

# Hierarchical clustering for gene expression data analysis

*Giorgio Valentini*

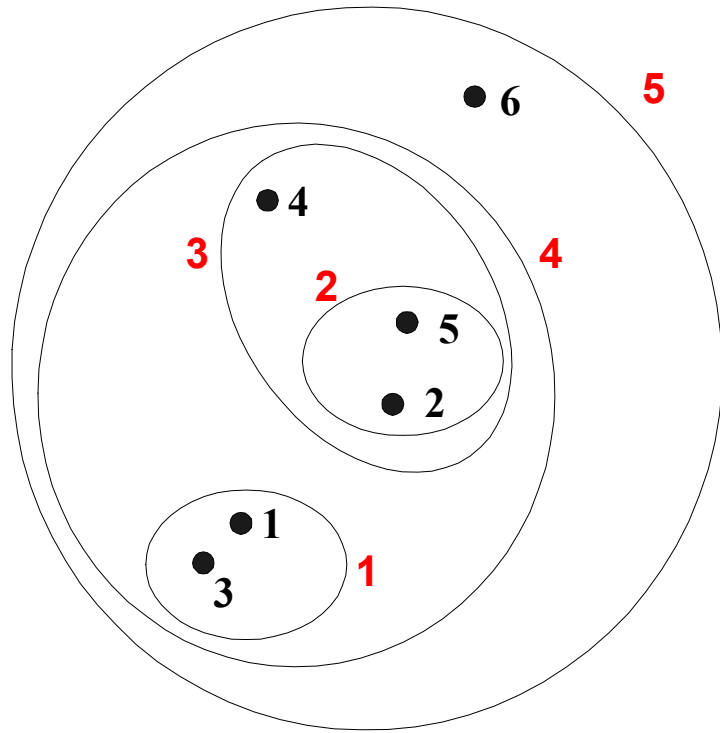
e-mail: [valentini@dsi.unimi.it](mailto:valentini@dsi.unimi.it)

**Dipartimento di Scienze dell'Informazione**  
Università degli Studi di Milano

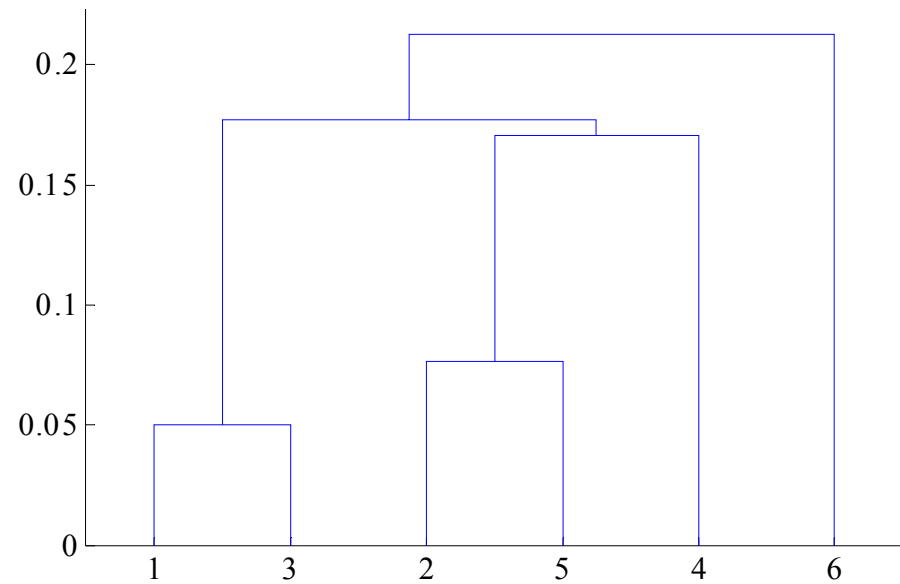
# Clustering of Microarray Data

1. *Clustering of gene expression profiles* (rows) => discovery of co-regulated and functionally related genes (or unrelated genes: different clusters)
2. *Clustering of samples* (columns) => identification of sub-types of related samples
3. *Two-way clustering* => combined sample clustering with gene clustering to identify which genes are the most important for sample clustering

# Hierarchical Clustering

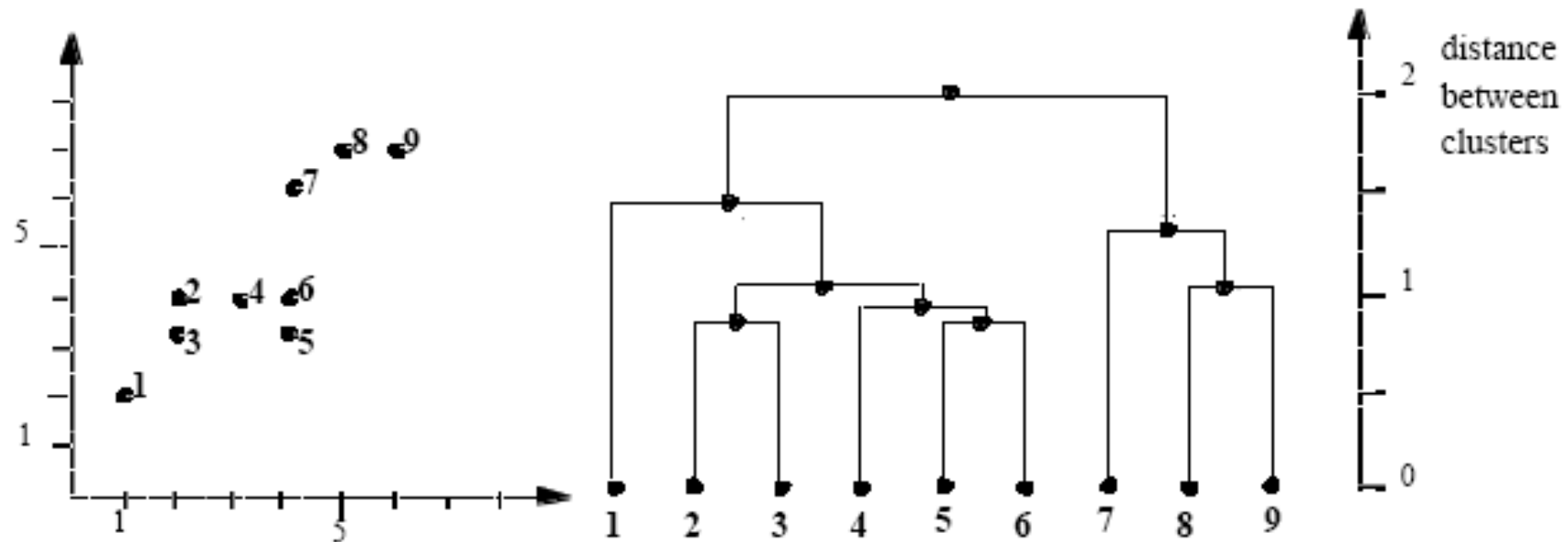


**Hierarchical Clustering**



**Dendrogram**

# Dendrograms



- The *root* represents the whole data set
- A *leaf* represents a single object in the data set
- An *internal node* represent the union of all objects in its sub-tree
- The *height* of an internal node represents the distance between its two child nodes

