CS6220: DATA MINING TECHNIQUES

Matrix Data: Classification: Part 1

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September 21, 2015

Matrix Data: Classification: Part 1

- Classification: Basic Concepts
- Decision Tree Induction
- Model Evaluation and Selection
- Summary

Supervised vs. Unsupervised Learning

- Supervised learning (classification)
 - Supervision: The training data (observations, measurements, etc.) are accompanied by labels indicating the class of the observations
 - New data is classified based on the training set
- Unsupervised learning (clustering)
 - The class labels of training data is unknown
 - Given a set of measurements, observations, etc. with the aim of establishing the existence of classes or clusters in the data

Prediction Problems: Classification vs. Numeric Prediction

- Classification
 - predicts categorical class labels
 - classifies data (constructs a model) based on the training set and the values (class labels) in a classifying attribute and uses it in classifying new data
- Numeric Prediction
 - models continuous-valued functions, i.e., predicts unknown or missing values
- Typical applications
 - Credit/loan approval:
 - Medical diagnosis: if a tumor is cancerous or benign
 - Fraud detection: if a transaction is fraudulent
 - Web page categorization: which category it is

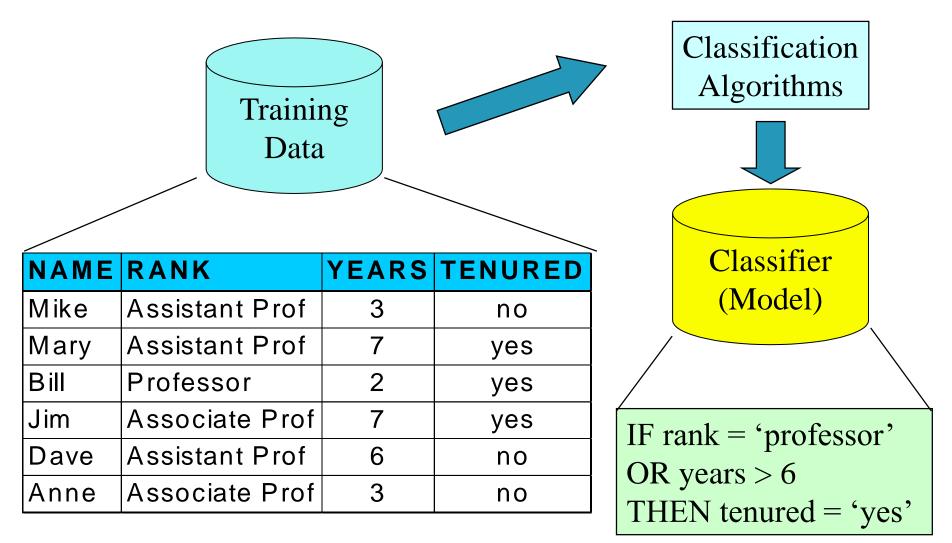
Classification—A Two-Step Process (1)

- Model construction: describing a set of predetermined classes
 - Each tuple/sample is assumed to belong to a predefined class, as determined by the class label attribute
 - For data point $i : \langle x_i, y_i \rangle$
 - Features: x_i; class label: y_i
 - The model is represented as classification rules, decision trees, or mathematical formulae
 - Also called classifier
 - The set of tuples used for model construction is training set

