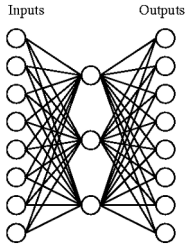


# Artificial Neural Networks (ANNs)

Biologically inspired computational model:

- (1) Simple computational units (neurons).
- (2) Highly interconnected - connectionist view
- (3) Vast parallel computation, consider:
  - Human brain has  $\sim 10^{11}$  neurons
  - Slow computational units, switching time  $\sim 10^{-3}$  sec (compared to the computer  $> 10^{-10}$  sec)
  - Yet, you can recognize a face in  $\sim 10^{-1}$  sec
  - This implies only about 100 sequential, computational neuron steps - this seems too low for something as complicated as recognizing a face
  - Parallel processing

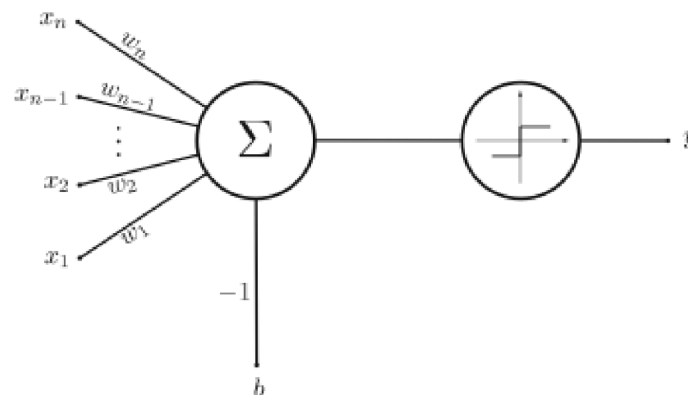
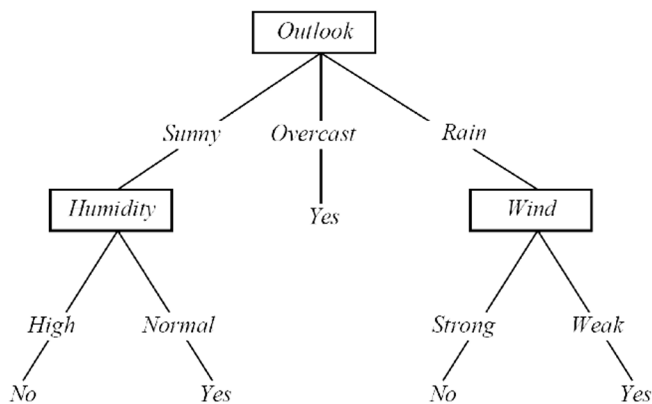


ANNs are naturally parallel - each neuron is a self-contained computational unit that depends only on its inputs.

# Learning

We have seen machine learning with different representations:

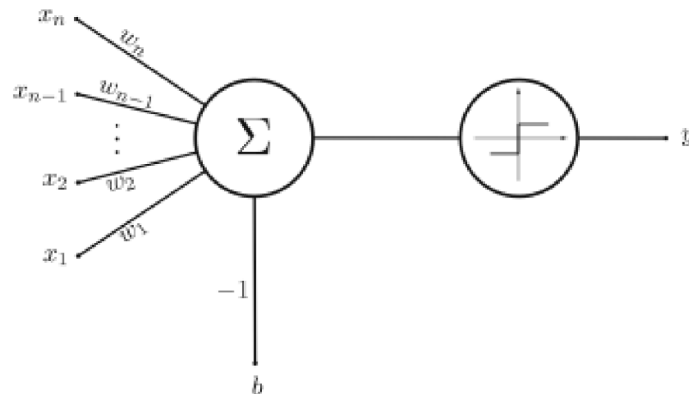
- (1) Decision trees -- symbolic representation of various decision rules -- “disjunction of conjunctions”
- (2) Perceptron -- learning of weights that represent a linear decision surface classifying a set of objects into two groups



