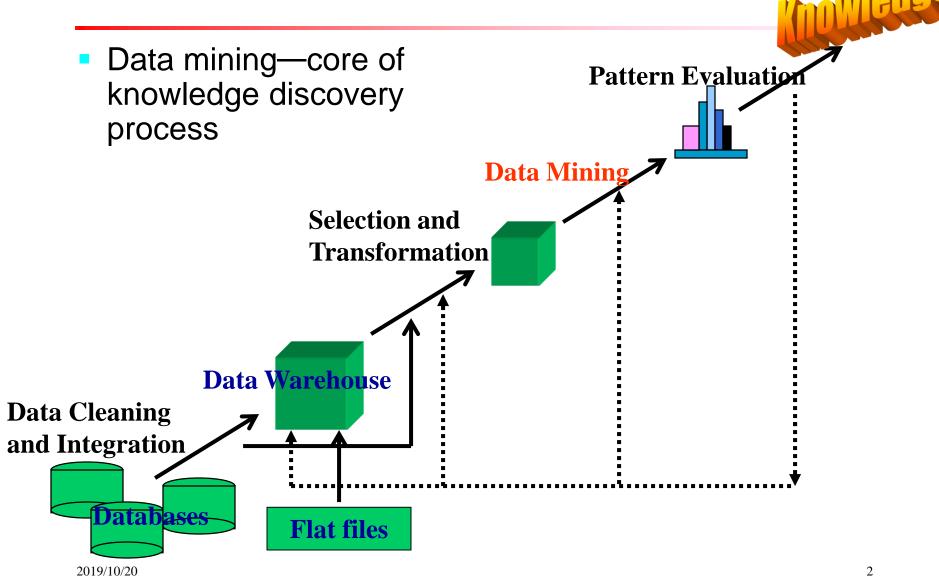
# **Data Mining**

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## **Review**



# **Cluster Analysis**

- What is Cluster Analysis?
- Types of Data in Cluster Analysis
- A Categorization of Major Clustering Methods
- Partitioning Methods
- Hierarchical Methods

- Density-Based Methods
- Grid-Based Methods
- Outlier Analysis
- Summary

## What is Cluster Analysis?

- Cluster: a collection of data objects
  - Similar to one another within the same cluster
  - Dissimilar to the objects in other clusters
- Cluster analysis
  - Finding similarities between data according to the characteristics in the data and grouping similar data objects into clusters
- Unsupervised learning: no predefined classes
- Typical applications
  - As a stand-alone tool to get insight into data distribution
  - As a preprocessing step for other algorithms

## **Examples of Clustering Applications**

- Marketing: Help marketers discover distinct groups according to their customer databases, and then use this knowledge to develop targeted marketing programs
- Land use: Identification of areas of similar land use in an earth observation database
- Insurance: Identifying groups of motor insurance policy holders with a high average claim cost
- City-planning: Identifying groups of houses according to their house type, value, and geographical location

## **Examples of Clustering Applications**

- Earth-quake studies: Observed earth quake epicenters are clustered along continent faults
- Biology: categorize genes with similar functionality
- - Document classification
  - Cluster Weblog data to discover groups of similar accessing patterns

# Clustering: Rich Applications and Multidisciplinary Efforts

- Pattern Recognition
- GIS
  - Create thematic maps in GIS by clustering feature spaces
  - Detect spatial clusters or for other spatial mining tasks
- Image Processing
- Economic Science (especially marketing research)
- Software package
  - S-Plus, SPSS, SAS, R

# What Is Good Clustering?

- A good clustering method will produce high quality clusters with
  - high intra-class similarity
  - low inter-class similarity
- The quality of a clustering result depends on both the similarity measure used by the method and its implementation
- The quality of a clustering method is also measured by its ability to discover some or all of the hidden patterns

## Measure the Quality of Clustering

- Dissimilarity/Similarity metric: Similarity is expressed in terms of a distance function, typically metric: d(i, j)
- The definitions of distance functions are usually very different for interval-scaled, boolean, categorical, ordinal ratio, and vector variables
- Weights may be assigned to different variables based on applications and data semantics
- It is hard to define "similar enough" or "good enough"
  - The answer is typically highly subjective

# **Requirements of Clustering**

- Scalability
- Ability to deal with different types of attributes
- Ability to handle dynamic data
- Discovery of clusters with arbitrary shape
- Minimal requirements for domain knowledge to determine input parameters
- Able to deal with noise and outliers
- Insensitive to order of input records
- High dimensionality
- Incorporation of user-specified constraints
- Interpretability and usability

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# **Cluster Analysis**

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## **Data Structures**

Data matrix

$$\begin{bmatrix} x_{11} & \cdots & x_{1f} & \cdots & x_{1p} \\ \cdots & \cdots & \cdots & \cdots \\ x_{i1} & \cdots & x_{if} & \cdots & x_{ip} \\ \cdots & \cdots & \cdots & \cdots \\ x_{n1} & \cdots & x_{nf} & \cdots & x_{np} \end{bmatrix}$$

Dissimilarity matrix 
$$\begin{bmatrix} 0 \\ d(2,1) & 0 \\ d(3,1) & d(3,2) & 0 \\ \vdots & \vdots & \vdots \\ d(n,1) & d(n,2) & \dots & 0 \end{bmatrix}$$