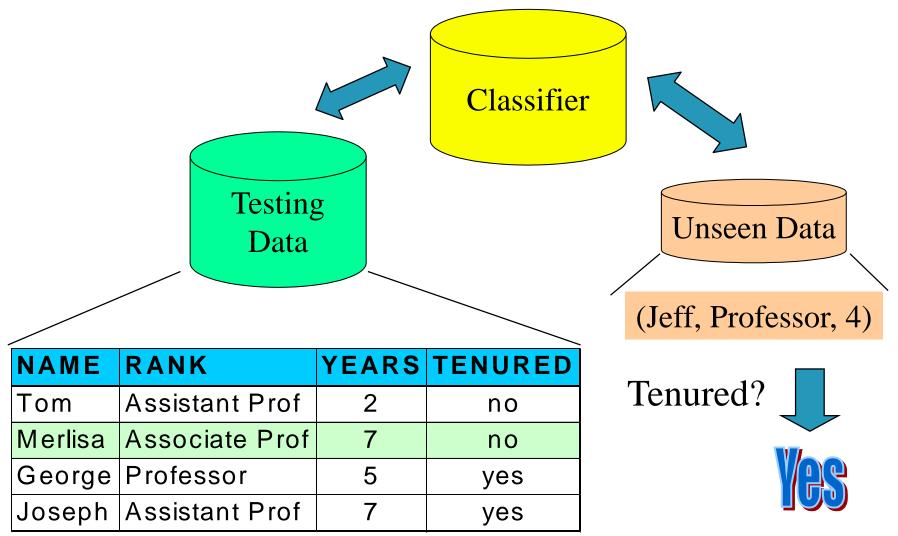
Classification—A Two-Step Process (2)

- Model usage: for classifying future or unknown objects
 - Estimate accuracy of the model
 - The known label of test sample is compared with the classified result from the model
 - Test set is independent of training set (otherwise overfitting)
 - Accuracy rate is the percentage of test set samples that are correctly classified by the model
 - Most used for binary classes
 - If the accuracy is acceptable, use the model to classify new data
- Note: If the test set is used to select models, it is called validation (test) set

Process (1): Model Construction



Process (2): Using the Model in Prediction



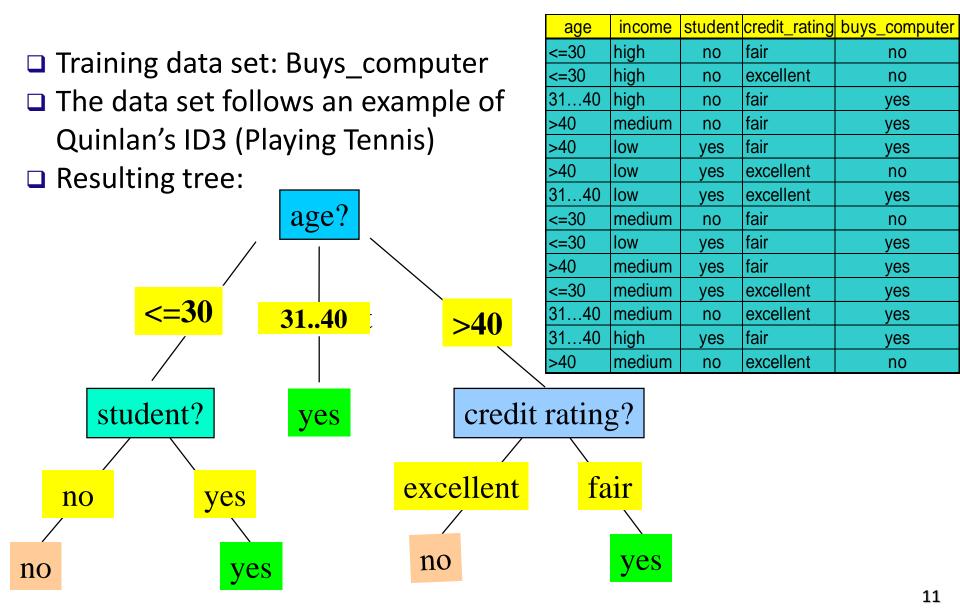
Classification Methods Overview

- Part 1
 - Decision Tree
 - Model Evaluation
- Part 2
 - Bayesian Learning: Naïve Bayes, Bayesian belief network
 - Logistic Regression
- Part 3
 - SVM
 - kNN
 - Other Topics

