NAG C Library Function Document

nag_anova_row_col (g04bcc)

1 Purpose

nag_anova_row_col (g04bcc) computes the analysis of variance for a general row and column design together with the treatment means and standard errors.

2 Specification

3 Description

In a row and column design the experimental material can be characterized by a two-way classification, nominally called rows and columns. Each experimental unit can be considered as being located in a particular row and column. It is assumed that all rows are of the same length and all columns are of the same length. Sets of equal numbers of rows and columns can be grouped together to form replicates, sometimes known as squares or rectangles, as appropriate.

If for a replicate, the number of rows, the number of columns and the number of treatments are equal and every treatment occurs once in each row and each column then the design is a Latin square. If this is not the case the treatments will be non-orthogonal to rows and columns. For example in the case of a lattice square each treatment occurs only once in each square.

For a row and column design, with t treatments in r rows and c columns and b replicates or squares with n = brc observations, the linear model is:

$$y_{ijk(l)} = \mu + \beta_i + \rho_j + \gamma_k + \tau_l + e_{ijk}$$

i = 1, 2, ..., b; j = 1, 2, ..., r; k = 1, 2, ..., c; l = 1, 2, ..., t, where β_i is the effect of the *i*th replicate, ρ_j is the effect of the *j*th row, γ_k is the effect of the *k*th column and the ijk(l) notation indicates that the *l*th treatment is applied to the unit in row *j*, column *k* of replicate *i*.

To compute the analysis of variance for a row and column design the mean is computed and subtracted from the observations to give, $y'_{ijk(l)} = y_{ijk(l)} - \hat{\mu}$. Since the replicates, rows and columns are orthogonal the estimated effects, ignoring treatment effects, $\hat{\beta}_i$, $\hat{\rho}_j$, $\hat{\gamma}_k$, can be computed using the appropriate means of the $y'_{ijk(l)}$, and the unadjusted sum of squares computed as the appropriate sum of squared totals for the $y'_{ijk(l)}$ divided by number of units per total. The observations adjusted for replicates, rows and columns can then be computed by subtracting the estimated effects from $y'_{ijk(l)}$ to give $y''_{ijk(l)}$.

In the case of a Latin square design the treatments are orthogonal to replicates, rows and columns and so the treatment effects, $\hat{\tau}_l$, can be estimated as the treatment means of the adjusted observations, $y''_{ijk(l)}$. The treatment sum of squares is computed as the sum of squared treatment totals of the $y''_{ij(l)}$ divided by the number of times each treatment is replicated. Finally the residuals, and hence the residual sum of squares, are given by, $r_{ij(l)} = y''_{ij(l)} - \hat{\tau}_l$.

For a design which is not orthogonal, for example a lattice square or an incomplete Latin square, the treatment effects adjusted for replicates, rows and columns need to be computed. The adjusted treatment effects are found as the solution to the equations:

$$A\hat{\tau} = (R - N_b N_b^T / (rc) - N_r N_r^T / (bc) - N_c N_c^T / (br))\hat{\tau} = q$$

where q is the vector of the treatment totals of the observations adjusted for replicates, rows and columns, $y''_{ijk(l)}$; R is a diagonal matrix with R_{ll} equal to the number of times the *l*th treatment is replicated, and N_b is the *t* by *b* incidence matrix, with $N_{l,i}$ equal to the number of times treatment *l* occurs in replicate *i*, with N_r and N_c being similarly defined for rows and columns. The solution to the equations can be written as:

$$\hat{\tau} = \Omega q$$

where, Ω is a generalized inverse of A. The solution is found from the eigenvalue decomposition of A. The residuals are first calculated by subtracting the estimated adjusted treatment effects from the adjusted observations to give $r'_{ij(l)} = y''_{ij(l)} - \hat{\tau}_l$. However, since only the unadjusted replicate, row and column effects have been removed and they are not orthogonal to treatments, the replicate, row and column means of the $r'_{ij(l)}$ have to be subtracted to give the correct residuals, $r_{ij(l)}$ and residual sum of squares.

