

# Deep Learning for Computer Vision

113-1/Fall 2024

<https://cool.ntu.edu.tw/courses/41702> (NTU COOL)

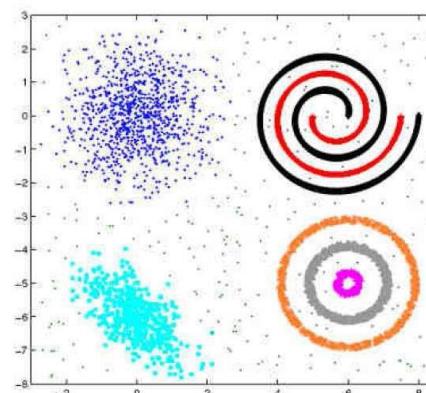
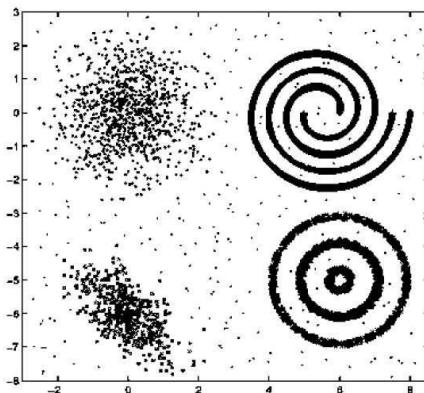
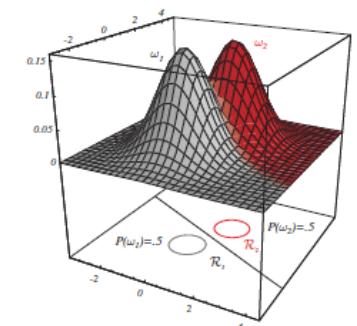
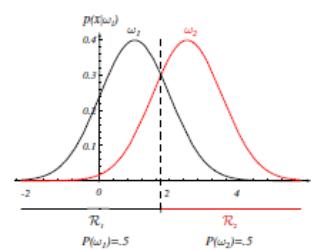
<http://vllab.ee.ntu.edu.tw/dlcv.html> (Public website)

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Dept. Electrical Engineering, National Taiwan University

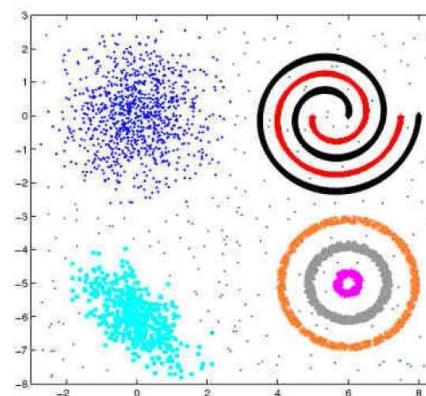
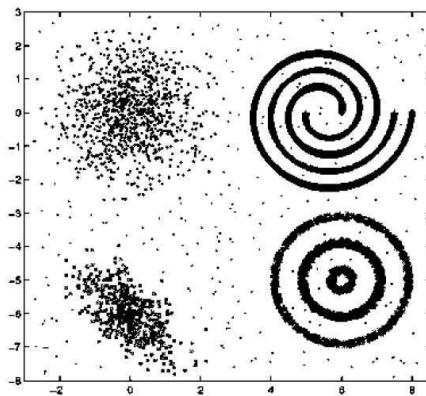
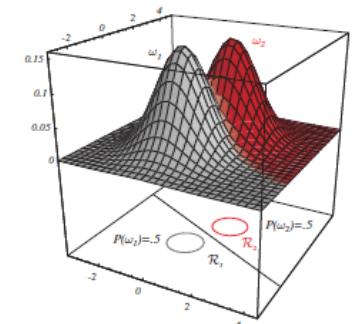
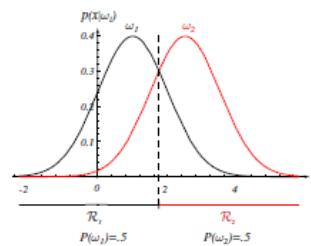
# What's to Be Covered in This Lecture...

- Unsupervised vs. Supervised Learning
  - Clustering & Dimension Reduction
  - Training, testing, & validation
  - Linear Classification
  - From Linear Classifier to Neural Nets



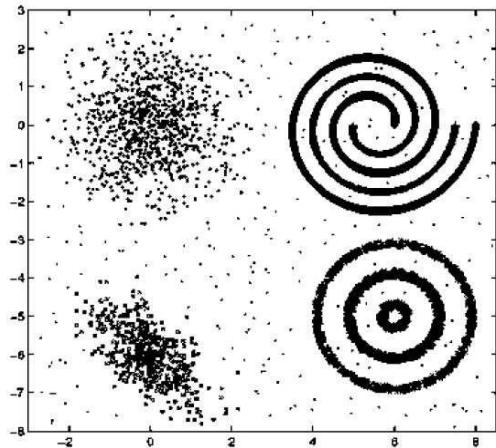
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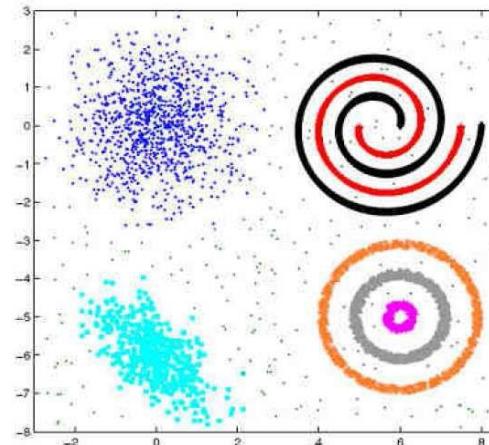


# Clustering

- Clustering is an unsupervised algorithm.
  - Given:  
a set of N unlabeled instances  $\{x_1, \dots, x_N\}$ ; # of clusters K
  - Goal: group the samples into K partitions
- Remarks:
  - High within-cluster (intra-cluster) similarity
  - Low between-cluster (inter-cluster) similarity
  - But...how to determine a proper similarity measure?



(a) Input data



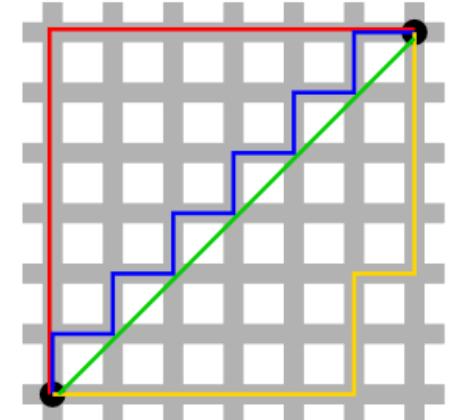
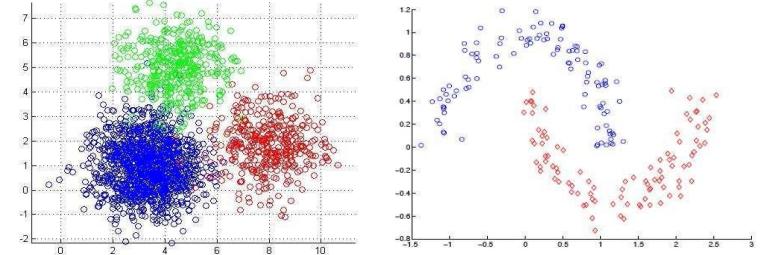
(b) Desired clustering

# Similarity is NOT Always Objective...



# Clustering (cont'd)

- Similarity:
  - A key component/measure to perform data clustering
  - Inversely proportional to distance
  - Example distance metrics:
    - Euclidean distance (L2 norm):  $d(x, z) = \|x - z\|_2 = \sqrt{\sum_{i=1}^D (x_i - z_i)^2}$
    - Manhattan distance (L1 norm):  $d(x, z) = \|x - z\|_1 = \sum_{i=1}^D |x_i - z_i|$
    - Note that  $p$ -norm of  $x$  is denoted as:

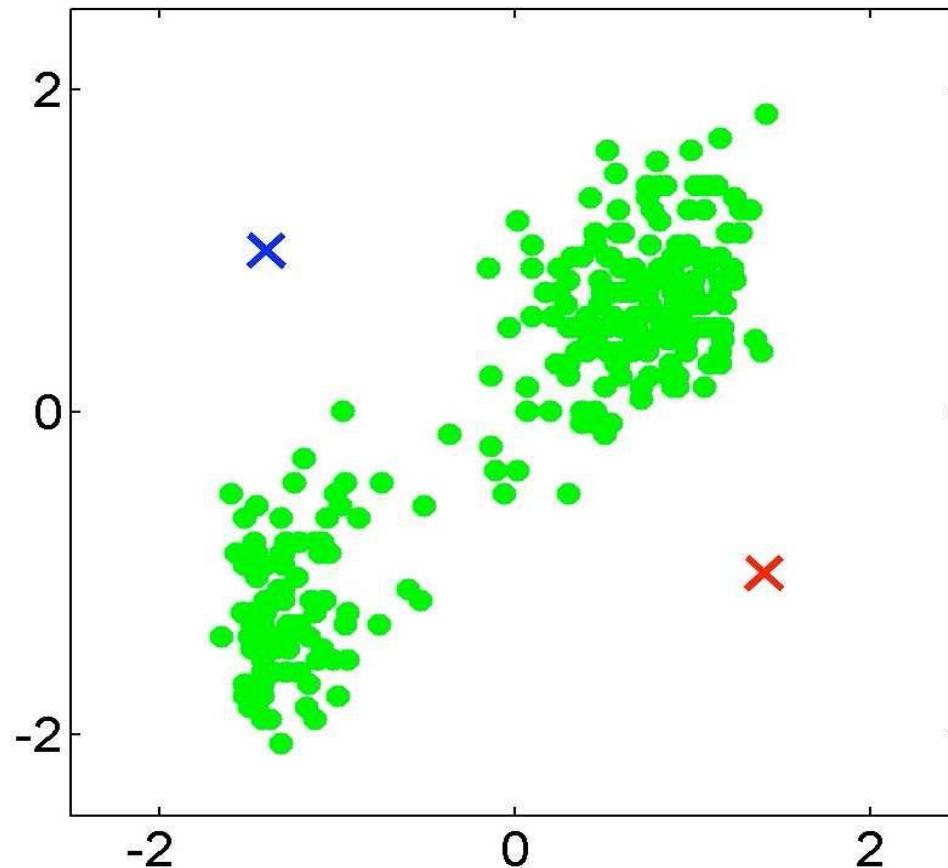


# K-Means Clustering

- **Input:**  $N$  examples  $\{\mathbf{x}_1, \dots, \mathbf{x}_N\}$  ( $\mathbf{x}_n \in \mathbb{R}^D$ ); number of partitions  $K$
- **Initialize:**  $K$  cluster centers  $\mu_1, \dots, \mu_K$ . Several initialization options:
  - Randomly initialize  $\mu_1, \dots, \mu_K$  anywhere in  $\mathbb{R}^D$
  - Or, simply choose any  $K$  examples as the cluster centers
- **Iterate:**
  - Assign each of example  $\mathbf{x}_n$  to its closest cluster center
  - Recompute the new cluster centers  $\mu_k$  (mean/centroid of the set  $C_k$ )
  - Repeat while not converge
- **Possible convergence criteria:**
  - Cluster centers do not change anymore
  - Max. number of iterations reached
- **Output:**
  - $K$  clusters (with centers/means of each cluster)