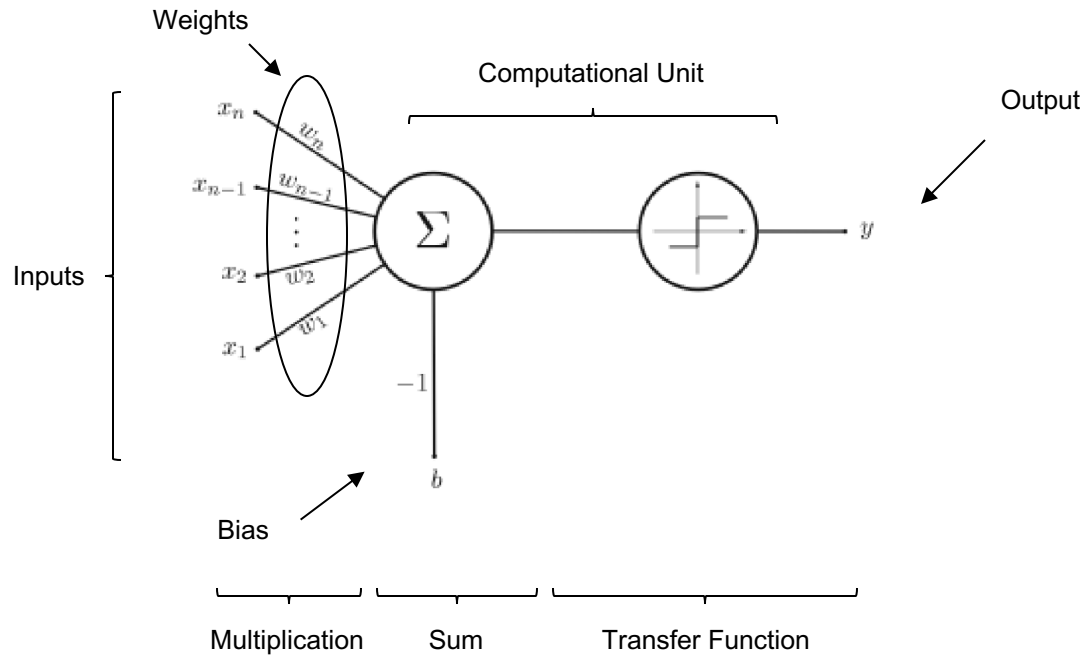
Different representations give rise to different hypothesis or model spaces. Machine learning algorithms search these model spaces for the best fitting model.

# The Perceptron

- A simple, single layered neural “network” - only has a single neuron.
- However, even this simple neural network is already powerful enough to perform (linear) classification tasks.



# The Architecture



Transfer Function:

$$\text{sgn}(k) = \begin{cases} +1 & \text{if } k \geq 0 \\ -1 & \text{otherwise} \end{cases}$$

Perceptron Computation: 
$$y = \text{sgn}\left(b + \sum_{i=1}^m w_i x_i\right)$$

Note:  $y \in \{+1, -1\}$

Binary Classification

# Computation

A perceptron computes the value,

$$y = \text{sgn}\left(b + \sum_{i=1}^m w_i x_i\right)$$

Ignoring the activation function  $\text{sgn}$  and setting  $m = 1$ , we obtain,

$$y' = b + w_1 x_1$$

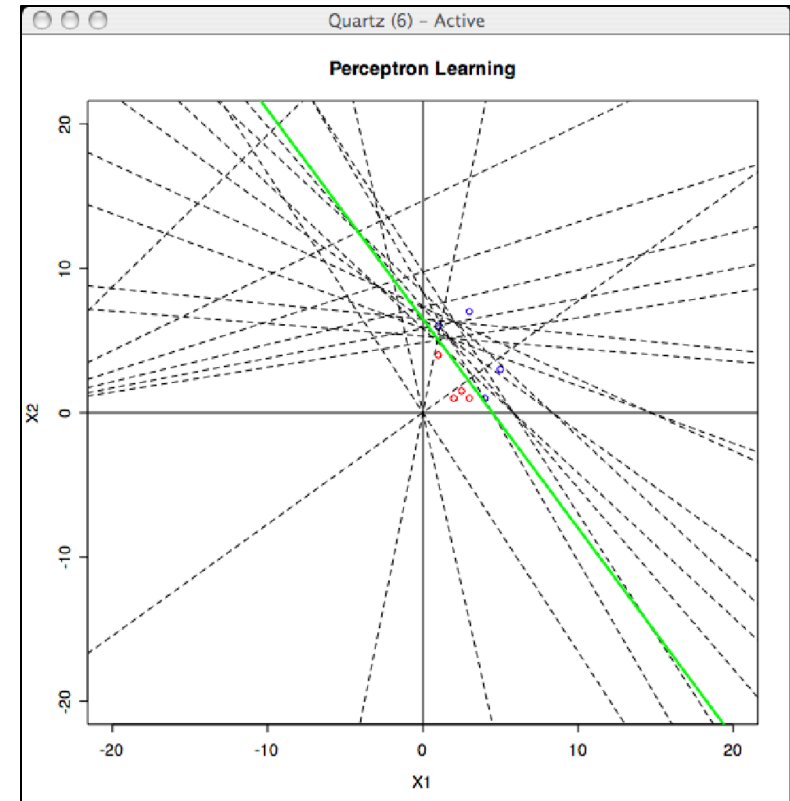
But this is the equation of a line with slope  $w$  and offset  $b$ .

