

# Cluster Analysis: Basic Concepts and Algorithms

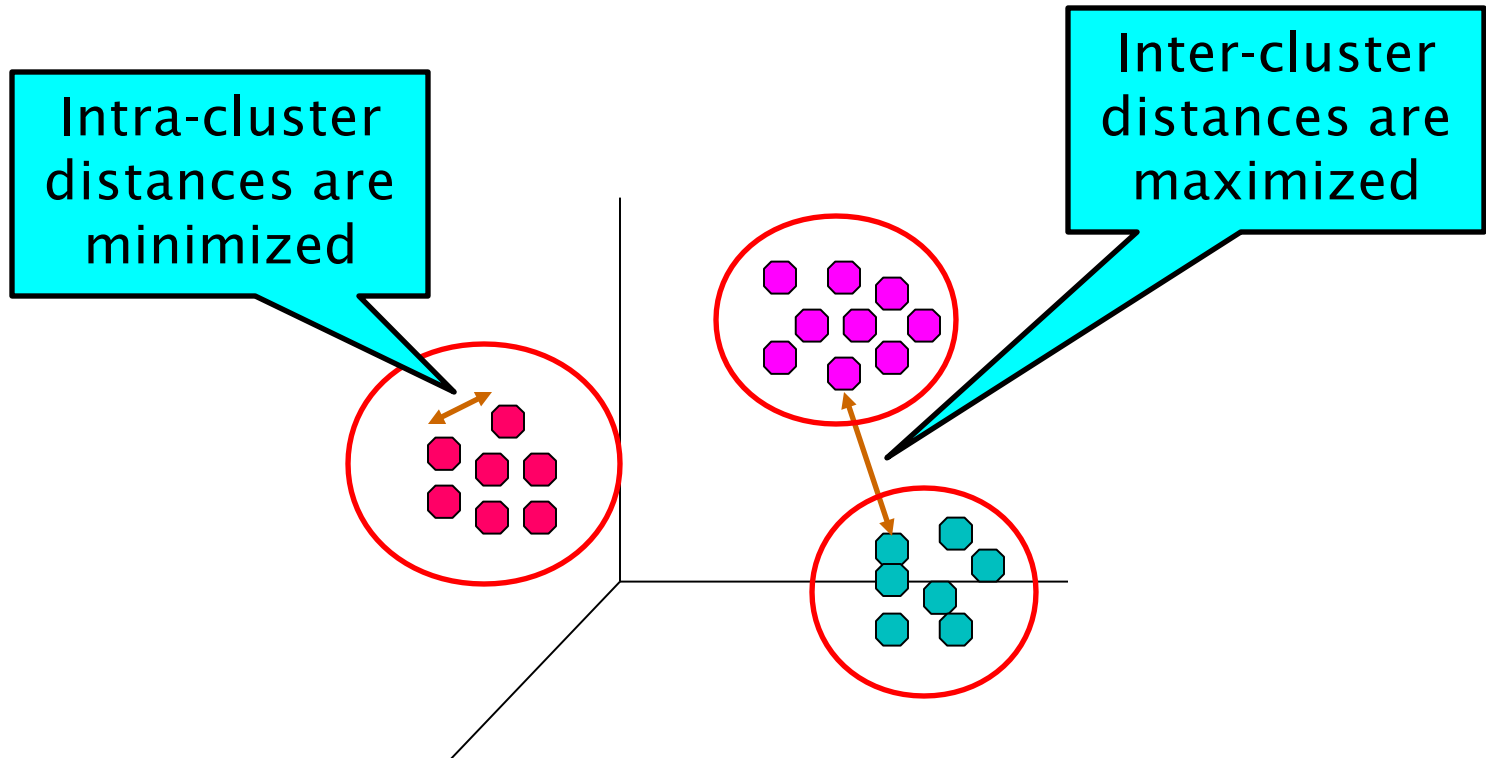


# Topics

- Introduction
- Types of Clustering
- Types of Clusters
- Clustering Algorithms
  - K-Means Clustering
  - Hierarchical Clustering
  - Density-based Clustering
- Cluster Validation

# What is Cluster Analysis?

- Finding groups of objects such that the objects in a group will be similar (or related) to one another and different from (or unrelated to) the objects in other groups



# Applications of Cluster Analysis

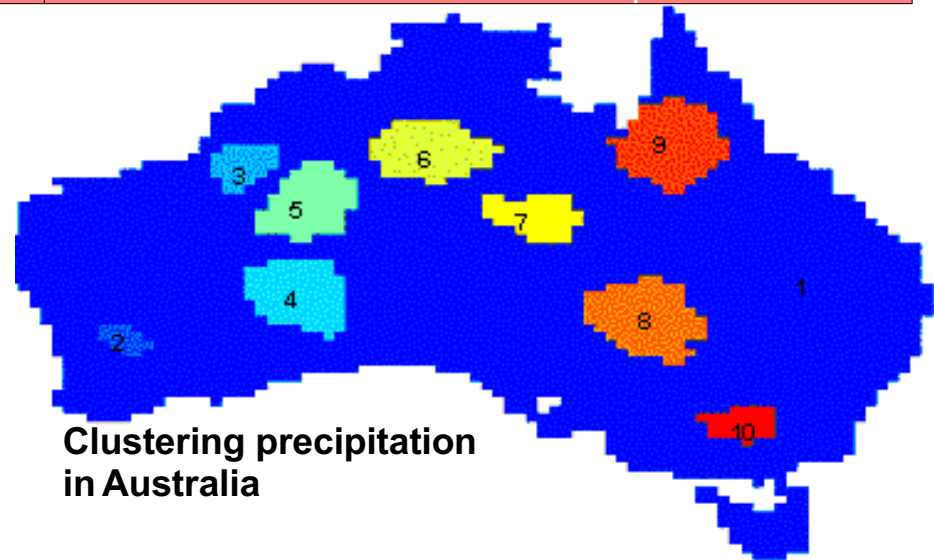
- **Understanding**

- Group related documents for browsing, group genes and proteins that have similar functionality, or group stocks with similar price fluctuations

	<i>Discovered Clusters</i>	<i>Industry Group</i>
<b>1</b>	Applied-Matl-DOWN,Bay-Network-DOWN,3-COM-DOWN,Cabletron-Sys-DOWN,CISCO-DOWN,HP-DOWN,DSC-Comm-DOWN,INTEL-DOWN,LSI-Logic-DOWN,Micron-Tech-DOWN,Texas-Inst-Down,Tellabs-Inc-Down, Natl-Semiconduct-DOWN,OracI-DOWN,SGI-DOWN, Sun-DOWN	Technology1-DOWN
<b>2</b>	Apple-Comp-DOWN,Autodesk-DOWN,DEC-DOWN,ADV-Micro-Device-DOWN,Andrew-Corp-DOWN,Computer-Assoc-DOWN,Circuit-City-DOWN,Compaq-DOWN,EMC-Corp-DOWN,Gen-Inst-DOWN,Motorola-DOWN,Microsoft-DOWN,Scientific-Atl-DOWN	Technology2-DOWN
<b>3</b>	Fannie-Mae-DOWN,Fed-Home-Loan-DOWN,MBNA-Corp-DOWN,Morgan-Stanley-DOWN	Financial-DOWN
<b>4</b>	Baker-Hughes-UP,Dresser-Inds-UP,Halliburton-HLD-UP,Louisiana-L and-UP,Phillips-Petro-UP,Unocal-UP, Schlumberger-UP	Oil-UP

- **Summarization**

- Reduce the size of large data sets



# What is not Cluster Analysis?

- Supervised classification
  - Uses class label information
- Simple segmentation
  - Dividing students into different registration groups alphabetically, by last name
- Results of a query
  - Groupings are a result of an external specification

→ **Clustering uses only the data**

# Similarity

**How do we measure  
similarity/proximity/dissimilarity/distance?**

## Examples

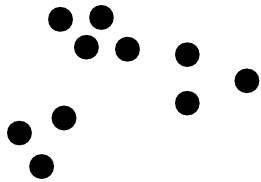
- Minkovsky distance: Manhattan distance, Euclidean Distance, etc.
- Jaccard index for binary data
- Gower's distance for mixed data (ratio/interval and nominal)
- Correlation coefficient as similarity between variables

# Notion of a Cluster can be Ambiguous

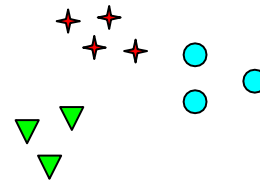
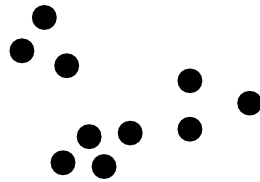


How many clusters?

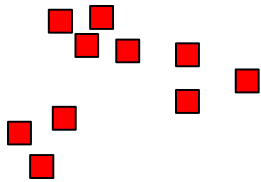
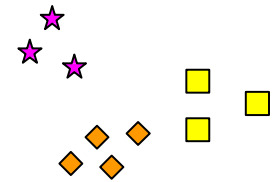
# Notion of a Cluster can be Ambiguous



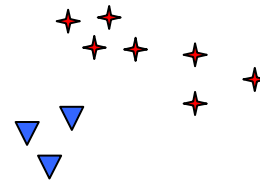
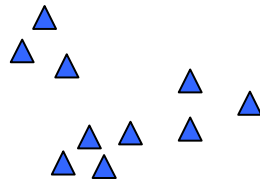
How many clusters?



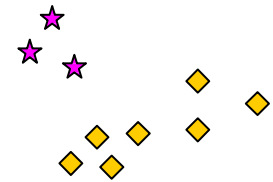
Six Clusters



Two Clusters



Four Clusters





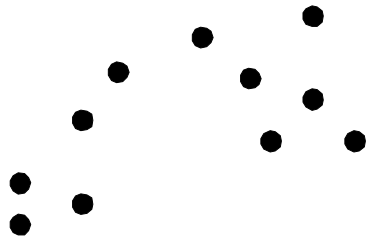
# Topics

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  - K-Means Clustering
  - Hierarchical Clustering
  - Density-based Clustering
- Cluster Validation

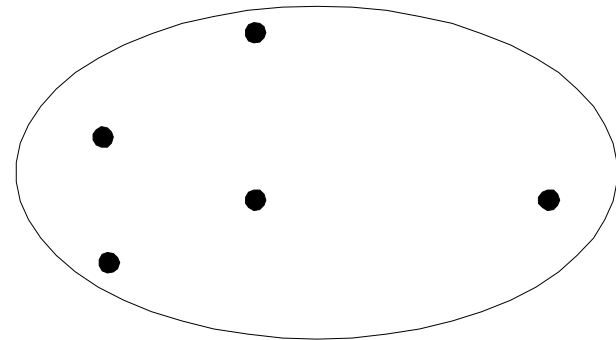
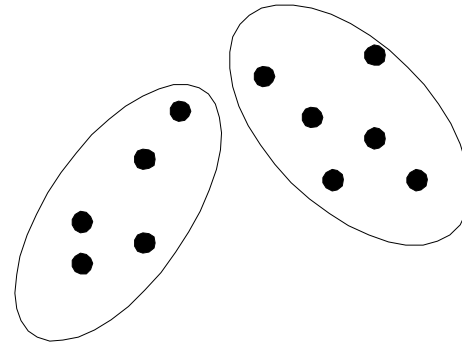
# Types of Clusterings

- A clustering is a **set of clusters**
- **Partitional Clustering**
  - A division data objects into non-overlapping subsets (clusters) such that each data object is in exactly one subset
- **Hierarchical clustering**
  - A set of nested clusters organized as a hierarchical tree

# Partitional Clustering

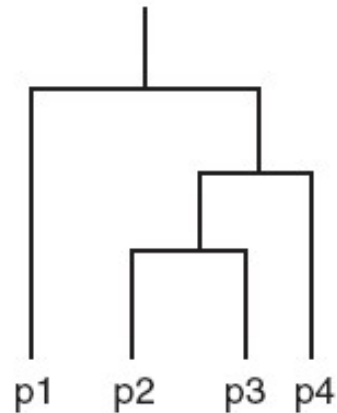


Original Points

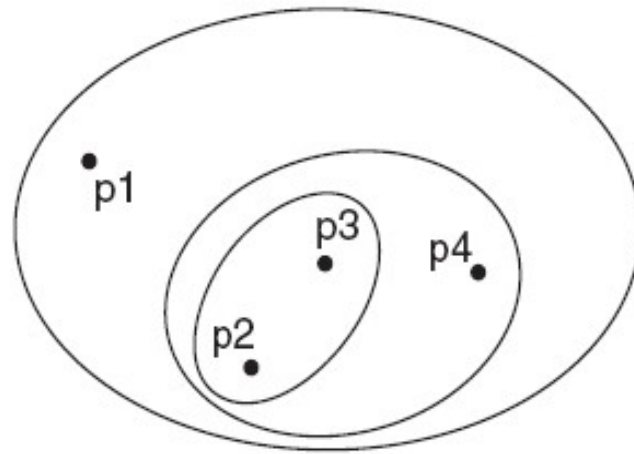


A Partitional Clustering

# Hierarchical Clustering



(a) Dendrogram.



(b) Nested cluster diagram.

**Figure 8.13.** A hierarchical clustering of four points shown as a dendrogram and as nested clusters.

# Other Distinctions Between Sets of Clusters

- **Exclusive versus non-exclusive**
  - In non-exclusive clusterings, points may belong to multiple clusters.
- **Fuzzy versus non-fuzzy**
  - In fuzzy clustering, a point belongs to every cluster with some membership weight between 0 and 1
  - Membership weights must sum to 1
  - Probabilistic clustering has similar characteristics
- **Partial versus complete**
  - In some cases, we only want to cluster some of the data
- **Heterogeneous versus homogeneous**
  - Cluster of widely different sizes, shapes, and densities

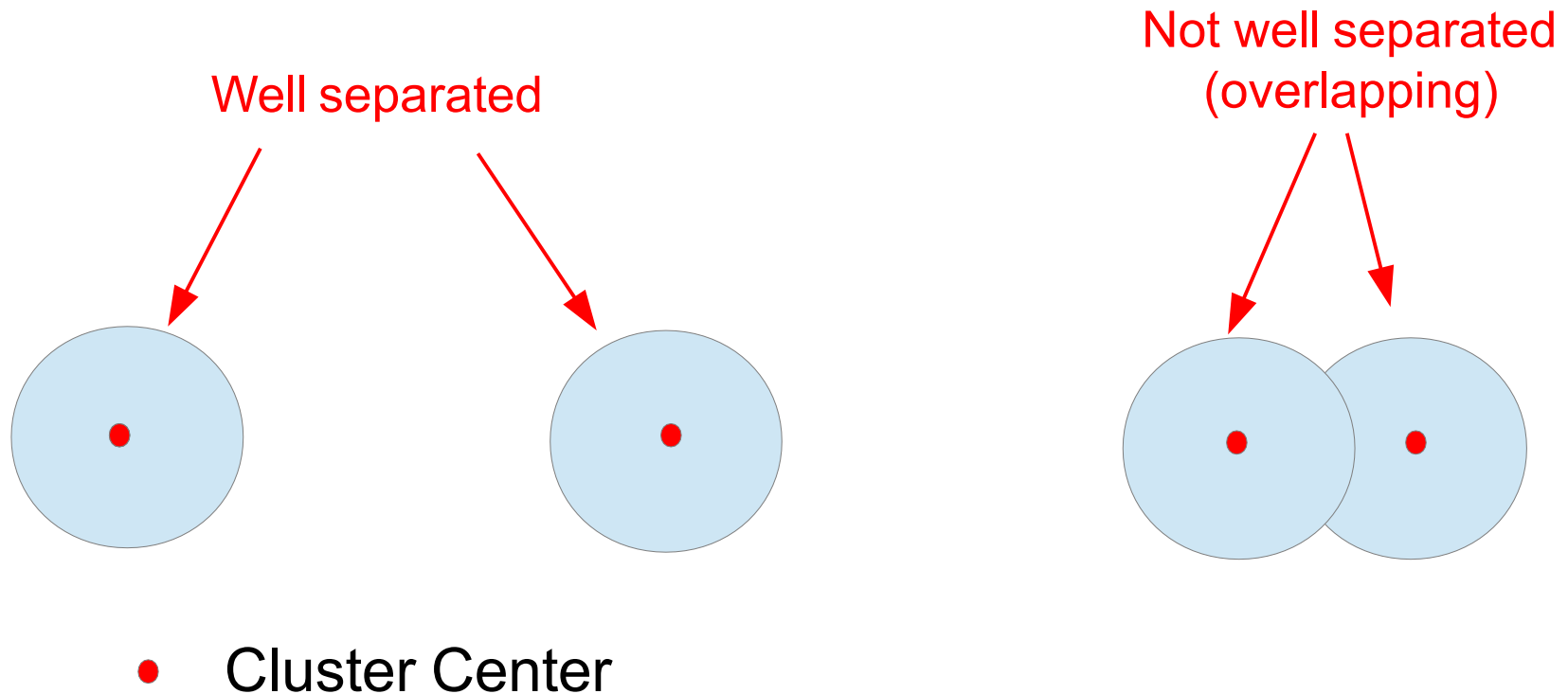
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# Types of Clusters

