# BBS654 Data Mining

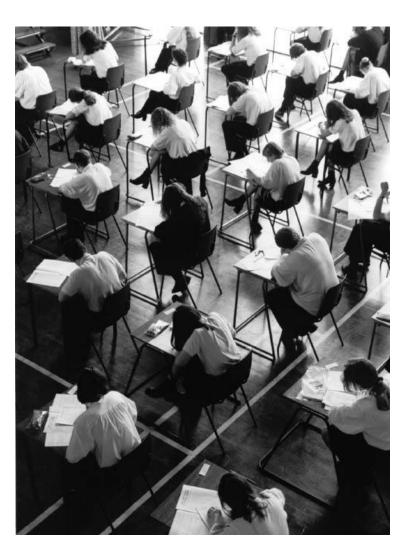
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Slides are adapted from Nazli Ikizler

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# Classification

- Classification systems:
  - Supervised learning
  - Make a rational prediction given evidence
  - There are several methods for this
  - Useful when you have labeled data (or can get it)



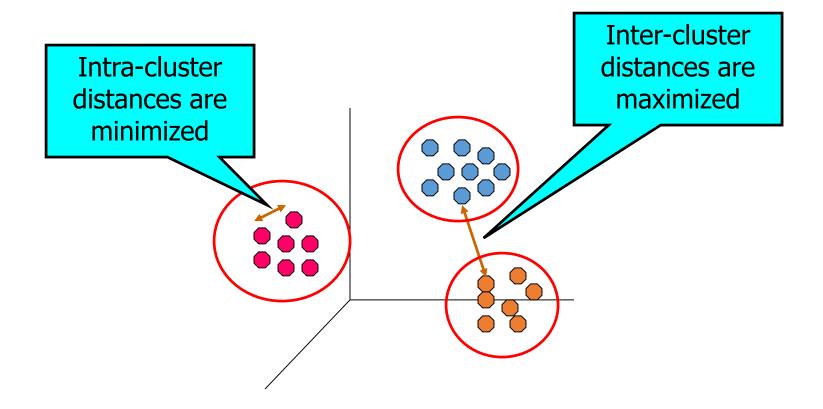
# Clustering

- Clustering systems:
  - Unsupervised learning
  - Detect patterns in unlabeled data
  - Useful when don't know what you're looking for
  - Requires data, but no labels
- Typical applications
  - As a stand-alone tool to get insight into data distribution
  - As a preprocessing step for other algorithms



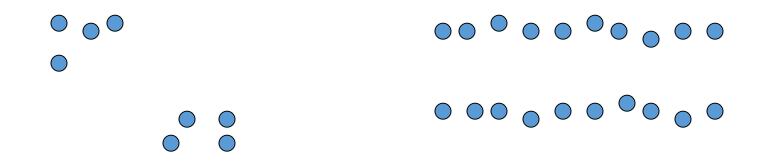
# What is Cluster Analysis?

 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



## Clustering

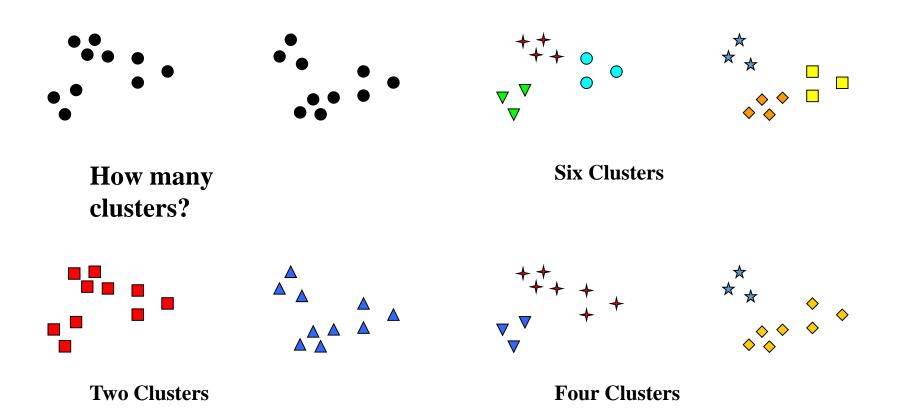
- Basic idea: group together similar instances
- Example: 2D point patterns



- What could "similar" mean?
  - One option: small (squared) Euclidean distance

dist
$$(x, y) = (x - y)^{\top} (x - y) = \sum_{i} (x_i - y_i)^2$$

Notion of a Cluster can be Ambiguous



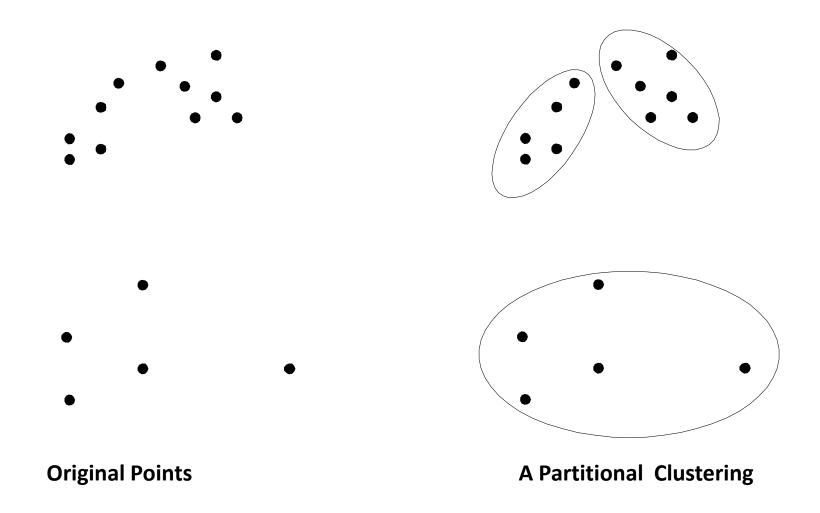
# Quality: What Is Good Clustering?

