



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
- Model-Based Clustering Methods
- Outlier Analysis
- Summary

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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
 - Grouping a set of data objects into clusters
- Clustering is unsupervised classification: no predefined classes
- Typical applications
 - As a stand-alone tool to get insight into data distribution
 - As a preprocessing step for other algorithms

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General Applications of Clustering



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- Pattern Recognition
- Spatial Data Analysis
 - create thematic maps in GIS by clustering feature spaces
 - detect spatial clusters and explain them in spatial data mining
- Image Processing
- Economic Science (especially market research)
- WWW
 - Document classification
 - Cluster Weblog data to discover groups of similar access patterns



Examples of Clustering Applications



- <u>Marketing:</u> Help marketers discover distinct groups in their customer bases, and then use this knowledge to develop targeted marketing programs
- <u>Land use:</u> Identification of areas of similar land use in an earth observation database
- <u>Insurance:</u> Identifying groups of motor insurance policy holders with a high average claim cost
- <u>City-planning:</u> Identifying groups of houses according to their house type, value, and geographical location
- <u>Earth-quake studies:</u> Observed earth quake epicenters should be clustered along continent faults

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What Is Good Clustering?



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



Requirements of Clustering in Data Mining



- Scalability
- · Ability to deal with different types of attributes
- · 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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Data Structures



- (two modes)

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

Dissimilarity matrix

- (one mode)

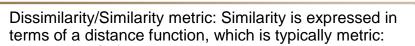
$$\begin{bmatrix} 0 \\ d(2,1) & 0 \\ d(3,1) & d(3,2) & 0 \\ \vdots & \vdots & \vdots \\ d(n,1) & d(n,2) & \dots & \dots & 0 \end{bmatrix}$$

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Measure the Quality of Clustering



d(i, j)

• There is a separate "quality" function that measures the "goodness" of a cluster.

- The definitions of distance functions are usually very different for interval-scaled, boolean, categorical, ordinal and ratio variables.
- Weights should be associated with different variables based on applications and data semantics.
- It is hard to define "similar enough" or "good enough"
 - the answer is typically highly subjective.

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Type of data in clustering analysis



- Interval-scaled variables:
- Binary variables:
- Nominal, ordinal, and ratio variables:
- Variables of mixed types:

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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

$$m_f = \frac{1}{n}(x_{1f} + x_{2f} + \dots + x_{nf}).$$

- 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

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Similarity and Dissimilarity Between Objects



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

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

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

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Similarity and Dissimilarity Between Objects (Cont.)



• If q = 2, d is Euclidean distance:

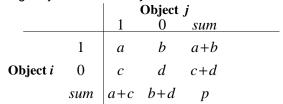
$$d(i,j) = \sqrt{(|x_{i1} - x_{j1}|^2 + |x_{i2} - x_{j2}|^2 + \dots + |x_{ip} - x_{jp}|^2)}$$

- Properties
 - $d(i,j) \ge 0$
 - d(i,i) = 0
 - d(i,j) = d(j,i)
 - $d(i,j) \le d(i,k) + d(k,j)$
- Also, one can use weighted distance, parametric
 Pearson product moment correlation, or other disimilarity measures



Binary Variables





• Simple matching coefficient (invariant, if the binary variable is

<u>symmetric</u>): $d(i, j) = \frac{b+c}{a+b+c+d}$ Jaccard coefficient (noninvariant if the binary variable is

Jaccard coefficient (noninvariant if the binary variable is asymmetric):

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

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Dissimilarity between Binary Variables



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

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

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

$$d (jim , mary) = \frac{1+2}{1+3+0} = 0.75$$



Nominal Variables



- A generalization of the 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 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

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Ordinal Variables



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- · 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 AeBt or AeBt
- Methods:
 - treat them like interval-scaled variables—not a good choice! (why?—the scale can be distorted)
 - apply logarithmic transformation

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

- treat them as continuous ordinal data treat their rank as intervalscaled

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Variables of Mixed **Types**



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- 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 $d\ (i,\ j) = \frac{\sum_{f=1}^{p} \delta_{ij}^{\ (f)} d_{ij}^{\ (f)}}{\sum_{f=1}^{p} \delta_{ij}^{\ (f)}}$

$$d(i, j) = \frac{\sum_{f=1}^{p} \delta_{ij}^{(f)} d_{ij}^{(f)}}{\sum_{f=1}^{p} \delta_{ii}^{(f)}}$$

− f is binary or nominal:

$$d_{ij}^{(f)} = 0$$
 if $x_{if} = x_{jf}$, or $d_{ij}^{(f)} = 1$ o.w.

