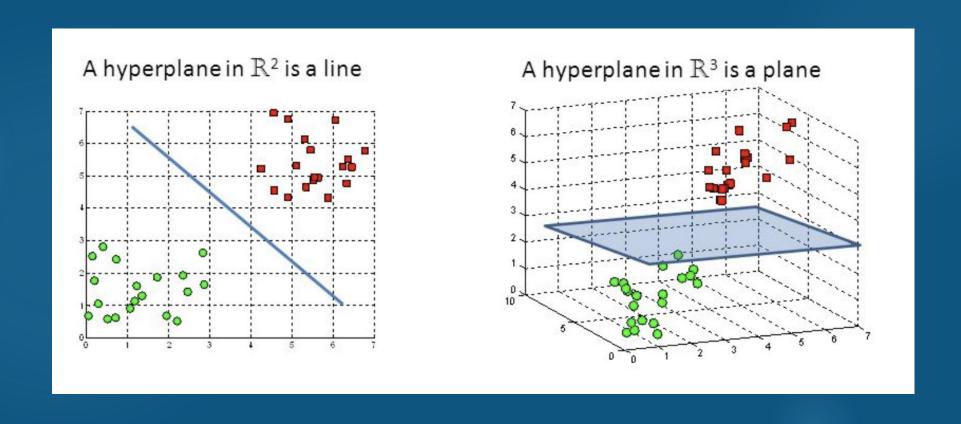
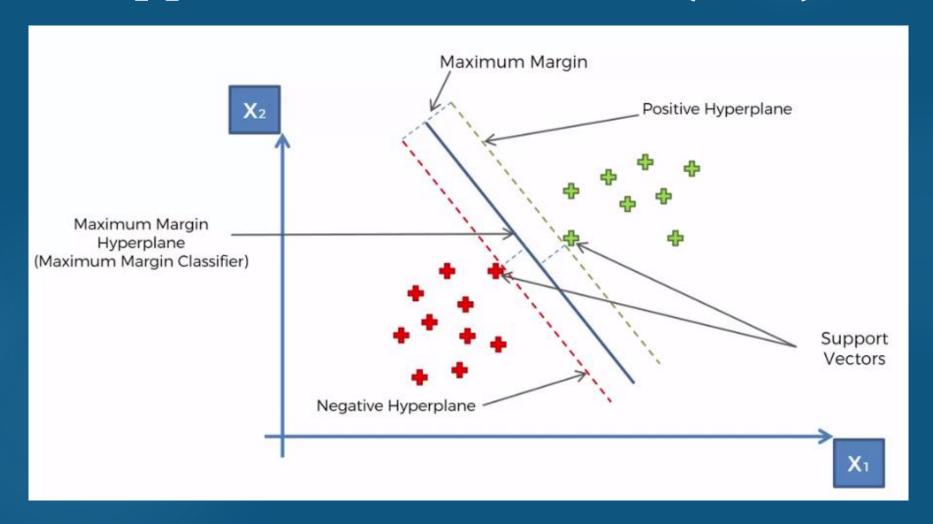
# SUPPORT VECTOR MACHINES (SVM)

## Support vector machines (SVM)

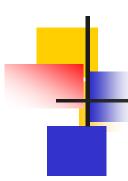
- ▶ **Support Vector Machines (SVM)** are supervised learning models used for classification and regression analysis.
- ▶ SVMs are based on the idea of finding a **hyperplane** that best divides a dataset into two classes



## Support vector machines (SVM)



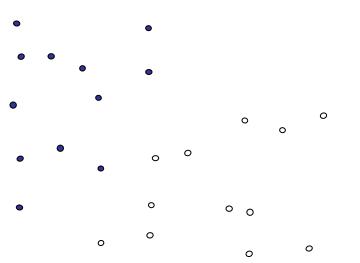
The goal is to choose a hyperplane with the greatest possible margin