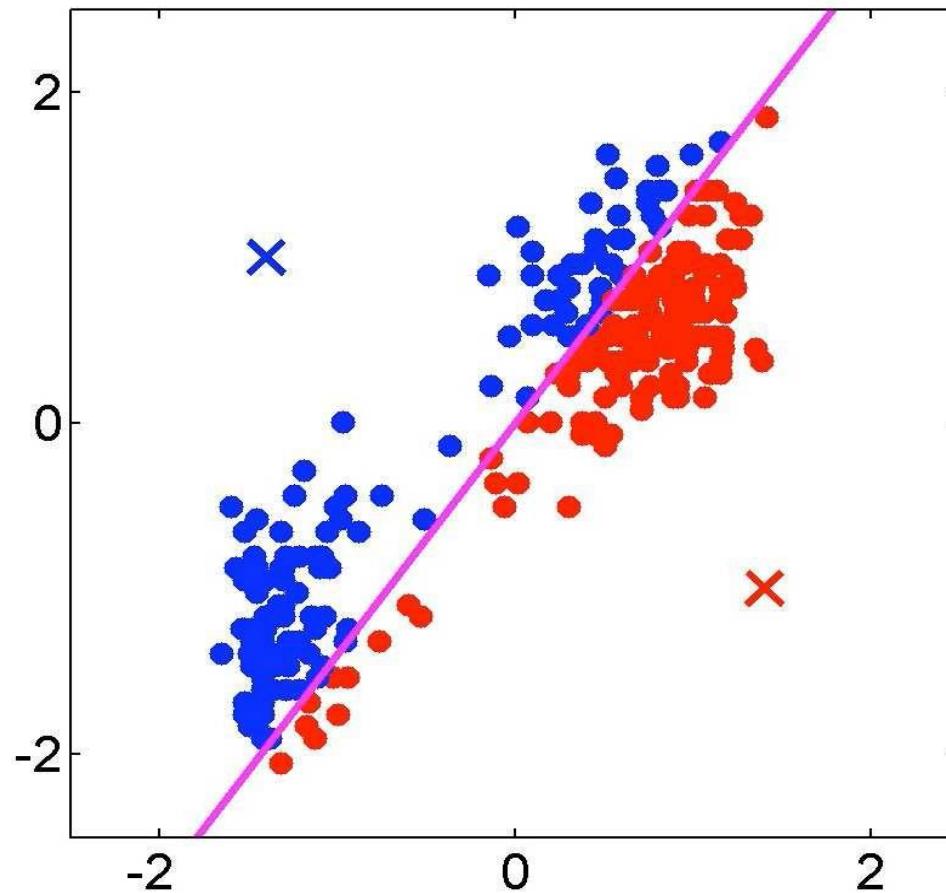
# K-Means Clustering

- Example ( $K = 2$ ): Initialization, iteration #1: pick cluster centers



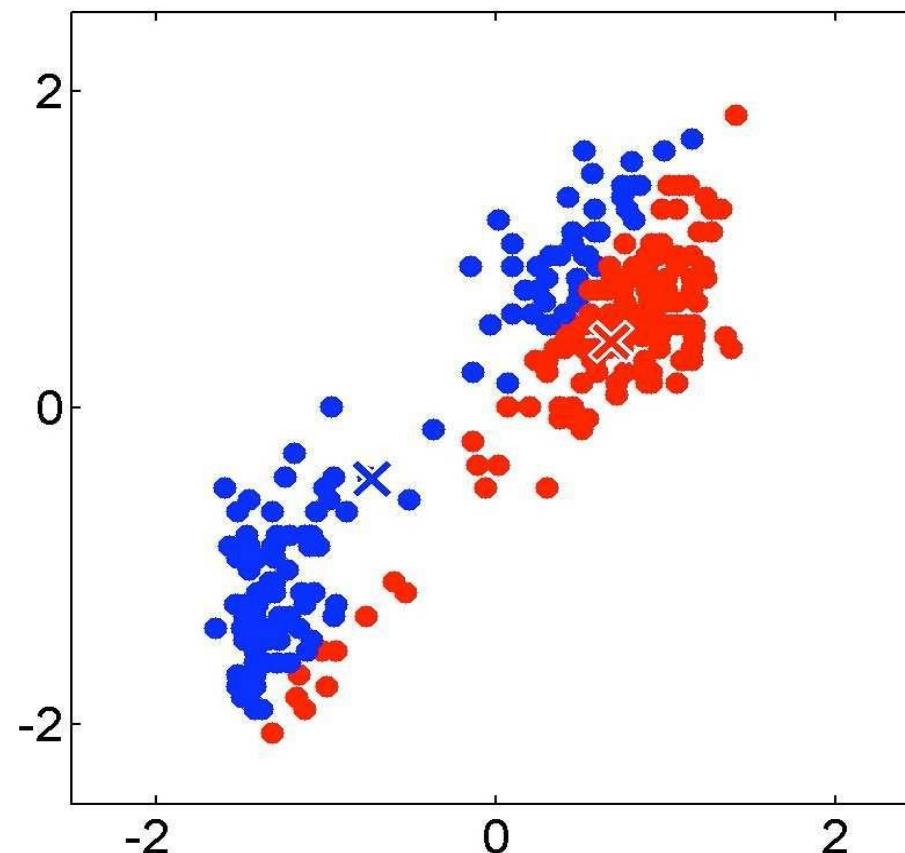
# K-Means Clustering

- Example ( $K = 2$ ): iteration #1-2, assign data to each cluster



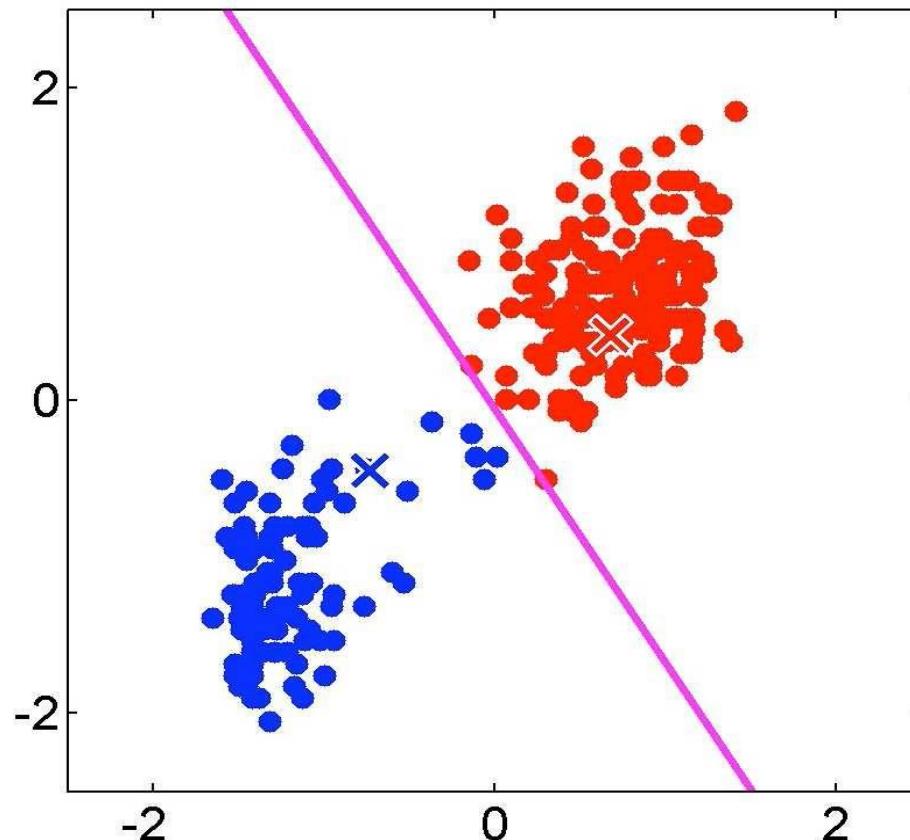
# K-Means Clustering

- Example ( $K = 2$ ): iteration #2-1, update cluster centers



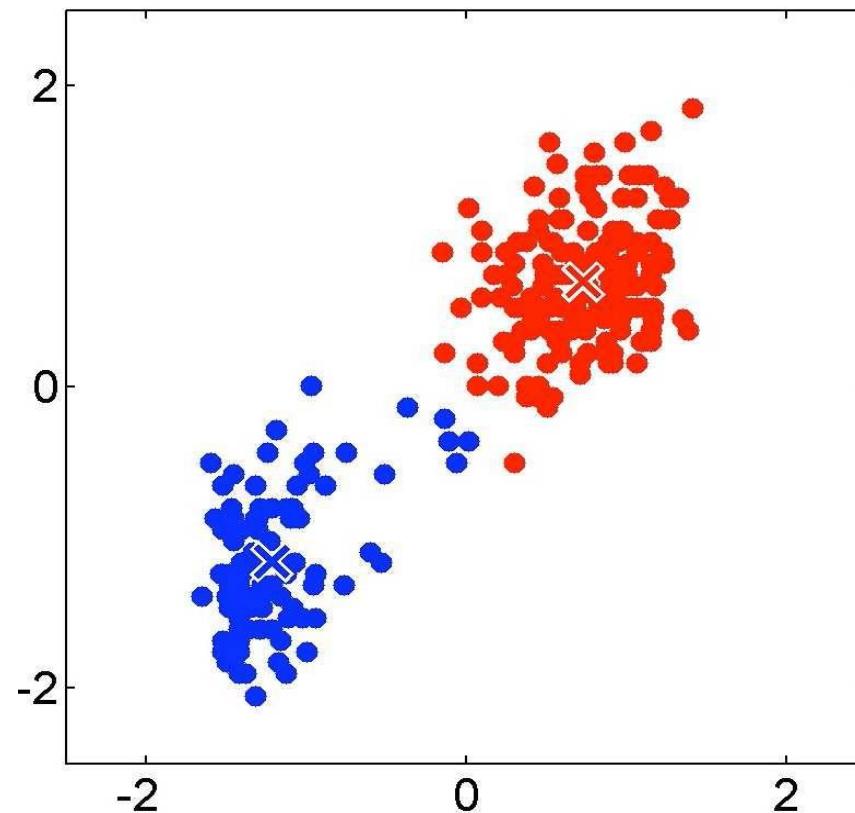
# K-Means Clustering

- Example ( $K = 2$ ): iteration #2, assign data to each cluster



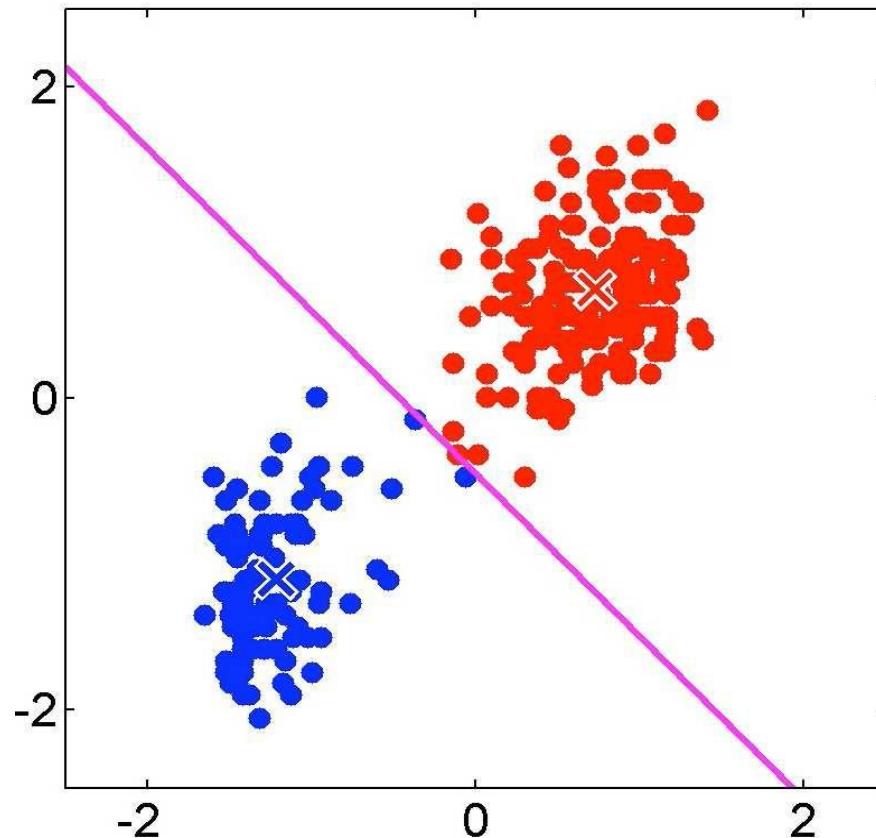
# K-Means Clustering

- Example ( $K = 2$ ): iteration #3-1



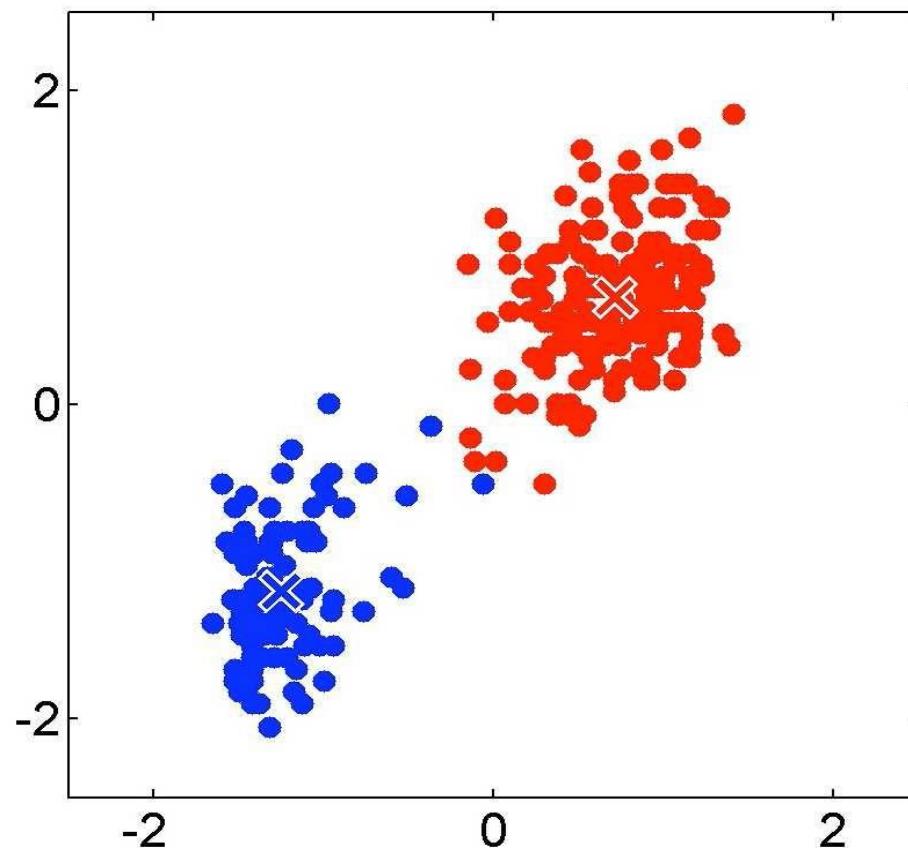
# K-Means Clustering

- Example ( $K = 2$ ): iteration #3-2



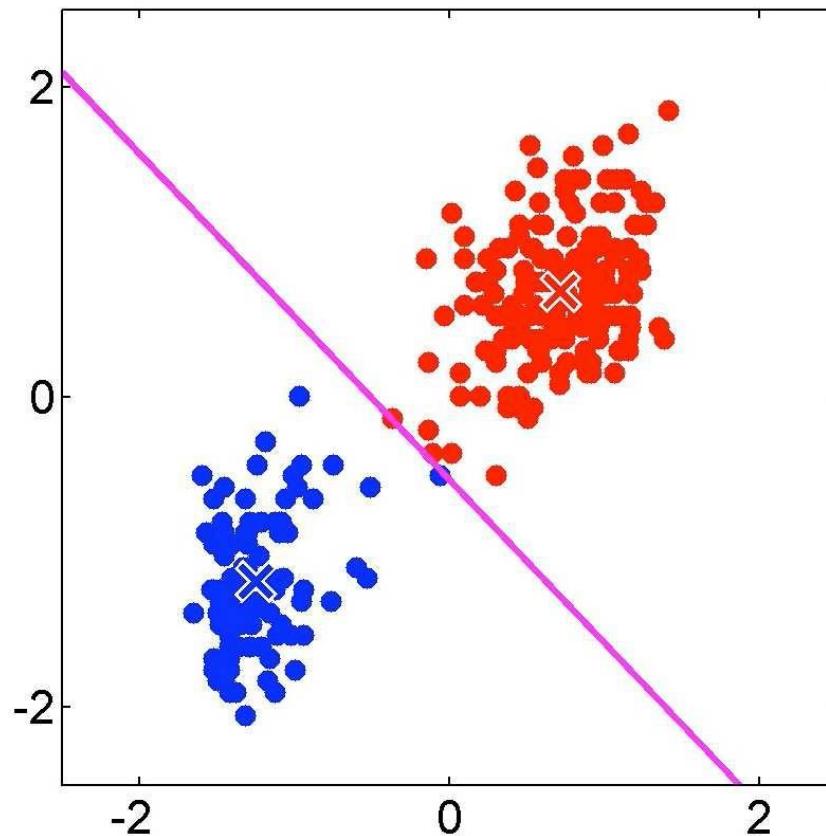
# K-Means Clustering

- Example ( $K = 2$ ): iteration #4-1



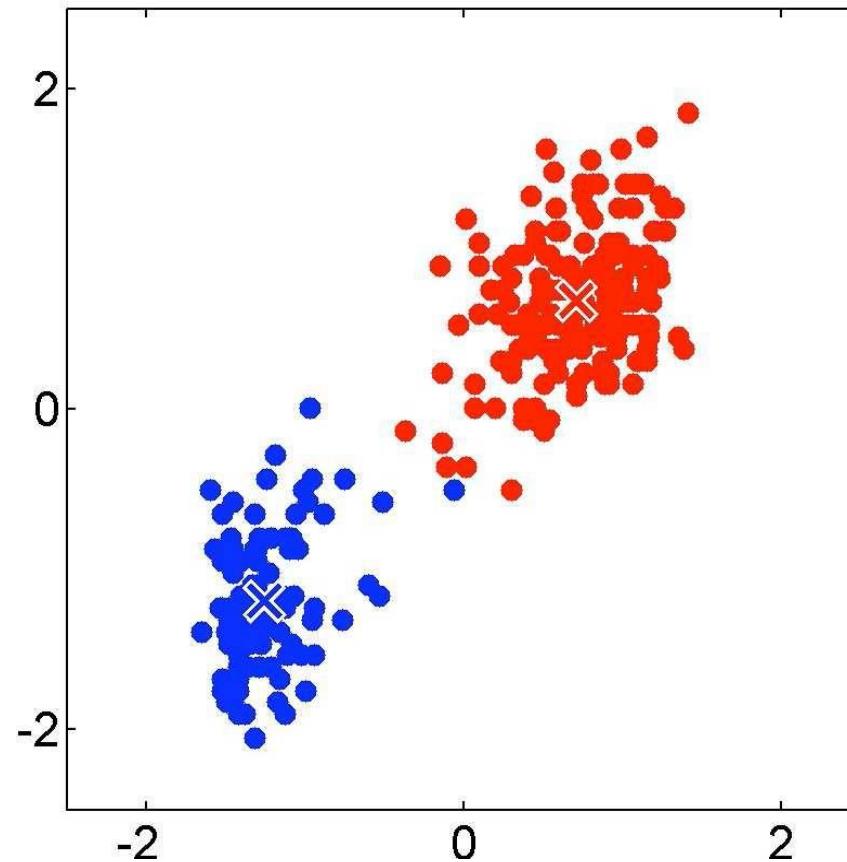
# K-Means Clustering

- Example ( $K = 2$ ): iteration #4-2



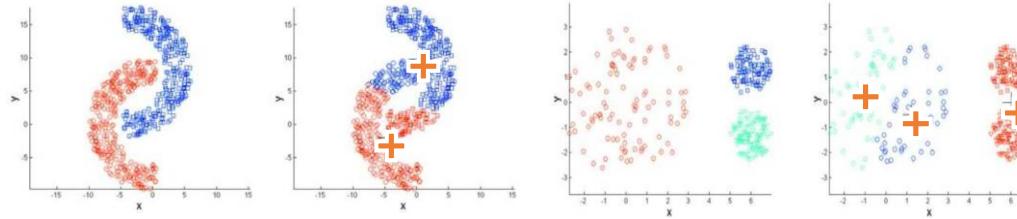