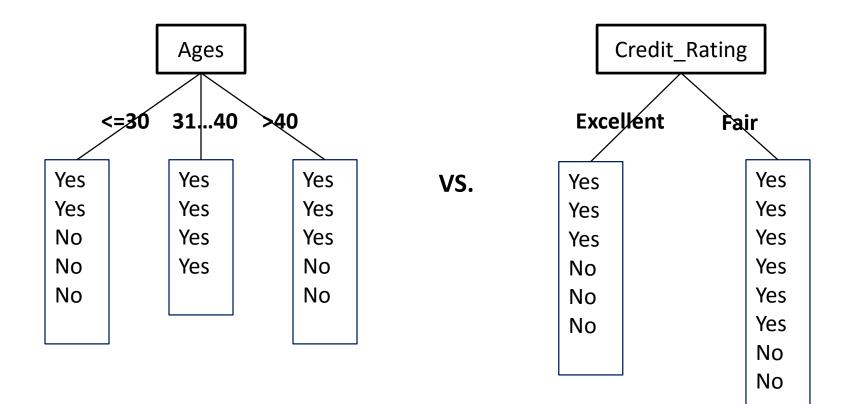
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Decision Tree Induction: An Example

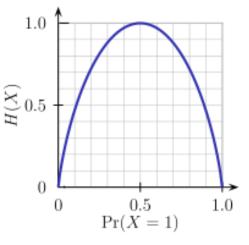


How to choose attributes?



Brief Review of Entropy

- Entropy (Information Theory)
 - A measure of uncertainty (impurity) associated with a random variable
 - Calculation: For a discrete random variable *Y* taking *m* distinct values {*y*₁, ..., *y_m*},
 - $H(Y) = -\sum_{i=1}^{m} p_i \log(p_i)$, where $p_i = P(Y = y_i)$
 - Interpretation:
 - Higher entropy => higher uncertainty
 - Lower entropy => lower uncertainty
- Conditional Entropy
 - $H(Y|X) = \sum_{x} p(x)H(Y|X=x)$



m = 2

Attribute Selection Measure: Information Gain (ID3/C4.5)

- Select the attribute with the highest information gain
- Let p_i be the probability that an arbitrary tuple in D belongs to class C_i, estimated by |C_{i, D}|/|D|
- Expected information (entropy) needed to classify a tuple in D: $Info(D) = -\sum_{i=1}^{m} p_i \log_2(p_i)$
- Information needed (after using A to split Dⁱ⁼¹ into v partitions) to classify D (conditional entropy): $Info_A(D) = \sum_{j=1}^{\nu} \frac{|D_j|}{|D|} \times Info(D_j)$
- Information gained by branching on attribute A

$$Gain(A) = Info(D) - Info_A(D)$$

Attribute Selection: Information Gain

| Class N: buys_computer = "no" | | | | | | | |
|-------------------------------|--------------|---------------------|---------------------|-------------------------------------|-----|--|--|
| Info(D) |) = I(9,5) = | $-\frac{9}{14}\log$ | $g_2(\frac{9}{14})$ | $-\frac{5}{14}\log_2(\frac{5}{14})$ |)=0 | | |
| | age | p _i | n _i | l(p _i , n _i) | | | |
| | <=30 | 2 | 3 | 0.971 | | | |
| | 3140 | 4 | 0 | 0 | | | |
| | >40 | 3 | 2 | 0.971 | | | |

Class P: buys_computer = "yes"

| age | income | student | credit_rating | buys_computer |
|------|--------|---------|---------------|---------------|
| <=30 | high | no | fair | no |
| <=30 | high | no | excellent | no |
| 3140 | high | no | fair | yes |
| >40 | medium | no | fair | yes |
| >40 | low | yes | fair | yes |
| >40 | low | yes | excellent | no |
| 3140 | low | yes | excellent | yes |
| <=30 | medium | no | fair | no |
| <=30 | low | yes | fair | yes |
| >40 | medium | yes | fair | yes |
| <=30 | medium | yes | excellent | yes |
| 3140 | medium | no | excellent | yes |
| 3140 | high | yes | fair | yes |
| >40 | medium | no | excellent | no |