# Type of Data in Clustering Analysis

- Interval-scaled variables
- Binary variables
- Nominal variables
- Ordinal variables
- Ratio-scaled variables
- Variables of mixed types

## **Interval-valued Variables**

- Standardize data
  - Calculate the mean absolute deviation:

$$s_f = \frac{1}{n}(|x_{1f} - m_f| + |x_{2f} - m_f| + ... + |x_{nf} - m_f|)$$
 where 
$$m_f = \frac{1}{n}(x_{1f} + x_{2f} + ... + x_{nf})$$

Calculate the standardized measurement (z-score)

$$z_{if} = \frac{x_{if} - m_f}{s_f}$$

 Using mean absolute deviation is more robust than using standard deviation

# Similarity and Dissimilarity Between Objects

- Distances are normally used to measure the similarity or dissimilarity between two data objects
- Some popular ones include: Minkowski distance:

$$d(i,j) = \sqrt[q]{(|x_{i_1} - x_{j_1}|^q + |x_{i_2} - x_{j_2}|^q + ... + |x_{i_p} - x_{j_p}|^q)}$$
 where  $i = (x_{i_1}, x_{i_2}, ..., x_{i_p})$  and  $j = (x_{j_1}, x_{j_2}, ..., x_{j_p})$  are two  $p$ -dimensional data objects, and  $q$  is a positive integer

If q = 1, d is Manhattan distance

$$d(i,j) = |x_{i_1} - x_{j_1}| + |x_{i_2} - x_{j_2}| + ... + |x_{i_p} - x_{j_p}|$$

# Similarity and Dissimilarity Between Objects

If q = 2, d is Euclidean distance:

$$d(i,j) = \sqrt{(|x_{i1} - x_{j1}|^2 + |x_{i2} - x_{j2}|^2 + ... + |x_{ip} - x_{jp}|^2)}$$

- Properties
  - $d(i,j) \geq 0$
  - d(i,i) = 0
  - d(i,j) = d(j,i)
  - $d(i,j) \leq d(i,k) + d(k,j)$
- Also, one can use weighted distance

# **Binary Variables**

- A contingency table for binary data
- Distance measure for symmetric binary variables:
- Distance measure for asymmetric binary variables:

	Object j				
		1	0	sum	
Ohiaati	1	a	b	a+b	
Object i	0	c	d	c+d	
	sum	a+c	b+d	p	

$$d(i,j) = \frac{b+c}{a+b+c+d}$$

$$d(i,j) = \frac{b+c}{a+b+c}$$

## Dissimilarity between Binary Variables

#### Example

Name	Gender	Fever	Cough	Test-1	Test-2	Test-3	Test-4
Jack	M	Y	N	P	N	N	N
Mary	F	Y	N	P	N	P	N
Jim	M	Y	P	N	N	N	N

- gender is a symmetric attribute
- the remaining attributes are asymmetric binary
- let the values Y and P be set to 1, and the value N be set to 0

$$d(jack,mary) = \frac{0+1}{2+0+1} = 0.33$$

$$d(jack,jim) = \frac{1+1}{1+1+1} = 0.67$$

$$d(jim,mary) = \frac{1+2}{1+1+2} = 0.75$$

## **Nominal Variables**

- A generalization of binary variable in that it can take more than 2 states, e.g., red, yellow, blue, green
- Method 1: Simple matching
  - m: # of matches, p: total # of nominal variables

$$d(i,j) = \frac{p-m}{p}$$

- Method 2: use a large number of binary variables
  - creating a new binary variable for each of the M nominal states

## **Ordinal Variables**

- An ordinal variable can be discrete or continuous
- Order is important, e.g., rank
- Can be treated like interval-scaled
  - replace  $x_{if}$  by their rank  $r_{if} \in \{1, ..., M_f\}$
  - map the range of each variable onto [0, 1] by replacing i-th object in the f-th variable by

$$z_{if} = \frac{r_{if} - 1}{M_f - 1}$$

compute the dissimilarity using methods for interval-scaled variables

## **Ratio-Scaled Variables**

Ratio-scaled variable: a positive measurement on a nonlinear scale, approximately at exponential scale, such as Ae<sup>Bt</sup> or Ae<sup>-Bt</sup>

#### Methods:

- treat them like interval-scaled variables not a good choice!
- apply logarithmic transformation

$$y_{if} = log(x_{if})$$

 treat them as continuous data and treat their rank as intervalscaled

## Variables of Mixed Types

- A database may contain all the six types of variables
  - symmetric binary, asymmetric binary, nominal, ordinal, interval and ratio
- One may use a weighted formula to combine their effects  $d(i,j) = \frac{\sum_{f=1}^{p} \delta_{ij}^{(f)} d_{ij}^{(f)}}{\sum_{f=1}^{p} \delta_{ij}^{(f)}}$ 
  - $\delta^{(f)}_{ij} = 0$  if  $x_{if}$  or  $x_{jf}$  is missing, or  $x_{if} = x_{jf} = 0$  and f is asymmetric attribute; otherwise,  $\delta^{(f)}_{ij} = 1$
  - f is binary or nominal:  $d_{ij}^{(f)} = 0$  if  $x_{if} = x_{jf}$ , or  $d_{ij}^{(f)} = 1$  otherwise
  - f is interval-based: use the normalized distance
  - f is ordinal
    - compute ranks r<sub>if</sub> and treat z<sub>if</sub> as interval-scaled
  - f is ratio-scaled