- f is interval-based: use the normalized distance
- f is ordinal or ratio-scaled
 - · compute ranks r_{if} and

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

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Major Clustering Approaches



- <u>Partitioning algorithms</u>: Construct various partitions and then evaluate them by some criterion
- <u>Hierarchy algorithms</u>: Create a hierarchical decomposition of the set of data (or objects) using some criterion
- <u>Density-based</u>: based on connectivity and density functions
- Grid-based: based on a multiple-level granularity structure
- Model-based: A model is hypothesized for each of the clusters and the idea is to find the best fit of that model to each other



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Partitioning Algorithms: Basic Concept



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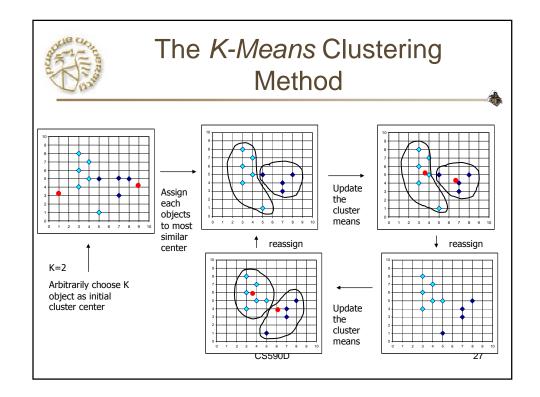
- <u>Partitioning method:</u> Construct a partition of a database
 D of **n** objects into a set of **k** clusters
- Given a k, find a partition of k clusters that optimizes the chosen partitioning criterion
 - Global optimal: exhaustively enumerate all partitions
 - Heuristic methods: k-means and k-medoids algorithms
 - <u>k-means</u> (MacQueen'67): Each cluster is represented by the center of the cluster
 - <u>k-medoids</u> 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 *k*, the *k-means* algorithm is implemented in four steps:
 - Partition objects into k nonempty subsets
 - Compute seed points as the centroids of the clusters of the current partition (the centroid is the center, i.e., *mean point*, of the cluster)
 - Assign each object to the cluster with the nearest seed point
 - Go back to Step 2, stop when no more new assignment





Comments on the *K-Means*Method



- <u>Strength:</u> Relatively efficient: O(tkn), where n is # objects, k is # clusters, and t is # iterations. Normally, k, t << n.
 - Comparing: PAM: O(k(n-k)²), CLARA: O(ks² + k(n-k))
- <u>Comment:</u> 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

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Variations of the *K-Means*Method



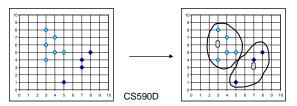
- A few variants of the k-means which differ in
 - Selection of the initial k means
 - Dissimilarity calculations
 - Strategies to calculate cluster means
- Handling categorical data: k-modes (Huang'98)
 - Replacing means of clusters with modes
 - Using new dissimilarity measures to deal with categorical objects
 - Using a frequency-based method to update modes of clusters
 - A mixture of categorical and numerical data: k-prototype method



What is the problem of k-Means Method?



- 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.



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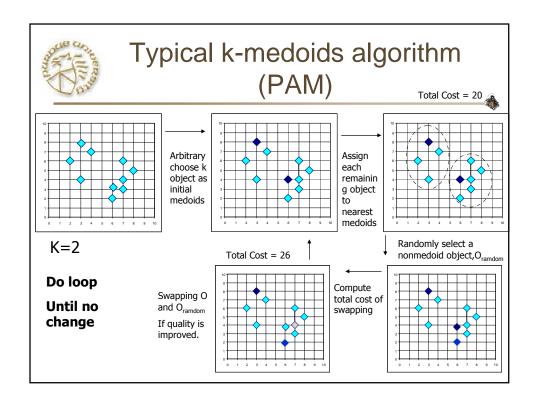


The K-Medoids Clustering Method



- Find representative objects, called medoids, in clusters
- PAM (Partitioning Around Medoids, 1987)
 - starts from an initial set of medoids and iteratively replaces one of the medoids by one of the non-medoids if it improves the total distance of the resulting clustering
 - PAM works effectively for small data sets, but does not scale well for large data sets
- CLARA (Kaufmann & Rousseeuw, 1990)
- CLARANS (Ng & Han, 1994): Randomized sampling
- Focusing + spatial data structure (Ester et al., 1995)

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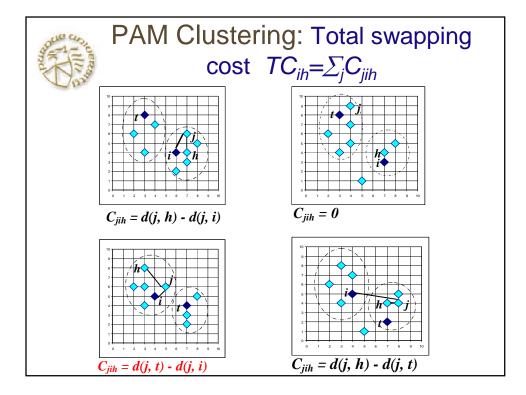




PAM (Partitioning Around Medoids) (1987)



- PAM (Kaufman and Rousseeuw, 1987), built in Splus
- Use real object to represent the cluster
 - Select **k** representative objects arbitrarily
 - For each pair of non-selected object h and selected object i,
 calculate the total swapping cost TC_{ih}
 - For each pair of *i* and *h*,
 - If $TC_{ih} < 0$, **i** is replaced by **h**
 - Then assign each non-selected object to the most similar representative object
 - repeat steps 2-3 until there is no change





What is the problem with PAM?



