### Data Mining - Classification I

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### Overview

- Introduction
  - Context
  - Overfitting

- Classification algorithms
  - k-neareast neighbor

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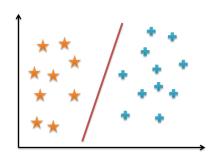
### Classification vs Clustering

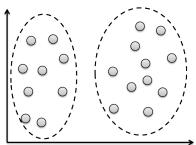
#### Classification

- Labeled data objects
- Assign a label to new objects

#### Clustering

- Data is not labeled
- Identify structure in data and group objects

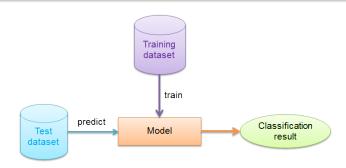




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#### Objectives

- Classification is the supervised learning task of data mining that predicts categorical class labels (discrete or nominal)
- A model is built based on the training set and the class labels.
  This model helps in classifying new data



### **Applications**

• Banking: card fraud detection



### **Applications**

 Pattern Recognition: Plate Recognition, Optimal Character Recognition





### **Applications**

• Security: Finger Print, Spam Filter.





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### **Applications**

• Transport: Self-driving vehicules.



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#### **Problematics**

#### **Difficulties**

- Training datasets: overfitting, imbalance of data, ...
- Information about data class
- Classification algorithms: complexity, training time, etc...
- Big data
- Performance

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#### **Notions**

```
Input set: (\mathcal{X}, \mathcal{Y}) = \{(\mathbf{x}_i, y_i)\}, \forall i = 1, ..., n: a set of n training objects
```

Object:  $\mathbf{x}_i \in \mathbb{R}^d$  with  $\mathbf{x}_i = (x_{i,1}, x_{i,2}, ..., x_{i,d}), y_i \in \mathcal{Y}$  ( $y_i$  is the class label of  $\mathbf{x}_i$ 

Model: Model  ${\mathcal M}$  allows to classify  ${\mathcal X}$  into many groups according to its  ${\mathcal Y}$ .

Problem: For new object (test data)  $\mathbf{x}_0 \notin \mathcal{X}$  and unknown  $y_0$ , the model  $\mathcal{M}$  is employed to predict  $y_0$ .

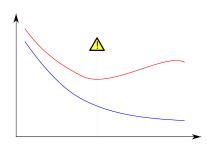
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## Overfitting

#### Definition

In data mining, overfitting problem occurs:

- A model is excessively complex, such as having too many parameters relative to the number of training objects.
- A model has poor predictive performance, as it overreacts due to the training data.



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# Overfitting

#### Overfitting

Many approaches to manage training datasets have been proposed to reduce variance and avoid overfitting:

- cross validation: leave-p-out, k-fold cross-validation
- bagging or bootstrap agrgregating
- and many more....

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#### Cross-validation

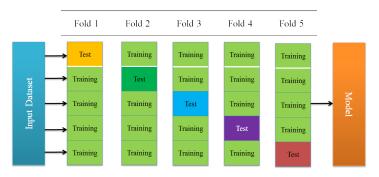
This technique is a statistics model:

- generalize to an independent data set
- involve partitioning a sample of data into complementary subsets: building the model from one subset (the training set), and validating the model on the other subset (the testing set)

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#### k-fold cross validation

- generalize to k subsets
- k-1 subsets for training and 1 subset for testing



5-fold cross validation

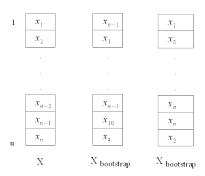
#### Leave-p-out

- Leave-p-out take p data out from the input dataset of n data, n-p for training and p for testing
- This cross-validation requires to learn and validate  $C_n^p$  times (if n=100 and p=30,  $C_{100}^{30}=3*10^{25}$ )
- Leave-one-out is often used as for  $C_n^1 = n$

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#### Bagging

- consider input set  $\mathcal{X}$  with n training objects
- create k new training sets  $\mathcal{X}_i$  by drawing n objects with replacement



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#### Definition

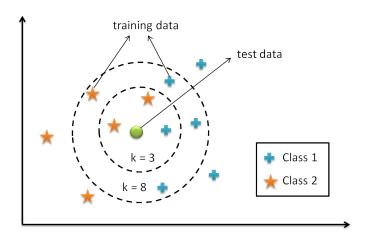
#### k-NN

k-Nearest Neighbors algorithm (or k-NN for short) is a non-parametric method often used for classification.