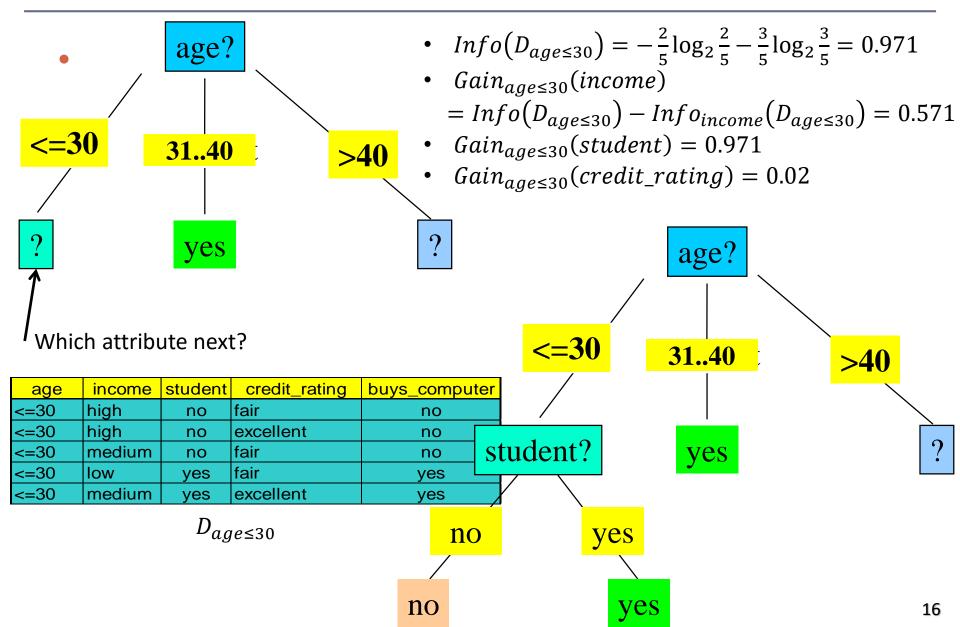
 $Info_{age}(D) = \frac{5}{14}I(2,3) + \frac{4}{14}I(4,0)$ 0.940 $+ \frac{5}{14}I(3,2) = 0.694$

 $\frac{5}{14}I(2,3)$ means "age <=30" has 5 out of 14 samples, with 2 yes'es and 3 no's. Hence

 $Gain(age) = Info(D) - Info_{age}(D) = 0.246$ Similarly,

Gain(income) = 0.029 Gain(student) = 0.151Gain(credit rating) = 0.048

Attribute Selection for a Branch



Algorithm for Decision Tree Induction

- Basic algorithm (a greedy algorithm)
 - Tree is constructed in a top-down recursive divide-and-conquer manner
 - At start, all the training examples are at the root
 - Attributes are categorical (if continuous-valued, they are discretized in advance)
 - Examples are partitioned recursively based on selected attributes
 - Test attributes are selected on the basis of a heuristic or statistical measure (e.g., information gain)
- Conditions for stopping partitioning
 - All samples for a given node belong to the same class
 - There are no remaining attributes for further partitioning majority voting is employed for classifying the leaf
 - There are no samples left use majority voting in the parent partition

Computing Information-Gain for Continuous-Valued Attributes

- Let attribute A be a continuous-valued attribute
- Must determine the *best split point* for A
 - Sort the value A in increasing order
 - Typically, the midpoint between each pair of adjacent values is considered as a possible *split point*
 - $(a_i + a_{i+1})/2$ is the midpoint between the values of a_i and a_{i+1}
 - The point with the *minimum expected information requirement* for A is selected as the split-point for A
- Split:
 - D1 is the set of tuples in D satisfying A ≤ split-point, and D2 is the set of tuples in D satisfying A > split-point

Gain Ratio for Attribute Selection (C4.5)

- Information gain measure is biased towards attributes with a large number of values
- C4.5 (a successor of ID3) uses gain ratio to overcome the problem (normalization to information gain)

