



## 18. Learning From Examples

# What is Learning?

- An **agent** is learning if it **improves its performance on future tasks after making observations** about the world
- Learning can range from trivial to profound:
  - jotting down a phone number (all of us can do this)
  - infer a new theory of the Universe (only a few such as Einstein can do this)

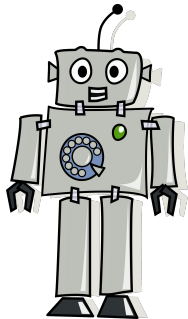
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- We will focus on one class of learning problem
  - “from a collection of input–output pairs, learn a function that predicts the output for new inputs”
- This class of learning task seems restricted but actually has vast applicability
  - Examples?



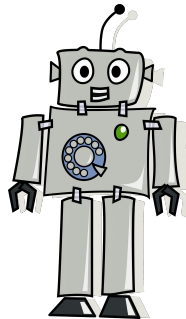
# 18.1 Forms of Learning

- Any component of an agent can be improved by learning from data
- The improvements, and the improvements techniques, depend on four major factors:
  - a. Which component is to be improved
    - workout improves muscles not mind power



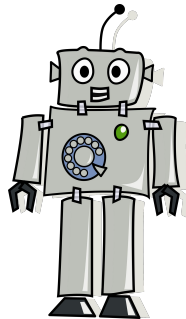
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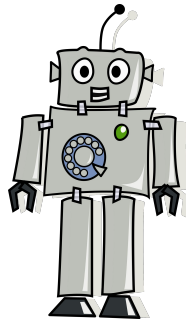
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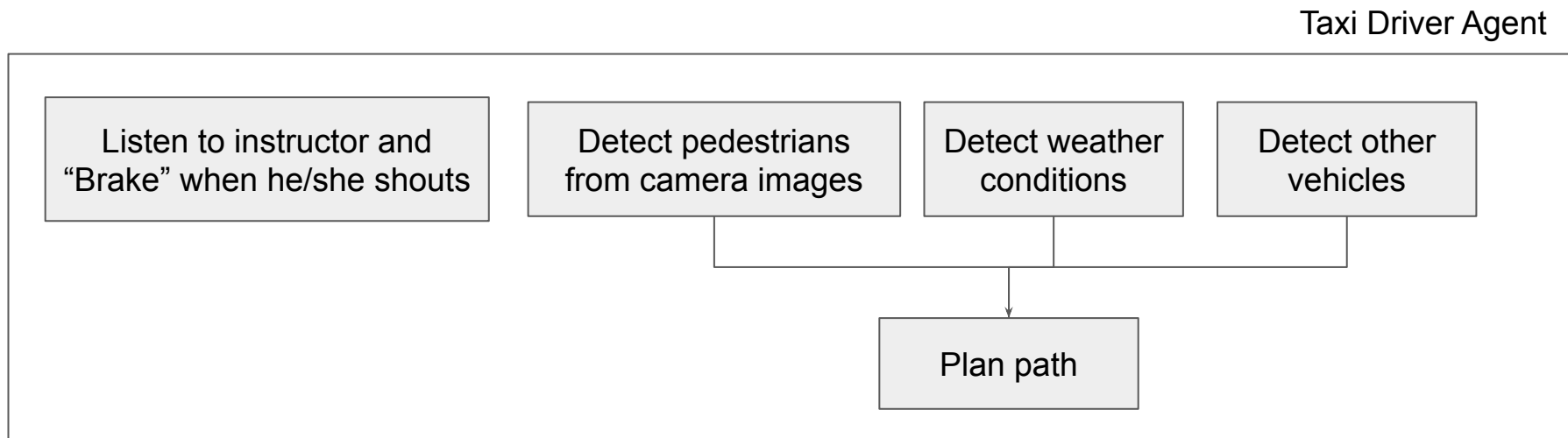
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    - model of human brain?
  - d. What feedback is available to learn from
    - reinforcement learning?



# 18.1 Components to be Learned

- (Although some agents can be end-to-end) An agent may have many components
- Which component do we want to learn at a time?





# 18.1 Representation

- Examples of representations for agent components:
  - Propositional and first-order logical sentences for the components in a logical agent
  - Bayesian networks for the inferential components of a decision-theoretic agent

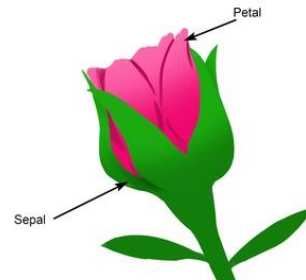
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- The Iris Flower Dataset (1936)
  - The data set consists of 50 samples from each of three species of Iris (Iris setosa, Iris virginica and Iris versicolor)
  - Four features were measured from each sample: the length and the width of the sepals and petals, in centimeters
  - <https://www.kaggle.com/uciml/iris>

	Sepal.Length	Sepal.Width	Petal.Length	Petal.Width	Species
1	5.1	3.5	1.4	0.2	setosa
2	4.9	3.0	1.4	0.2	setosa
3	4.7	3.2	1.3	0.2	setosa
4	4.6	3.1	1.5	0.2	setosa
5	5.0	3.6	1.4	0.2	setosa
6	5.4	3.9	1.7	0.4	setosa
7	4.6	3.4	1.4	0.3	setosa
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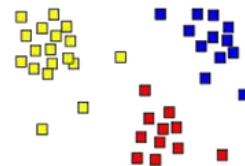


# 18.1 Feedback to Learn From

There are **three** types of feedback that determine the **three main types of learning**:

1. **Unsupervised learning** - the agent learns patterns in the input even though no explicit feedback is supplied

The most common unsupervised learning task is clustering: detecting potentially useful clusters of input examples. For example, a taxi agent might gradually develop a concept of “good traffic days” and “bad traffic days” without ever being given labeled examples of each by a teacher



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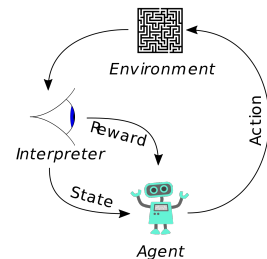
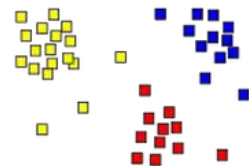
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2. **Reinforcement learning** - the agent learns from a series of reinforcements - rewards or punishments

For example, the lack of a tip at the end of the journey gives the taxi agent an indication that it did something wrong. The two points for a win at the end of a chess game tells the agent it did something right. It is up to the agent to decide which of the actions prior to the reinforcement were most responsible for it.



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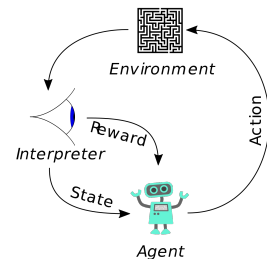
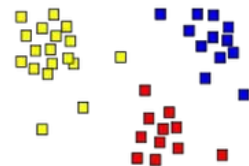
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3. **Supervised learning** - the agent observes some example input–output pairs and learns a function that maps from input to output

The iris dataset example



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- Example:
  - Imagine that you are trying to build a system to guess a person's age from a photo
  - **Supervised Learning:** You gather some labeled examples by snapping pictures of people and asking their age
  - **Reality:** Some of the people lied about their age
  - It's not just that there is random noise in the data; rather the inaccuracies are systematic, and to uncover them is an unsupervised learning problem involving images, self-reported ages, and true (unknown) ages. Thus, both noise and lack of labels create a continuum between supervised and unsupervised learning.

