Aggregation of pairwise comparison matrices: A clustering approach

Kolos Csaba Ágoston^a

Sándor Bozóki^b

László Csató^c

21st February 2024

"There can be many reasons for searching for representative objects. Not only can these objects provide a characterization of the clusters, but they can often be used for further work or research, especially when it is more economical or convenient to use a small set of k objects instead of the large set one started off with."¹

Abstract

We consider clustering in group decision making where the opinions are given by pairwise comparison matrices. In particular, the k-medoids model is suggested to classify the matrices as it has a linear programming problem formulation. Its objective function depends on the measure of dissimilarity between the matrices but not on the weights derived from them. With one cluster, our methodology provides an alternative to the conventional aggregation procedures. It can also be used to quantify the reliability of the aggregation. The proposed theoretical framework is applied to a large-scale experimental dataset, on which it is able to automatically detect some mistakes made by the decision-makers.

Keywords: Decision analysis; aggregation; clustering; large-scale group decision making; pairwise comparison matrix

MSC class: 90-10, 90B50, 91B08

JEL classification number: C38, C44

^a Email: kolos.agoston@uni-corvinus.hu

Corvinus University of Budapest (BCE), Institute of Operations and Decision Sciences, Department of Operations Research and Actuarial Sciences, Budapest, Hungary

 $^{^{\}rm b}$ Email: sandor.bozoki@sztaki.hun-ren.hu

Institute for Computer Science and Control (SZTAKI), Hungarian Research Network (HUN-REN), Laboratory on Engineering and Management Intelligence, Research Group of Operations Research and Decision Systems, Budapest, Hungary

Corvinus University of Budapest (BCE), Institute of Operations and Decision Sciences, Department of Operations Research and Actuarial Sciences, Budapest, Hungary

^c Corresponding author. Email: laszlo.csato@sztaki.hun-ren.hu

Institute for Computer Science and Control (SZTAKI), Hungarian Research Network (HUN-REN), Laboratory on Engineering and Management Intelligence, Research Group of Operations Research and Decision Systems, Budapest, Hungary

Corvinus University of Budapest (BCE), Institute of Operations and Decision Sciences, Department of Operations Research and Actuarial Sciences, Budapest, Hungary

¹ Source: Kaufman and Rousseeuw (1990, p. 71).

1 Introduction

The fast development of information technology has enabled large-scale group decision making (LGDSM): in many decision making problems, the possible number of decision-makers (DMs) now easily reaches thousands (García-Zamora et al., 2022). However, because the cognitive abilities of humans have not evolved parallel to the exponential growth of data, it is necessary to aggregate the opinions of DMs (Tang and Liao, 2021; Fan et al., 2024). For this purpose, one of the most widely used techniques is *clustering*: allocating the individual judgements into groups such that judgements in the same group (called a cluster) are more similar to each other than to those in other groups (clusters).

The Analytic Hierarchy Process (AHP) (Saaty, 1977, 1980) is a popular multi-criteria decision making (MCDM) methodology, hence, solving group AHP (GAHP) models is almost as old as AHP itself (Aczél and Saaty, 1983). Traditionally, there are two main approaches to aggregating individual preferences and creating a group consensus: (a) aggregating the individual pairwise comparison matrices (PCMs) and deriving priorities from the common matrix (Aczél and Saaty, 1983); and (b) deriving priorities from the individual PCMs and aggregating these priorities (Basak and Saaty, 1993). Ossadnik et al. (2016) recommend the second option as no other aggregation technique satisfies a comparable number of evaluation criteria. However, this approach is sensitive to extreme opinions, the supposed consensus may not accurately portray the true group preference (Amenta et al., 2020). Furthermore, Duleba and Szádoczki (2022) find that distance-based aggregation outperforms conventional methods for a high number of DMs.

On the other hand, we do not know of any studies where individual PCMs are clustered in order to aggregate them. The current paper aims to fill this research gap, which is our main contribution to the extant literature.

The proposed clustering methodology can be used for several purposes since:

- it provides an alternative aggregation procedure if the number of clusters is restricted to one;
- it gives information on the ability of aggregated preferences to represent the individual preferences;
- it makes it possible to detect some mistakes in the original data, which is far from trivial in LGDSM problems.

In addition, a crucial advantage of our aggregation method is the lack of any requirement on the number of missing entries in the PCMs, which is an especially useful property in LGDSM problems where the underlying pairwise comparisons are obtained via a questionnaire that can be finished at every point (Fan et al., 2024). Naturally, a complete PCM will have a stronger effect on clustering, but this seems to be a fair compensation for the DMs who have devoted more effort and time to reveal their preferences.

The proposed clustering approach guarantees that the cluster centres are PCMs given by at least one DM. This is also an attractive property because a cluster centre is difficult to interpret if it strongly differs from individual opinions in a MCDM problem.