SplitInfo_A(D) =
$$-\sum_{j=1}^{\nu} \frac{|D_j|}{|D|} \times \log_2(\frac{|D_j|}{|D|})$$

- GainRatio(A) = Gain(A)/SplitInfo(A)
- Ex. $SplitInfo_{income}(D) = -\frac{4}{14} \times \log_2\left(\frac{4}{14}\right) - \frac{6}{14} \times \log_2\left(\frac{6}{14}\right) - \frac{4}{14} \times \log_2\left(\frac{4}{14}\right) = 1.557$
 - gain_ratio(income) = 0.029/1.557 = 0.019
- The attribute with the maximum gain ratio is selected as the splitting attribute

Gini Index (CART, IBM IntelligentMiner)

• If a data set *D* contains examples from *n* classes, gini index, gini(*D*) is defined as $gini(D) = 1 - \sum_{j=1}^{\nu} p_j^2$

where p_j is the relative frequency of class j in D

- If a data set *D* is split on A into two subsets D_1 and D_2 , the gini index gini(*D*) is defined as $gini_A(D) = \frac{|D_1|}{|D|}gini(D_1) + \frac{|D_2|}{|D|}gini(D_2)$
- Reduction in Impurity:

$$\Delta gini(A) = gini(D) - gini_A(D)$$

 The attribute provides the smallest gini_{split}(D) (or the largest reduction in impurity) is chosen to split the node (need to enumerate all the possible splitting points for each attribute)

Computation of Gini Index

- Ex. D has 9 tuples in buys_computer = "yes" and 5 in "no" $gini(D) = 1 - \left(\frac{9}{14}\right)^2 - \left(\frac{5}{14}\right)^2 = 0.459$
- Suppose the attribute income partitions D into 10 in D₁: {low, medium} and 4 in D₂

$$\begin{split} gini_{income \in \{low, medium\}}(D) &= \left(\frac{10}{14}\right) Gini(D_1) + \left(\frac{4}{14}\right) Gini(D_2) \\ &= \frac{10}{14} \left(1 - \left(\frac{7}{10}\right)^2 - \left(\frac{3}{10}\right)^2\right) + \frac{4}{14} \left(1 - \left(\frac{2}{4}\right)^2 - \left(\frac{2}{4}\right)^2\right) \\ &= 0.443 \\ &= Gini_{income \in \{high\}}(D). \end{split}$$

Gini_{low,high} is 0.458; Gini_{medium,high} is 0.450. Thus, split on the {low,medium} (and {high}) since it has the lowest Gini index

Comparing Attribute Selection Measures

- The three measures, in general, return good results but
 - Information gain:
 - biased towards multivalued attributes
 - Gain ratio:
 - tends to prefer unbalanced splits in which one partition is much smaller than the others (why?)
 - Gini index:
 - biased to multivalued attributes

***Other Attribute Selection Measures**

- <u>CHAID</u>: a popular decision tree algorithm, measure based on χ^2 test for independence
- <u>C-SEP</u>: performs better than info. gain and gini index in certain cases
- <u>G-statistic</u>: has a close approximation to χ^2 distribution
- MDL (Minimal Description Length) principle (i.e., the simplest solution is preferred):
 - The best tree as the one that requires the fewest # of bits to both (1) encode the tree, and (2) encode the exceptions to the tree
- Multivariate splits (partition based on multiple variable combinations)
 - <u>CART</u>: finds multivariate splits based on a linear comb. of attrs.
- Which attribute selection measure is the best?
 - Most give good results, none is significantly superior than others

Overfitting and Tree Pruning

- <u>Overfitting</u>: An induced tree may overfit the training data
 - Too many branches, some may reflect anomalies due to noise or outliers
 - Poor accuracy for unseen samples
- Two approaches to avoid overfitting
 - <u>Prepruning</u>: *Halt tree construction early*-do not split a node if this would result in the goodness measure falling below a threshold
 - Difficult to choose an appropriate threshold
 - <u>Postpruning</u>: *Remove branches* from a "fully grown" tree—get a sequence of progressively pruned trees
 - Use a set of data different from the training data to decide which is the "best pruned tree"