- The input consists of the k closest training examples in the feature space.
- The output is a class label. An object is classified by a majority vote of its neighbors, with the object being assigned to the class most common among its k nearest neighbors.
- ullet k is a positive integer, typically small for example k = 1, 3, 5, 10, etc.

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### k-NN



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### Pros and Cons

#### Advantages

- No input parametric
- No training step required
- Simple to understand and to interpret

#### Disadvantages

- Require many operations for testing
- Low accuracy comparing to Decision Tree, Random Forest, SVM, Neural Networks, etc...

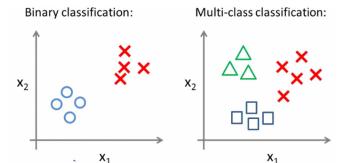
```
Input set: (\mathcal{X}, \mathcal{Y}) = \{(\mathbf{x}_i, y_i)\}, \forall i = 1, ..., n: a set of N training objects
```

Object: 
$$\mathbf{x}_i \in \mathbb{R}^d$$
 with  $\mathbf{x}_i = (x_{i,1}, x_{i,2}, ..., x_{i,d}), y_i \in \mathcal{Y}$  ( $y_i$  is  $\{-1; +1\}$ 

Model: A learn classifier  $f(\mathbf{x}_i)$  such that

$$f(\mathbf{x}_i) = \begin{cases} \geq 0 & y_i = +1 \\ < 0 & y_i = -1 \end{cases}$$

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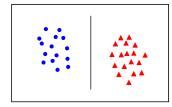


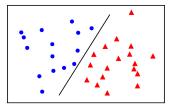
#### Binary vs Multi-class classification

A multi-class classifier is able to classify into more 2 classes. Binary classifier can only deal with 2-class problems

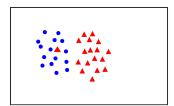
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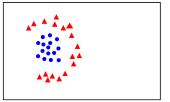
linearly separable

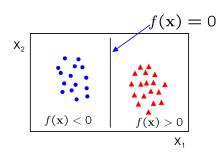




not linearly separable







A linear classifier has the form:

$$f(\mathbf{x}) = \mathbf{w}\mathbf{x}^T + b$$

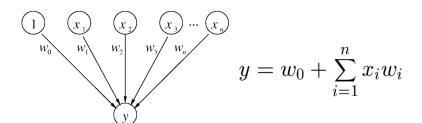
- in 2D the discriminant is a line, in kD  $(k \ge 3)$ , is a hyperplane denoted by  $\Delta(\mathbf{v}, a) = \{\mathbf{x} \in \mathbb{R}^d | \mathbf{v}\mathbf{x}^T + a = 0\}$
- w is the normal to the line, and b the bias
- w is known as the weight vector

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#### Perceptron

- An algorithm for supervised learning of binary classifiers.
- A type of linear classifier.



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Given linearly separable data  $\mathcal{X}$  labelled into two categories  $y_i = \{-1, 1\}$ , find a weight vector  $\mathbf{w}$  such that the discriminant function:

$$f(\mathbf{x}) = \mathbf{w}\mathbf{x}^T$$

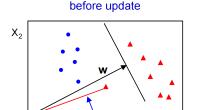
#### Perceptron Algorithm

- intialize  $\mathbf{w}_0 = \mathbf{0}$  (or close to  $\mathbf{0}$ )
- for all  $\mathbf{x}_i \in \mathcal{X}$ , if  $\mathbf{x}_i$  is misclassified

