

## NAG Library Function Document

### **nag\_zhegvd (f08sqc)**

## 1 Purpose

nag\_zhegvd (f08sqc) computes all the eigenvalues and, optionally, the eigenvectors of a complex generalized Hermitian-definite eigenproblem, of the form

$$Az = \lambda Bz, \quad ABz = \lambda z \quad \text{or} \quad BAz = \lambda z,$$

where  $A$  and  $B$  are Hermitian and  $B$  is also positive definite. If eigenvectors are desired, it uses a divide-and-conquer algorithm.

## 2 Specification

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

void nag_zhegvd (Nag_OrderType order, Integer itype, Nag_JobType job,
                 Nag_UptoType uplo, Integer n, Complex a[], Integer pda, Complex b[],
                 Integer pdb, double w[], NagError *fail)
```

## 3 Description

nag\_zhegvd (f08sqc) first performs a Cholesky factorization of the matrix  $B$  as  $B = U^H U$ , when **uplo** = Nag\_Upper or  $B = LL^H$ , when **uplo** = Nag\_Lower. 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

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

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

$$Z^{-1} A Z^{-H} = \Lambda \quad \text{and} \quad Z^H B Z = I,$$

and for  $BAz = \lambda z$  we have

$$Z^H A Z = \Lambda \quad \text{and} \quad Z^H B^{-1} Z = I.$$

## 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 <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 3.2.1.3 in the Essential Introduction for a more detailed explanation of the use of this argument.  
*Constraint:* **order** = Nag\_RowMajor or Nag\_ColMajor.
- 2: **itype** – Integer *Input*  
*On entry:* specifies the problem type to be solved.  
**itype** = 1  
 $Az = \lambda z.$   
**itype** = 2  
 $ABz = \lambda z.$   
**itype** = 3  
 $BAz = \lambda z.$   
*Constraint:* **itype** = 1, 2 or 3.
- 3: **job** – Nag\_JobType *Input*  
*On entry:* indicates whether eigenvectors are computed.  
**job** = Nag\_EigVals  
Only eigenvalues are computed.  
**job** = Nag\_DoBoth  
Eigenvalues and eigenvectors are computed.  
*Constraint:* **job** = Nag\_EigVals or Nag\_DoBoth.
- 4: **uplo** – Nag\_UptoType *Input*  
*On entry:* if **uplo** = Nag\_Upper, the upper triangles of  $A$  and  $B$  are stored.  
If **uplo** = Nag\_Lower, the lower triangles of  $A$  and  $B$  are stored.  
*Constraint:* **uplo** = Nag\_Upper or Nag\_Lower.
- 5: **n** – Integer *Input*  
*On entry:*  $n$ , the order of the matrices  $A$  and  $B$ .  
*Constraint:* **n**  $\geq 0$ .
- 6: **a**[*dim*] – Complex *Input/Output*  
**Note:** the dimension, *dim*, of the array **a** must be at least  $\max(1, \mathbf{pda} \times \mathbf{n})$ .  
*On entry:* the  $n$  by  $n$  Hermitian matrix  $A$ .  
If **order** = 'Nag\_ColMajor',  $A_{ij}$  is stored in **a**[(*j* - 1)  $\times$  **pda** + *i* - 1].  
If **order** = 'Nag\_RowMajor',  $A_{ij}$  is stored in **a**[(*i* - 1)  $\times$  **pda** + *j* - 1].  
If **uplo** = 'Nag\_Upper', the upper triangular part of  $A$  must be stored and the elements of the array below the diagonal are not referenced.  
If **uplo** = 'Nag\_Lower', the lower triangular part of  $A$  must be stored and the elements of the array above the diagonal are not referenced.  
*On exit:* if **job** = Nag\_DoBoth, **a** contains the matrix  $Z$  of eigenvectors. The eigenvectors are normalized as follows:

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

If **job** = Nag\_EigVals, the upper triangle (if **uplo** = Nag\_Upper) or the lower triangle (if **uplo** = Nag\_Lower) of **a**, including the diagonal, is overwritten.

7: **pda** – Integer *Input*

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

*Constraint:*  $\mathbf{pda} \geq \max(1, \mathbf{n})$ .

8: **b[dim]** – Complex *Input/Output*

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