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- Pam is more robust than k-means in the presence of noise and outliers because a medoid is less influenced by outliers or other extreme values than a mean
- Pam works efficiently for small data sets but does not scale well for large data sets.
 - O(k(n-k)2) for each iteration

where n is # of data,k is # of clusters

→ Sampling based method, CLARA(Clustering LARge Applications)



CLARA (Clustering Large Applications) (1990)



- CLARA (Kaufmann and Rousseeuw in 1990)
 - Built in statistical analysis packages, such as S+
- It draws *multiple samples* of the data set, applies *PAM* on each sample, and gives the best clustering as the output
- Strength: deals with larger data sets than PAM
- Weakness:
 - Efficiency depends on the sample size
 - A good clustering based on samples will not necessarily represent a good clustering of the whole data set if the sample is biased

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CLARANS ("Randomized" CLARA) (1994)



- CLARANS (A Clustering Algorithm based on Randomized Search) (Ng and Han'94)
- CLARANS draws sample of neighbors dynamically
- The clustering process can be presented as searching a graph where every node is a potential solution, that is, a set of *k* medoids
- If the local optimum is found, CLARANS starts with new randomly selected node in search for a new local optimum
- It is more efficient and scalable than both PAM and CLARA
- Focusing techniques and spatial access structures may further improve its performance (Ester et al.'95)



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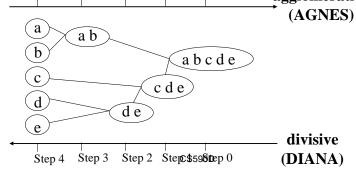
Hierarchical Clustering



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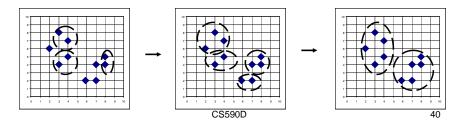
Use distance matrix as clustering criteria. This method does not require the number of clusters
 k as an input, but needs a termination condition
 Step 0 Step 1 Step 2 Step 3 Step 4 agglomerative





AGNES (Agglomerative Nesting)

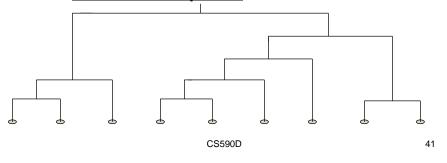
- Introduced in Kaufmann and Rousseeuw (1990)
- · Implemented in statistical analysis packages, e.g., Splus
- Use the Single-Link method and the dissimilarity matrix.
- · Merge nodes that have the least dissimilarity
- · Go on in a non-descending fashion
- · Eventually all nodes belong to the same cluster





A Dendrogram Shows How the Clusters are Merged Hierarchically

- Decompose data objects into a several levels of nested partitioning (<u>tree</u> of clusters), called a <u>dendrogram</u>.
- A <u>clustering</u> of the data objects is obtained by <u>cutting</u> the dendrogram at the desired level, then each connected component forms a cluster.

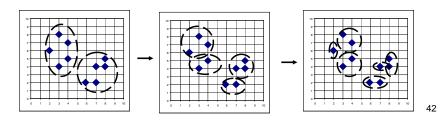




DIANA (Divisive Analysis)



- Implemented in statistical analysis packages, e.g., Splus
- Inverse order of AGNES
- · Eventually each node forms a cluster on its own





More on Hierarchical Clustering Methods



- Major weakness of agglomerative clustering methods
 - <u>do not scale</u> 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
 - <u>CURE (1998)</u>: selects well-scattered points from the cluster and then shrinks them towards the center of the cluster by a specified fraction
 - CHAMELEON (1999): hierarchical clustering using dynamic modeling



BIRCH (1996)



- Birch: Balanced Iterative Reducing and Clustering using Hierarchies, by Zhang, Ramakrishnan, Livny (SIGMOD'96)
- 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 an arbitrary clustering algorithm to cluster the leaf nodes of the CF-tree
- Scales linearly: finds a good clustering with a single scan and improves the quality with a few additional scans
- Weakness: handles only numeric data, and sensitive to the order of the data record.

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Clustering Feature Vector

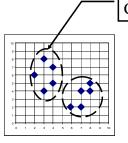


Clustering Feature: $CF = (N, \overrightarrow{LS}, SS)$

N: Number of data points

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

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



CF = (5, (16,30), (54,190))

(3,4)

(2,6) (4,5)

(4,7)

(3,8)

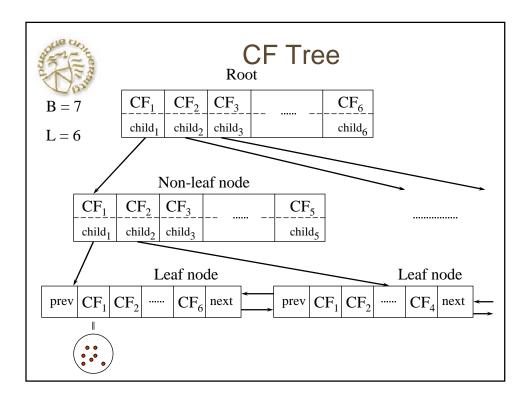
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CF-Tree in BIRCH



- summary of the statistics for a given subcluster: the 0-th, 1st and 2nd moments of the subcluster from the statistical point of view.
- 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





CURE (Clustering Using REpresentatives)









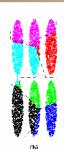
- CURE: proposed by Guha, Rastogi & Shim, 1998
 - Stops the creation of a cluster hierarchy if a level consists of k clusters
 - Uses multiple representative points to evaluate the distance between clusters, adjusts well to arbitrary shaped clusters and avoids single-link effect

