NAG Toolbox for MATLAB

f08sn

1 Purpose

f08sn computes all the eigenvalues and, optionally, the eigenvectors of a complex generalized Hermitian-definite eigenproblem, of the form

$$Az = \lambda Bz$$
, $ABz = \lambda z$ or $BAz = \lambda z$,

where A and B are Hermitian and B is also positive-definite.

2 Syntax

$$[a, b, w, info] = f08sn(itype, jobz, uplo, a, b, 'n', n)$$

3 Description

f08sn first performs a Cholesky factorization of the matrix B as $B = U^{H}U$, when **uplo** = 'U' or $B = LL^{H}$, when **uplo** = 'L'. The generalized problem is then reduced to a standard symmetric eigenvalue problem

$$Cx = \lambda x$$
,

which is solved for the eigenvalues and, optionally, the eigenvectors; the eigenvectors are then backtransformed to give the eigenvectors of the original problem.

For the problem $Az = \lambda Bz$, the eigenvectors are normalized so that the matrix of eigenvectors, z, satisfies

$$Z^{H}AZ = \Lambda$$
 and $Z^{H}BZ = I$,

where Λ is the diagonal matrix whose diagonal elements are the eigenvalues. For the problem $ABz = \lambda z$ we correspondingly have

$$Z^{-1}AZ^{-H} = \Lambda$$
 and $Z^{H}BZ = I$.

and for $BAz = \lambda z$ we have

$$Z^{H}AZ = \Lambda$$
 and $Z^{H}B^{-1}Z = I$.

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

5 Parameters

5.1 Compulsory Input Parameters

1: itype – int32 scalar

Specifies the problem type to be solved.

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itype = 1

$$Az = \lambda Bz$$
.
itype = 2
 $ABz = \lambda z$.
itype = 3
 $BAz = \lambda z$.

2: **jobz** – **string**

If jobz = 'N', compute eigenvalues only.

If jobz = 'V', compute eigenvalues and eigenvectors.

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

3: **uplo – string**

If uplo = 'U', the upper triangles of A and B are stored.

If $\mathbf{uplo} = 'L'$, the lower triangles of A and B are stored.

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

4: a(lda,*) – complex array

The first dimension of the array \mathbf{a} must be at least $\max(1, \mathbf{n})$

The second dimension of the array must be at least $max(1, \mathbf{n})$

The n by n Hermitian matrix A.

If $\mathbf{uplo} = 'U'$, the upper triangular part of A must be stored and the elements of the array below the diagonal are not referenced.

If $\mathbf{uplo} = 'L'$, the lower triangular part of A must be stored and the elements of the array above the diagonal are not referenced.

5: b(ldb,*) – complex array

The first dimension of the array **b** must be at least $max(1, \mathbf{n})$

The second dimension of the array must be at least $max(1, \mathbf{n})$

The Hermitian positive-definite matrix *B*:

if $\mathbf{uplo} = 'U'$, the leading n by n upper triangular part of \mathbf{b} contains the upper triangular part of the matrix B;

if $\mathbf{uplo} = 'L'$, the leading n by n lower triangular part of \mathbf{b} contains the lower triangular part of the matrix B.

5.2 Optional Input Parameters

1: n - int32 scalar

Default: The first dimension of the arrays **a**, **b** and the second dimension of the arrays **a**, **b**. (An error is raised if these dimensions are not equal.)

n, the order of the matrices A and B.

Constraint: $\mathbf{n} \geq 0$.

5.3 Input Parameters Omitted from the MATLAB Interface

lda, ldb, work, lwork, rwork

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5.4 Output Parameters

1: a(lda,*) - complex array

The first dimension of the array **a** must be at least $max(1, \mathbf{n})$

The second dimension of the array must be at least $max(1, \mathbf{n})$

If $\mathbf{jobz} = \mathbf{V}$, then if $\mathbf{info} = 0$, a contains the matrix Z of eigenvectors. The eigenvectors are normalized as follows:

if **itype** = 1 or 2,
$$Z^HBZ = I$$
; if **itype** = 3, $Z^HB^{-1}Z = I$.

If **jobz** = 'N', the upper triangle (if **uplo** = 'U') or the lower triangle (if **uplo** = 'L') of **a**, including the diagonal, is destroyed.