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• Transform f, and treat f as interval-scaled.  $y_{if} = log(x_{if})$ 

 $Z_{if} = \frac{r_{if} - 1}{M_{c} - 1}$ 

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## **Exercise**

1. Please compute the dissimilarity matrix for the data set.

ID	Test-1	Test-2	Test-3
	(categorical)	(ordinal)	(ratio-scaled)
1	Α	excellent	445
2	В	fair	22
3	С	good	164
4	Α	excellent	1,210

## **Solution**

## For test-1, use simple matching

0			
d(2,1)	0		
d(3,1)	d(3,2)	0	
d(4,1)	d(4,2)	d(4,3)	0

=

0			
1	0		
1	1	0	
0	1	1	0

For test-2

0			
1	0		
0.5	0.5	0	
0	1	0.5	0

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## **Solution**

## For test-3, use log transformation

- Convert test-3 to 2.65, 1.34, 2.21, 3.08
- Normalize to 0.75, 0, 0.5,1

0			
0.75	0		
0.25	0.5	0	
0.25	1	0.5	0

## Dissimilarity matrix

0			
0.92	0		
0.58	0.67	0	
0.08	1	0.67	0

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## **Major Clustering Approaches (I)**

#### Partitioning approach:

- Construct various partitions and then evaluate them by some criterion, e.g., minimizing the sum of square errors
- Typical methods: k-means, k-medoids, CLARANS

#### Hierarchical approach:

- Create a hierarchical decomposition of the set of data (or objects) using some criterion
- Typical methods: Diana, Agnes, BIRCH, ROCK, CHAMELEON

#### Density-based approach:

- Based on connectivity and density functions
- Typical methods: DBSACN, OPTICS, DenClue

## **Major Clustering Approaches (II)**

- Grid-based approach:
  - based on a multiple-level granularity structure
  - Typical methods: STING, WaveCluster, CLIQUE
- Probabilistic Model-based approach:
  - Typical methods: EM

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# Centroid, Radius and Diameter of a Cluster (for numerical data sets)

Centroid: the "middle" of a cluster

$$C = \frac{\sum_{i=1}^{N} (t_i)}{N}$$

Radius: square root of average distance from any point of the cluster to its centroid

$$R = \sqrt{\frac{\sum_{i=1}^{N} (t_i - c)^2}{N}}$$

 Diameter: square root of average mean squared distance between all pairs of points in the cluster

$$D = \sqrt{\frac{\sum_{i=1}^{N} \sum_{j=1}^{N} (t_i - t_j)^2}{N(N-1)}}$$

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## **Partitioning Algorithms: Basic Concept**

Partitioning method: Construct a partition of a database D of n objects into a set of k clusters, s.t., min sum of squared distance

$$\sum_{m=1}^{k} \sum_{t_{mi} \in Km} (C_m - t_{mi})^2$$

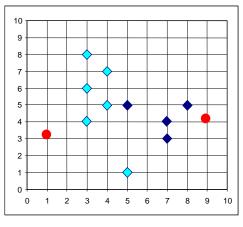
- Given a k, find a partition of k clusters that optimizes the chosen partitioning criterion
  - k-means (MacQueen'67): Each cluster is represented by the center of the cluster
  - k-medoids or PAM (Partition around medoids) (Kaufman & Rousseeuw'87): Each cluster is represented by one of the objects in the cluster

## The K-Means Clustering Method

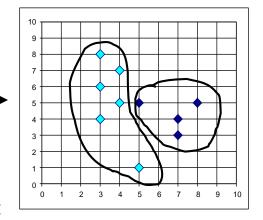
- Given a k, the k-means algorithm is implemented in four steps:
  - Give k random seeds as the initial centroids
  - Compute the centroid of each cluster of the current partition (the centroid is the center, i.e., mean point)
  - For each object, compute its distance to the centroids
    - Assign it to the cluster with the nearest centroid
  - Go back to Step 2, stop when no more new assignment

# The K-Means Clustering Method

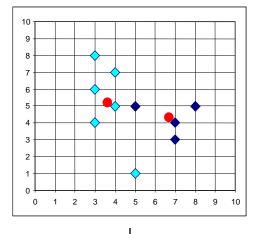
## Example



Assign each objects to most similar center

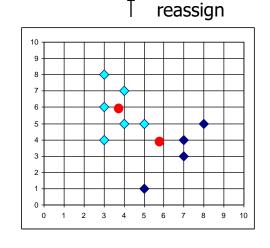


Update the cluster means

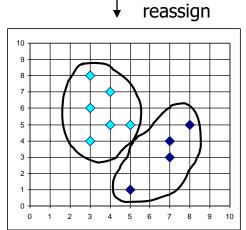


K=2

Arbitrarily choose K object as initial cluster center



Update the cluster means



## Comments on the *K-Means* Method

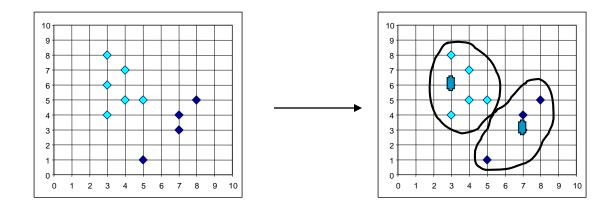
- Strength: Relatively efficient: O(tkn), where n is # objects, k is # clusters, and t is # iterations. Normally, k, t << n</p>
- Comment: Often terminates at a local optimum. The global optimum may be found using techniques such as: deterministic annealing and genetic algorithms
- Weakness
  - Applicable only when mean is defined, then what about categorical data?
  - Need to specify k, the number of clusters, in advance
  - Unable to handle noisy data and outliers
  - Not suitable to discover clusters with non-convex shapes