# Unsupervised, Supervised, and Reinforcement Learning



Match the following:

Unsupervised Learning

We have a dataset but there is no target to be predicted. Rather, we want to learn a model that might have generated that set.

Supervised Learning

You teach your kid about different kinds of fruits that are available in world by showing the image of each fruit(X) and its name (Y).

Reinforcement Learning

You ask your child to put apples into different buckets based on size or color.

Unsupervised Learning

This is a setting where we have a sequential decision problem. Making a decision now influences what decisions we can make in the future. A reward function is provided that tells us how “good” certain states are.

Supervised Learning

We have a data set that includes the target values (the values we wish to predict). We try to learn a function that correctly predict the target values from the other features, which can then be used to make predictions about other examples.

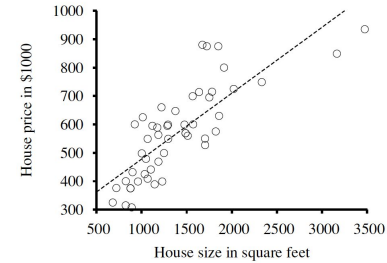
Reinforcement Learning

You give apples to your kid in the morning only after brushing the teeth.

<https://www.quora.com/What-is-the-difference-between-supervised-unsupervised-reinforcement-and-deep-learning>

# 18.6.1 Univariate Linear Regression

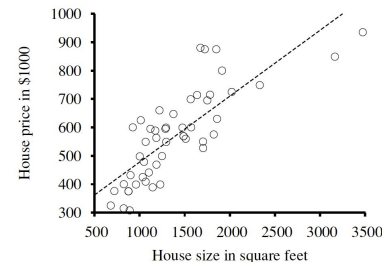
- A univariate linear function (a straight line) with input  $x$  and output  $y$  has the form
  - $y = w_1 x + w_0$ , where  $w_0$  and  $w_1$  are real-valued coefficients to be learned
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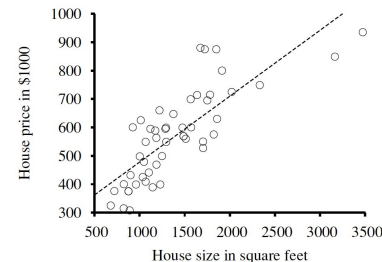
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  - A training set of  $n$  points in the  $x, y$  plane, each point representing the size in square feet and the price of a house offered for sale

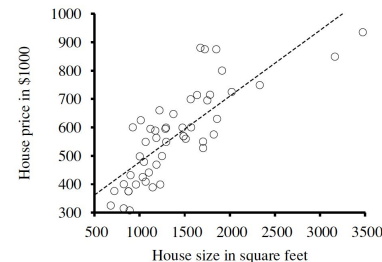


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- The task of finding the  $h_w$  that best fits these data is called **linear regression**
  - To fit a line to the data, all we have to do is find the values of the weights  $[w_0, w_1]$  that minimize the empirical loss



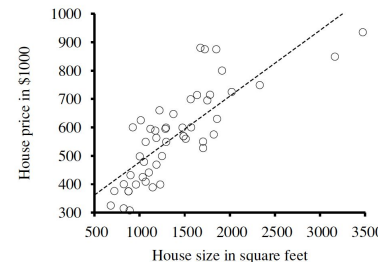
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- It is traditional to use the squared loss function,  $L_2$ , summed over all the training examples:

$$Loss(h_w) = \sum_{j=1}^N (y_j - (w_1 x_j + w_0))^2$$



## 18.6.1 Univariate Linear Regression

- We would like to find  $w^* = \operatorname{argmin}_w \operatorname{Loss}(h_w)$
- The sum  $\operatorname{Loss}(h_w)$  is minimized when its partial derivatives with respect to  $w_0$  and  $w_1$  are zero

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- These equations have a unique solution:

$$w_1 = \frac{N(\sum x_j y_j) - (\sum x_j)(\sum y_j)}{N(\sum x_j^2) - (\sum x_j)^2}; \quad w_0 = (\sum y_j - w_1(\sum x_j))/N$$

- For example,  $w_1 = 0.232$ ,  $w_0 = 246$

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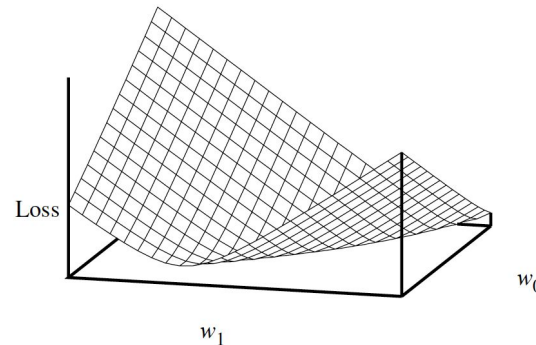
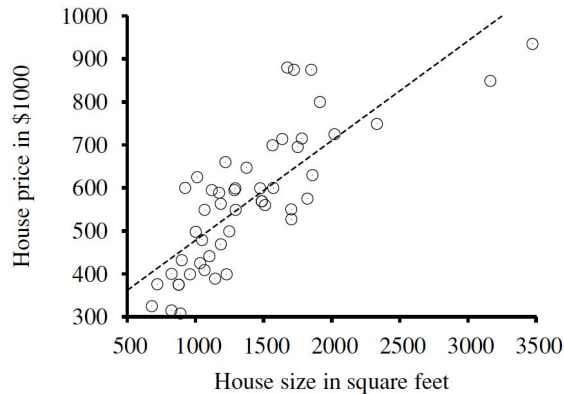
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- For univariate ('only 1 variable as input') linear regression, the weight space defined by  $w_0$  and  $w_1$  is two-dimensional
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## Univariate Linear Regression

- We would like to predict "sepal width" using "sepal length" on the Iris Flower Dataset
- Dataset: [rawdata](#) and [metadata](#)

```
In [0]: from keras.models import Sequential
        from keras.layers import Dense
        import numpy as np
        import matplotlib.pyplot as plt
```

```
In [16]: datapath = 'https://raw.githubusercontent.com/badriadhikari/2019-Fall-AI/master/MODULE-I/iris.data'

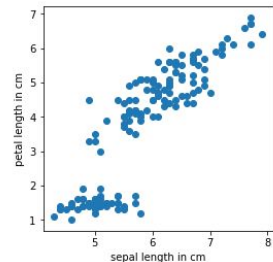
        # Column 1. sepal length in cm (load as col 0)
        # Column 2. petal length in cm (load as col 1)
        # Use loadtxt instead when there are non-numeric values as well
        dataset = np.genfromtxt(datapath, delimiter=",", usecols=(0, 2))

        print('')
        print(dataset.shape)
        print('')
        print(dataset[0:5])

        (150, 2)

        [[5.1 1.4]
         [4.9 1.4]
         [4.7 1.3]
         [4.6 1.5]
         [5.  1.4]]
```