- A good clustering method will produce high quality clusters
  - high <u>intra-class</u> similarity: cohesive within clusters
  - low <u>inter-class</u> similarity: <u>distinctive</u> between clusters
- The <u>quality</u> of a clustering method depends on
  - the similarity measure used by the method
  - its implementation, and
  - Its ability to discover some or all of the hidden patterns

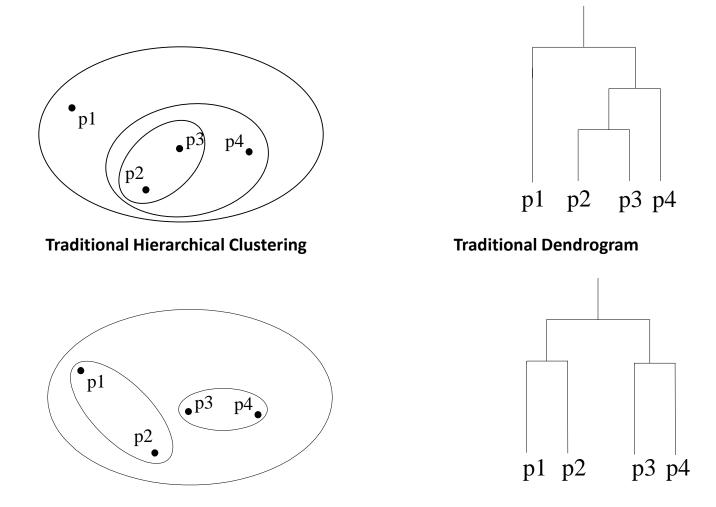
## Types of Clusterings

- A clustering is a set of clusters
- Important distinction between hierarchical and partitional sets 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



## Hierarchical Clustering



Non-traditional Hierarchical Clustering

Non-traditional Dendrogram

# Clustering Algorithms

- K-means and its variants
- Hierarchical clustering
- Density-based clustering

## 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
  - 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

#### Objective function

• 
$$J = \sum_{j=1}^{k} \sum_{C(i)=j} ||x_i - c_j||^2$$

- Total within-cluster variance
- Re-arrange the objective function

• 
$$J = \sum_{j=1}^{k} \sum_{i} w_{ij} ||x_i - c_j||^2$$

• 
$$w_{ij} \in \{0,1\}$$

- $w_{ij} = 1$ , if  $x_i$  belongs to cluster j;  $w_{ij} = 0$ , otherwise
- Looking for:
  - The best assignment  $w_{ij}$
  - The best center  $c_j$

## Solution of K-Means

$$J = \sum_{j=1}^{k} \sum_{i} w_{ij} ||x_i - c_j||^2$$

• Iterations

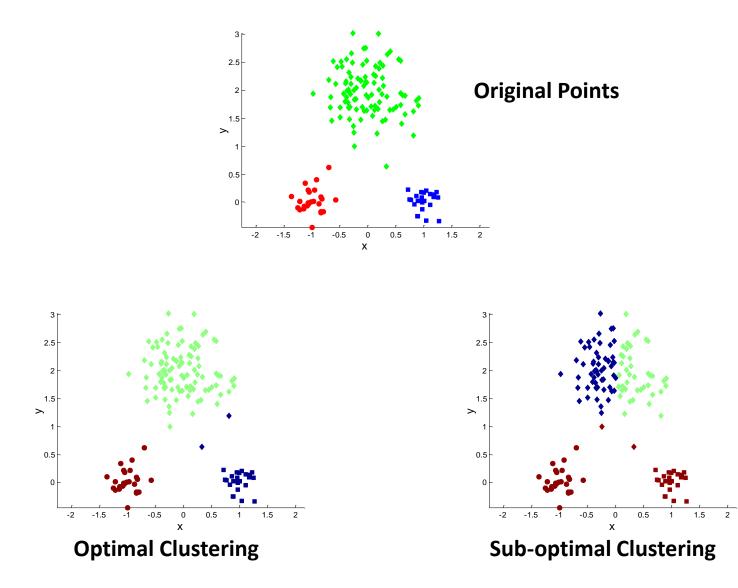
- Step 1: Fix centers  $c_j$ , find assignment  $w_{ij}$  that minimizes J• =>  $w_{ij} = 1$ ,  $if ||x_i - c_j||^2$  is the smallest
- Step 2: Fix assignment w<sub>ij</sub>, find centers that minimize J
  => first derivative of J = 0

• => 
$$\frac{\partial J}{\partial c_j} = -2 \sum_i w_{ij} (x_i - c_j) = 0$$
  
• => $c_j = \frac{\sum_i w_{ij} x_i}{\sum_i w_{ij}}$   
• Note  $\sum_i w_{ij}$  is the total number of objects in cluster j

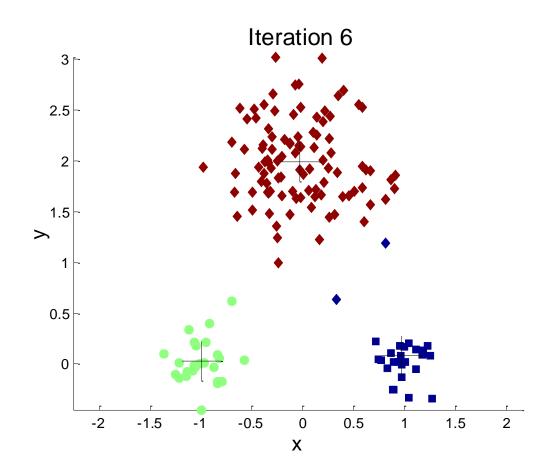
#### K-means Clustering – Details

- Initial centroids are often chosen randomly.
  - Clusters produced vary from one run to another.
- The centroid is (typically) the mean of the points in the cluster.
- 'Closeness' is measured by Euclidean distance, cosine similarity, correlation, etc.
- K-means will converge for common similarity measures
- Most of the convergence happens 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