*On entry:* the *n* by *n* Hermitian matrix *B*.

If **order** = 'Nag\_ColMajor',  $B_{ij}$  is stored in **b**[(*j* – 1) ×  **pdb** + *i* – 1].

If **order** = 'Nag\_RowMajor',  $B_{ij}$  is stored in **b**[(*i* – 1) ×  **pdb** + *j* – 1].

If **uplo** = 'Nag\_Upper', the upper triangular part of *B* must be stored and the elements of the array below the diagonal are not referenced.

If **uplo** = 'Nag\_Lower', the lower triangular part of *B* must be stored and the elements of the array above the diagonal are not referenced.

*On exit:* the triangular factor *U* or *L* from the Cholesky factorization  $B = U^H U$  or  $B = LL^H$ .

9: **pdb** – Integer *Input*

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

*Constraint:*  $\mathbf{pdb} \geq \max(1, \mathbf{n})$ .

10: **w[n]** – double *Output*

*On exit:* the eigenvalues in ascending order.

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

The NAG error argument (see Section 3.6 in the Essential Introduction).

## 6 Error Indicators and Warnings

### NE\_ALLOC\_FAIL

Dynamic memory allocation failed.

### NE\_BAD\_PARAM

On entry, argument  $\langle\text{value}\rangle$  had an illegal value.

### NE\_CONVERGENCE

The algorithm failed to converge;  $\langle\text{value}\rangle$  off-diagonal elements of an intermediate tridiagonal form did not converge to zero.

### NE\_INT

On entry, **itype** =  $\langle\text{value}\rangle$ .

*Constraint:* **itype** = 1, 2 or 3.

On entry, **n** =  $\langle value \rangle$ .  
 Constraint: **n**  $\geq 0$ .

On entry, **pda** =  $\langle value \rangle$ .  
 Constraint: **pda** > 0.

On entry, **pdb** =  $\langle value \rangle$ .  
 Constraint: **pdb** > 0.

## NE\_INT\_2

On entry, **pda** =  $\langle value \rangle$  and **n** =  $\langle value \rangle$ .  
 Constraint: **pda**  $\geq \max(1, n)$ .

On entry, **pdb** =  $\langle value \rangle$  and **n** =  $\langle value \rangle$ .  
 Constraint: **pdb**  $\geq \max(1, n)$ .

## 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 contact NAG for assistance.

## NE\_MAT\_NOT\_POS\_DEF

If **fail.errnum** = **n** +  $\langle value \rangle$ , for  $1 \leq \langle value \rangle \leq \mathbf{n}$ , then the leading minor of order  $\langle value \rangle$  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 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 Parallelism and Performance

`nag_zhegvd` (f08sqc) is threaded by NAG for parallel execution in multithreaded implementations of the NAG Library.

`nag_zhegvd` (f08sqc) 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 Users' Note for your implementation for any additional implementation-specific information.

## 9 Further Comments

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

The real analogue of this function is `nag_dsygvd` (f08scc).

## 10 Example

This example finds all the eigenvalues and eigenvectors of the generalized Hermitian eigenproblem  $ABz = \lambda z$ , where

$$A = \begin{pmatrix} -7.36 & 0.77 - 0.43i & -0.64 - 0.92i & 3.01 - 6.97i \\ 0.77 + 0.43i & 3.49 & 2.19 + 4.45i & 1.90 + 3.73i \\ -0.64 + 0.92i & 2.19 - 4.45i & 0.12 & 2.88 - 3.17i \\ 3.01 + 6.97i & 1.90 - 3.73i & 2.88 + 3.17i & -2.54 \end{pmatrix}$$

and

$$B = \begin{pmatrix} 3.23 & 1.51 - 1.92i & 1.90 + 0.84i & 0.42 + 2.50i \\ 1.51 + 1.92i & 3.58 & -0.23 + 1.11i & -1.18 + 1.37i \\ 1.90 - 0.84i & -0.23 - 1.11i & 4.09 & 2.33 - 0.14i \\ 0.42 - 2.50i & -1.18 - 1.37i & 2.33 + 0.14i & 4.29 \end{pmatrix},$$

together with an estimate of the condition number of  $B$ , and approximate error bounds for the computed eigenvalues and eigenvectors.

The example program for nag\_zhegv (f08snc) illustrates solving a generalized Hermitian eigenproblem of the form  $Az = \lambda Bz$ .

### 10.1 Program Text

