

## NAG Library Function Document

### nag\_nearest\_correlation\_bounded (g02abc)

#### 1 Purpose

nag\_nearest\_correlation\_bounded (g02abc) computes the nearest correlation matrix, in the Frobenius norm or weighted Frobenius norm, and optionally with bounds on the eigenvalues, to a given square, input matrix.

#### 2 Specification

```
#include <nag.h>
#include <nagg02.h>

void nag_nearest_correlation_bounded (Nag_OrderType order, double g[],
    Integer pdg, Integer n, Nag_NearCorr_ProbType opt, double alpha,
    double w[], double errtol, Integer maxits, Integer maxit, double x[],
    Integer pdx, Integer *iter, Integer *feval, double *nrmgrd,
    NagError *fail)
```

#### 3 Description

Finds the nearest correlation matrix  $X$  by minimizing  $\frac{1}{2}\|G - X\|^2$  where  $G$  is an approximate correlation matrix.

The norm can either be the Frobenius norm or the weighted Frobenius norm  $\frac{1}{2}\|W^{\frac{1}{2}}(G - X)W^{\frac{1}{2}}\|_F^2$ .

You can optionally specify a lower bound on the eigenvalues,  $\alpha$ , of the computed correlation matrix, forcing the matrix to be positive definite,  $0 < \alpha < 1$ .

Note that if the weights vary by several orders of magnitude from one another the algorithm may fail to converge.

#### 4 References

Borsdorf R and Higham N J (2010) A preconditioned (Newton) algorithm for the nearest correlation matrix *IMA Journal of Numerical Analysis* **30(1)** 94–107

Qi H and Sun D (2006) A quadratically convergent Newton method for computing the nearest correlation matrix *SIAM J. Matrix AnalAppl* **29(2)** 360–385

#### 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: **g[*pdg* × *n*]** – double *Input/Output*

**Note:** the (*i*, *j*)th element of the matrix  $G$  is stored in

**g**[(*j* – 1) × **pdg** + *i* – 1] when **order** = Nag\_ColMajor;  
**g**[(*i* – 1) × **pdg** + *j* – 1] when **order** = Nag\_RowMajor.

- On entry:*  $G$ , the initial matrix.  
*On exit:*  $G$  is overwritten.
- 3: **pdg** – Integer *Input*  
*On entry:* the stride separating row or column elements (depending on the value of **order**) in the array **g**.  
*Constraint:* **pdg**  $\geq$  **n**.
- 4: **n** – Integer *Input*  
*On entry:* the order of the matrix  $G$ .  
*Constraint:* **n**  $>$  0.
- 5: **opt** – Nag\_NearCorr\_ProbType *Input*  
*On entry:* indicates the problem to be solved.  
**opt** = Nag\_LowerBound  
 The lower bound problem is solved.  
**opt** = Nag\_WeightedNorm  
 The weighted norm problem is solved.  
**opt** = Nag\_Both  
 Both problems are solved.  
*Constraint:* **opt** = Nag\_LowerBound, Nag\_WeightedNorm or Nag\_Both.
- 6: **alpha** – double *Input*  
*On entry:* the value of  $\alpha$ .  
 If **opt** = Nag\_WeightedNorm, **alpha** need not be set.  
*Constraint:*  $0.0 < \mathbf{alpha} < 1.0$ .
- 7: **w[n]** – double *Input/Output*  
*On entry:* the square roots of the diagonal elements of  $W$ , that is the diagonal of  $W^{\frac{1}{2}}$ .  
 If **opt** = Nag\_LowerBound, **w** need not be set.  
*On exit:* if **opt** = Nag\_WeightedNorm or Nag\_Both, the array is scaled so  $0 < \mathbf{w}[i-1] \leq 1$ , for  $i = 1, 2, \dots, n$ .  
*Constraint:*  $\mathbf{w}[i-1] > 0.0$ , for  $i = 1, 2, \dots, n$ .
- 8: **errtol** – double *Input*  
*On entry:* the termination tolerance for the Newton iteration. If **errtol**  $\leq$  0.0 then  $\mathbf{n} \times \sqrt{\mathit{machine\ precision}}$  is used.
- 9: **maxits** – Integer *Input*  
*On entry:* specifies the maximum number of iterations to be used by the iterative scheme to solve the linear algebraic equations at each Newton step.  
 If **maxits**  $\leq$  0,  $2 \times \mathbf{n}$  is used.
- 10: **maxit** – Integer *Input*  
*On entry:* specifies the maximum number of Newton iterations.  
 If **maxit**  $\leq$  0, 200 is used.