**Estimation:** 



- denotes +1
- ° denotes -1



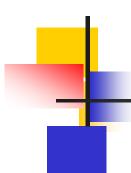
0

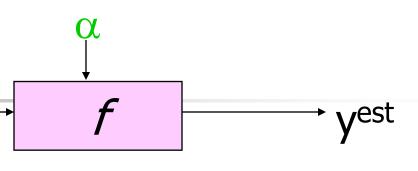
0 0

$$f(x, w, b) = sign(w. x - b)$$

w: weight vector

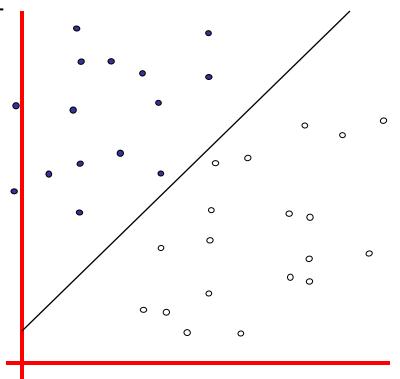
x: data vector

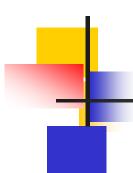


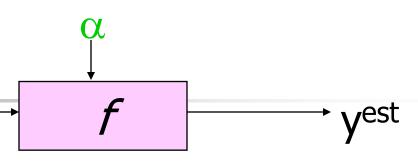


f(x, w, b) = sign(w. x - b)

° denotes -1

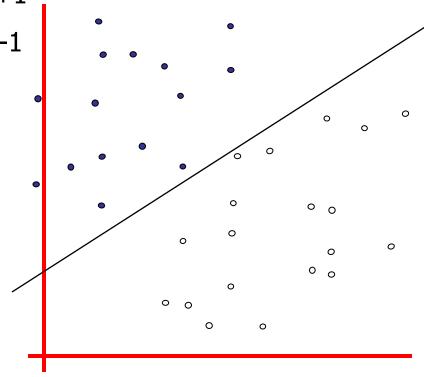


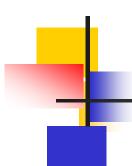


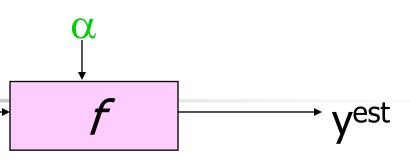


f(x, w, b) = sign(w. x - b)

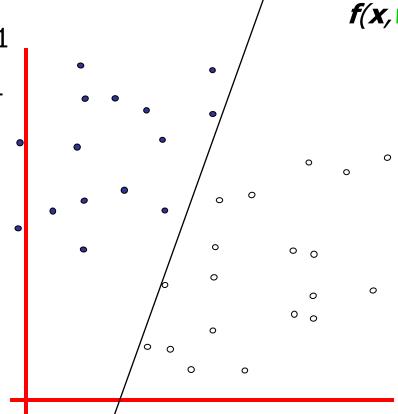
° denotes -1



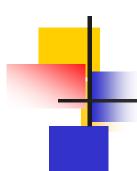


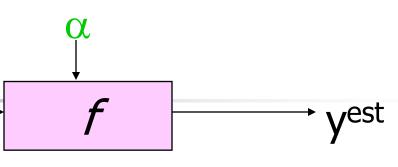


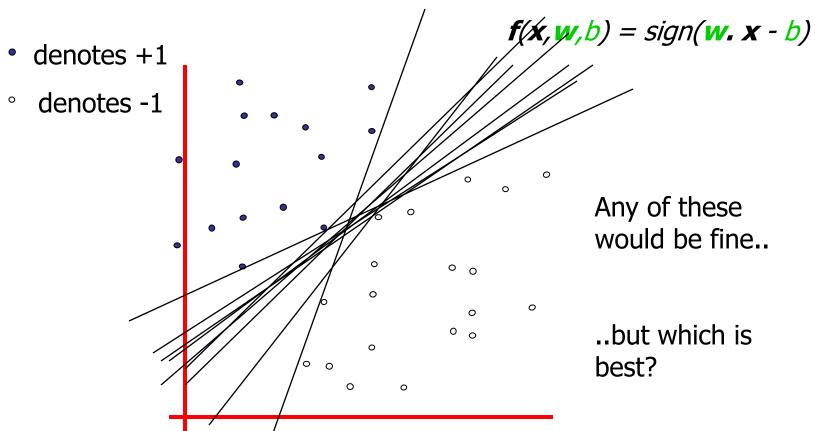
- denotes +1
- ° denotes -1



f(x, w, b) = sign(w. x - b)