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Drawbacks of Distance-Based Method







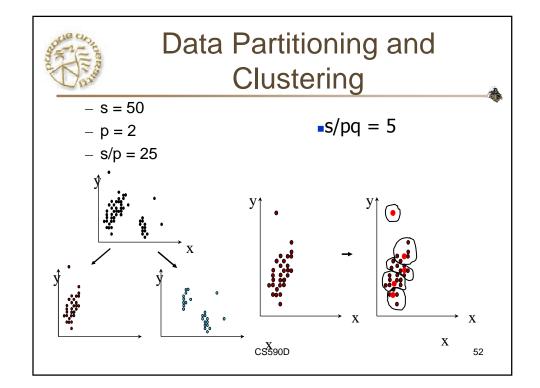
- Drawbacks of square-error based clustering method
 - Consider only one point as representative of a cluster
 - Good only for convex shaped, similar size and density, and if k can be reasonably estimated

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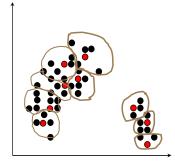
Cure: The Algorithm

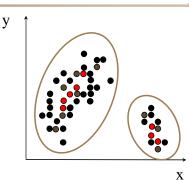
- Draw random sample s.
- Partition sample to p partitions with size s/p
- Partially cluster partitions into s/pq clusters
- Eliminate outliers
 - By random sampling
 - If a cluster grows too slow, eliminate it.
- Cluster partial clusters.
- Label data in disk





Cure: Shrinking Representative Points





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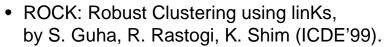
- Shrink the multiple representative points towards the gravity center by a fraction of α .
- Multiple representatives capture the shape of the cluster

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Clustering Categorical Data: ROCK



- Use links to measure similarity/proximity
- Not distance based
- Computational complexity: $O(n^2 + nm_m m_a + n^2 \log n)$
- · Basic ideas:
 - Similarity function and neighbors: $Sim(T_1,T_2) = \frac{|T_1 \cap T_2|}{|T_1 \cup T_2|}$ Let $T_1 = \{1,2,3\}, \ T_2 = \{3,4,5\}$

$$Sim(T1, T2) = \frac{|\{3\}|}{|\{1,2,3,4,5\}|} = \frac{1}{5} = 0.2$$



Rock: Algorithm



 Links: The number of common neighbours for the two points.

$$\{1,2,3\}, \{1,2,4\}, \{1,2,5\}, \{1,3,4\}, \{1,3,5\}$$

 $\{1,4,5\}, \{2,3,4\}, \{2,3,5\}, \{2,4,5\}, \{3,4,5\}$
 $\{1,2,3\} \xrightarrow{3} \{1,2,4\}$

- Algorithm
 - Draw random sample
 - Cluster with links
 - Label data in disk

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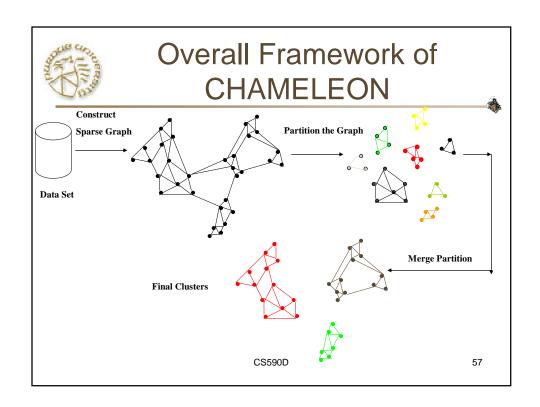


CHAMELEON (Hierarchical clustering using dynamic modeling)



- CHAMELEON: by G. Karypis, E.H. Han, and V. Kumar'99
- · 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
 - Cure ignores information about interconnectivity of the objects,
 Rock ignores information about the closeness of two clusters
- A two-phase algorithm
 - 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

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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
 - One scan
 - Need density parameters as termination condition
- Several interesting studies:
 - DBSCAN: Ester, et al. (KDD'96)
 - OPTICS: Ankerst, et al (SIGMOD'99).
 - DENCLUE: Hinneburg & D. Keim (KDD'98)
 - CLIQUE: Agrawal, et al. (SIGMOD'98)

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Density Concepts

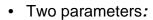


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- Core object (CO)—object with at least 'M' objects within a radius 'E-neighborhood'
- Directly density reachable (DDR)—x is CO, y is in x's 'Eneighborhood'
- Density reachable—there exists a chain of DDR objects from x to y
- Density based cluster–density connected objects maximum w.r.t. reachability

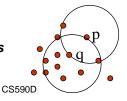


Density-Based Clustering: Background



- Eps: Maximum radius of the neighbourhood
- MinPts: Minimum number of points in an Eps-neighbourhood of that point
- $N_{Eps}(p)$: {q belongs to D | dist(p,q) <= Eps}
- Directly density-reachable: A point p is directly density-reachable from a point q wrt. Eps, MinPts if
 - 1) p belongs to N_{Eps}(q)
 - 2) core point condition:

$$|N_{Eps}(q)| >= MinPts$$



MinPts = 5

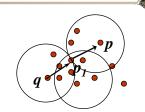
Eps = 1 cm

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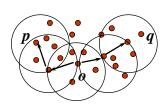


Density-Based Clustering: Background (II)

- Density-reachable:
 - A point p is density-reachable from a point q wrt. *Eps*, *MinPts* if there is a chain of points $p_1, ..., p_n, p_1 = q, p_n = p$ such that p_{i+1} is directly density-reachable from p_i



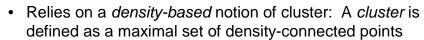
- Density-connected
 - A point p is density-connected to a point q wrt. Eps, MinPts if there is a point o such that both, p and q are density-reachable from o wrt. Eps and MinPts.