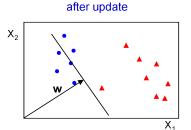
$$\mathbf{w}_t = \mathbf{w}_{t-1} + \alpha sign(f(\mathbf{x}_i))\mathbf{x}_i$$

• until all the data is correctly classified

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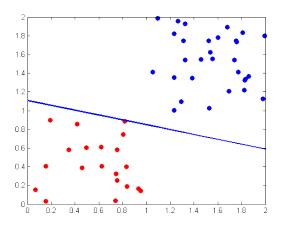
 $\mathbf{x}_i$ 



 $\mathbf{w} \leftarrow \mathbf{w} - \alpha \mathbf{x}_i$ 

if the data is linearly separable, then the algorithm will converge. After the convergence,  $\mathbf{w} = \alpha \sum \mathbf{x}_i$ .

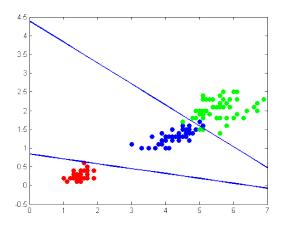
 $X_1$ 



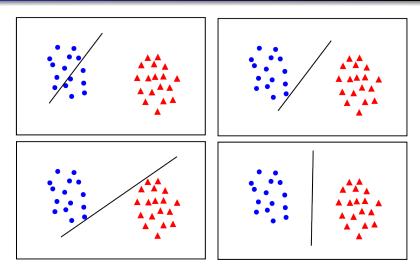
Perceptron Classifier for 2-class dataset

- Convergence can be slow and Perceptron algorithm depends heavily on input parameters  $\mathbf{w}_0$  and  $\alpha$
- Perceptron is able to handle 2-class problems.
- How to use Perceptron for multi-class problems?
- What if we have 3 classes such as Iris dataset?

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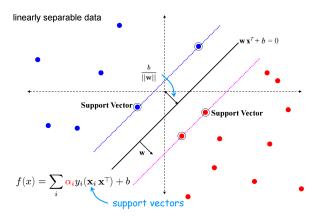
Perceptron Classifier for Iris dataset



Are all decision boundaries equally good? or what is the best  $\mathbf{w}$ ?

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The decision boundary or the hyperplan should be as far away from the data of both classes as possible.



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A support vector machine (SVM) is a supervised learning technique from the field of machine learning applicable to both classification and regression.

- Rooted in the Statistical Learning Theory developed by Vladimir Vapnik and co-workers at AT&T Bell Laboratories in 1995
- Very famous due to its success for hand writting recognitions (or image classification in general)

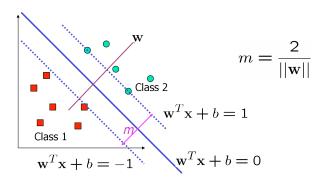
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- SVMs are based on the principle of Structural Risk Minimization.
- Non-linearly map the input space into a very high dimensional feature space in order to construct an optimal separating hyperplane in this space (a maximal margin classifier)

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The margin, the smallest distance between the decision boundary  $\Delta(\mathbf{v}, a)$  and the examples  $\mathbf{x}_i$ , must be maximized.

$$m = min_{i=1..n} dist(\mathbf{x}_i, \Delta(\mathbf{v}, a))$$



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Let  $\mathbf{x}$  be a vector in  $\mathbb{R}^d$  and  $\Delta(\mathbf{v}, a) = {\mathbf{s} \in \mathbb{R}^d | \mathbf{v}\mathbf{s}^T + a = 0}$  an hyperplane. The distance between  $\mathbf{x}$  and the hyperplan  $\Delta(\mathbf{v}, a)$  is

$$dist(\mathbf{x}, \Delta(\mathbf{v}, a)) = \frac{|\mathbf{v}\mathbf{x}^T + a|}{||v||}$$