# Hierarchical Clustering

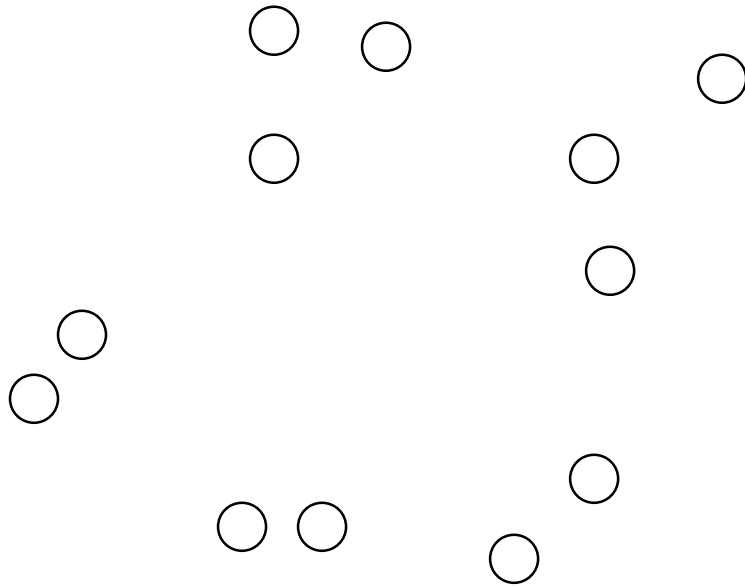
- Two main types of hierarchical clustering.
  - **Agglomerative:**
    - Start with the points as individual clusters
    - At each step, merge the closest pair of clusters.
    - Until only one cluster (or  $k$  clusters) left
    - This requires defining the notion of cluster proximity.
  - **Divisive:**
    - Start with one, all-inclusive cluster
    - At each step, split a cluster
    - Until each cluster contains a point (or there are  $k$  clusters)
    - Need to decide which cluster to split at each step.

# Basic Agglomerative Hierarchical Clustering Algorithm

1. Initially, each object forms its own cluster
  2. Compute all pairwise distances between the initial clusters (objects)
- repeat**
3. Merge the closest pair (A, B) in the set of the current clusters into a new cluster  $C = A \cup B$
  4. Remove A and B from the set of current clusters; insert C into the set of current clusters
  5. Determine the distance between the new cluster C and all other clusters in the set of current clusters
- until** only a single cluster remains

# Agglomerative Hierarchical Clustering: Starting Situation

- For agglomerative hierarchical clustering we start with clusters of individual points and a proximity matrix.

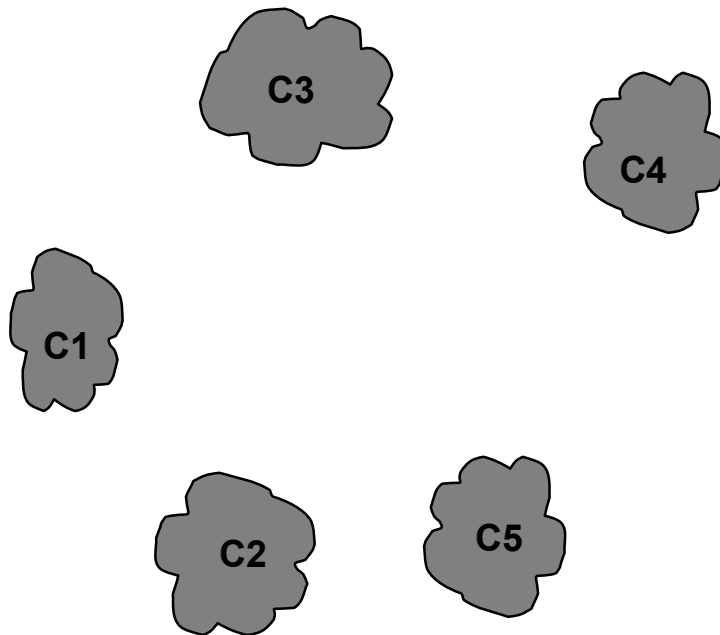


	p1	p2	p3	p4	p5	...
p1						
p2						
p3						
p4						
p5						
.						
.						
.						

**Proximity Matrix**

# Agglomerative Hierarchical Clustering: Intermediate Situation

- After some merging steps, we have some clusters.



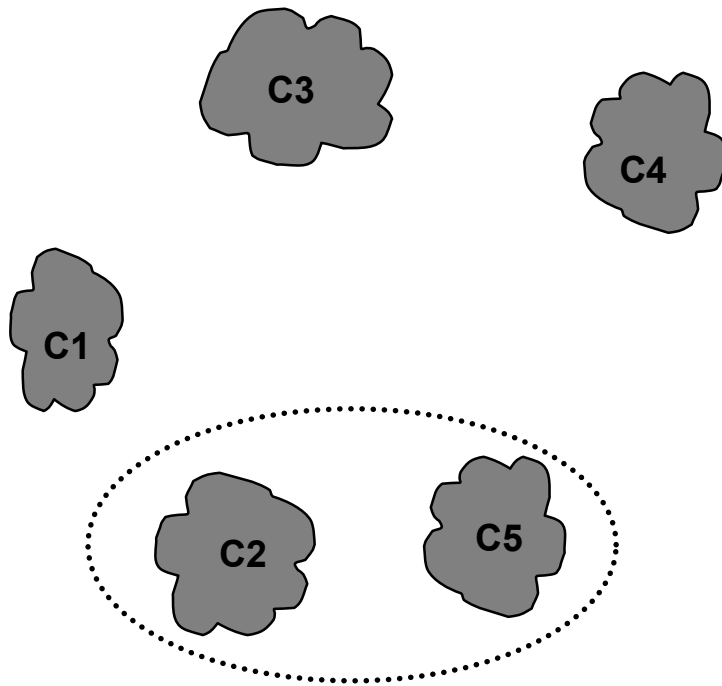
	C1	C2	C3	C4	C5
C1					
C2					
C3					
C4					
C5					

**Proximity Matrix**



# Agglomerative Hierarchical Clustering: Intermediate Situation

- We want to merge the two closest clusters (C2 and C5) and update the proximity matrix.