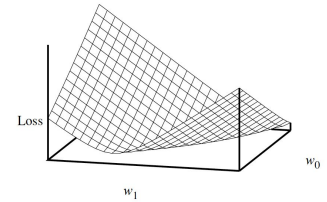
```
In [17]: plt.figure(figsize=(4,4))
        plt.scatter(dataset[:, 0], dataset[:, 1])
        plt.xlabel('sepal length in cm')
        plt.ylabel('petal length in cm')
        plt.show()
```



# 18.6.1 Beyond Linear Models (for Regression)

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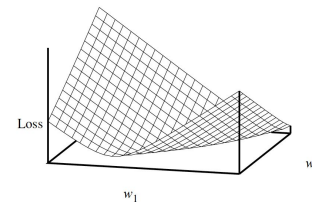
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- Here, we face a general optimization search problem in a continuous weight space
  - Such problems can be addressed using algorithms such as Hill-Climbing algorithm that follows the **gradient** of the function to be optimized.
  - Specifically, because we are trying to minimize loss, we can use **gradient descent**



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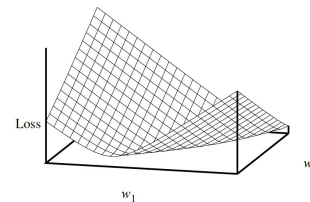
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$\mathbf{w} \leftarrow$  any point in the parameter space  
loop until convergence do

for each  $w_i$  in  $\mathbf{w}$  do

$$w_i \leftarrow w_i - \alpha \frac{\partial}{\partial w_i} Loss(\mathbf{w})$$



The parameter  $\alpha$ , is called the **learning rate** when we are trying to minimize loss in a learning problem. It can be a fixed constant, or it can decay over time as the learning process proceeds.



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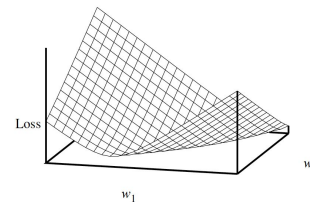
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For univariate regression, this reduces to:  $w_0 \leftarrow w_0 + \alpha \sum_j (y_j - h_{\mathbf{w}}(x_j))$ ;  $w_1 \leftarrow w_1 + \alpha \sum_j (y_j - h_{\mathbf{w}}(x_j)) \times x_j$

# Calculating Weights in Univariate Linear Regression

- Solving the equations with derivative of loss function equal to zero, we saw that we can calculate the two weights

$$\frac{\partial}{\partial w_0} \sum_{j=1}^N (y_j - (w_1 x_j + w_0))^2 = 0 \text{ and } \frac{\partial}{\partial w_1} \sum_{j=1}^N (y_j - (w_1 x_j + w_0))^2 = 0$$
$$\longrightarrow w_1 = \frac{N(\sum x_j y_j) - (\sum x_j)(\sum y_j)}{N(\sum x_j^2) - (\sum x_j)^2}; \quad w_0 = (\sum y_j - w_1(\sum x_j))/N$$

- The gradient descent method also gives us an equation to calculate the weights

$$w_i \leftarrow w_i - \alpha \frac{\partial}{\partial w_i} \text{Loss}(\mathbf{w})$$
$$\longrightarrow w_0 \leftarrow w_0 + \alpha \sum_j (y_j - h_{\mathbf{w}}(x_j)); \quad w_1 \leftarrow w_1 + \alpha \sum_j (y_j - h_{\mathbf{w}}(x_j)) \times x_j$$

- Both methods provide us a technique to calculate the weights
  - What is the difference between the two methods?
  - Which is more general?



# Classroom Discussion & Demo



 Open in Colab

## Linear Regression with Two Input Variables using the Iris Flower Dataset

- We would like to predict "petal width" (column 4 in the original dataset) using sepal width (column 2) and petal length (column 3)
- Dataset: [rawdata](#) and [metadata](#)

```
In [2]: from keras.models import Sequential
        from keras.layers import Dense
        import numpy as np
        import matplotlib.pyplot as plt

        # Column 2. sepal width in cm (load as col 0)
        # Column 3. petal length in cm (load as col 1)
        # Column 4. petal width in cm (load as col 2)
        datapath = 'https://raw.githubusercontent.com/badriadhikari/2019-Fall-AI/master/MODULE-I/iris.dat
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        dataset = np.genfromtxt(datapath, delimiter=",", usecols=(1, 2, 3))

        print('')
        print(dataset.shape)
        print('')
        print(dataset[0:5])

        (150, 3)

        [[3.5 1.4 0.2]
         [3.  1.4 0.2]
         [3.2 1.3 0.2]
         [3.1 1.5 0.2]
         [3.6 1.4 0.2]]
```

```
In [5]: # Q1. Why is shuffling important before splitting?
        np.random.shuffle(dataset)
        print('')
        print(dataset[0:5])
        train = dataset[:100]
        valid = dataset[100:]
        print('')
        print(train.shape)
        print('')
        print(valid.shape)

        [[3.7 1.5 0.2]
         [3.  4.2 1.5]
         [2.8 6.7 2. ]
         [2.4 3.7 1. ]
         [3.  1.4 0.2]]

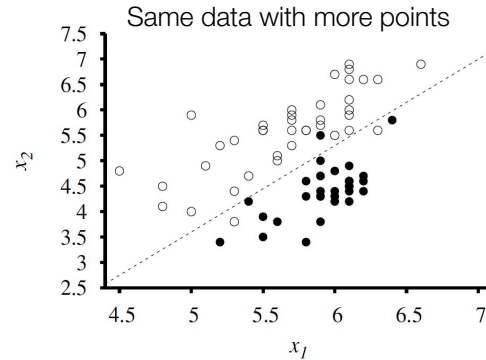
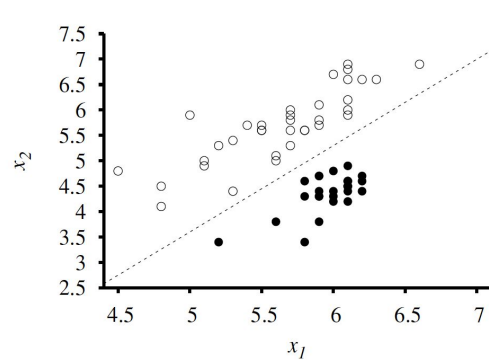
        (100, 3)

        (50, 3)
```

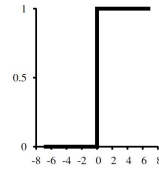
Find a partner!  
Answer the questions!

XIIT PAIR  
SHARE

## 18.6.3 Linear Classifiers with a Hard Threshold

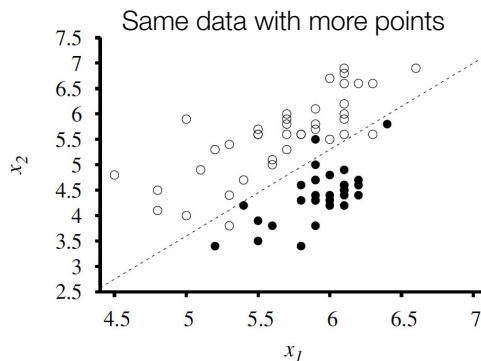
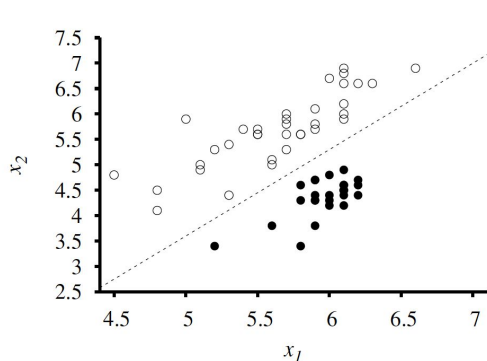


Linear functions can be used to do classification as well (with the help of a threshold).