## Variations of the *K-Means* Method

- Handling categorical data: k-modes
  - Replacing means of clusters with modes
  - Using new dissimilarity measures to deal with categorical objects
  - Using a frequency-based method to update modes of clusters
  - A mixture of categorical and numerical data: k-prototype method
- Expectation Maximization: an extension to k-means
  - Assign each object to a cluster according to a weight (prob.)
  - New means are computed based on weighted measures

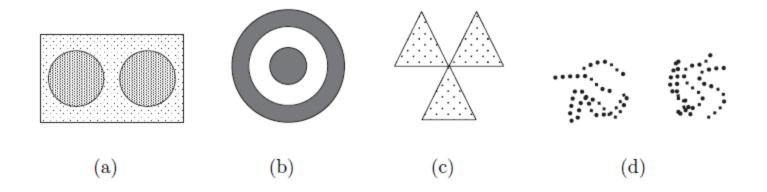
#### What Is the Problem of the K-Means Method?

- 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 objects in a cluster as a reference point, medoids can be used, which is the most centrally located object in a cluster



#### **Exercise**

1. Identify the clusters using the K-means (using squared error as the objective function). Note that darkness or the number of dots indicates density.



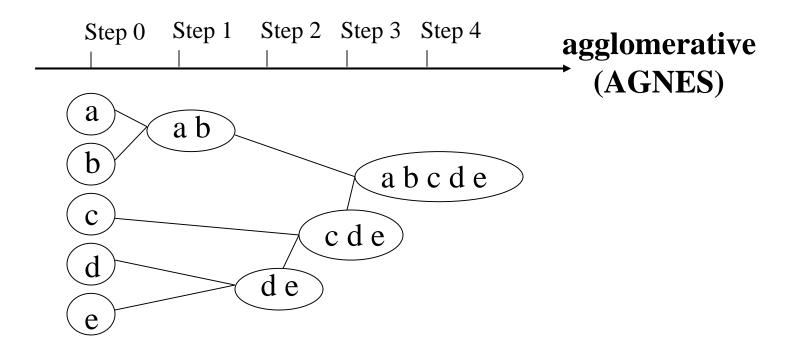
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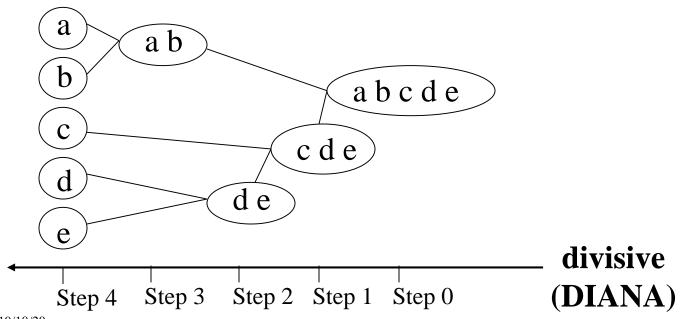
# **Hierarchical Clustering**

This method does not require the number of clusters k as an input, but needs a termination condition



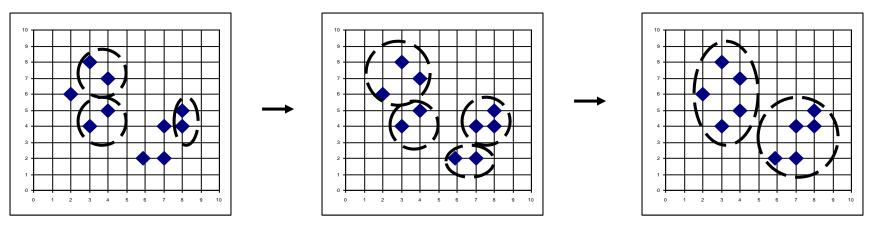
## **Hierarchical Clustering**

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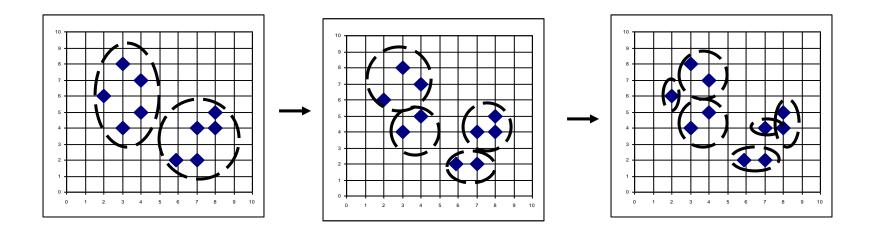
## **AGNES** (Agglomerative Nesting)

- Introduced by Kaufmann and Rousseeuw (1990)
- Use the Single-Link method and the dissimilarity matrix
- Merge nodes that have the least dissimilarity
- Go on in a non-descending fashion
- Eventually all nodes belong to the same cluster



## **DIANA** (Divisive Analysis)

- Introduced by Kaufmann and Rousseeuw (1990)
- Inverse order of AGNES
- Eventually each node forms a cluster on its own



