

# Supplementary materials to TITLE

Martin Papenberg

## Standard algorithm for anticlustering

To assign samples to batches, we use anticlustering. Anticlustering is an optimization method that is characterized by (a) an objective function that quantifies the balance among batches, and (b) an algorithm that conducts the batch assignment in such a way that balance among batches is maximized via the objective function. Anticlustering owes its name to the fact that the objective functions it uses are the reversal of criteria used in cluster analysis. For example, Späth (1986) already recognized that by maximizing instead of minimizing the k-means criterion (the “variance”), he was able to create groups that are similar to each other, and presented it as an improvement over the more intuitive random assignment.

In our application, we optimize the (a) diversity objective using (b) the exchange algorithm by Papenberg and Klau (2021). This constellation of objective and algorithm is also the default method implemented in the R package `anticlust` (Papenberg 2019). The diversity is defined as the total sum the within-cluster sums of dissimilarities (Brusco, Cradit, and Steinley 2020). When assigning samples in such a way that the diversity is maximized, we simultaneously maximize similarity between batches. In Papenberg and Klau (2021), we referred to the maximization of the diversity as “anticluster editing” because the minimization of the diversity is also well-known from the area of cluster analysis—under the term “cluster editing” (Böcker, Briesemeister, and Klau 2011; Shamir, Sharan, and Tsur 2004).

The anticlustering exchange algorithm takes as input (1) a distance matrix, quantifying the pairwise dissimilarity among samples, and (2) the number of batches  $K$ . As measure of dissimilarity, we use the common Euclidean distance, which transfers the features of our samples to pairwise distances. It is defined as

$$d(x, y) = \sqrt{\sum_{i=1}^M (x_i - y_i)^2}$$

where  $M$  is the number of numeric features describing two samples  $x = (x_1, \dots, x_M)$  and  $y = (y_1, \dots, y_M)$ . When samples are described by two features, the Euclidean distance corresponds to the geometric, “straightline” distance between two points in a two-dimensional space; more similar items are closer to each other (see Figure 1). For categorical variables, we use binary (or „dummy“) coding before computing the Euclidean distance matrix. **example! best if related to the current application**

The exchange algorithm consists of two steps: an initialization step and an optimization step. As initialization, it randomly assigns samples to  $K$  equal-sized batches. In principle, unequal-sized batches would also be possible with our algorithm, but equal-sized batches were required in the current application. After initialization, the algorithm selects the first sample and checks how the diversity would change if the sample were swapped with each sample that is currently assigned to a different batch. After simulating each exchange—that is  $N - \frac{N}{K}$  exchanges—it realizes the one exchange that increases the diversity the most. It does not conduct an exchange if no improvement in diversity is possible. This procedure is repeated for each sample, leading to the evaluation of  $N \cdot \frac{N}{K}$  exchanges in total; it terminates after the last sample was processed.<sup>1</sup>

<sup>1</sup>The procedure might also restart at the first element and reiterate through all samples until no exchange leads to an improvement any more, i.e., until a local maximum is found. In `anticlust`, we also implemented this local maximum search, which corresponds to the algorithm LCW by Weitz and Lakshminarayanan (1998). However, the default method in `anticlust` is the exchange method that terminates after a fixed number of exchanges.