- 11: **x**[**pdx** × **n**] – double *Output*  
**Note:** the  $(i, j)$ th element of the matrix  $X$  is stored in  
 $\mathbf{x}[(j - 1) \times \mathbf{pdx} + i - 1]$  when **order** = Nag\_ColMajor;  
 $\mathbf{x}[(i - 1) \times \mathbf{pdx} + j - 1]$  when **order** = Nag\_RowMajor.  
*On exit:* contains the nearest correlation matrix.
- 12: **pdx** – Integer *Input*  
*On entry:* the stride separating row or column elements (depending on the value of **order**) in the array **x**.  
*Constraint:* **pdx** ≥ **n**.
- 13: **iter** – Integer \* *Output*  
*On exit:* the number of Newton steps taken.
- 14: **feval** – Integer \* *Output*  
*On exit:* the number of function evaluations of the dual problem.
- 15: **nrmgrd** – double \* *Output*  
*On exit:* the norm of the gradient of the last Newton step.
- 16: **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  $\langle value \rangle$  had an illegal value.

### NE\_CONVERGENCE

Newton iteration fails to converge in  $\langle value \rangle$  iterations. Increase **maxit** or check the call to the function.

The *machine precision* is limiting convergence. In this instance the returned matrix  $X$  may be useful.

### NE\_EIGENPROBLEM

Failure to solve intermediate eigenproblem.

### NE\_INT

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

### NE\_INT\_2

On entry, **pdg** =  $\langle value \rangle$  and **n** =  $\langle value \rangle$ .  
Constraint: **pdg** ≥ **n**.

On entry, **pdx** =  $\langle value \rangle$  and **n** =  $\langle value \rangle$ .  
Constraint: **pdx** ≥ **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\_REAL**

On entry, **alpha** = *value*.  
Constraint:  $0.0 < \mathbf{alpha} < 1.0$ .

**NE\_WEIGHTS\_NOT\_POSITIVE**

On entry, all elements of **w** were not positive.

**7 Accuracy**

The returned accuracy is controlled by **errtol** and limited by *machine precision*.

**8 Parallelism and Performance**

nag\_nearest\_correlation\_bounded (g02abc) is threaded by NAG for parallel execution in multithreaded implementations of the NAG Library.

nag\_nearest\_correlation\_bounded (g02abc) 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**

Arrays are internally allocated by nag\_nearest\_correlation\_bounded (g02abc). The total size of these arrays is  $12 \times \mathbf{n} + 3 \times \mathbf{n} \times \mathbf{n} + \max(2 \times \mathbf{n} \times \mathbf{n} + 6 \times \mathbf{n} + 1, 120 + 9 \times \mathbf{n})$  double elements and  $5 \times \mathbf{n} + 3$  Integer elements. All allocated memory is freed before return of nag\_nearest\_correlation\_bounded (g02abc).

**10 Example**

This example finds the nearest correlation matrix to:

$$G = \begin{pmatrix} 2 & -1 & 0 & 0 \\ -1 & 2 & -1 & 0 \\ 0 & -1 & 2 & -1 \\ 0 & 0 & -1 & 2 \end{pmatrix}$$

weighted by  $W^{\frac{1}{2}} = \text{diag}(100, 20, 20, 20)$  with minimum eigenvalue 0.02.

**10.1 Program Text**

```
/* nag_nearest_correlation_bounded (g02abc) Example Program.
 *
 * Copyright 2011, Numerical Algorithms Group.
 *
 * Mark 23, 2011.
 */

#include <nag.h>
#include <nag_stdlib.h>
#include <nagf08.h>
#include <nagg02.h>
#include <nagx04.h>
```

```

int main(void)
{
    /* Scalars */
    Integer          exit_status = 0;
    double           alpha, errtol, nrmgrd;
    Integer          feval, i, iter, j, maxit, maxits, n, pdeig, pdg, pdx;

    /* Arrays */
    char             nag_enum_arg[100];
    double           *eig = 0, *g = 0, *w = 0, *x = 0;

    /* Nag Types */
    Nag_OrderType   order;
    Nag_NearCorr_ProbType opt;
    NagError        fail;

    INIT_FAIL(fail);

#ifdef NAG_COLUMN_MAJOR
#define G(I, J) g[(J-1)*pdg + I-1]
#define X(I, J) x[(J-1)*pdx + I-1]
    order = Nag_ColMajor;
#else
#define G(I, J) g[(I-1)*pdg + J-1]
#define X(I, J) x[(I-1)*pdx + J-1]
    order = Nag_RowMajor;
#endif

    /* Output preamble */
    printf("nag_nearest_correlation_bounded (g02abc)");
    printf(" Example Program Results\n\n");
    fflush(stdout);

    /* Skip heading in data file */
    scanf("%*[\n] ");
    /* Read in the problem size, opt and alpha */
    scanf("%ld", &n);
    scanf("%39s", nag_enum_arg);
    /*
     * nag_enum_name_to_value (x04nac).
     * Converts NAG enum member name to value
     */
    opt = (Nag_NearCorr_ProbType) nag_enum_name_to_value(nag_enum_arg);
    scanf("%lf%*[\n]", &alpha);

    pdg = n;
    pdx = n;
    if( order == Nag_ColMajor)
        pdeig = 1;
    else
        pdeig = n;

    if (
        !(g = NAG_ALLOC((pdg)*(n), double)) ||
        !(w = NAG_ALLOC((n), double)) ||
        !(x = NAG_ALLOC((pdx)*(n), double)) ||
        !(eig = NAG_ALLOC((n), double))
        )
    {
        printf("Allocation failure\n");
        exit_status = -1;
        goto END;
    }

    /* Read in the matrix g */
    for (i = 1; i <= n; i++)
        for (j = 1; j <= n; j++)
            scanf("%lf", &G(i, j));
    scanf("%*[\n] ");