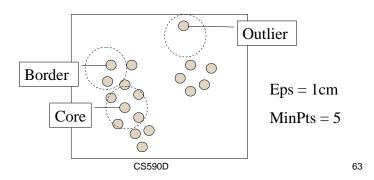
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DBSCAN: Density Based Spatial Clustering of Applications with Noise



Discovers clusters of arbitrary shape in spatial databases with noise





DBSCAN: The Algorithm



- Arbitrary select a point p
- Retrieve all points density-reachable from p wrt Eps and MinPts.
- If p is a core point, a cluster is formed.
- If p is a border point, no points are density-reachable from p and DBSCAN visits the next point of the database.
- Continue the process until all of the points have been processed.



OPTICS: A Cluster-Ordering Method (1999)

- OPTICS: Ordering Points To Identify the **Clustering Structure**
 - Ankerst, Breunig, Kriegel, and Sander (SIGMOD'99)
 - Produces a special order of the database wrt its density-based clustering structure
 - This cluster-ordering contains info equiv to the density-based clusterings corresponding to a broad range of parameter settings
 - Good for both automatic and interactive cluster analysis, including finding intrinsic clustering structure
 - Can be represented graphically or using visualization techniques

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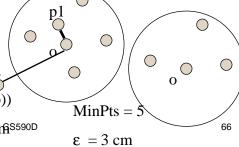
OPTICS: Some Extension from **DBSCAN**

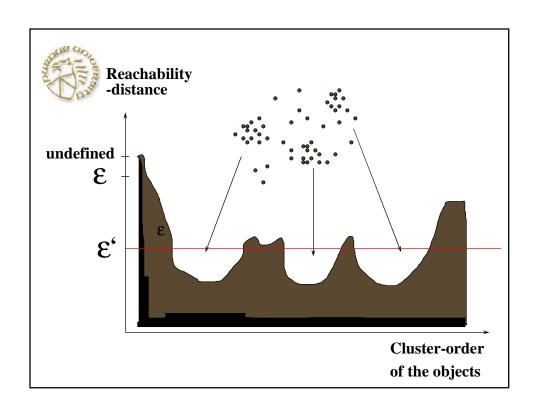
- Index-based:

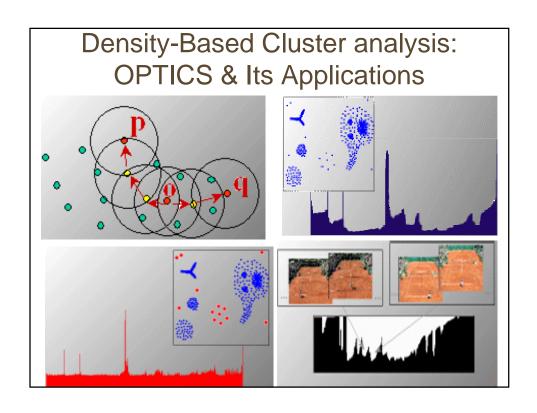
 - N = 20
 - p = 75%
 - M = N(1-p) = 5
 - Complexity: $O(kN^2)$
- Core Distance
- Reachability Distance

Max (core-distance (o), d (o, p))

r(p1, o) = 2.8cm. $r(p2, o) = 4crf^{S590D}$









DENCLUE: Using density functions



- DENsity-based CLUstEring by Hinneburg & Keim (KDD'98)
- · Major features
 - Solid mathematical foundation
 - Good for data sets with large amounts of noise
 - Allows a compact mathematical description of arbitrarily shaped clusters in high-dimensional data sets
 - Significant faster than existing algorithm (faster than DBSCAN by a factor of up to 45)
 - But needs a large number of parameters

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Denclue: Technical Essence



- Uses grid cells but only keeps information about grid cells that do actually contain data points and manages these cells in a tree-based access structure.
- Influence function: describes the impact of a data point within its neighborhood.
- Overall density of the data space can be calculated as the sum of the influence function of all data points.
- Clusters can be determined mathematically by identifying density attractors.
- Density attractors are local maximal of the overall density function.