Observation: For the general case the perceptron computes a hyperplane in order to accomplish its classification task,

$$y' = b + \sum_{i=1}^m w_i x_i = b + \vec{w} \cdot \vec{x}$$

# Perceptron Learning Revisited

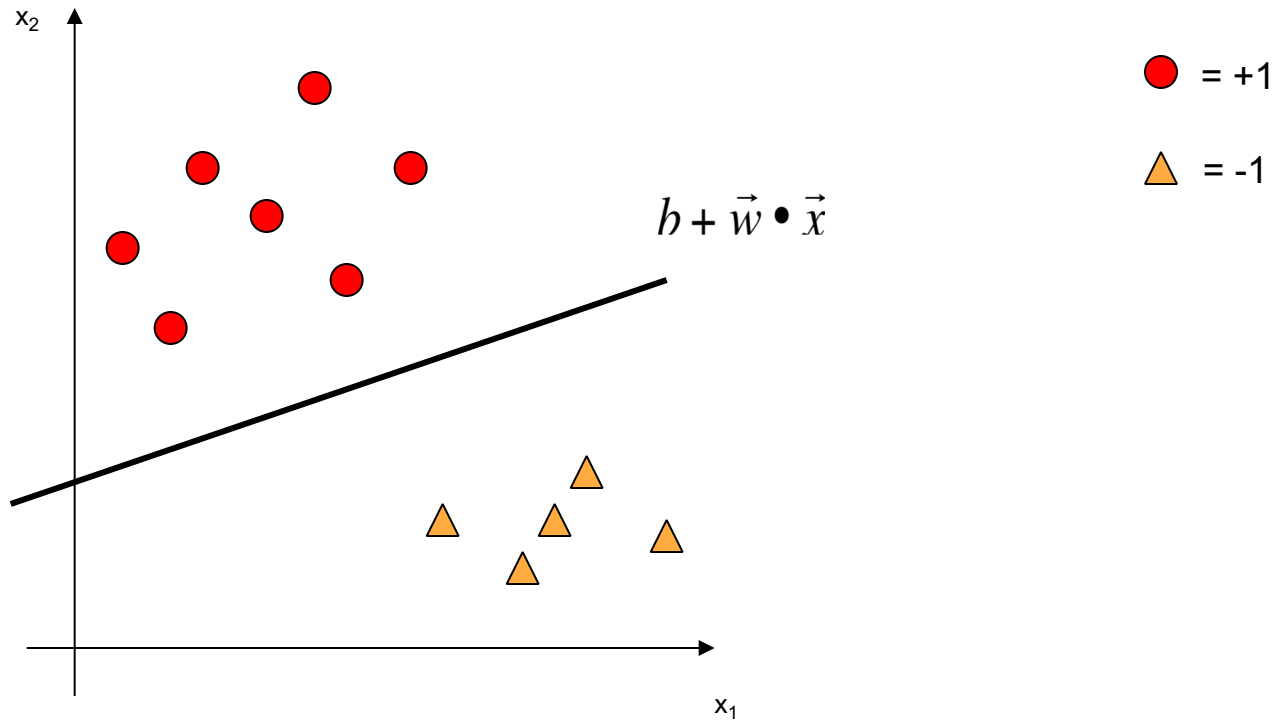
```
Initialize  $\bar{w}$  and  $b$  to random values.  
repeat  
  for each  $(\bar{x}_i, y_i) \in D$  do  
    if  $\hat{f}(\bar{x}_i) \neq y_i$  then  
      Update  $\bar{w}$  and  $b$  incrementally.  
    end if  
  end for  
until  $D$  is perfectly classified.  
return  $\bar{w}$  and  $b$ 
```