```

```

/* Read in the vector w */
for (i = 0; i < n; i++)
    scanf("%lf", &w[i]);
scanf("%*[\n] ");

/* Use the defaults for errtol, maxits and maxit */
errtol = 0.0;
maxits = 0;
maxit = 0;

/*
 * nag_nearest_correlation_bounded (g02abc).
 * Computes the nearest correlation matrix incorporating weights
 * and/or bounds
 */
nag_nearest_correlation_bounded(order, g, pdg, n, opt, alpha, w, errtol,
                                maxits, maxit, x, pdx, &iter, &feval,
                                &nrmgrd, &fail);

if (fail.code != NE_NOERROR)
{
    printf("%s\n", fail.message);
    exit_status = 1;
    goto END;
}

/*
 * nag_gen_real_mat_print (x04cac).
 * Print real general matrix (easy-to-use)
 */
nag_gen_real_mat_print(order, Nag_GeneralMatrix, Nag_NonUnitDiag, n, n, x,
                        pdx, "Nearest Correlation Matrix x", NULL, &fail);
if (fail.code != NE_NOERROR)
{
    printf("%s\n", fail.message);
    exit_status = 1;
    goto END;
}

printf("\nNumber of Newton steps taken: %lld\n", iter);
printf("Number of function evaluations: %9ld\n\n", feval);
printf("alpha: %37.3f \n\n", alpha);

/* nag_dsyev (f08fac).
 * Computes all eigenvalues and, optionally, eigenvectors of a real
 * symmetric matrix
 */
nag_dsyev(order, Nag_EigVals, Nag_Upper, n, x, pdx, eig, &fail);
if (fail.code != NE_NOERROR)
{
    printf("%s\n", fail.message);
    exit_status = 1;
    goto END;
}

/* nag_gen_real_mat_print (x04cac).
 * Print real general matrix (easy-to-use)
 */
nag_gen_real_mat_print(order, Nag_GeneralMatrix, Nag_NonUnitDiag, 1, n,
                        eig, pdeig, "Eigenvalues of x", NULL, &fail);
if (fail.code != NE_NOERROR)
{
    printf("%s\n", fail.message);
    exit_status = 1;
}

END:
NAG_FREE(eig);
NAG_FREE(g);
NAG_FREE(w);
NAG_FREE(x);
return exit_status;
}

```

**10.2 Program Data**

```
nag_nearest_correlation_bounded (g02abc) Example Program Data
4 Nag_Both 0.02      :: n, opt, alpha
  2.0   -1.0   0.0   0.0
 -1.0   2.0   -1.0   0.0
  0.0   -1.0   2.0  -1.0
  0.0   0.0   -1.0   2.0  :: End of g
100.0  20.0  20.0  20.0  :: w
```

**10.3 Program Results**

```
nag_nearest_correlation_bounded (g02abc) Example Program Results
```

```
Nearest Correlation Matrix x
      1      2      3      4
1  1.0000 -0.9187  0.0257  0.0086
2 -0.9187  1.0000 -0.3008  0.2270
3  0.0257 -0.3008  1.0000 -0.8859
4  0.0086  0.2270 -0.8859  1.0000
```

```
Number of Newton steps taken:      5
Number of function evaluations:    6
```

```
alpha:                             0.020
```

```
Eigenvalues of x
      1      2      3      4
1  0.0392  0.1183  1.6515  2.1910
```

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