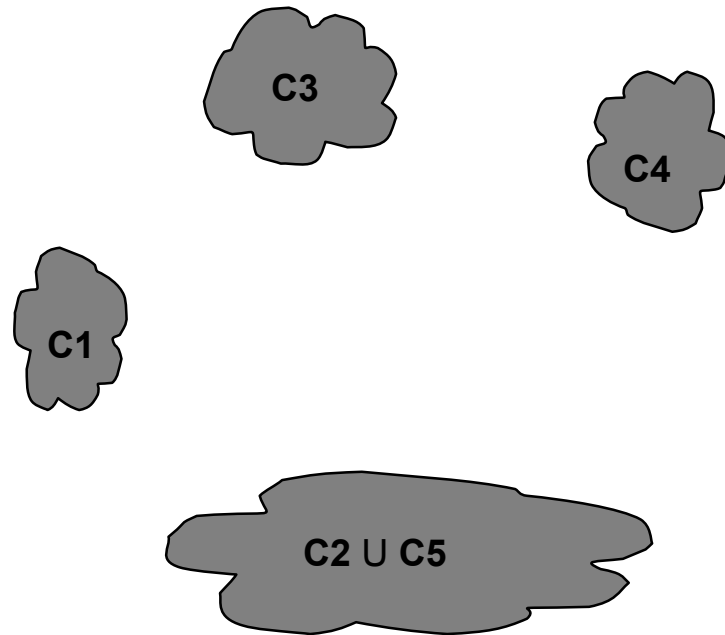


	C1	C2	C3	C4	C5
C1					
C2					
C3					
C4					
C5					

**Proximity Matrix**

# Agglomerative Hierarchical Clustering: after Merging

- The question is “How do we update the proximity matrix?”



	C1	$\begin{matrix} \text{C2} \\ \cup \\ \text{C5} \end{matrix}$	C3	C4
C1		?		
$\text{C2} \cup \text{C5}$	?	?	?	?
C3		?		
C4		?		

**Distance Matrix**

- Key operation is the computation of the distance of two clusters.
- Different approaches to defining the distance between clusters distinguishes the different algorithms

# Inter-cluster distances

- Four widely used ways of defining the **inter-cluster distance**, i.e., the distance between two separate clusters  $C_i$  and  $C_j$ , are

- o **single linkage method** (nearest neighbor):

$$d(C_i, C_j) = \min_{x \in C_i, y \in C_j} \{ d(x, y) \}$$

- o **complete linkage method** (furthest neighbor):

$$d(C_i, C_j) = \max_{x \in C_i, y \in C_j} \{ d(x, y) \}$$

- o **average linkage method** (unweighted pair-group average):

$$d(C_i, C_j) = \text{avg}_{x \in C_i, y \in C_j} \{ d(x, y) \}$$

- o **centroid linkage method** (distance between cluster centroids  $c_i$  and  $c_j$ ):

$$d(C_i, C_j) = d(c_i, c_j)$$

# Single linkage (minimum distance) method

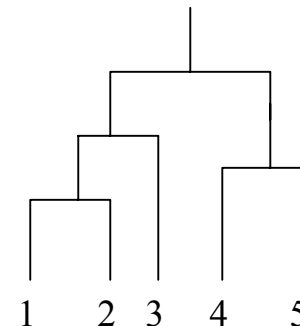
- Distance (dissimilarity) of two clusters is based on the two most similar (closest) points in the different clusters  $C_i$  and  $C_j$  :

$$d(C_i, C_j) = \min_{x \in C_i, y \in C_j} \{ d(x, y) \}$$

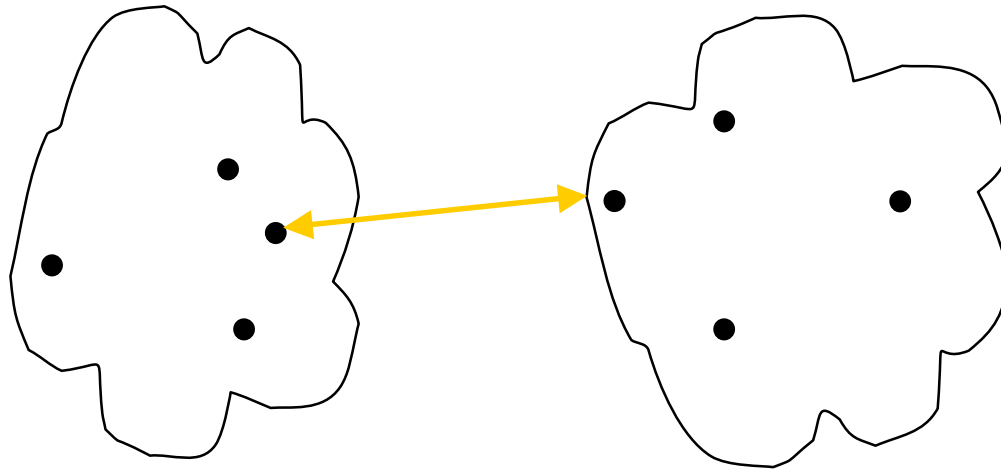
- Determined by one pair of points, i.e., by one link in the proximity graph.
- Can handle non-elliptical shapes.
- Sensitive to noise and outliers.

Similarity matrix

	I1	I2	I3	I4	I5
I1	1.00	0.90	0.10	0.65	0.20
I2	0.90	1.00	0.70	0.60	0.50
I3	0.10	0.70	1.00	0.40	0.30
I4	0.65	0.60	0.40	1.00	0.80
I5	0.20	0.50	0.30	0.80	1.00

