NAG C Library Function Document

nag_zhbevd (f08hqc)

1 Purpose

nag_zhbevd (f08hqc) computes all the eigenvalues and, optionally, all the eigenvectors of a complex Hermitian band matrix. If the eigenvectors are requested, then it uses a divide-and-conquer algorithm to compute eigenvalues and eigenvectors. However, if only eigenvalues are required, then it uses the Pal–Walker–Kahan variant of the *QL* or *QR* algorithm.

2 Specification

```
#include <nag.h>
#include <nagf08.h>

void nag_zhbevd (Nag_OrderType order, Nag_JobType job, Nag_UploType uplo,
     Integer n, Integer kd, Complex ab[], Integer pdab, double w[], Complex z[],
     Integer pdz, NagError *fail)
```

3 Description

nag_zhbevd (f08hqc) computes all the eigenvalues and, optionally, all the eigenvectors of a complex Hermitian band matrix A. In other words, it can compute the spectral factorization of A as

$$A = Z\Lambda Z^{\mathrm{H}},$$

where Λ is a real diagonal matrix whose diagonal elements are the eigenvalues λ_i, and Z is the (complex) unitary matrix whose columns are the eigenvectors z_i. Thus

$$Az_i = \lambda_i z_i, \quad i = 1, 2, \ldots, n.$$

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 Arguments

1: **order** – Nag_OrderType *Input*

On entry: the **order** argument specifies the two-dimensional storage scheme being used, i.e., row-major ordering or column-major ordering. C language defined storage is specified by **order** = **Nag_RowMajor**. See Section 2.2.1.4 of the Essential Introduction for a more detailed explanation of the use of this argument.

Constraint: **order** = **Nag_RowMajor** or **Nag_ColMajor**.

2: **job** – Nag_JobType *Input*

On entry: indicates whether eigenvectors are computed.

job = **Nag_DoNothing**

Only eigenvalues are computed.

job = **Nag_EigVecs**

Eigenvalues and eigenvectors are computed.

Constraint: **job** = **Nag_DoNothing** or **Nag_EigVecs**.

3: **uplo** – Nag_UploType *Input*

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

uplo = **Nag_Upper**

The upper triangular part of A is stored.

uplo = **Nag_Lower**

The lower triangular part of A is stored.

Constraint: **uplo** = **Nag_Upper** or **Nag_Lower**.

4: **n** – Integer *Input*

On entry: n, the order of the matrix A.

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

5: **kd** – Integer *Input*

On entry: if **uplo** = **Nag_Upper**, the number of superdiagonals, k, of the matrix A.

If **uplo** = **Nag_Lower**, the number of subdiagonals.

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

6: **ab**[dim] – Complex *Input/Output*

Note: the dimension, *dim*, of the array **ab** must be at least $\max(1, \mathbf{pdab} \times \mathbf{n})$.

On entry: the n by n Hermitian band matrix A with k sub- or super-diagonals. This is stored as a notional two-dimensional array with row elements or column elements stored contiguously. Just the upper or lower triangular part of the array is held depending on the value of **uplo**. The storage of elements a_{ij} depends on the **order** and **uplo** arguments as follows:

if **order** = **Nag_ColMajor** and **uplo** = **Nag_Upper**,
a_{ij} is stored in $\mathbf{ab}[k + i - j + (j-1) \times \mathbf{pdab}]$, for $i = 1, \ldots, n$ and $j = i, \ldots, \min(n, i + k)$;
if **order** = **Nag_ColMajor** and **uplo** = **Nag_Lower**,
a_{ij} is stored in $\mathbf{ab}[i - j + (j-1) \times \mathbf{pdab}]$, for $i = 1, \ldots, n$ and $j = \max(1, i - k), \ldots, i$;
if **order** = **Nag_RowMajor** and **uplo** = **Nag_Upper**,
a_{ij} is stored in $\mathbf{ab}[j - i + (i-1) \times \mathbf{pdab}]$, for $i = 1, \ldots, n$ and $j = i, \ldots, \min(n, i + k)$;
if **order** = **Nag_RowMajor** and **uplo** = **Nag_Lower**,
a_{ij} is stored in $\mathbf{ab}[k + j - i + (i-1) \times \mathbf{pdab}]$, for $i = 1, \ldots, n$ and $j = \max(1, i - k), \ldots, i$.

