F01BVF - 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

F01BVF transforms the generalized symmetric-definite eigenproblem $Ax = \lambda Bx$ to the equivalent standard eigenproblem $Cy = \lambda y$, where A, B and C are symmetric band matrices and B is positive-definite. B must have been decomposed by F01BUF.

2 Specification

3 Description

A is a symmetric band matrix of order n and bandwidth $2m_A+1$. The positive-definite symmetric band matrix B, of order n and bandwidth $2m_B+1$, must have been previously decomposed by F01BUF as $ULDL^TU^T$. F01BVF applies U, L and D to A, m_A rows at a time, restoring the band form of A at each stage by plane rotations. The parameter k defines the change-over point in the decomposition of B as used by F01BUF and is also used as a change-over point in the transformations applied by this routine. For maximum efficiency, k should be chosen to be the multiple of m_A nearest to n/2. The resulting symmetric band matrix C is overwritten on A. The eigenvalues of C, and thus of the original problem, may be found using F08HEF (SSBTRD/DSBTRD) and F08JFF (SSTERF/DSTERF). For selected eigenvalues, use F08HEF (SSBTRD/DSBTRD) and F02BFF.

4 References

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

5 Parameters

1: N — INTEGER

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

2: MA1 — INTEGER Input

On entry: $m_A + 1$, where m_A is the number of non-zero super-diagonals in A. Normally MA1 \ll N.

3: MB1 — INTEGER Input

On entry: $m_B + 1$, where m_B is the number of non-zero super-diagonals in B.

Constraint: $MB1 \leq MA1$.

4: M3 — INTEGER Input

On entry: the value of $3m_A + m_B$.

5: K — INTEGER Input

On entry: k, the change-over point in the transformations. It must be the same as the value used by F01BUF in the decomposition of B.

Suggested value: the optimum value is the multiple of m_A nearest to n/2.

Constraint: $MB1 - 1 \le K \le N$.

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6: A(IA,N) - real array

Input/Output

On entry: the upper triangle of the n by n symmetric band matrix A, with the diagonal of the matrix stored in the $(m_A + 1)$ th row of the array, and the m_A super-diagonals within the band stored in the first m_A rows of the array. Each column of the matrix is stored in the corresponding column of the array. For example, if n = 6 and $m_A = 2$, the storage scheme is

Elements in the top left corner of the array need not be set. The following code assigns the matrix elements within the band to the correct elements of the array:

On exit: A is overwritten by the corresponding elements of C.

7: IA — INTEGER Input

On entry: the first dimension of the array A as declared in the (sub)program from which F01BVF is called.

Constraint: IA \geq MA1.

8: B(IB,N) - real array

Input

On entry: the elements of the decomposition of matrix B as returned by F01BUF.

9: IB — INTEGER Input

On entry: the first dimension of the array B as declared in the (sub)program from which F01BVF is called.

Constraint: $IB \ge MB1$.

10: V(IV,M3) - real array

Work space

11: IV — INTEGER

Input

On entry: the first dimension of the array V as declared in the (sub)program from which F01BVF is called.

Constraint: IV $\geq m_A + m_B$.

12: W(M3) - real array

Workspace

13: IFAIL — INTEGER

Input/Output

On entry: IFAIL must be set to 0, -1 or 1. For users not familiar with this parameter (described in Chapter P01) the recommended value is 0.

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

6 Error Indicators and Warnings

Errors detected by the routine:

IFAIL = 1

On entry, MB1 > MA1.

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7 Accuracy

In general the computed system is exactly congruent to a problem $(A+E)x = \lambda(B+F)x$, where ||E|| and ||F|| are of the order of $\epsilon\kappa(B)||A||$ and $\epsilon\kappa(B)||B||$ respectively, where $\kappa(B)$ is the condition number of B with respect to inversion and ϵ is the **machine precision**. This means that when B is positive-definite but not well-conditioned with respect to inversion, the method, which effectively involves the inversion of B, may lead to a severe loss of accuracy in well-conditioned eigenvalues.

8 Further Comments

The time taken by the routine is approximately proportional to $n^2 m_B^2$ and the distance of k from n/2, e.g., k = n/4 and k = 3n/4 take 502% longer.

When B is positive-definite and well-conditioned with respect to inversion, the generalized symmetric eigenproblem can be reduced to the standard symmetric problem $Py = \lambda y$ where $P = L^{-1}AL^{-T}$ and $B = LL^{T}$, the Cholesky factorization.

When A and B are of band form, especially if the bandwidth is small compared with the order of the matrices, storage considerations may rule out the possibility of working with P since it will be a full matrix in general. However, for any factorization of the form $B = SS^T$, the generalized symmetric problem reduces to the standard form

$$S^{-1}AS^{-T}(S^Tx) = \lambda(S^Tx)$$

and there does exist a factorization such that $S^{-1}AS^{-T}$ is still of band form (see Crawford [1]). Writing

$$C = S^{-1}AS^{-T}$$
 and $y = S^Tx$

the standard form is $Cy = \lambda y$ and the bandwidth of C is the maximum bandwidth of A and B.