#### Two different K-means Clusterings

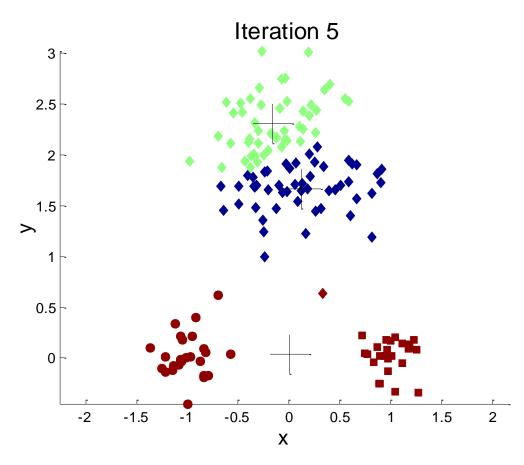


#### Importance of Choosing Initial Centroids



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## Importance of Choosing Initial Centroids



## Evaluating K-means Clusters

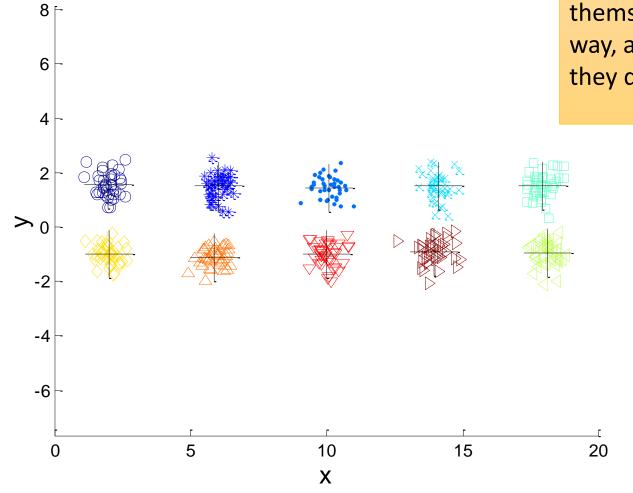
- Most common measure is Sum of Squared Error (SSE)
  - For each point, the error is the distance to the nearest cluster
  - To get SSE, we square these errors and sum them.

$$SSE = \sum_{i=1}^{K} \sum_{x \in C_i} dist^2(m_i, x)$$

- *x* is a data point in cluster *C*<sub>i</sub> and *m*<sub>i</sub> is the representative point for cluster *C*<sub>i</sub>
  - can show that *m<sub>i</sub>* corresponds to the center (mean) of the cluster
- Given two clusters, we can choose the one with the smallest error
- One easy way to reduce SSE is to increase K, the number of clusters
  - A good clustering with smaller K can have a lower SSE than a poor clustering with higher K

### 10 Clusters Example

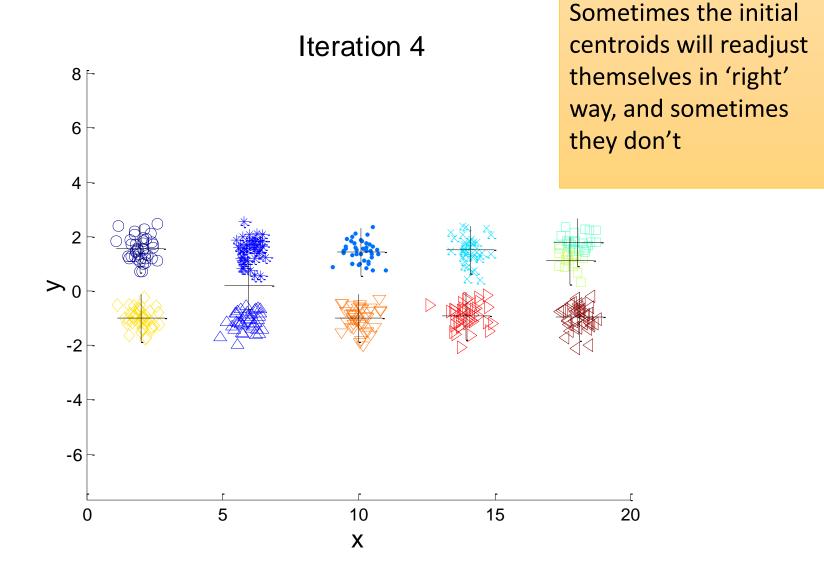
Sometimes the initial centroids will readjust themselves in 'right' way, and sometimes they don't



Iteration 4

Starting with two initial centroids in one cluster of each pair of clusters

#### 10 Clusters Example



Starting with some pairs of clusters having three initial centroids, while other have only one.

## Solutions to Initial Centroids Problem

- Multiple runs
  - Helps, but probability is not on your side
- 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
- Postprocessing
- Bisecting K-means
  - Not as susceptible to initialization issues

## **Bisecting K-means**

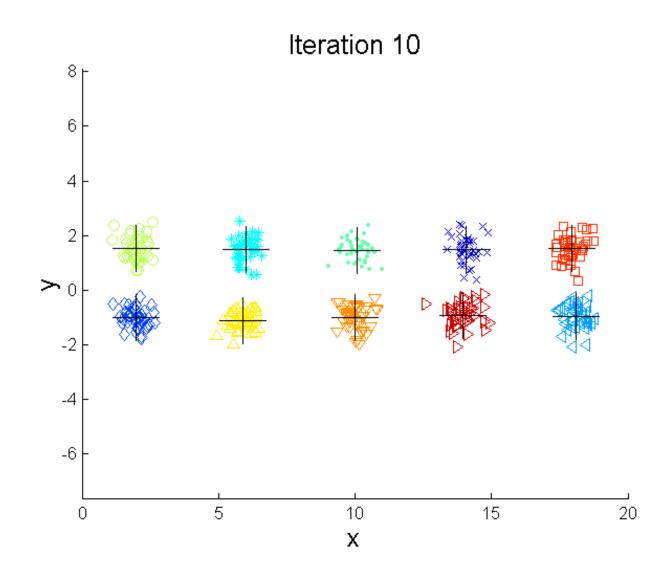
- Bisecting K-means algorithm
  - Variant of K-means that can produce a partitional or a hierarchical clustering