- Center-based clusters
- Contiguous clusters
- Density-based clusters
- Conceptual clusters
- Described by an Objective Function

# Center-based Clusters

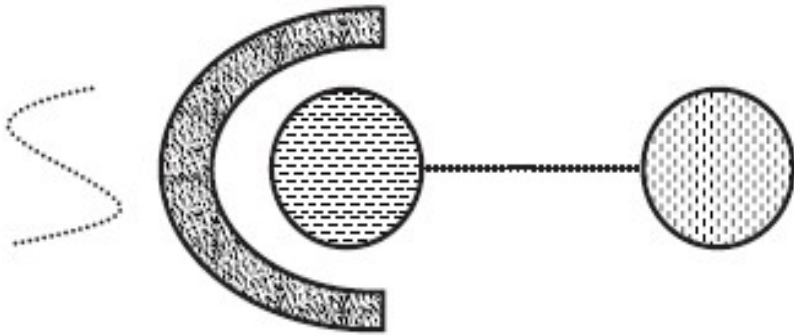


A cluster is a set of objects such that an object in a cluster is closer (more similar) to the “center” of a cluster, than to the center of any other cluster

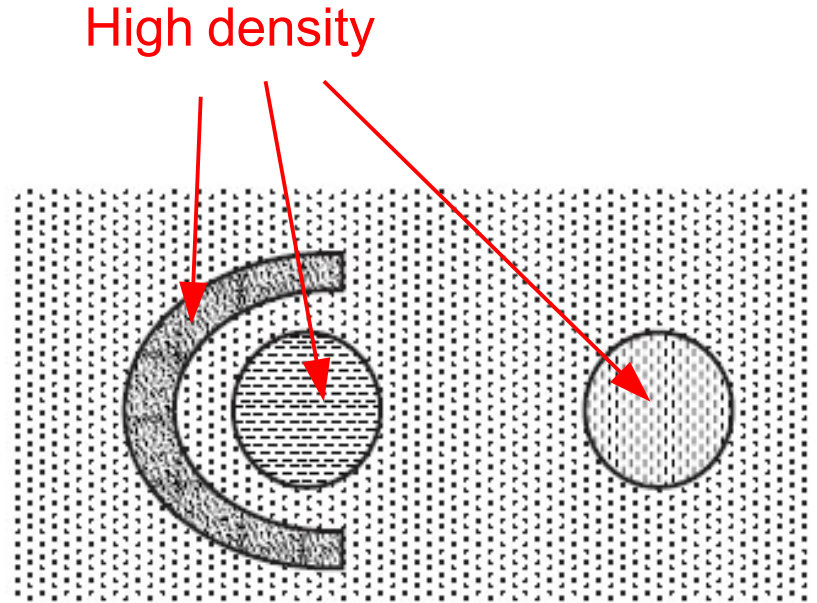
The center of a cluster is often a **centroid**, the average of all the points in the cluster, or a **medoid**, the most “representative” point of a cluster



# Contiguous and Density-based Clusters

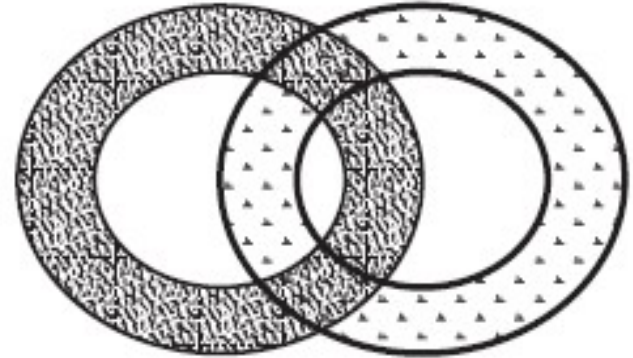
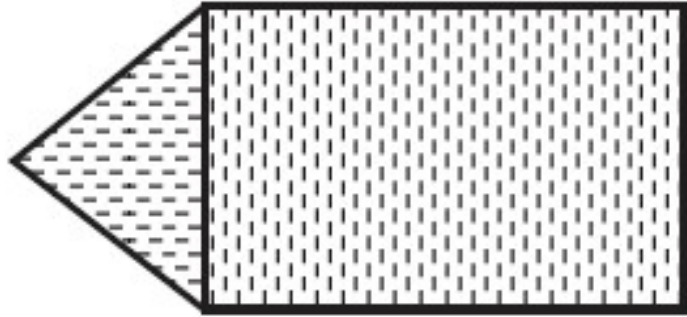


(c) Contiguity-based clusters. Each point is closer to at least one point in its cluster than to any point in another cluster.



(d) Density-based clusters. Clusters are regions of high density separated by regions of low density.

# Conceptual Clusters



(e) Conceptual clusters. Points in a cluster share some general property that derives from the entire set of points. (Points in the intersection of the circles belong to both.)

Conceptual clusters are hard to detect since they are often not:

- Center-based
- Contiguity-based
- Density-based

# Topics

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- Types of Clusters
- **Objective Functions**
- Clustering Algorithms
  - K-Means Clustering
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# Objective Functions

The best clustering minimizes or maximizes an objective function.

- Example: Minimize the Sum of Squared Errors

$$SSE = \sum_{i=1}^K \sum_{\mathbf{x} \in C_i} \|\mathbf{x} - \mathbf{m}_i\|^2$$

- $\mathbf{x}$  is a data point in cluster  $C_i$ ,  $\mathbf{m}_i$  is the center for cluster  $C_i$  as the mean of all points in the cluster and  $\|\cdot\|$  is the L2 norm (= Euclidean distance).

**Problem:** Enumerate all possible ways of dividing the points into clusters and evaluate the 'goodness' of each potential set of clusters by using the given objective function. (NP Hard)

# Objective Functions

Global objective function

- Typically used in **partitional clustering** (k-means uses SSE)
- **Mixture Models** assume that the data is a 'mixture' of a number of parametric statistical distributions (e.g., a mixture of Gaussians).

Local objective function

- **Hierarchical clustering** algorithms typically have local objectives
- **Density-based clustering** is based on local density estimates
- **Graph based approaches.** Graph partitioning and shared nearest neighbors

We will talk about the objective functions when we talk about individual clustering algorithms.

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# K-means Clustering

- **Partitional** clustering approach
- Each cluster is associated with a centroid (center point)
- Each point is assigned to the cluster with the closest centroid
- Number of clusters,  $K$ , must be specified

The basic algorithm is very simple:

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- 1: Select  $K$  points as the initial centroids.
  - 2: **repeat**
  - 3:     Form  $K$  clusters by assigning all points to the closest centroid.
  - 4:     Recompute the centroid of each cluster.
  - 5: **until** The centroids don't change
-

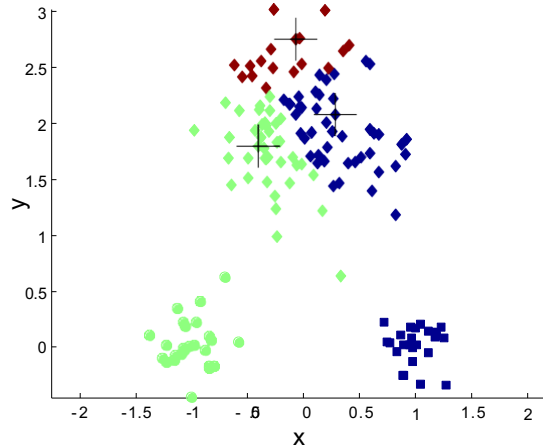
# K-means Clustering – Details

- **Initial centroids** are often chosen randomly.
  - Clusters produced vary from one run to another.
- The centroid is the mean of the points in the cluster.
- ‘Closeness’ is measured by **Euclidean distance**
- K-means will converge (points stop changing assignment) typically in the first few iterations.
  - Often the stopping condition is changed to ‘Until relatively few points change clusters’
- Complexity is  $O(n * K * I * d)$ 
  - $n$  = number of points,  $K$  = number of clusters,  
 $I$  = number of iterations,  $d$  = number of attributes

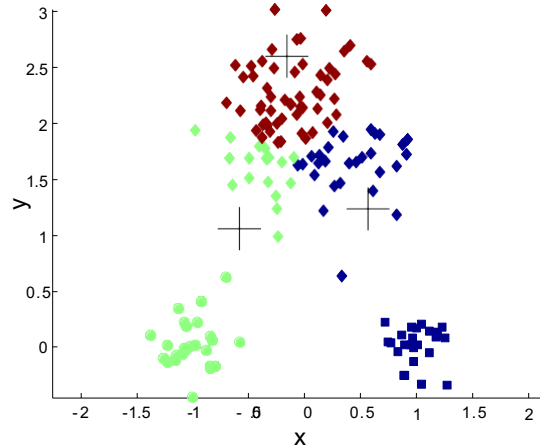


# K-Means Example

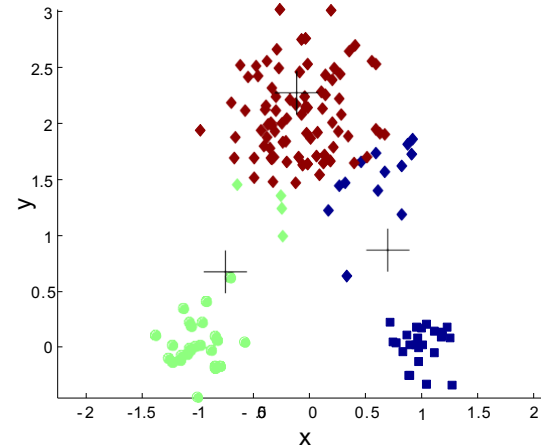
Iteration1



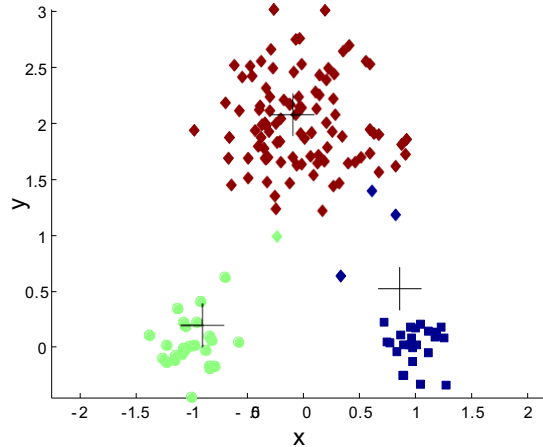
Iteration2



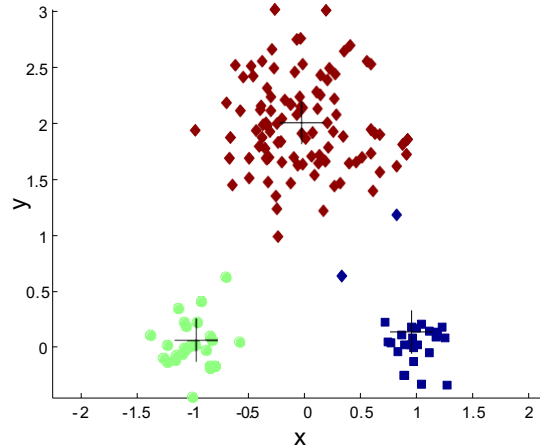
Iteration3



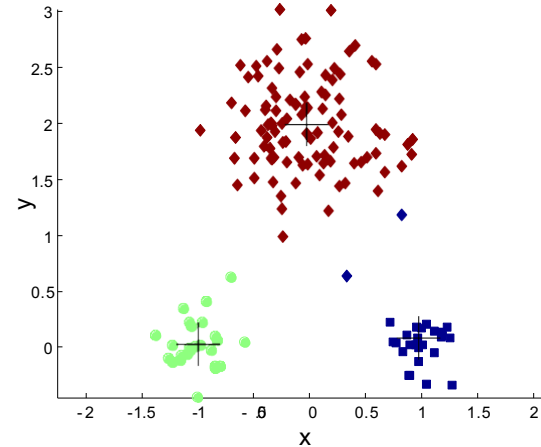
Iteration4



Iteration5



Iteration6



# Problems with Selecting Initial Points

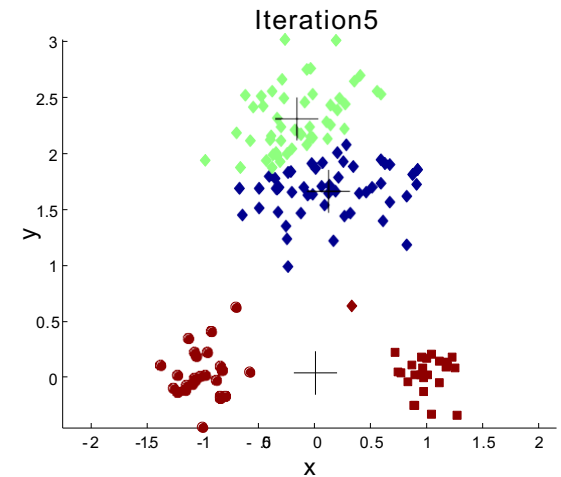
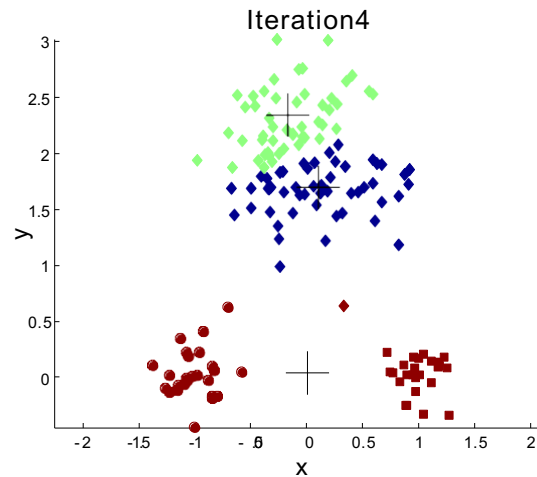
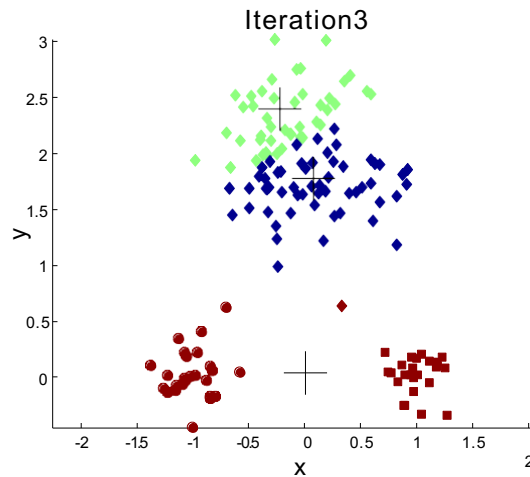
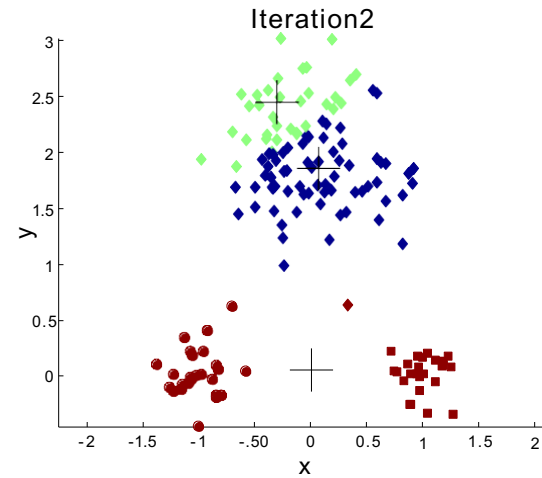
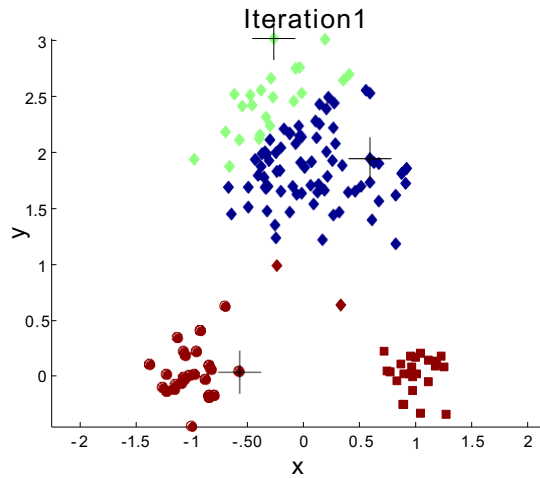
If there are  $K$  'real' clusters then the chance of selecting one centroid from each cluster is small.

