

Data Mining Steps		
1. Data Cleaning	Removal of noise and inconsistent records	
2. Data Integration	Combing multiple sources	
3. Data Selection	Only data relevant for the task are retrieved from the database	
4. Data Transformation	Converting data into a form more appropriate for mining	
5. Data Mining	Application of intelligent methods to extract data patterns	
6. Model Evaluation	Identification of truly interesting patterns representing knowledge	
7. Knowledge Presentation	Visualization or other knowledge presentation techniques	
Data mining could also be called Knowledge Discovery in Databases (see kdnuggets.com)		

Types of Attributes		
Nomial	e.g., ID numbers, eye color, zip codes	
Ordinal	e.g., rankings, grades, height	
Interval	e.g., calendar dates, temperatures	
Ratio	e.g., length, time, counts	

#### Distance Measures

# **Euclidean Distance:**

$$dist = \sqrt{\sum_{k=1}^{n} (p_k - q_k)^2}$$

# Minkowski Distance:

$$dist = \left(\sum_{k=1}^{n} |p_k - q_k|^r\right)^{\frac{1}{r}}$$

r=1, City Block r=2, Euclidean r->inf., Chebyshev

Manhattan = City Block

Jaccard coefficient, Hamming, Cosine are a similarity / dissimilarity measures



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# Measures of Node Impurity

GAIN = measure before split - measure after split

$$GINI(t) = 1 - \sum_{j} [p(j | t)]^{2}$$

 $p(j \mid t)$  is the relative frequency of class j at node t

$$GINI_{split} = \sum_{i=1}^{k} \frac{n_i}{n} GINI(i)$$

 $\begin{array}{ll} \mbox{where,} & \mbox{$n_i$ = number of records at child i,} \\ \mbox{$n = number of records at node p.} \\ \mbox{Pick the smallest} \end{array}$ 

$$Entropy(t) = -\sum_{j} p(j \mid t) \log p(j \mid t)$$

Information Gain.

$$GAIN_{split} = Entropy(p) - \left(\sum_{i=1}^{k} \frac{n_i}{n} Entropy(i)\right)$$

Parent Node, p is split into k partitions;  $n_i$  is number of records in partition i

$$GainRATIO_{quit} = \frac{GAIN_{squit}}{SplitINFO}$$

$$SplitINFO = -\sum_{i=1}^{k} \frac{n_i}{n} \log \frac{n_i}{n}$$

Parent Node, p is split into k partitions; n<sub>i</sub> is number of records in partition i

$$Error(t) = 1 - \max_{i} P(i \mid t)$$

# Model Evaluation

	PREDICTED CLASS		
ACTUAL		Class=Yes	Class=No
	Class=Yes	a (TP)	b (FN)
CLASS	Class=No	c (FP)	d (TN)

Accuracy =  $\frac{TP+TN}{TP+FN+TN+FP}$ 

Precision =  $\frac{TP}{TP+FP}$ 

 $Recall = \frac{TP}{TP + FN}$ 

F-measure =  $\frac{2 \cdot TP}{2 \cdot TP + FN + FP}$ 

 $\mathrm{Cost} = TP \times Cost_{TP} + FN \times Cost_{FN}$ 

 $+\,TN\times Cost_{TN}+FP\times Cost_{FP}$ 

Sensitivity = Recall

Specificity =  $1 - \frac{FP}{FP + TN} = \frac{TN}{TN + FP}$ 

False Positive Rate = 1 - Specificity

Kappa = (observed agreement - chance agreement) / (1- chance agreement)

Kappa = (Dreal – Drandom) / (Dperfect – Drandom), where D indicates the sum of values in diagonal of the confusion matrix

### K-Nearest Neighbor

- \* Compute distance between two points
- \* Determine the class from nearest neighbor list
  - \* Take the majority vote of class labels  $\label{eq:majority} \text{among the $k$-nearest neighbors}$
  - \* Weigh the vote according to distance

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# K-Nearest Neighbor (cont)

\* weight factor, w = 1 / d^2

Classify records by using a collection of

"if...then..." rules

Rule: (Condition) --> y

where:

\* Condition is a conjunction of attributes

\* y is the class label

LHS: rule antecedent or condition

RHS: rule consequent

Examples of classification rules:

(Blood Type=Warm) ^ (Lay Eggs=Yes) --> Birds

(Taxable Income < 50K) ^ (Refund=Yes) --> Evade=No

Sequential covering is a rule-based classifier.