Constructs a line (hyperplane) as a classifier

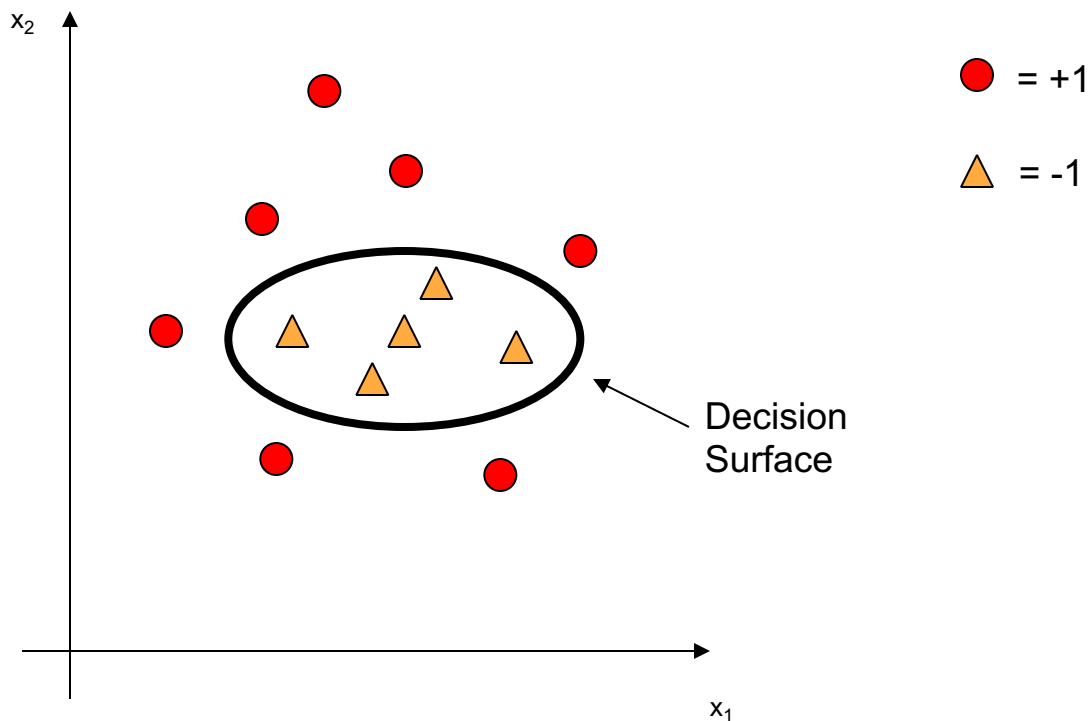
$$h + \vec{w} \cdot \vec{x}$$

# Classification



In order for the hyperplane to become a classifier we need to find  $b$  and  $w \Rightarrow$  learning!

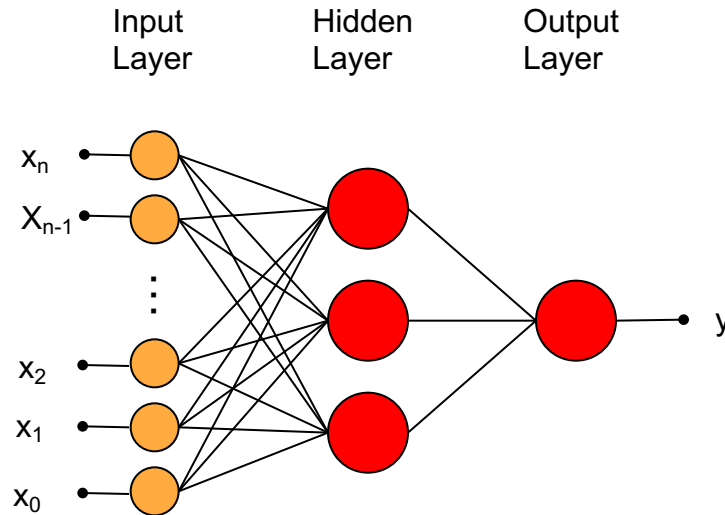
# What About Non-Linearity?



Can we learn this decision surface? ...Yes! *Multi-Layer Perceptrons*

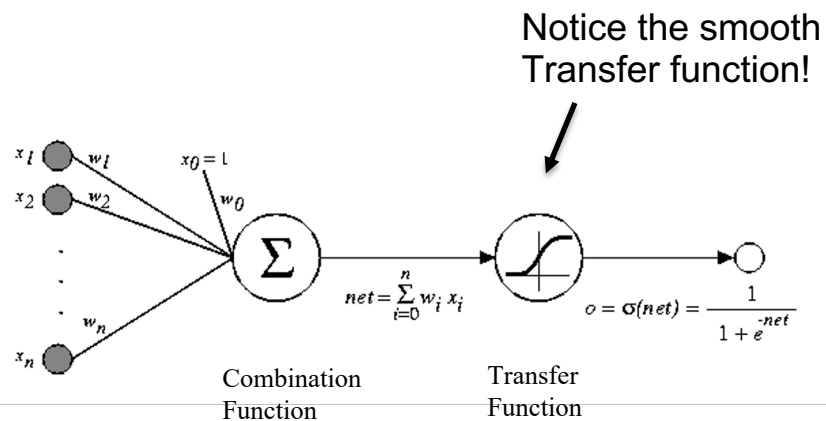


# Multi-Layer Perceptrons (ANNs)



○ ≡ Linear/Input Unit

● ≡

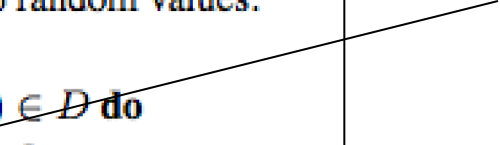


# How do we train?

Perceptron was easy:

```
Initialize  $\bar{w}$  and  $b$  to random values.  
repeat  
  for each  $(\bar{x}_i, y_i) \in D$  do  
    if  $\hat{f}(\bar{x}_i) \neq y_i$  then  
      Update  $\bar{w}$  and  $b$  incrementally.  
    end if  
  end for  
until  $D$  is perfectly classified.  
return  $\bar{w}$  and  $b$ 
```