- Chance is relatively small when  $K$  is large
- If clusters are the same size,  $n$ , then

$$P = \frac{\text{number of ways to select one centroid from each cluster}}{\text{number of ways to select } K \text{ centroids}} = \frac{K!n^K}{(Kn)^K} = \frac{K!}{K^K}$$

- For example, if  $K = 10$ , then probability =  $10!/10^{10} = 0.00036$
- Sometimes the initial centroids will readjust themselves in 'right' way, and sometimes they don't
- Consider an example of five pairs of clusters

# Importance of Choosing Initial Centroids ...



# Solutions to Initial Centroids Problem

- Multiple runs
- Sample and use hierarchical clustering to determine initial centroids
- Select more than  $k$  initial centroids and then select among these initial centroids (Select most widely separated)

# Evaluating K-means Clusters

- Most common measure is Sum of Squared Error (SSE)
  - For each point, the error is the distance to the nearest cluster center

$$SSE = \sum_{i=1}^K \sum_{\mathbf{x} \in C_i} \|\mathbf{x} - \mathbf{m}_i\|^2$$

- $\mathbf{x}$  is a data point in cluster  $C_i$ ,  $\mathbf{m}_i$  is the center for cluster  $C_i$  as the mean of all points in the cluster and  $\|\cdot\|$  is the L2 norm (= Euclidean distance).
  - Given two clusterings, we can choose the one with the smallest error
  - Only compare clusterings with the same K! One easy way to reduce SSE is to increase K, the number of clusters
- K-Means is a heuristic to minimize SSE.

# Pre-processing and Post-processing

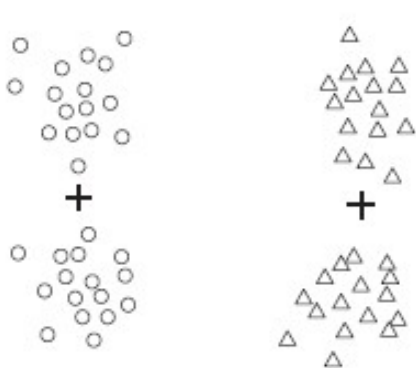
- Pre-processing
  - **Normalize** the data (e.g., scale to unit standard deviation)
  - **Eliminate outliers**
- Post-processing
  - **Eliminate** small clusters that may represent outliers
  - **Split** 'loose' clusters, i.e., clusters with relatively high SSE
  - **Merge** clusters that are 'close' and that have relatively low SSE

# Bisecting K-means

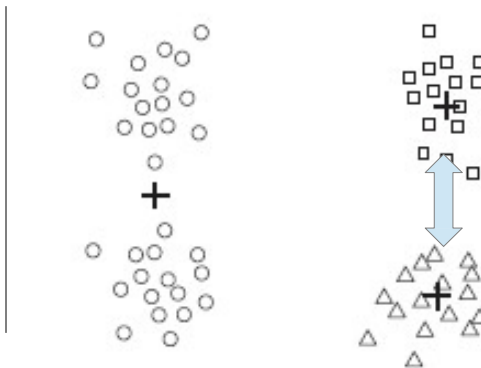
Variant of K-means that can produce a partitional or a hierarchical clustering

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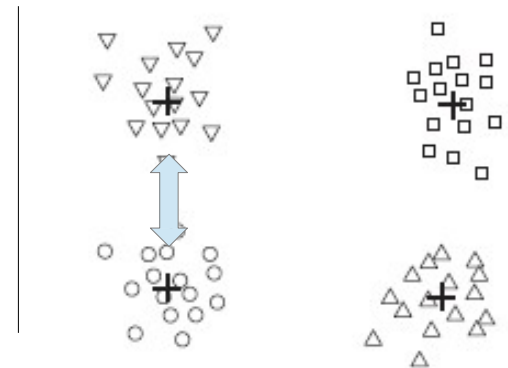
- 1: Initialize the list of clusters to contain the cluster containing all points.
  - 2: **repeat**
  - 3:   Select a cluster from the list of clusters
  - 4:   **for**  $i = 1$  to *number\_of\_iterations* **do**
  - 5:     Bisect the selected cluster using basic K-means
  - 6:   **end for**
  - 7:   Add the two clusters from the bisection with the lowest SSE to the list of clusters.
  - 8: **until** Until the list of clusters contains  $K$  clusters
- 



(a) Iteration 1.



(b) Iteration 2.



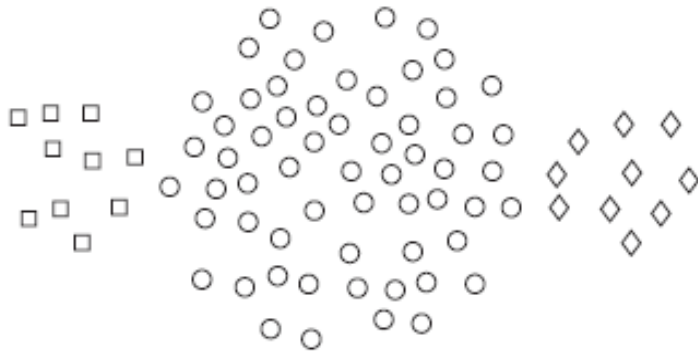
(c) Iteration 3.

# Limitations of K-means

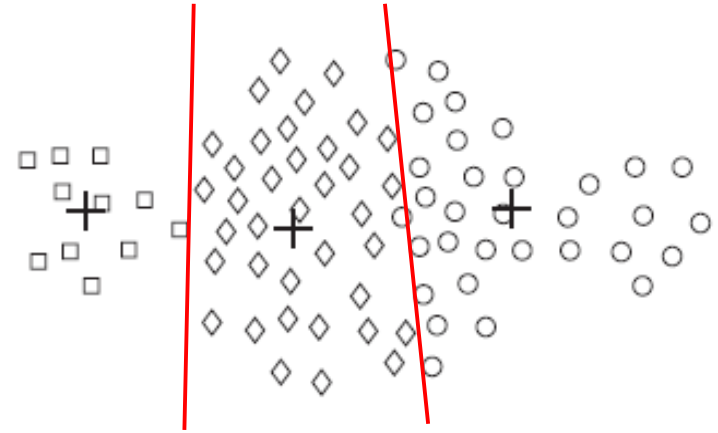
- K-means has problems when clusters are of differing
  - Sizes
  - Densities
  - Non-globular shapes
- K-means has problems when the data contains outliers.



# Limitations of K-means: Differing Sizes



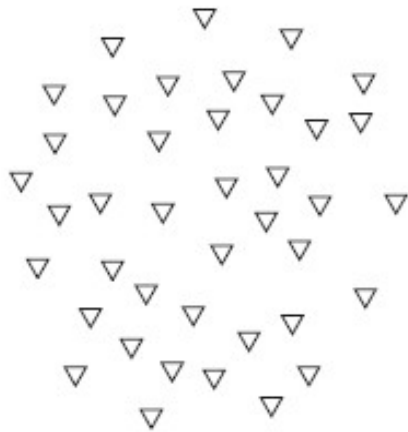
(a) Original points.



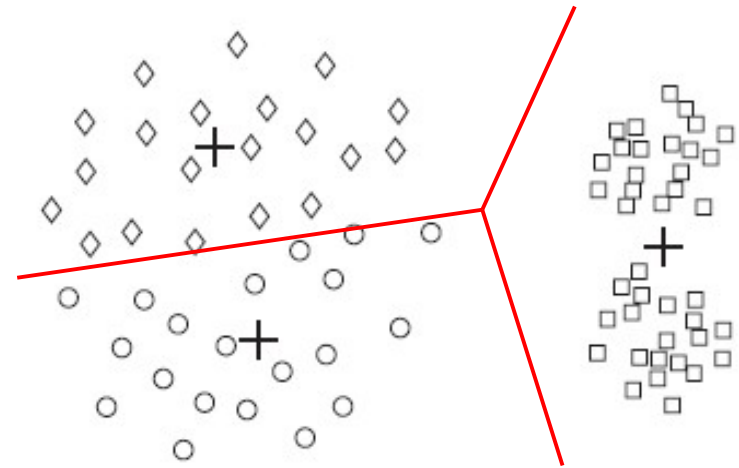
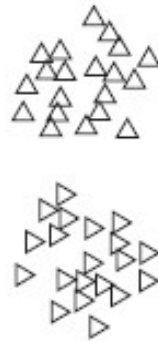
(b) Three K-means clusters.

**Figure 8.9.** K-means with clusters of different size.

# Limitations of K-means: Differing Density



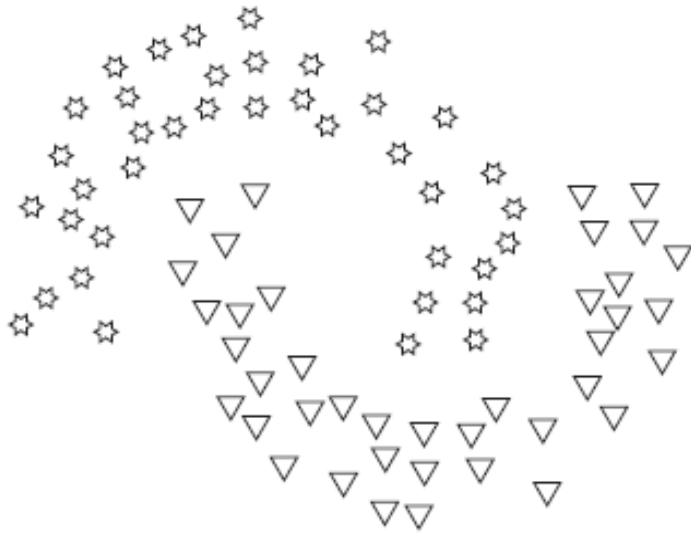
(a) Original points.



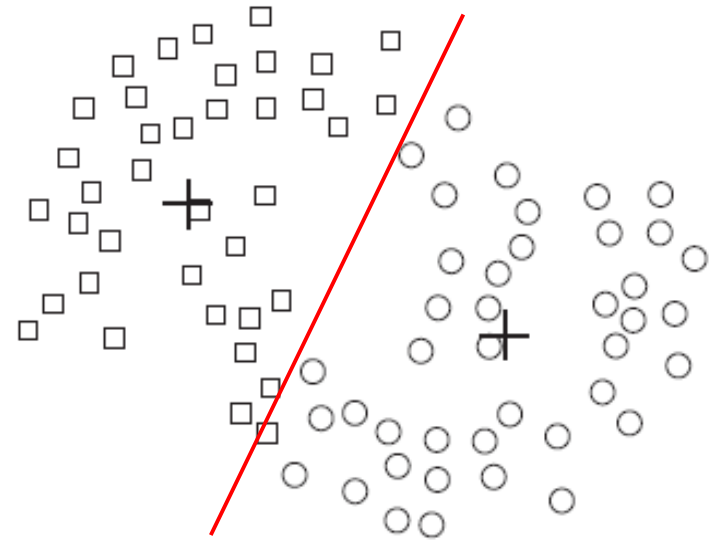
(b) Three K-means clusters.

**Figure 8.10.** K-means with clusters of different density.

# Limitations of K-means: Non-globular Shapes



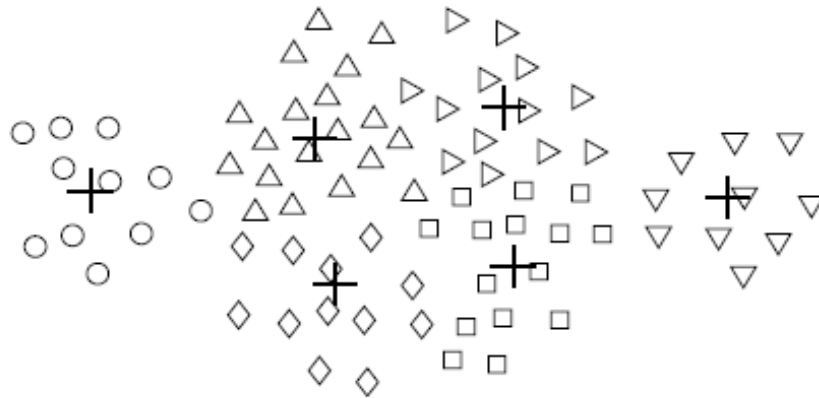
(a) Original points.



(b) Two K-means clusters.

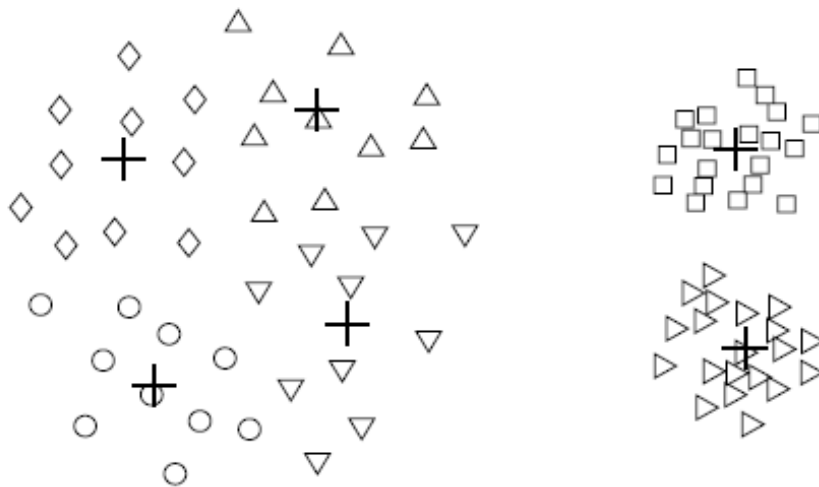
**Figure 8.11.** K-means with non-globular clusters.