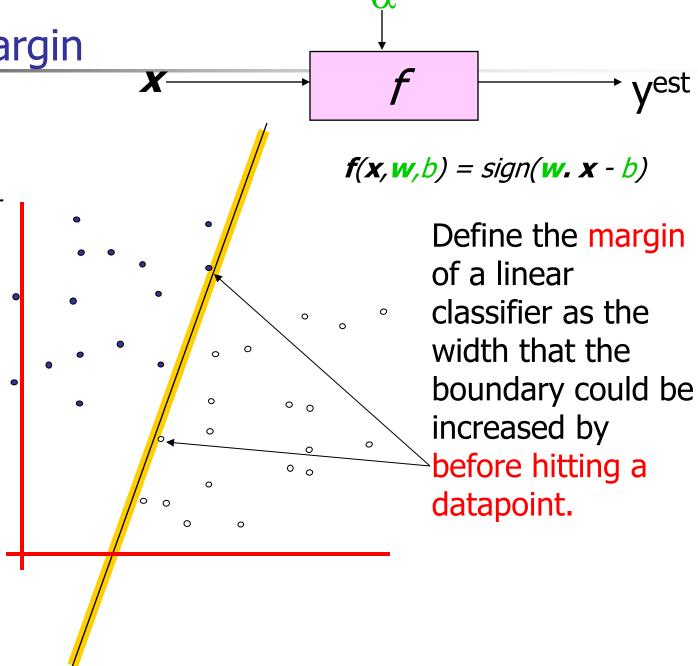


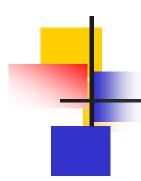


#### Classifier Margin

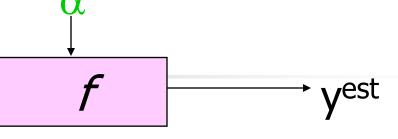
• denotes +1

° denotes -1



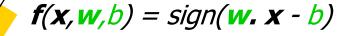


#### Maximum Margin



• denotes +1

° denotes -1



The maximum margin linear classifier is the linear classifier with the, um, maximum margin.

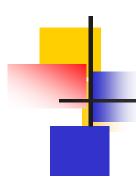
This is the simplest kind of SVM (Called an LSVM)

Linear SVM

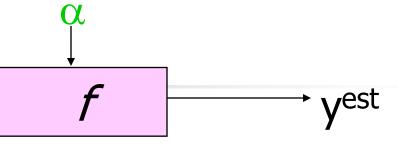
0 0

0 0

0



#### Maximum Margin

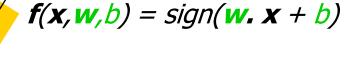


• denotes +1

° denotes -1

#### **Support Vectors**

are those datapoints that the margin pushes up against



The maximum margin linear classifier is the linear classifier with the, um, maximum margin.

This is the simplest kind of SVM (Called an LSVM)

Linear SVM

0 0

0 0

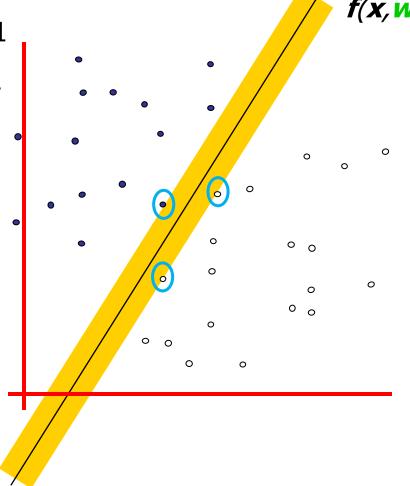
## Why Maximum Margin?