# Recent Hierarchical Clustering Methods

- Major weakness of hierarchical clustering methods
  - do not scale well: time complexity of at least  $O(n^2)$ , where n is the number of total objects
  - can never undo what was done previously
- Integration of hierarchical with distance-based clustering
  - BIRCH (1996): uses CF-tree and incrementally adjusts the quality of sub-clusters
  - ROCK (1999): clustering categorical data by neighbor and link analysis
  - CHAMELEON (1999): hierarchical clustering using dynamic modeling

## **BIRCH (1996)**

- BIRCH: integrated hierarchical clustering
- Clustering feature, Clustering feature tree
- Incrementally construct a CF (Clustering Feature) tree, a hierarchical data structure for multiphase clustering
  - Phase 1: scan DB to build an initial in-memory CF tree (a multi-level compression of the data that tries to preserve the inherent clustering structure of the data)
  - Phase 2: use a clustering algorithm to cluster the leaf nodes of the CF-tree

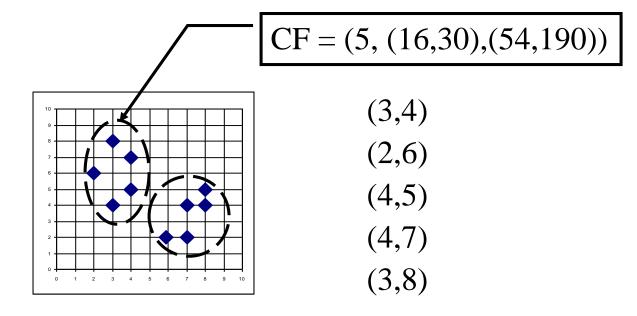
# **Clustering Feature Vector in BIRCH**

Clustering Feature: CF = (N, LS, SS), summarize the cluster members

#### N: Number of data points

LS: 
$$\sum_{i=1}^{N} = X_i$$

SS: 
$$\sum_{i=1}^{N} = \overline{X_i^2}$$



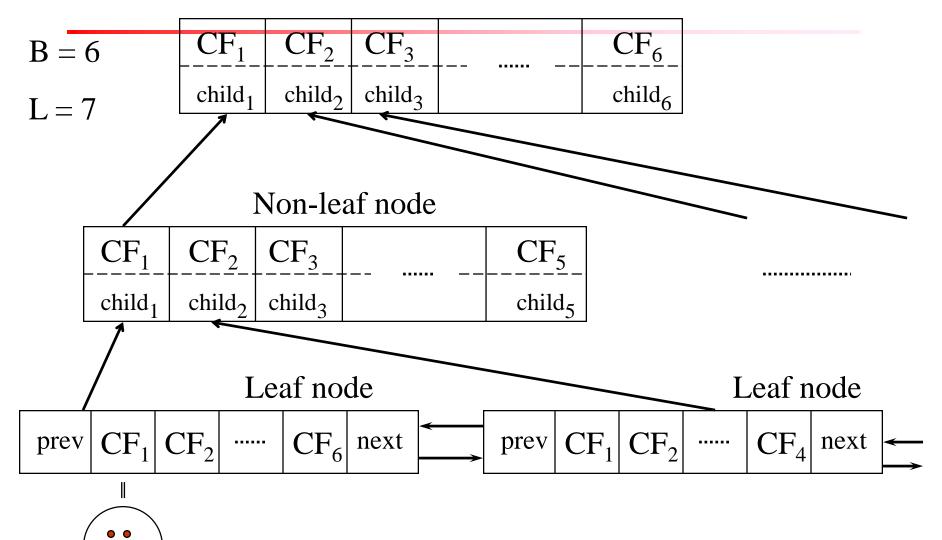
### **CF-Tree in BIRCH**

#### Clustering feature:

- summary of the statistics for a given subcluster
- registers crucial measurements for computing cluster and utilizes storage efficiently
- A CF tree is a height-balanced tree that stores the clustering features for a hierarchical clustering
  - A nonleaf node in a tree has descendants or "children"
  - The nonleaf nodes store sums of the CFs of their children
- A CF tree has two parameters
  - Branching factor: specify the maximum number of children
  - Threshold: max diameter of sub-clusters stored at the leaf nodes

### The CF Tree Structure

Root



#### **BIRCH**

#### Phase 1

- Insert each object to its closest leaf entry
- If the diameter of a leaf is larger than a threshold, the leaf will be split
- Update the CF and its ancestor's CF
- If the size of the CF tree is too big, re-build the tree from the leaf node, no re-scan the original objects
- Two parameters (branching factor, threshold), control the size of the tree

#### **BIRCH**

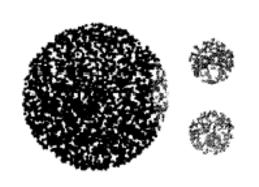
#### Scales linearly

- Complexity: O(n)
- Scalable for large database
- Incremental clustering
- Finds a good clustering with a single scan, I/O cost small

#### Weakness

- Handles only numeric data, and sensitive to the order of the data record
- Not good at arbitrary shaped cluster

Centroid-based clustering All-points agglomerative clustering





CURE: middle ground between centroid-based clustering and all-points agglomerative clustering

- Start with each individual point as a separate cluster
- Merge closest clusters till each cluster contains more than c points
- For each cluster, use c scattered points as representatives
- If more than k clusters
  - Clusters with the closest pair of representative points are merged
  - Update the representative points of merged clusters