2: b(ldb,*) – complex array

The first dimension of the array **b** must be at least $max(1, \mathbf{n})$

The second dimension of the array must be at least $max(1, \mathbf{n})$

If $\inf \mathbf{o} \leq \mathbf{n}$, the part of \mathbf{b} containing the matrix contains the triangular factor U or L from the Cholesky factorization $\mathbf{b} = U^{\mathrm{H}}U$ or $\mathbf{b} = LL^{\mathrm{H}}$.

3: $\mathbf{w}(*)$ – double array

Note: the dimension of the array w must be at least $max(1, \mathbf{n})$.

If info = 0, the eigenvalues in ascending order.

4: info – int32 scalar

info = 0 unless the function detects an error (see Section 6).

6 Error Indicators and Warnings

Errors or warnings detected by the function:

$$info = -i$$

If info = -i, parameter i had an illegal value on entry. The parameters are numbered as follows:

1: itype, 2: jobz, 3: uplo, 4: n, 5: a, 6: lda, 7: b, 8: ldb, 9: w, 10: work, 11: lwork, 12: rwork, 13: info.

It is possible that **info** refers to a parameter that is omitted from the MATLAB interface. This usually indicates that an error in one of the other input parameters has caused an incorrect value to be inferred.

info > 0

f07fr or f08fn returned an error code:

- \leq **n** if **info** = i, f08fn failed to converge; i off-diagonal elements of an intermediate tridiagonal form did not converge to zero;
- > **n** if **info** = **n** + i, for $1 \le i \le$ **n**, then the leading minor of order i of B is not positive-definite. The factorization of B could not be completed and no eigenvalues or eigenvectors were computed.

7 Accuracy

If B is ill-conditioned with respect to inversion, then the error bounds for the computed eigenvalues and vectors may be large, although when the diagonal elements of B differ widely in magnitude the

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eigenvalues and eigenvectors may be less sensitive than the condition of *B* would suggest. See Section 4.10 of Anderson *et al.* 1999 for details of the error bounds.

The example program below illustrates the computation of approximate error bounds.

8 Further Comments

The total number of floating-point operations is proportional to n^3 .

The real analogue of this function is f08sa.

9 Example

```
itype = int32(1);
jobz = 'Vectors';
uplo = 'Upper';
a = [complex(-7.36, +0), complex(0.77, -0.43), complex(-0.64, -0.92),
complex(3.01, -6.97);
     complex(0, 0), complex(3.49, +0), complex(2.19, +4.45), complex(1.9, +0)
+3.73);
      complex(0, +0), complex(0, 0), complex(0.12, +0), complex(2.88, -
3.17):
    complex(0, 0), complex(0, 0), complex(0, 0), complex(-2.54, +0)];
  = [complex(3.23, +0), complex(1.51, -1.92), complex(1.9, +0.84),
complex(0.42, +2.5);
     complex(0, +0), complex(3.58, +0), complex(-0.23, +1.11), complex(-
1.18, +1.37);
      complex(0, 0), complex(0, 0), complex(4.09, +0), complex(2.33, -
0.14):
     complex(0, 0), complex(0, 0), complex(0, 0), complex(4.29, 0)];
[aOut, bOut, w, info] = f08sn(itype, jobz, uplo, a, b)
aOut =
  -0.7721 + 1.5598i
                     -0.3504 + 0.6060i
                                            0.2835 - 0.5806i
                                                                 0.2310 -
1.2161i
    0.6038 - 0.1627i
                      -0.0993 + 0.0631i
                                           -0.3769 - 0.3194i
                                                                -0.4710 +
0.4814i
    0.5954 - 0.6430i
                      0.6851 - 0.5987i -0.3338 - 0.0134i
                                                                -0.2242 +
0.6335i
  -0.6810
                    -0.8127
                                        0.6663
                                                          0.8515
bOut =
   1.7972
                         0.8402 - 1.0683i
                                            1.0572 + 0.4674i
                                                                 0.2337 +
1.3910i
                                             -0.4702 - 0.3131i
                         1.3164
                                                                  0.0834 -
0.0368i
         0
                              0
                                             1.5604
                                                                  0.9360 -
0.9900i
                          0
                                            0
                                                          0.6603
       0
   -5.9990
   -2.9936
   0.5047
    3.9990
info =
           0
```

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