° denotes -1

#### **Support Vectors**

are those datapoints that the margin pushes up against

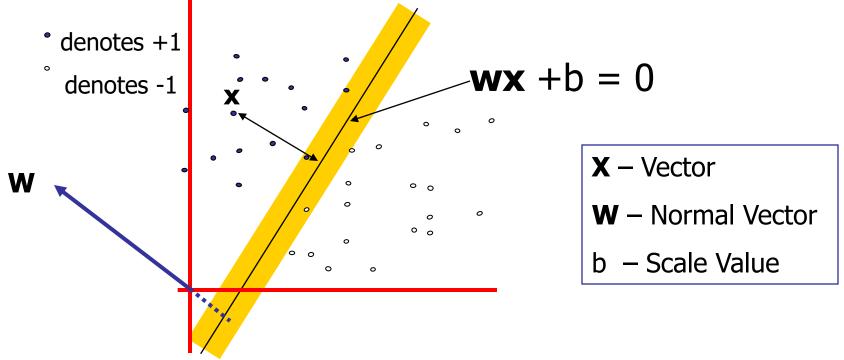


 $f(x, \mathbf{w}, b) = sign(\mathbf{w}, \mathbf{x} - b)$ 

The maximum margin linear classifier is the linear classifier with the maximum margin.

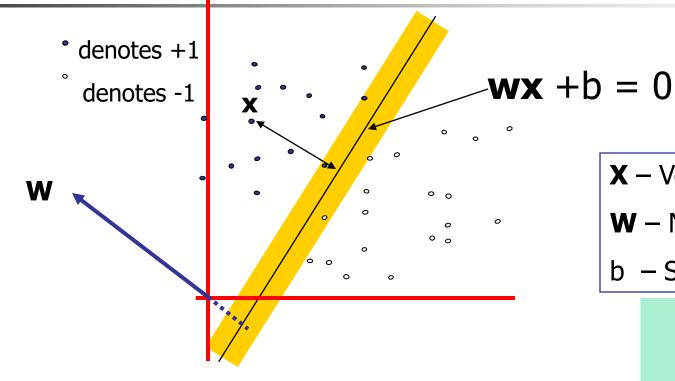
This is the simplest kind of SVM (Called an LSVM)

#### How to calculate the distance from a point to a line?



- http://mathworld.wolfram.com/Point-LineDistance2-Dimensional.html
- In our case,  $w_1 * x_1 + w_2 * x_2 + b = 0$ ,
- thus,  $\mathbf{w} = (w_1, w_2), \mathbf{x} = (x_1, x_2)$

### Estimate the Margin



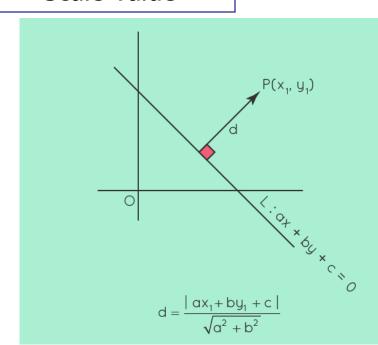
What is the distance expression for a point  $\mathbf{x}$  to a line  $\mathbf{w}\mathbf{x}+\mathbf{b}=0$ ?

$$d(\mathbf{x}) = \frac{\left|\mathbf{x} \cdot \mathbf{w} + b\right|}{\sqrt{\left\|\mathbf{w}\right\|_{2}^{2}}} = \frac{\left|\mathbf{x} \cdot \mathbf{w} + b\right|}{\sqrt{\sum_{i=1}^{d} w_{i}^{2}}}$$

**X** – Vector

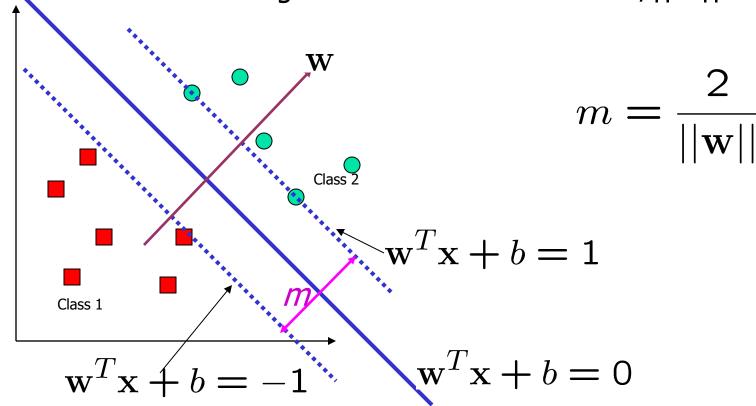
**W** – Normal Vector

b - Scale Value



## Large-margin Decision Boundary

- The decision boundary should be as far away from the data of both classes as possible
  - We should maximize the margin, *m*
  - Distance between the origin and the line  $\mathbf{w}^{t}\mathbf{x}=-\mathbf{b}$  is  $\mathbf{b}/||\mathbf{w}||$