# K-Means Clustering

- Example ( $K = 2$ ): iteration #5, cluster means are not changed.



# K-Means Clustering (cont'd)

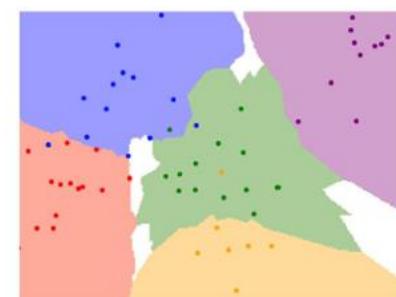
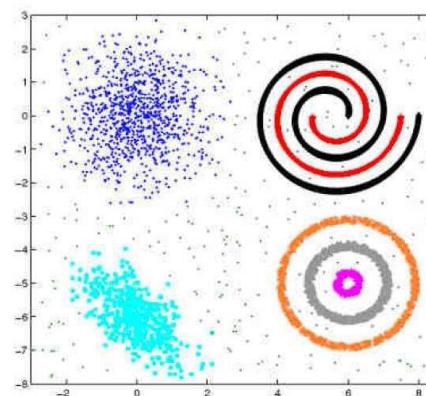
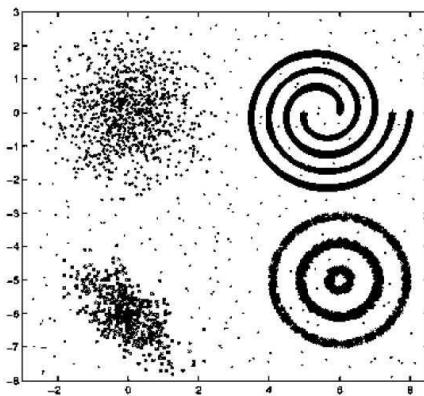
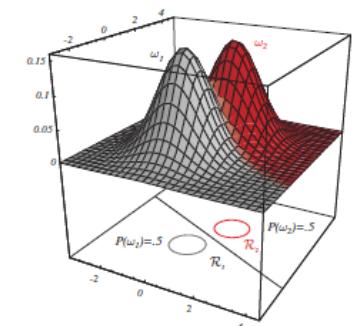
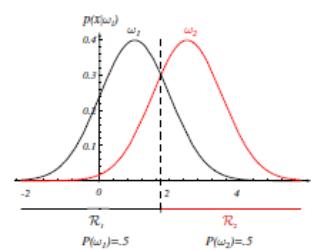
- Easy to implement, but...
  - Preferable for round shaped clusters with similar sizes



- Limitations
  - Sensitive to initialization →
  - Sensitive to outliers →
  - Hard assignment only →
- Remarks
  - Expectation-maximization (EM) algorithm
  - Speed-up possible by hierarchical clustering (e.g.,  $100 = 10^2$  clusters)

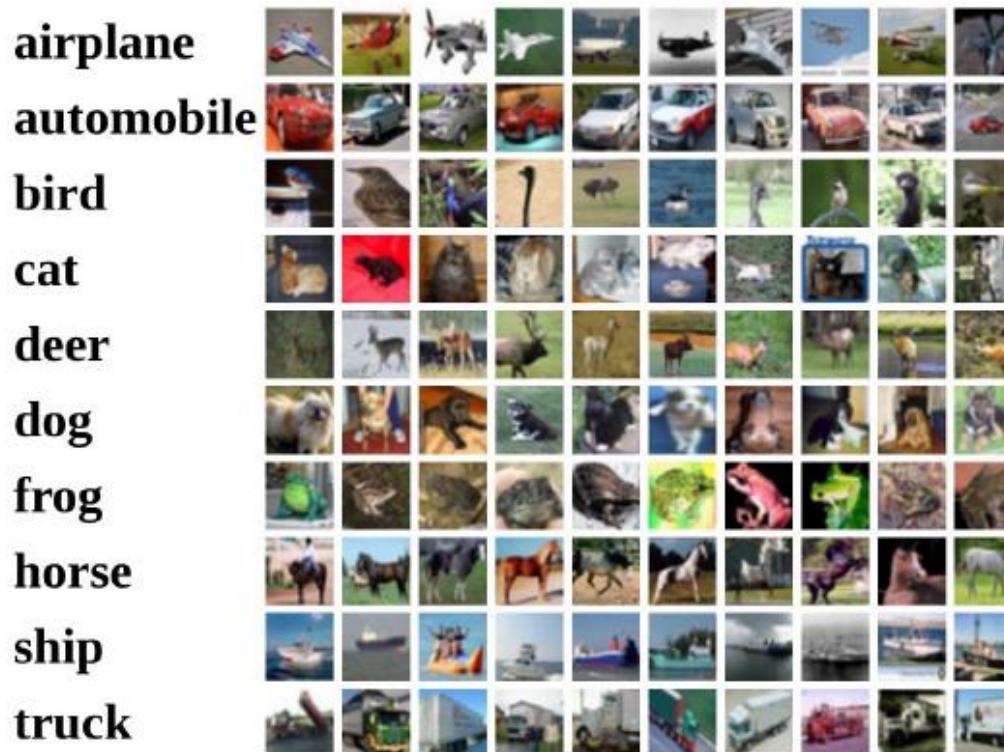
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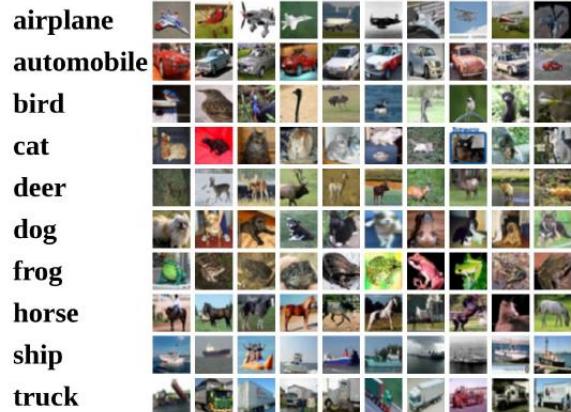


# Linear Classification

- Linear Classifier
  - Consider that we have 10 object categories of interest
    - E.g., CIFAR10 with 50K training & 10K test images of 10 categories.  
And, each image is of size 32 x 32 x 3 pixels.