# Overcoming K-means Limitations



(a) Unequal sizes.

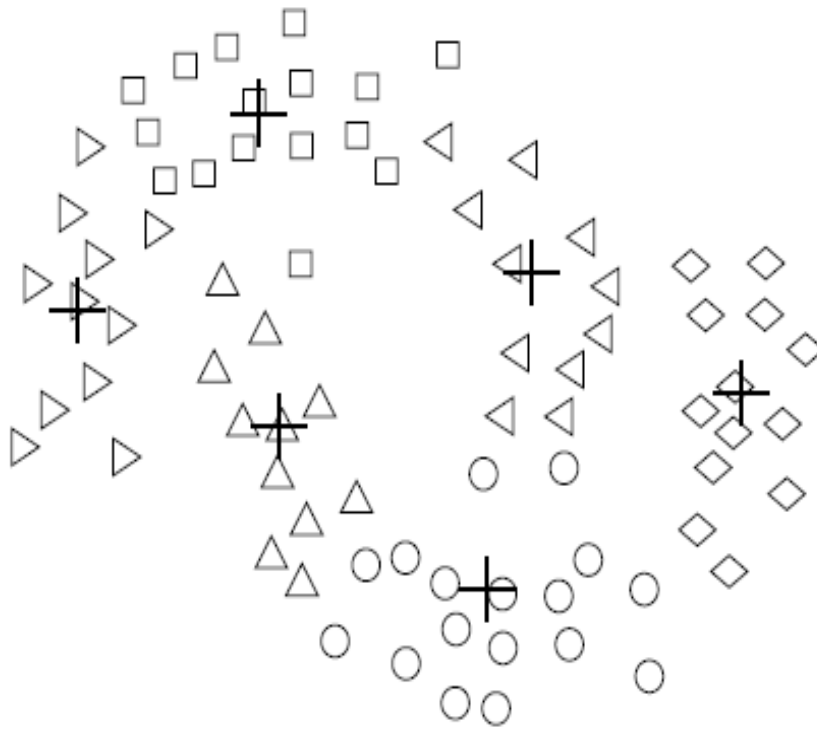
Use a larger  
number of clusters



(b) Unequal densities.

Several clusters  
represent a true  
cluster

# Overcoming K-means Limitations



(c) Non-spherical shapes.

Use a larger  
number of clusters

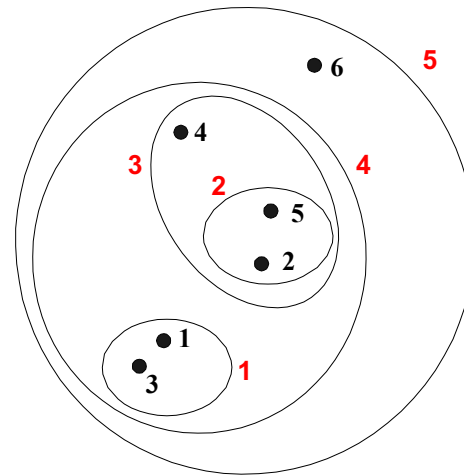
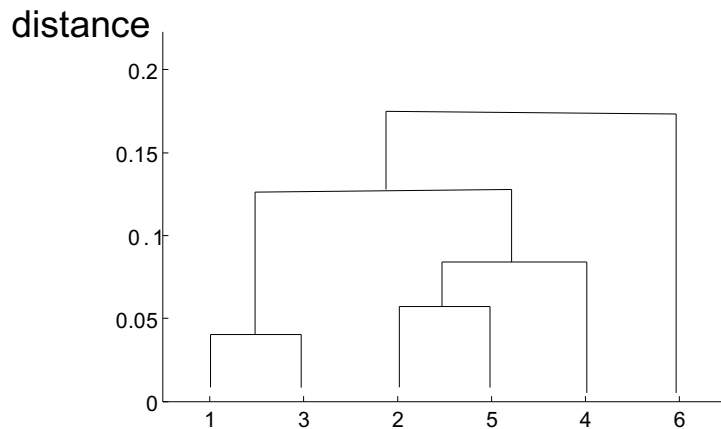
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# Hierarchical Clustering

- Produces a set of nested clusters organized as a hierarchical tree
- Can be visualized as a **dendrogram**
  - A tree like diagram that records the sequences of merges or splits



# Strengths of Hierarchical Clustering

- You do not have to assume any particular number of clusters
  - Any desired number of clusters can be obtained by 'cutting' the dendrogram at the proper level
- They may correspond to meaningful taxonomies
  - Example in biological sciences (e.g., animal kingdom, phylogeny reconstruction, ...)



# 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
  - Divisive:
    - Start with one, all-inclusive cluster
    - At each step, split a cluster until each cluster contains a point (or there are  $k$  clusters)
- Traditional hierarchical algorithms
  - use a similarity or distance matrix
  - merge or split one cluster at a time

# Agglomerative Clustering Algorithm

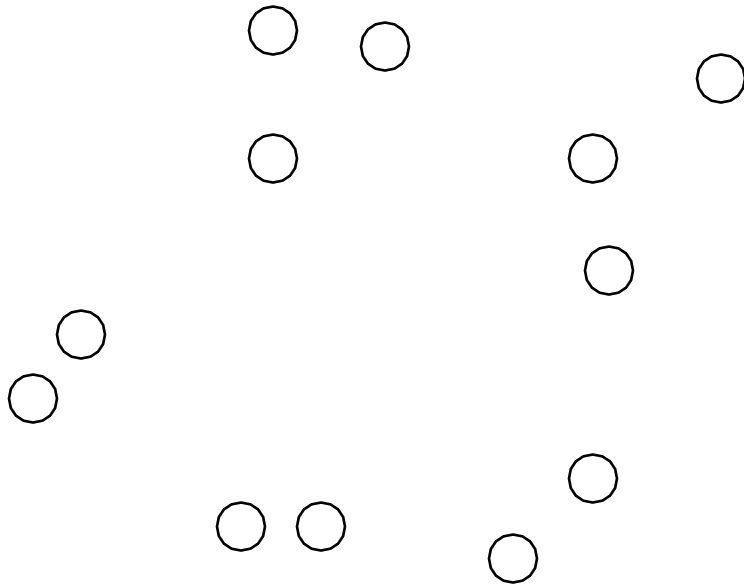
- More popular hierarchical clustering technique
- Basic algorithm is straightforward

1. Compute the proximity matrix
2. Let each data point be a cluster
3. Repeat
- 4        Merge the two closest clusters
- Update the proximity matrix
5. Until only a single cluster remains

- Key operation is the computation of the proximity of two clusters → Different approaches to **defining the distance between clusters** distinguish the different algorithms

# Starting Situation

Start with clusters of individual points and a proximity matrix



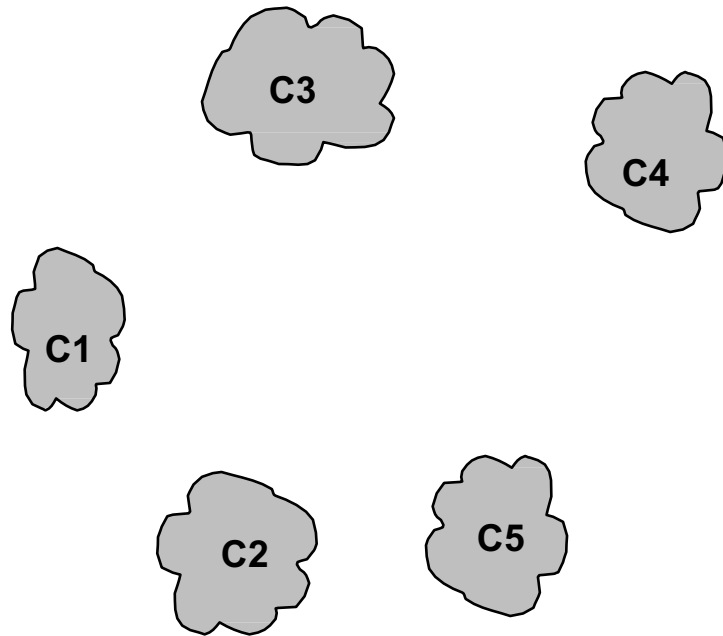
	p1	p2	p3	p4	p5	. . .
p1						
p2						
p3						
p4						
p5						
.						
.						
.						

**Proximity Matrix**



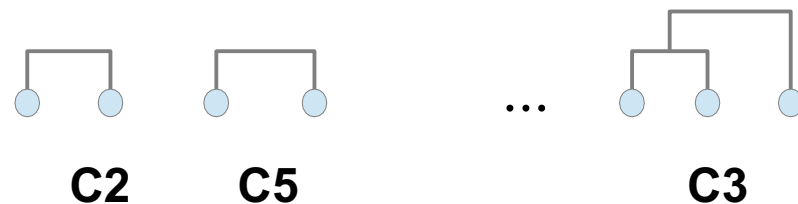
# Intermediate Situation

After some merging steps, we have some clusters



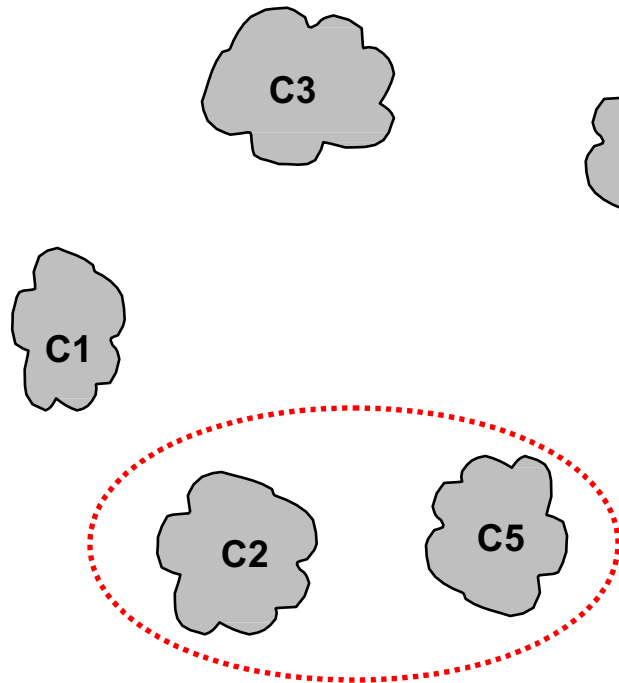
	C1	C2	C3	C4	C5
C1					
C2					
C3					
C4					
C5					

**Proximity Matrix**



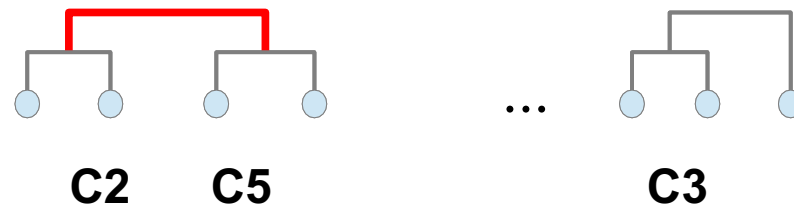
# Intermediate Situation

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



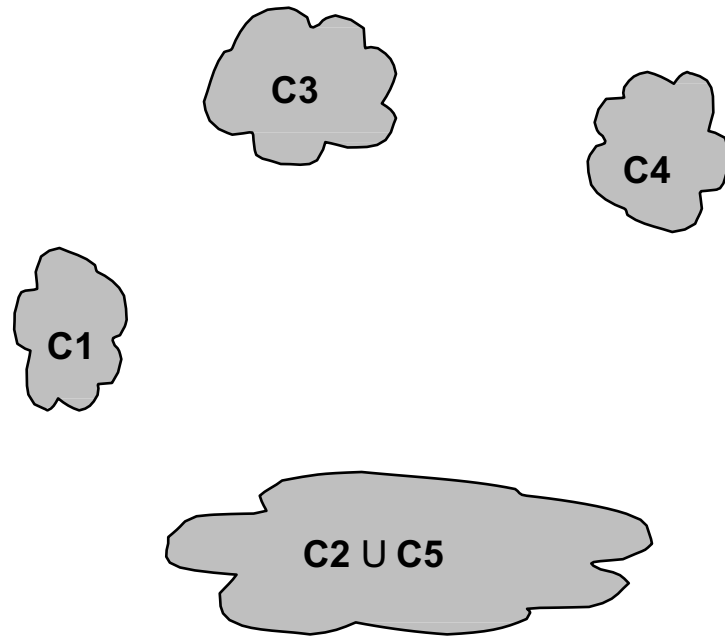
	C1	C2	C3	C4	C5
C1					
C2					
C3					
C4					
C5					

Proximity Matrix



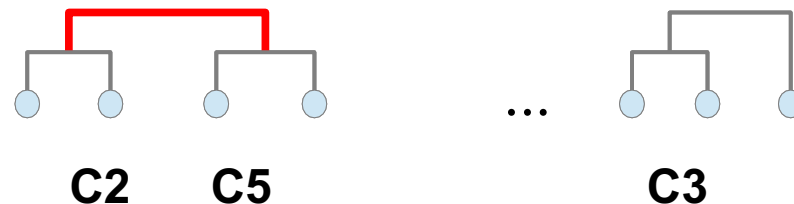
# After Merging

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

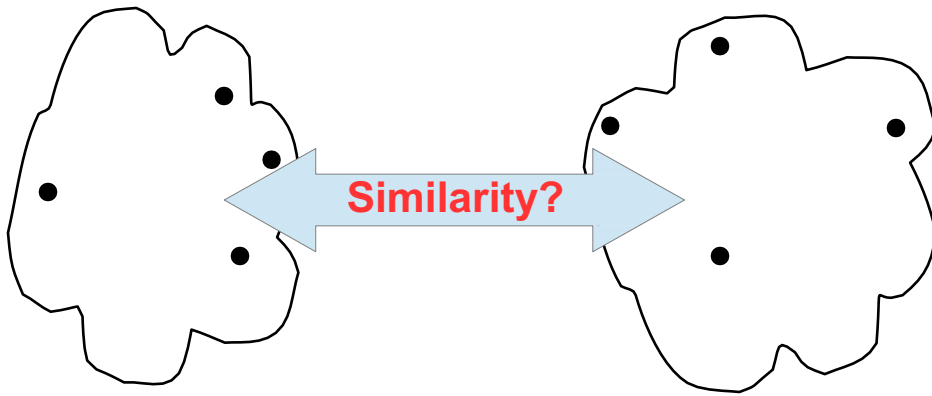


		$C2 \cup C5$		
	C1	C5	C3	C4
C1		?		
C2 U C5	?	?	?	?
C3		?		
C4		?		