# Finding the Decision Boundary

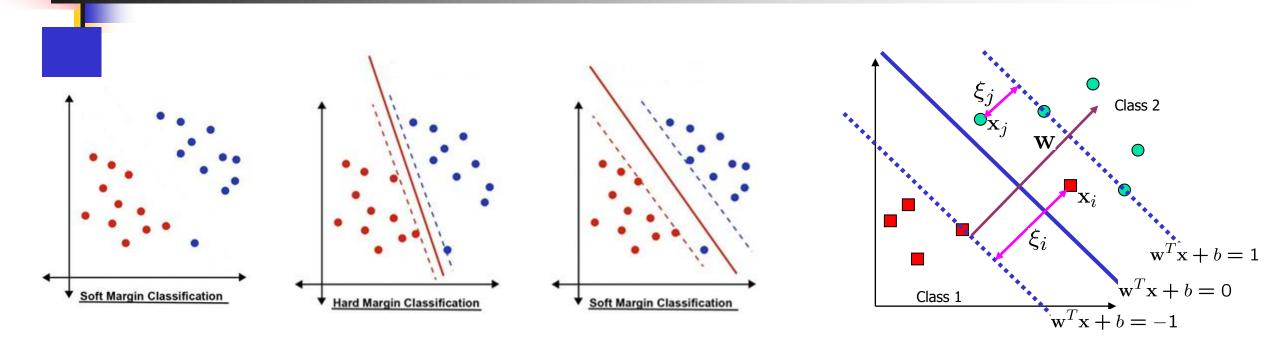
- Let  $\{x_1, ..., x_n\}$  be our data set and let  $y_i \in \{1,-1\}$  be the class label of  $x_i$
- The decision boundary should classify all points correctly ⇒

$$y_i(\mathbf{w}^T\mathbf{x}_i + b) \ge 1, \quad \forall i$$

- To see this: when y=-1, we wish (wx+b)<1, when y=1, we wish (wx+b)>1. For support vectors, we wish y(wx+b)=1.
- The decision boundary can be found by solving the following constrained optimization problem

Minimize 
$$\frac{1}{2}||\mathbf{w}||^2$$
 subject to  $y_i(\mathbf{w}^T\mathbf{x}_i+b)\geq 1$   $\forall i$ 

## Allowing errors in our solutions



- We allow "error"  $\xi_i$  in classification; it is based on the output of the discriminant function  $\mathbf{w}^\mathsf{T}\mathbf{x} + \mathbf{b}$
- $\xi_i$  approximates the number of misclassified samples

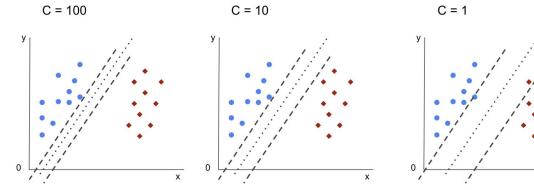
## Soft Margin Hyperplane

• If we minimize  $\sum_i \xi_i$ ,  $\xi_i$  can be computed by

$$\begin{cases} \mathbf{w}^T \mathbf{x}_i + b \ge 1 - \xi_i & y_i = 1 \\ \mathbf{w}^T \mathbf{x}_i + b \le -1 + \xi_i & y_i = -1 \\ \xi_i \ge 0 & \forall i \end{cases}$$

- ξ<sub>i</sub> are "slack variables" in optimization
- Note that  $\xi_i$ =0 if there is no error for  $\mathbf{x}_i$
- $\xi_i$  is an upper bound of the number of errors
- We want to minimize
  - C: tradeoff parameter between error and margin

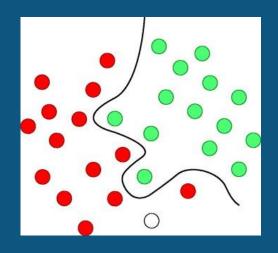
$$\frac{1}{2}||\mathbf{w}||^2 + C\sum_{i=1}^n \xi_i$$