Last but not least, the suggested k-medoids clustering has a linear programming (LP) formulation, which allows for imposing additional restrictions. For example, the inconsistency of the cluster centres can be required to remain below an exogenous threshold.

The theoretical framework is applied to the experimental data of Bozóki et al. (2013).

The paper is structured as follows. The methodology is described in Section 2, and the numerical results are presented in Section 3. Finally, Section 4 offers concluding remarks.

2 Theoretical background

First, the mathematics underlying the proposed approach is discussed. Section 2.1 introduces the necessary concepts from the literature of pairwise comparison matrices, Section 2.2 summarises the details of cluster analysis. Section 2.3 overviews some measures to quantify the (dis)similarity of pairwise comparison matrices, while Section 2.4 provides a LP formulation for our k-medoids clustering model.

2.1 Pairwise comparison matrices

Pairwise comparison matrices are often used to determine the cardinal preferences of DMs, who answer questions like "How many times is a criterion more important than another one?" or "How many times is a given alternative better than another one with respect to a fixed criterion?" These pairwise ratios are collected into the $n \times n$ pairwise comparison matrix $A = [a_{ij}]$, where $a_{ij} > 0$ and $a_{ij} = 1/a_{ji}$ hold for all $1 \le i, j \le n$. A is called consistent if $a_{ij}a_{jk} = a_{ik}$ is satisfied for all $1 \le i, j, k \le n$. Otherwise, when cardinal transitivity is violated, the matrix is called inconsistent.

From a practical point of view, it is natural and reasonable to allow for some degree of inconsistency. Several inconsistency indices have been proposed to measure the contradictions in a pairwise comparison matrix (Bozóki and Rapcsák, 2008; Brunelli, 2018; Csató, 2019a). In particular, Saaty (1977) has suggested the so-called *inconsistency index* CI in his celebrated theory of AHP, which is an affine transformation of the dominant eigenvalue λ_{max} of the pairwise comparison matrix A:

$$CI(A) = \frac{\lambda_{\max} - n}{n - 1}.$$

This is divided by the random index RI_n , the average CI of a large number of randomly generated $n \times n$ pairwise comparison matrices to obtain the *inconsistency ratio* $CR(A) = CI(A)/RI_n$. According to Saaty, CR should remain below 0.1 to accept the matrix as a reasonable representation of consistent preferences.

Analogously, many weighting methods are available to find a positive weight vector $\mathbf{w} = (w_1, w_2, \ldots, w_n)$ such that the ratios w_i/w_j give good approximations to the pairwise comparisons a_{ij} (Choo and Wedley, 2004). The *Logarithmic Least Squares Method (LLSM)* (Crawford and Williams, 1985; Csató, 2019b; De Graan, 1980; Fichtner, 1986; Rabinowitz, 1976) is one of the most popular:

$$\min \sum_{i,j} \left[\log a_{ij} - \log \left(\frac{w_i}{w_j} \right) \right]^2$$
$$\sum_{i=1}^n w_i = 1,$$
$$w_i > 0, \qquad i = 1, 2, \dots, n.$$

The unique solution to this optimisation problem is given by the geometric means of the entries in the rows:

$$\frac{w_i}{w_j} = \frac{\sqrt[n]{\prod_{k=1}^n a_{ik}}}{\sqrt[n]{\prod_{k=1}^n a_{jk}}}$$

Therefore, it is often called the *geometric mean* method.

2.2 Cluster analysis

Clustering, the problem of dividing a set of objects into groups (clusters) such that any object is more similar to an arbitrary object in its group than to any object in a distinct group, is one of the most popular methods in exploratory data science. Although no "perfect" clustering algorithm exists, perhaps the most natural solution is k-means clustering, when each object belongs to the cluster with the nearest mean.

However, this leads to an NP-hard non-convex discrete optimisation problem. Therefore, most state-of-the-art statistical and data science software implement a heuristic iterative algorithm that alternates two steps: i) assigning all points to the nearest cluster centre; and ii) recalculating the cluster centres. The operation in step ii) is quite trivial in Euclidean spaces (take the average in every dimension), but it may be challenging if the objects do not belong to a multidimensional Euclidean space (Majstorović et al., 2018). Furthermore, the heuristic iterative algorithm above can converge to a local rather than global optimum. Hence, Ágoston and E.-Nagy (2024) provide a mixed integer linear programming (MILP) formulation to obtain an exact solution for the minimum sum-of-clustering problem that allows for adding arbitrary linear constraints and can be solved up to 150 objects.

An alternative approach to the k-means clustering is the k-medoids problem (Kaufman and Rousseeuw, 1986; Schubert and Rousseeuw, 2019). Here, all cluster centres should be objects themselves, allowing for easier interpretability of the cluster centres. This seems to be an important advantage in LGDSM since the DMs may not be willing to accept a solution that has not been suggested by any of them. In addition, k-medoids can be used with an arbitrary dissimilarity measure, while k-means generally requires Euclidean distance. Last but not least, minimising the sum of pairwise dissimilarities instead of the sum of squared Euclidean distances makes the k-medoids problem less sensitive to outliers than the k-means problem.