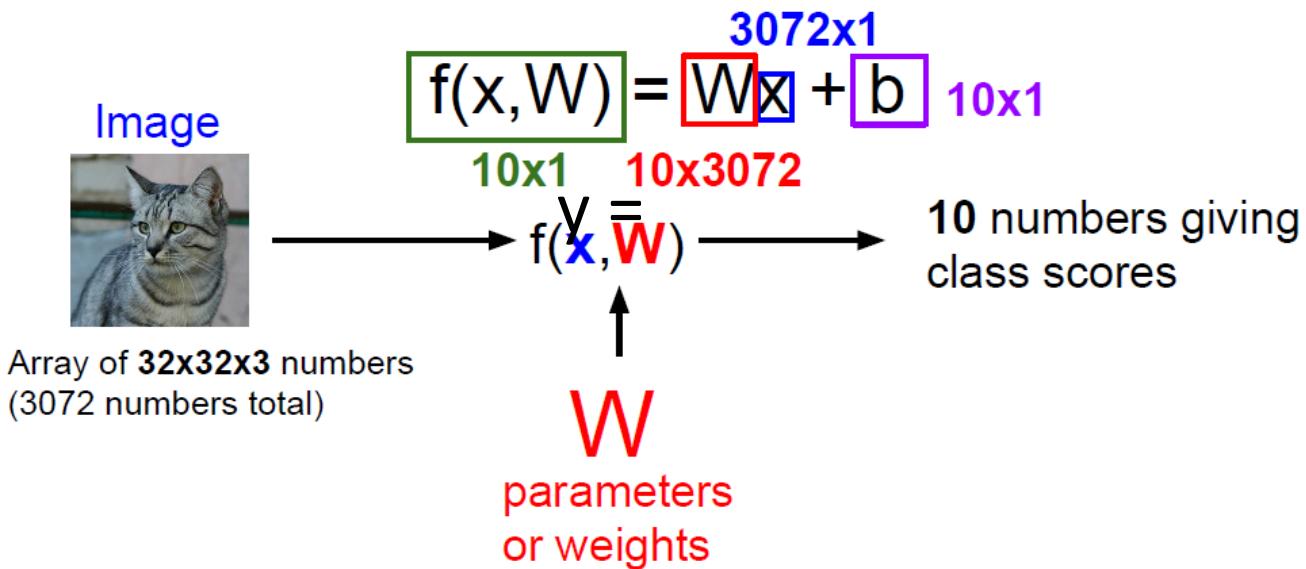


# Linear Classification (cont'd)



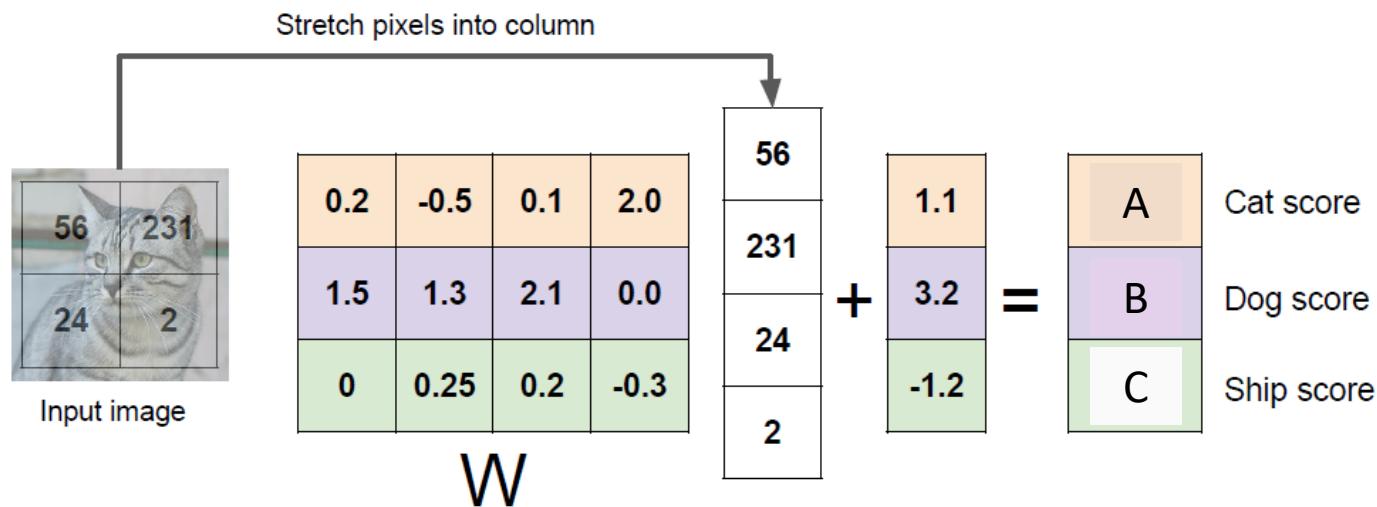
- Linear Classifier

- Let's take the input image as  $\mathbf{x}$ , and the linear classifier as  $\mathbf{W}$ .  
We need  $\mathbf{y} = \mathbf{W}\mathbf{x} + \mathbf{b}$  as a 10-dimensional output vector, indicating the score for each class.



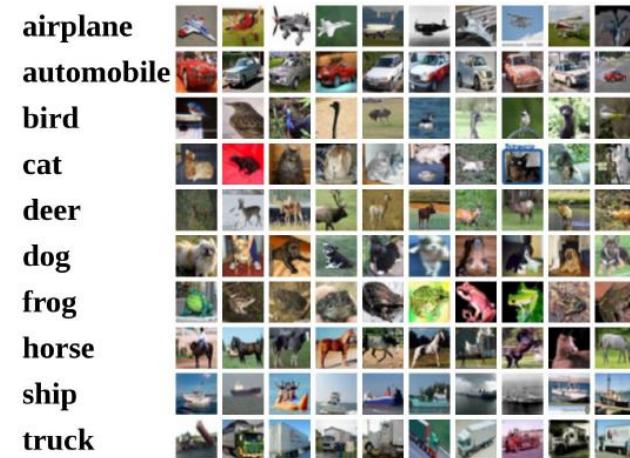
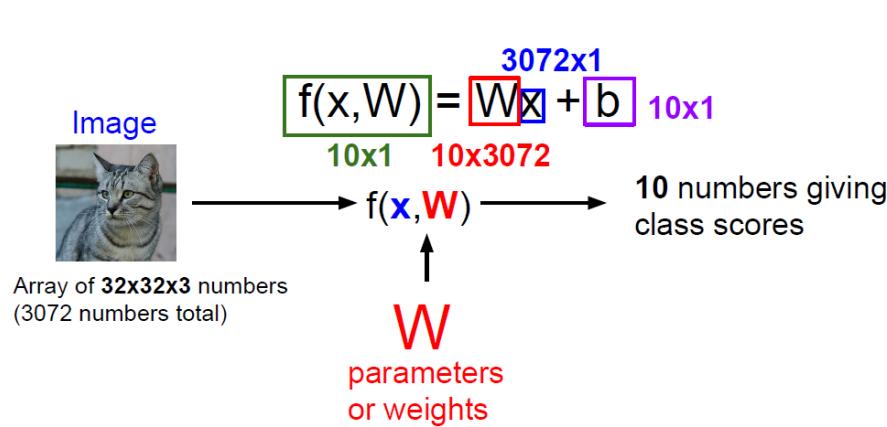
# Linear Classification (cont'd)

- Linear Classifier
  - For example, an image with  $2 \times 2$  pixels & 3 classes of interest we need to learn a linear classifier  $\mathbf{W}$  (plus a bias  $\mathbf{b}$ ), so that desirable outputs  $\mathbf{y} = \mathbf{Wx} + \mathbf{b}$  can be expected.



# Final Remarks

- Interpreting  $\mathbf{W}$  in  $\mathbf{y} = \mathbf{Wx} + \mathbf{b}$ 
  - Once  $\mathbf{W}$  is learned for predicting  $\mathbf{y}$ , each row in  $\mathbf{W}$  can be viewed as an **exemplar** of the corresponding class.
  - Recall that,  $\mathbf{Wx}$  basically performs **inner product** (or **correlation**) between the input  $\mathbf{x}$  and the exemplar of each class -> **similarity** matters!
  - Any potential problem or limitation?



# Loss Function (or Cost/Objective Function)

- **Loss is a function of model parameter  $\mathbf{W}$**

- Tells us how **good/bad** our learned model  $\mathbf{W}$  in  $\mathbf{y} = \mathbf{Wx} + \mathbf{b}$  is.  
If loss function, the lower, the better!
- Given a labeled dataset  $\{(x_i, y_i)\}_{i=1}^N$   
where  $\mathbf{x}$  and  $\mathbf{y}$  indicate the input instance and its label, respectively,

Loss of a single input instance is denoted as  $L_i(f(x_i, W), y_i)$ ,

and that for the entire dataset is the sum or average of per-instance losses:

$$L = \frac{1}{N} \sum_i L_i(f(x_i, W), y_i) .$$

- In practice, calculating full sum for  $L$  is expensive.
  - Approximate sum using a **minibatch** of instances (e.g., 32, 64, 128 samples, etc.)

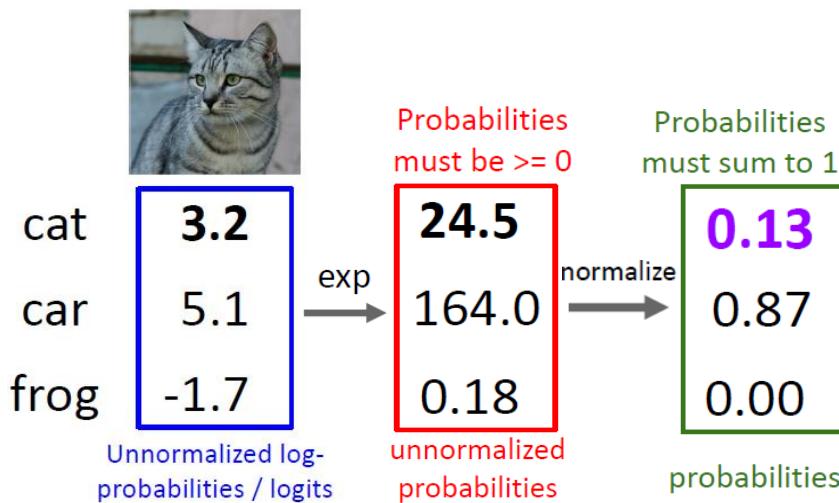
# Loss Function (cont'd)

- **Cross-Entropy Loss (Multinomial Logistic Regression)**

- Interpret the classifier scores as **probabilities**
- **Softmax function:**

$$P(Y = k \mid X = x_i) = \frac{\exp(s_k)}{\sum_j \exp(s_j)} \quad \text{with } s = f(x_i; W) \text{ as the classifier output for input } \mathbf{x}_i$$

- See example below:



$$L_i = -\log P(Y = y_i \mid X = x_i)$$

$$L_{\text{cat}} = -\log(0.13) = 2.04$$

What about this L?

What are its possibly min/max value??

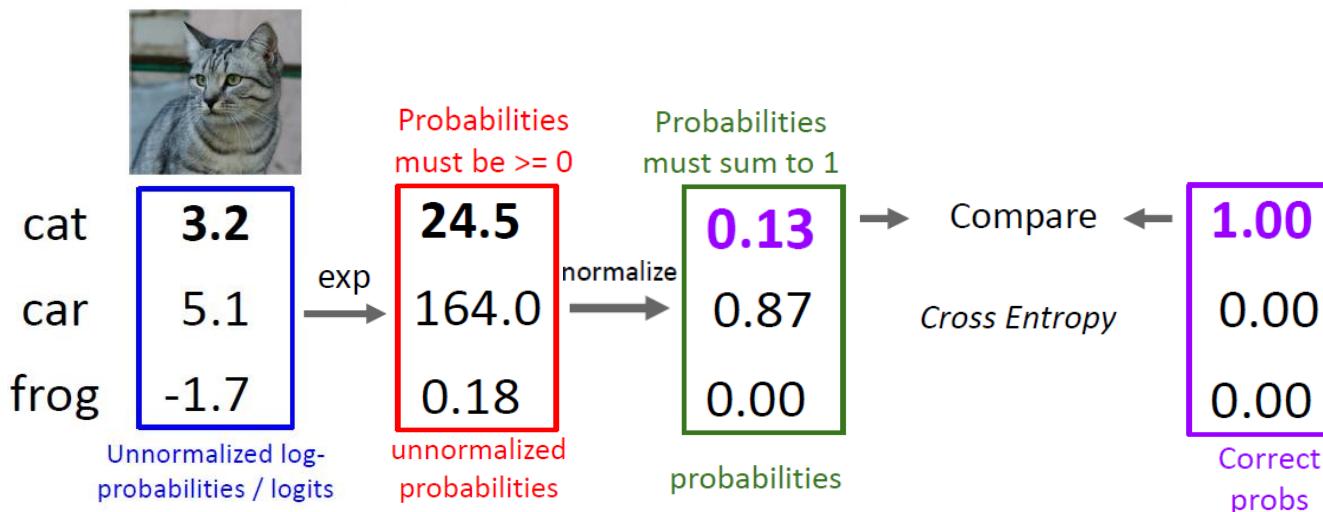
- **Cross-Entropy Loss (cont'd)**

- **Softmax function:**

$$P(Y = k \mid X = x_i) = \frac{\exp(s_k)}{\sum_j \exp(s_j)} \quad \text{with } s = f(x_i; W) \text{ as the classifier output for input } x_i$$

→  $L_i = -\log P(Y = y_i \mid X = x_i)$  or  $L_i = -\log \left( \frac{\exp(s_{y_i})}{\sum_j \exp(s_j)} \right)$

- **(Binary) Cross Entropy Loss (or  $L_{BCE}$ ; see example below):**



- **Learning  $W$  from  $L_{BCE}$**

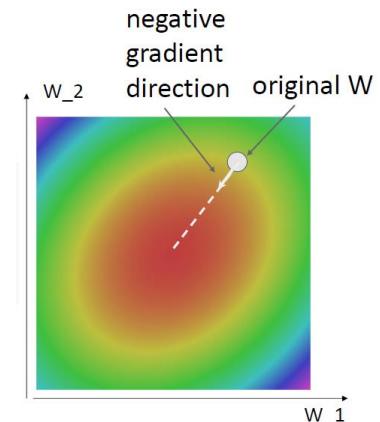
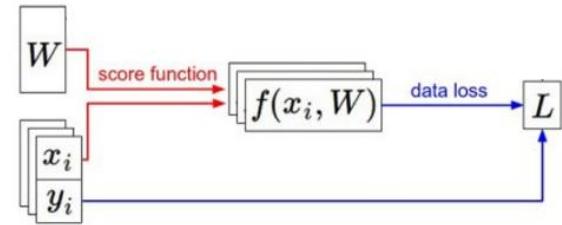
- Computing **gradients**:

Following the slope to reach the (hopefully global) minimum for  $W$ .