Proximity Matrix



# How to Define Inter-Cluster Similarity

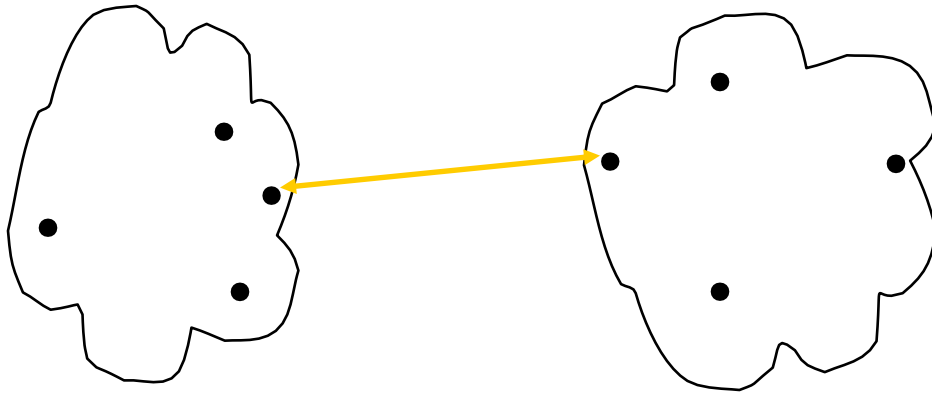


- MIN
- MAX
- Group Average
- Distance Between Centroids
- Other methods driven by an objective function
  - Ward's Method uses squared error

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

• **Proximity Matrix**

# How to Define Inter-Cluster Similarity



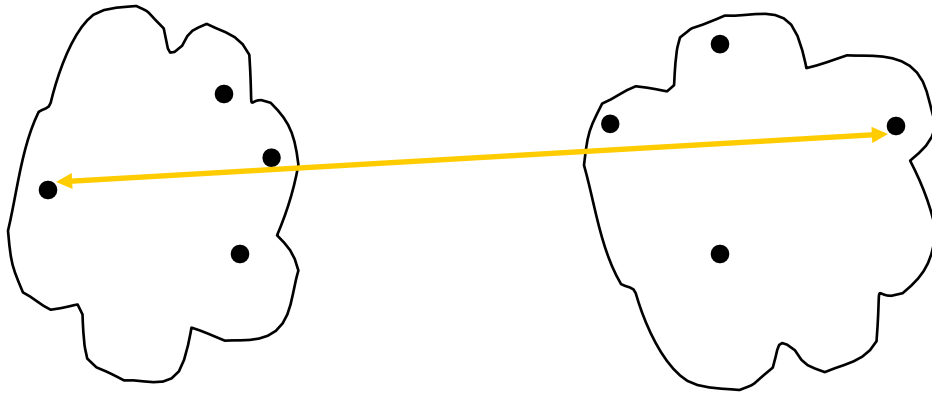
- **MIN (Single Link)**
- MAX
- Group Average
- Distance Between Centroids
- Other methods driven by an objective function
  - Ward's Method uses squared error

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

• **Proximity Matrix**



# How to Define Inter-Cluster Similarity

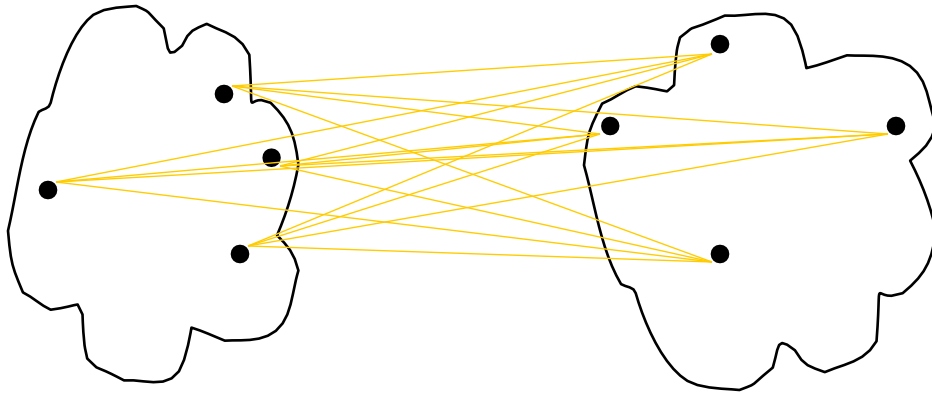


- MIN
- **MAX (Complete Link)**
- Group Average
- Distance Between Centroids
- Other methods driven by an objective function
  - Ward's Method uses squared error

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

• **Proximity Matrix**

# How to Define Inter-Cluster Similarity



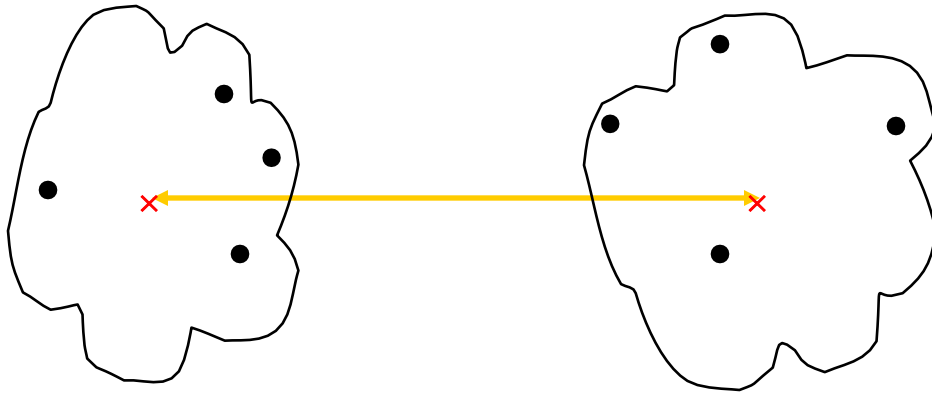
- MIN
- MAX
- **Group Average (Average Link)**
- Distance Between Centroids
- Other methods driven by an objective function
  - Ward's Method uses squared error

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

.

• Proximity Matrix

# How to Define Inter-Cluster Similarity

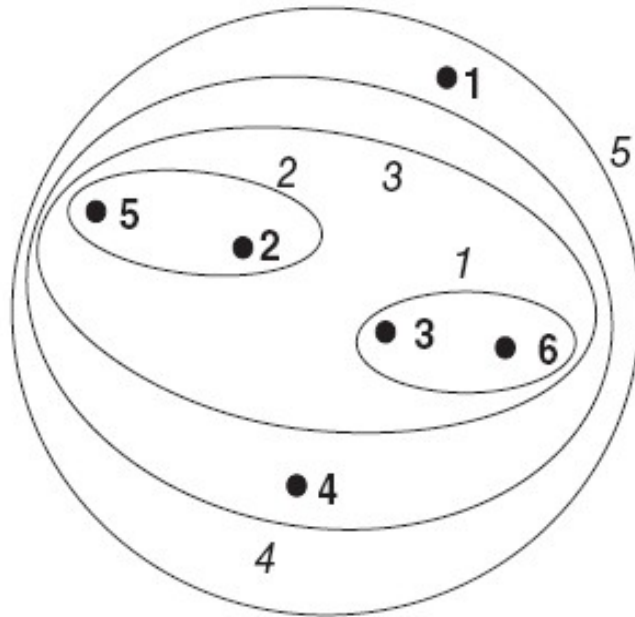


- MIN
- MAX
- Group Average
- **Distance Between Centroids**
- Other methods driven by an objective function
  - Ward's Method uses squared error

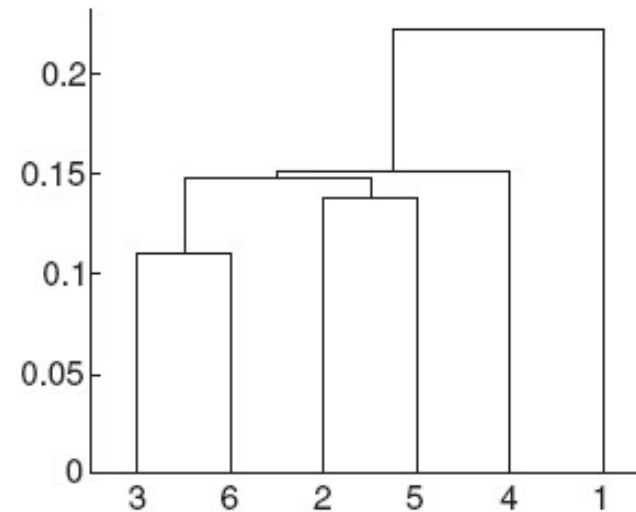
	p1	p2	p3	p4	p5	...
p1						
p2						
p3						
p4						
p5						
.						

• **Proximity Matrix**

# Single Link



(a) Single link clustering.



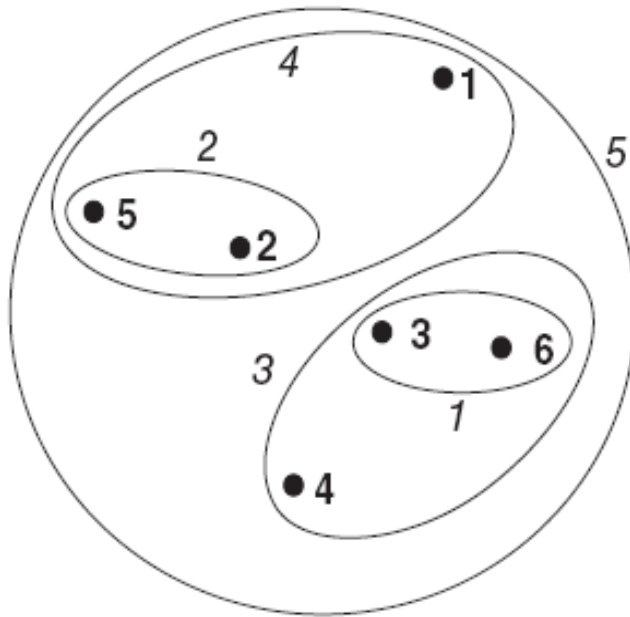
(b) Single link dendrogram.

Figure 8.16. Single link clustering of the six points shown in Figure 8.15.

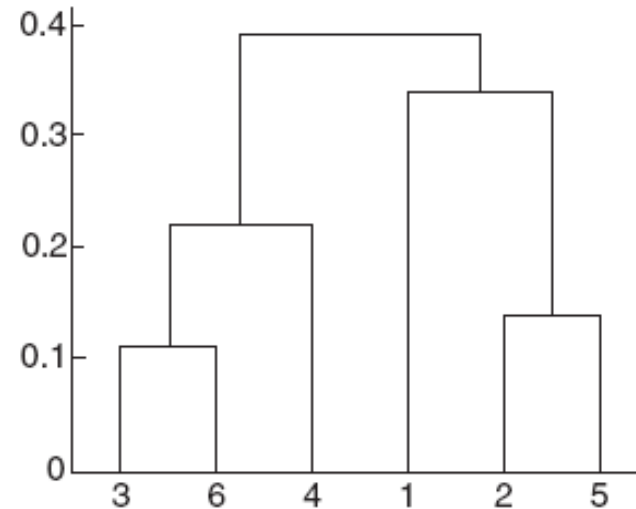
**Advantage:** Non-spherical, non-convex clusters

**Problem:** Chaining

# Complete Link



(a) Complete link clustering.



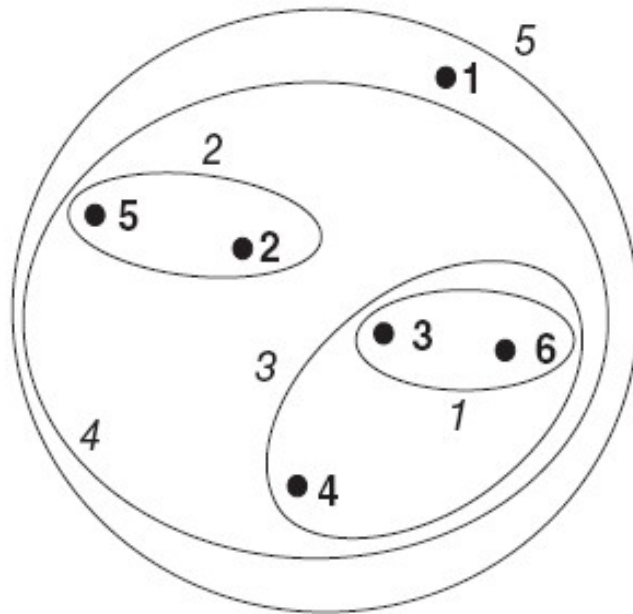
(b) Complete link dendrogram.

Figure 8.17. Complete link clustering of the six points shown in Figure 8.15.

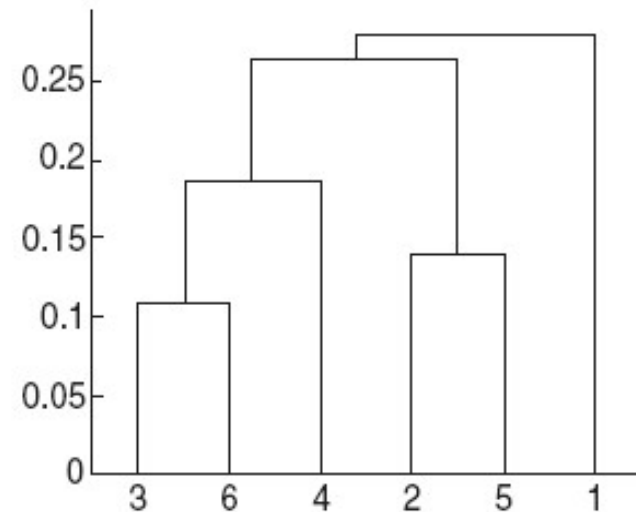
**Advantage:** more robust against noise (no chaining)

**Problem:** Tends to break large clusters,  
Biased towards globular clusters

# Average Link



(a) Group average clustering.



(b) Group average dendrogram.

Figure 8.18. Group average clustering of the six points shown in Figure 8.15.

## Compromise between Single and Complete Link

# Cluster Similarity: Ward's Method

- Similarity of two clusters is based on the increase in squared error when two clusters are merged
- Less susceptible to noise and outliers
- Biased towards globular clusters
- Hierarchical analogue of K-means
  - Can be used to initialize K-means

# Hierarchical Clustering: Complexity

- $O(N^2)$  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 proximity matrix of size  $N^2$  must be updated and searched
  - Complexity can be reduced to  $O(N^2 \log(N))$  time for some approaches



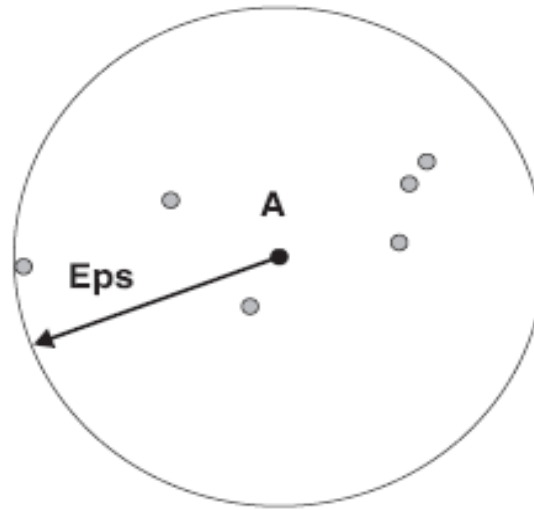
# Hierarchical Clustering: Limitations

- **Greedy:** Once a decision is made to combine two clusters, it cannot be undone
- **No global 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
  - Chaining, breaking large clusters

# Topics

- Introduction
- Types of Clustering
- Types of Clusters
- **Clustering Algorithms**
  - K-Means Clustering
  - Hierarchical Clustering
  - **Density-based Clustering**
- Cluster Validation

# DBSCAN



Density = 7 points

Figure 8.20. Center-based density.