For unknown reasons, the k-medoids model is not widely used and is not implemented in usual statistical software packages. Nonetheless, the problem can be formulated as an LP model according to Section 2.4, and solved with standard LP solvers in small and medium size instances.

Originally, clustering was intended to form groups in Euclidean spaces, but the underlying idea can be applied in other fields. Clustering algorithms appear in network science, where the goal is to classify nodes in a graph such that the nodes in a given cluster are connected more strongly to each other than to nodes in other clusters. Other objects can also be clustered, for example, curves or mortality tables.

To conclude, cluster analysis requires a clustering algorithm and a dissimilarity measure, which does not necessarily be a distance. This paper considers the k-medoids problem due to its advantages discussed above.

2.3 Quantifying the similarity of pairwise comparison matrices

Several ideas exist in the literature for computing the (dis)similarity of two pairwise comparison matrices A and B. Crawford and Williams (1985, p. 389) essentially introduces the following metric:

$$D_1(A,B) = \sqrt{\left(\sum_{i=1}^n \sum_{j=1}^n \left(\log(a_{ij}) - \log(b_{ij})\right)^2\right)}.$$

The original definition takes the sum only for the entries above the diagonal, but this is equivalent to D_1 except for a constant factor. Obviously, D_1 is symmetric and equals 0 if and only if A = B. It also satisfies the triangle inequality (Crawford and Williams, 1985).

Tekile et al. (2023) measure the closeness of two complete pairwise comparison matrices by another formula called Manhattan or L^1 distance:

$$D_2(A, B) = \sum_{i=1}^n \sum_{j=1}^n |\log(a_{ij}) - \log(b_{ij})|.$$

 D_2 also satisfies the triangle inequality.

Kułakowski et al. (2022) use another indicator, the so-called compatibility index, which was originally defined by Saaty (2008):

$$D_{3u}(A,B) = \frac{1}{n^2} \left(\sum_{i=1}^n \sum_{j=1}^n a_{ij} b_{ji} \right).$$

While D_{3u} is symmetric, $D_{3u}(A, B) \ge n^2$. Hence, it is worth normalising D_{3u} as has been done in Ágoston and Csató (2024):

$$D_3(A,B) = \frac{1}{n^2} \left(\sum_{i=1}^n \sum_{j=1}^n (a_{ij}b_{ji} - 1) \right).$$

 D_3 remains symmetric, and $D_3(A, B) = 0$ if and only if A = B. However, D_3 is not a distance.

Lemma 1. Dissimilarity index D_3 does not satisfy the triangle inequality.

Proof. It is sufficient to give a counterexample. Consider the following three pairwise comparison matrices:

$$A = \begin{bmatrix} 1 & 2 & 1 \\ 1/2 & 1 & 1 \\ 1 & 1 & 1 \end{bmatrix}, \qquad B = \begin{bmatrix} 1 & 3 & 1 \\ 1/3 & 1 & 1 \\ 1 & 1 & 1 \end{bmatrix}, \qquad C = \begin{bmatrix} 1 & 4 & 1 \\ 1/4 & 1 & 1 \\ 1 & 1 & 1 \end{bmatrix}.$$

It can be checked that $D_3(A, B) = 1/6$, $D_3(B, C) = 1/12$, but $D_3(A, C) = 1/2 > 1/4 = D_3(A, B) + D_3(B, C)$.

Kułakowski et al. (2022) consider three other versions of the compatibility index (all of them are modified to ensure D(A, A) = 0 and $D(A, B) \ge 0$):

$$D_4(A,B) = \frac{2}{n(n-1)} \left(\sum_{i=1}^{n-1} \sum_{j=i+1}^n \max\left\{ a_{ij}b_{ji}, a_{ji}b_{ij} \right\} - 1 \right);$$

$$D_5(A,B) = \max\left\{ a_{ij}b_{ji} - 1 : 1 \le i, j \le n \right\};$$

$$D_6(A,B) = -\frac{2}{n(n-1)} \left(\sum_{i=1}^{n-1} \sum_{j=i+1}^n \min\left\{ a_{ij}b_{ji}, a_{ji}b_{ij} \right\} - 1 \right).$$

Based on D_4 – D_6 , another reasonable index could be

 $D_7(A, B) = -\min \{a_{ij}b_{ji} - 1 : 1 \le i, j \le n\}.$

Lemma 2. Indices D_4 and D_5 do not satisfy the triangle inequality.

Proof. The PCMs from the proof of Lemma 1 can be used: $D_4(A, B) = 1/6$, $D_4(B, C) = 1/9$ but $D_4(A, C) = 1/3 > 5/18 = D_4(A, B) + D_4(B, C)$, and $D_5(A, B) = 1/2$, $D_5(B, C) = 1/3$ but $D_5(A, C) = 1 > 5/6 = D_5(A, B) + D_5(B, C)$.