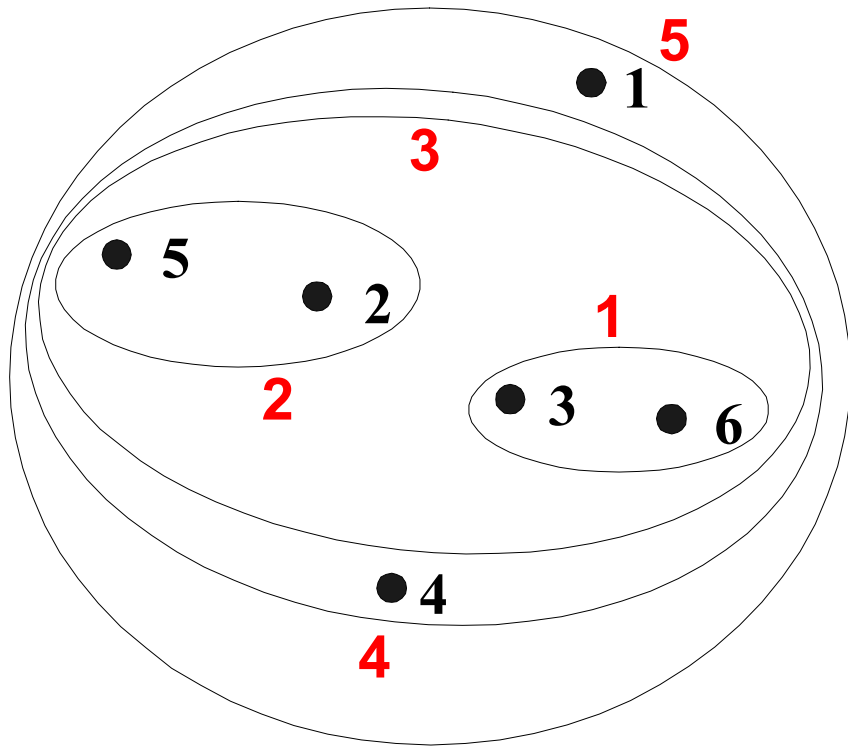


# Single linkage

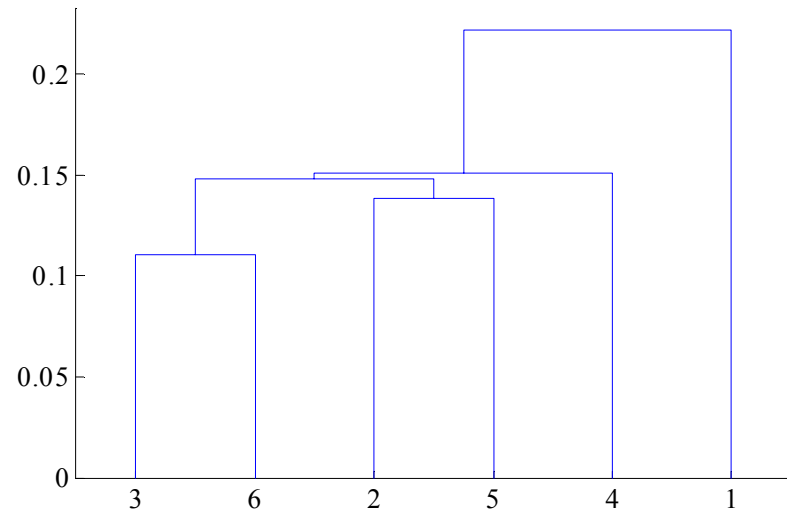


$$d(C_i, C_j) = \min_{x \in C_i, y \in C_j} \{ d(x, y) \}$$

# Hierarchical Clustering: minimum distance

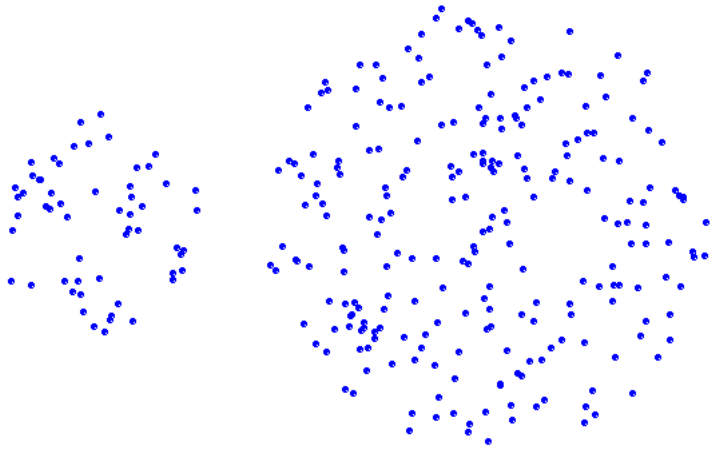


**Nested Clusters**

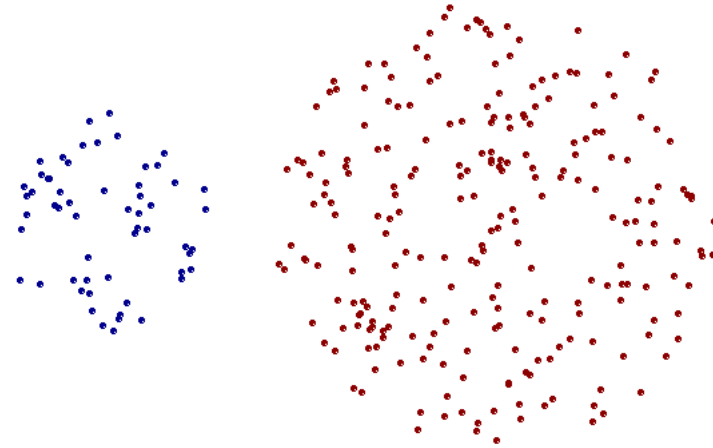


**Dendrogram**

# Strength of minimum distance

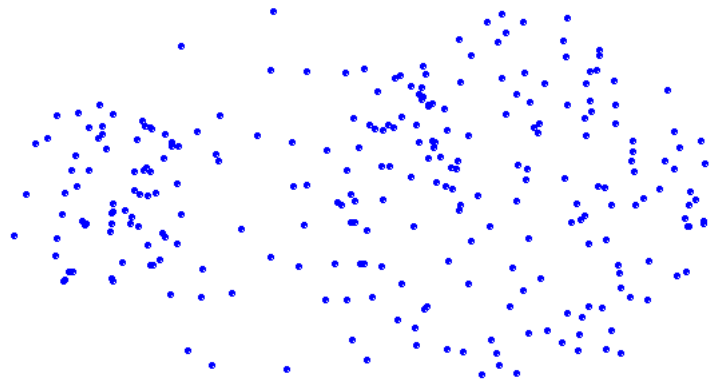


**Original Points**

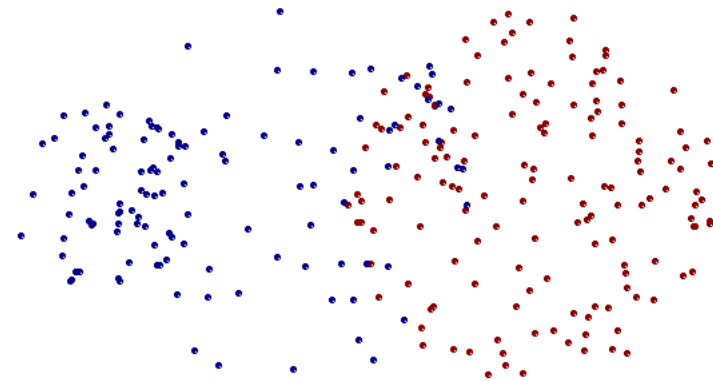


**Two Clusters**

# Limitation of minimum distance



**Original Points**



**Two Clusters**



# Complete Linkage (minimum distance) method

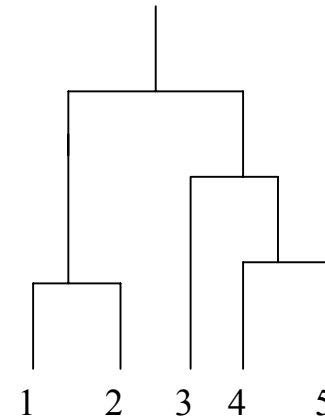
- Distance of two clusters is based on the two least similar (most distant) points in the different clusters  $C_i$  and  $C_j$ :

$$d(C_i, C_j) = \max_{x \in C_i, y \in C_j} \{ d(x, y) \}$$

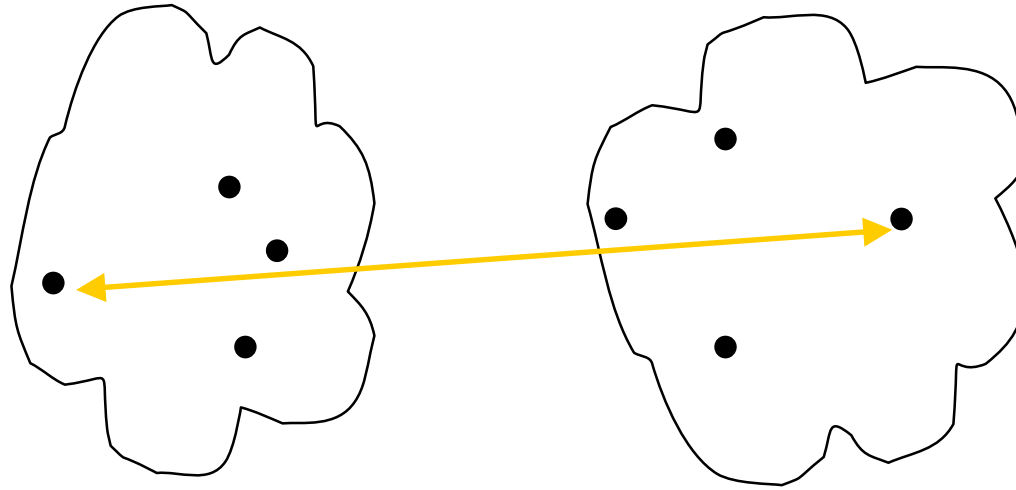
- Determined by all pairs of points in the two clusters.
- Tends to break large clusters.
- Less susceptible to noise and outliers.