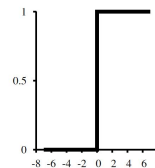
Plot of two seismic data parameters, body wave magnitude  $x_1$  and surface wave magnitude  $x_2$ , for earthquakes (white circles) and nuclear explosions (black circles) occurring between 1982 and 1990 in Asia and the Middle East

# 18.6.3 Linear Classifiers with a Hard Threshold



Plot of two seismic data parameters, body wave magnitude  $x_1$  and surface wave magnitude  $x_2$ , for earthquakes (white circles) and nuclear explosions (black circles) occurring between 1982 and 1990 in Asia and the Middle East

Linear functions can be used to do classification as well (with the help of a threshold).



A **decision boundary** is a line (or a surface, in higher dimensions) that separates the two classes.

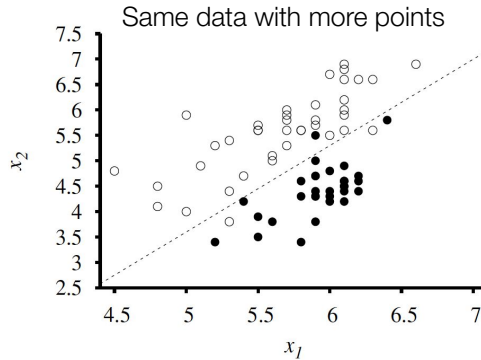
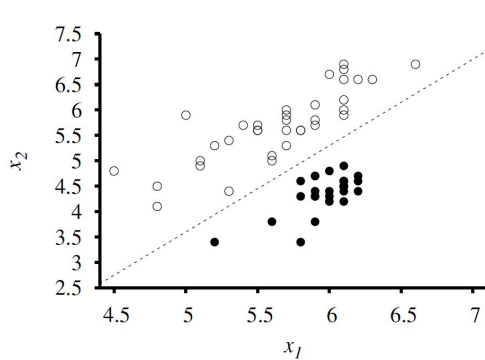
The decision boundary is a straight line in the first figure.

A linear decision boundary is called a **linear separator** and data that admit such a separator are called **linearly separable**.

What are the differences between the two pairs of plots?



# 18.6.3 Linear Classifiers with a Hard Threshold

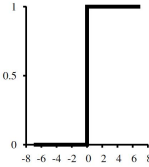


Plot of two seismic data parameters, body wave magnitude  $x_1$  and surface wave magnitude  $x_2$ , for earthquakes (white circles) and nuclear explosions (black circles) occurring between 1982 and 1990 in Asia and the Middle East

$$x_2 = 1.7 x_1 - 4.9 \quad \text{or} \quad -4.9 + 1.7 x_1 - x_2 = 0$$

The explosions, which we want to classify with value 1, are to the right of this line with higher values of  $x_1$  and lower values of  $x_2$ , so they are points for which  $-4.9 + 1.7x_1 - x_2 > 0$ , while earthquakes have  $-4.9 + 1.7x_1 - x_2 < 0$ .

Linear functions can be used to do classification as well (with the help of a threshold).



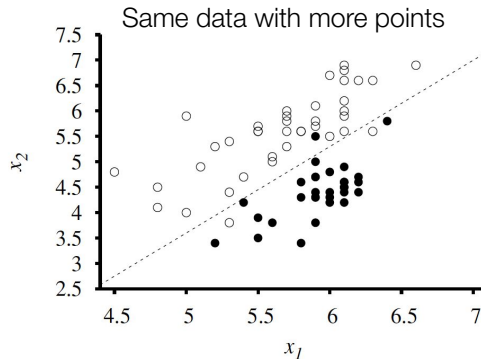
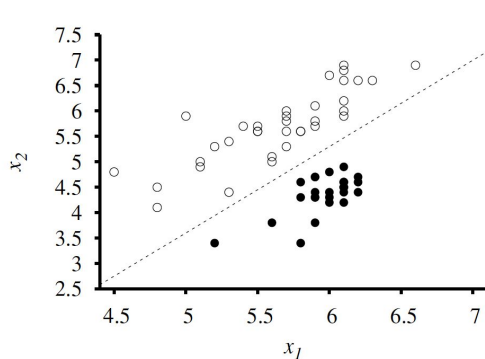
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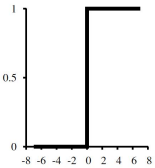
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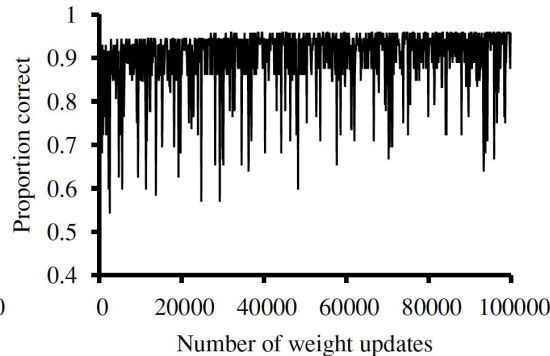
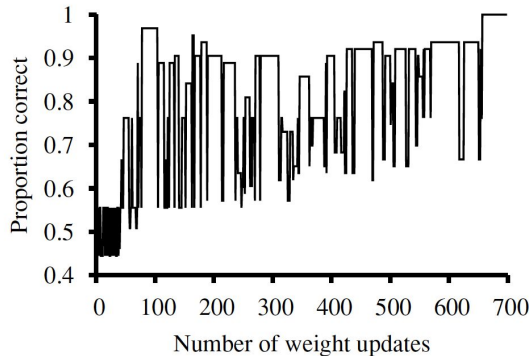
# 18.6.3 Linear Classifiers with a Hard Threshold



Linear functions can be used to do classification as well (with the help of a threshold).



Plot of two seismic data parameters, body wave magnitude  $x_1$  and surface wave magnitude  $x_2$ , for earthquakes (white circles) and nuclear explosions (black circles) occurring between 1982 and 1990 in Asia and the Middle East



Plot of total training-set accuracy vs. number of iterations through the training set

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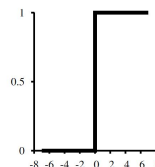
## 18.6.4 Linear Classification with Logistic Regression

Passing the output of a linear function through the threshold function creates a linear classifier, yet the hard nature of the threshold (e.g. 0) causes some **problems**:

**1)** The hypothesis  $h_w(x)$  is a discontinuous function of its inputs and weights

- $h_w(x) = \text{Threshold}(w \cdot x)$ , where  $\text{Threshold}(z) = 1$  if  $z \geq 0$  and 0 otherwise
- Discontinuous functions are not differentiable
- This makes learning with the perceptron rule (gradient descent) very unpredictable

$$w_i \leftarrow w_i - \alpha \frac{\partial}{\partial w_i} \text{Loss}(\mathbf{w})$$





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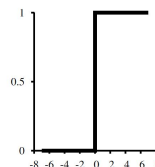
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**2)** The linear classifier always announces a completely confident prediction of 1 or 0, even for examples that are very close to the boundary