Proposition 1. Indices D_6 and D_7 satisfy the triangle inequality, thus, they are distances on the set of pairwise comparison matrices.

Proof. For both D_6 and D_7 , it is sufficient to show that the triangle inequality holds elementwise for any pairwise comparison matrices $A = [a_{ij}]$, $B = [b_{ij}]$, and $C = [c_{ij}]$. $1 = a_{ij} \leq b_{ij} \leq c_{ij}$ can be assumed without loss of generality. Then, by elementary calculus, the following three inequalities hold:

$1 - \frac{a_{ij}}{b_{ij}}$	+	$1 - \frac{b_{ij}}{c_{ij}}$	\geq	$1 - \frac{a_{ij}}{c_{ij}};$
$1 - \frac{a_{ij}}{c_{ij}}$	+	$1 - \frac{b_{ij}}{c_{ij}}$	\geq	$1 - \frac{a_{ij}}{b_{ij}};$
$1 - \frac{a_{ij}}{b_{ij}}$	+	$1 - \frac{a_{ij}}{c_{ij}}$	\geq	$1 - \frac{b_{ij}}{c_{ij}}.$

Fichtner (1984) verifies that the popular eigenvector method (Saaty, 1977) can be defined by a metric, too. However, although it satisfies the requirements of a distance, it is discontinuous.

All of the above dissimilarity measures D_1-D_7 are invariant under the relabeling of the alternatives. Consequently, they are invariant to transposition, that is,

$$D_i(A,B) = D_i(A^{\top}, B^{\top}),$$

where A^{\top} is the transpose of the pairwise comparison matrix A.

Since the input of the k-medoids problem is the dissimilarity matrix of the objects, any of the indices D_1-D_7 can be used in the proposed methodology. Our numerical experiment will consider D_1 and D_3 .

2.4 An LP formulation of the *k*-medoids clustering problem

Assume that the $m \times m$ matrix $\Delta = [\delta_{ij}]$ contains the (pairwise) dissimilarities between any two of the *m* objects (individual PCMs). For example, $\delta_{ij} = D_1 \left(A^{(i)}, A^{(j)} \right)$ if measure D_1 is used to determine the dissimilarity of the preferences given by matrices $A^{(i)}$ and $A^{(j)}$ of decision makers *i* and *j*. Let *M* denote the set of integers $\{1, \ldots, m\}$.

Label	Туре	Size (n)	Number of PCMs (m)
M4	Objective (map)	4	66
M6	Objective (map)	6	77
M8	Objective (map)	8	82
$\mathbf{S4}$	Subjective (summer house)	4	68
$\mathbf{S6}$	Subjective (summer house)	6	77
S8	Subjective (summer house)	8	78

Table 1: Sample sizes in the experimental database

The k-medoids problem is as follows (see, for instance, Vinod (1969)):

$$\sum_{i=1}^{m} \sum_{j=1}^{m} \delta_{ij} x_{ij} \to \min$$
(1)

s.t.
$$\sum_{j=1}^{m} x_{ij} = 1 \qquad \forall i \in M \qquad (2)$$

$$x_{ij} \le y_j \qquad \qquad \forall \ i, j \in M \tag{3}$$

$$\sum_{j=1}^{m} y_j = k \tag{4}$$

$$\begin{aligned} x_{ij} &\geq 0 & \forall i, j \in M \\ y_j &\in \{0, 1\} & \forall j \in M \end{aligned}$$

In this formulation, the binary variable y_j equals one if object j is a cluster centre (and zero otherwise), while x_{ij} equals one if object i is placed in the cluster whose centre is object j (and zero otherwise). Constraint (4) ensures that there are k different cluster centres. According to (2), each object is classified into exactly one cluster. Finally, due to the constraints (3), object i can be associated with object j ($x_{ij} = 1$) only if object j is a cluster centre, namely, $y_j = 1$.

3 Numerical results

Bozóki et al. (2013) have collected 454 pairwise comparison matrices in a controlled experiment where even the questioning order is known for each matrix. However, here we use only the final complete pairwise comparison matrices. The matrices are distinguished by their size (n = 4, n = 6, n = 8 alternatives) and type (objective, where the students compared maps; subjective, where the students compared summer houses). The sample sizes are reported in Table 1.

