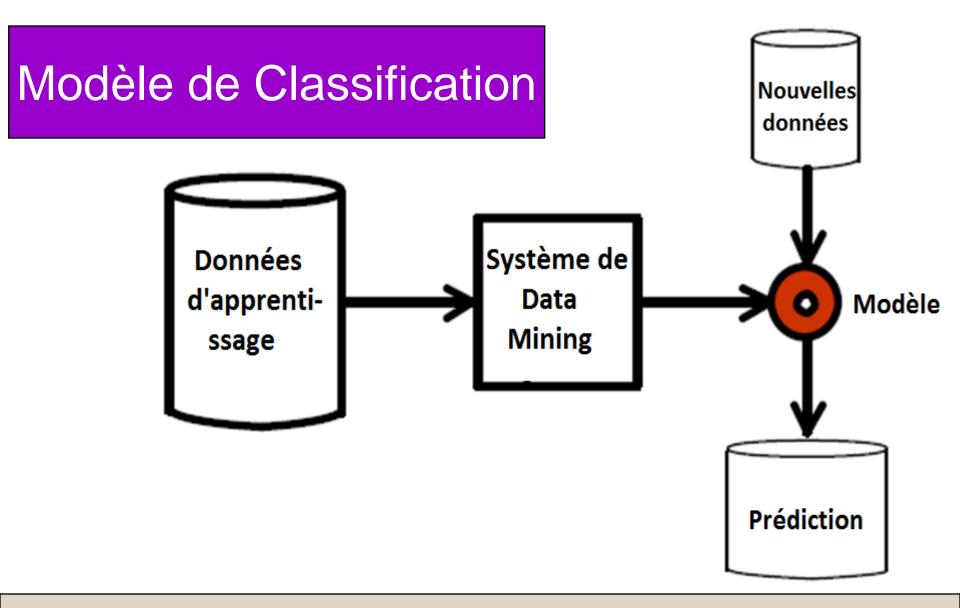
Classification



Bayesian Classification: Why?

- <u>A statistical classifier</u>: performs *probabilistic prediction*, *i.e.*, predicts class membership probabilities
- Foundation: Based on Bayes' Theorem.
- <u>Performance</u>: A simple Bayesian classifier, *naïve* Bayesian classifier, has comparable performance with decision tree and selected neural network classifiers
- Incremental: Each training example can incrementally increase/decrease the probability that a hypothesis is correct — prior knowledge can be combined with observed data
- <u>Standard</u>: Even when Bayesian methods are computationally intractable, they can provide a standard of optimal decision making against which other methods can be measured

Bayesian Theorem: Basics

- Let **X** be a data sample ("evidence"): class label is unknown
- Let H be a *hypothesis* that X belongs to class C
- Classification is to determine P(H|X), the probability that the hypothesis holds given the observed data sample X
- P(H) (*prior probability*), the initial probability
 - E.g., X will buy computer, regardless of age, income, ...
- P(X): probability that sample data is observed
- P(X|H) (*posteriori probability*), the probability of observing the sample X, given that the hypothesis holds
 - E.g., Given that X will buy computer, the prob. that X is 31..40, medium income

Bayesian Theorem

Given training data X, posteriori probability of a hypothesis
 H, P(H|X), follows the Bayes theorem

$$P(H|\mathbf{X}) = \frac{P(\mathbf{X}|H)P(H)}{P(\mathbf{X})}$$

- Informally, this can be written as
 posteriori = likelihood x prior/evidence
- Predicts **X** belongs to C_2 iff the probability $P(C_i|\mathbf{X})$ is the highest among all the $P(C_k|X)$ for all the *k* classes
- Practical difficulty: require initial knowledge of many probabilities, significant computational cost

Towards Naïve Bayesian Classifier

- Let D be a training set of tuples and their associated class labels, and each tuple is represented by an n-D attribute vector X = (x₁, x₂, ..., x_n)
- Suppose there are *m* classes C₁, C₂, ..., C_m.
- Classification is to derive the maximum posteriori, i.e., the maximal P(C_i|X)
- This can be derived from Bayes' theorem

$$P(C_i | \mathbf{X}) = \frac{P(\mathbf{X} | C_i) P(C_i)}{P(\mathbf{X})}$$

 Since P(X) is constant for all classes, only needs to be maximized

$$P(C_i | \mathbf{X}) = P(\mathbf{X} | C_i) P(C_i)$$

Derivation of Naïve Bayes Classifier

- A simplified assumption: attributes are conditionally independent (i.e., no dependence relation between attributes): $P(\mathbf{X}|_{C_{i}}) = \prod_{k=1}^{n} P(x_{k}|_{C_{i}}) = P(x_{1}|_{C_{i}}) \times P(x_{2}|_{C_{i}}) \times ... \times P(x_{n}|_{C_{i}})$
- This greatly reduces the computation cost: Only counts
 the class distribution
- If A_k is categorical, P(x_k|C_i) is the # of tuples in C_i having value x_k for A_k divided by |C_{i, D}| (# of tuples of C_i in D)
- If A_k is continous-valued, $P(x_k|C_i)$ is usually computed based on Gaussian distribution with a mean μ and standard deviation σ 1 $-\frac{(x-\mu)^2}{2}$

$$g(x,\mu,\sigma) = \frac{1}{\sqrt{2\pi\sigma}} e^{-\frac{(x-\mu)^2}{2\sigma^2}}$$

and $P(x_k|C_i)$ is

$$P(\mathbf{X} | C_i) = g(x_k, \mu_{C_i}, \sigma_{C_i})$$

Naïve Bayesian Classifier: Training Dataset

Class:

C1:buys_computer = 'yes' C2:buys_computer = 'no'

Data sample X = (age <=30, Income = medium, Student = yes Credit_rating = Fair)

age	income	student	credit_rating	_com
<=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

Naïve Bayesian Classifier: An Example

- P(C_i): P(buys_computer = "yes") = 9/14 = 0.643
 P(buys_computer = "no") = 5/14 = 0.357
- Compute P(X|C_i) for each class
 P(age = "<=30" | buys_computer = "yes") = 2/9 = 0.222</p>
 P(age = "<= 30" | buys_computer = "no") = 3/5 = 0.6</p>
 P(income = "medium" | buys_computer = "yes") = 4/9 = 0.444
 P(income = "medium" | buys_computer = "no") = 2/5 = 0.4
 P(student = "yes" | buys_computer = "yes) = 6/9 = 0.667
 P(student = "yes" | buys_computer = "no") = 1/5 = 0.2
 P(credit_rating = "fair" | buys_computer = "yes") = 6/9 = 0.667
 P(credit_rating = "fair" | buys_computer = "no") = 2/5 = 0.4
- X = (age <= 30, income = medium, student = yes, credit_rating = fair)

Therefore, X belongs to class ("buys_computer = yes")

Avoiding the 0-Probability Problem

 Naïve Bayesian prediction requires each conditional prob. be nonzero. Otherwise, the predicted prob. will be zero

$$P(X \mid C_i) = \prod_{k=1}^{n} P(x_k \mid C_i)$$

- Ex. Suppose a dataset with 1000 tuples, income=low (0), income= medium (990), and income = high (10),
- Use Laplacian correction (or Laplacian estimator)
 - Adding 1 to each case

Prob(income = low) = 1/1003

Prob(income = medium) = 991/1003

Prob(income = high) = 11/1003

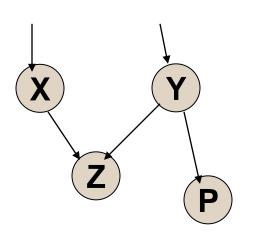
The "corrected" prob. estimates are close to their "uncorrected" counterparts

Naïve Bayesian Classifier: Comments

- Advantages
 - Easy to implement
 - Good results obtained in most of the cases
- Disadvantages
 - Assumption: class conditional independence, therefore loss of accuracy
 - Practically, dependencies exist among variables
 - E.g., hospitals: patients: Profile: age, family history, etc. Symptoms: fever, cough etc., Disease: lung cancer, diabetes, etc.
 - Dependencies among these cannot be modeled by Naïve Bayesian Classifier
- How to deal with these dependencies?
 - Bayesian Belief Networks

Bayesian Belief Networks

- Bayesian belief network allows a *subset* of the variables conditionally independent
- A graphical model of causal relationships
 - Represents <u>dependency</u> among the variables
 - Gives a specification of joint probability distribution



- □ Nodes: random variables
- □ Links: dependency
- □ X and Y are the parents of Z, and Y is the parent of P
- $\hfill\square$ No dependency between Z and P
- □ Has no loops or cycles

Classification by Backpropagation

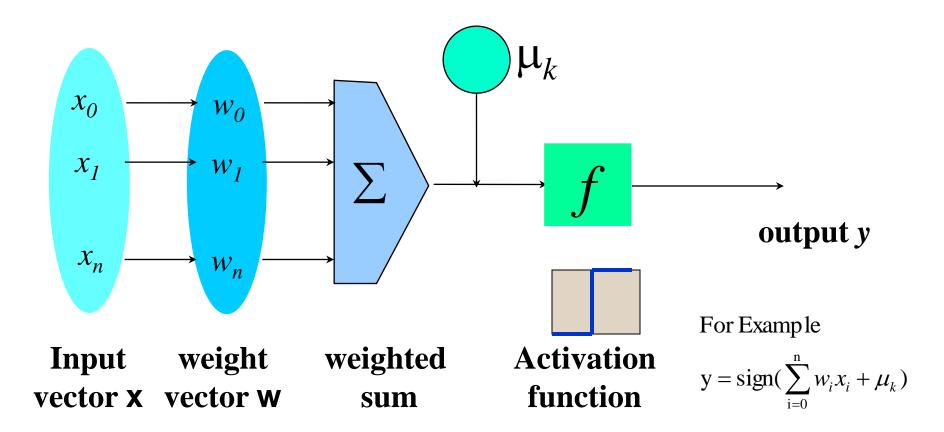
- Backpropagation: A **neural network** learning algorithm
- Started by psychologists and neurobiologists to develop and test computational analogues of neurons
- A neural network: A set of connected input/output units where each connection has a weight associated with it
- During the learning phase, the network learns by adjusting the weights so as to be able to predict the correct class label of the input tuples
- Also referred to as connectionist learning due to the connections between units