### Rule Evaluation

Accuracy = 
$$\frac{n_c}{n_c}$$

Laplace = 
$$\frac{n_c + 1}{n_c + b}$$

$$\text{M-estimate} = \frac{n_c + kp}{n + k}$$

n: Number of instances covered by rule  $n_c$ : Number of instances of class c covered by rule k: Number of classes p: Prior probability (for the positive class)

# Bayesian Classification

Conditional Probability:  $P(C|A) = \frac{P(A,C)}{P(C|A)}$ 

$$P(A \mid C) = \frac{P(A,C)}{P(A \mid C)}$$

Baye's theorem:

$$P(C \mid A) = \frac{P(A \mid C)P(C)}{P(A)}$$

 $posterior = \frac{likelihood \times prior}{normalizing constant}$ 

Naïve Bayes Classifer:

Original: 
$$P(A_r | C) = \frac{N_{tc}}{N_c}$$

Laplace: 
$$P(A_i \mid C) = \frac{N_{ic} + 1}{N_c + c}$$

m - estimate : 
$$P(A_{\varepsilon} \mid C) = \frac{N_{\varepsilon} + mp}{N_{\varepsilon} + m}$$

c: number of classes, p: prior probability, m: parameter

P(B|A), read as the probability of B given A.

$$P(B|A) = \frac{P(A \text{ and } B)}{P(A)} P(A \text{ and } B) = P(A) \cdot P(B|A)$$

p(a,b) is the probability that both a and b happen.

p(a|b) is the probability that a happens, knowing that b has already happened.

Association Min-Apriori, LIFT, Simpson's Paradox, Anti-m-

Analysis onotone property

Ensemble Staking, Random Forest

Methods

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Terms (cont)	
Decision Trees	C4.5, Pessimistic estimate, Occam's Razor, Hunt's Algorithm
Model Evaluation	Cross-validation, Bootstrap, Leave-one out (C-V), Misclassification error rate, Repeated holdout, Stratification
Bayes	Probabilistic classifier
Data Visual- ization	Chernoff faces, Data cube, Percentile plots, Parallel coordinates
Nonlinear Dimensionality Reduction	Principal components, ISOMAP, Multidimensional scaling

# Ensemble Techniques

### AdaBoost Algorithm:

$$\alpha_t = \frac{1}{2} \ln \left( \frac{1 - \epsilon_t}{\epsilon_t} \right)$$

error ( $\epsilon_t$ ) = # of misclassified divided by total

$$w_i = w_1 = w_2 = \dots = w_{10} = \frac{1}{10} = 0.1$$

Re-weighting:

 $misclassified = w_i \times e^{+\alpha_t}$ 

 $correct\ classified = w_i \times e^{-\alpha_t}$ 

**Manipulate training data:** bagging and boosting ensemble of "experts", each specializing on different portions of the instance space

Manipulate output values: error-correcting output coding (ensemble of "experts", each predicting 1 bit of the {multibit} full class label)

Methods: BAGGing, Boosting, AdaBoost

# Rules Analysis

$$support = \frac{P(A,B)}{Total}$$

$$confidence = \frac{P(A,B)}{P(A)}$$

$$Lift = \frac{Confidence}{P(B)}$$
Example:
$$Rule \{b\} \rightarrow \{c\} \qquad \qquad \boxed{\begin{array}{c|c} & \boxed{b} & \boxed{b} \\ \hline b & \boxed{3} & \boxed{4} \\ \hline b & \boxed{2} & \boxed{1} \\ \hline \end{array}}$$

$$support = 3/10 = 0.3$$

$$confidence = 3/7 = 0.4286$$

$$lift = \frac{3/7}{5/10}$$

### Apriori Algorithm

Let k=1

Generate frequent itemsets of length 1
Repeat until no new frequent itemsets are identified

