j,j),j=i,min(n,i+kb))
        End Do
      Else If (uplo=='L') Then
        Do i = 1, n
          Read (nin,*)(ab(1+i-j,j),j=max(1,i-ka),i)
        End Do
        Do i = 1, n
          Read (nin,*)(bb(1+i-j,j),j=max(1,i-kb),i)
        End Do
      End If
      Compute the split Cholesky factorization of B The NAG name equivalent of zpbstf is f08utf \,
!
      Call zpbstf(uplo,n,kb,bb,ldbb,info)
      Write (nout,*)
      If (info>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
!
        The NAG name equivalent of zhbgst is f08usf
        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
1
        The NAG name equivalent of zhbtrd is f08hsf
        Call zhbtrd('N',uplo,n,ka,ab,ldab,d,e,x,ldx,work,info)
!
        Calclate the eigenvalues of T (same as C)
        The NAG name equivalent of dsterf is f08jff
!
        Call dsterf(n,d,e,info)
        If (info>0) Then
          Write (nout,*) 'Failure to converge.'
        Else
          Print eigenvalues
          Write (nout,*) 'Eigenvalues'
          Write (nout, 99999) d(1:n)
        End If
      End If
99999 Format (3X, (8F8.4))
    End Program f08usfe
```

10.2 Program Data

10.3 Program Results

```
F08USF Example Program Results

Eigenvalues
-6.6089 -2.0416 0.1603 1.7712
```

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