error



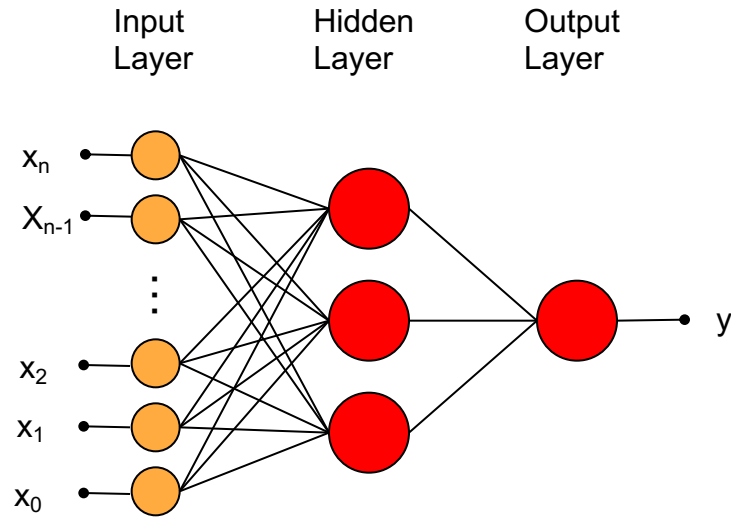
Every time we found an error of the predicted value  $f(x_i)$  compared to the label in the training set  $y_i$ , we update  $w$  and  $b$ .

Not so easy in multi-layer neural networks – the error can occur deep in the network!

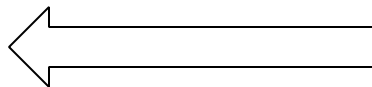
# Artificial Neural Networks

## Feed-forward with Backpropagation

We have to be a bit smarter in the case of ANNs: compute the signal (feed forward) and then use the error at the output to update all the weights by propagating the error back through the network.