On exit: A is overwritten by the values generated during the reduction to tridiagonal form.

If **uplo** = **Nag_Upper**, the first superdiagonal and the diagonal of the tridiagonal matrix are returned in rows **kd** and $\mathbf{kd} + 1$ of the array **ab**, respectively.

If **uplo** = **Nag_Lower**, the diagonal and the first subdiagonal of the tridiagonal matrix are returned in the first two rows of the array **ab**.

7: **pdab** – Integer *Input*

On entry: the stride separating row or column elements (depending on the value of **order**) of the matrix A in the array **ab**.

Constraint: $\mathbf{pdab} \geq \mathbf{kd} + 1$.

8: **w**[*dim*] – double *Output*

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

On exit: the eigenvalues of the matrix A in ascending order.

9: **z**[*dim*] – Complex *Output*

Note: the dimension, *dim*, of the array **z** must be at least

$\max(1, \mathbf{pdz} \times \mathbf{n})$ when **job** = **Nag_EigVecs**;
1 when **job** = **Nag_DoNothing**.

If **order** = **Nag_ColMajor**, the (i,j)th element of the matrix Z is stored in $\mathbf{z}[(j-1) \times \mathbf{pdz} + i - 1]$.

If **order** = **Nag_RowMajor**, the (i,j)th element of the matrix Z is stored in $\mathbf{z}[(i-1) \times \mathbf{pdz} + j - 1]$.

On exit: if **job** = **Nag_EigVecs**, **z** is overwritten by the unitary matrix Z which contains the eigenvectors of A. The ith column of Z contains the eigenvector which corresponds to the eigenvalue **w**[i].

If **job** = **Nag_DoNothing**, **z** is not referenced.

10: **pdz** – Integer *Input*

On entry: the stride separating matrix row or column elements (depending on the value of **order**) in the array **z**.

Constraints:

if **job** = **Nag_EigVecs**, **pdz** $\geq \max(1, \mathbf{n})$;
if **job** = **Nag_DoNothing**, **pdz** ≥ 1.

11: **fail** – NagError * *Input/Output*

The NAG error argument (see Section 2.6 of the Essential Introduction).

6 Error Indicators and Warnings

NE_ALLOC_FAIL

Dynamic memory allocation failed.

NE_BAD_PARAM

On entry, argument ⟨*value*⟩ had an illegal value.

NE_CONVERGENCE

The algorithm failed to converge, ⟨*value*⟩ elements of an intermediate tridiagonal form did not converge to zero.

NE_ENUM_INT_2

On entry, **job** = ⟨*value*⟩, **n** = ⟨*value*⟩, **pdz** = ⟨*value*⟩.
Constraint: if **job** = **Nag_DoNothing**, **pdz** ≥ 1.

On entry, **job** = ⟨*value*⟩, **n** = ⟨*value*⟩, **pdz** = ⟨*value*⟩.
Constraint: if **job** = **Nag_EigVecs**, **pdz** $\geq \max(1, \mathbf{n})$.

On entry, **pdz** = ⟨*value*⟩, **job** = ⟨*value*⟩, **n** = ⟨*value*⟩.
Constraint: if **job** = **Nag_EigVecs**, **pdz** $\geq \max(1, \mathbf{n})$;
if **job** = **Nag_DoNothing**, **pdz** ≥ 1.

NE_INT

On entry, **kd** = ⟨*value*⟩.
Constraint: **kd** ≥ 0.

On entry, **n** = ⟨*value*⟩.
Constraint: **n** ≥ 0.

On entry, **pdab** = ⟨*value*⟩.
Constraint: **pdab** > 0.