Similarity matrix

	I1	I2	I3	I4	I5
I1	1.00	0.90	0.10	0.65	0.20
I2	0.90	1.00	0.70	0.60	0.50
I3	0.10	0.70	1.00	0.40	0.30
I4	0.65	0.60	0.40	1.00	0.80
I5	0.20	0.50	0.30	0.80	1.00



# Complete linkage

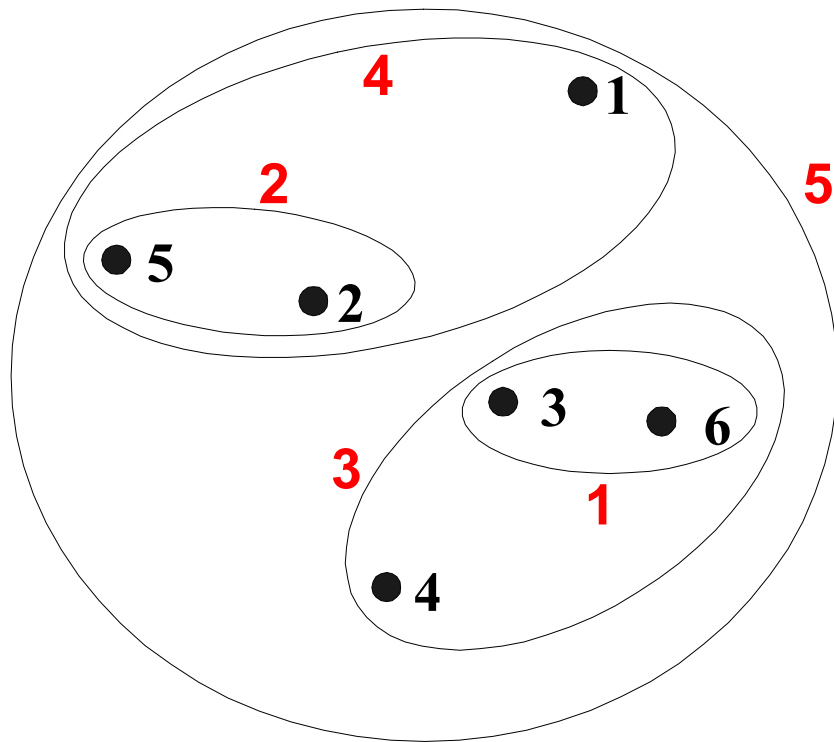


$$d(C_i, C_j) = \max_{x \in C_i, y \in C_j} \{ d(x, y) \}$$

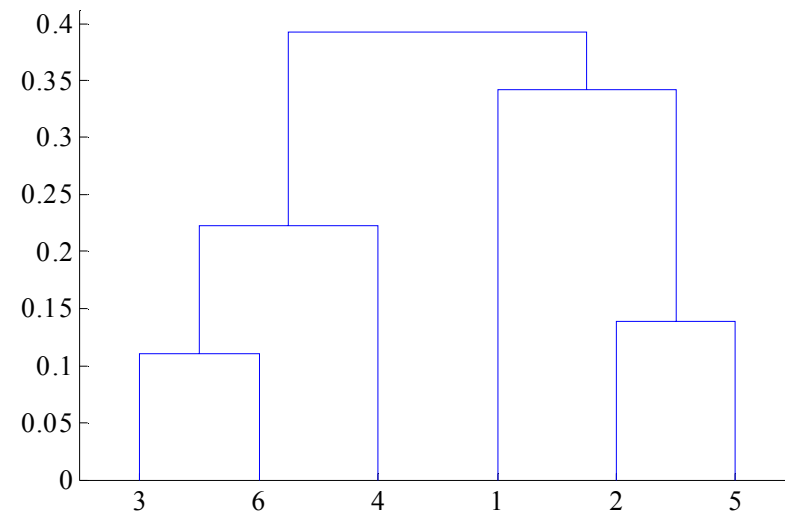
## Cluster Similarity: maximum distance or Complete Linkage

- Similarity of two clusters is based on the two most distant points in the different clusters.
- Tends to break large clusters.
- Less susceptible to noise and outliers.
- Biased towards globular clusters.