#### Choose representatives

- the point farthest from the mean of the cluster
- for 2 to c do
   the point farthest from the previously chosen point
- Shrink the scattered points toward the mean by a fraction  $\alpha$

for each scattered point p do representative =  $p + \alpha$  \* (mean – p)

#### Merge

Euclidian distance

$$dist(u, v) = \min_{p \in u. \text{rep}, q \in v. \text{rep}} dist(p, q)$$

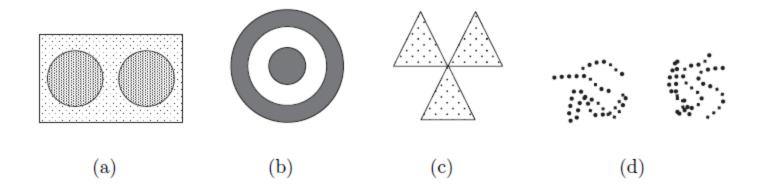
 Closest clusters — minimum distance between representative points from two clusters

- Multiple representative points allow CURE to discover arbitrary shaped clusters
- Less sensitive to outliers
  - Shrink scattered points toward the mean, weaken the effects of outliers
- Time complexity  $O(n^2 \log(n))$
- For large-scale database, do sampling and partitioning

- Draw a random sampling S from original objects
- Cluster the sampled objects (basic CURE)
- Eliminate outliers
- Each unsampled original object is assigned to the cluster containing the closest representative point to it

## **Exercise**

1. Identify the clusters using the Single-Link method. Note that darkness or the number of dots indicates density.



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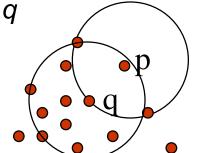
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## **Density-Based Clustering Methods**

- Clustering based on density (local cluster criterion), such as density-connected points
- Major features:
  - Discover clusters of arbitrary shape
  - Handle noise
  - Need density parameters as termination condition
  - Complexity is O(n²)

# **Density-Based Clustering: Basic Concepts**

- Two parameters:
  - $\varepsilon$ -neighborhood: neighborhood within a radius  $\varepsilon$  of a point
  - *MinPts*: Min number of points in  $\varepsilon$ -neighborhood of a point
- core object: If the number of points in  $\varepsilon$ -neighborhood of point p exceeds MinPts
- Directly density-reachable: A point p is directly density-reachable from a point q w.r.t.,  $\varepsilon$ , MinPts if
  - p belongs to ε-neighborhood of q
  - q is core object



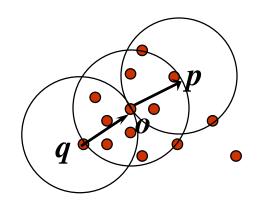
MinPts = 5

$$\varepsilon = 1 \text{ cm}$$

## Density-Reachable and Density-Connected

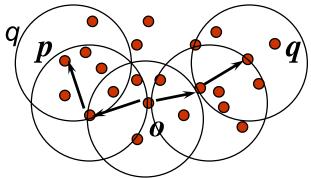
#### Density-reachable:

• A point p is density-reachable from a point q w.r.t.  $\varepsilon$ , MinPts if there is a chain of points  $p_1, \ldots, p_n, p_1 = q, p_n = p$  such that  $p_{i+1}$  is directly density-reachable from  $p_i$ 



#### Density-connected

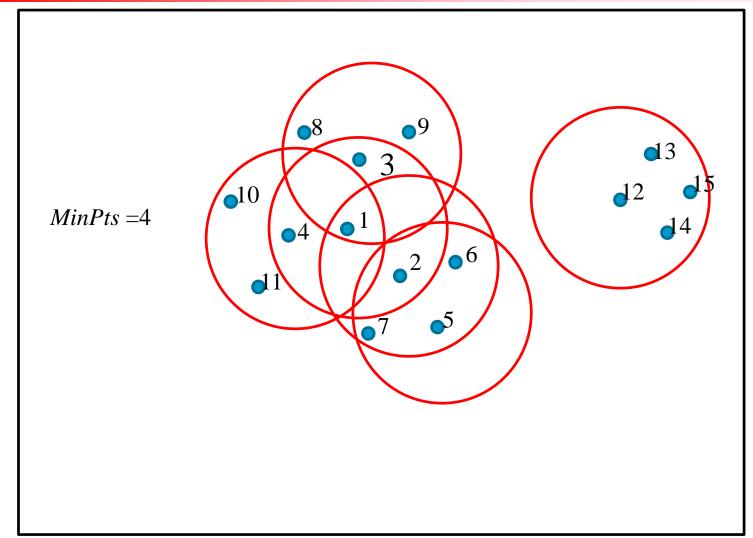
A point p is density-connected to a point q w.r.t. ε, MinPts if there is a point o such that both p and q are density-reachable from o w.r.t. ε and MinPts



# **DBSCAN: The Algorithm**

```
(1) mark all the objects as unvisited;
(2) do
(3)
        randomly select an unvisited object p;
(4)
        mark p as visited;
(5)
        if the \varepsilon-neighborhood of p has at least MinPts objects
             create a new cluster C, and add p to C;
(6)
(7)
            let N be the set of objects in the \varepsilon-neighborhood of \boldsymbol{p};
(8)
            for each point p' in N
                 if p' is unvisited
(9)
(10)
                      mark p' as visited;
(11)
                      if the \varepsilon-neighborhood of p' has at least MinPts points, add
those points to N;
(12)
                 if p' is not yet a member of any cluster, add p' to C;
            end for
(13)
(14)
            output C;
       else mark p as noise;
(15)
(16) until no object is unvisited.
```