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

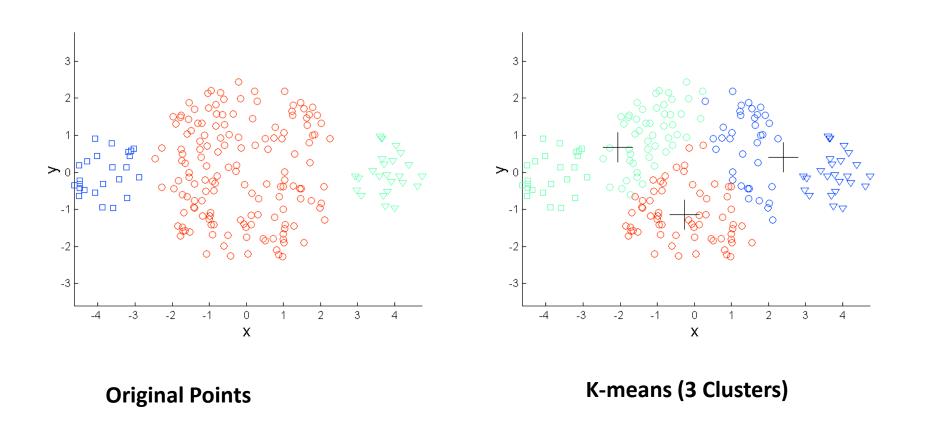
### Bisecting K-means Example



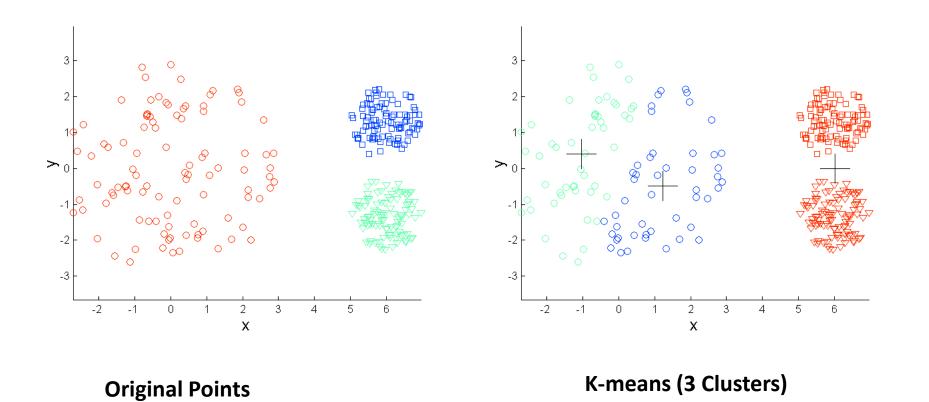
## Limitations of K-means

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

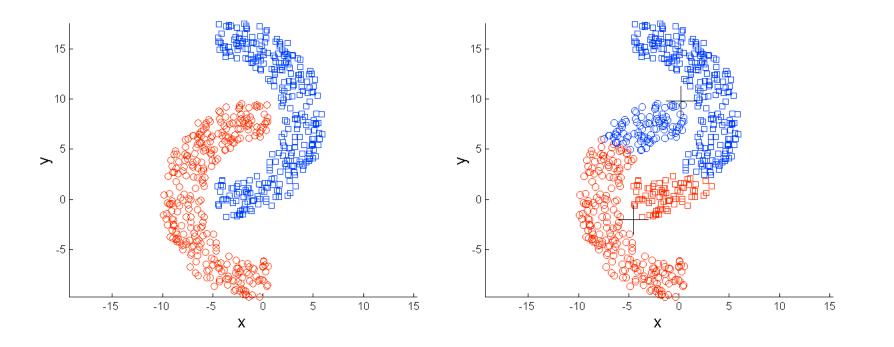
#### Limitations of K-means: Differing Sizes



# Limitations of K-means: Differing Density



#### Limitations of K-means: Non-globular Shapes

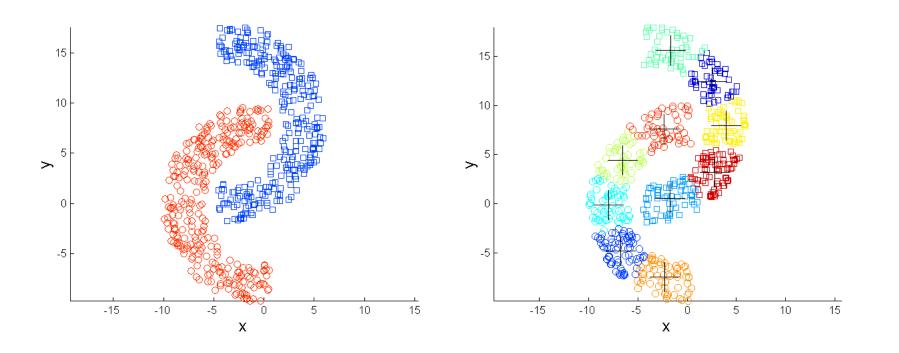


**Original Points** 

K-means (2 Clusters)

One solution is to use many clusters. Find parts of clusters, but need to put together.

#### **Overcoming K-means Limitations**



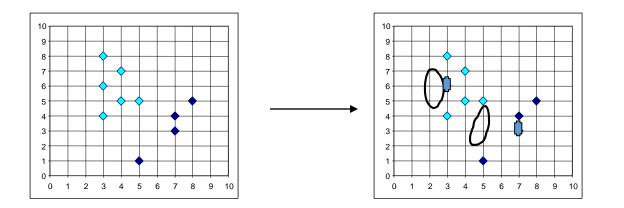
**Original Points** 

**K-means Clusters** 

One solution is to use many clusters. Find parts of clusters, but need to put together.

