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TRGRPS - AN INTERACTIVE ALGORITHM FOR GROUP RECOGNITION WITH AN EXAMPLE FROM SPARTINETEA*

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Introduction

I know that group recognition is a subject of great interest to this group. I realize, of course, that there are preferences for different methods, according to different objectives.

The algorithms used by phytosociologists appear to be of two generic types:

1. The first focuses on homogeneity within the groups. Groups are sought in which the member objects are as similar as possible in terms of given characters. The clustering algorithm, which I described in an earlier paper (Orlóci 1967) with group homogeneity measured in terms of the sum of squares, represents an example in this connection.
2. The second type shifts focus from homogeneity within the groups to discontinuities in the entire sample. The clustering algorithm TRGRPS is of this type. It is formulated around the definition that one group of relevés is discontinuous with another group if none of its neighbourhoods, with given radius, overlap with any of the neighbourhoods with the same radius in the other. The definition and related concepts are explained in the text.

The algorithm is programed in BASIC under file name TRGRPS.BAS. When it is hypothesized that discontinuities subdivide the sample into a given number of groups and the hypothesis is presented to the program with data, TRGRPS. BAS performs the necessary tests automatically and, if groups are found, it produces group descriptors.

*Contribution from the Working Group for Data Processing in Phytosociology, International Society for Vegetation Science. For nomenclature of species see Lausi, Beefink & Kortekaas (1975).

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Relevant Concepts

Sample space

It will be assumed that there are N relevés in the sample labelled by symbols A, B, C, etc. It will be assumed also that in the associated sample space the relevés are represented by points in the line,



The relative placements of the points to one another are important. The assumption of such a simple unidimensional arrangement is not necessary, but convenient, because it can simplify matters when the concepts are explained without detracting from the generality of the conclusions.

Nearest neighbour relevés

The length of the line segment, connecting any two points in sample space, is proportional to differences in the species composition of the relevés. The line segments represent distances in p -dimensions. Each dimension in p -space corresponds to a different species in the sample. It is noted that differences need not be measured in terms of a distance function. Any symmetric measure of dissimilarity is permissible. The point nearest to another point in terms of any such distance or dissimilarity is said to be that point's nearest neighbour. In the example shown, A is nearest neighbour to B, E to F, etc.

The neighbourhood of a relevé

Each relevé may or may not have other relevés in its neighbourhood within a given radius (r) for which the neighbourhood is defined. At $r = d(B, C)$, where $d(B, C)$ is the distance of point B from C, the following neighbourhoods of relevés exist:

{AB}	for relevé A
{ABC}	for relevé B
{BC}	for relevé C
{DE}	for relevé D and E
{F}	for relevé F

It is noted that {AB} and {BC} are overlapping neighbourhoods; *i.e.* relevé B is possessed in common by both. {ABC}, {DE} and {F} are non-overlapping. Neighbourhoods that are non-overlapping are said to be discontinuous.

Clustering by Neighbourhoods

Given $r = 0$ as the initial state of the neighbourhood radius in a clustering procedure, as many discontinuous neighbourhoods can be recognized in the sample A, B, C, D, E, F as there are non-identical relevés.

{A}, {B}, {C}, {D}, {E}, {F}

Now, let r be incremented by a constant, small quantity in steps until the condition $d(A, B) \leq r < d(B, C)$ is satisfied. Here, $d(A, B)$ represents the smallest distance other than zero in the sample and $d(B, C)$ the second smallest distance. At such a radius,

{AB}, {C}, {DE}, {F}

are recognized as discontinuous neighbourhoods. If r is further incremented, so that it will satisfy the condition $d(B, C) \leq r < d(E, F)$, where $d(E, F)$ is the third smallest distance in the sample, the discontinuous neighbourhoods are

{ABC}, {DE}, {F}

Further increments in r , until $d(E, F) \leq r < d(C, D)$ is satisfied, yield the discontinuous neighbourhoods,

{ABC}, {DEF}

Finally, when $r \geq d(C, D)$, only a single group {ABCDEF}

is recognized. All neighbourhoods at a given r under the given constraints satisfy the original definition of discontinuous groups stated in the introduction. In other words, none of the neighbourhoods in any one group overlaps with any neighbourhood in the other. It is evident from what has been said that the procedure of clustering by neighbourhoods uses single link fusions similarly to Sneath's (1957) method of single linkage clustering.