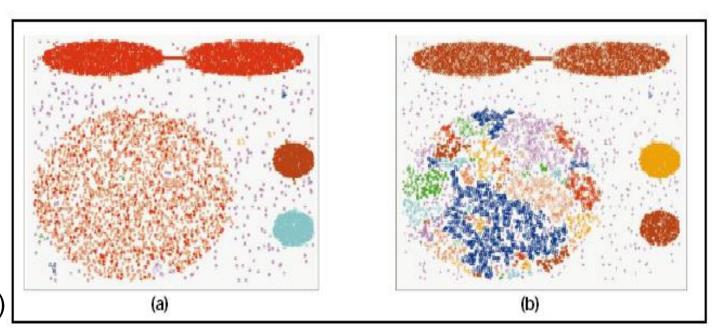
# **DBSCAN: The Algorithm**

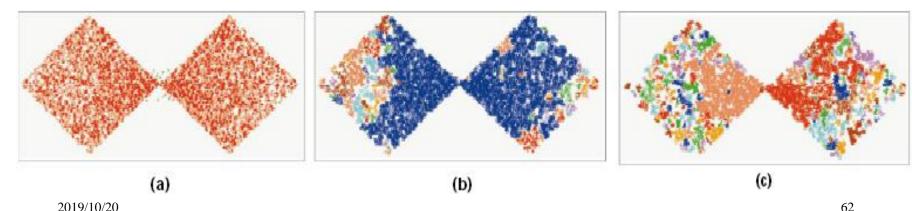


### **DBSCAN: Sensitive to Parameters**

DBScan results with MinPts = 4 and  $\epsilon$ = (a) 0.5, (b) 0.4

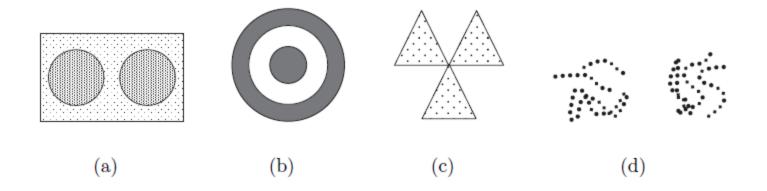
DBScan results with MinPts = 4 and  $\epsilon$ = (a) 5.0, (b) 4.0, (c) 3.0





## **Exercise**

1. Identify the clusters using DBSCAN. Note that darkness or the number of dots indicates density.



# **Cluster Analysis**

- What is Cluster Analysis?
- Types of Data in Cluster Analysis
- A Categorization of Major Clustering Methods
- Partitioning Methods
- Hierarchical Methods

- Density-Based Methods
- Grid-Based Methods
- Outlier Analysis
- Summary

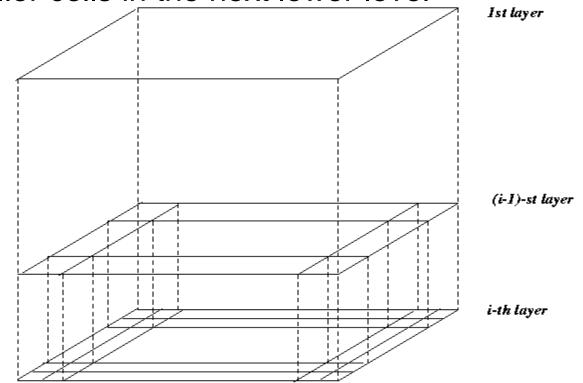
# **Grid-Based Clustering Method**

- Using multi-resolution grid data structure
- Several interesting methods
  - STING (a STatistical INformation Grid approach) by Wang,
     Yang and Muntz (1997)
  - WaveCluster by Sheikholeslami, Chatterjee, and Zhang (VLDB'98)
    - A multi-resolution clustering approach using wavelet method
  - CLIQUE: Agrawal, et al. (SIGMOD'98)
    - On high-dimensional data

# STING: A Statistical Information Grid Approach

- The spatial area is divided into rectangular cells
- There are several levels of cells corresponding to different levels of resolution

Each cell at a high level is partitioned into a number of smaller cells in the next lower level



# The STING Clustering Method

- Statistical info of each cell is calculated and stored beforehand and is used to answer queries
  - count, mean, s, min, max
  - type of distribution—normal, uniform, NONE, etc.
- Clusters are identified based on count, cell size, etc.

$$n = \sum_{t} n_{t}$$

$$m = \frac{\sum_{t} m_{t} n_{t}}{n}$$

$$s = \sqrt{\frac{\sum_{t} (s_{t}^{2} + m_{t}^{2}) n_{t}}{n} - m^{2}}$$

$$min = \min_{t} (min_{t})$$

$$max = \max_{t} (max_{t})$$

# The STING Query Method

- Use a top-down approach to answer spatial data queries
- Start from a pre-selected layer typically with a small number of cells
- For each cell in the current level compute the confidence interval
- Remove the irrelevant cells from further consideration
- When finish examining the current layer, proceed to the next lower level of the relevant cells
- Repeat this process until the bottom layer is reached

#### **Exercise**

- Please give some comments on STING in the following aspects:
- (1) cluster shape
- (2) computational complexity
- (3) cluster quality
- (4) incremental clustering

## **Comments on STING Clustering**

#### Advantages:

- Query-independent
- incremental update
- O(K) for query, where K is the number of grid cells at the lowest level
- O(n) for generating clusters

#### Disadvantages:

- All the cluster boundaries are either horizontal or vertical, and no diagonal boundary is detected
- Processing time depends on the size of each grid

# **Cluster Analysis**

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Density-Based

#### Methods

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# What Is Outlier Discovery?

