NAG Fortran Library Routine Document

G05GBF

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

G05GBF generates a random correlation matrix with given eigenvalues.

2 Specification

```
SUBROUTINE G05GBF(N, D, C, LDC, EPS, WK, IFAIL)
INTEGER           N, LDC, IFAIL
real              D(N), C(LDC,N), EPS, WK(2*N)
```

3 Description

Given n eigenvalues, $\lambda_1, \lambda_2, \ldots, \lambda_n$, such that

$$\sum_{i=1}^{n} \lambda_i = n$$

and $\lambda_i \geq 0$, for $i = 1, 2, \ldots, n$.

G05GBF will generate a random correlation matrix, C, of dimension n, with eigenvalues $\lambda_1, \lambda_2, \ldots, \lambda_n$.

The method used is based on that described by Lin and Bendel (1985). Let D be the diagonal matrix with values $\lambda_1, \lambda_2, \ldots, \lambda_n$ and let A be a random orthogonal matrix generated by G05GAF then the matrix $C_0 = ADA^T$ is a random covariance matrix with eigenvalues $\lambda_1, \lambda_2, \ldots, \lambda_n$. The matrix C_0 is transformed into a correlation matrix by means of $n-1$ elementary rotation matrices P_i such that $C = P_{n-1}P_{n-2}\ldots P_1 C_0 P_1^T \ldots P_{n-2}^T P_{n-1}^T$. The restriction on the sum of eigenvalues implies that for any diagonal element of $C_0 > 1$, there is another diagonal element < 1. The P_i are constructed from such pairs, chosen at random, to produce a unit diagonal element corresponding to the first element. This is repeated until all diagonal elements are 1 to within a given tolerance ϵ.

The randomness of C should be interpreted only to the extent that A is a random orthogonal matrix and C is computed from A using the P_i which are chosen as arbitrarily as possible.

4 References

Lin S P and Bendel R B (1985) Algorithm AS213: Generation of population correlation on matrices with specified eigenvalues *Appl. Statist.* **34** 193–198

5 Parameters

1: N – INTEGER *Input*

On entry: the dimension of the correlation matrix to be generated, n.

Constraint: N ≥ 1.

2: D(N) – ***real*** array *Input*

On entry: the n eigenvalues, λ_i, for $i = 1, 2, \ldots, n$.

Constraints:

$$\mathrm{D}(i) \geq 0.0, \text{ for } i = 1, 2, \ldots, n, \text{ and } \sum_{i=1}^{n} \mathrm{D}(i) = n \text{ to within EPS.}$$

3: C(LDC,N) – ***real*** array *Output*

On exit: a random correlation matrix, C, of dimension n.

4: LDC – INTEGER *Input*

On entry: the first dimension of the array C as declared in the (sub)program from which G05GBF is called.

Constraint: LDC $\geq$ N.

5: EPS – ***real*** *Input*

On entry: the maximum acceptable error in the diagonal elements, ϵ.

Constraint: EPS $\geq$ N$\times$***machine precision***.

Suggested value: EPS=0.00001.

6: WK(2$*$N) – ***real*** array *Workspace*

7: IFAIL – INTEGER *Input/Output*

On entry: IFAIL must be set to 0, -1 or 1. Users who are unfamiliar with this parameter should refer to Chapter P01 for details.

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

For environments where it might be inappropriate to halt program execution when an error is detected, the value -1 or 1 is recommended. If the output of error messages is undesirable, then the value 1 is recommended. Otherwise, for users not familiar with this parameter the recommended value is 0. **When the value -1 or 1 is used it is essential to test the value of IFAIL on exit.**

6 Error Indicators and Warnings

If on entry IFAIL $= 0$ or -1, explanatory error messages are output on the current error message unit (as defined by X04AAF).

Errors or warnings detected by the routine:

IFAIL $= 1$

On entry, N < 0,
or LDC $<$ N,
or EPS $<$ N$\times$ ***machine precision***.

IFAIL $= 2$

On entry, D$(i) < 0.0$ for some i,
or $\displaystyle\sum_{i=1}^{n} \mathrm{D}(i) \neq n$ to within EPS.

IFAIL $= 3$

The error in a diagonal element is greater than EPS. The value of EPS should be increased. Otherwise the program could be re-run with a different value used for the seed of the random number generator, see G05CBF or G05CCF.

7 Accuracy

The maximum error in a diagonal element is given by EPS.

8 Further Comments

The time taken by the routine is approximately proportional to n^2.

9 Example

A 3 by 3 correlation matrix with eigenvalues of 0.7, 0.9 and 1.4 is generated and printed.

The generator mechanism used is selected by an initial call to G05ZAF.

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.

```
*     G05GBF Example Program Text
*     Mark 20 Revised. NAG Copyright 2001.
*     .. Parameters ..
      INTEGER          NIN, NOUT
      PARAMETER        (NIN=5,NOUT=6)
      INTEGER          NMAX
      PARAMETER        (NMAX=10)
*     .. Local Scalars ..
      real             EPS
      INTEGER          I, IFAIL, J, LDC, N
*     .. Local Arrays ..
      real             C(NMAX,NMAX), D(NMAX), WK(2*NMAX)
*     .. External Subroutines ..
      EXTERNAL         G05CBF, G05GBF, G05ZAF
*     .. Executable Statements ..
      CALL G05ZAF('O')
      WRITE (NOUT,*) 'G05GBF Example Program Results'
*     Skip heading in data file
      READ (NIN,*)
      READ (NIN,*) N
      IF (N.LE.NMAX) THEN
         READ (NIN,*) (D(I),I=1,N)
*
         WRITE (NOUT,*)
*
         LDC = NMAX
         CALL G05CBF(0)
         EPS = 0.0001e0
*
         IFAIL = 0
*
         CALL G05GBF(N,D,C,LDC,EPS,WK,IFAIL)
*
         DO 20 I = 1, N
            WRITE (NOUT,99999) (C(I,J),J=1,N)
   20    CONTINUE
      END IF
      STOP
*
99999 FORMAT (1X,3F9.3)
      END
```

9.2 Program Data

```
G05GBF Example Program Data
3
0.7 0.9 1.4
```

9.3 Program Results

```
 G05GBF Example Program Results

     1.000    0.100   -0.251
     0.100    1.000   -0.239
    -0.251   -0.239    1.000
```