Let  $\mathbf{s}_{x}$  be the closest point to  $\mathbf{x} \in \Delta, \mathbf{s}_{x} = argmin||\mathbf{x} - \mathbf{s}||$ ,

$$\mathbf{x} = \mathbf{s}_{x} + m \frac{\mathbf{v}}{||\mathbf{v}||} \iff \mathbf{x} - \mathbf{s}_{x} = m \frac{\mathbf{v}}{||\mathbf{v}||}$$

So that, taking the scalar product with vector  $\mathbf{v}$  we have:

$$\mathbf{v} \times m \frac{\mathbf{v}^T}{\mathbf{v}} = \mathbf{v}(\mathbf{x}^T - \mathbf{s}_x^T) = \mathbf{v}\mathbf{x}^T - \mathbf{v}\mathbf{s}_x^T = \mathbf{v}\mathbf{x}^T + a - (\mathbf{v}\mathbf{s}_x^T + a) = \mathbf{v}\mathbf{x}^T + a$$

$$\mathbf{vs}_{\times}^T + a = 0$$
 because  $\mathbf{s}_{\times} \in \Delta$ 

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So, therefore

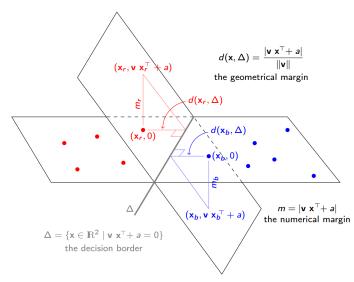
$$\mathbf{v} \times m \frac{\mathbf{v}^T}{||\mathbf{v}||} = m \frac{||\mathbf{v}||^2}{||\mathbf{v}||} = m||\mathbf{v}|| = \mathbf{v}\mathbf{x}^T + a$$

Thus,

$$m = \frac{\mathbf{v}\mathbf{x}^T + a}{||\mathbf{v}||}$$

and

$$dist(\mathbf{x}, \Delta(\mathbf{v}, a)) = \frac{|\mathbf{v}\mathbf{x}^T + a|}{||v||}$$



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The decision hyperplane  $\Delta(\mathbf{v}, a) = \{\mathbf{s} \in \mathbb{R}^d | \mathbf{v}\mathbf{s}^T + a = 0\}$ , we have to maximize the margin

$$\max_{\mathbf{v},a} \min_{i \in [1..n]} dist(\mathbf{x}_i, \Delta(\mathbf{v}, a))$$

Maximize the confidence

$$\begin{cases} \max_{\mathbf{v},a} m & \max_{i \in [1..n]} \frac{|\mathbf{v}\mathbf{x}^T + a|}{||\mathbf{v}||} \ge m & \text{with } \frac{y_i(\mathbf{v}\mathbf{x}^T + a)}{||\mathbf{v}||} \ge m \end{cases}$$

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Change variable  $\mathbf{w} = \frac{\mathbf{v}}{m||\mathbf{v}||} \implies ||\mathbf{w}|| = \frac{1}{m}$  and  $b = \frac{a}{m||\mathbf{v}||}$  Maximize the confidence

$$\begin{cases} \max_{\mathbf{v}, a} & m \\ \text{with} & \frac{y_i(\mathbf{v}\mathbf{x}^T + a)}{||v||} \ge m \end{cases}$$

$$\begin{cases} \max_{\mathbf{w},b} m \\ \text{with } y_i(\mathbf{wx}^T + b) \ge 1; i = 1..n \end{cases} \begin{cases} \min_{\mathbf{w},a} ||\mathbf{w}|| \\ \text{with } y_i(\mathbf{wx}^T + b) \ge 1; i = 1..n \end{cases}$$

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#### Linear SVMs are the solution of the following problem (called primal)