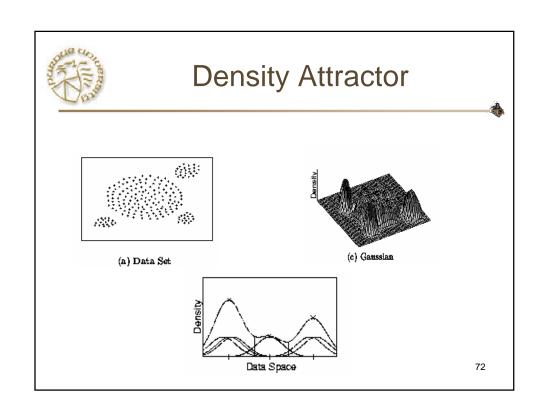
Gradient: The steepness of a slope

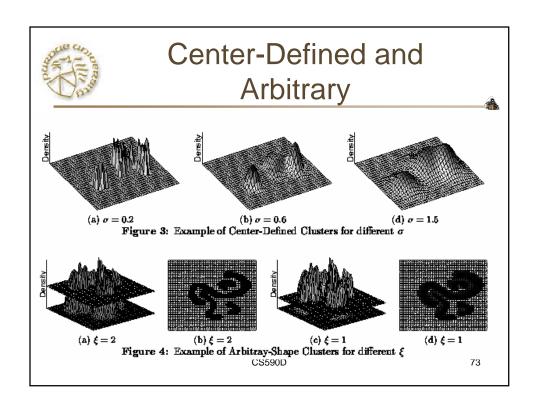
• Example

$$f_{Gaussian}(x,y) = e^{-\frac{d(x,y)^{2}}{2\sigma^{2}}}$$

$$f_{Gaussian}^{D}(x) = \sum_{i=1}^{N} e^{-\frac{d(x,x_{i})^{2}}{2\sigma^{2}}}$$

$$\nabla f_{Gaussian}^{D}(x,x_{i}) = \sum_{i=1}^{N} (x_{i}-x) \cdot e^{-\frac{d(x,x_{i})^{2}}{2\sigma^{2}}}$$







Cluster Analysis



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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)

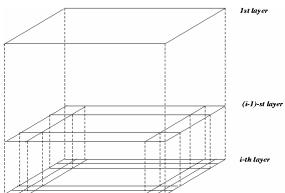
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STING: A Statistical Information Grid Approach



- Wang, Yang and Muntz (VLDB'97)
- The spatial area area is divided into rectangular cells
- There are several levels of cells corresponding to different levels of resolution



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STING: A Statistical Information Grid Approach (2)



- Each cell at a high level is partitioned into a number of smaller cells in the next lower level
- Statistical info of each cell is calculated and stored beforehand and is used to answer queries
- Parameters of higher level cells can be easily calculated from parameters of lower level cell
 - count, mean, s, min, max
 - type of distribution—normal, *uniform*, etc.
- 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

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STING: A Statistical Information Grid Approach (3)

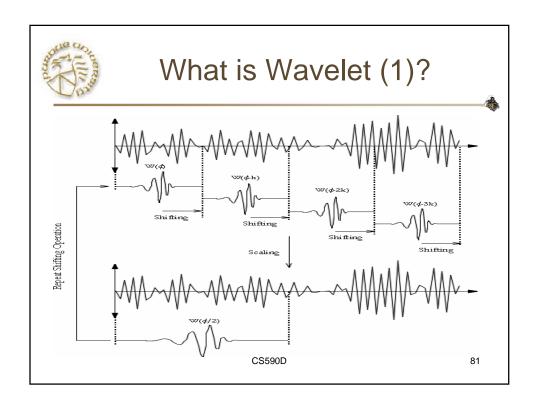


- Remove the irrelevant cells from further consideration
- When finish examining the current layer, proceed to the next lower level
- Repeat this process until the bottom layer is reached
- Advantages:
 - Query-independent, easy to parallelize, incremental update
 - O(K), where K is the number of grid cells at the lowest level
- Disadvantages:
 - All the cluster boundaries are either horizontal or vertical, and no diagonal boundary is detected



WaveCluster (1998)

- Sheikholeslami, Chatterjee, and Zhang (VLDB'98)
- A multi-resolution clustering approach which applies wavelet transform to the feature space
 - A wavelet transform is a signal processing technique that decomposes a signal into different frequency sub-band.
- · Both grid-based and density-based
- Input parameters:
 - # of grid cells for each dimension
 - the wavelet, and the # of applications of wavelet transform.





WaveCluster (1998)

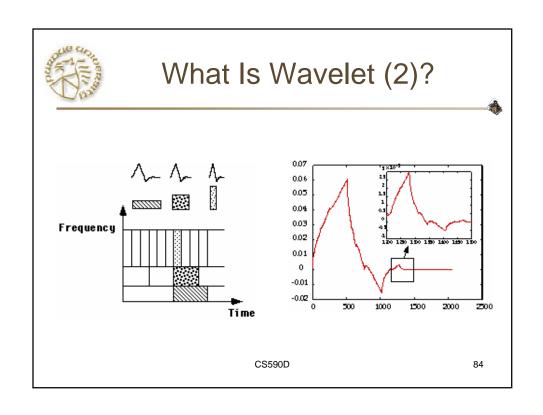
- How to apply wavelet transform to find clusters
 - Summaries the data by imposing a multidimensional grid structure onto data space
 - These multidimensional spatial data objects are represented in a n-dimensional feature space
 - Apply wavelet transform on feature space to find the dense regions in the feature space
 - Apply wavelet transform multiple times which result in clusters at different scales from fine to coarse