Given the sums of squares, the mean squares are computed as the sums of squares divided by the degrees of freedom. The degrees of freedom for the unadjusted replicates, rows and columns are b-1, r-1 and c-1 respectively and for the Latin square designs the degrees of freedom for the treatments is t-1. In the general case the degrees of freedom for treatments is the rank of the matrix Ω . The *F*-statistic given by the ratio of the treatment mean square to the residual mean square tests the hypothesis:

$$H_0: \tau_1 = \tau_2 = \ldots = \tau_t = 0.$$

The standard errors for the difference in treatment effects, or treatment means, for Latin square designs, are given by:

$$se(\hat{\tau}_j - \hat{\tau}_{j*}) = \sqrt{2s^2/(bt)}$$

where s^2 is the residual mean square. In the general case the variances of the treatment effects are given by:

$$\operatorname{Var}(\hat{\tau}) = \Omega s^2$$

from which the appropriate standard errors of the difference between treatment effects or the difference between adjusted means can be calculated.

The analysis of a row-column design can be considered as consisting of different strata: the replicate stratum, the rows within replicate and the columns within replicate strata and the units stratum. In the Latin square design all the information on the treatment effects is given at the units stratum. In other designs there may be a loss of information due to the non-orthogonality of treatments and replicates, rows and columns and information on treatments may be available in higher strata. The efficiency of the estimation at the units stratum is given by the (canonical) efficiency factors, these are the non-zero eigenvalues of the matrix, A, divided by the number of replicates in the case of equal replication, or by the mean of the number of replicates in the unequally replicated case, (see John (1987)). If more than one eigenvalue is zero then the design is said to be disconnected and information on some treatment comparisons can only be obtained from higher strata.

4 Parameters

1:	nrep – Integer	Input
	On entry: the number of replicates, b.	
	Constraint: $nrep \ge 1$.	
2:	nrow – Integer	Input
	On entry: the number of rows per replicate, r.	
	Constraint: $nrow \ge 2$.	
3:	ncol – Integer	Input
	On entry: the number of columns per replicate, c.	
	Constraint: $ncol \ge 2$.	

4: **y[nrep*nrow*ncol]** – const double

On entry: the n = brc observations ordered by columns within rows within replicates. That is $\mathbf{y}[rc(i-1) + r(j-1) + k - 1]$ contains the observation from the k column of the jth row of the *i*th replicate, i = 1, 2, ..., b; j = 1, 2, ..., r; k = 1, 2, ..., c.

5: **nt** – Integer

On entry: the number of treatments. If only replicates, rows and columns are required in the analysis then set $\mathbf{nt} = 1$.

Constraint: $\mathbf{nt} \geq 1$.

6: **it**[*dim1*] – const Integer

Note: the dimension, dim1, of the array it must be at least nrep*nrow*ncol if nt > 1, and 1 otherwise.

On entry: if nt > 1, it[i-1] indicates which of the **nt** treatments unit *i* received, i = 1, 2, ..., n. If nt = 1, it is not referenced.

Constraint: if $\mathbf{nt} \geq 2$, $1 \leq \mathbf{it}[i-1] \leq \mathbf{nt}$, for $i = 1, 2, \ldots, n$.

7: **gmean** – double *

On exit: the grand mean, $\hat{\mu}$.

8: **tmean[nt]** – double

On exit: if $\mathbf{nt} \geq 2$, tmean[l-1] contains the (adjusted) mean for the *l*th treatment, $\hat{\mu}^* + \hat{\tau}_l$, l = 1, 2, ..., t, where $\hat{\mu}^*$ is the mean of the treatment adjusted observations $y_{ijk(l)} - \hat{\tau}_l$. Otherwise tmean is not referenced.

9: **table[6][5]** – double

On exit: the analysis of variance table. Column 1 contains the degrees of freedom, column 2 the sum of squares, and where appropriate, column 3 the mean squares, column 4 the *F*-statistic and column 5 the significance level of the *F*-statistic. Row 1 is for replicates, row 2 for rows, row 3 for columns, row 4 for treatments (if nt > 1), row 5 for residual and row 6 for total. Mean squares are computed for all but the total row, *F*-statistics and significance are computed for treatments, replicates, rows and columns. Any unfilled cells are set to zero.