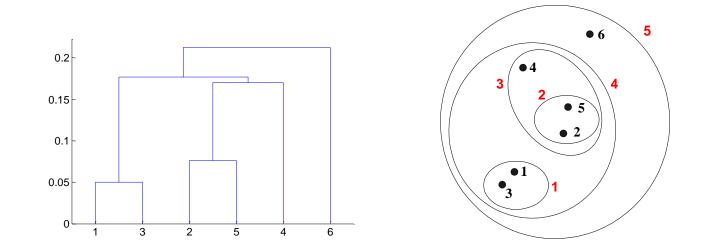
#### K-means is sensitive to outliers

- The k-means algorithm is sensitive to outliers !
  - Since an object with an extremely large value may substantially distort the distribution of the data
- K-Medoids: Instead of taking the mean value of the object in a cluster as a reference point, medoids can be used, which is the most centrally located object in a cluster



## **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

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

# **Hierarchical Clustering**

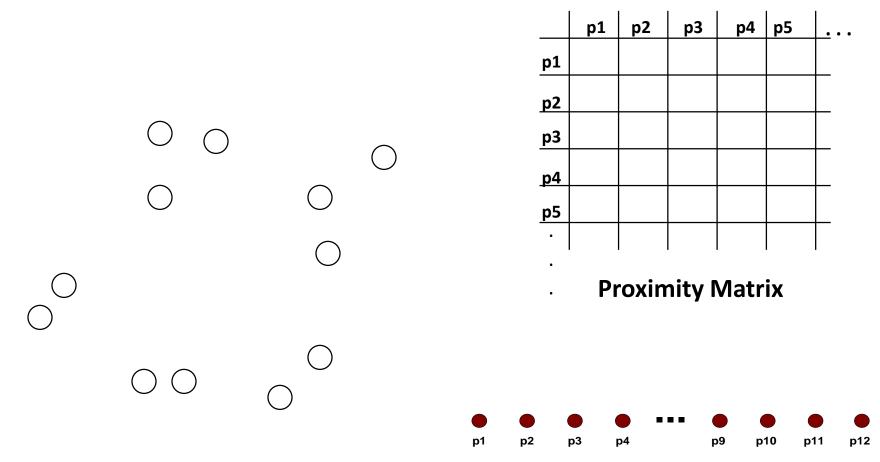
- Two main types of hierarchical clustering
  - Agglomerative:
    - Start with the points as individual clusters
    - At each step, merge the closest pair of clusters until only one cluster (or k clusters) left
  - Divisive:
    - Start with one, all-inclusive cluster
    - At each step, split a cluster 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
  - 5. Update the proximity matrix
  - 6. 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

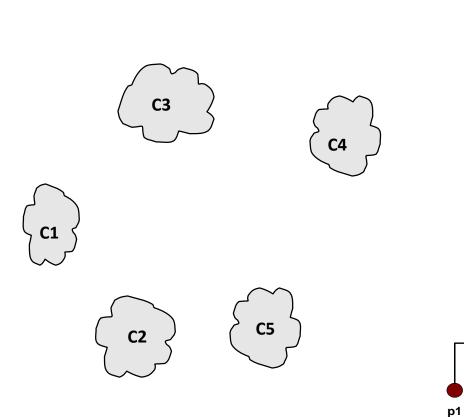
# Starting Situation

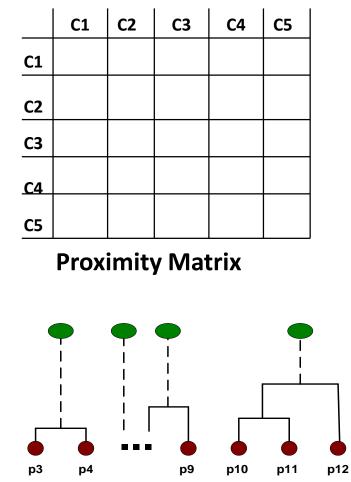
• Start with clusters of individual points and a proximity matrix



## Intermediate Situation

• After some merging steps, we have some clusters





p2

# Intermediate Situation

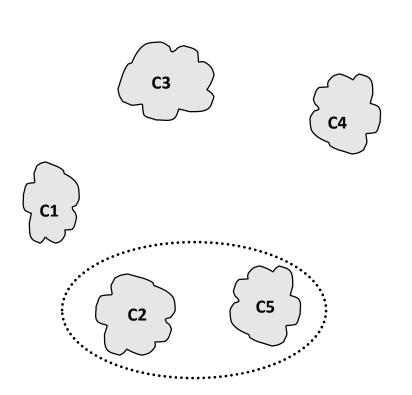
We want to merge the two closest clusters (C2 and C5) and update the proximity matrix.