Let  $\{(\mathbf{x}_i, y_i); i = 1..n\}$  be a set of labelled data with  $\mathbf{x} \in \mathbb{R}^d, y_i \in \{1, -1\}$ . A support vector machine is a linear classifier associated with the following decision function:  $D(x) = sign(\mathbf{w}\mathbf{x}^T + b)$  where  $\mathbf{w} \in \mathbb{R}^d$  and  $b \in \mathbb{R}$  a given thought the solution of the following problem:

$$\begin{cases} \min_{\mathbf{w} \in \mathbf{R}^d, b \in \mathbb{R}} & \frac{1}{2} ||\mathbf{w}||^2 \\ \text{with} & y_i(\mathbf{w}\mathbf{x}^T + b) \ge 1; i = 1..n \end{cases}$$

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Minimize  $\frac{1}{2}||\mathbf{w}||$  subject to  $1 - y_i(\mathbf{w}\mathbf{x}^T + b) \le 0$  for i = 1..n The Lagrangian is:

$$\mathcal{L} = \frac{1}{2} \mathbf{w} \mathbf{w}^T + \sum_{i=1}^n \alpha_i (1 - y_i (\mathbf{w} \mathbf{x}^T + b))$$

Note that  $||\mathbf{w}|| = \mathbf{w}\mathbf{w}^T$  and  $\alpha_I \geq 0$ 

$$\mathcal{L} = \frac{1}{2} \mathbf{w} \mathbf{w}^T + \sum_{i=1}^n \alpha_i (1 - y_i (\mathbf{w} \mathbf{x}^T + b))$$

Setting the gradient of  $\mathcal{L}$  w.r.t.  $\mathbf{w}$  and  $\mathbf{b}$  to zero, we have:

$$\frac{\delta \mathcal{L}}{\delta \mathbf{w}} =$$

$$\frac{\delta \mathcal{L}}{\delta h} =$$

If we substitue  $\mathbf{w} = \sum_{i=1}^{n} \alpha_i y_i \mathbf{x}_i$  to  $\mathcal{L}$ , we have

$$\mathcal{L} = \sum_{i=1}^{n} \alpha_i - \frac{1}{2} \sum_{i=1}^{n} \sum_{j=1}^{n} \alpha_i \alpha_j y_i y_j \mathbf{x}_i \mathbf{x}_j^T$$

- Note that  $\sum_{i=1}^{n} \alpha_i y_i = 0$
- ullet This is a function of lpha
- It is known as the dual problem: if we know  $\mathbf{w}$ , we know all  $\alpha$ ; or if we know all  $\alpha$ , we know  $\mathbf{w}$

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Thue dual problem is thus:

$$\max \sum_{i=1}^{n} \alpha_i - \frac{1}{2} \sum_{i=1}^{n} \sum_{j=1}^{n} \alpha_i \alpha_j y_i y_j \mathbf{x}_i \mathbf{x}_j^T$$

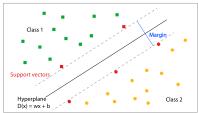
subject to  $\alpha_i \geq 0$  and  $\sum_{i=1}^n \alpha_i y_i = 0$ 

- $\mathbf{x}_i$  with non-zero  $\alpha_i$  are called support vectors (SV)
- $\mathbf{w} = \sum_{i=1}^{n} \alpha_i y_i \mathbf{x}_i$  is linear combination of a small number of data points
- To test new data z
  - compute  $\mathbf{wz}^T + b = \sum_{i=1}^s \alpha_i y_i(\mathbf{x}_i \mathbf{z}^T) + b$
  - if the sum is positive, z is classified as class 1 or class 2 otherwise.

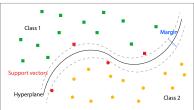
Classification

 Up to now, we have only seen large-margin classifier with a linear decision boundary. How to do if our problem is nonlinear?