- **Density** = number of points within a specified radius (Eps)

# DBSCAN

MinPts = 4

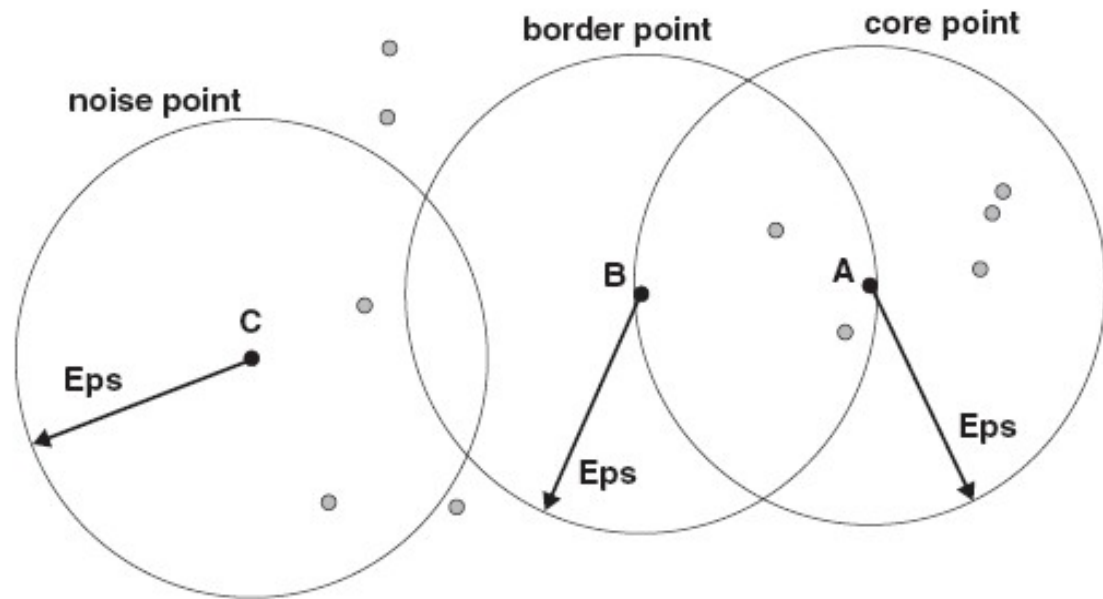


Figure 8.21. Core, border, and noise points.

- A point is a **core point** if it has more than a specified number of points (MinPts) within Eps. These are points that are at the interior of a cluster
- A **border point** has fewer than MinPts within Eps, but is in the neighborhood of a core point
- A **noise point** is any point that is not a core point or a border point.

# DBSCAN Algorithm

**DBSCAN**(D, eps, MinPts)

C = 0

for each unvisited point P in dataset D

mark P as visited

NeighborPts = regionQuery(P, eps)

if sizeof(NeighborPts) < MinPts

mark P as NOISE

else

C = next cluster

expandCluster(P, NeighborPts, C, eps, MinPts)

**expandCluster**(P, NeighborPts, C, eps, MinPts)

add P to cluster C

for each point P' in NeighborPts

if P' is not visited

mark P' as visited

NeighborPts' = regionQuery(P', eps)

if sizeof(NeighborPts') >= MinPts

NeighborPts = NeighborPts joined with NeighborPts'

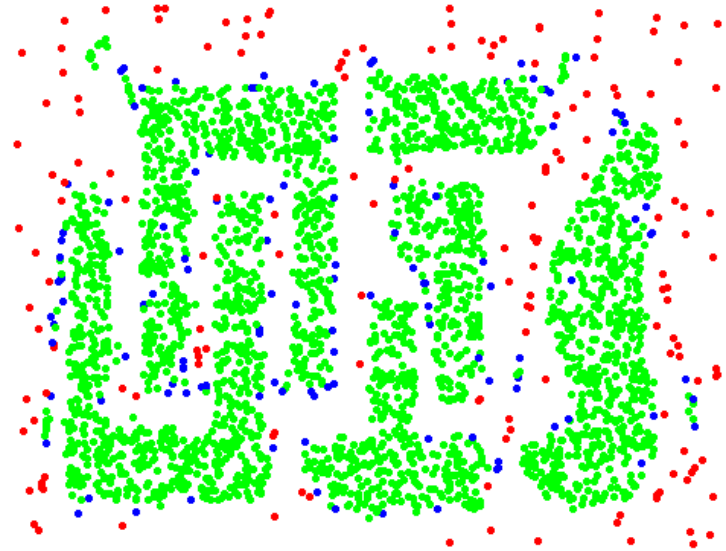
if P' is not yet member of any cluster

add P' to cluster C

# DBSCAN: Core, Border and Noise Points



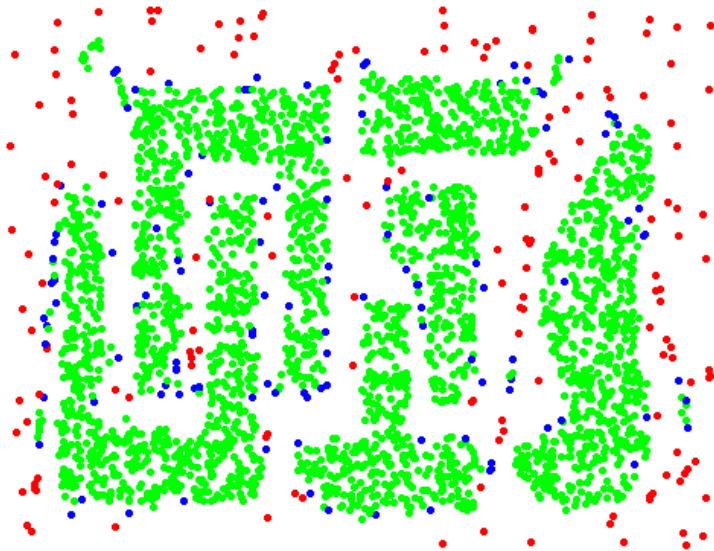
Original Points



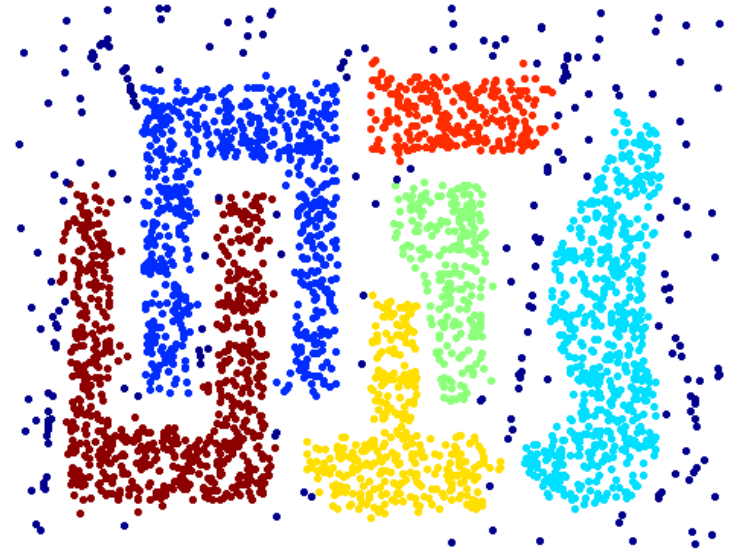
Point types: **core**,  
**border** and **noise**

Eps = 10, MinPts = 4

# DBSCAN: Determine Clusters



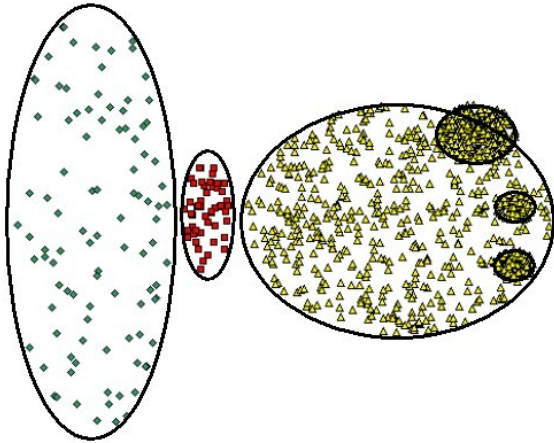
Point types: **core**,  
**border** and **noise**



Clusters

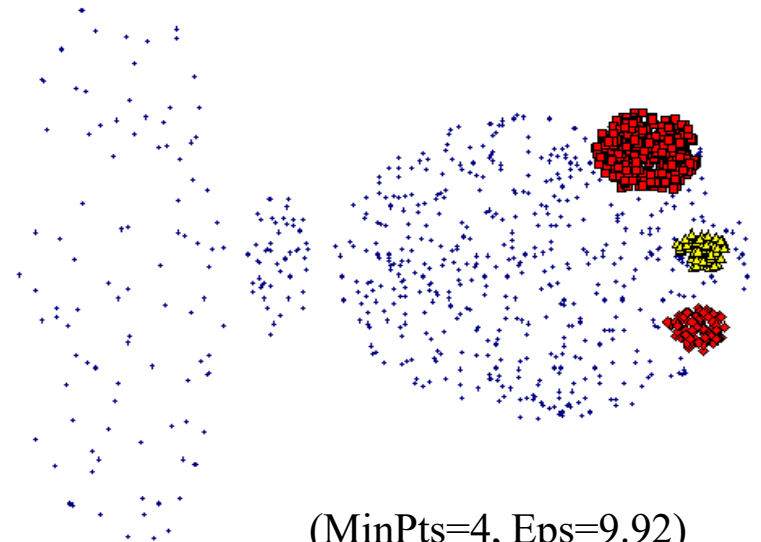
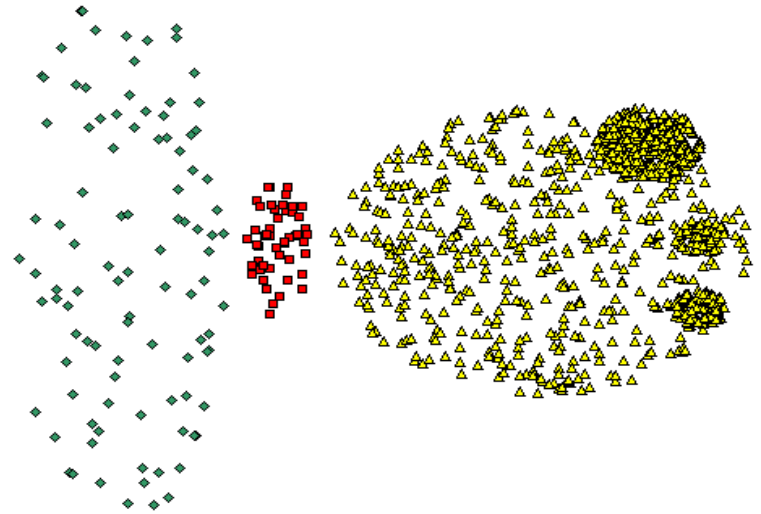
- Resistant to Noise
- Can handle clusters of different shapes and sizes
- Eps and MinPts depend on each other and can be hard to specify

# When DBSCAN Does NOT Work Well



Original Points

- Varying densities
- High-dimensional data

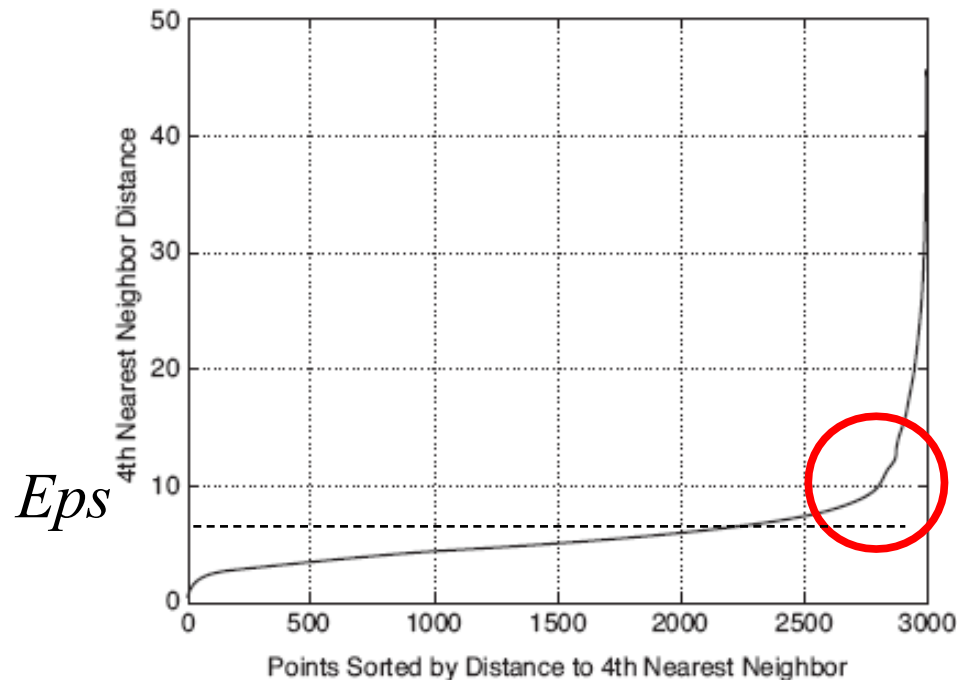


(MinPts=4, Eps=9.92)



# DBSCAN: Determining EPS and MinPts

- Idea is that for points in a cluster, their  $k$ th nearest neighbors are at roughly the same distance
- Noise points have the  $k$ th nearest neighbor at farther distance
- So, plot sorted distance of every point to its  $k$ th nearest neighbor



$$MinPts = k$$

# Some Other Clustering Algorithms

- **Center-based Clustering**

- Fuzzy c-means
- PAM (Partitioning Around Medoids)

- **Mixture Models**

- Expectation-maximization (EM) algorithm

- **Hierarchical**

- CURE (Clustering Using Representatives): shrinks points toward center
- BIRCH (balanced iterative reducing and clustering using hierarchies)

- **Graph-based Clustering**

- Graph partitioning on a sparsified proximity graph
- Shared nearest-neighbor (SNN graph)

- **Subspace Clustering**

- **Data Stream Clustering**

# Topics

- Introduction
- Types of Clustering
- Types of Clusters
- Clustering Algorithms
  - K-Means Clustering
  - Hierarchical Clustering
  - Density-based Clustering
- Cluster Validation

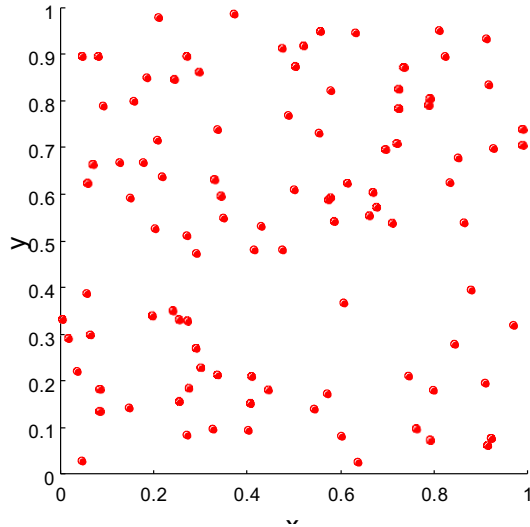
# Cluster Validity

- For supervised classification (= we have a class label) we have a variety of measures to evaluate how good our model is: Accuracy, precision, recall
- For cluster analysis (=unsupervised learning), the analogous question is:

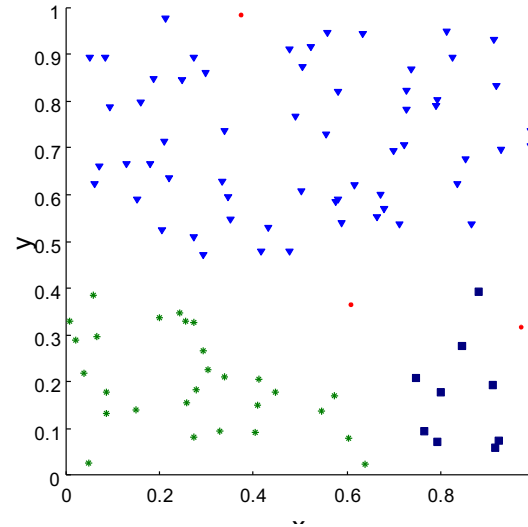
**How to evaluate the “goodness” of the resulting clusters?**

# Clusters found in Random Data (Overfitting)

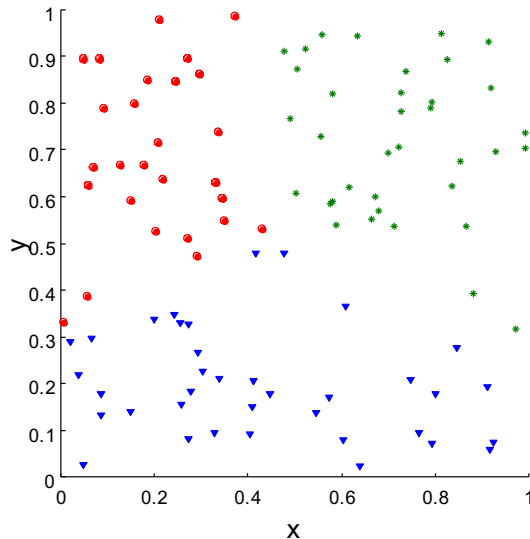
Random  
Points



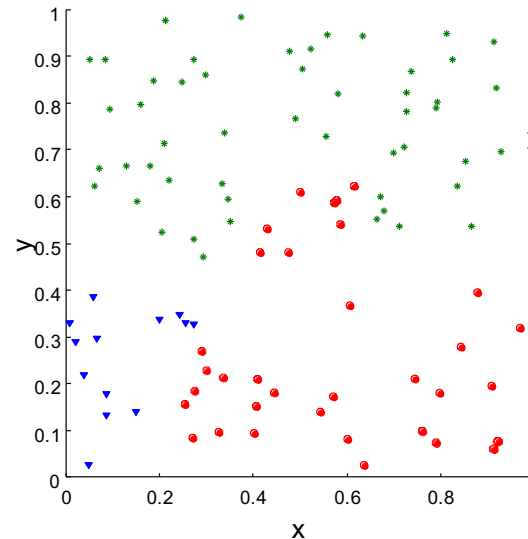
DBSCAN



K-means



Complete  
Link



**If you tell a clustering algorithm to find clusters then it will!**

# Different Aspects of Cluster Validation

1. Determining the **clustering tendency** of a set of data, i.e., distinguishing whether non-random structure actually exists in the data (e.g., to avoid overfitting).
2. **External Validation:** Compare the results of a cluster analysis to externally known class labels (ground truth).
3. **Internal Validation:** Evaluating how well the results of a cluster analysis fit the data *without* reference to external information.
4. **Compare clusterings** to determine which is better.
5. Determining the **'correct' number of clusters**.