- **Gradient Descent** via numeric or analytic gradients:

- Iteratively step in the direction of the **negative gradient** & search for  $W$
- Hyperparameters: weight initialization, # of steps, learning rate, etc.

```
# Vanilla gradient descent
w = initialize_weights()
for t in range(num_steps):
    dw = compute_gradient(loss_fn, data, w)
    w -= learning_rate * dw
```



- **Stochastic Gradient Descent**

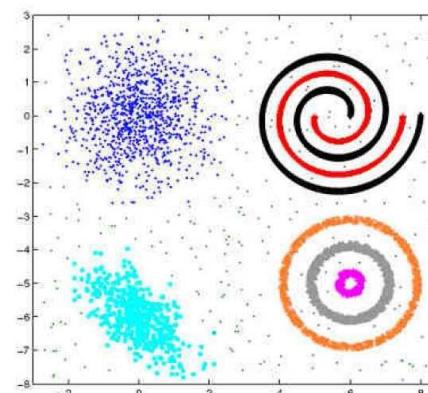
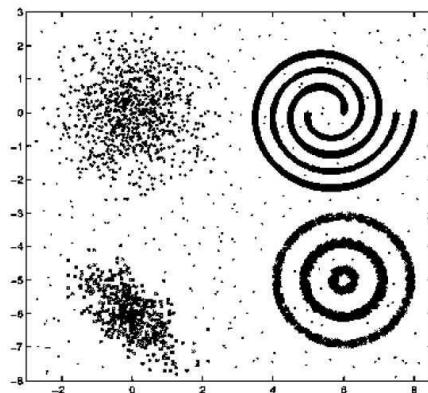
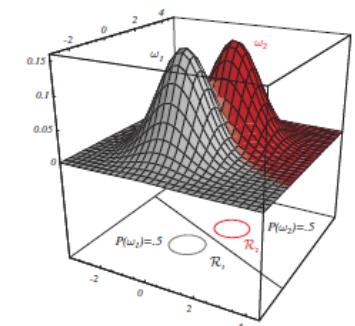
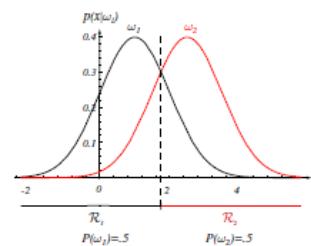
- Full sum in  $L$  is **expensive when large  $N$**
- **Approximate** sum using a minibatch of instances (e.g., 32, 64, 128, etc.)
- Additional hyperparameters of batch size and data sampling

```
# Stochastic gradient descent
w = initialize_weights()
for t in range(num_steps):
    minibatch = sample_data(data, batch_size)
    dw = compute_gradient(loss_fn, minibatch, w)
    w -= learning_rate * dw
```



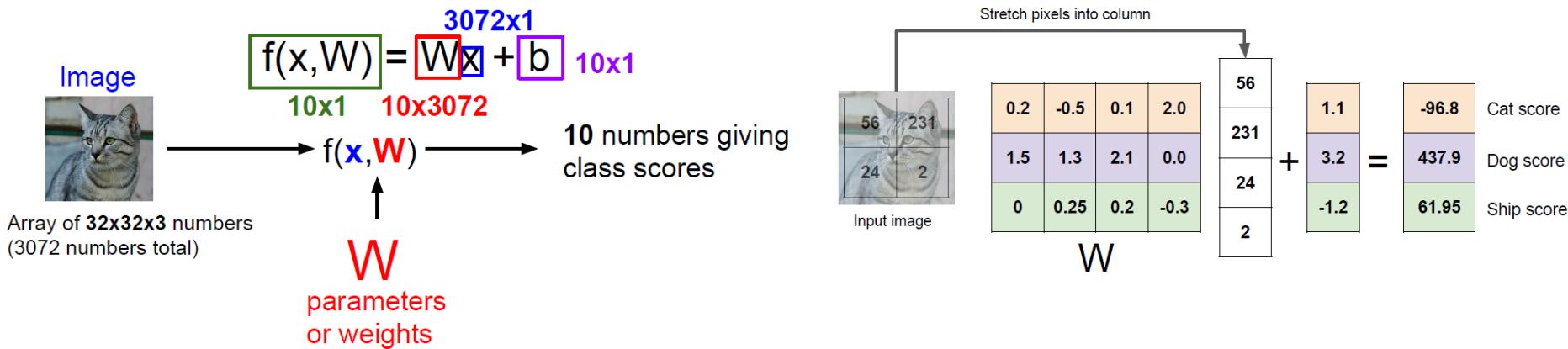
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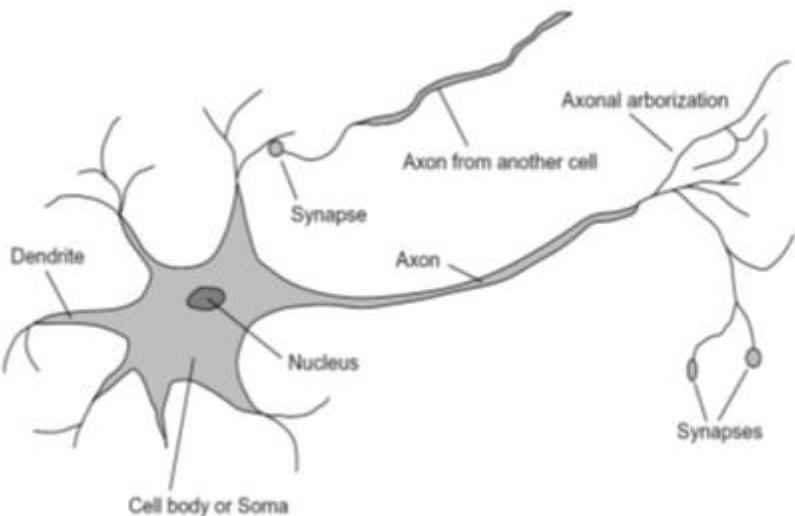


# Recap: Linear Classification

- Linear Classifier
  - Let's take the input image as  $x$ , and the linear classifier as  $W$ . We need  $y = Wx + b$  as a 10-dimensional output vector, indicating the score for each class.
  - For example, an image with  $2 \times 2$  pixels & 3 classes of interest we need to learn a linear classifier  $W$  (plus a bias  $b$ ), so that desirable outputs  $y = Wx + b$  can be expected.



# Biological neuron and Perceptrons



A biological neuron

Input

Weights

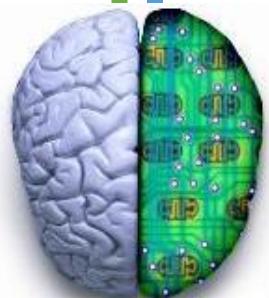
$$\begin{array}{l} x_1 \xrightarrow{w_1} \\ x_2 \xrightarrow{w_2} \\ x_3 \xrightarrow{w_3} \\ \vdots \\ \vdots \\ x_d \xrightarrow{w_d} \end{array}$$

Output:  $\sigma(w \cdot x + b)$

Sigmoid function:

$$\sigma(t) = \frac{1}{1 + e^{-t}}$$

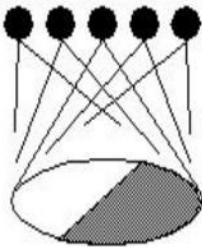
An artificial neuron (Perceptron)  
- a linear classifier



# Hubel/Wiesel Architecture and Multi-layer Neural Network

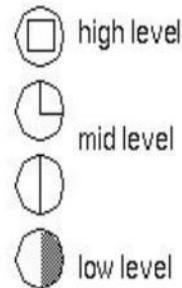
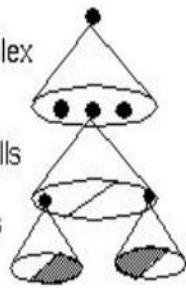
Hubel & Weisel