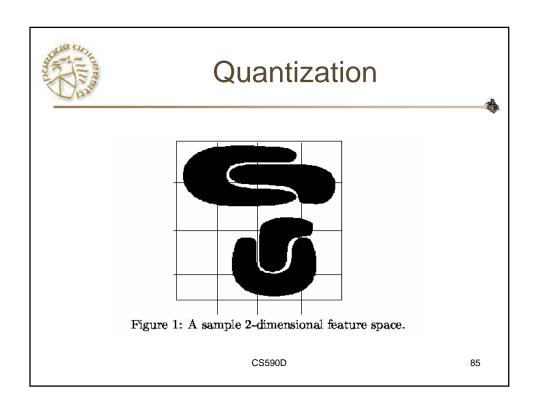
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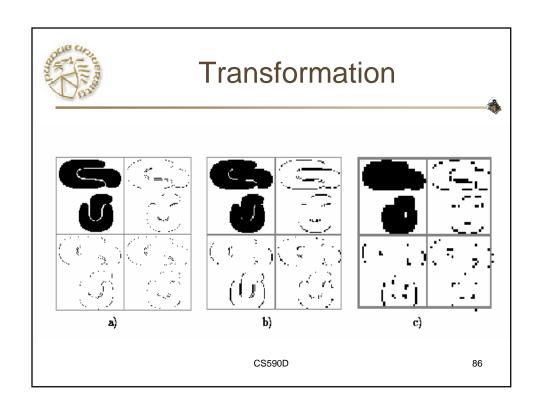


Wavelet Transform

- Decomposes a signal into different frequency subbands. (can be applied to ndimensional signals)
- Data are transformed to preserve relative distance between objects at different levels of resolution.
- Allows natural clusters to become more distinguishable









WaveCluster (1998)



- · Why is wavelet transformation useful for clustering
 - Unsupervised clustering
 It uses hat-shape filters to emphasize region where points cluster, but simultaneously to suppress weaker information in their boundary
 - Effective removal of outliers
 - Multi-resolution
 - Cost efficiency
- · Major features:
 - Complexity O(N)
 - Detect arbitrary shaped clusters at different scales
 - Not sensitive to noise, not sensitive to input order
 - Only applicable to low dimensional data



CLIQUE (Clustering In QUEst)



- Agrawal, Gehrke, Gunopulos, Raghavan (SIGMOD'98).
- Automatically identifying subspaces of a high dimensional data space that allow better clustering than original space
- CLIQUE can be considered as both density-based and grid-based
 - It partitions each dimension into the same number of equal length interval
 - It partitions an m-dimensional data space into non-overlapping rectangular units
 - A unit is dense if the fraction of total data points contained in the unit exceeds the input model parameter
 - A cluster is a maximal set of connected dense units within a subspace

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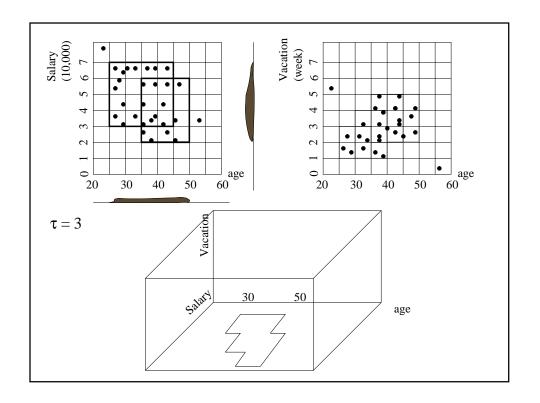


CLIQUE: The Major Steps



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- Partition the data space and find the number of points that lie inside each cell of the partition.
- Identify the subspaces that contain clusters using the Apriori principle
- Identify clusters:
 - Determine dense units in all subspaces of interests
 - Determine connected dense units in all subspaces of interests.
- Generate minimal description for the clusters
 - Determine maximal regions that cover a cluster of connected dense units for each cluster
 - Determination of minimal cover for each cluster





Strength and Weakness of *CLIQUE*



- It <u>automatically</u> finds <u>subspaces</u> of the <u>highest</u> dimensionality such that high density clusters exist in those subspaces
- It is insensitive to the order of records in input and does not presume some canonical data distribution
- It scales *linearly* with the size of input and has good scalability as the number of dimensions in the data increases

Weakness

 The accuracy of the clustering result may be degraded at the expense of simplicity of the method



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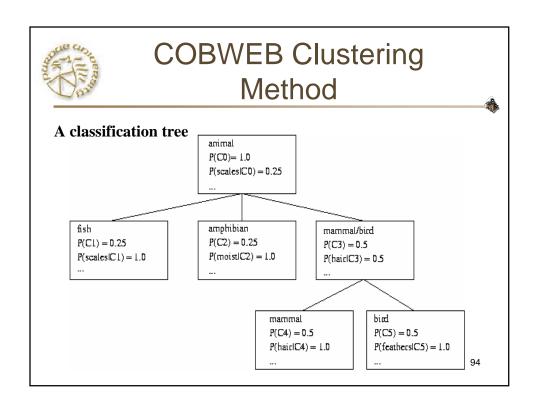


Model-Based Clustering Methods



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- Attempt to optimize the fit between the data and some mathematical model
- Statistical and AI approach
 - Conceptual clustering
 - · A form of clustering in machine learning
 - Produces a classification scheme for a set of unlabeled objects
 - Finds characteristic description for each concept (class)
 - COBWEB (Fisher'87)
 - A popular a simple method of incremental conceptual learning
 - Creates a hierarchical clustering in the form of a classification tree
 - Each node refers to a concept and contains a probabilistic description of that concept