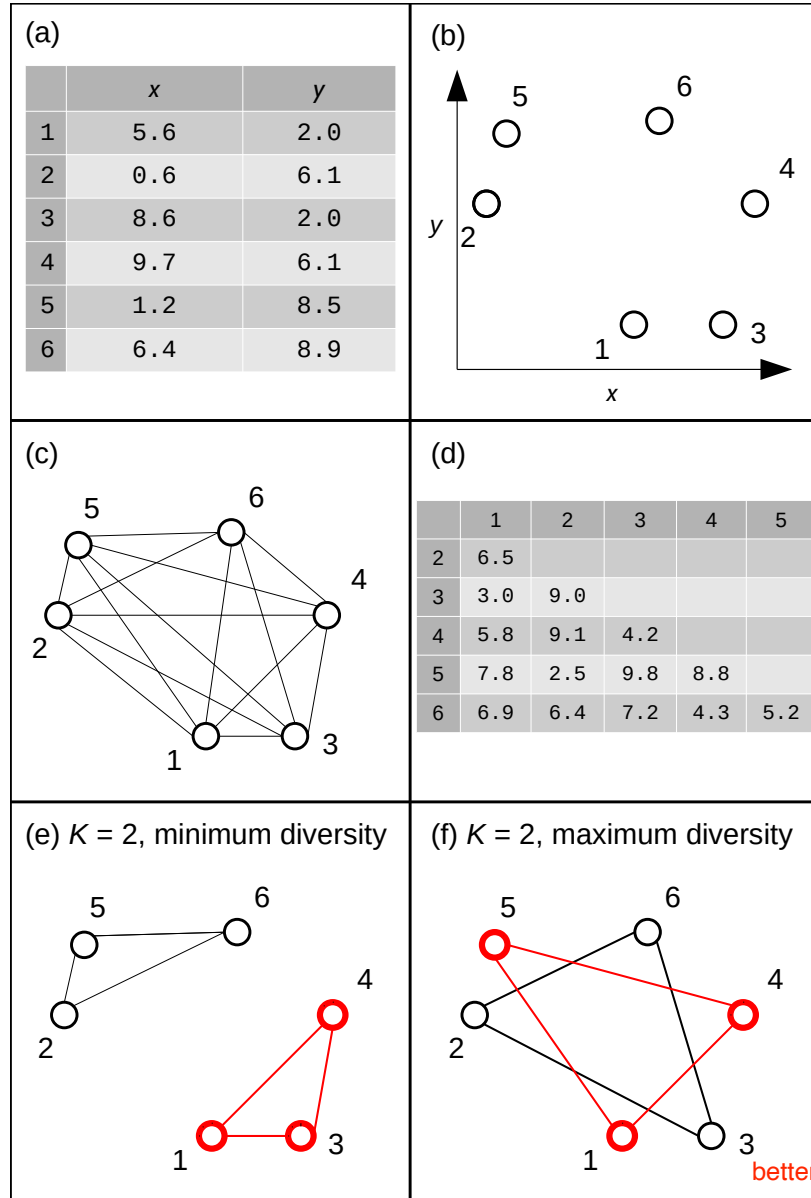


Figure 1: Illustrates the conversion from numeric features to Euclidean distance, and (anti)clustering assignments based on minimum and maximum diversity using the Euclidean distance. Panel (a) shows the numeric values on two hypothetical features  $x$  and  $y$  for six items; panel (b) shows their representation in the two-dimensional space; panel (c) represents the Euclidean distances between items as straight lines in the two-dimensional space; panel (d) illustrates the Euclidean distances in a dissimilarity matrix; these values are proportional to the length of the connecting lines in panel (c). Panel (e) illustrates a clustering assignment of the 6 data points to  $K = 2$  equal-sized groups via minimum diversity. Panel (f) illustrates an anticlustering assignment of 6 data points to  $K = 2$  equal-sized groups via maximum diversity. This Figure is an adaptation of an illustration of the Preprint (Version 1) at <https://doi.org/10.31234/osf.io/3razc>, and it is published with a CC-BY Attribution 4.0 International license.

## Including must-link constraints with anticlustering

To include must-link constraints with anticlustering, we adjusted both the initialization phase as well as the optimization phase of the exchange algorithm. For initialization, we recognize that we now may have to assign multiple samples simultaneously (“must-link groups”) to a batch—not all samples can be considered independently of all other samples. Therefore, we first collect all samples that have at least one other sample (“must-link partner”) that has to be assigned to the same group.

- We obtain a list of all must-link groups together and store them with their cardinality.
- We must assign these must-link groups to the  $K$  batches.
- This corresponds to a bin packing problem!
  - We have  $n$  elements –  $n$  = number of must-link groups – where each element has a size, corresponding to the number of elements in that must-link group.
  - Each bin corresponds to a batch and has target capacity of  $\frac{N}{K}$
  - We use “randomized fit” heuristic to fill the batches: Select first batches that fits, and try out batches in random order (leads to rather even distribution in all batches). A “worst fit” heuristic might also be used, but some random element is good if we restart our optimization heuristic multiple times for better results.
- The remaining elements (that are not part of must-link groups) can be assigned randomly to batches (while respecting the cardinality constraints / equal-sized batches) – they are not part of the bin packing problem.

### Optimization algorithm:

- “Merge” samples: A must-link group is treated as a single unit during optimization.
- New distance matrix has to be computed: New distances are given by summing up all distances between all elements that are part of must-link groups (this works for the diversity objective because it is a mere sum)
- Exchange heuristic is conducted on the merged elements. We only exchange must-link groups of the same size to ensure that the cardinality constraints are still respected in the end (equal-sized groups).

## References

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