```
/* nag_zhegvd (f08sdc) Example Program.
*
* Copyright 2011 Numerical Algorithms Group.
*
* Mark 23, 2011.
*/
#include <stdio.h>
#include <nag.h>
#include <nagx04.h>
#include <nag_stdlib.h>
#include <nagf07.h>
#include <nagf08.h>
#include <nagf16.h>
#include <nagx02.h>
#include <naga02.h>

int main(void)
{
    /* Scalars */
    double      anorm, bnorm, eps, rcond, rcondb, t1, t2, t3;
    Integer     i, j, n, pda, pdb;
    Integer     exit_status = 0;
    /* Arrays */
    Complex    *a = 0, *b = 0;
    double      *eerbnd = 0, *rcondz = 0, *w = 0, *zerbnd = 0;
    char        nag_enum_arg[40];

    /* Nag Types */
    NagError    fail;
    Nag_OrderType order;
    Nag_UptoType uplo;

#ifdef NAG_COLUMN_MAJOR
#define A(I, J) a[(J-1)*pda + I - 1]
#define B(I, J) b[(J-1)*pdb + I - 1]
    order = Nag_ColMajor;
#else
#define A(I, J) a[(I-1)*pda + J - 1]
#define B(I, J) b[(I-1)*pdb + J - 1]
    order = Nag_RowMajor;
#endif
}
```

```

INIT_FAIL(fail);

printf("nag_zhegvd (f08sqc) Example Program Results\n\n");

/* Skip heading in data file */
scanf("%*[^\n]");
scanf("%ld%*[^\n]", &n);
if (n < 0)
{
    printf("Invalid n\n");
    exit_status = 1;
    goto END;;
}
scanf(" %39s%*[^\n]", nag_enum_arg);
/* nag_enum_name_to_value (x04nac).
 * Converts NAG enum member name to value
 */
uplo = (Nag_UptoType) nag_enum_name_to_value(nag_enum_arg);

pda = n;
pdb = n;
/* Allocate memory */
if (!(a      = NAG_ALLOC(n * n, Complex)) ||
    !(b      = NAG_ALLOC(n * n, Complex)) ||
    !(eerbnd = NAG_ALLOC(n, double)) ||
    !(rcondz = NAG_ALLOC(n, double)) ||
    !(w      = NAG_ALLOC(n, double)) ||
    !(zerbnd = NAG_ALLOC(n, double)))
{
    printf("Allocation failure\n");
    exit_status = -1;
    goto END;
}

/* Read the triangular parts of the matrices A and B */
if (uplo == Nag_Upper)
{
    for (i = 1; i <= n; ++i)
        for (j = i; j <= n; ++j)
            scanf("( %lf , %lf ) ", &A(i, j).re, &A(i, j).im);
    scanf("%*[^\n]");
    for (i = 1; i <= n; ++i)
        for (j = i; j <= n; ++j)
            scanf("( %lf , %lf ) ", &B(i, j).re, &B(i, j).im);
}
else
{
    for (i = 1; i <= n; ++i)
        for (j = 1; j <= i; ++j)
            scanf("( %lf , %lf ) ", &A(i, j).re, &A(i, j).im);
    scanf("%*[^\n]");
    for (i = 1; i <= n; ++i)
        for (j = 1; j <= i; ++j)
            scanf("( %lf , %lf ) ", &B(i, j).re, &B(i, j).im);
}
scanf("%*[^\n]");

/* Compute the one-norms of the symmetric matrices A and B
 * using nag_zhe_norm (f16ucc).
 */
nag_zhe_norm(order, Nag_OneNorm, uplo, n, a, pda, &anorm, &fail);
nag_zhe_norm(order, Nag_OneNorm, uplo, n, b, pdb, &bnorm, &fail);
if (fail.code != NE_NOERROR)
{
    printf("Error from nag_zhe_norm (f16ucc).\n%s\n", fail.message);
    exit_status = 1;
    goto END;
}

/* Solve the generalized Hermitian eigenvalue problem A*B*x = lambda*x

```

```

* using nag_zhegvd (f08sqc).
*/
nag_zhegvd(order, 2, Nag_DoBoth, uplo, n, a, pda, b, pdb, w, &fail);
if (fail.code != NE_NOERROR)
{
    printf("Error from nag_zhegvd (f08sqc).\\n%s\\n", fail.message);
    exit_status = 1;
    goto END;
}

/* Normalize the eigenvectors */
for(j=1; j<=n; j++)
    for(i=n; i>=1; i--) A(i, j) = nag_complex_divide(A(i, j),A(1, j));

/* Print eigensolution */
printf(" Eigenvalues\\n   ");
for (j = 0; j < n; ++j) printf(" %11.4f%s", w[j], j%6 == 5?"\\n":"");
printf("\\n");

/* Prnit normalized vectors using nag_gen_complx_mat_print (x04dac). */
fflush(stdout);
nag_gen_complx_mat_print(order, Nag_GeneralMatrix, Nag_NonUnitDiag, n, n,
                           a, pda, "Eigenvectors", 0, &fail);
if (fail.code != NE_NOERROR)
{
    printf("Error from nag_gen_complx_mat_print (x04dac).\\n%s\\n",
           fail.message);
    exit_status = 1;
    goto END;
}