#### Linear separation



#### Non-linear separation



• To handle non-linearly separable problems, we use error  $\xi$ 

$$\begin{cases} \mathbf{w}^{T} \mathbf{x}_{i} + b \geq 1 - \xi_{i} & y_{i} = 1 \\ \mathbf{w}^{T} \mathbf{x}_{i} + b \leq -1 + \xi_{i} & y_{i} = -1 \\ \xi_{i} \geq 0 & \forall i \end{cases}$$

•  $\xi_i$  are slack variables in optimization,  $\xi = 0$  if there is no error for  $\mathbf{x}_i$ 

We want to minimize

$$\frac{1}{2}||\mathbf{w}||^2 + C\sum_{i=1}^n \xi_i$$

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• The dual of this new problem is:

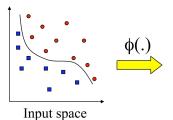
$$\max \sum_{i=1}^{n} \alpha_i - \frac{1}{2} \sum_{i=1}^{n} \sum_{j=1}^{n} \alpha_i \alpha_j y_i y_j \mathbf{x}_i \mathbf{x}_j^T$$

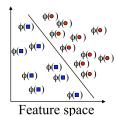
subject to 
$$C \ge \alpha_i \ge 0, \sum_{i=1}^n \alpha_i y_i = 0$$

• This is similar to the problem in the linear separable case.

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- Up to now, we have only seen large-margin classifier with a linear decision boundary. How to do if our problem is nonlinear?
- Idea: transform x<sub>i</sub> to a higher dimensional space to "make life easier"





Note: feature space is of higher dimension than the input space in practice

- Computation in the feature space is costly because it is high dimensional
- Kernel functions are used to overcome this.

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• The dual of this new problem is:

$$\max \sum_{i=1}^{n} \alpha_i - \frac{1}{2} \sum_{i=1}^{n} \sum_{j=1}^{n} \alpha_i \alpha_j y_i y_j \mathbf{x}_i \mathbf{x}_j^T$$

subject to 
$$C \ge \alpha_i \ge 0, \sum_{i=1}^n \alpha_i y_i = 0$$

- We just calculate the inner product without explicite the mapping of original data to the feature space.
- Geometric similarity measures can be expressed by inner products.
- Kernal function is often defined as:

$$K(\mathbf{x}_i, \mathbf{x}_j) = \phi(\mathbf{x}_i)\phi(\mathbf{x}_j)^T$$

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#### Examples of kernel functions

Linear kernels

$$K(\mathbf{x}, \mathbf{y}) = \mathbf{x}\mathbf{y}^T$$

Polynomial kernels with degree d

$$K(\mathbf{x}, \mathbf{y}) = (\mathbf{x}\mathbf{y}^T + 1)^d$$

Gaussian kernels

$$K(\mathbf{x},\mathbf{y}) = e^{\frac{-||\mathbf{x}-\mathbf{y}||^2}{2\sigma^2}}$$

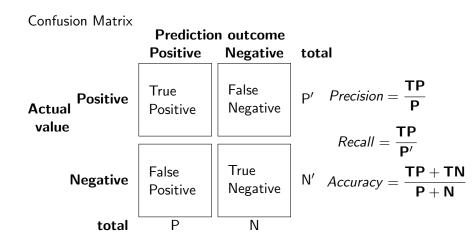
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Special transformation from  $\mathbb{R}^2 o \mathbb{R}^3$ 

$$\phi(\mathbf{x}) = \phi : (\mathbf{x}_1, \mathbf{x}_2) \to (\mathbf{x}_1^2, \mathbf{x}_2^2, \sqrt{2}\mathbf{x}_1\mathbf{x}_2)$$
$$\phi(\mathbf{x})\phi(\mathbf{y})^T = (\mathbf{x}\mathbf{y}^T)^2$$

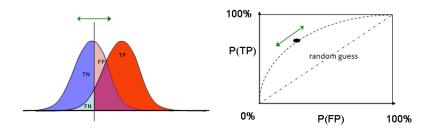
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#### **Evaluation**



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#### **Evaluation**



A receiver operating characteristic (ROC), or ROC curve, is a graphical plot that illustrates the performance of a binary classifier system.

 True positive rate (TPR) vs False positive rate (FPR) at various threshold settings.

Ocst vs benefit analysis of decision making Nhat-Quang Doan Classification

#### Evaluation

To compare different binary classifiers, AUC (Area under ROC curve) becomes a good criterion.