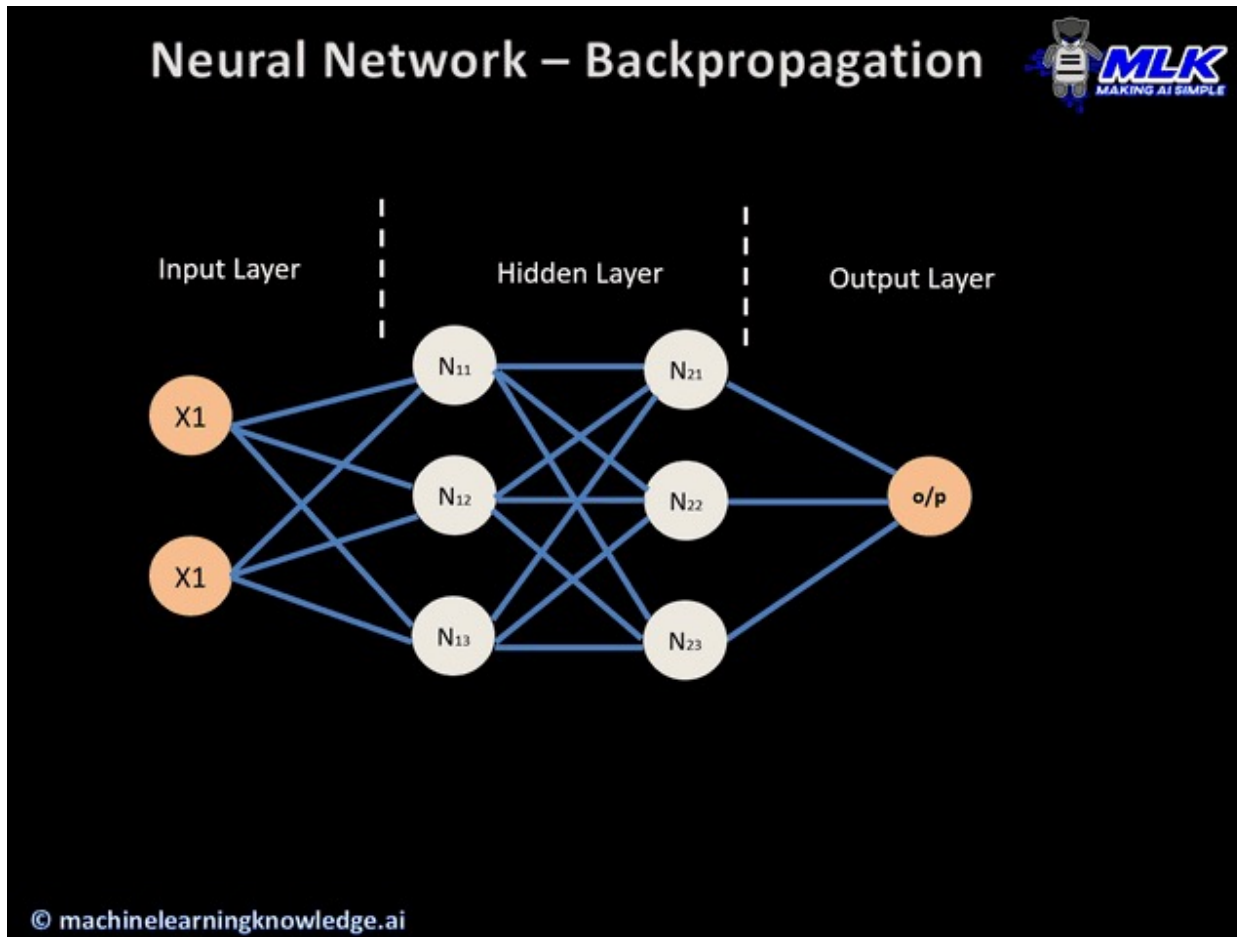


Signal Feed-forward



Error  
Backpropagation

# Back Propagation Training



# Backpropagation

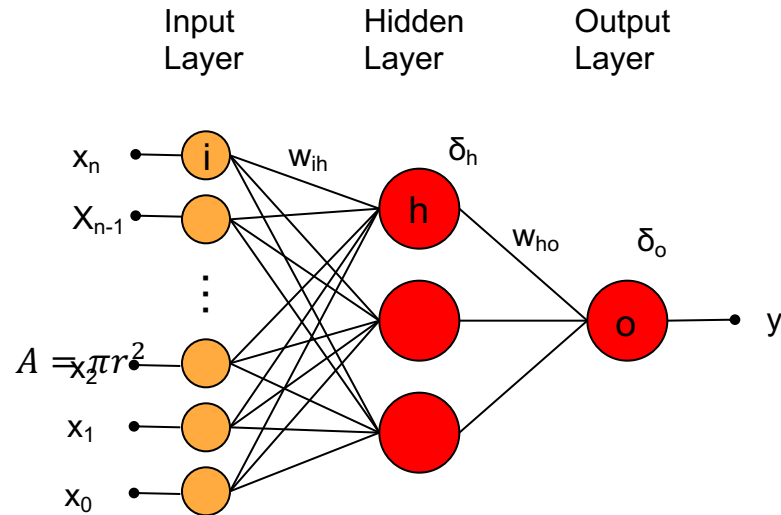
$$E = (y' - y)^2 \quad (\text{output error})$$

$$\delta_o = y(1 - y)E \quad (\text{output node error})$$

$$w_{ho} \leftarrow w_{ho} + \alpha \delta_o \quad (\text{weight update})$$

$$\delta_h = y(1 - y)w_{ho}\delta_o \quad (\text{hidden node error})$$

$$w_{ih} \leftarrow w_{ih} + \alpha \delta_h \quad (\text{weight update})$$



This only works because

$$\delta_o = y(1 - y)E = \frac{\partial E}{\partial w \cdot x} = \frac{\partial (y' - y)^2}{\partial w \cdot x} = 2(y' - y) \left( \frac{\partial y'}{\partial w \cdot x} - \frac{\partial y}{\partial w \cdot x} \right)$$

and the output  $y$  is differentiable because the transfer function is differentiable. Also note, everything is based on the *rate of change* of the error...we are searching in the direction where the rate of change will minimize the output error.

For this to work transfer Function has to be smooth!!

# Backpropagation Algorithm

Note: this algorithm is for a NN with a single output node  $o$  and a single hidden layer. It can easily be generalized.