10: c[nt][tdc] - double

On exit: the upper triangular part of **c** contains the variance-covariance matrix of the treatment effects, the strictly lower triangular part contains the standard errors of the difference between two treatment effects (means), i.e., $\mathbf{c}[i-1][j-1]$ contains the covariance of treatment *i* and *j* if $j \ge i$ and the standard error of the difference between treatment *i* and *j* if j < i, i = 1, 2, ..., t; j = 1, 2, ..., t.

11: **tdc** – Integer

On entry: the second dimension of the array \mathbf{c} as declared in the function from which nag_anova_row_col is called.

Constraint: $tdc \ge nt$.

12: **irep[nt]** – Integer

On exit: if nt > 1, irep[l-1] contains the treatment replications, R_{ll} , l = 1, 2, ..., nt. Otherwise irep is not referenced.

13: **rpmean[nrep]** – double

On exit: if nrep > 1, rpmean[i-1] contains the mean for the *i*th replicate, $\hat{\mu} + \hat{\beta}_i$, i = 1, 2, ..., b. Otherwise rpmean is not referenced.

Input

Input

Output

Output

Output

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Input

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Output

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14:	rmean[nrep*nrow] – double Output On exit: rmean $[j-1]$ contains the mean for the <i>j</i> th row, $\hat{\mu} + \hat{\rho}_i$, $j = 1, 2,, r$.	
15:	cmean[nrep*ncol] – double <i>On exit:</i> cmean [$k - 1$] contains the mean for the <i>k</i> th column, $\hat{\mu} + \hat{\gamma}_k$, $k = 1, 2,, c$.	
16:	$\mathbf{r}[\mathbf{nrep*nrow*ncol}]$ – double Output On exit: $\mathbf{r}[i-1]$ contains the residuals, r_i , $i = 1, 2,, n$.	
17:	$ef[nt]$ - doubleOutputOn exit: if $nt \ge 2$, the canonical efficiency factors. Otherwise ef is not referenced.	
18:	tol – double Input On entry: the tolerance value used to check for zero eigenvalues of the matrix Ω . If tol = 0.0 a default value of 0.00001 is used. Constraint: tol ≥ 0.0	
19:	irdf – IntegerInputOn entry: an adjustment to the degrees of freedom for the residual and total. If $irdf \ge 1$ the degrees	
	of freedom for the total is set to $n - irdf$ and the residual degrees of freedom adjusted accordingly. If $irdf = 0$, the total degrees of freedom for the total is set to $n - 1$, as usual. Constraint: $irdf \ge 0$.	
20:	fail - NagError *Input/OutputThe NAG error parameter (see the Essential Introduction).	
5	Error Indicators and Warnings	
NE_	INT_ARG_LT	
	On entry, nrep must not be less than 1: $nrep = \langle value \rangle$.	
	On entry, nrow must not be less than 2: $nrow = \langle value \rangle$.	
	On entry, ncol must not be less than 2: $ncol = \langle value \rangle$.	
	On entry, nt must not be less than 1: $\mathbf{nt} = \langle value \rangle$.	
	On entry, irdf must not be less than 0: irdf = $\langle value \rangle$.	

NE_2_INT_ARG_LT

On entry, $\mathbf{tdc} = \langle value \rangle$ while $\mathbf{nt} = \langle value \rangle$. These parameters must satisfy $\mathbf{tdc} \geq \mathbf{nt}$.

NE_REAL_ARG_LT

On entry, tol must not be less than 0.0: tol = $\langle value \rangle$.

NE_ARRAY_CONS

The contents of array it are not valid. Constraint: if $\mathbf{nt} \ge 2$, $1 \le \mathbf{it}[i] \le \mathbf{nt}$, $i = 0, 1, 2, \dots, \mathbf{nrep} \ast \mathbf{nrow} \ast \mathbf{ncol}$.