The Algorithm TRGRPS

TRGRPS incorporates clustering by neighbourhoods and various other devices which enable the user to test hypotheses about the existence of discontinuous groups of relevés. The algorithm regards a group of relevés as discontinuous if none of its neighbourhoods overlap with any neighbourhood in other groups. The flow pattern of operations and decisions is described below:

- Present data; specify a hypothesis for the number of groups R ; compute distances between relevés.
- Specify initial value of neighbourhood radius and its increments.
- Cluster relevés and count the number of groups. Designate this number by k .
- If $k = R$ then accept the hypothesis.
If $k < R$ then modify the hypothesis to less than R groups and continue with step b.
If $k > R$ then increment neighbourhood radius and continue at c.

I realize that based on sampling considerations tests of the kind as in this algorithm should preferably be probabilistic. However, the algorithm as I present it here does not incorporate probabilistic considerations. I plan to discuss these aspects in a separate essay presently in the preparatory stage.

A Related Algorithm

TRGRPS evolved through modifications of the original Jancey (1974) algorithm which I abstracted (Orlói 1975) under the name TRGRUP. I stressed a nearest neighbour property of the groups in the original descriptions of TRGRUP, although I incorporated the idea of clustering by neighbourhoods in the program TRGRUP.BAS. The present paper corrects this anomaly by recognizing that although the condition that every member in a discontinuous group must have its nearest neighbour within the group is necessary it is by no means sufficient because when this condition is satisfied neighbourhoods can overlap between the groups. The groups with contents

{AB}, {C}, {DEF}

represent an éclatant example in this connection. In these, the nearest neighbour property is satisfied; every member in each group has its nearest neighbour within its own group. Yet two groups, {AB} and {C}, appear continuous because the neighbourhoods of relevés A, B and C overlap at the implied radius $d(E, F)$.

TRGRPS.BAS differs from TRGRUP.BAS by the speci-

fic constraint that none of the relevés are allowed to remain unclustered when the hypothesis is true. This means that for the hypothesis of R groups to be accepted in TRGRPS, the sample must divide into exactly R groups, and not just R nodes plus a number of points that float unclustered among the groups. Programs TRGRUP.BAS and TRGRPS.BAS yield identical results when the group structure is strong in the sample.

Running the Program TRGRPS.BAS

The program uses the BASIC programming language, written in conversational mode for processing from a teletype console. After the RUN command, the program requests values for the following parameters:

Number of species (P)Number of relevés (N)Minimum group size (S)Number of groups (R)Initial value of neighbourhood radius (C)Increment in neighbourhood radius (D)

Values for P and N are requested first, and after P and N , values for S and R . S can be set to one if no hypothesis specifies *a priori* minimum group size. An initial value for R derives from the hypothesis. Alternate values, small or large, may be assigned to R in successive runs until a solution is accepted. It is noted that running time and R are inversely related.

Data input is from DISK file RAWD which contains P sets of N numbers. The data in computer core are stored in array **D** which has P rows and N columns. Data input is followed by computation of relevé distances. The program uses the Euclidean distance. However, if so desired, any symmetric function of a relevé dissimilarity may replace the Euclidean distance. The distances are stored in array **A** which has N^2 elements.

After the elements of **A** are generated, a search begins for the smallest value. The message initial value of linkage parameter is printed followed by display of the smallest distance value. A request to confirm this value is displayed next. The user responds by typing either the same value, or another value, which then serves as the initial value for **C**. The program determines a maximum of 50 values, each representing a difference between pairs of the smallest distance values in **A**, ordered according to size, to assist in the choice of a value for the increment **D** of **C**. Any value can be chosen. However,

it must be kept in mind that a large value of D shortens the computation time required, while a small value increases precision. A zero is not permissible for D .

Once the program gets underway, the sample is searched for discontinuous groups. When exactly R groups are found, the group descriptors are printed and the execution of the program stops. If R groups are not found a new search begins for $R - 1$ groups.

Printout of Results

Examples are given in the next section. When a solution is found under a current hypothesis for R , as modified by parameters S , C , D , the following descriptors are printed:

- A value of C representing the current size of the neighbourhood radius r at the last fusion. This can be used as a global measure for coherence in the groups.
- A label for each group.
- Relevé labels in each group.

Results of repeated runs at different R can be used to trace a hierarchical group structure of relevés in the sample.

Results from Spartinetea

The results presented here derive from the data in Table 1 representing saltmarsh relevé section f (Van der Maarel, Orlóci & Pignatti 1976, Lausi, Beefink & Kortekaas 1976). Inspection of the raw data reveals that an initial hypothesis of 5 relevé groups may be tenable with S set equal to 5. When TRGRPS is run a solution is indeed found at this value of R (see Table 2).