For 2, 3, and 4, we can further distinguish whether we want to evaluate the entire clustering or just individual clusters.

# Measures of Cluster Validity

Numerical measures that are applied to judge various aspects of cluster validity, are classified into the following three types.

- **External Index:** Used to measure the extent to which cluster labels match externally supplied class labels.
  - Entropy, Purity, Rand index
- **Internal Index:** Used to measure the goodness of a clustering structure *without* respect to external information.
  - Sum of Squared Error (SSE), Silhouette coefficient
- **Relative Index:** Used to compare two different clusterings or clusters.
  - Often an external or internal index is used for this function, e.g., SSE or entropy

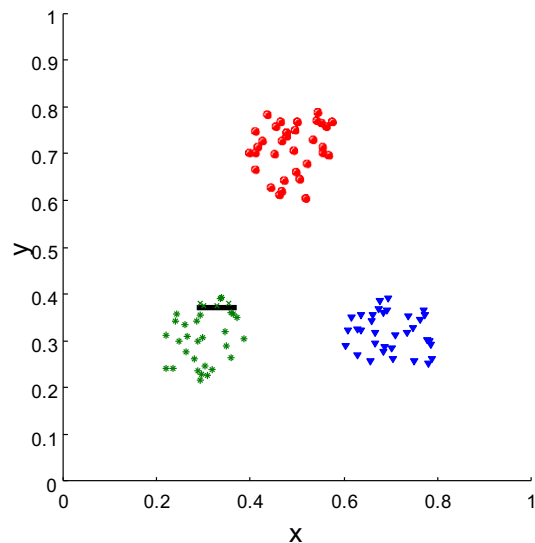
# Measuring Cluster Validity Via Correlation

- Two matrices
  - Proximity Matrix
  - “Incidence” Matrix
    - One row and one column for each data point
    - An entry is 1 if the associated pair of points belong to the same cluster
    - An entry is 0 if the associated pair of points belongs to different clusters
- Compute the correlation between the two matrices
  - Since the matrices are symmetric, only the correlation between  $n(n-1) / 2$  entries needs to be calculated.
- High correlation indicates that points that belong to the same cluster are close to each other.
- Not a good measure for some density or contiguity based clusters (e.g., single link HC).

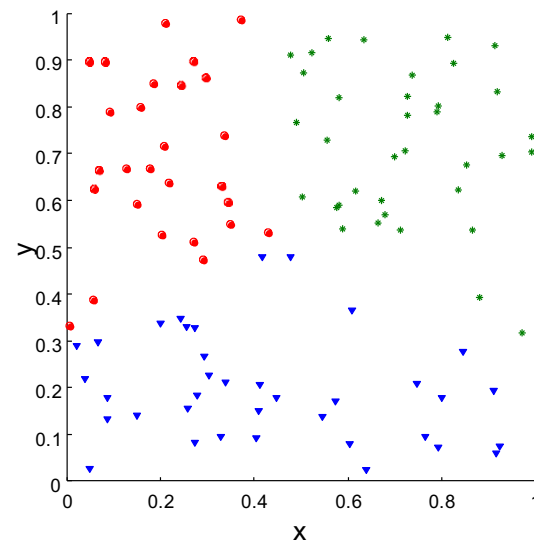


# Measuring Cluster Validity Via Correlation

- Correlation of incidence and proximity matrices for the K-means clusterings of the following two data sets.



**Corr = -0.9235**

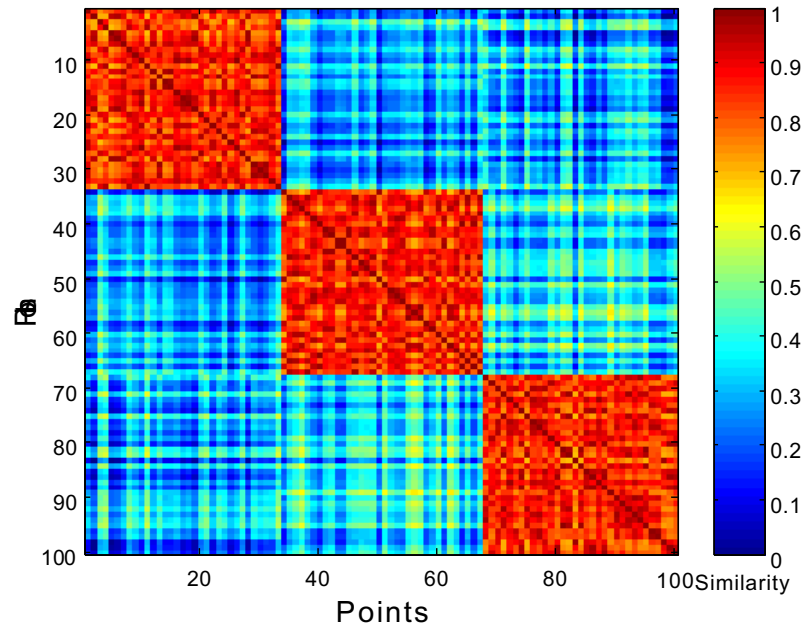
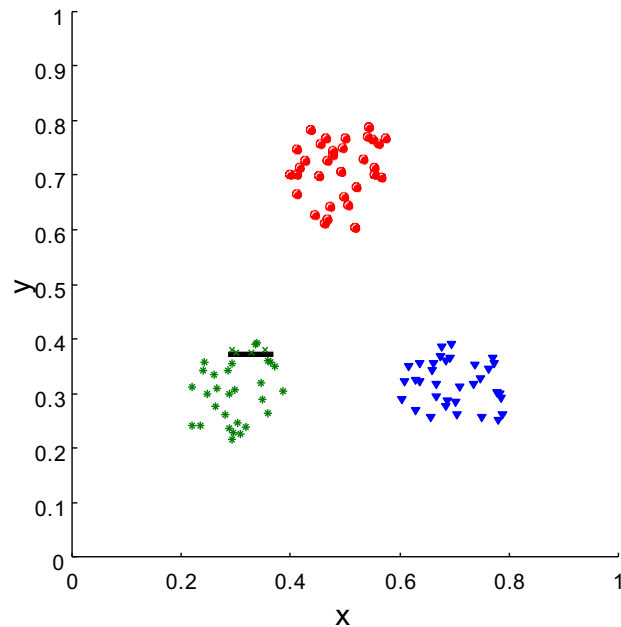


**Corr = -0.5810**

Note: Correlation is negative between distance matrix and incidence matrix

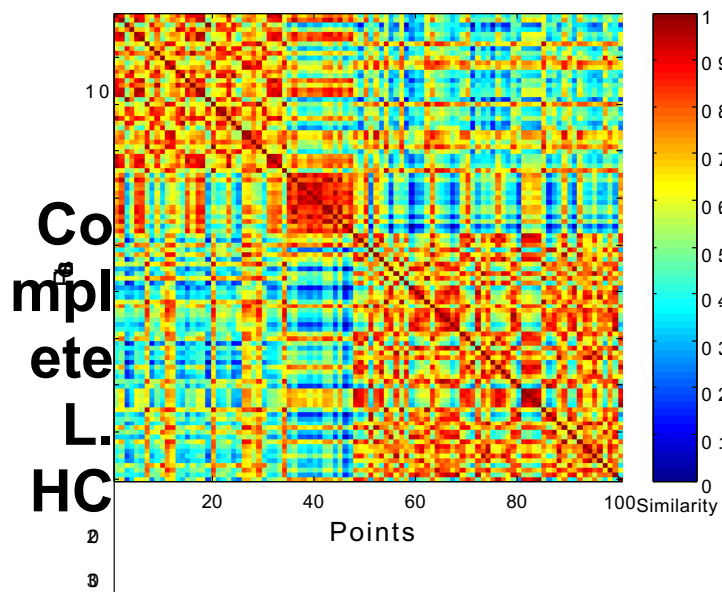
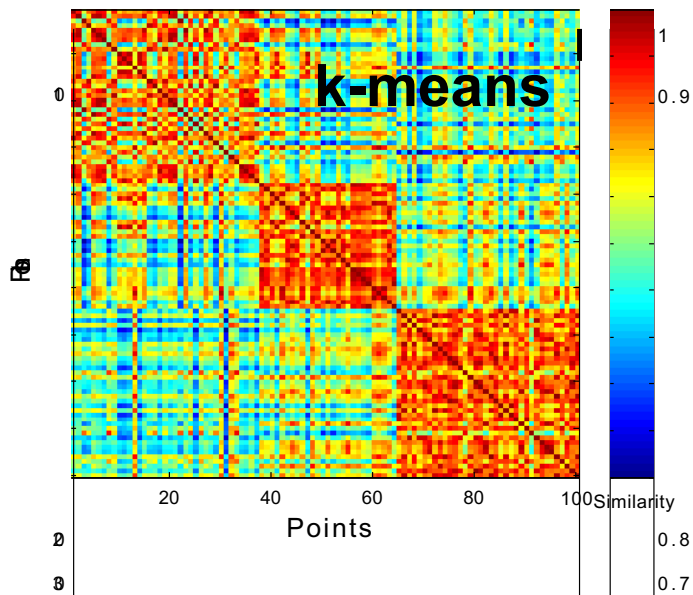
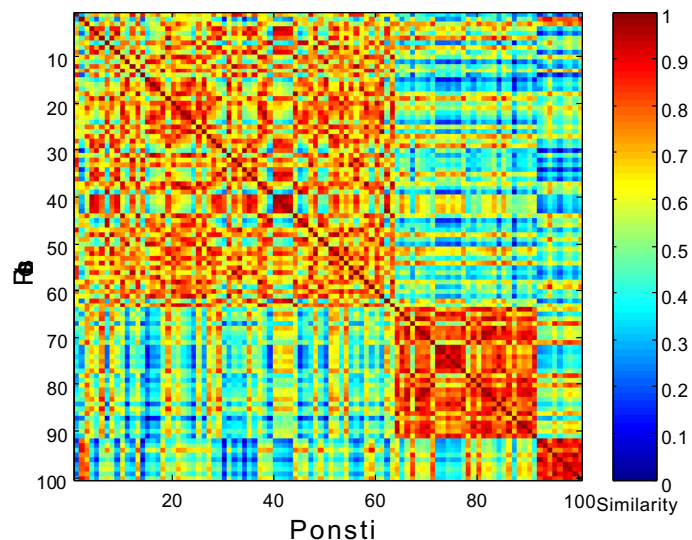
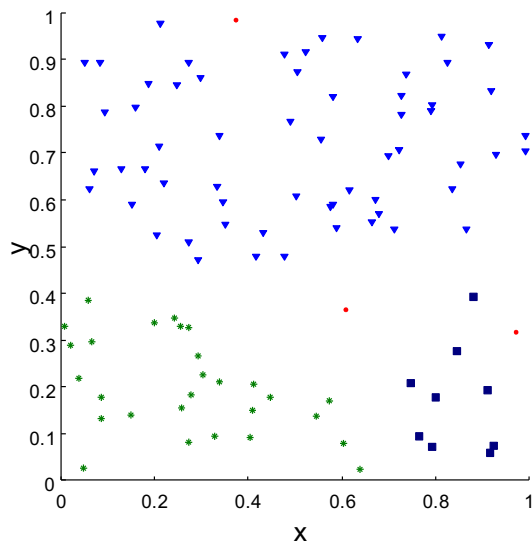
# Using Similarity Matrix for Cluster Validation

Order the similarity matrix with respect to cluster labels and inspect visually.