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Model Evaluation and Selection

- Evaluation metrics: How can we measure accuracy?
 Other metrics to consider?
- Use validation test set of class-labeled tuples instead of training set when assessing accuracy
- Methods for estimating a classifier's accuracy:
 - Holdout method, random subsampling
 - Cross-validation

Classifier Evaluation Metrics: Confusion Matrix

Confusion Matrix:

| Actual class\Predicted class | C ₁ | ¬ C ₁ | |
|------------------------------|----------------------|----------------------|--|
| C ₁ | True Positives (TP) | False Negatives (FN) | |
| ¬ C ₁ | False Positives (FP) | True Negatives (TN) | |

Example of Confusion Matrix:

| Actual class\Predicted class | buy_computer = yes | buy_computer = no | Total |
|---------------------------------|-----------------------|----------------------|-------|
| buy_computer = yes | 6954 | 46 | 7000 |
| buy_computer = no | 412 | 2588 | 3000 |
| Total | 7366 | 2634 | 10000 |

- Given *m* classes, an entry, *CM*_{i,j} in a confusion matrix indicates # of tuples in class *i* that were labeled by the classifier as class *j*
- May have extra rows/columns to provide totals

Classifier Evaluation Metrics: Accuracy, Error Rate, Sensitivity and Specificity

| A\P | С | ¬C | |
|-----|----|----|-----|
| С | ΤР | FN | Ρ |
| −C | FP | ΤN | Ν |
| | P' | N' | All |

 Classifier Accuracy, or recognition rate: percentage of test set tuples that are correctly classified

Accuracy = (TP + TN)/All

Error rate: 1 – accuracy, or
 Error rate = (FP + FN)/All

- Class Imbalance Problem:
 - One class may be *rare*, e.g. fraud, or HIV-positive
 - Significant majority of the negative class and minority of the positive class
 - Sensitivity: True Positive recognition rate

Sensitivity = TP/P

- Specificity: True Negative recognition rate
 - Specificity = TN/N

Classifier Evaluation Metrics: Precision and Recall, and F-measures Precision: exactness – what % of tuples that the classifier labeled as positive are actually positive $\overline{TP + FP}$ precision Recall: completeness – what % of positive tuples did the classifier label as positive? TPrecall $\overline{TP + FN}$ Perfect score is 1.0 Inverse relationship between precision & recall • F measure (F₁ or F-score): harmonic mean of precision and $2 \times precision \times recall$ recall, precision + recall• F_{β} : weighted measure of precision and recall • assigns ß times as much weight to recall as to precision $F_{\beta} = \frac{(1+\beta^2) \times precision \times recall}{\beta^2 \times precision + recall}$

Classifier Evaluation Metrics: Example

| Actual Class\Predicted class | cancer = yes | cancer = no | Total | Recognition(%) |
|------------------------------|--------------|-------------|-------|---------------------------|
| cancer = yes | 90 | 210 | 300 | 30.00 (sensitivity) |
| cancer = no | 140 | 9560 | 9700 | 98.56 (specificity) |
| Total | 230 | 9770 | 10000 | 96.40 (<i>accuracy</i>) |

• *Precision* = 90/230 = 39.13%

Recall = 90/300 = 30.00%

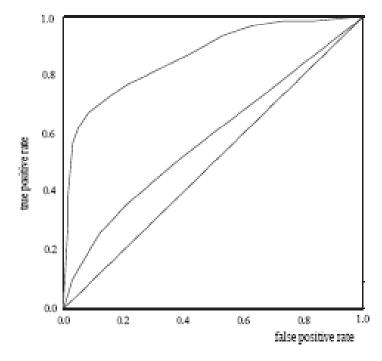
Evaluating Classifier Accuracy: Holdout & Cross-Validation Methods