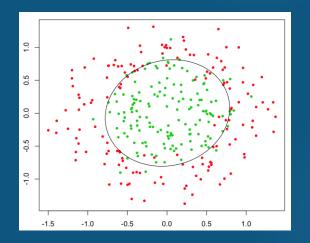


The optimization problem becomes

Minimize 
$$\frac{1}{2}||\mathbf{w}||^2 + C\sum_{i=1}^n \xi_i$$
  
subject to  $y_i(\mathbf{w}^T\mathbf{x}_i + b) \ge 1 - \xi_i, \quad \xi_i \ge 0$ 

## Support Vector machine (SVM)



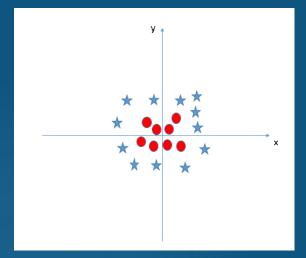


#### Support Vector Machines:

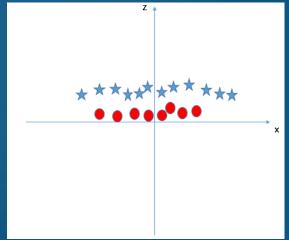
When non-linear boundary is used for the classification purpose, it is called Support Vector Machines (SVMs). It is an extension of the Support Vector Classifier and done by enlarging the feature space in a specific way, using *kernels*.

## Find the Hyperplane

In the scenario below, we can't have linear hyperplane between the two classes, so how does SVM classify these two classes? Till now, we have only looked at the linear hyperplane.



SVM can solve this problem. It solves this problem by introducing additional feature. Here, we will add a new feature  $z=x^2+y^2$ .

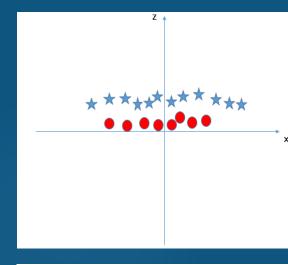


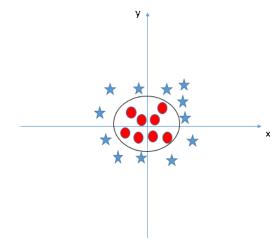
In the original plot, red circles appear close to the origin of x and y axes, leading to lower value of z and star relatively away from the origin result to higher value of z.

#### KERNEL TRICK

Should we need to add this feature manually (inefficient with much computation) to have a hyperplane?

No, SVM has a technique called the kernel trick. These are functions which takes low dimensional input space and transform it efficiently (trick) to a higher dimensional space i.e. it converts non separable problem to separable problem, these functions are called kernels.





#### KERNEL FUNCTIONS

Let's look at an example:

$$\mathbf{x}=(x_1,x_2,x_3)^T \ \mathbf{y}=(y_1,y_2,y_3)^T$$

Here x and y are two data points in 3 dimensions. Let's assume that we need to map x and y to 9-dimensional space. We need to do the following calculations to get the final result, which is just a scalar. The computational complexity, in this case, is  $O(n^2)$ .

$$egin{aligned} \phi(\mathbf{x}) &= (x_1^2, x_1 x_2, x_1 x_3, x_2 x_1, x_2^2, x_2 x_3, x_3 x_1, x_3 x_2, x_3^2)^T \ \phi(\mathbf{y}) &= (y_1^2, y_1 y_2, y_1 y_3, y_2 y_1, y_2^2, y_2 y_3, y_3 y_1, y_3 y_2, y_3^2)^T \end{aligned}$$

$$\phi(\mathbf{x})^T\phi(\mathbf{y}) = \sum_{i,j=1}^3 x_i x_j y_i y_j$$

#### KERNEL FUNCTIONS

However, if we use the kernel function, which is denoted as k(x, y), instead of doing the complicated computations in the 9-dimensional space, we reach the same result within the 3-dimensional space by calculating the dot product of x-transpose and y. The computational complexity, in this case, is O(n).

$$egin{align} k(\mathbf{x},\mathbf{y}) &= (\mathbf{x}^T\mathbf{y})^2 \ &= (x_1y_1 + x_2y_2 + x_3y_3)^2 \ &= \sum_{i,j=1}^3 x_i x_j y_i y_j \ \end{array}$$

In essence, what the kernel trick does for us is to offer a more efficient and less expensive way to transform data into higher dimensions. With that saying, the application of the kernel trick is not limited to the SVM algorithm. Any computations involving the dot products (x, y) can utilize the kernel trick.