- In many situations, we really need more gradated predictions



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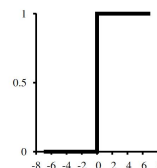
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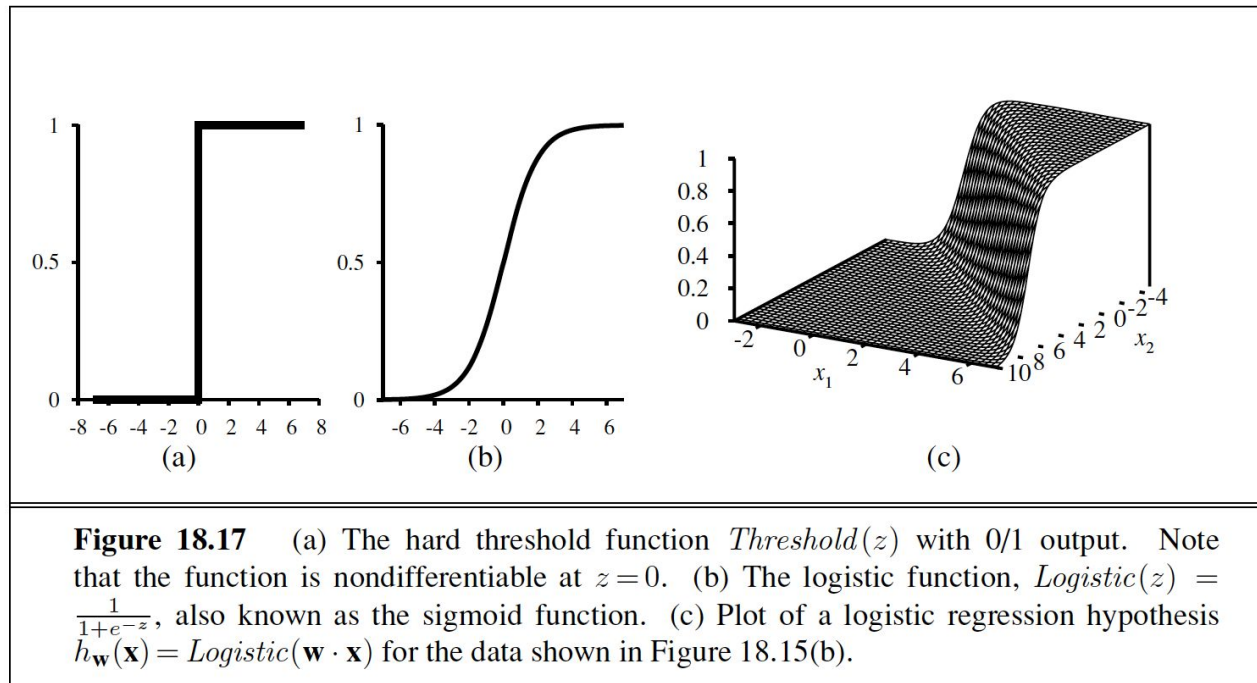
These issues can be resolved to a large extent by softening the threshold function—approximating the hard threshold with a continuous, differentiable function

# 18.6.4 Linear Classification with Logistic Regression

$$\text{Logistic}(z) = \frac{1}{1 + e^{-z}}$$

$$h_{\mathbf{w}}(\mathbf{x}) = \text{Logistic}(\mathbf{w} \cdot \mathbf{x}) = \frac{1}{1 + e^{-\mathbf{w} \cdot \mathbf{x}}}$$

The process of fitting the weights of this model to minimize loss on a data set is **logistic regression**.



# 18.6.4 Updating Weights in Logistic Regression (multivariate)

$$h_{\mathbf{w}}(\mathbf{x}) = \text{Logistic}(\mathbf{w} \cdot \mathbf{x}) = \frac{1}{1 + e^{-\mathbf{w} \cdot \mathbf{x}}}$$

Loss function for Logistic Regression

$$w_i \leftarrow w_i - \alpha \frac{\partial}{\partial w_i} \text{Loss}(\mathbf{w})$$

Perceptron Learning rule (gradient descent)

$$\frac{\partial}{\partial w_i} \text{Loss}(\mathbf{w}) = \frac{\partial}{\partial w_i} (y - h_{\mathbf{w}}(\mathbf{x}))^2$$

Derivative of the Loss function

How to calculate this?

Power function rule:

$$\frac{d(g^n)}{dx} = n g^{n-1} \frac{dg}{dx}$$

Chain rule:

$$\partial g(f(x)) / \partial x = g'(f(x)) \partial f(x) / \partial x$$

For logistic function:

$$L'(x) = L(x) (1 - L(x))$$

# 18.6.4 Updating Weights in Logistic Regression (multivariate)

$$\begin{aligned}
 \frac{\partial}{\partial w_i} \text{Loss}(\mathbf{w}) &= \frac{\partial}{\partial w_i} (y - h_{\mathbf{w}}(\mathbf{x}))^2 \\
 &= 2(y - h_{\mathbf{w}}(\mathbf{x})) \frac{\partial}{\partial w_i} (y - h_{\mathbf{w}}(\mathbf{x})) \quad \text{from power function rule } \frac{\partial}{\partial x} g(x)^n = n[g(x)]^{n-1} \frac{\partial}{\partial x} (g(x)) \\
 &= 2(y - h_{\mathbf{w}}(\mathbf{x})) \frac{\partial}{\partial w_i} (-h_{\mathbf{w}}(\mathbf{x})) \\
 &= -2(y - h_{\mathbf{w}}(\mathbf{x})) \frac{\partial}{\partial w_i} h_{\mathbf{w}}(\mathbf{x}) \\
 &= -2(y - h_{\mathbf{w}}(\mathbf{x})) \frac{\partial}{\partial w \cdot \mathbf{x}} g(\mathbf{w} \cdot \mathbf{x}) \cdot \frac{\partial}{\partial w_i} (\mathbf{w} \cdot \mathbf{x}) \quad \text{from chain rule } \frac{\partial}{\partial x} g(f(x)) = g'(f(x)) \cdot \frac{\partial}{\partial x} f(x) \\
 &= -2(y - h_{\mathbf{w}}(\mathbf{x})) [g'(\mathbf{w} \cdot \mathbf{x})] \cdot \frac{\partial}{\partial w_i} (w_1 x_1 + w_2 x_2 + \dots + w_n x_n) \\
 &= -2(y - h_{\mathbf{w}}(\mathbf{x})) [g(\mathbf{w} \cdot \mathbf{x}) (1 - g(\mathbf{w} \cdot \mathbf{x}))] x_i \quad \text{derivative of a logistic function } \frac{\partial}{\partial x} \left( \frac{1}{1+e^{-x}} \right) = \frac{1}{1+e^{-x}} \left( 1 - \frac{1}{1+e^{-x}} \right) \\
 &\therefore \frac{\partial}{\partial w_i} \text{Loss}(\mathbf{w}) = -2(y - h_{\mathbf{w}}(\mathbf{x})) h_{\mathbf{w}}(\mathbf{x}) (1 - h_{\mathbf{w}}(\mathbf{x})) x_i
 \end{aligned}$$

$$w_i \leftarrow w_i - \alpha \frac{\partial}{\partial w_i} \text{Loss}(\mathbf{w})$$

$$w_i \leftarrow w_i + \alpha (y - h_{\mathbf{w}}(\mathbf{x})) \times h_{\mathbf{w}}(\mathbf{x}) (1 - h_{\mathbf{w}}(\mathbf{x})) \times x_i$$