More on Statistical-Based Clustering



Limitations of COBWEB

- The assumption that the attributes are independent of each other is often too strong because correlation may exist
- Not suitable for clustering large database data skewed tree and expensive probability distributions

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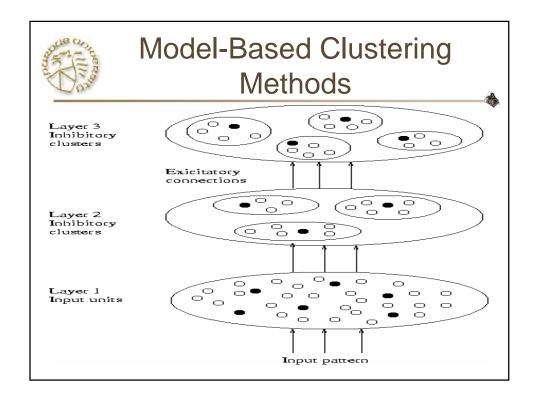
- an extension of COBWEB for incremental clustering of continuous data
- suffers similar problems as COBWEB
- AutoClass (Cheeseman and Stutz, 1996)
 - Uses Bayesian statistical analysis to estimate the number of clusters
 - Popular in industry



Other Model-Based Clustering Methods



- Neural network approaches
 - Represent each cluster as an exemplar, acting as a "prototype" of the cluster
 - New objects are distributed to the cluster whose exemplar is the most similar according to some dostance measure
- Competitive learning
 - Involves a hierarchical architecture of several units (neurons)
 - Neurons compete in a "winner-takes-all" fashion for the object currently being presented





Self-organizing feature maps (SOMs)



- Clustering is also performed by having several units competing for the current object
- The unit whose weight vector is closest to the current object wins
- The winner and its neighbors learn by having their weights adjusted
- SOMs are believed to resemble processing that can occur in the brain
- Useful for visualizing high-dimensional data in 2or 3-D space

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Cluster Analysis



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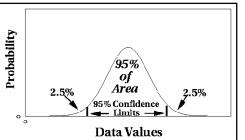
What Is Outlier Discovery?



- The set of objects are considerably dissimilar from the remainder of the data
- Example: Sports: Michael Jordon, Wayne Gretzky, ...
- Problem
 - Find top n outlier points
- Applications:
 - Credit card fraud detection
 - Telecom fraud detection
 - Customer segmentation
 - Medical analysis

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Outlier Discovery Education Statistical Approaches



- Use discordancy tests depending on
 - data distribution
 - distribution parameter (e.g., mean, variance)
 - number of expected outliers
- Drawbacks
 - most tests are for single attribute
 - In many cases, data distribution may not be known



Outlier Discovery: Distance-Based Approach



- Introduced to counter the main limitations imposed by statistical methods
 - We need multi-dimensional analysis without knowing data distribution.
- 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
 - Nested-loop algorithm
 - Cell-based algorithm

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Outlier Discovery: 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
- OLAP data cube technique
 - uses data cubes to identify regions of anomalies in large multidimensional data



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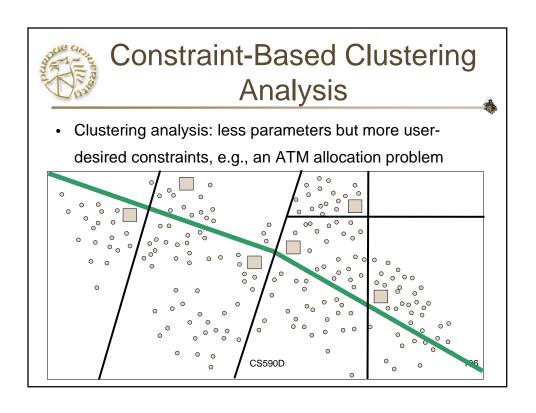
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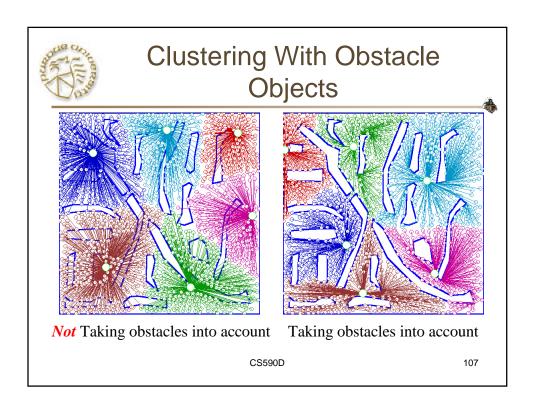


Problems and Challenges



- Considerable progress has been made in scalable clustering methods
 - Partitioning: k-means, k-medoids, CLARANS
 - Hierarchical: BIRCH, CURE
 - Density-based: DBSCAN, CLIQUE, OPTICS
 - Grid-based: STING, WaveCluster
 - Model-based: Autoclass, Denclue, Cobweb
- Current clustering techniques do not <u>address</u> all the requirements adequately
- Constraint-based clustering analysis: Constraints exist in data space (bridges and highways) or in user queries







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, density-based methods, grid-based methods, and model-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
- There are still lots of research issues on cluster analysis, such as constraint-based clustering

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