Neural Network as a Classifier

• Weakness

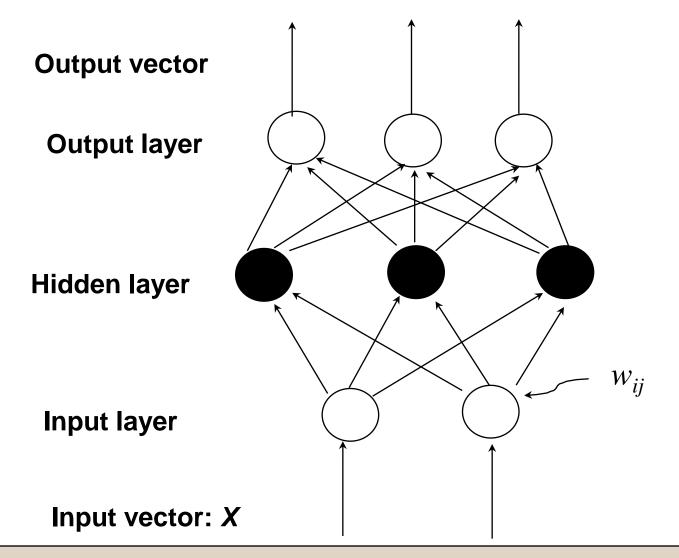
- Long training time
- Require a number of parameters typically best determined empirically, e.g., the network topology or ``structure."
- Poor interpretability: Difficult to interpret the symbolic meaning behind the learned weights and of ``hidden units" in the network
- Strength
 - High tolerance to noisy data
 - Ability to classify untrained patterns
 - Well-suited for continuous-valued inputs and outputs
 - Successful on a wide array of real-world data
 - Algorithms are inherently parallel
 - Techniques have recently been developed for the extraction of rules from trained neural networks

A Neuron (= a perceptron)



• The *n*-dimensional input vector **x** is mapped into variable y by means of the scalar product and a nonlinear function mapping

A Multi-Layer Feed-Forward Neural Network



How A Multi-Layer Neural Network Works?

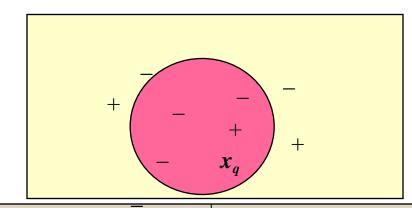
- The **inputs** to the network correspond to the attributes measured for each training tuple
- Inputs are fed simultaneously into the units making up the input layer
- They are then weighted and fed simultaneously to a hidden layer
- The number of hidden layers is arbitrary, although usually only one
- The weighted outputs of the last hidden layer are input to units making up the **output layer**, which emits the network's prediction
- The network is **feed-forward** in that none of the weights cycles back to an input unit or to an output unit of a previous layer
- From a statistical point of view, networks perform **nonlinear regression**: Given enough hidden units and enough training samples, they can closely approximate any function

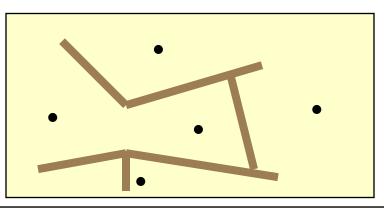
Backpropagation

- Iteratively process a set of training tuples & compare the network's prediction with the actual known target value
- For each training tuple, the weights are modified to minimize the mean squared error between the network's prediction and the actual target value
- Modifications are made in the "backwards" direction: from the output layer, through each hidden layer down to the first hidden layer, hence "backpropagation"
- Steps
 - Initialize weights (to small random #s) and biases in the network
 - Propagate the inputs forward (by applying activation function)
 - Backpropagate the error (by updating weights and biases)
 - Terminating condition (when error is very small, etc.)

The *k*-Nearest Neighbor Algorithm

- All instances correspond to points in the n-D space
- The nearest neighbor are defined in terms of Euclidean distance, dist(X₁, X₂)
- Target function could be discrete- or real- valued
- For discrete-valued, *k*-NN returns the most common value among the *k* training examples nearest to *x*_q
- Vonoroi diagram: the decision surface induced by 1-NN for a typical set of training examples





Discussion on the k-NN Algorithm

- k-NN for real-valued prediction for a given unknown tuple
 - Returns the mean values of the k nearest neighbors
- Distance-weighted nearest neighbor algorithm
 - Weight the contribution of each of the k neighbors according to their distance to the query x_a $w \equiv \frac{1}{d(x_q, x_i)^2}$
 - Give greater weight to closer neighbors
- Robust to noisy data by averaging k-nearest neighbors
- Curse of dimensionality: distance between neighbors could be dominated by irrelevant attributes
 - To overcome it, axes stretch or elimination of the least relevant attributes