## 18.6.4 Linear Classification with Logistic Regression

Additional **advantages of Logistic Regression** (compared to hard threshold):

1. In a linearly separable case, logistic regression is somewhat slower to converge, but behaves much more predictably

## 18.6.4 Linear Classification with Logistic Regression

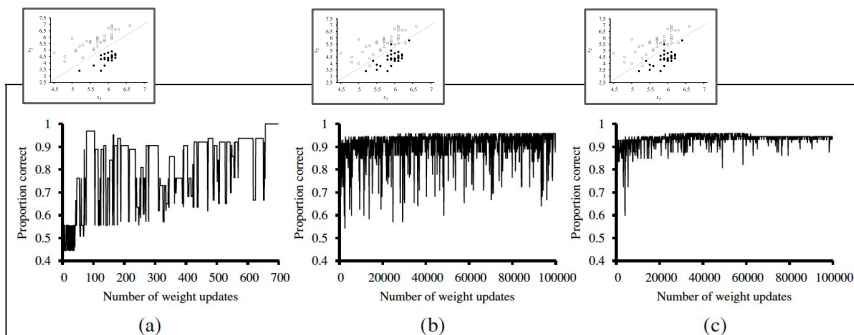
Additional **advantages of Logistic Regression** (compared to hard threshold):

1. In a linearly separable case, logistic regression is somewhat slower to converge, but behaves much more predictably
2. Where the data are noisy and nonseparable, logistic regression converges far more quickly and reliably; these advantages tend to carry over into real-world applications

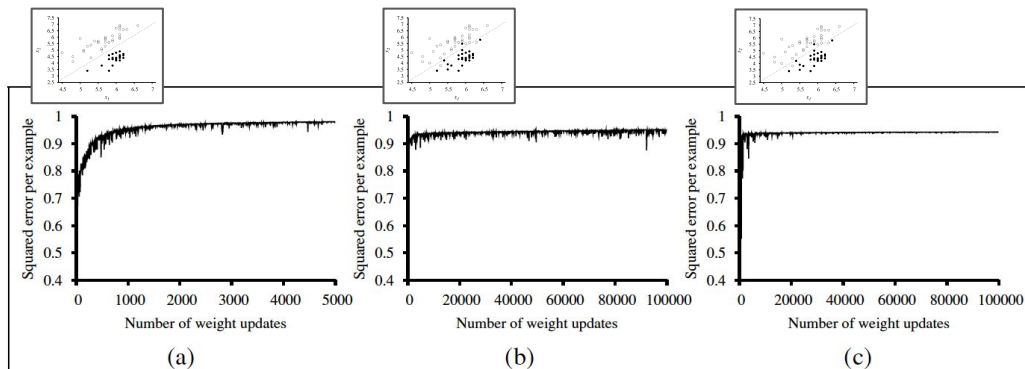
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Additional **advantages of Logistic Regression** (compared to hard threshold):

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**Figure 18.16** (a) Plot of total training-set accuracy vs. number of iterations through the training set for the perceptron learning rule, given the earthquake/explosion data in Figure 18.15(a). (b) The same plot for the noisy, non-separable data in Figure 18.15(b); note the change in scale of the  $x$ -axis. (c) The same plot as in (b), with a learning rate schedule  $\alpha(t) = 1000/(1000 + t)$ .



**Figure 18.18** Repeat of the experiments in Figure 18.16 using logistic regression and squared error. The plot in (a) covers 5000 iterations rather than 1000, while (b) and (c) use the same scale.



### Logistic Regression Example using the Iris Flower Dataset

- We would like to predict if a given data point (one row) belongs to the 'Iris-setosa' class
- Dataset: [rawdata](#) and [metadata](#)

```
In [25]: from keras.models import Sequential
from keras.layers import Dense
import numpy as np
import matplotlib.pyplot as plt

datapath = 'https://raw.githubusercontent.com/badriadhikari/2019-Fall-AI/master/MODULE-I/iris.dat'
dataset = np.genfromtxt(datapath, delimiter=";", dtype = str)

print('')
print(dataset.shape)
print('')
print(dataset[0:5])

dataset[:, 4] = np.where(dataset[:, 4] == 'Iris-setosa', 1, 0)
#dataset[:, 4] = np.where(dataset[:, 4] == 'Iris-versicolor', 1, 0)
#dataset[:, 4] = np.where(dataset[:, 4] == 'Iris-virginica', 1, 0)

dataset = dataset.astype(float)
print('')
print(dataset[0:5])

np.random.shuffle(dataset)
print('')
print(dataset[0:5])

train = dataset[:100]
valid = dataset[100:]
print('')
print(train.shape)
print('')
print(valid.shape)

(150, 5)

[['5.1' '3.5' '1.4' '0.2' 'Iris-setosa']
 ['4.9' '3.0' '1.4' '0.2' 'Iris-setosa']
 ['4.7' '3.2' '1.3' '0.2' 'Iris-setosa']
 ['4.6' '3.1' '1.5' '0.2' 'Iris-setosa']
 ['5.0' '3.6' '1.4' '0.2' 'Iris-setosa']]

[[5.1 3.5 1.4 0.2 1.]
 [4.9 3.  1.4 0.2 1.]
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 [5.  3.6 1.4 0.2 1.]]

[[6.1 2.8 4.  1.3 0.]
 [5.  3.3 1.4 0.2 1.]
 [6.4 2.7 5.3 1.9 0.]
 [5.7 2.5 5.  2.  0.]
 [4.8 3.4 1.9 0.2 1.]]

(100, 5)

(50, 5)
```

# 18.7 Artificial Neural Networks (Motivation)

In 1953, Professor Theodor Erismann devised an experiment

- performing it upon his assistant and student, Ivo Kohler

He made Kohler wear a pair of hand-engineered goggles

- Specially arranged mirrors flipped the light that would reach eyes, top becoming bottom, and bottom top.

After 10 days, Kohler had grown accustomed to the invariably upside-down world

- everything seemed to him normal, rightside-up
- He could do everyday activities in public perfectly well: walk along a crowded sidewalk, even ride a bicycle



<https://www.theguardian.com/education/2012/nov/12/improbable-research-seeing-upside-down>

# 18.7 Artificial Neural Networks (Motivation)

Science

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## Eye-specific termination bands in tecta of three-eyed frogs

M Constantine-Paton, MI Law

+ See all authors and affiliations

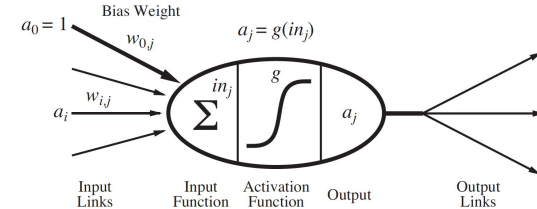
Science 10 Nov 1978:  
Vol. 202, Issue 4368, pp. 639-641  
DOI: 10.1126/science.309179