On entry, **pdz** = ⟨*value*⟩.
Constraint: **pdz** > 0.

NE_INT_2

On entry, **pdab** = ⟨*value*⟩, **kd** = ⟨*value*⟩.
Constraint: **pdab** ≥ **kd** + 1.

NE_INTERNAL_ERROR

An internal error has occurred in this function. Check the function call and any array sizes. If the call is correct then please consult NAG for assistance.

7 Accuracy

The computed eigenvalues and eigenvectors are exact for a nearby matrix $(A + E)$, where

$$\|E\|_2 = \mathrm{O}(\epsilon)\|A\|_2,$$

and ϵ is the ***machine precision***. See Section 4.7 of Anderson *et al.* (1999) for further details.

8 Further Comments

The real analogue of this function is nag_dsbevd (f08hcc).

9 Example

To compute all the eigenvalues and eigenvectors of the Hermitian band matrix A, where

$$A = \begin{pmatrix} 1+0i & 2-1i & 3-1i & 0+0i & 0+0i \\ 2+1i & 2+0i & 3-2i & 4-2i & 0+0i \\ 3+1i & 3+2i & 3+0i & 4-3i & 5-3i \\ 0+0i & 4+2i & 4+3i & 4+0i & 5-4i \\ 0+0i & 0+0i & 5+3i & 5+4i & 5+0i \end{pmatrix}.$$

9.1 Program Text

```
/* nag_zhbevd (f08hqc) Example Program.
 *
 * Copyright 2001 Numerical Algorithms Group.
 *
 * Mark 7, 2001.
 */

#include <stdio.h>
#include <nag.h>
#include <nag_stdlib.h>
#include <nagf08.h>
#include <nagx04.h>

int main(void)
{
  /* Scalars */
  Integer  i, j, k, kd, n, pdab, pdz, w_len;
  Integer  exit_status=0;
  NagError fail;
  Nag_JobType    job;
  Nag_UploType   uplo;
  Nag_OrderType  order;
  /* Arrays */
```

```
  char    uplo_char[2], job_char[2];
  Complex *ab=0, *z=0;
  double  *w=0;

#ifdef NAG_COLUMN_MAJOR
#define AB_UPPER(I,J) ab[(J-1)*pdab + k + I - J - 1]
#define AB_LOWER(I,J) ab[(J-1)*pdab + I - J]
  order = Nag_ColMajor;
#else
#define AB_UPPER(I,J) ab[(I-1)*pdab + J - I]
#define AB_LOWER(I,J) ab[(I-1)*pdab + k + J - I - 1]
  order = Nag_RowMajor;
#endif

  INIT_FAIL(fail);
  Vprintf("nag_zhbevd (f08hqc) Example Program Results\n\n");

  /* Skip heading in data file */
  Vscanf("%*[^\n] ");
  Vscanf("%ld%ld%*[^\n] ", &n, &kd);
  pdab = kd + 1;
  pdz = n;
  w_len = n;

  /* Allocate memory */
  if ( !(ab = NAG_ALLOC(pdab * n, Complex)) ||
       !(w = NAG_ALLOC(w_len, double)) ||
       !(z = NAG_ALLOC(n * n, Complex)) )
    {
      Vprintf("Allocation failure\n");
      exit_status = -1;
      goto END;
    }
  /* Read whether Upper or Lower part of A is stored */
  Vscanf(" ' %1s '%*[^\n] ", uplo_char);
  if (*(unsigned char *)uplo_char == 'L')
    uplo = Nag_Lower;
  else if (*(unsigned char *)uplo_char == 'U')
    uplo = Nag_Upper;
  else
    {
      Vprintf("Unrecognised character for Nag_UploType type\n");
      exit_status = -1;
      goto END;
    }
  /* Read A from data file */
  k = kd + 1;
  if (uplo == Nag_Upper)
    {
      for (i = 1; i <= n; ++i)
        {
          for (j = i; j <= MIN(i+kd,n); ++j)
            {
              Vscanf(" ( %lf , %lf )", &AB_UPPER(i,j).re,
                     &AB_UPPER(i,j).im);
            }
        }
      Vscanf("%*[^\n] ");
    }
  else
    {
      for (i = 1; i <= n; ++i)
        {
          for (j = MAX(1,i-kd); j <= i; ++j)
            {
              Vscanf(" ( %lf , %lf )", &AB_LOWER(i,j).re,
                     &AB_LOWER(i,j).im);
            }
        }
      Vscanf("%*[^\n] ");
    }
```

```
  /* Read type of job to be performed */
  Vscanf(" ' %1s '%*[^\n] ", job_char);
  if (*(unsigned char *)job_char == 'V')
    job = Nag_EigVecs;
  else
    job = Nag_DoNothing;
  /* Calculate all the eigenvalues and eigenvectors of A */
  /* nag_zhbevd (f08hqc).
   * All eigenvalues and optionally all eigenvectors of
   * complex Hermitian band matrix (divide-and-conquer)
   */
  nag_zhbevd(order, job, uplo, n, kd, ab, pdab, w, z, pdz, &fail);
  if (fail.code != NE_NOERROR)
    {
      Vprintf("Error from nag_zhbevd (f08hqc).\n%s\n", fail.message);
      exit_status = 1;
      goto END;
    }
  /* Print eigenvalues and eigenvectors */
  Vprintf(" Eigenvalues\n");
  for (i = 0; i < n; ++i)
    Vprintf("   %5ld     %8.4f\n", i+1, w[i]);
  Vprintf("\n");
  /* nag_gen_complx_mat_print_comp (x04dbc).
   * Print complex general matrix (comprehensive)
   */
  nag_gen_complx_mat_print_comp(order, Nag_GeneralMatrix, Nag_NonUnitDiag, n, n,
                                z, pdz, Nag_AboveForm, "%7.4f", "Eigenvectors",
                                Nag_IntegerLabels, 0, Nag_IntegerLabels, 0, 80,
                                0, 0, &fail);
  if (fail.code != NE_NOERROR)
    {
      Vprintf("Error from nag_gen_complx_mat_print_comp (x04dbc).\n%s\n",
              fail.message);
      exit_status = 1;
      goto END;
    }
 END:
  if (ab) NAG_FREE(ab);
  if (w) NAG_FREE(w);
  if (z) NAG_FREE(z);
  return exit_status;
}
```