Gen erate length (k+1) candidate itemsets

from

length k frequent itemsets

Prune candidate itemsets containing

subsets

of length k that are infrequent

Count the support of each candidate by

sca nning the DB

Eli minate candidates that are infreq -

uent,

leaving only those that are frequent

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### K-means Clustering

Select K points as the initial centroids repeat

Rec ompute the centroid of each cluster until the centroids don't change

**Closeness** is measured by distance (e.g., Euclidean), similarity (e.g., Cosine), correlation.

Centroid is typically the mean of the points in the cluster

#### **Hierarchical Clustering**

### Single-Link or MIN

Similarity of two clusters is based on the two most similar (closest / minimum) points in the different clusters

Determined by one pair of points, i.e., by one link in the proximity graph.

#### Complete or MAX

Similarity of two clusters is based on the two least similar (most distant, maximum) points in the different clusters

Determined by all pairs of points in the two clusters

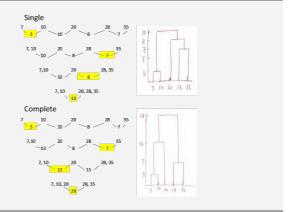
#### **Group Average**

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

**Agglomerative** clustering starts with points as individual clusters and merges closest clusters until only one cluster left.

**Divisive** clustering starts with one, all-inclusive cluster and splits a cluster until each cluster only has one point.

# Dendrogram Example



Dataset: {7, 10, 20, 28, 35}

# **Density-Based Clustering**

current\_cluster\_label <-- 1</pre>

for all core points do

if the core point has no cluster label then

cur ren t\_c lus ter \_label <--

curren t\_c lus ter \_label +1

Label the current core point with

the cluster label

end if

 $\begin{tabular}{ll} \textbf{for all points in the Eps-ne igh bor hood,} \\ \textbf{except i-th the point itself } \textbf{do} \end{tabular}$ 

if the point does not have a cluster

label then

Label the point with

cluster label

end if

end for



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### **Density-Based Clustering (cont)**

#### > end for

DBSCAN is a popular algorithm

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

A point is a core point if it has more than a specified number of points (MinPts) within Eps

These are points that are at the interior of a cluster

A border point has fewer than MinPts within Eps, but is in the neighborhood of a core point

A noise point is any point that is not a core point or a border point

# Other Clustering Methods

**Fuzzy** is a partitional clustering method. **Fuzzy clustering** (also referred to as **soft clustering**) is a form of clustering in that each data point can belong to more than one cluster.

**Graph-based** methods: Jarvis-Patrick, Shared-Near Neighbor (SNN, Density), Chameleon

Model-based methods: Expectation-Maximization

### Regression Analysis

- \* Linear Regression
- | Least squares
- \* Subset selection
- \* Stepwise selection
- \* Regularized regression

| Ridge

| Lasso

# Regression Analysis (cont)

| Elastic Net

### **Anomaly Detection**

Anomaly is a pattern in the data that does not conform to the expected behavior (e.g., outliers, exceptions, peculiarities, surprise)

#### Types of Anomaly

Point: An individual data instance is anomalous w.r.t. the data Contextual: An individual data instance is anomalous within a context

Collective: A collection of related data instances is anomalous

#### **Approaches**

- \* Graphical (e.g., boxplots, scatter plots)
- \* Statistical (e.g., normal distribution, likelihood)
- | Parametric Techniques
- | Non-parametric Techniques
- \* Distance (e.g., nearest-neighbor, density, clustering)

Local outlier factor (LOF) is a density-based distance approach

Mahalanobis Distance is a clustering-based distance approach



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