/* Estimate the reciprocal condition number of the Cholesky factor of B.
 * to estimate the reciprocal condition.
 * nag_ztrcon (f07tuc)
 * Note that: cond(B) = 1/rcond**2
 */
nag_ztrcon(order, Nag_OneNorm, uplo, Nag_NonUnitDiag, n, b, pdb,
            &rcond, &fail);
if (fail.code != NE_NOERROR)
{
    printf("Error from nag_ztrcon (f07tuc).\\n%s\\n", fail.message);
    exit_status = 1;
    goto END;
}

/* Print the reciprocal condition number of B */
rcondb = rcond * rcond;

printf("\\nEstimate of reciprocal condition number for B\\n      %11.1e\\n",
       rcondb);

/* Get the machine precision, using nag_machine_precision (x02ajc) */
eps = nag_machine_precision;
if (rcond < eps)
{
    printf("\\nB is very ill-conditioned, error estimates have not been "
           "computed\\n");
    goto END;
}

/* Estimate reciprocal condition numbers for the eigenvectors of A*B-lambda*I
 * nag_ddisna (f08flc)
 */
nag_ddisna(Nag_EigVecs, n, n, w, rcondz, &fail);
if (fail.code != NE_NOERROR)
{
    printf("Error from nag_ddisna (f08flc).\\n%s\\n", fail.message);
    exit_status = 1;
    goto END;
}

```

```

/* Compute the error estimates for the eigenvalues and eigenvectors. */
t1 = 1.0 / rcond;
t2 = eps * t1;
t3 = anorm * bnorm;
for (i = 0; i < n; ++i)
{
    eerbnd[i] = eps * (t3 + abs(w[i])/rcondb);
    zerbnd[i] = t2 * (t3/rcondz[i] + t1);
}
/* Print the approximate error bounds for the eigenvalues and vectors. */
printf("\nError estimates for the eigenvalues\n");
for (i = 0; i < n; ++i) printf(" %11.1e%s", eerbnd[i], i%6 == 5?"\n":"");
printf("\n\nError estimates for the eigenvectors\n");
for (i = 0; i < n; ++i) printf(" %11.1e%s", zerbnd[i], i%6 == 5?"\n":"");
printf("\n");

END:
NAG_FREE(a);
NAG_FREE(b);
NAG_FREE(eerbnd);
NAG_FREE(rcondz);
NAG_FREE(w);
NAG_FREE(zerbnd);

return exit_status;
}

```

## 10.2 Program Data

nag\_zhegvd (f08sqc) Example Program Data

```

4 : n

Nag_Upper : uplo

(-7.36, 0.00) ( 0.77, -0.43) (-0.64, -0.92) ( 3.01, -6.97)
( 3.49, 0.00) ( 2.19, 4.45) ( 1.90, 3.73)
( 0.12, 0.00) ( 2.88, -3.17)
(-2.54, 0.00) : matrix A

( 3.23, 0.00) ( 1.51, -1.92) ( 1.90, 0.84) ( 0.42, 2.50)
( 3.58, 0.00) (-0.23, 1.11) (-1.18, 1.37)
( 4.09, 0.00) ( 2.33, -0.14)
( 4.29, 0.00) : matrix B

```

## 10.3 Program Results

nag\_zhegvd (f08sqc) Example Program Results

|              |         |         |         |         |
|--------------|---------|---------|---------|---------|
| Eigenvalues  |         |         |         |         |
| -61.7321     | -6.6195 | 0.0725  | 43.1883 |         |
| Eigenvectors |         |         |         |         |
|              | 1       | 2       | 3       |         |
| 1            | 1.0000  | 1.0000  | 1.0000  | 1.0000  |
|              | 0.0000  | -0.0000 | -0.0000 | 0.0000  |
| 2            | -0.4648 | 1.5001  | -0.2201 | -1.7750 |
|              | 0.0291  | 1.9518  | -0.3108 | 0.6305  |
| 3            | 0.1123  | -3.2220 | -0.5544 | 0.0399  |
|              | 0.0866  | 0.8343  | -0.1985 | 1.0638  |
| 4            | -0.5690 | 1.0723  | 0.2491  | -1.2264 |
|              | -0.5820 | 0.9277  | 0.5732  | 0.7828  |

Estimate of reciprocal condition number for B  
2.5e-03

```
Error estimates for the eigenvalues
  2.7e-12    2.8e-13    2.3e-14    1.9e-12

Error estimates for the eigenvectors
  5.2e-14    1.1e-13    1.1e-13    5.4e-14
```

---