#### **EXAMPLES**

For example let x and y be defined as x = (x1, x2, x3) and y = (y1, y2, y3).

The mapping to 9 dimensions would be

$$f(x) = (x1x1, x1x2, x1x3, x2x1, x2x2, x2x3, x3x1, x3x2, x3x3)$$

We can define a kernel which would be equivalent to the above equation

$$K(x, y) = dot(x, y) = x.y = (x^{T}y)^{2}$$

Let,

$$x = (1, 2, 3)$$
  
 $y = (4, 5, 6).$ 

Then:

$$f(x) = (1, 2, 3, 2, 4, 6, 3, 6, 9)$$
  
 $f(y) = (16, 20, 24, 20, 25, 30, 24, 30, 36)$ 

Calculating  $\langle f(x), f(y) \rangle$ , gives us

$$16 + 40 + 72 + 40 + 100 + 180 + 72 + 180 + 324 = 1024$$

Instead of doing so many calculations, if we apply the kernel instead:

$$K(x, y) = (4 + 10 + 18)^{2} = 32^{2} = 1024$$

#### KERNELS

- Linear Kernel:  $K(x, y) = x \cdot y$
- Polynomial Kernel:  $K(x, y) = (x \cdot y + c)^d$
- Radial Basis Function (RBF) Kernel:

$$K(x, y) = \exp(-\gamma ||x - y||^2)$$

Sigmoid Kernel:

$$K(x, y) = \tanh(\alpha x \cdot y + c)$$

#### Polynomial Kernel:

Suppose we have two features X1 and X2 and output variable as Y, so using polynomial kernel we can write it as:

$$X1^{T}.X2 = \begin{bmatrix} X1 \\ X2 \end{bmatrix}.[X1 \ X2]$$

$$= \begin{bmatrix} X1^{2} & X1.X2 \\ X1.X2 & X2^{2} \end{bmatrix}$$

So, we basically need to find  $X_1^2$ ,  $X_2^2$  and  $X_1$ .  $X_2$ , and now we can see that 2 dimensions got converted into 5 dimensions.

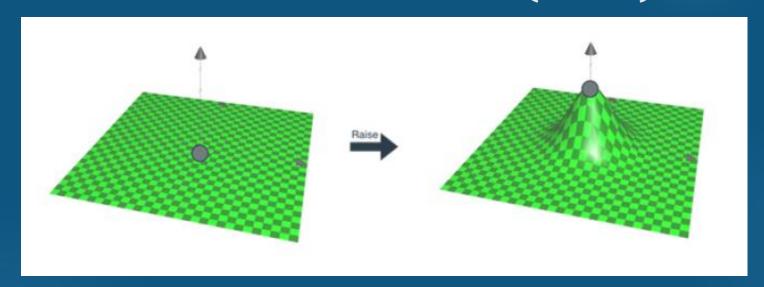
#### Linear Kernel:

Linear kernel is the dot product of the input samples. The dot product of the two points determines the cosine similarity, The higher the value, the more similar the points are.

#### Sigmoid Kernel:

It is just taking your input, mapping them to a value of 0 and 1 so that they can be separated by a simple straight line.