3.1 Analysing a sample of objective type

In the case of dataset M4, the "correct" consistent pairwise comparison matrix provided by the ratios of country areas that appear on the map are known:

$$\begin{bmatrix} 1 & 1.691 & 0.282 & 0.770 \\ 0.591 & 1 & 0.167 & 0.455 \\ 3.544 & 5.991 & 1 & 2.725 \\ 1.300 & 2.198 & 0.367 & 1 \end{bmatrix}.$$

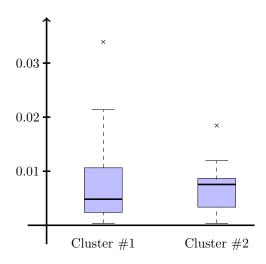


Figure 1: Distribution of inconsistency ratios CR, dataset M4, k = 2 clusters

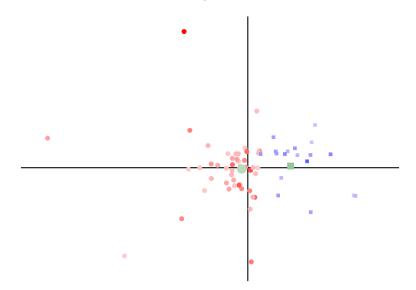


Figure 2: Approximated coordinates of PCMs, dataset M4, measure D_3

Notes: The two clusters are represented by red dots and blue squares. A darker mark is associated with a higher inconsistency. The cluster centers are green and have double size.

For k = 2 clusters, the cluster centres coincide for both dissimilarity measures D_1 and D_3 :

$$CC_1^{(1)} = CC_3^{(1)} = \begin{bmatrix} 1 & 1.500 & 0.286 & 0.833 \\ 0.667 & 1 & 0.154 & 0.435 \\ 3.500 & 6.500 & 1 & 2.300 \\ 1.200 & 2.300 & 0.435 & 1 \end{bmatrix},$$

and

$$CC_1^{(2)} = CC_3^{(2)} = \begin{bmatrix} 1 & 1.200 & 0.400 & 0.909 \\ 0.833 & 1 & 0.200 & 0.500 \\ 2.500 & 5.000 & 1 & 2.300 \\ 1.100 & 2.000 & 0.500 & 1 \end{bmatrix}.$$

The two clusters are the same: 47 matrices belong to first and 19 to the second group.

Figure 1 shows boxplots for the inconsistency ratios CR in the two clusters.

We have used multidimensional scaling (MDS) to visualise the pairwise comparison matrices (Kruskal, 1964; Kruskal and Wish, 1978). This technique places each matrix

into a lower-dimensional space such that the original distances are preserved to the extent possible. In Figure 2, red dots and blue squares represent the two clusters. A darker mark is associated with a higher inconsistency. The two cluster centres are green and have double size. Clearly, the clusters are not determined by the inconsistency of the matrices.

The main findings can be summarised as follows:

and

- The clusters are quite balanced with respect to the number of matrices in them;
- The results are insensitive to the dissimilarity measure $(D_1 \text{ or } D_3)$ used;
- The two groups are not distinguished by the level of inconsistency, although the variance of CR is somewhat higher for the first group.

The analysis has been repeated with higher numbers of clusters. A strong relationship remains between the resulting groups for different values of k, as well as for the two dissimilarity measures D_1 and D_3 .

In sample M8 with measure D_3 and k = 2 clusters, an interesting example has been found that uncovers a possible application of our methodology. The cluster centres are:

	[1	1.800	0.769	0.556	1.800	5.000	1.100	3.000	
	0.556	1	0.455	0.313	1.100	2.700	0.909	1.500	
	1.300	2.200	1	0.833	3.000	7.000	1.800	3.300	
$CC_{3}^{(1)} =$	1.800	3.200	1.200	1	5.000	10.100	2.200	6.000	
$CC_3^{+} =$	0.556	0.909	0.333	0.200	1	2.300	0.625	1.400	,
	0.200	0.370	0.143	0.099	0.435	1	0.250	0.769	
	0.909	1.100	0.556	0.455	1.600	4.000	1	2.200	
	0.333	0.667	0.303	0.167	0.714	1.300	0.455	1	
	Γ 1	1.500	0.500	0.500	3.000	6.000	1.200	3.000	
	0.667	1	0.500	0.333	1.500	3.000	0.667	1.500	
	2.000	2.000	1	0.667	5.000	8.000	1.500	4.000	
$CC_{3}^{(2)} =$	2.000	3.000	1.500	1	4.500	0.100	2.000	5.000	
$CC_3 =$	0.333	0.667	0.200	0.222	1	2.000	0.667	1.200	•
	0.167	0.333	0.125	10.000	0.500	1	0.286	0.500	
	0.833	1.500	0.667	0.500	1.500	3.500	1	2.500	
	0.333	0.667	0.250	0.200	0.833	2.000	0.400	1	

The first cluster contains 81 matrices and the second cluster consists of only one. The two cluster centres are similar except for the preference between the fourth and sixth alternatives (countries), highlighted by bold font. Since the pairwise comparisons show the ratio of the area of two countries, the only PCM in the second cluster is almost certain to contain a typo: the student has thought that country 4 is ten times larger than country 6 but mistakenly written the reciprocal 1/10, which is a standard mistake. Unsurprisingly, the associated PCM fundamentally differs from all other matrices, and the presented clustering approach is able to detect the outlier without any other specification. This can be highly advantageous in LGDSM problems.

3.2 Analysing a sample of subjective type

When the students have compared summer houses, no "natural" PCM exists around which the opinions are centred.