# Hierarchical Clustering: maximum distance

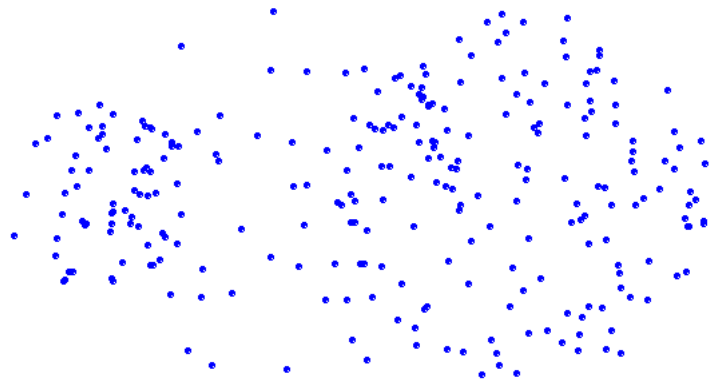


**Nested Clusters**

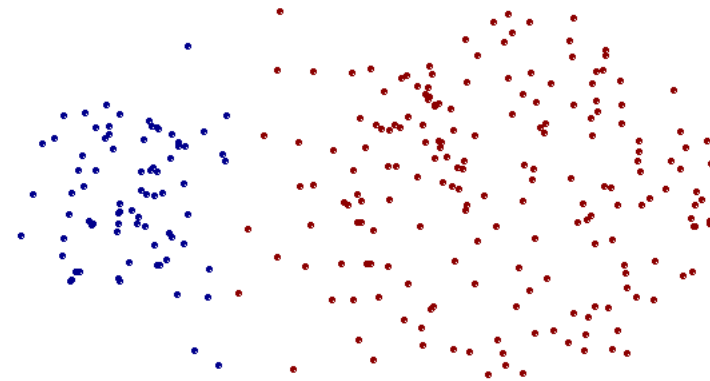


**Dendrogram**

# Strength of maximum distance

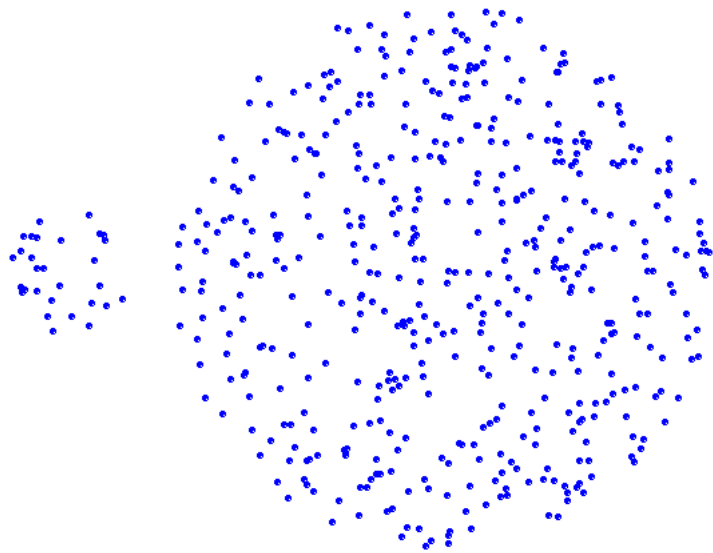


**Original Points**

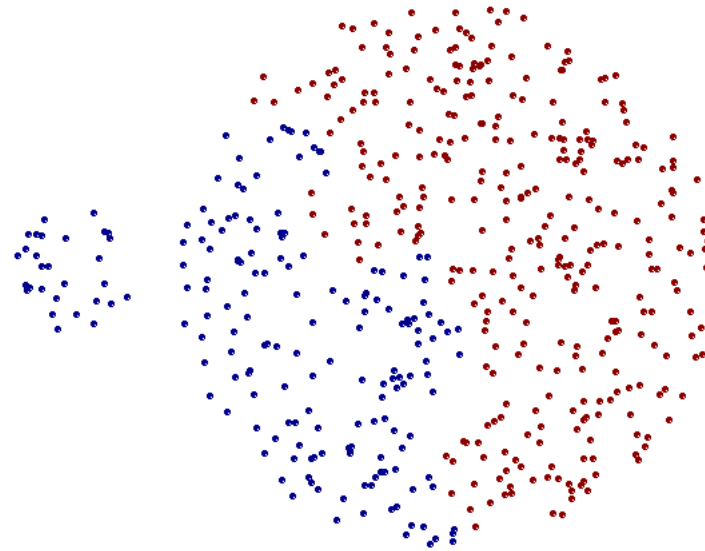


**Two Clusters**

# Limitations of maximum distance



**Original Points**



**Two Clusters**

# Average linkage (average distance) method

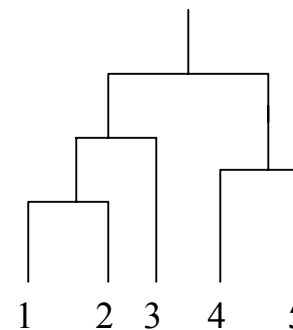
- Distance of two clusters is the average of pairwise distances between points in the two clusters  $C_i$  and  $C_j$ :

$$d(C_i, C_j) = \frac{1}{|C_i| |C_j|} \sum_{x \in C_i} \sum_{y \in C_j} d(x, y)$$

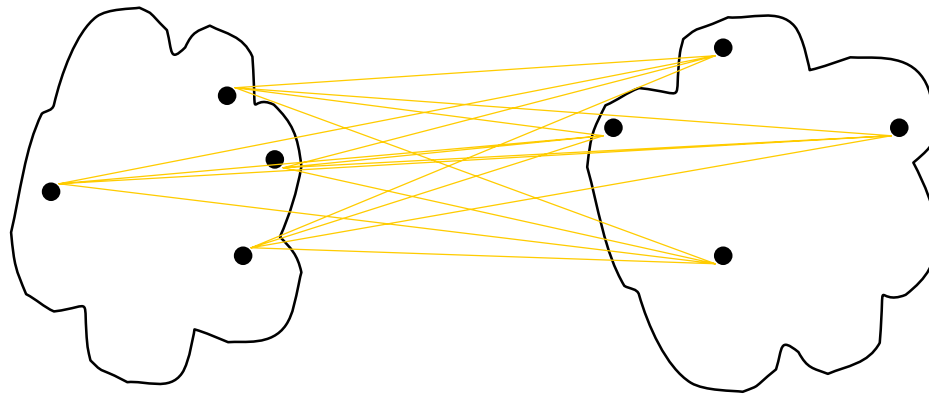
- Compromise between Single and Complete Link.
- Need to use average connectivity for scalability since total connectivity favors large clusters.
- Less susceptible to noise and outliers.
- Biased towards globular clusters.

Similarity matrix

	I1	I2	I3	I4	I5
I1	1.00	0.90	0.10	0.65	0.20
I2	0.90	1.00	0.70	0.60	0.50
I3	0.10	0.70	1.00	0.40	0.30
I4	0.65	0.60	0.40	1.00	0.80
I5	0.20	0.50	0.30	0.80	1.00