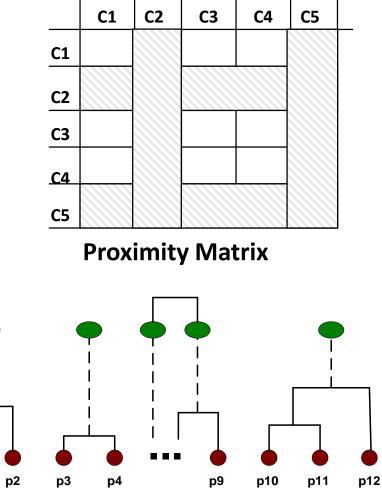
 <u>c1</u>

 <u>c2</u>

 <u>c3</u>

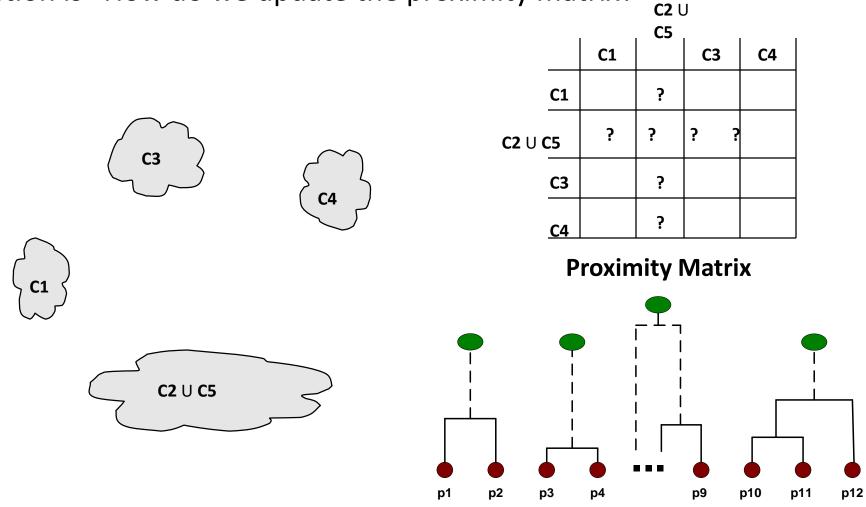
p1

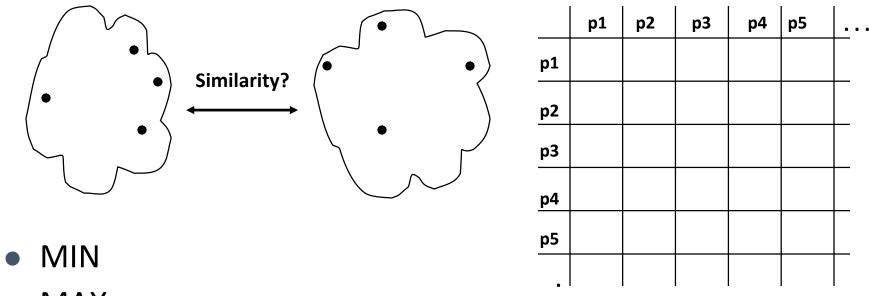




# After Merging

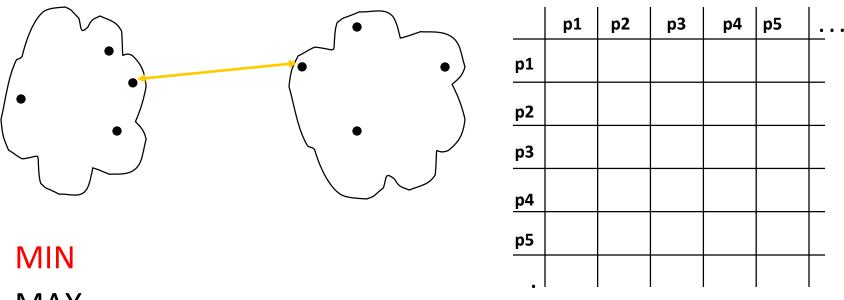
• The question is "How do we update the proximity matrix?"



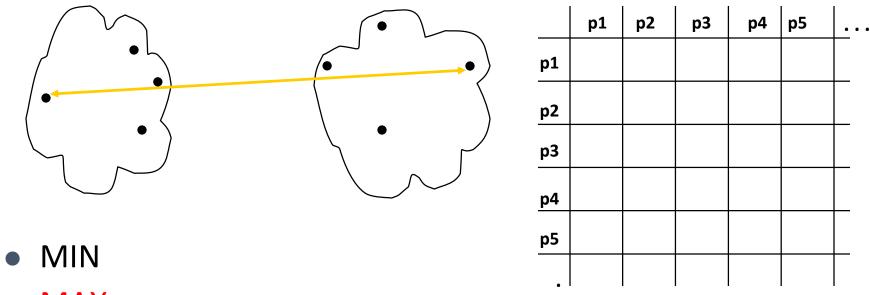


- MAX
- Group Average

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

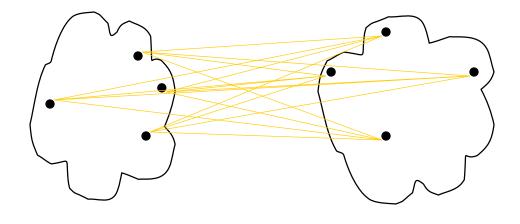


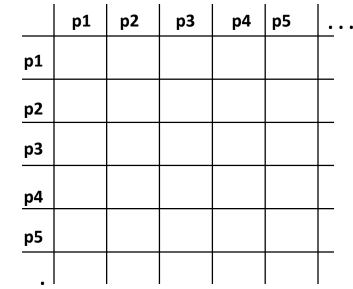
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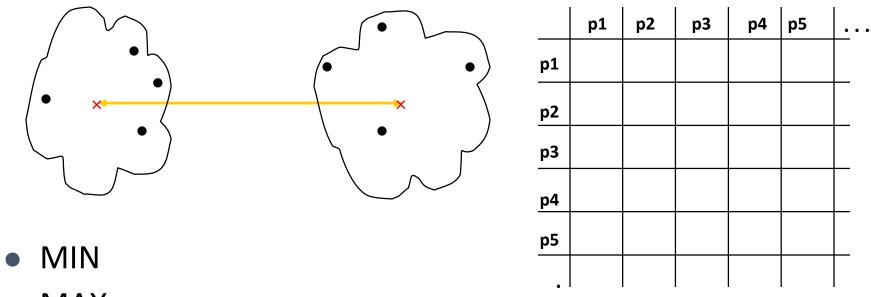
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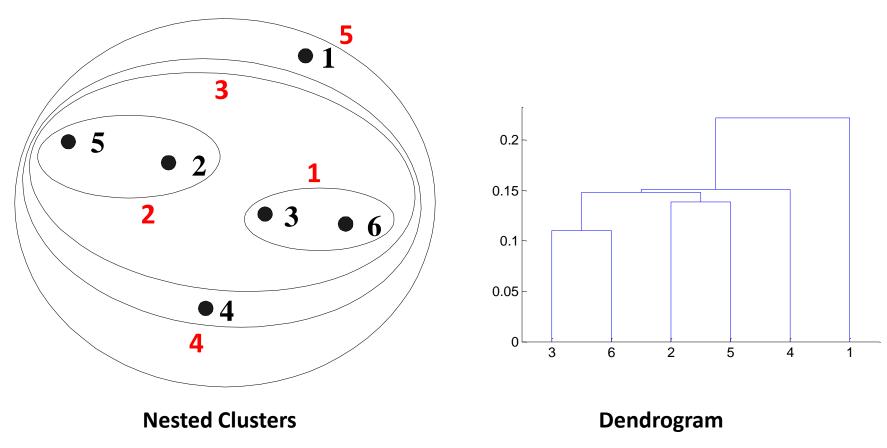
- MIN
- MAX
- Group Average
- Distance Between Centroids
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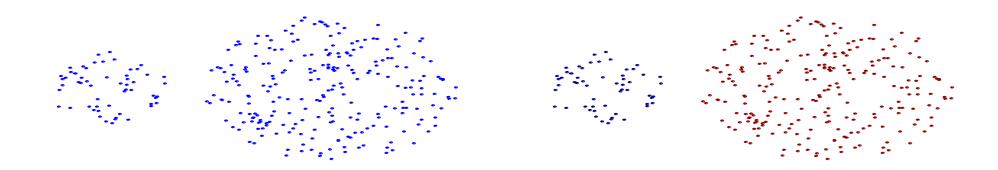
- MAX
- Group Average
- Distance Between Centroids
- Other methods driven by an objective function
  - Ward's Method uses squared error