Table 1. Sample of 50 reprints from *Spartinetwa*. The Braun-Blanquet scale is replaced by the following scores in the table: 1 = +, 2 = *, 3 = 1, 5 = 2, 7 = 3, 8 = 4, 9 = 5. Releve order follows syntaxonomic affiliation.

[illegible]

Table 2. Results from program TRGRPS.BAS. The number of groups hypothesized is 5. See the main text for description of program.

```

TRGRPS      16:15      22-Nov-75
PROGRAM NAME -- TRGRPS
-----
NUMBER OF SPECIES 227
NUMBER OF QUADRATS 750
MINIMUM GROUP SIZE REQUIRED 75
NUMBER OF GROUPS REQUIRED 75
INITIAL VALUE OF NEIGHBOURHOOD RADIUS = 1
CONFIRM BY TYPING 1 ELSE TYPE DESIRED VALUE: 72
SELECT A VALUE FROM THE LIST BELOW
TO SERVE AS THE INCREMENT IN LINKAGE PARAMETER.
1 2
TYPE SELECTED VALUE:
72
5 - GROUPS
MAXIMUM NEIGHBOURHOOD RADIUS 67
GROUP 1
1 2 3 5 6
GROUP 2
5 7 8 10 9
GROUP 3
11 12 14 15 13
GROUP 4
16 19 20 17 18
GROUP 5
21 22 25 24 23 27 28 29 30 34 31 32 35 33 36 37 40
50 46 47 48 38 40 19 26 41 47 42 45 44
TIME: 177.74 SECS.

```

The correspondence of the groups at the 5-group level in TRGRPS and the units in conventional syntaxonomy is most remarkable. The first group, and the second, are subassociations (*typicum* and *salicornietosum*) of *Spartinetum maritimae*. The third group represents *Spartinetum alterniflorae*, the fourth *Spartinetum patensis*, and the fifth *Spartinetum townsendii*.

A hierarchical pattern of fusions, derived by repeated runs of TRGRPS. BAS with different values of *R*, is traced in Fig. 1. Interestingly, the first fusion does not occur between the subassociations of *Spartinetum maritimae*, notwithstanding expectations, but between *Spartinetum maritimae salicornietosum* and *Spartinetum townsendii*. With these the other subassociation, *typicum*, fuses only in the second step, fol-

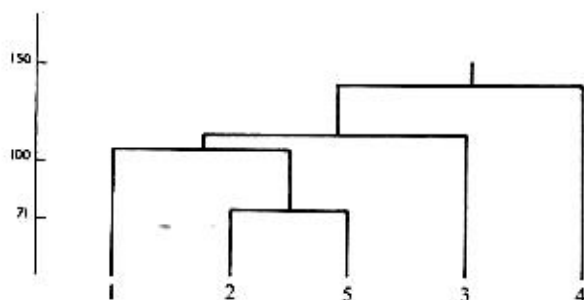


Figure 1. Hierarchical structure in Spartinetea (Table 1) at the level of 5 groups. Vertical scale indicates values of the neighbourhood radius. Symbols at the base designate syntaxonomic units recognized in TRGRPS. Legend to symbols: 1 - *Spartinetum maritimae typicum*, 2 - *Spartinetum maritimae salicornietosum*, 5 - *Spartinetum townsendii*, 3 - *Spartinetum alterniflorae*, 4 - *Spartinetum batensis*. Numbers 1 to 5 identify groups in the computer printouts at the 5 groups level in Table 2.

lowed by fusions with *Spartinetum alterniflorae*, and lastly, *Spartinetum patensis*.

Discussion

It is evident intuitively that the relevés in a discrete group must exhibit a density pattern that permits their recognition as members in a coherent group. Coherence, the discreteness of groups, is associated with criteria of non-overlapping neighbourhoods in algorithm TRGRPS. When applied to a sample of *Spartinetea*, the results of a comparison of the groups suggest that the method of group recognition in TRGRPS resembles the method of group recognition in traditional syntaxonomy.

The algorithm TRGRPS should be a useful tool in the hands of phytosociologists, interested in the construction of syntaxonomic systems and in the identification of a parsimonious path through neighbouring groups for which hierarchical relationships are to be shown.

Summary

TRGRPS can detect groups or signal if discrete groups cannot be found in a sample. The present paper elaborates on the concepts, describes the algorithm and provides illustrations from syntaxonomy. A computer program (TRGRPS.BAS) is available from the author upon request.

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