# Using Similarity Matrix for Cluster Validation

Clusters in random data are not so crisp



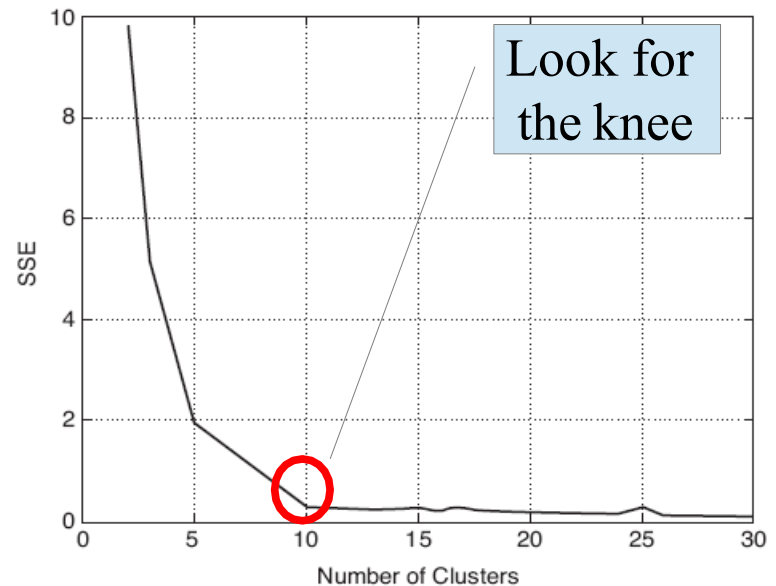
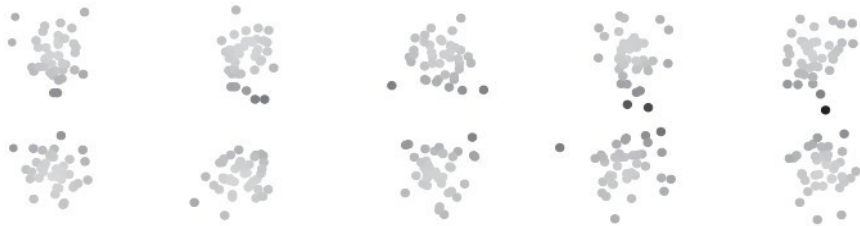
# Internal Measures: SSE

- SSE is good for comparing two clusterings or two clusters (average SSE).

$$SSE = \sum_{i=1}^K \sum_{\mathbf{x} \in C_i} \|\mathbf{x} - \mathbf{m}_i\|^2$$

- Can also be used to estimate the number of clusters

10 clusters



**Figure 8.32.** SSE versus number of clusters for the data of Figure 8.29.

# Internal Measures: Cohesion and Separation

- **Cluster Cohesion:** Measures how closely related are objects in a cluster
  - Example: Within cluster sum of squares (WSS=SSE)

$$WSS = \sum_i \sum_{x \in C_i} \|x - m_i\|^2$$

- **Cluster Separation:** Measure how distinct or well-separated a cluster is from other clusters
  - Example: Between cluster sum of squares (BSS)

$$BSS = \sum_i |C_i| \|m - m_i\|^2$$

Where  $|C_i|$  is the size of cluster  $i$

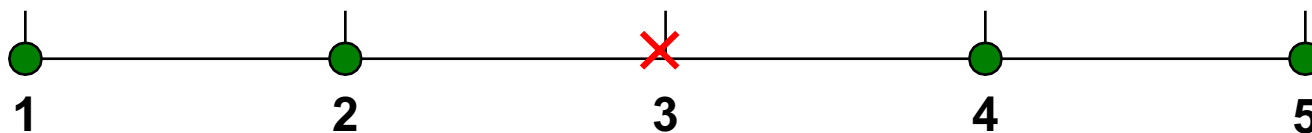
- **Total sum of squares**  $TSS = WSS + BSS$

# Internal Measures: Cohesion and Separation

Example:  $TSS = BSS + WSS = \text{constant}$  for a given data set

m

**K=1 cluster:**

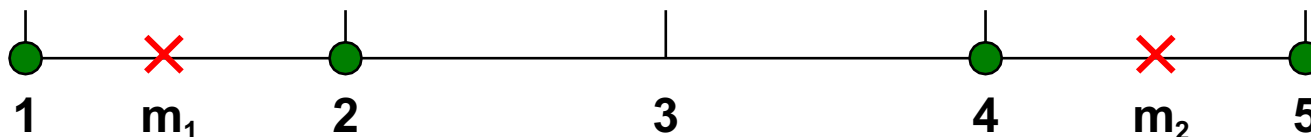


$$WSS = (1-3)^2 + (2-3)^2 + (4-3)^2 + (5-3)^2 = 10$$

$$BSS = 4 \times (3-3)^2 = 0$$

$$Total = 10 + 0 = 10$$

**K=2 clusters:**



$$WSS = (1-1.5)^2 + (2-1.5)^2 + (4-4.5)^2 + (5-4.5)^2 = 1$$

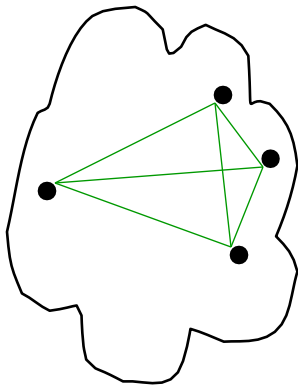
$$BSS = 2 \times (3-1.5)^2 + 2 \times (4.5-3)^2 = 9$$

$$Total = 1 + 9 = 10$$

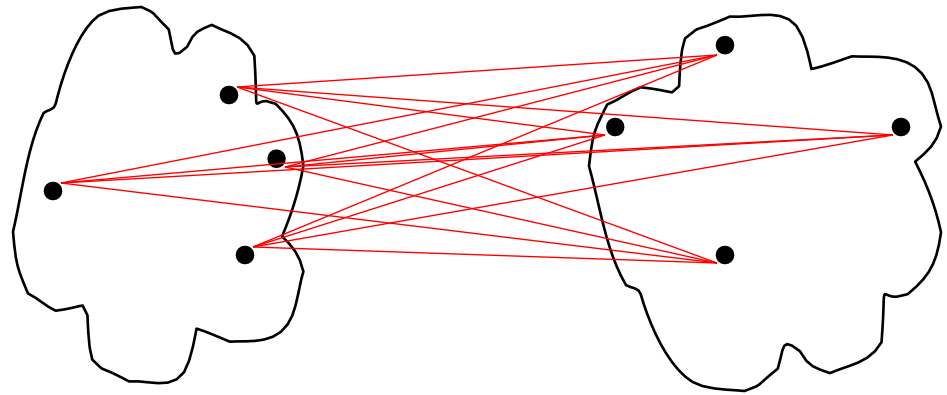
# Internal Measures: Cohesion and Separation

A proximity graph based approach can also be used for cohesion and separation.

- Cluster cohesion is the sum of the weight of all links within a cluster.
- Cluster separation is the sum of the weights between nodes in the cluster and nodes outside the cluster.



cohesion



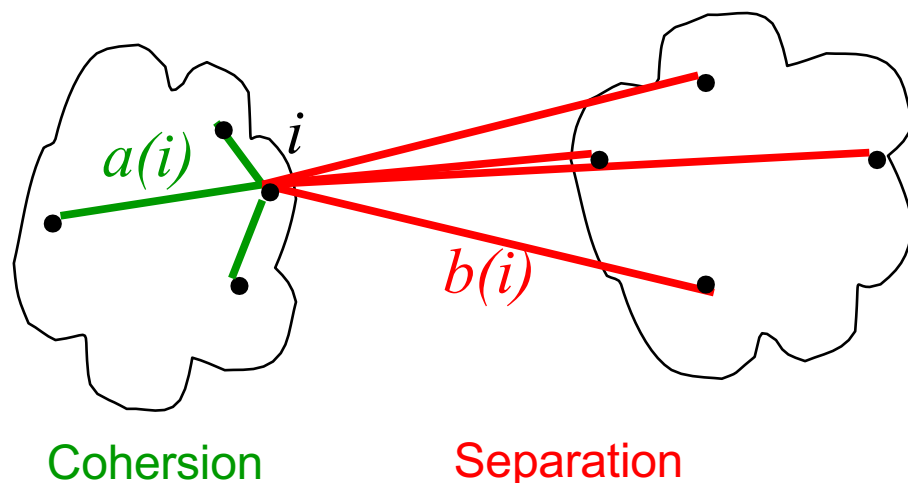
separation

# Internal Measures: Silhouette Coefficient

- Silhouette Coefficient combine ideas of both cohesion and separation, but for individual points. For an individual point  $i$ :
  - Calculate  $a(i)$  = average dissimilarity of  $i$  to all other points in its cluster
  - Calculate  $b(i)$  = lowest average dissimilarity of  $i$  to any other

$$s(i) = \begin{cases} 1 - a(i)/b(i), & \text{if } a(i) < b(i) \\ 0, & \text{if } a(i) = b(i) \\ b(i)/a(i) - 1, & \text{if } a(i) > b(i) \end{cases}$$

$$-1 \leq s(i) \leq 1$$



- The closer to 1 the better.
- Can calculate the Average Silhouette width for a cluster or a clustering



# Internal Measures: Silhouette Plot

Silhouette plot of pam(x = dis.bc, k = 5)

n = 160

5 clusters  $C_j$

$j : n_j \mid \text{ave}_{i \in C_j} s_i$

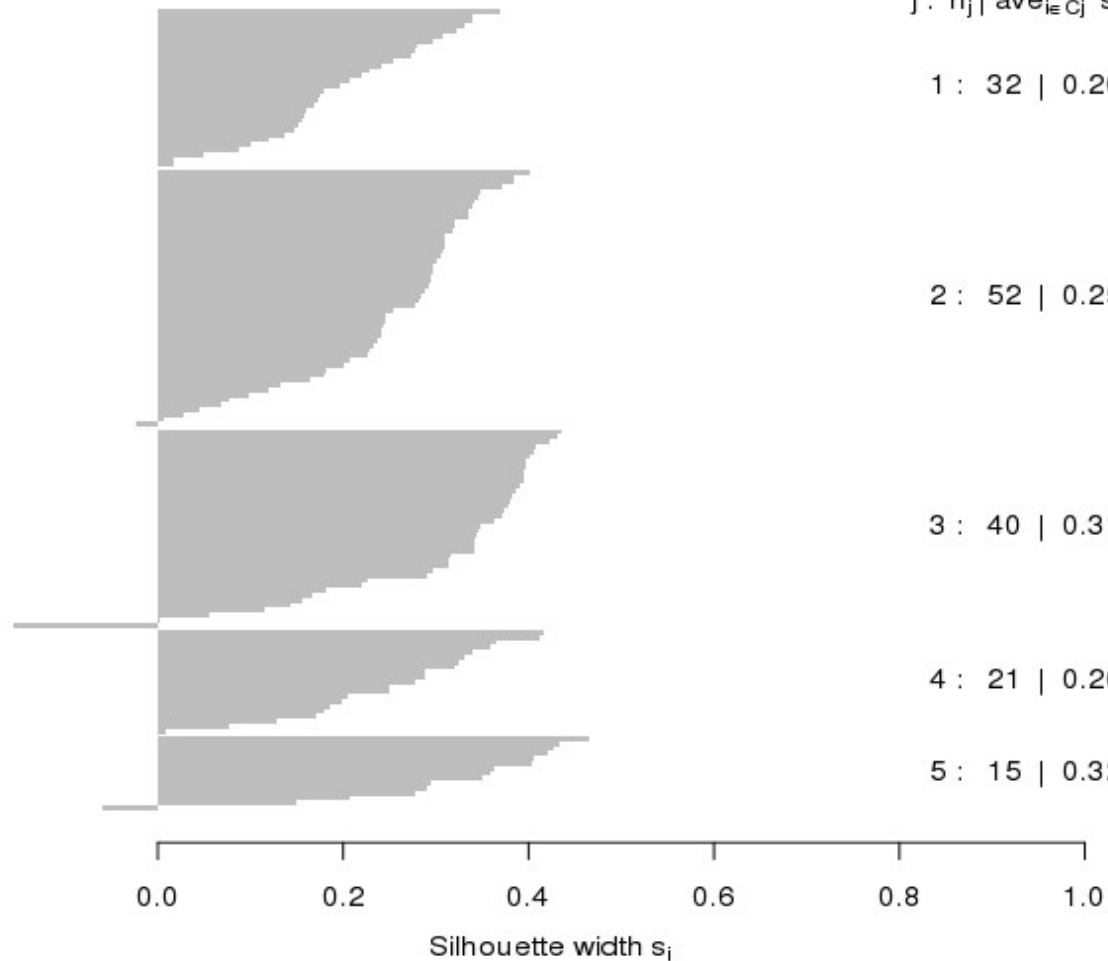
1 : 32 | 0.20

2 : 52 | 0.25

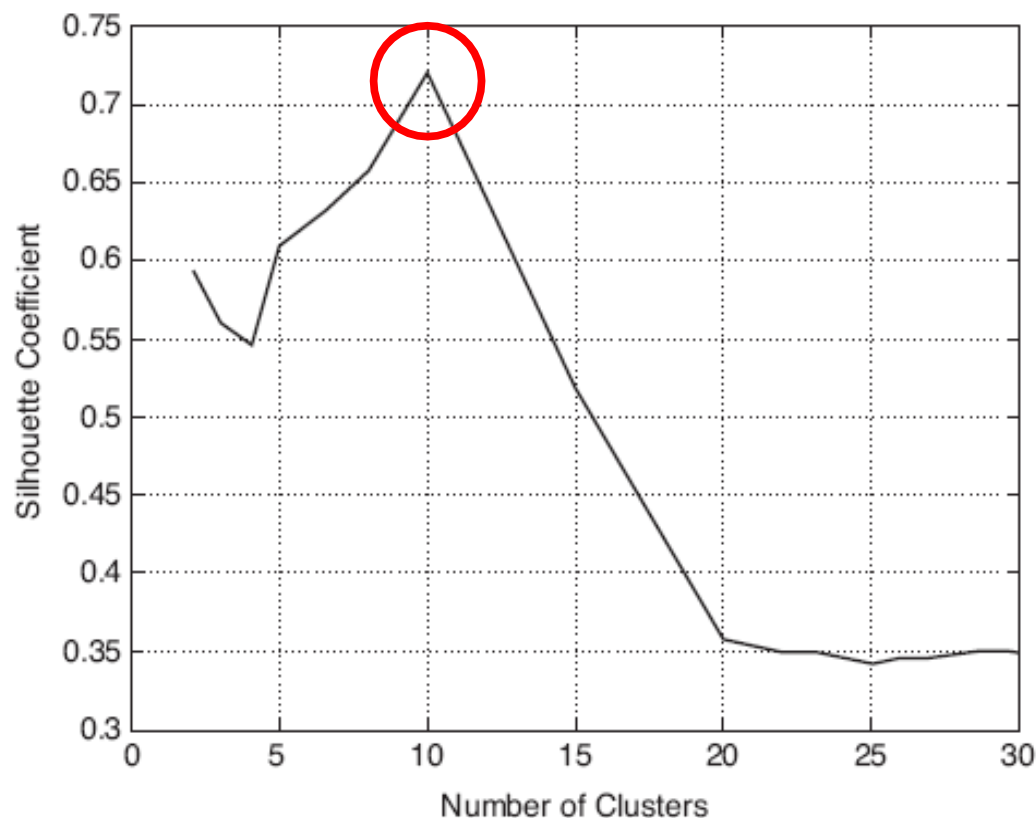
3 : 40 | 0.31

4 : 21 | 0.26

5 : 15 | 0.32



# Internal Measures: Choosing k with Silhouette



**Figure 8.33.** Average silhouette coefficient versus number of clusters for the data of Figure 8.29.

# External Measures of Cluster Validity: Entropy and Purity

**Table 5.9.** K-means Clustering Results for LA Document Data Set

Cluster	Entertainment	Financial	Foreign	Metro	National	Sports	Entropy	Purity
1	3	5	40	506	96	27	1.2270	0.7474
2	4	7	280	29	39	2	1.1472	0.7756
3	1	1	1	7	4	671	0.1813	0.9796
4	10	162	3	119	73	2	1.7487	0.4390
5	331	22	5	70	13	23	1.3976	0.7134
6	5	358	12	212	48	13	1.5523	0.5525
Total	354	555	341	943	273	738	1.1450	0.7203

**entropy** For each cluster, the class distribution of the data is calculated first, i.e., for cluster  $j$  we compute  $p_{ij}$ , the ‘probability’ that a member of cluster  $j$  belongs to class  $i$  as follows:  $p_{ij} = m_{ij}/m_j$ , where  $m_j$  is the number of values in cluster  $j$  and  $m_{ij}$  is the number of values of class  $i$  in cluster  $j$ . Then using this class distribution, the entropy of each cluster  $j$  is calculated using the standard formula  $e_j = \sum_{i=1}^L p_{ij} \log_2 p_{ij}$ , where the  $L$  is the number of classes. The total entropy for a set of clusters is calculated as the sum of the entropies of each cluster weighted by the size of each cluster, i.e.,  $e = \sum_{j=1}^K \frac{m_j}{m} e_j$ , where  $m_j$  is the size of cluster  $j$ ,  $K$  is the number of clusters, and  $m$  is the total number of data points.

**purity** Using the terminology derived for entropy, the purity of cluster  $j$ , is given by  $purity_j = \max_i p_{ij}$  and the overall purity of a clustering by  $purity = \sum_{j=1}^K \frac{m_j}{m} purity_j$ .

Other measures: Precision, Recall, F-measure, Rand, Adj. Rand

# Final Comment on Cluster Validity

“The validation of clustering structures is the most difficult and frustrating part of cluster analysis.

Without a strong effort in this direction, cluster analysis will remain a black art accessible only to those true believers who have experience and great courage.”

*Algorithms for Clustering Data, Jain and Dubes*