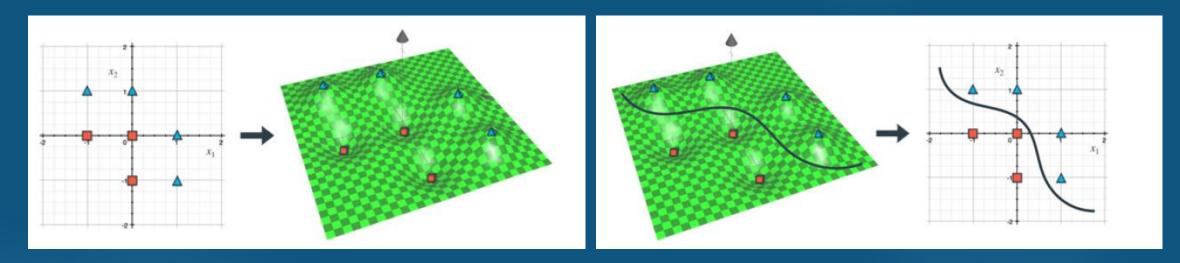
## Radial basis Function (RBF) kernel



Imagine if you had a point in the plane, and the plane was like a blanket. Then you pinch the blanket at that point, and raise it. This is how a *radial basis function* looks like.

We can raise the blanket at any point we like, and that gives us one different radial basis function. The *radial basis function kernel* (also called rbf kernel) is precisely the set of all radial basis functions for every point in the plane.

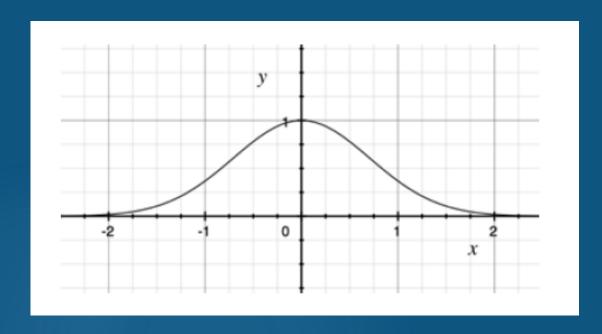
## Radial basis Function (RBF) kernel



Lift the plane at every triangle, and push it down at every square. Then simply draw a plane at height 0, and intersect it with our surface. This is the same as looking at the curve formed by the points at height 0.

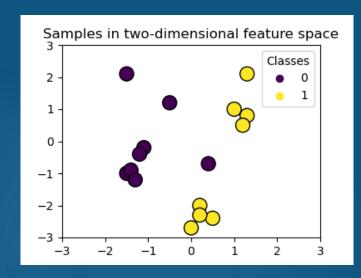
Imagine if there is a landscape with mountains and with the sea. The curve will correspond to the coastline, namely, where the water and the land meet. This gives us the curves (when we project everything back to the plane), and we obtain our desired classifier.

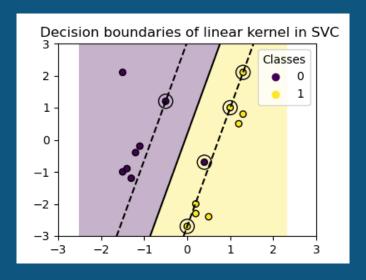
## Radial basis Function (RBF) kernel

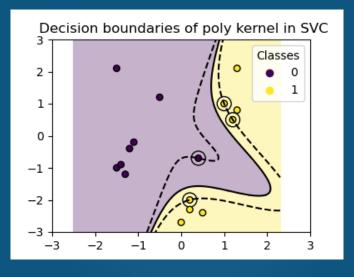


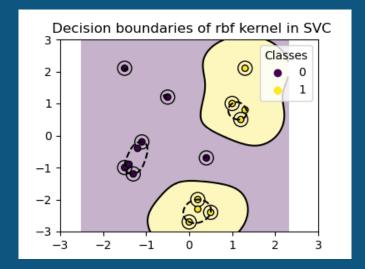
- 1. One of the simplest radial basis function has the formula:  $y = e^{-x^2}$
- 2. Notice that this bump happens at 0. If we wanted it at any different point, say p, we simply translate the formula to  $y = e^{-(x-p)^2}$
- 3. Thus, if we want to obtain the radial basis function centered at the point 5:  $y = e^{-(x-5)^2}$

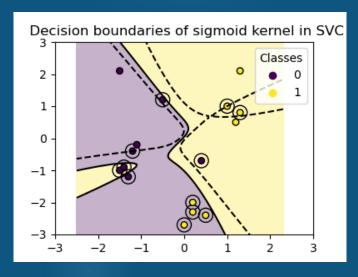
#### DIFFERENT KERNELS











# Thank You