Holdout method

- Given data is randomly partitioned into two independent sets
 - Training set (e.g., 2/3) for model construction
 - Test set (e.g., 1/3) for accuracy estimation
- <u>Random sampling</u>: a variation of holdout
 - Repeat holdout k times, accuracy = avg. of the accuracies obtained
- Cross-validation (k-fold, where k = 10 is most popular)
 - Randomly partition the data into *k mutually exclusive* subsets, each approximately equal size
 - At *i*-th iteration, use D_i as test set and others as training set
 - <u>Leave-one-out</u>: *k* folds where *k* = # of tuples, for small sized data
 - <u>*Stratified cross-validation</u>*: folds are stratified so that class dist. in each fold is approx. the same as that in the initial data

Model Selection: ROC Curves

- ROC (Receiver Operating Characteristics) curves: for visual comparison of classification models
- Originated from signal detection theory
- Shows the trade-off between the true positive rate and the false positive rate
- The area under the ROC curve is a measure of the accuracy of the model
- Rank the test tuples in decreasing order: the one that is most likely to belong to the positive class appears at the top of the list
- Area under the curve: the closer to the diagonal line (i.e., the closer the area is to 0.5), the less accurate is the model



- Vertical axis represents the true positive rate
- Horizontal axis rep. the false positive rate
- The plot also shows a diagonal line
- A model with perfect accuracy will have an area of 1.0

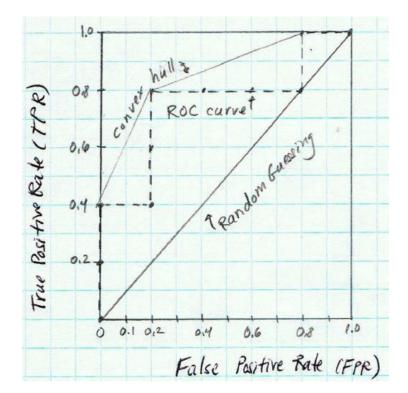
Plotting an ROC Curve

- True positive rate: TPR = TP/P (sensitivity)
- False positive rate: FPR = FP/N (1-specificity)

- Rank tuples according to how likely they will be a positive tuple
 - Idea: when we include more tuples in, we are more likely to make mistakes, that is the trade-off!
 - Nice property: not threshold (cut-off) need to be specified, only rank matters

| $Tuple \ \#$ | Class | Prob. | TP | FP | TN | FN | TPR | FPR |
|--------------|-------|-------|----|----|----|----|-----|-----|
| 1 | р | 0.9 | 1 | 0 | 5 | 4 | 0.2 | 0 |
| 2 | р | 0.8 | 2 | 0 | 5 | 3 | 0.4 | 0 |
| 3 | n | 0.7 | 2 | 1 | 4 | 3 | 0.4 | 0.2 |
| 4 | р | 0.6 | 3 | 1 | 4 | 2 | 0.6 | 0.2 |
| 5 | р | 0.55 | 4 | 1 | 4 | 1 | 0.8 | 0.2 |
| 6 | n | 0.54 | 4 | 2 | 3 | 1 | 0.8 | 0.4 |
| 7 | n | 0.53 | 4 | 3 | 2 | 1 | 0.8 | 0.6 |
| 8 | n | 0.51 | 4 | 4 | 1 | 1 | 0.8 | 0.8 |
| 9 | р | 0.50 | 5 | 4 | 0 | 1 | 1.0 | 0.8 |
| 10 | n | 0.4 | 5 | 5 | 0 | 0 | 1.0 | 1.0 |

Example



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Summary

- Classification is a form of data analysis that extracts models describing important data classes.
- decision tree induction
- Evaluation
 - Evaluation metrics include: accuracy, sensitivity, specificity, precision, recall, F measure, and F_{β} measure.
 - k-fold cross-validation is recommended for accuracy estimation.
 - ROC curves are useful for model selection.

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