	$CL_{3}^{(1)}$	$CL_{3}^{(2)}$	$CL_{3}^{(3)}$	$CL_{3}^{(4)}$	\sum
$CL_{1}^{(1)}$	25	0	1	0	26
$CL_{1}^{(2)}$	0	9	5	1	15
$CL_{1}^{(3)}$	2	0	0	7	9
$CL_{1}^{(4)}$	0	0	6	12	18
\sum	27	9	12	20	68

Table 2: Contingency table for sample S4

Four clusters, measures D_1 (row) and D_3 (column)

Since sample S4 contains four different houses, we have started the analysis with four clusters, supposing that each alternative will be the best in one cluster.

For the measure D_1 , the cluster centres are as follows:

$CC_{1}^{(1)} =$	$\begin{bmatrix} 1\\ 0.500\\ 0.143\\ 0.333 \end{bmatrix}$	$2.000 \\ 1 \\ 0.200 \\ 0.500$	$7.000 \\ 5.000 \\ 1 \\ 2.000$	$\begin{array}{c} 3.000 \\ 2.000 \\ 0.500 \\ 1 \end{array} \right]$,	$CC_{1}^{(2)} = $	$ \begin{array}{c} 1\\ 0.667\\ 2.000\\ 3.000 \end{array} $	$1.500 \\ 1 \\ 3.000 \\ 3.000$	$0.500 \\ 0.333 \\ 1 \\ 1.500$	$\begin{array}{c} 0.333 \\ 0.333 \\ 0.667 \\ 1 \end{array}$],
$CC_1^{(3)} =$	$\begin{bmatrix} 1\\ 0.333\\ 0.200\\ 2.500 \end{bmatrix}$	3.000 1 0.333 7.000	$5.000 \\ 3.000 \\ 1 \\ 9.000$	$\begin{array}{c} 0.400 \\ 0.143 \\ 0.111 \\ 1 \end{array} \right]$,	$CC_{1}^{(4)} = $	1 0.167 0.333 0.333	6.000 1 5.000 5.000	$3.000 \\ 0.200 \\ 1 \\ 2.000$	3.000 0.200 0.500 1].

On the other hand, for the measure D_3 :

$$CC_{3}^{(1)} = \begin{bmatrix} 1 & 2.000 & 7.000 & 3.000 \\ 0.500 & 1 & 5.000 & 2.000 \\ 0.143 & 0.200 & 1 & 0.500 \\ 0.333 & 0.500 & 2.000 & 1 \end{bmatrix}, \qquad CC_{3}^{(2)} = \begin{bmatrix} 1 & 1.500 & 0.333 & 0.200 \\ 0.667 & 1 & 0.333 & 0.200 \\ 3.000 & 3.000 & 1 & 0.500 \\ 5.000 & 5.000 & 2.000 & 1 \end{bmatrix}.$$
$$CC_{3}^{(3)} = \begin{bmatrix} 1 & 4.000 & 1.500 & 2.000 \\ 0.250 & 1 & 0.333 & 0.667 \\ 0.667 & 3.000 & 1 & 3.000 \\ 0.500 & 1.500 & 0.333 & 1 \end{bmatrix}, \qquad CC_{3}^{(4)} = \begin{bmatrix} 1 & 5.000 & 3.000 & 1.000 \\ 0.200 & 1 & 0.500 & 0.143 \\ 0.333 & 2.000 & 1 & 0.200 \\ 1.000 & 7.000 & 5.000 & 1 \end{bmatrix}.$$

Note that the centres of the first clusters (matrices $CC_1^{(1)}$ and $CC_3^{(1)}$) coincide. The cluster sizes are 26, 15, 9, 18 for index D_1 and 27, 9, 12, 20 for index D_3 .

The contingency table of the two groupings is shown in Table 2. The similarity of the clusters is lower than in the objective task (Section 3.1), but the largest cluster almost coincides.

The inconsistency ratios of the cluster centres are 0.007, 0.008, 0.028, 0.063 in the case of D_1 , while 0.007, 0.009, 0.022, 0.013 in the case of D_3 . Similar to the results of Section 3.1, the clusters are relatively uniform with respect to the level of inconsistency as can be seen in Figure 3. Even though the cluster centres are slightly less inconsistent for D_3 , each centre has an acceptable inconsistency according to the famous 10% rule of thumb. Naturally, this is not guaranteed in other datasets but the LP model of the *k*-medoids problem (Section 2.4) might contain any (linear) restriction on the inconsistency of the cluster centres.