The contents of array **it** are not valid. Constraint: some value of $\mathbf{it} = j$ for all $j = 1, 2, ..., \mathbf{nt}$.

NE_ALLOC_FAIL

Memory allocation failed.

NE_ARRAY_CONSTANT

On entry, the elements of the array y are constant.

NE_G04BC_ST_ERR

A computed standard error is zero due to rounding errors, or the eigenvalue computation failed to converge. Both are unlikely errors.

NE_G04BC_REPS

The treatments are totally confounded with replicates, rows and columns, so the treatment sum of squares and degrees of freedom are zero. The analysis of variance table is not computed, except for replicate, row, column, total sum of squares and degrees of freedom.

NE_G04BC_RESD

The residual degrees of freedom or the residual sum of squares are zero, columns 3, 4 and 5 of the analysis of variance table will not be computed and the matrix of standard errors and covariances, \mathbf{c} , will not be scaled.

NE_G04BC_DISCON

The design is disconnected, the standard errors may not be valid. The design may have a nested structure.

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.

6 Further Comments

To estimate missing values the Healy and Westmacott procedure or its derivatives may be used (see John and Quenouille (1977)). This is an iterative procedure in which estimates of the missing values are adjusted by subtracting the corresponding values of the residuals. The new estimates are then used in the analysis of variance. This process is repeated until convergence. A suitable initial value may be the grand mean. When using this procedure **irdf** should be set to the number of missing values plus one to obtain the correct degrees of freedom for the residual sum of squares.

For analysis of covariance the residuals are obtained from an analysis of variance of both the response variable and the covariates. The residuals from the response variable are then regressed on the residuals from the covariates using, say, nag_regress_confid_interval (g02cbc) or nag_regsn_mult_linear (g02dac). The results from those routines can be used to test for the significance of the covariates. To test the significance of the treatment effects after fitting the covariate, the residual sum of squares from the regression should be compared with the residual sum of squares obtained from the equivalent regression but using the residuals from fitting replicates, rows and columns only.

6.1 Accuracy

The algorithm used in this routine, described in Section 3, achieves greater accuracy than the traditional algorithms based on the subtraction of sums of squares.

6.2 References

Cochran W G and Cox G M (1957) *Experimental Designs* Wiley Davis O L (1978) *The Design and Analysis of Industrial Experiments* Longman John J A (1987) *Cyclic Designs* Chapman and Hall

[NP3652/1]

John J A and Quenouille M H (1977) *Experiments: Design and Analysis* Griffin Searle S R (1971) *Linear Models* Wiley

7 See Also

nag_regress_confid_interval (g02cbc)
nag_regsn_mult_linear (g02dac)

8 Example

The data for a 5×5 Latin square is input and the ANOVA and treatment means computed and printed. Since the design is orthogonal only one standard error need be printed