## Hierarchical Clustering: MIN or Single Link

- Similarity of two clusters is based on the two most similar (closest) points in the different clusters
  - Determined by one pair of points, i.e., by one link in the proximity graph.



## Strength of MIN

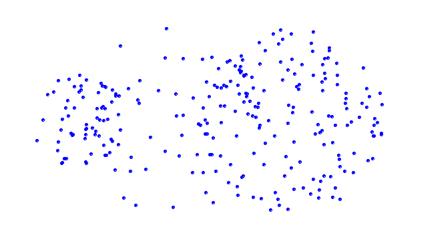


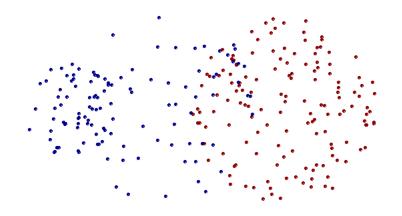
**Original Points** 

**Two Clusters** 

• Can handle non-elliptical shapes

## Limitations of MIN





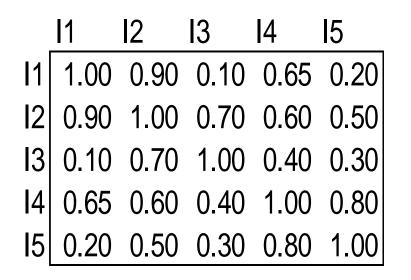
**Original Points** 

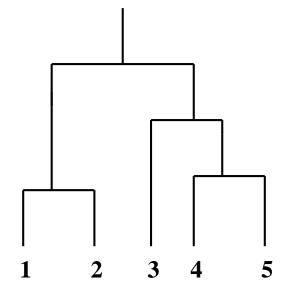
**Two Clusters** 

• Sensitive to noise and outliers

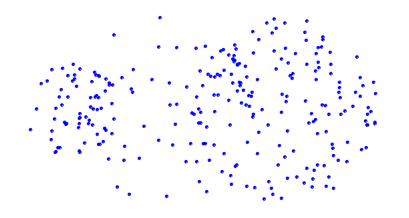
Cluster Similarity: MAX or Complete Linkage

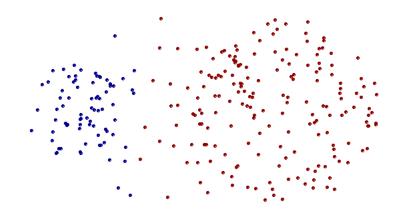
- Similarity of two clusters is based on the two least similar (most distant) points in the different clusters
  - Determined by all pairs of points in the two clusters





## Strength of MAX



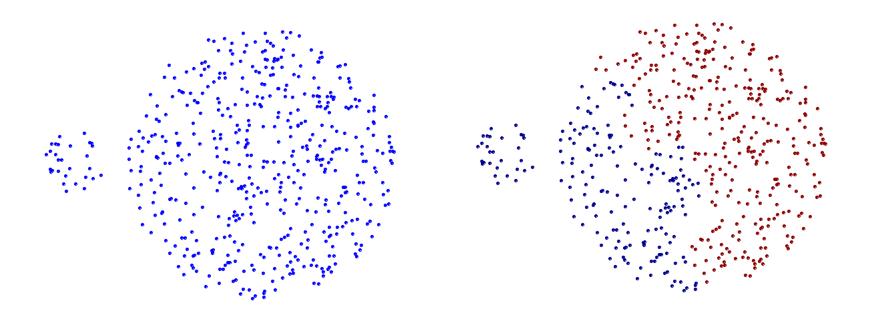


**Original Points** 

**Two Clusters** 

• Less susceptible to noise and outliers

## Limitations of MAX



**Original Points** 

#### **Two Clusters**

- •Tends to break large clusters
- Biased towards globular clusters

# Cluster Similarity: Group Average

• Proximity of two clusters is the average of pairwise proximity between points in the two clusters.