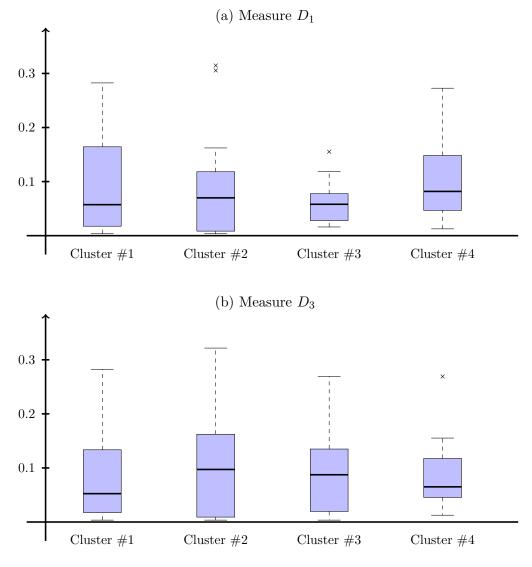


Figure 3: Distribution of inconsistency ratios CR, dataset S4, k = 4 clusters

As the cluster centres are PCMs, the associated priority weights can help the interpretation of the clusters. To that end, we have used the geometric mean method (Section 2.1). For the dissimilarity measure D_1 , they are as follows:

$\left[\begin{array}{c} 0.495\\ 0.291\\ 0.067\\ 0.148 \end{array}\right],$	$\left[\begin{array}{c} 0.155\\ 0.114\\ 0.310\\ 0.420 \end{array}\right],$	$\begin{bmatrix} 0.260\\ 0.102\\ 0.049\\ 0.589 \end{bmatrix},$	$\begin{bmatrix} 0.511\\ 0.054\\ 0.180\\ 0.254 \end{bmatrix}.$
[0.148]	0.420	0.589	$\left[\begin{array}{c} 0.254 \end{array} \right]$

On the other hand, for index D_3 :

$\left[\begin{array}{c} 0.495\\ 0.291 \end{array}\right]$	$\left[\begin{array}{c} 0.109\\ 0.089\end{array}\right]$	$\left[\begin{array}{c} 0.403\\ 0.105\end{array}\right]$	$\left[\begin{array}{c} 0.368\\ 0.065\end{array}\right]$
$\left[\begin{array}{c} 0.067\\ 0.148 \end{array}\right],$	$\left[\begin{array}{c} 0.284\\ 0.517 \end{array}\right]'$	$\left[\begin{array}{c} 0.339\\ 0.153 \end{array}\right]'$	$\left[\begin{array}{c} 0.113\\ 0.455 \end{array}\right]^{-1}$

Against our conjecture, the second and the third summer houses do not have the highest priority in any cluster; the first and the last alternatives are the best in two clusters, respectively. Aggregating the individual pairwise comparison matrices by the geometric means of the entries (Aczél and Saaty, 1983) and applying LLSM to the common matrix results in the following priority vector:

$$\begin{bmatrix} 0.410\\ 0.164\\ 0.146\\ 0.279 \end{bmatrix}.$$
(5)

The implied ranking $1 \succ 4 \succ 2 \succ 3$ differs from the ranking in any cluster centre, thus, the aggregated ranking does not correspond to the preferences of any group.

Clustering may provide an aggregation procedure if there is only one cluster, whose centre represents "best" the whole set of PCMs. The cluster centres are

$$CC_1 = \begin{bmatrix} 1 & 2.000 & 3.000 & 1.500 \\ 0.500 & 1 & 2.000 & 0.500 \\ 0.333 & 0.500 & 1 & 0.250 \\ 0.667 & 2.000 & 4.000 & 1 \end{bmatrix} \text{ and } CC_3 = \begin{bmatrix} 1 & 2.000 & 1.500 & 1.000 \\ 0.500 & 1 & 0.667 & 0.667 \\ 0.667 & 1.500 & 1 & 0.667 \\ 1.000 & 1.500 & 1.500 & 1 \end{bmatrix}$$

for measures D_1 and D_3 , respectively, and the corresponding weight vectors are

[0.381]		[0.319]	
0.185		0.166	
0.099	and	0.219	,
0.334		0.296	

which lead to the rankings $1 \succ 4 \succ 2 \succ 3$ and $1 \succ 4 \succ 3 \succ 2$, respectively. The rankings are identical to the ranking derived from the weights (5) associated with the aggregated matrix, except for a rank reversal at the bottom in the case of dissimilarity measure D_3 . On the other hand, the relative weights of the first and the last alternatives are closer to each other according to both clusters than according to the result obtained by a reasonable aggregation of the individual matrices given in (5).

3.3 The appropriate number of clusters

There is no unique or standard procedure to determine the appropriate number of clusters. Nonetheless, a popular choice is the so-called "elbow" method: the sum of distances from the cluster centres (the objective function of the LP in Section 2.4) is plotted for all relevant numbers of clusters, and the value for which the decreasing trend flattens is chosen. The corresponding chart for sample S4 is shown in Figure 4, suggesting that k = 4 or k = 5 clusters is the best option as a further increase in k does not lead to a substantial reduction in the optimal value of the objective function.