topographical mapping

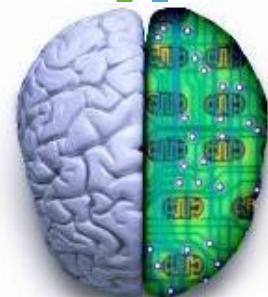


featural hierarchy

hyper-complex  
cells  
complex cells  
simple cells



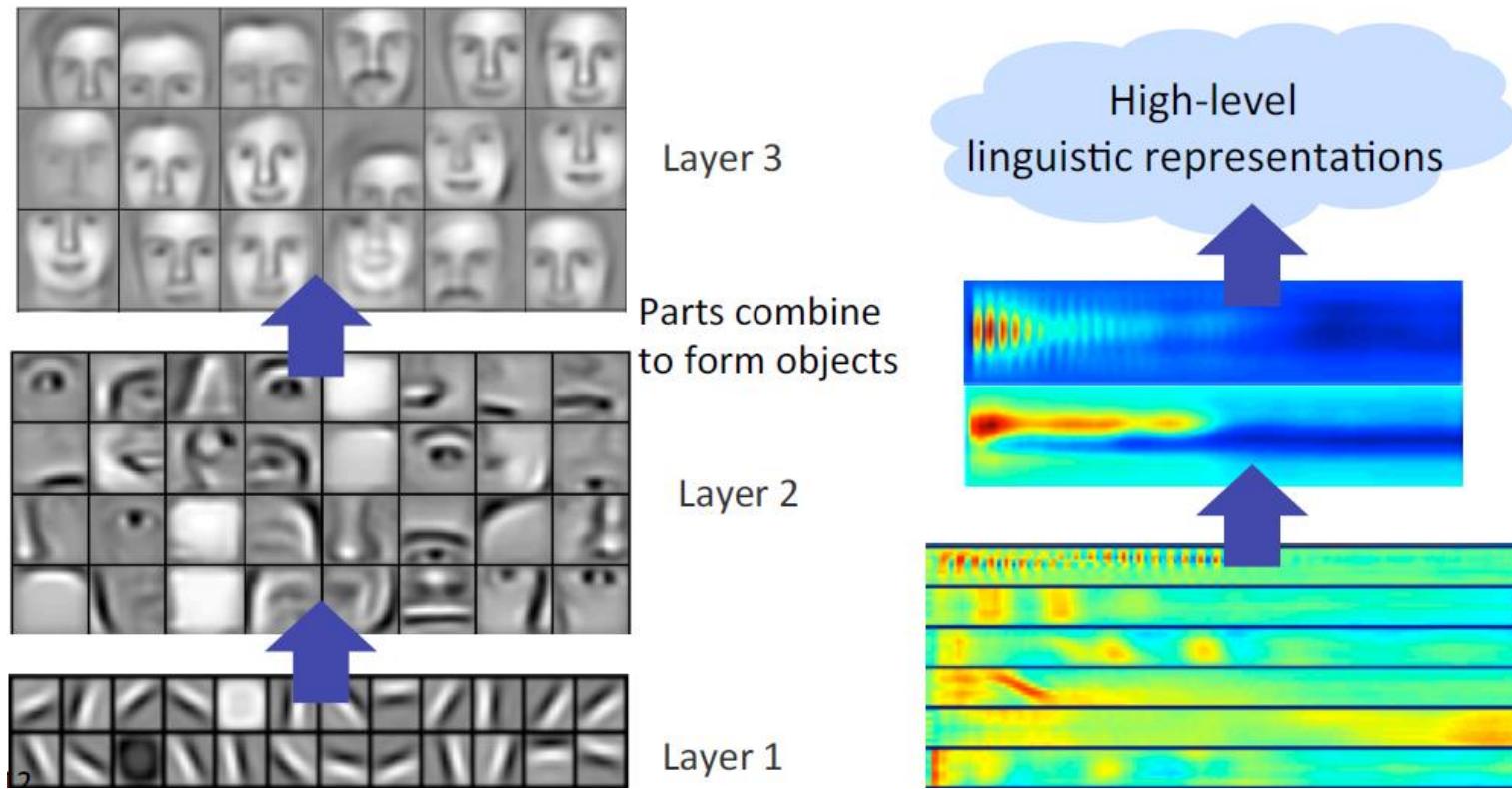
Hubel and Weisel's architecture



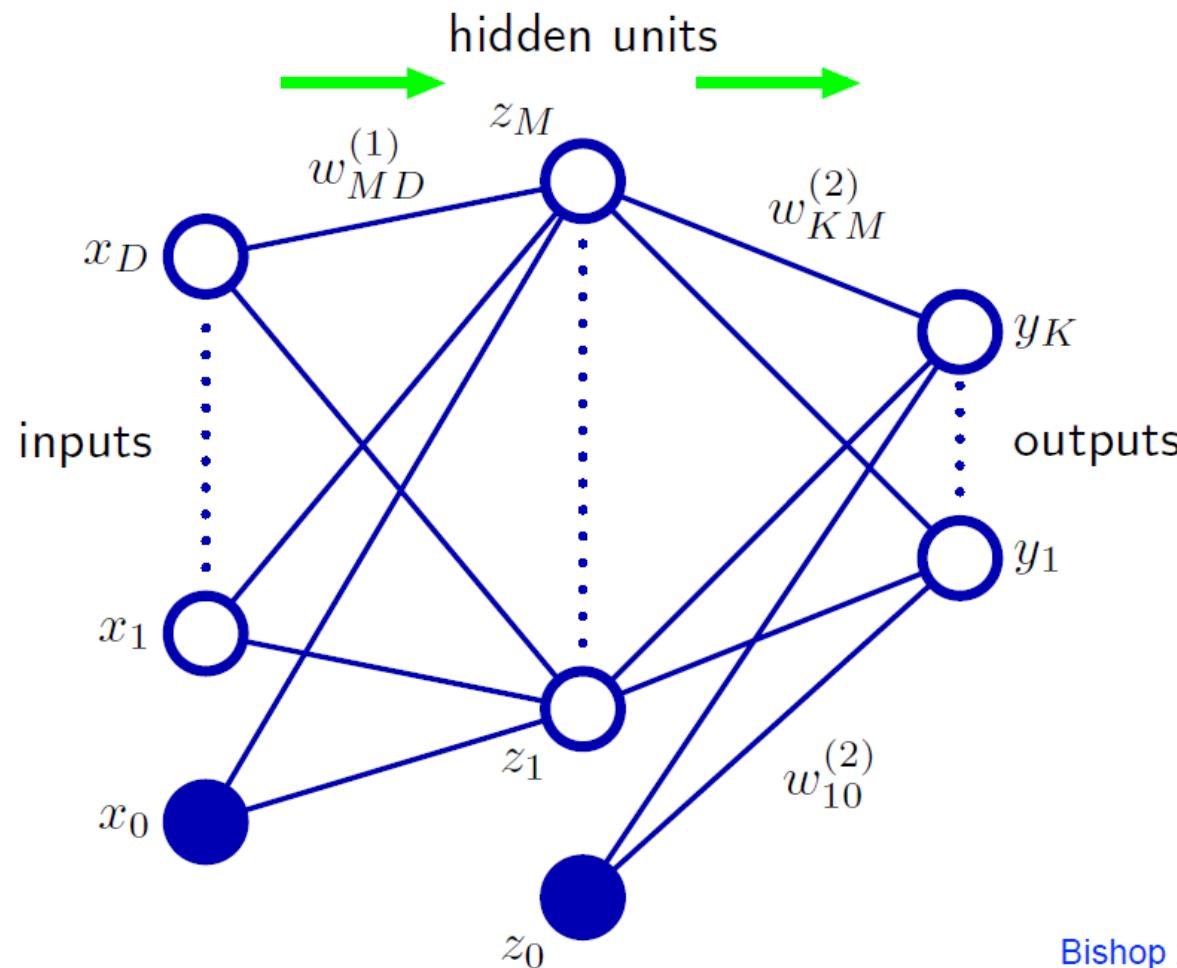
**Multi-Layer Perception or Neural Network**  
- A *non-linear* classifier

# Hierarchical Representation Learning

- Successive model layers learn deeper intermediate representations.

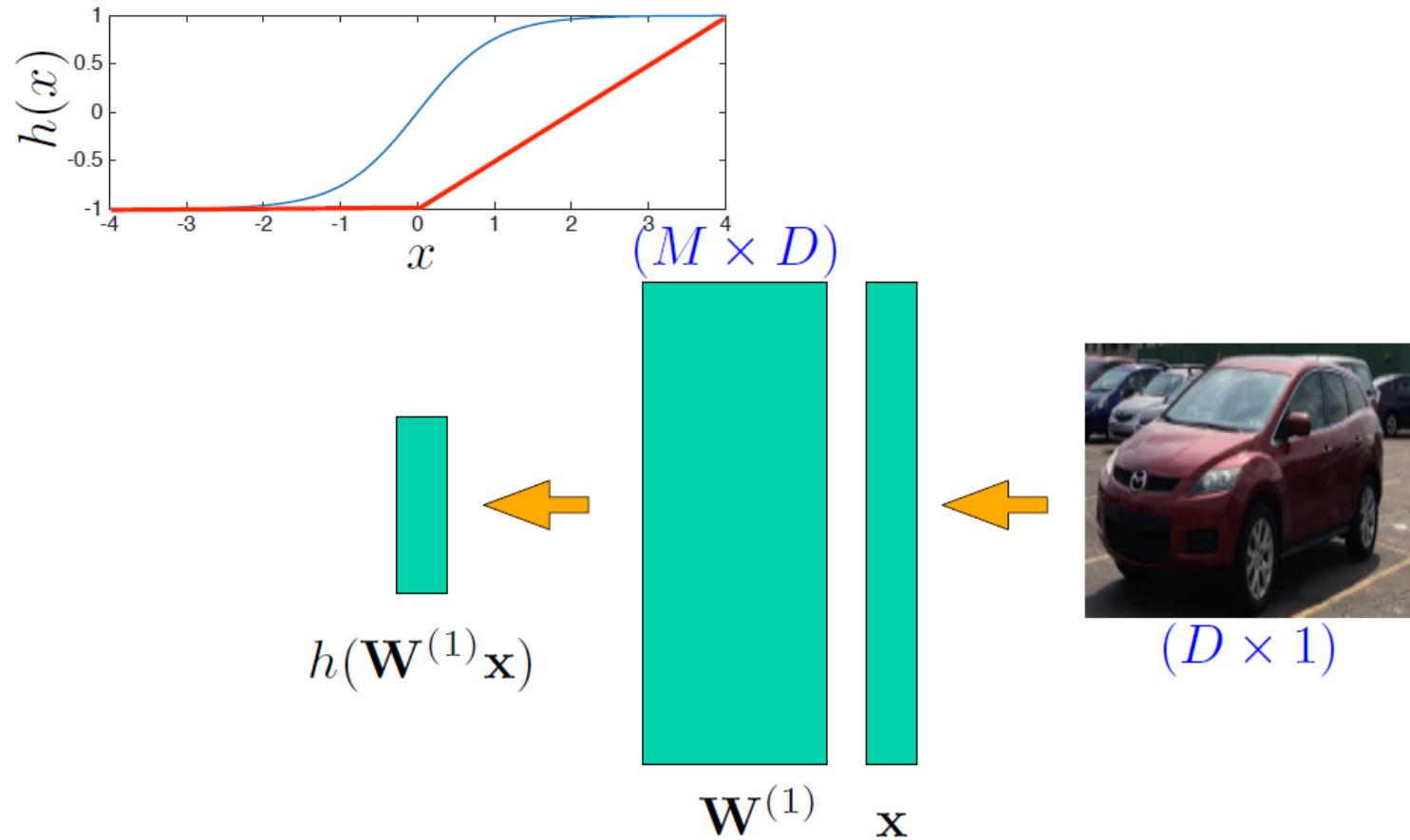


# Multi-Layer Perceptron: A Nonlinear Classifier (cont'd)



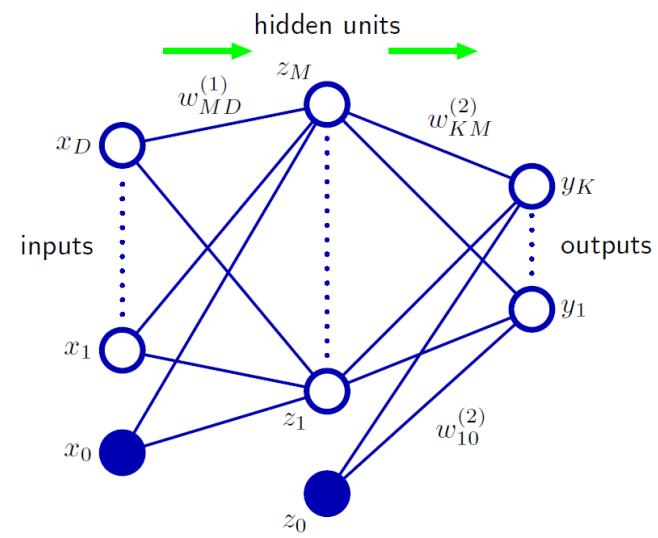
Bishop 2006

# Multi-Layer Perceptron: A Nonlinear Classifier



# Layer 1 in MLP

$$\mathbf{z} = \begin{bmatrix} z_1 \\ \vdots \\ z_M \end{bmatrix} \leftarrow \begin{bmatrix} h[\mathbf{x}^T \mathbf{w}_1^{(1)}] \\ \vdots \\ h[\mathbf{x}^T \mathbf{w}_M^{(1)}] \end{bmatrix}$$

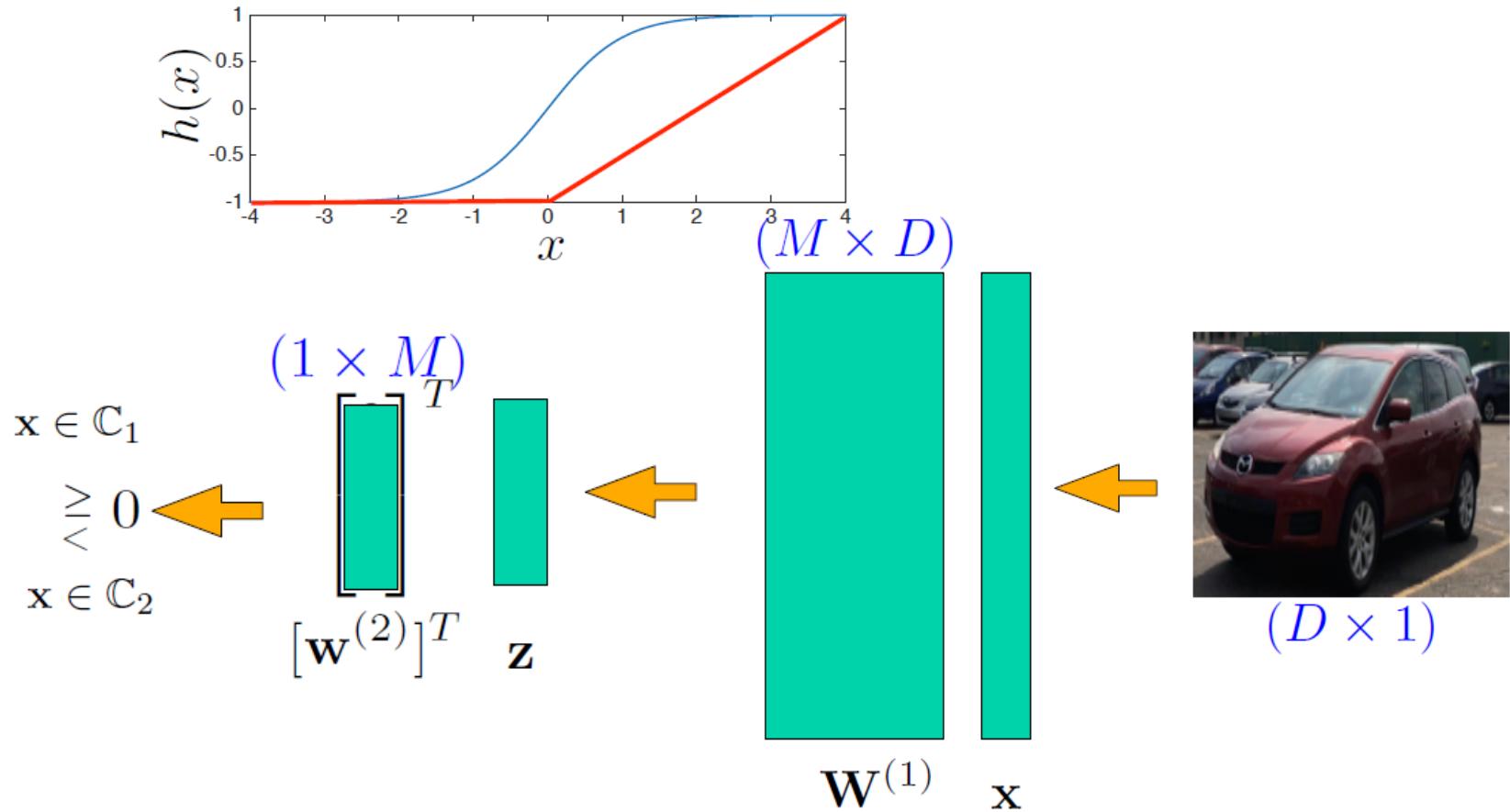


$h()$  = non-linear function

$[\mathbf{w}_1^{(1)}, \dots, \mathbf{w}_M^{(1)}]$  = 1st layer's  $D \times M$  weights