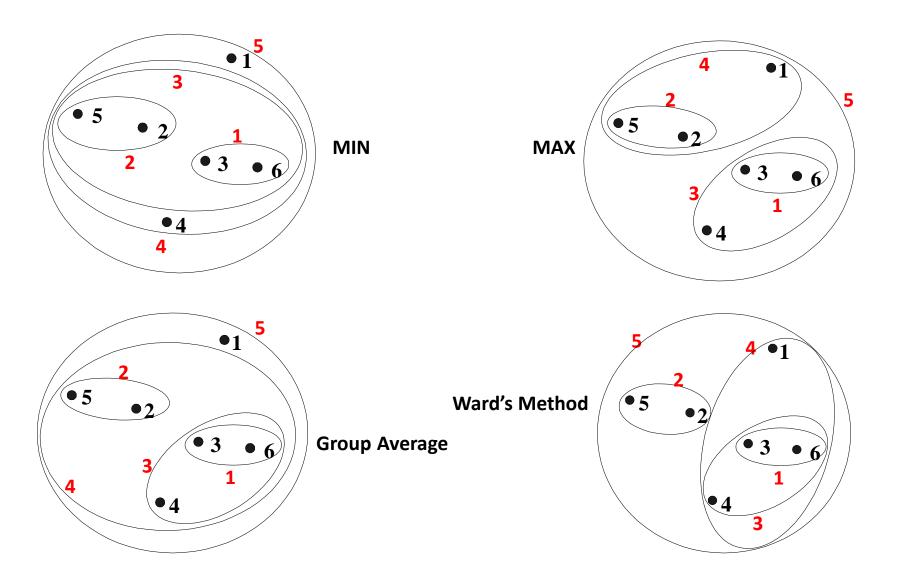
$$proximity(Cluster_{i}, Cluster_{j}) = \frac{\sum_{\substack{p_{i} \in Cluster_{i} \\ p_{j} \in Cluster_{j}}} \sum_{\substack{p_{i} \in Cluster_{i} \\ p_{j} \in Cluster_{j}}} \sum_{\substack{p_{i} \in Cluster_{i} \\ P_{i} \in Cluster_{i}}} \sum_{\substack{p_{i} \in$$

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

# 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
- Can be used to initialize K-means

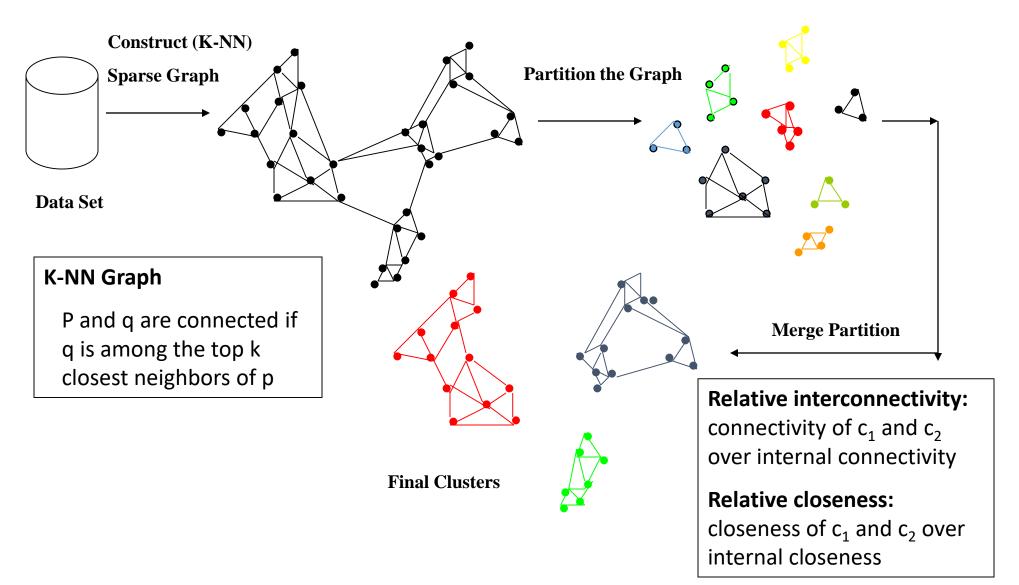
### **Hierarchical Clustering: Comparison**



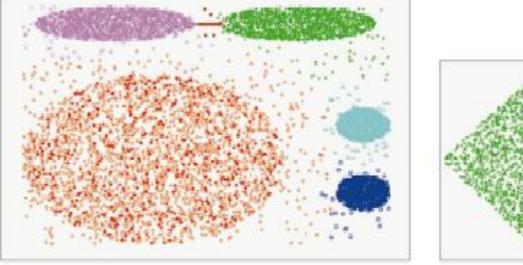
### CHAMELEON: Hierarchical Clustering Using Dynamic Modeling (1999)

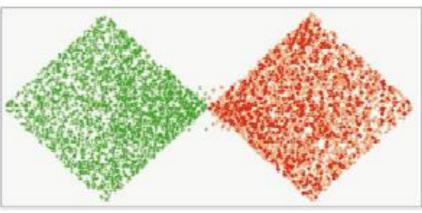
- CHAMELEON: G. Karypis, E. H. Han, and V. Kumar, 1999
- Measures the similarity based on a dynamic model
  - Two clusters are merged only if the *interconnectivity* and *closeness (proximity)* between two clusters are high *relative to* the internal interconnectivity of the clusters and closeness of items within the clusters
- Graph-based, and a two-phase algorithm
  - 1. Use a graph-partitioning algorithm: cluster objects into a large number of relatively small sub-clusters
  - 2. Use an agglomerative hierarchical clustering algorithm: find the genuine clusters by repeatedly combining these sub-clusters

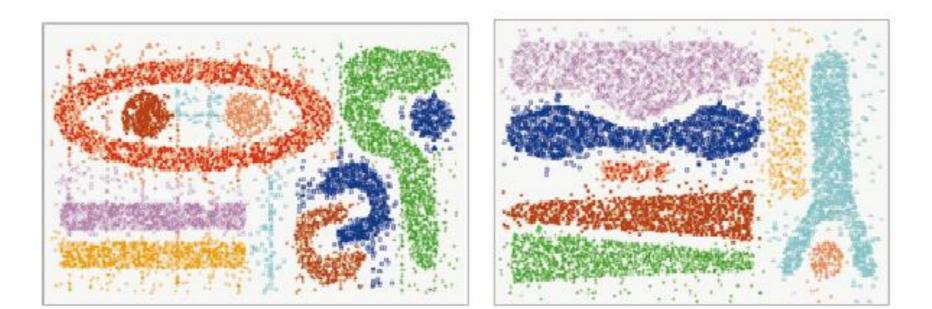
# **Overall Framework of CHAMELEON**



### **CHAMELEON (Clustering Complex Objects)**







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