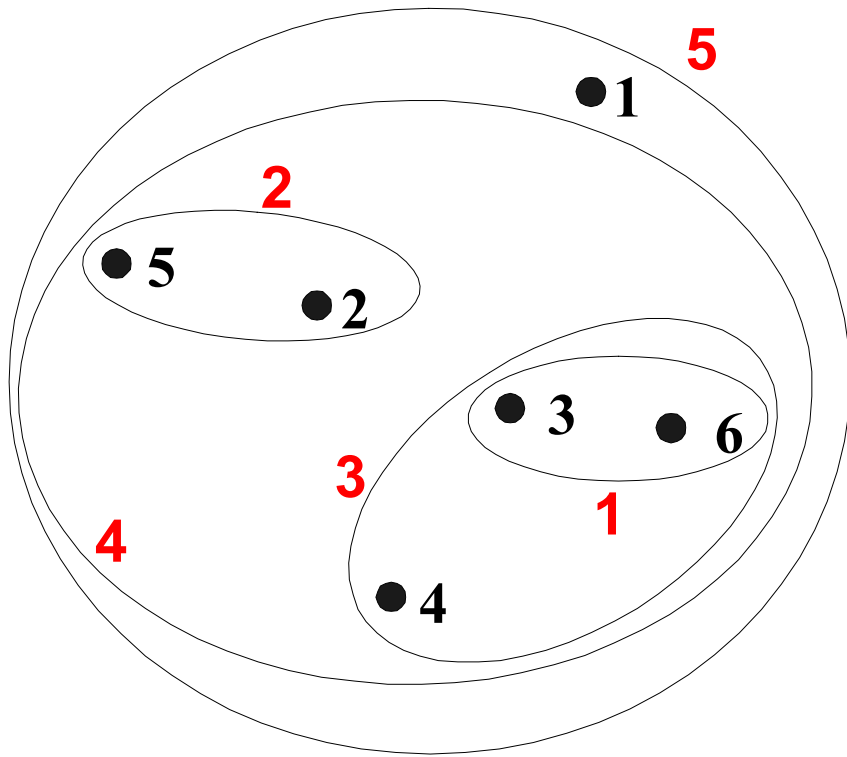
# Average linkage



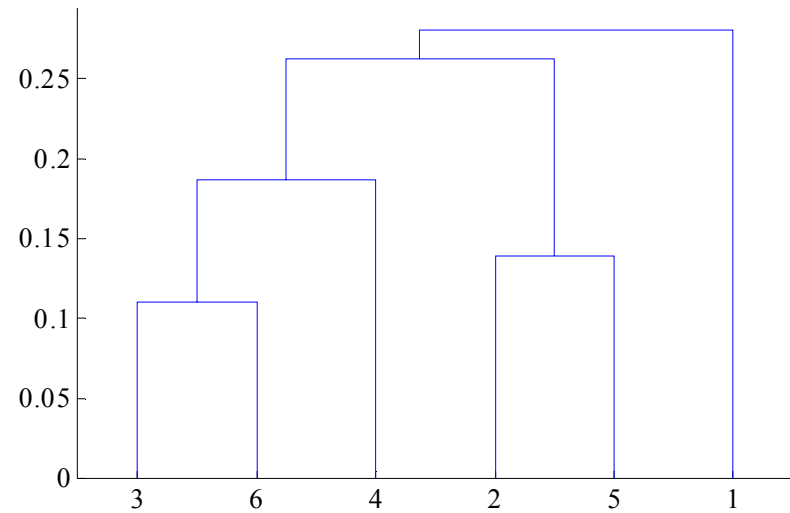
$$d(C_i, C_j) = \frac{1}{|C_i| |C_j|} \sum_{x \in C_i} \sum_{y \in C_j} d(x, y)$$



# Hierarchical Clustering: Average distance



**Nested Clusters**



**Dendrogram**

# Centroid linkage (centroid distance) method

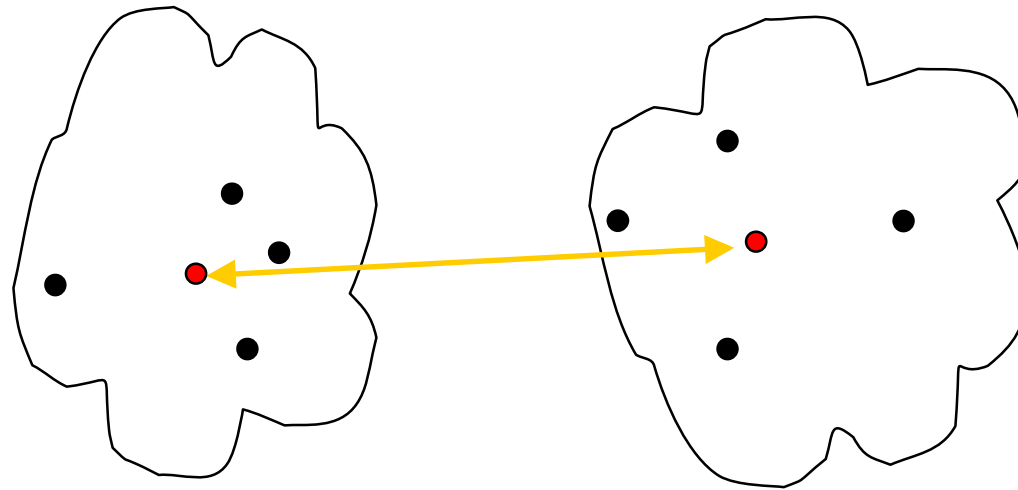
- Distance of two clusters is distance of the two centroids  $c_i$  and  $c_j$  of the two clusters  $C_i$  and  $C_j$ :

$$d(C_i, C_j) = d(c_i, c_j)$$

$$c_i = \frac{1}{|C_i|} \sum_{x \in C_i} x \quad c_j = \frac{1}{|C_j|} \sum_{x \in C_j} x$$

- Compromise between Single and Complete Link.
- Less computationally intensive with respect to average linkage.

# Centroid linkage



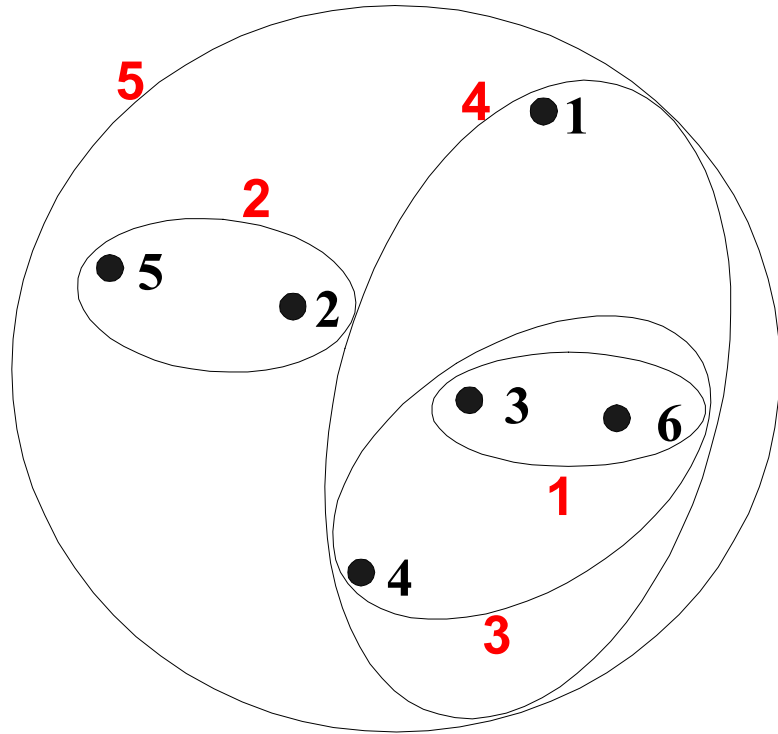
$$d(C_i, C_j) = d(c_i, c_j)$$

$$c_i = \frac{1}{|C_i|} \sum_{x \in C_i} x \quad c_j = \frac{1}{|C_j|} \sum_{x \in C_j} x$$