8.1 Program Text

```
/* nag_anova_row_col (g04bcc) Example Program.
* Copyright 2000 Numerical Algorithms Group.
* Mark 6, 2000.
*/
#include <stdio.h>
#include <nag.h>
#include <nag_stdlib.h>
#include <nagg04.h>
int main (void)
{
  const char *fmt_99999[]={"%3.0f ", "%10.4f ", "%10.4f
                                                                ", "%10.4f
                                                                                ۳,
"%8.4f"};
 double *c=0, c_b20=le-5, *cmean=0, *ef=0, gmean, *r=0, *rmean=0; *rpmean=0;
 double *table=0, *tmean=0, *y=0;
 Integer c__0=0, i, *irep, *it, j, n, ncol, nrep, nrow, nt;
 Integer exit_status=0;
 NagError fail;
#define TABLE(I,J) table[((I)-1)*5 + (J)-1]
#define C(I,J) c[((I)-1)*nt + (J)-1]
 INIT_FAIL(fail);
 Vprintf("g04bcc Example Program Results\n");
  /* Skip heading in data file */
 Vscanf("%*[^\n]");
 Vscanf("%ld %ld %ld %ld", &nrep, &nrow, &ncol, &nt);
  if (!(c = NAG_ALLOC(nt*nt, double))
      || !(cmean = NAG_ALLOC(nrep*ncol, double))
      || !(ef = NAG_ALLOC(nt, double))
      || !(r = NAG_ALLOC(nrep*nrow*ncol, double))
      || !(y = NAG_ALLOC(nrep*nrow*ncol, double))
      || !(rmean = NAG_ALLOC(nrep*nrow, double))
      || !(rpmean = NAG_ALLOC(nrep, double))
      || !(tmean = NAG_ALLOC(nt, double))
      || !(table = NAG_ALLOC(30, double))
      || !(irep = NAG_ALLOC(nt, Integer))
```

```
|| !(it = NAG_ALLOC(nrep*nrow*ncol, Integer)))
    {
      Vprintf("Allocation failure\n");
     exit_status = -1;
     goto END;
    }
 n = nrep * nrow * ncol;
 for (i = 1; i <= n; ++i)
   Vscanf("%lf", &y[i - 1]);
 for (i = 1; i <= n; ++i)
   Vscanf("%ld", &it[i - 1]);
 gO4bcc(nrep, nrow, ncol, y, nt, it, &gmean, tmean, table,
 c, nt, irep, rpmean, rmean, cmean, r, ef, c_b20, c_0,
 &fail):
 if (fail.code != NE NOERROR)
      Vprintf("Error from g04bcc.\n%s\n", fail.message);
     exit_status = 1;
     goto END;
   }
 Vprintf("\n");
 Vprintf("%s\n"," ANOVA TABLE");
 Vprintf("\n");
 if (nrep > 1)
    {
     Vprintf("\n%s", " Reps
                                   ");
     for (j = 1; j <= 5; ++j)
Vprintf(fmt_999999[j-1], TABLE(1,j));
   }
 Vprintf("\n%s"," Rows
                               ");
 for (j = 1; j <= 5; ++j)
   Vprintf(fmt_999999[j-1], TABLE(2,j));
 Vprintf("\n%s", " Columns
                                ");
 for (j = 1; j <= 5; ++j)
   Vprintf(fmt_99999[j-1], TABLE(3,j));
 Vprintf("\n\n%s", " Treatments ");
 for (j = 1; j <= 5; ++j)
   Vprintf(fmt_99999[j-1], TABLE(4,j));
 Vprintf("\n%s", " Residual
                                ");
 for (j = 1; j <= 3; ++j)
   Vprintf(fmt_99999[j-1], TABLE(5,j));
 Vprintf("\n%s", " Total
                                ");
 for (j = 1; j <= 2; ++j)
   Vprintf(fmt_999999[j-1], TABLE(6,j));
 Vprintf("\n%s\n\n", " Treatment means");
 for (i = 1; i <= nt; ++i)</pre>
   Vprintf("%10.4f%s", tmean[i - 1], i%6?"":"\n");
   Vprintf("\n\n%s%10.4f\n", " S.E. of difference (orthogonal design) = ",
C(2,1));
END:
 if (c) NAG_FREE(c);
 if (cmean) NAG_FREE(cmean);
```

g04bcc

```
if (ef) NAG_FREE(ef);
if (r) NAG_FREE(r);
if (y) NAG_FREE(y);
if (rmean) NAG_FREE(rmean);
if (rpmean) NAG_FREE(rmean);
if (tmean) NAG_FREE(tmean);
if (table) NAG_FREE(table);
if (irep) NAG_FREE(irep);
if (it) NAG_FREE(it);
return exit_status;
}
```

8.2 Program Data

gO4bcc Example Program Data

1 5 5 5

8.3 Program Results

g04bcc Example Program Results

ANOVA TABLE

Rows 4 29.4231 7.3558 9.0266 0.0013 Columns 22.9950 7.0545 0.0037 4 5.7487 Treatments 4 0.5423 0.1356 0.1664 0.9514 Residual 12 9.7788 0.8149 Total 24 62.7392 Treatment means 7.3180 7.2440 7.2060 6.9000 7.2600 S.E. of difference (orthogonal design) = 0.5709