9.2 Program Data

```
nag_zhbevd (f08hqc) Example Program Data
  5  2                                                    :Values of N and KD
  'L'                                                     :Value of UPLO
  (1.0, 0.0)
  (2.0, 1.0) (2.0, 0.0)
  (3.0, 1.0) (3.0, 2.0) (3.0, 0.0)
             (4.0, 2.0) (4.0, 3.0) (4.0, 0.0)
                        (5.0, 3.0) (5.0, 4.0) (5.0, 0.0) :End of matrix A
  'V'                                                     :Value of JOB
```

9.3 Program Results

```
nag_zhbevd (f08hqc) Example Program Results

 Eigenvalues
       1      -6.4185
       2      -1.4094
       3       1.4421
       4       4.4856
       5      16.9002

 Eigenvectors

         1       2       3       4       5
```

```
 1 -0.2591  0.6367  0.4516  0.5503  0.1439
   -0.0000 -0.0000 -0.0000 -0.0000 -0.0000

 2  0.0245 -0.2578 -0.3029  0.4785  0.3060
    0.4344  0.2413 -0.4402  0.2759  0.0411

 3  0.5159 -0.3039  0.3160  0.2128  0.4681
   -0.1095 -0.3481  0.2978  0.0465  0.2306

 4  0.0004  0.3450 -0.4088 -0.1707  0.4098
   -0.5093 -0.0832 -0.3213  0.0200  0.3832

 5 -0.4333 -0.2469  0.0204  0.0175  0.1819
    0.1353  0.2634  0.2262 -0.5611  0.5136
```