$\mathbf{x} = D \times 1$  raw input

# Multi-Layer Perceptron: A Nonlinear Classifier (cont'd)



# Layer 2 in MLP

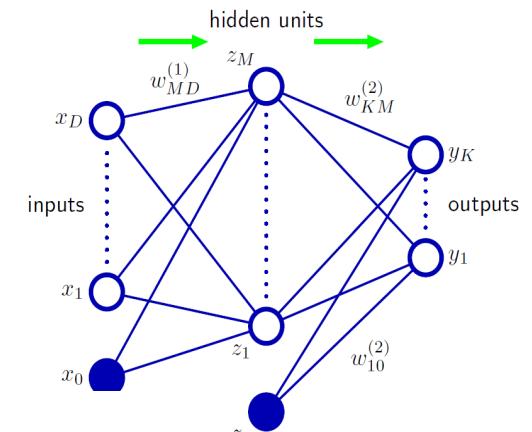


$$\mathbf{x} \in \mathbb{R}^D$$

$$\mathbf{z} \in \mathbb{C}_1$$

$$\mathbf{z}^T \mathbf{w}^{(2)} \geq 0$$

$$\mathbf{z} \in \mathbb{C}_2$$

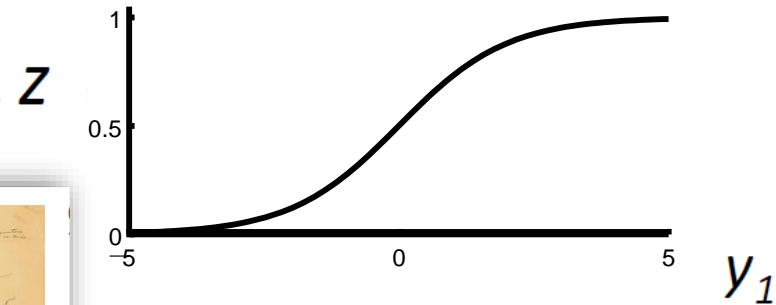
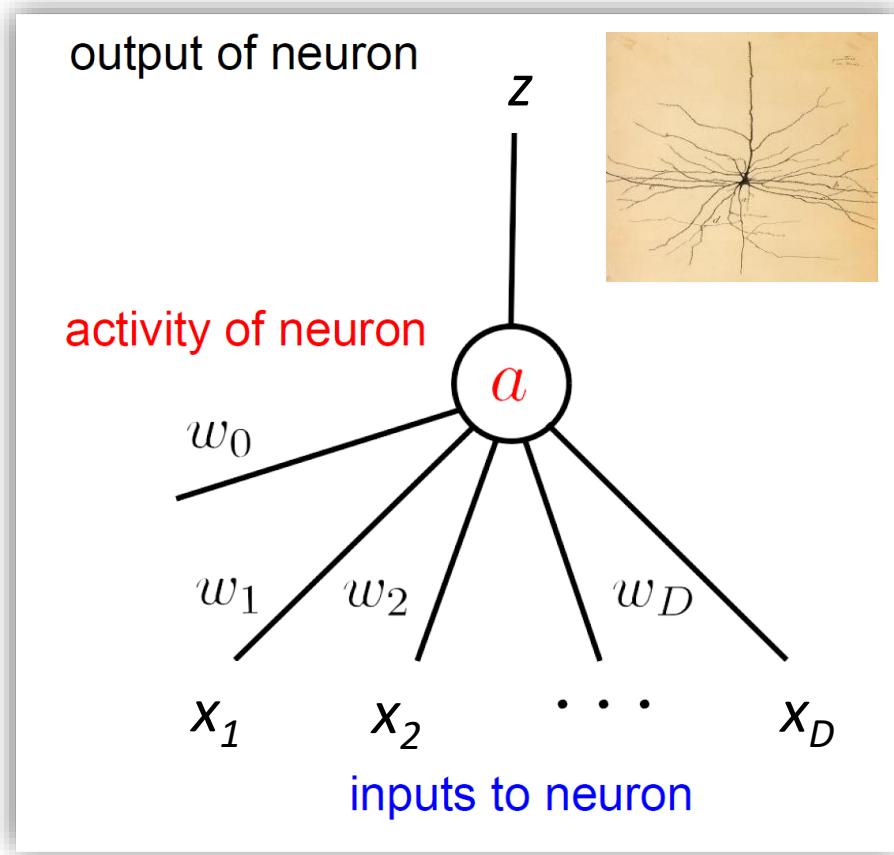


$\mathbf{z} = M \times 1$  output of layer 1

$\mathbf{w}^{(2)} = 2\text{nd layer's } M \times 1 \text{ weight vector}$

# Let's Take a Closer Look... Output $y = Wx + b$

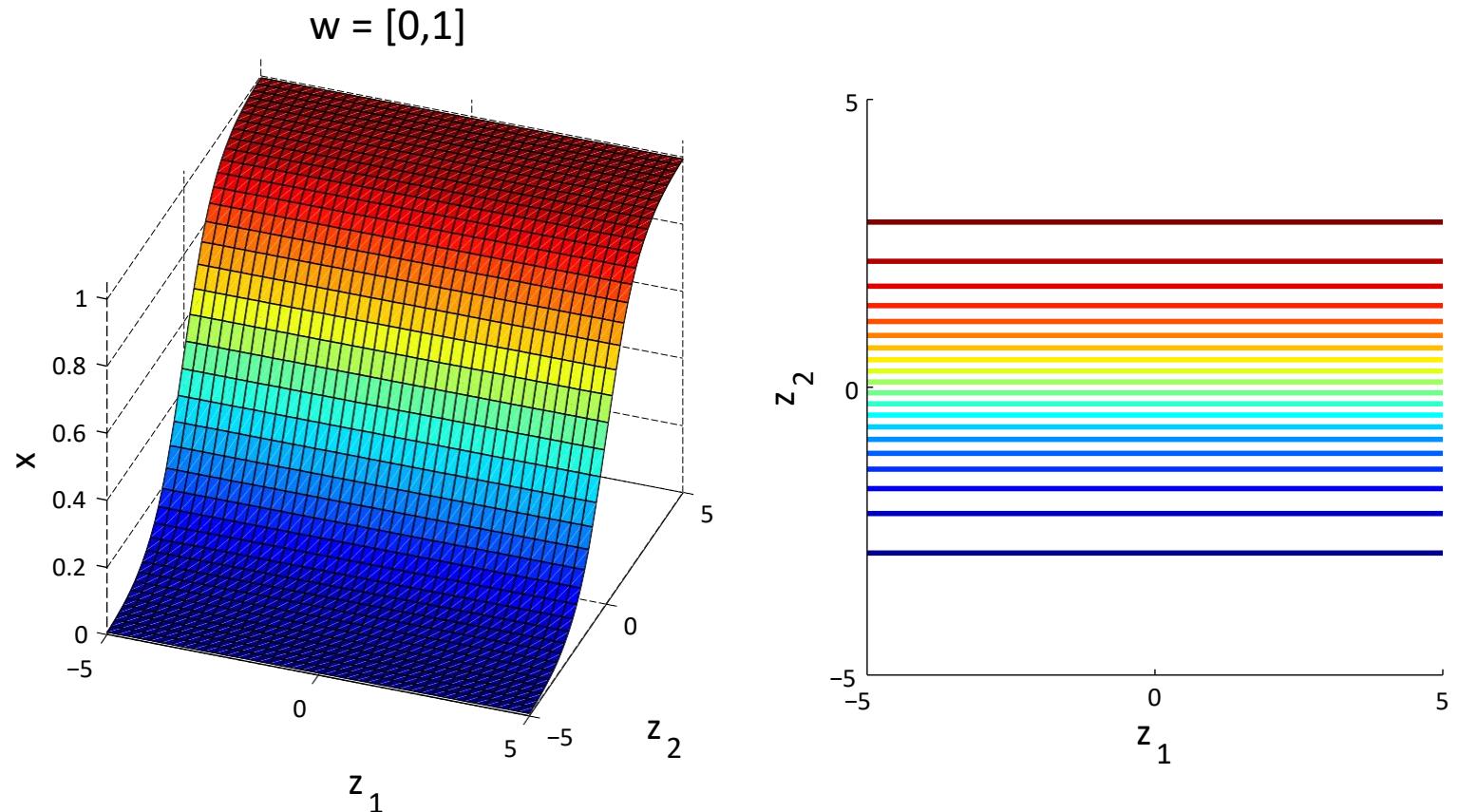
- A single neuron



$$z(y_1) = \frac{1}{1 + \exp(-y_1)}, z \in (0,1)$$

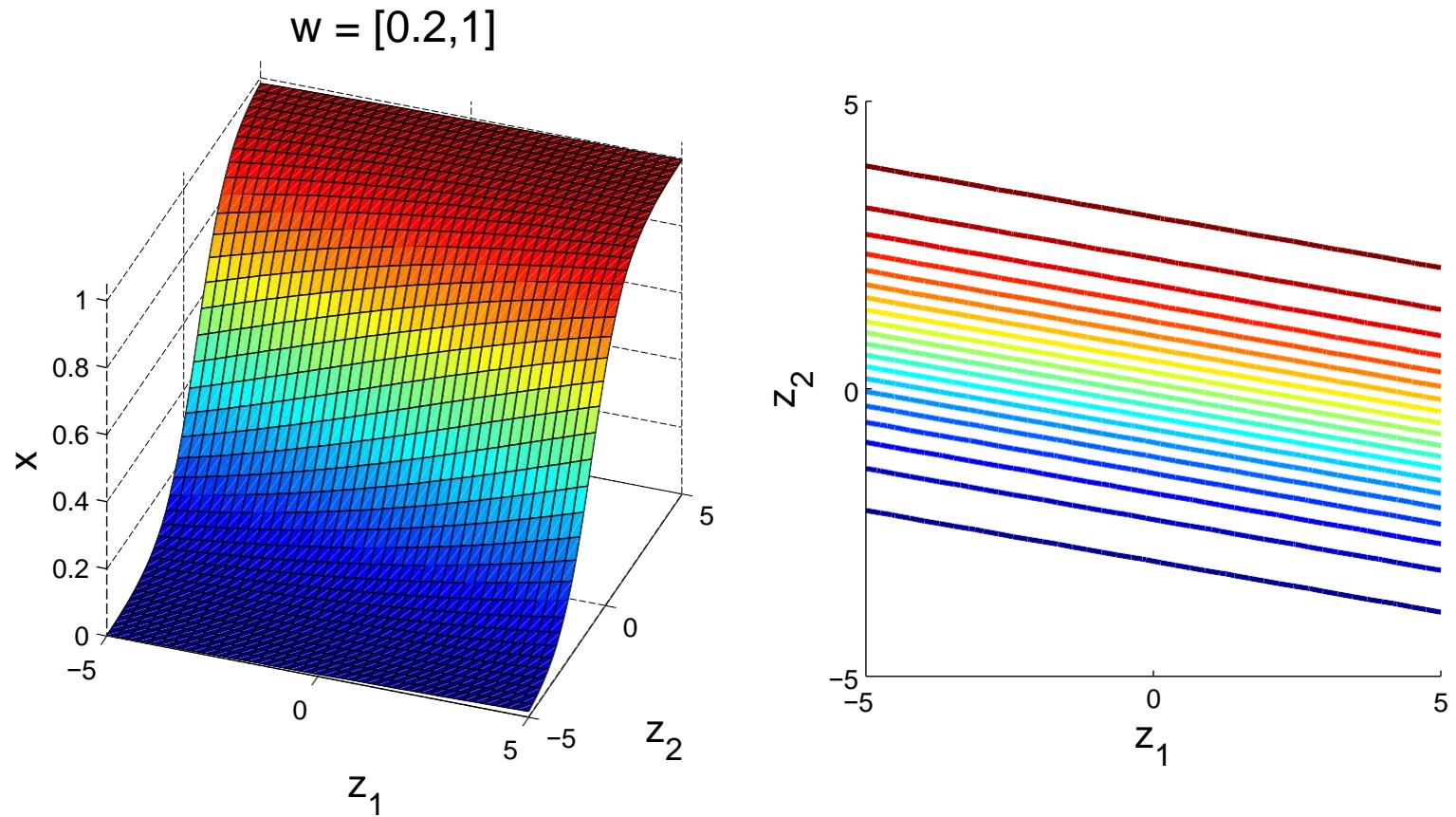
$$\begin{aligned}y_1 &= w_0 + \sum_{d=1}^D w_d z_d \\&= \sum_{d=0}^D w_d z_d\end{aligned}$$

# Input-Output Function of a Single Neuron



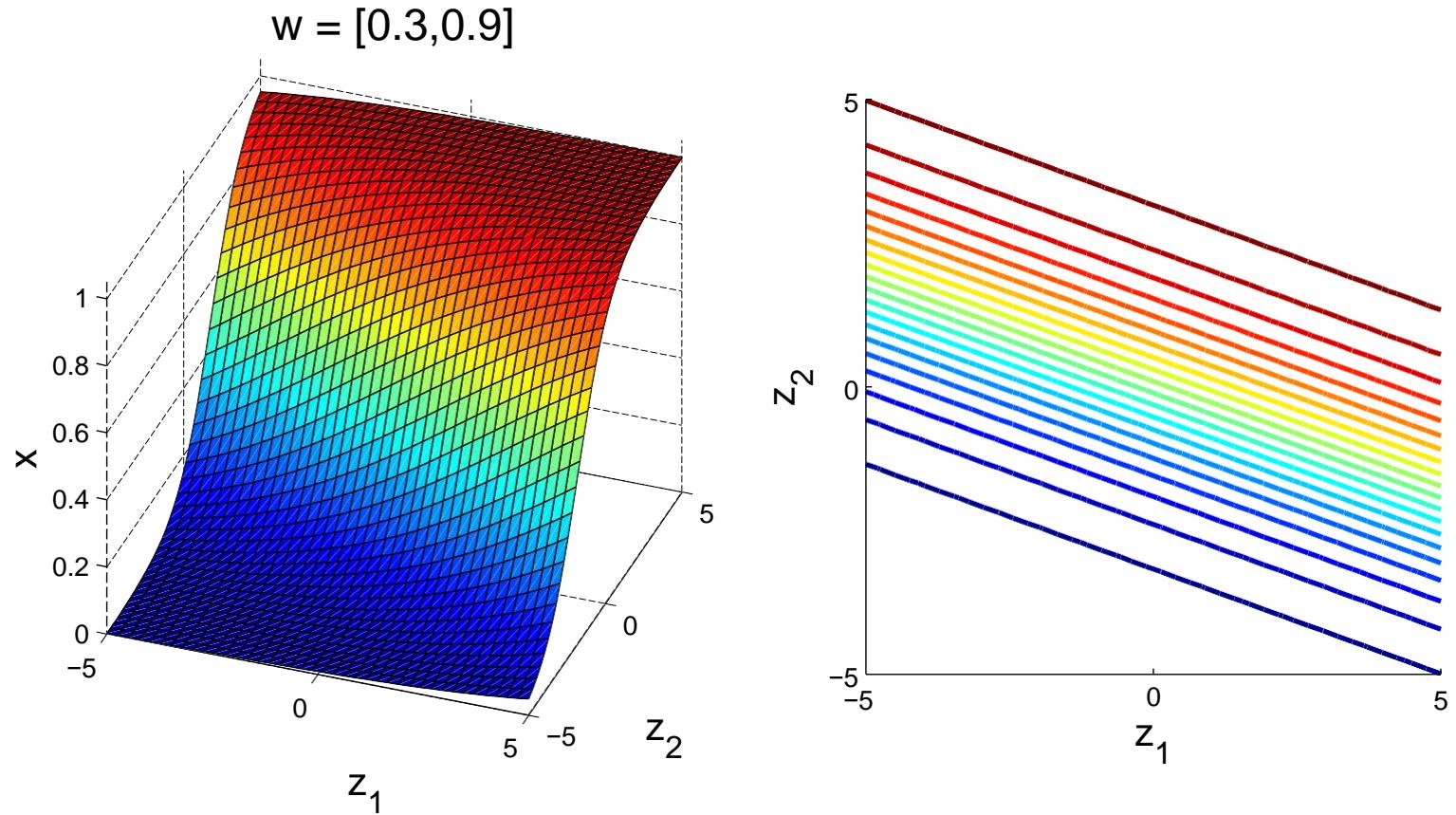
$$x(z_1, z_2) = \frac{1}{1 + \exp(-w_1 z_1 - w_2 z_2)}$$