Initialize the weights in the network (often randomly)

Do

For each example  $e$  in the training set

// forward pass

$y$  = compute neural net output

$y'$  = label for  $e$  from training data

Calculate error  $E = (y' - y)^2$  at the output units

// backward pass

Compute error  $\delta_o$  for weights from a hidden node  $h$  to the output node  $o$  using  $E$

Compute error  $\delta_h$  for weights from an input node  $i$  to hidden node  $h$  using  $\delta_o$

Update the weights in the network

Until all examples classified correctly or stopping criterion satisfied

Return the network

# Neural Network Learning

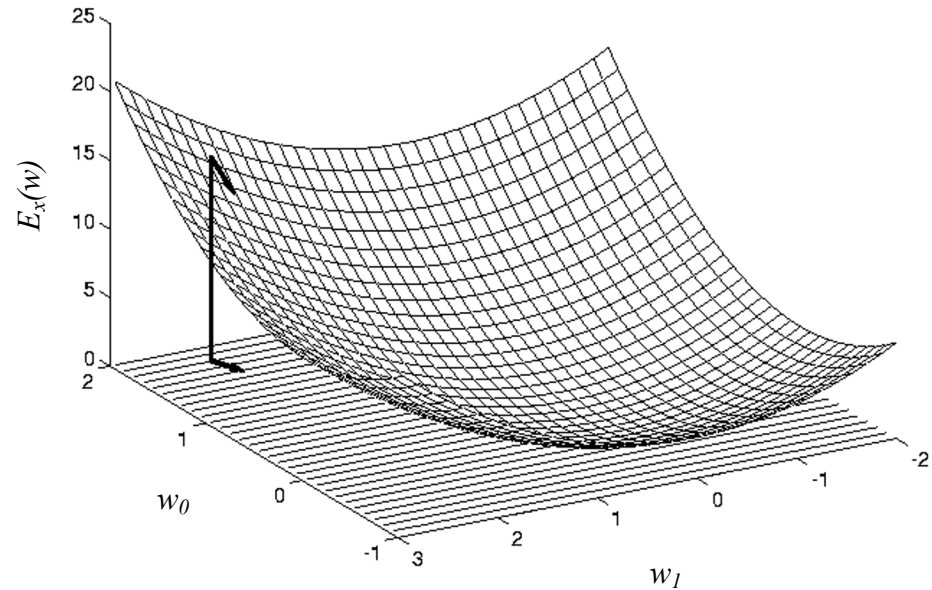
- Define the network error in terms of weights  $w$  as

$$E_x(w) = (y' - y)^2$$

for some training instance  $x$ .

- Use the gradient (slope) of the error surface to guide the search towards appropriate weights:

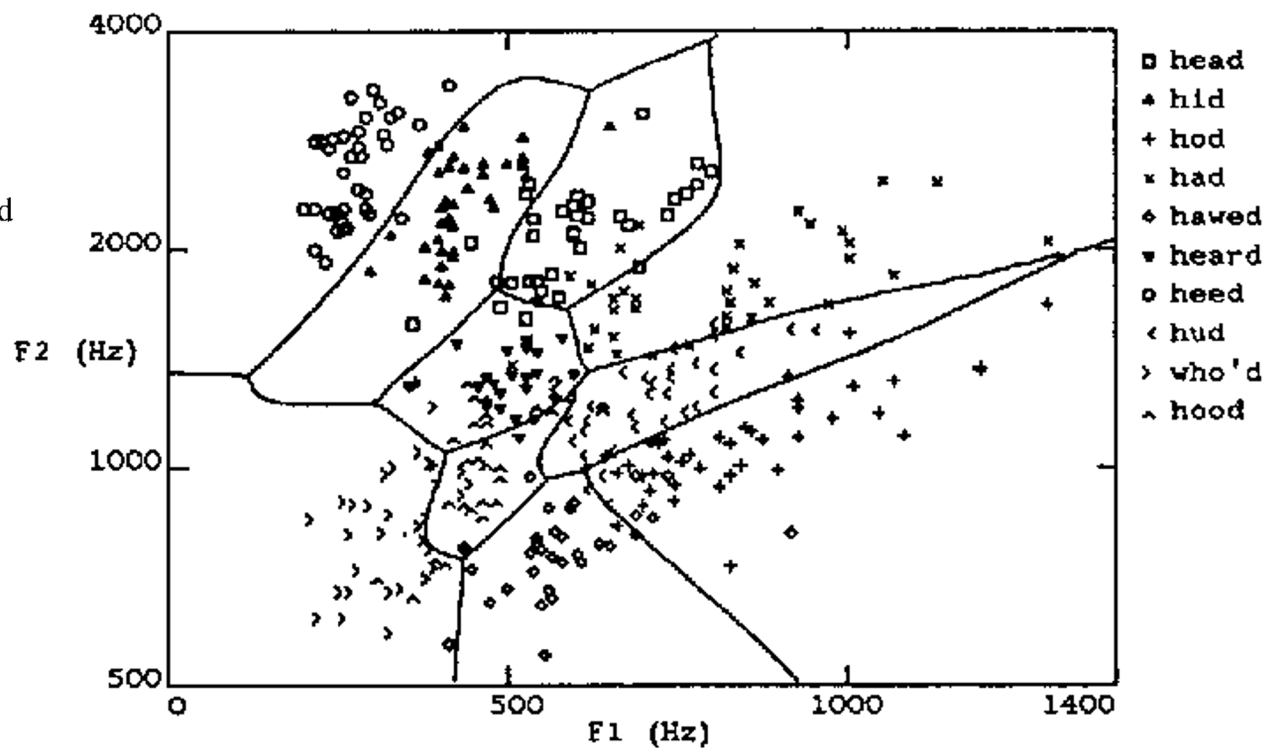
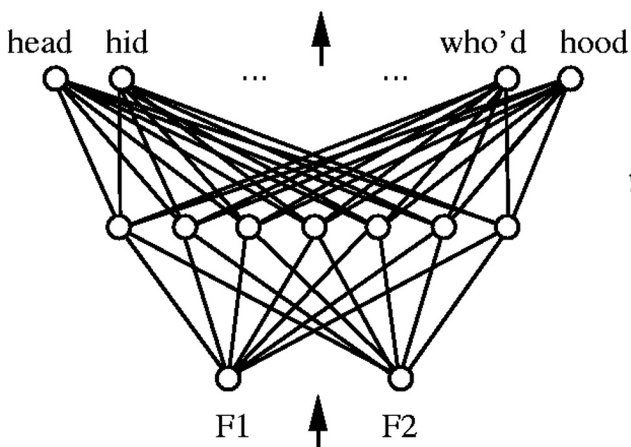
$$\Delta w_k = -\eta \frac{\partial E_x}{\partial w_k}$$