- What are outliers?
  - The set of objects are considerably dissimilar from the remaining of the data
  - Caused by
    - Measurement or execution errors
    - Result of inherent variability
- Mining outliers is valuable
- Applications:
  - Credit card fraud detection
  - Customer segmentation
  - Medical analysis

#### **Outlier Detection**

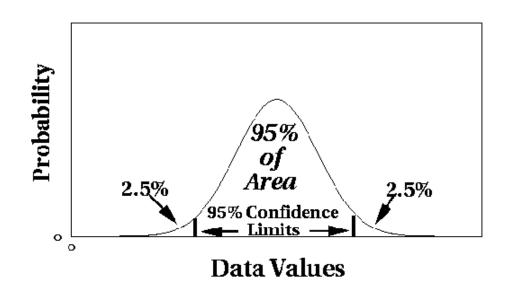
#### Visualization

- Weak in data with categorical data, high dimensional data
- Good at numerical data of 2 or 3 dimensions
- Clustering
  - Byproduct of clustering may be outliers
- Computer-based methods
  - Statistical-based outlier detection
  - Distance-based outlier detection
  - Deviation-based outlier detection

## **Outlier Detection: Statistical Approaches**

- Assume a distribution (e.g. normal distribution) for the data set and then use discordancy test to find outliers
- Discordancy tests depends on knowledge
  - data distribution
  - two hypothesis
  - distribution parameter (e.g., mean, variance)
  - number of expected outliers

## **Outlier Detection: Statistical Approaches**



#### Drawbacks

- Most tests are for single attribute
- In many cases, data distribution may not be known
- Require input parameters

## **Outlier Detection: Distance-Based Approach**

- Introduced to overcome the main limitations imposed by statistical methods
  - We need multi-dimensional analysis without knowing data distribution, no statistical test
- Distance-based outlier: A DB(p, d)-outlier is an object
   O in a dataset T such that at least a fraction p of the objects in T lies at a distance greater than d from O
- Algorithms for mining distance-based outliers
  - Index-based algorithm
  - Cell-based algorithm

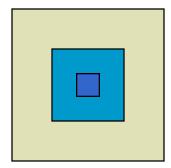
## **Index-based Algorithm**

- Search for neighbors of each object O within radius d around the object
- Multi-dimensional index structure, e.g. kd tree
- Max number of objects within d-neighborhood of each outlier
- The worst case O(kn²)
  k dimensionality, n number of objects
- Drawbacks:
  - Tree building is computational intensive

## **Cell-based Algorithm**

#### Cell partition

- Partition data space into cells, side length d/2k<sup>1/2</sup>
- Each cell has two layers around it
  - First layer one cell thick
  - Second layer (2k<sup>1/2</sup>-1) cells thick



#### Outlier detection

- If count of the first layer >M, no outlier in this cell
- If count of the second layer<=M, all objects are outliers</p>
- Otherwise, examine every object in the cell
- Good for large-scale data set

## **Cell-based Algorithm**

2	2	2	2	2	2	2
2 /	2	2	2	2	2	2
2	2	1	2 1 C 1 2	1	2	2
2	2	1	C	1,	2	2
2	2	1 1 2	1	1,	2	2
2 2 2 2 2 2	2		2	2	2	2
2	2	2	2	2	2	2

- *k*=2, *k* dimensionality
- Layer-1 property: given any point x in cell C, and any point y in layer-1 cell, dist(x,y) ≤ d
- Layer-2 property: given any point x in cell C, and any point y outside layer-2 cell, dist(x,y) > d

## **Outlier Detection: Deviation-Based Approach**

- Identifies outliers by examining the main characteristics of objects in a group
- Objects that "deviate" from this description are considered outliers
- Sequential exception technique
  - Simulates the way in which humans can distinguish unusual objects from among a series of supposedly like objects
  - A sequence of subsets,  $\{S_1, S_2, ..., S_m\}$ ,  $S_{j-1} \subset S_j$
  - Calculate the dissimilarity difference between the current subset with the proceeding subset in the sequence

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## **Summary**

- Cluster analysis groups objects based on their similarity and has wide applications
- Measure of similarity can be computed for various types of data
- Clustering algorithms can be categorized into partitioning methods, hierarchical methods, densitybased methods, grid-based methods
- Outlier detection and analysis are very useful for fraud detection, etc. and can be performed by statistical, distance-based or deviation-based approaches

## Summary

- Considerable progress has been made in scalable clustering methods
  - Partitioning: k-means, k-medoids, CLARANS
  - Hierarchical: BIRCH, ROCK, CHAMELEON
  - Density-based: DBSCAN, OPTICS, DenClue
  - Grid-based: STING, WaveCluster, CLIQUE
  - Model-based: EM, Fuzzy K-Means
  - Frequent pattern-based: pCluster
  - Constraint-based: COD, constrained-clustering