Each stage in the transformation consists of two phases. The first reduces a leading principal sub-matrix of B to the identity matrix and this introduces non-zero elements outside the band of A. In the second, further transformations are applied which leave the reduced part of B unaltered and drive the extra elements upwards and off the top left corner of A. Alternatively, B may be reduced to the identity matrix starting at the bottom right-hand corner and the extra elements introduced in A can be driven downwards.

The advantage of the $ULDL^TU^T$ decomposition of B is that no extra elements have to be pushed over the whole length of A. If k is taken as approximately n/2, the shifting is limited to halfway. At each stage the size of the triangular bumps produced in A depends on the number of rows and columns of B which are eliminated in the first phase and on the bandwidth of B. The number of rows and columns over which these triangles are moved at each step in the second phase is equal to the bandwidth of A.

In this routine, A is defined as being at least as wide as B and must be filled out with zeros if necessary as it is overwritten with C. The number of rows and columns of B which are effectively eliminated at each stage is m_A .

9 Example

To find the 3 smallest eigenvalues of $Ax = \lambda Bx$, where

$$A = \begin{pmatrix} 11 & 12 & & & & & \\ 12 & 12 & 13 & & & & & \\ & 13 & 13 & 14 & & & & \\ & & 14 & 14 & 15 & & & \\ & & & 15 & 15 & 16 & & \\ & & & & 16 & 16 & 17 & & \\ & & & & & 17 & 17 & 18 & \\ & & & & & & 18 & 19 & \\ & & & & & & & 19 & 19 \end{pmatrix}$$

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```
22
102
       23
      103
             24
                  25
           104
       24
             25
                  105
                         26
                   26
                        106
                               27
                         27
                             107
                                     28
                                           29
                               28
                                    108
                                     29
                                         109
```

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.

```
F01BVF Example Program Text
Mark 18 Revised. NAG Copyright 1997.
.. Parameters ..
INTEGER
                 NMAX, MA1MAX, MB1MAX, M3, IA, IB, IV
PARAMETER
                  (NMAX=20,MA1MAX=8,MB1MAX=8,M3=3*MA1MAX+MB1MAX-4,
                 IA=MA1MAX,IB=MB1MAX,IV=MA1MAX+MB1MAX-2)
INTEGER
                 NIN, NOUT
                  (NIN=5, NOUT=6)
PARAMETER
.. Local Scalars ..
real
                 ABSTOI.
INTEGER
                 I, IFAIL, INFO, J, K, M, M1, M2, MA1, MB1, N,
.. Local Arrays ..
real
                 A(IA,NMAX), B(IB,NMAX), D(NMAX), E(NMAX),
                 R(NMAX), V(IV,M3), W(M3), WORK(4*NMAX)
INTEGER
                 IBLOCK(NMAX), ISPLIT(NMAX), IWORK(3*NMAX)
.. External Subroutines ...
                  ssbtrd, sstebz, F01BUF, F01BVF
EXTERNAL
.. Intrinsic Functions ..
INTRINSIC
                 MAX
.. Executable Statements ..
WRITE (NOUT,*) 'F01BVF Example Program Results'
Skip heading in data file
READ (NIN,*)
READ (NIN,*) N, MA1, MB1
IF (N.GT.O .AND. N.LE.NMAX .AND. MA1.GE.O .AND. MA1.LE.
    MA1MAX .AND. MB1.GE.O .AND. MB1.LE.MB1MAX) THEN
   READ (NIN,*) ((A(J,I),J=MAX(1,MA1+1-I),MA1),I=1,N)
   READ (NIN,*) ((B(J,I),J=MAX(1,MB1+1-I),MB1),I=1,N)
   K = N/2
   IFAIL = 0
   CALL FO1BUF(N, MB1, K, B, IB, W, IFAIL)
   CALL FO1BVF(N,MA1,MB1,M3,K,A,IA,B,IB,V,IV,W,IFAIL)
   CALL ssbtrd('N', 'U', N, MA1-1, A, IA, D, E, W, 1, WORK, INFO)
   IF (INFO.NE.O) THEN
      WRITE (NOUT,99999) 'ssbtrd', INFO
   ELSE
      ABSTOL = 0.0e0
      READ (NIN,*) M1, M2
      CALL sstebz('I', 'E', N, 0.0e0, 0.0e0, M1, M2, ABSTOL, D, E, M, NSPLIT,
```

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9.2 Program Data

```
F01BVF Example Program Data
  9 2 2
   11
   12
         12
   13
         13
   14
   15
         15
   16
         16
   17
         17
   18
         18
   19
         19
   101
   22
        102
    23
        103
    24
        104
    25
        105
    26
        106
    27
        107
   28
        108
   29
        109
  1 3
```

9.3 Program Results

```
F01BVF Example Program Results

Selected eigenvalues
-0.2643 -0.1530 -0.0418
```

 $[NP3390/19/pdf] F01BVF.5 \; (last)$