☞ Backpropagation can be understood as a stochastic gradient search on the error surface of the network.

# Representational Power

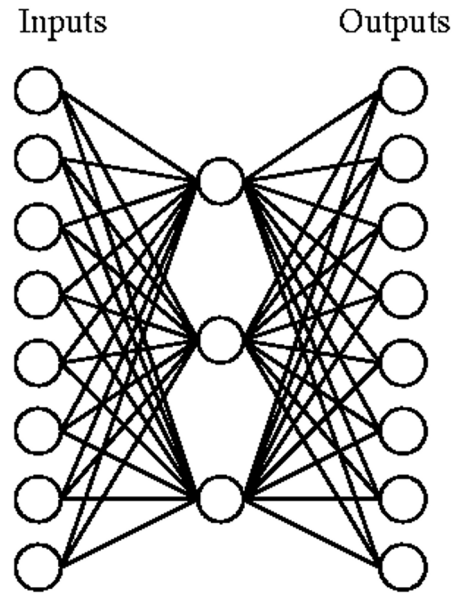
- Every bounded continuous function can be approximated with arbitrarily small error by a network with one hidden layer.
- Any function can be approximated to arbitrary accuracy by a network with two hidden layers.





# Hidden Layer Representations

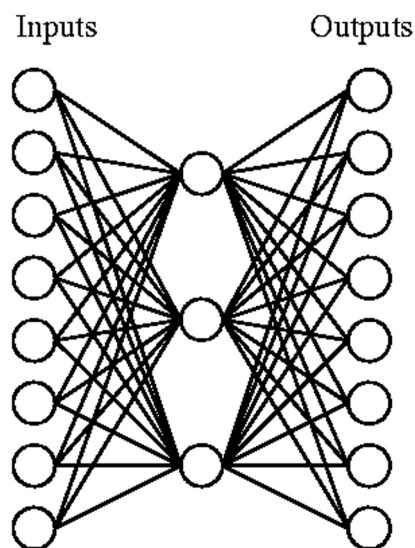
Target Function:



Input	Output
10000000	→ 10000000
01000000	→ 01000000
00100000	→ 00100000
00010000	→ 00010000
00001000	→ 00001000
00000100	→ 00000100
00000010	→ 00000010
00000001	→ 00000001

Can this be learned?

# Hidden Layer Representations



Input	Hidden Values	Output
10000000	→ .89 .04 .08	→ 10000000
01000000	→ .01 .11 .88	→ 01000000
00100000	→ .01 .97 .27	→ 00100000
00010000	→ .99 .97 .71	→ 00010000
00001000	→ .03 .05 .02	→ 00001000
00000100	→ .22 .99 .99	→ 00000100
00000010	→ .80 .01 .98	→ 00000010
00000001	→ .60 .94 .01	→ 00000001

1 0 0
0 0 1
0 1 0
1 1 1
0 0 0
0 1 1
1 0 1
1 1 0

☞ This neural network architecture is sometimes also called autoencoder because of its ability to invent new representations of the input data and is a popular building block in deep-learning.