Dr. Martha Constantine-Paton is a neuroscientist at MIT



## 18.7.1 Neural Network Structures

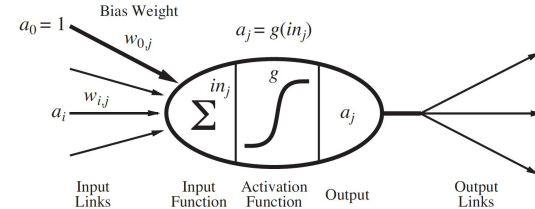
- Neural networks are composed of nodes or units connected by directed links
- A link from unit **i** to unit **j** serves to propagate the activation  **$a_i$**  from **i** to **j**
- Each link also has a numeric weight  **$w_{i,j}$**  associated with it, which determines the strength and sign of the connection
- Just as in linear regression models, each unit has a dummy input  $a_0 = 1$  with an associated weight  **$w_{0,j}$**



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- Each unit **j** first computes a weighted sum of its inputs:

$$in_j = \sum_{i=0}^n w_{i,j} a_i$$



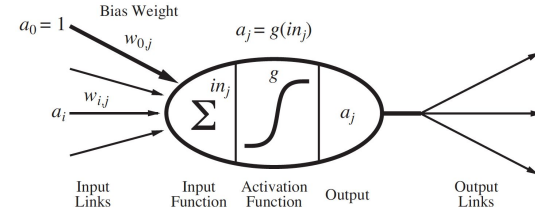
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- Each unit **j** first computes a weighted sum of its inputs:

$$in_j = \sum_{i=0}^n w_{i,j} a_i$$

- Then, we apply an **activation function g** to this sum to derive the output:

$$a_j = g(in_j) = g\left(\sum_{i=0}^n w_{i,j} a_i\right)$$



# 18.7.1 Neural Network Structures

There are distinct ways to connect neurons:

1. A feed-forward network has connections only in one direction—that is, it forms a directed acyclic graph
  - a. Every node receives input from “upstream” nodes and delivers output to “downstream” nodes; there are **no loops**
  - b. A feed-forward network represents a function of its current input; thus, it has **no internal state** other than the weights themselves

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  - b. A feed-forward network represents a function of its current input; thus, it has **no internal state** other than the weights themselves
2. A recurrent network, on the other hand, feeds its outputs back into its own inputs
  - a. This means that the activation levels of the network form a dynamical system that may **reach a stable state or exhibit oscillations** or even chaotic behavior
  - b. Moreover, the response of the network to a given input depends on its initial state, which may depend on previous inputs
  - c. Hence, recurrent networks (unlike feed-forward networks) can support **short-term memory**
  - d. This makes them more **interesting as models of the brain**, but also more **difficult to understand**

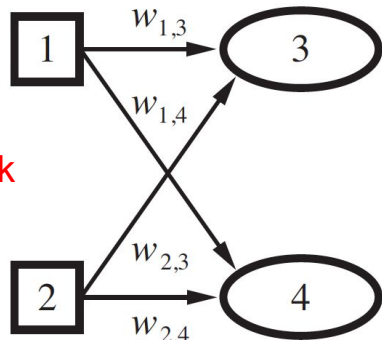
We will focus on feed-forward networks!



## 18.7.2 Single-layer feed-forward neural networks (Perceptrons)

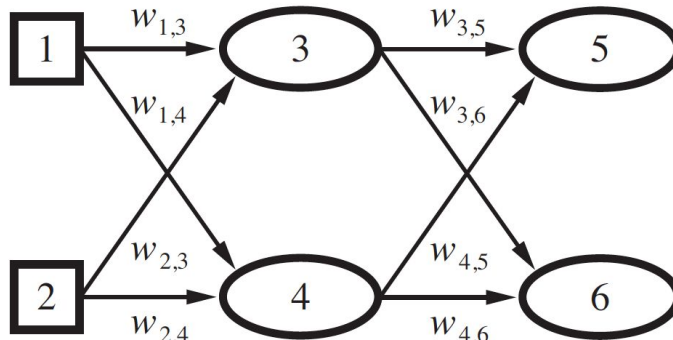
A network with all the inputs connected directly to the outputs is called a single-layer neural network, or a perceptron network

Perceptron Network



(a)

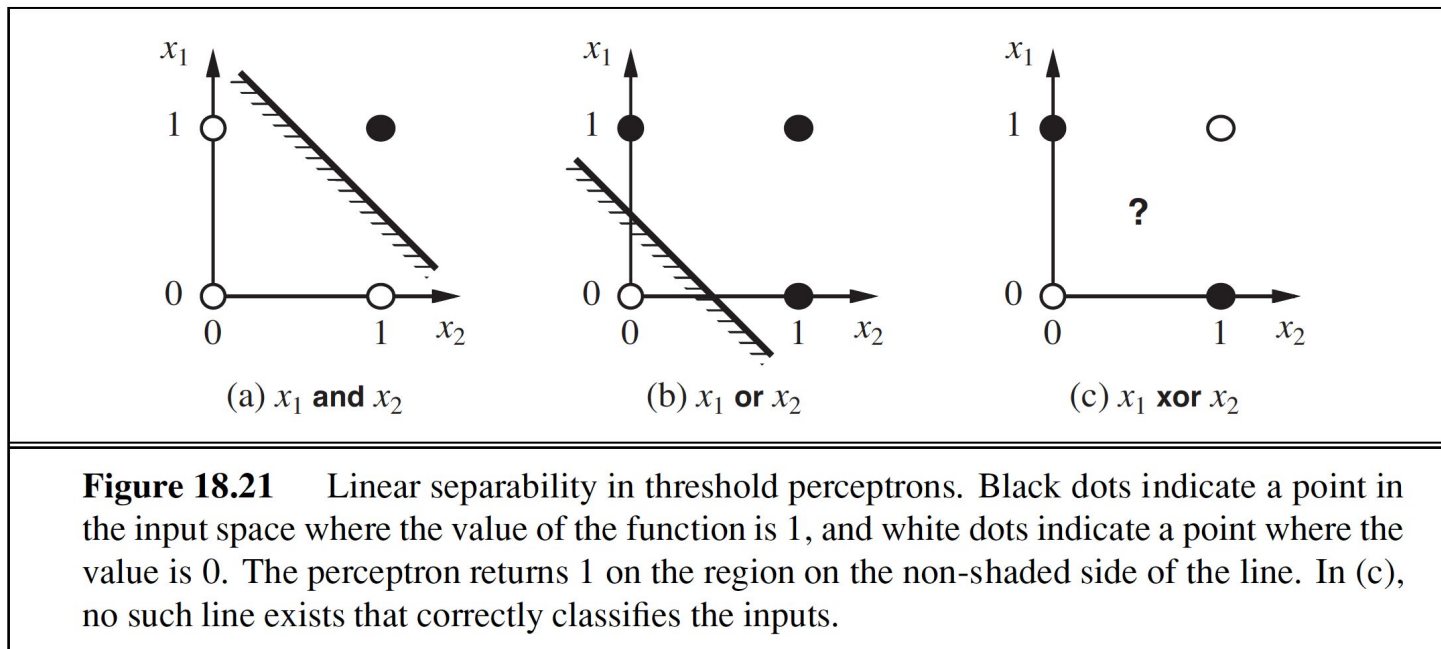
Neural Network



(b)