A more sophisticated version of the "elbow" method is the gap statistic (Tibshirani et al., 2001). Although it "is designed to be applicable to any clustering method and distance measure" (Tibshirani et al., 2001, p. 411), during the implementation uniformly distributed samples are generated in a multivariate Euclidean space. Therefore, this method does not seem to be suitable for our problem.

Several other approaches are available to obtain the optimal number of clusters but the majority of them are based on the sum of squares. Although the variance could be calculated, we think it has no common meaning for the dissimilarity measures considered.

One exception is the silhouette method (Kaufman and Rousseeuw, 1990), which uses the dissimilarity matrix Δ but does not rely on any specific property of Euclidean spaces.

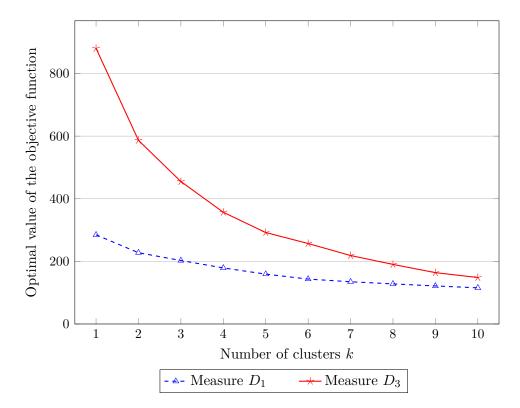


Figure 4: The optimum as a function of the number of clusters, sample S4

Number of clusters	Measure D_1	Measure D_3
2	0.286	0.467
3	0.285	0.407
4	0.260	0.503
5	0.310	0.510
6	0.293	0.494
7	0.302	0.477
8	0.321	0.488
9	0.337	0.535
10	0.345	0.568
11	0.348	0.556
12	0.325	0.571
13	0.325	0.553
14	0.330	0.545
15	0.327	0.538
16	0.324	0.514
17	0.308	0.512
18	0.310	0.523
19	0.306	0.521
20	0.310	0.518

Table 3: Mean silhouette values, sample S4

The average silhouette values are given in Table 3. They tend to grow for small values of k, the maximum is reached at 11 (12) clusters in the case of index D_1 (D_3). However,

these numbers seem to be unreasonably high.

Another common method is using a hierarchical clustering method and choosing the number of clusters based on dendrograms, which are presented in Figure 5. Note that Figure 5.a has been derived by the square of D_1 to make the structure more visible. The dendrogram shows the level of dissimilarity at which the corresponding objects are merged into the same cluster. According to these dendrograms, the number of clusters k should be between three and six in our case.

To conclude, we currently could not recommend any method that immediately gives the number of clusters. The problem is somewhat analogous to the choice of the dissimilarity measure, where there is no perfect solution, too.

4 Conclusions

This paper has provided a new perspective on group decision making, especially large-scale group decision making, by proposing the k-medoids clustering for pairwise comparison matrices. Our method has some advantages over other solutions:

- it is independent of the specification of the weighting method;
- it can handle incomplete data such that the impact of incomplete pairwise comparison matrices, which contain less information, is inherently reduced;
- it is more robust to outliers than the *k*-means clustering algorithm;
- the cluster centres are guaranteed to be individual pairwise comparison matrices, making them easier to accept by the decision-makers;
- its LP formulation allows for adding various restrictions, for example, regarding the inconsistency of the cluster centres.

The suggested approach has been used to analyse the experimental data of Bozóki et al. (2013), which demonstrated that it (i) can provide an alternative aggregation technique by choosing one cluster; (b) is able to automatically detect outliers.

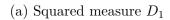
Hopefully, all practitioners dealing with data from large-scale group decision-making may benefit from using the clustering model presented here. Further investigations are especially welcome because, at the moment, there are few results on choosing the dissimilarity measure underlying the k-medoids algorithm or the appropriate number of clusters k.

Acknowledgements

The research was supported by the National Research, Development and Innovation Office under Grants FK 145838 and TKP2021-NKTA-01 NRDIO.

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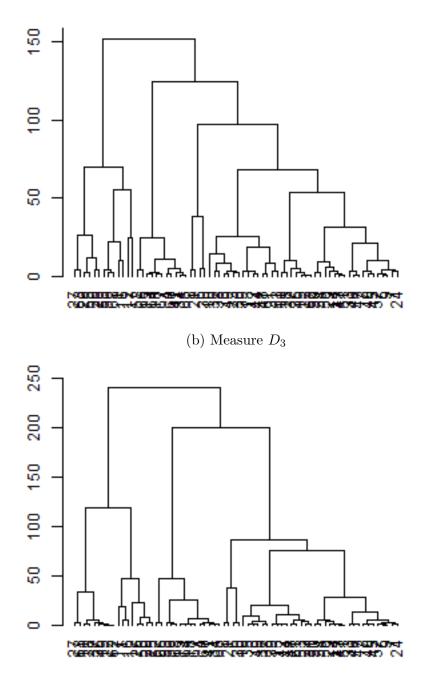


Figure 5: Dendrograms, sample S4

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