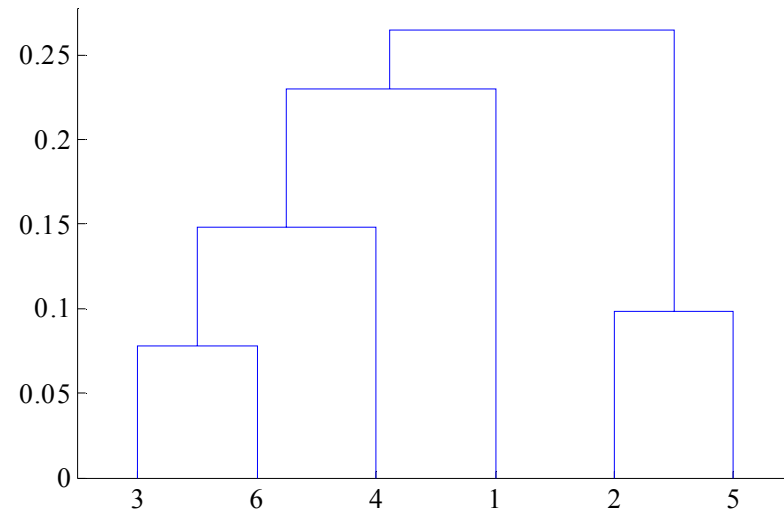
# Cluster Similarity: Ward's Method

- Similarity of two clusters is based on the increase in squared error when two clusters are merged.
  - Similar to group average if distance between points is distance squared.
- Less susceptible to noise and outliers.
- Biased towards globular clusters.
- Hierarchical analogue of K-means
  - But Ward's method does not correspond to a local minimum
  - Can be used to initialize K-means

# Hierarchical Clustering: Ward's method

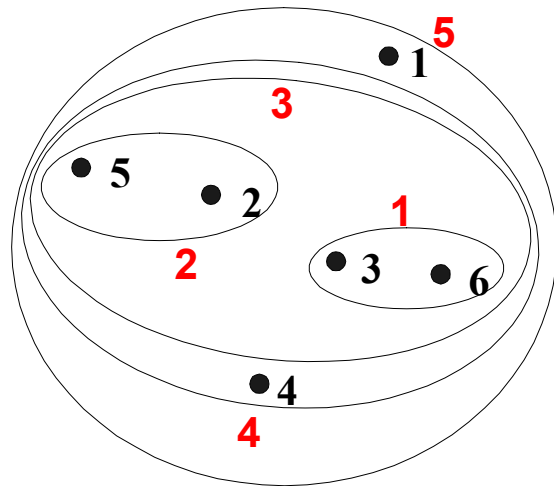


**Nested Clusters**

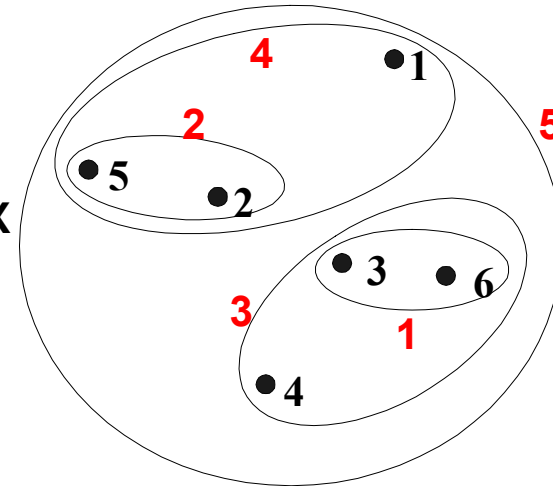


**Dendrogram**

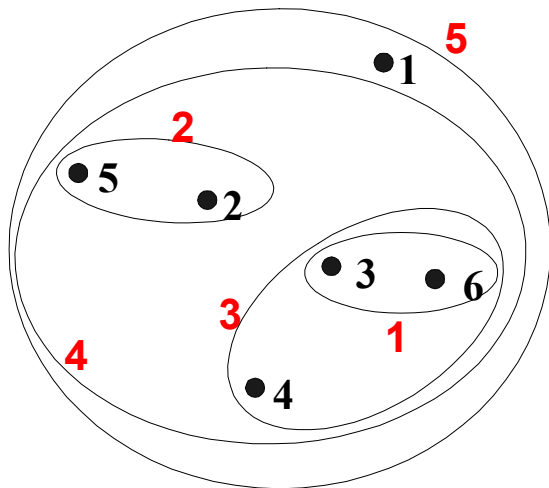
# Hierarchical Clustering: comparison



MIN

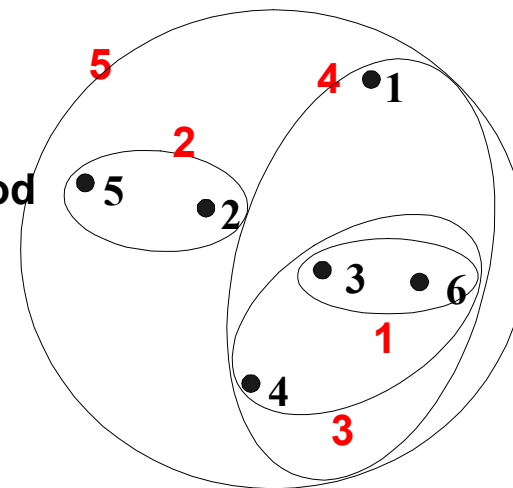


MAX



Average

Ward's Method



# Comparison of minimum, maximum, average and centroid distance

## Minimum distance

- When  $d_{\min}$  is used to measure distance between clusters, the algorithm is called the nearest-neighbor or single-linkage clustering algorithm
- If the algorithm is allowed to run until only one cluster remains, the result is a minimum spanning tree (MST)
- This algorithm favors elongated classes

## Maximum distance

- When  $d_{\max}$  is used to measure distance between clusters, the algorithm is called the farthest-neighbor or complete-linkage clustering algorithm
- From a graph-theoretic point of view, each cluster constitutes a complete sub-graph
- This algorithm favors compact classes

## Average and centroid distance

- The minimum and maximum distance are extremely sensitive to outliers since their measurement of between-cluster distance involves minima or maxima
- The average and centroid distance approaches are more robust to outliers
- Of the two, the centroid distance is computationally more attractive
- Notice that the average distance approach involves the computation of  $|C_i||C_j|$  distances for each pair of clusters

## Hierarchical Clustering: Time and Space requirements

- $O(N^2)$  space since it uses the proximity matrix.
  - $N$  is the number of points.
- $O(N^3)$  time in many cases.
  - There are  $N$  steps and at each step the size,  $N^2$ , proximity matrix must be updated and searched.
  - By being careful, the complexity can be reduced to  $O(N^2 \log(N))$  time for some approaches.



## Hierarchical Clustering: problems and limitations

- Once a decision is made to combine two clusters, it cannot be undone.
- No objective function is directly minimized.
- Different schemes have problems with one or more of the following:
  - Sensitivity to noise and outliers.
  - Difficulty handling different sized clusters and convex shapes.
  - Breaking large clusters.

# Advantages and disadvantages of Hierarchical clustering

## Advantages

- Does not require the number of clusters to be known in advance
- No input parameters (besides the choice of the (dis)similarity)
- Computes a complete hierarchy of clusters
- Good result visualizations integrated into the methods

## Disadvantages

- May not scale well: runtime for the standard methods:  $O(n^2 \log n)$
- No explicit clusters: a “flat” partition can be derived afterwards (e.g. via a cut through the dendrogram or termination condition in the construction)
- No automatic discovering of “optimal clusters”

# Hierarchical clustering of tissues and genes:

Alizadeh et al. 2000, Distinct types of diffuse large B-cell lymphoma identified by gene expression profiling, Nature 403:3.