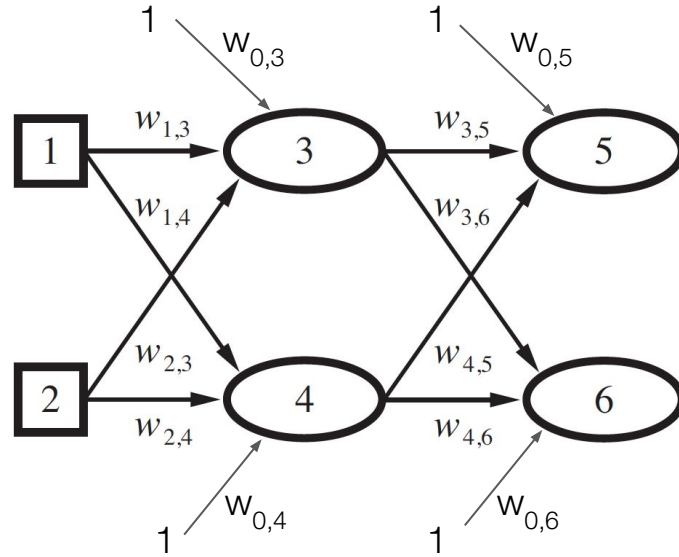
**Figure 18.20** (a) A perceptron network with two inputs and two output units. (b) A neural network with two inputs, one hidden layer of two units, and one output unit. Not shown are the dummy inputs and their associated weights.

## 18.7.2 Single-layer feed-forward neural networks (Perceptrons)



What is the key limitation of a perceptron network? How to overcome the limitation?

## 18.7.3 Multilayer Feed-forward Neural Networks



What do  $a_3$  and  $a_4$  stand for?



What are  $w_{0,5}$ ,  $w_{0,3}$ ,  $w_{0,4}$ , and  $w_{0,6}$ ?

$$\begin{aligned} a_5 &= g(w_{0,5} + w_{3,5} a_3 + w_{4,5} a_4) \\ &= g(w_{0,5} + w_{3,5} g(w_{0,3} + w_{1,3} a_1 + w_{2,3} a_2) + w_{4,5} g(w_{0,4} + w_{1,4} a_1 + w_{2,4} a_2)) \\ &= g(w_{0,5} + w_{3,5} g(w_{0,3} + w_{1,3} x_1 + w_{2,3} x_2) + w_{4,5} g(w_{0,4} + w_{1,4} x_1 + w_{2,4} x_2)) \end{aligned}$$

# Neural Network Structures

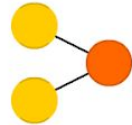
*A mostly complete chart of*

## Neural Networks

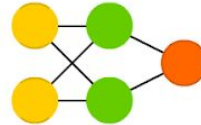
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- Backfed Input Cell
- Input Cell
- △ Noisy Input Cell
- Hidden Cell
- Probabilistic Hidden Cell
- △ Spiking Hidden Cell
- Output Cell
- Match Input Output Cell
- Recurrent Cell

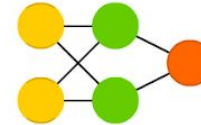
Perceptron (P)



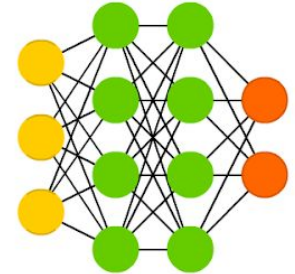
Feed Forward (FF)



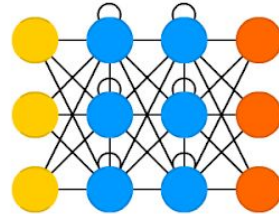
Radial Basis Network (RBF)



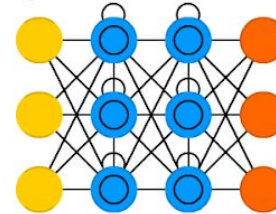
Deep Feed Forward (DFF)



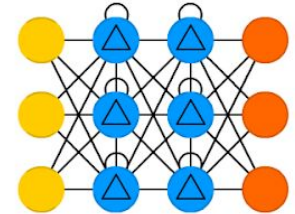
Recurrent Neural Network (RNN)



Long / Short Term Memory (LSTM)

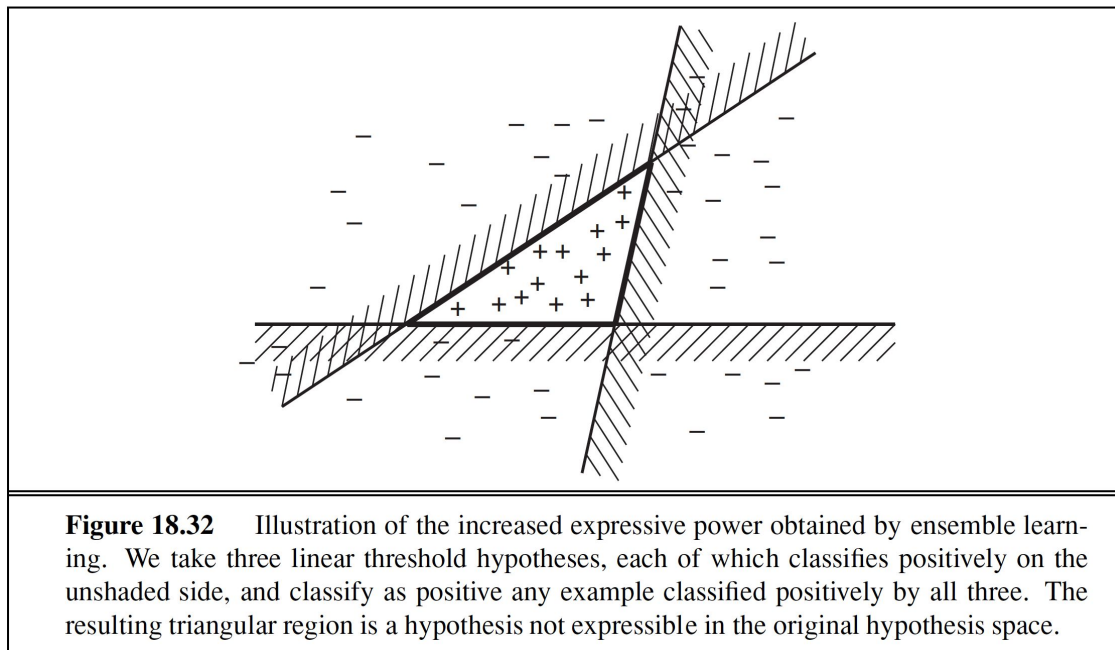


Gated Recurrent Unit (GRU)



## 18.10 Ensemble Learning

- The idea of ensemble learning methods is to select a collection, or ensemble, of hypotheses from the hypothesis space and combine their predictions; i.e. use multiple models; e.g. boosting



# Summary

- If the available feedback provides the correct answer for example inputs, then the learning problem is called **supervised learning**. The task is to learn a function  $y = h(x)$ .
- Learning a discrete-valued function is called **classification**; learning a continuous function is called **regression**.
- Sometimes not all errors are equal. A **loss function** tells us how bad each error is.
- **Logistic regression** replaces the perceptron's hard threshold with a soft threshold defined by a logistic function. Gradient descent works well even for noisy data that are not linearly separable.
- Neural networks represent complex nonlinear functions with a network of linear threshold units.
- Ensemble methods such as **boosting** often perform better than individual methods.