# Input-Output Function of a Single Neuron



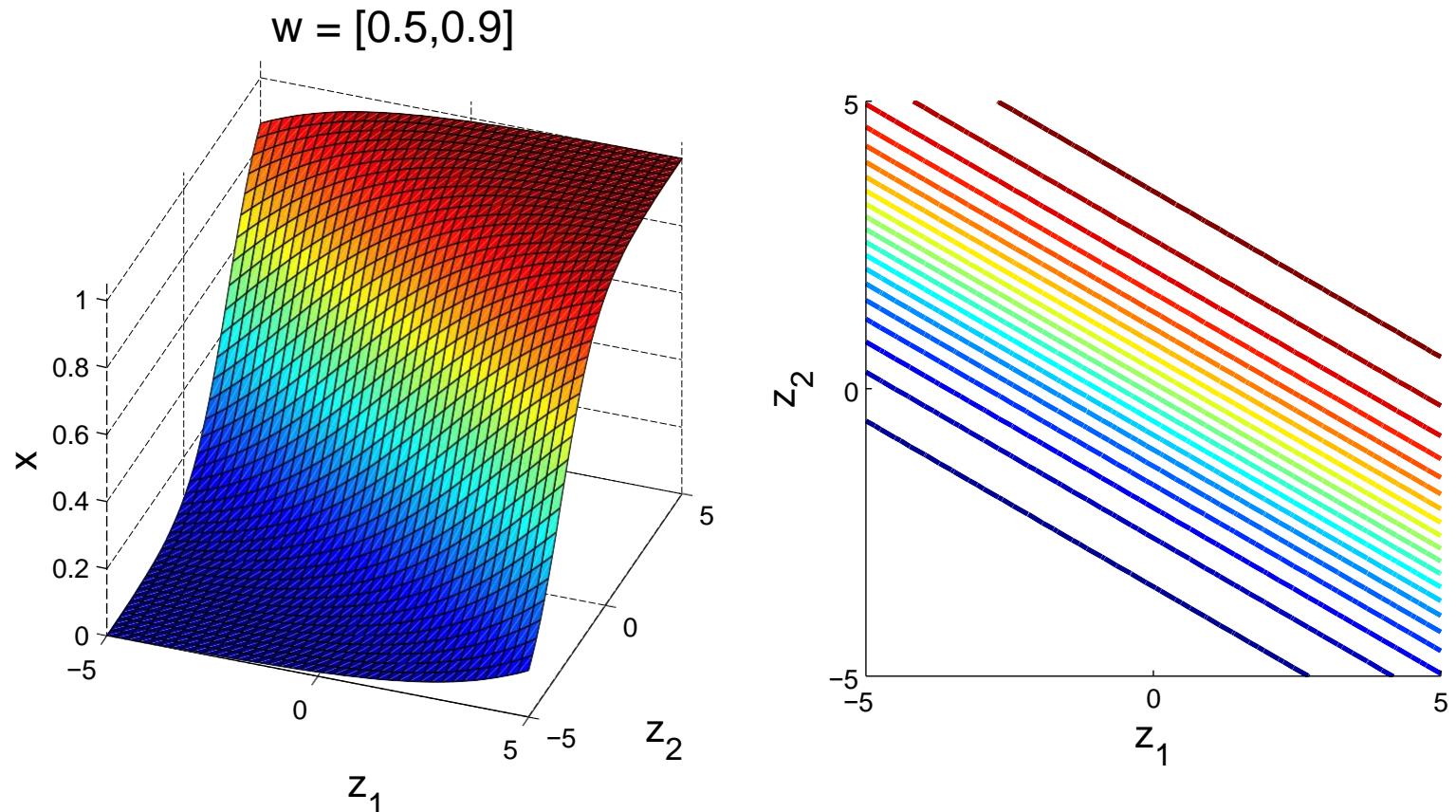
$$x(z_1, z_2) = \frac{1}{1 + \exp(-w_1 z_1 - w_2 z_2)}$$

# Input-Output Function of a Single Neuron



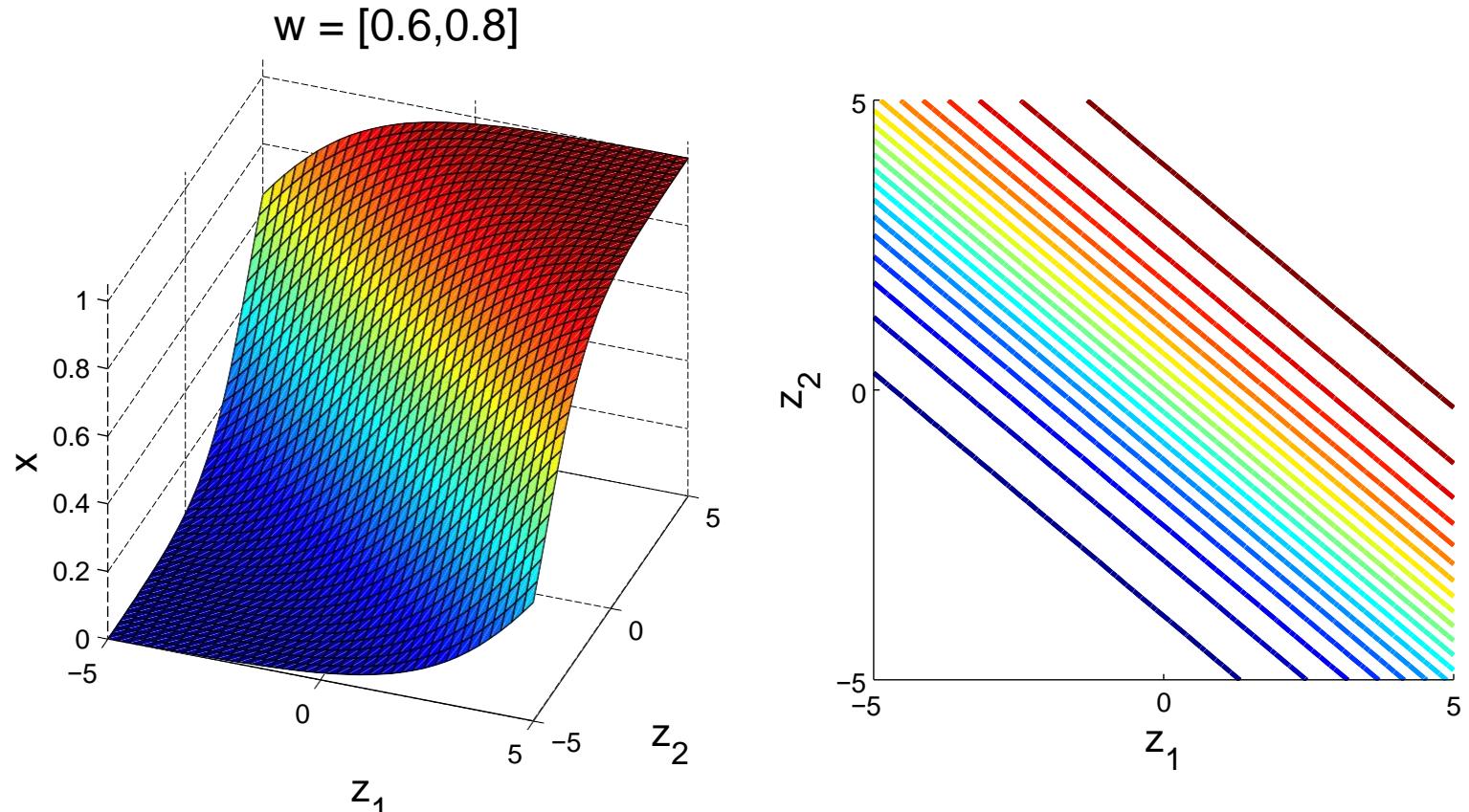
$$x(z_1, z_2) = \frac{1}{1 + \exp(-w_1 z_1 - w_2 z_2)}$$

# Input-Output Function of a Single Neuron



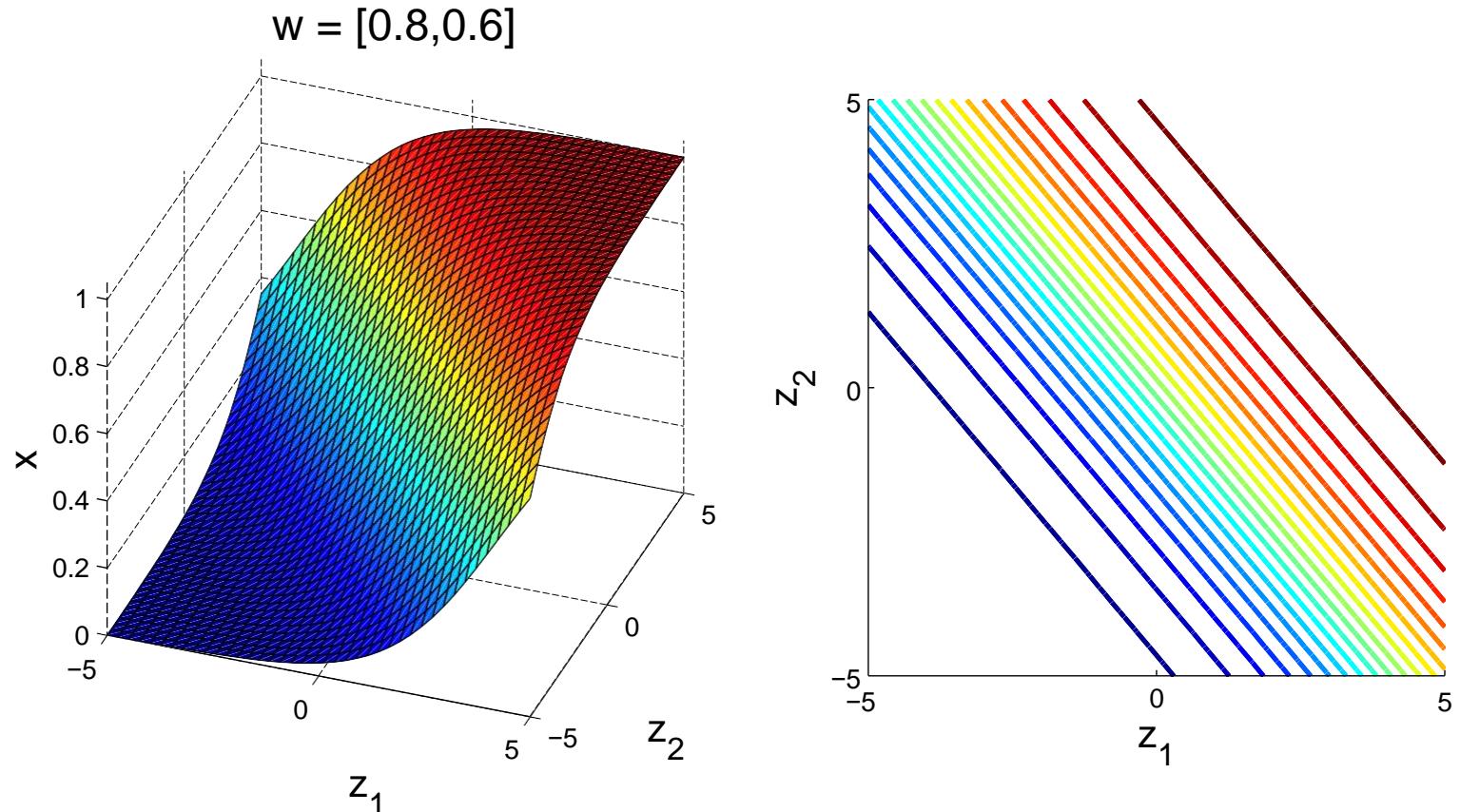
$$x(z_1, z_2) = \frac{1}{1 + \exp(-w_1 z_1 - w_2 z_2)}$$

# Input-Output Function of a Single Neuron



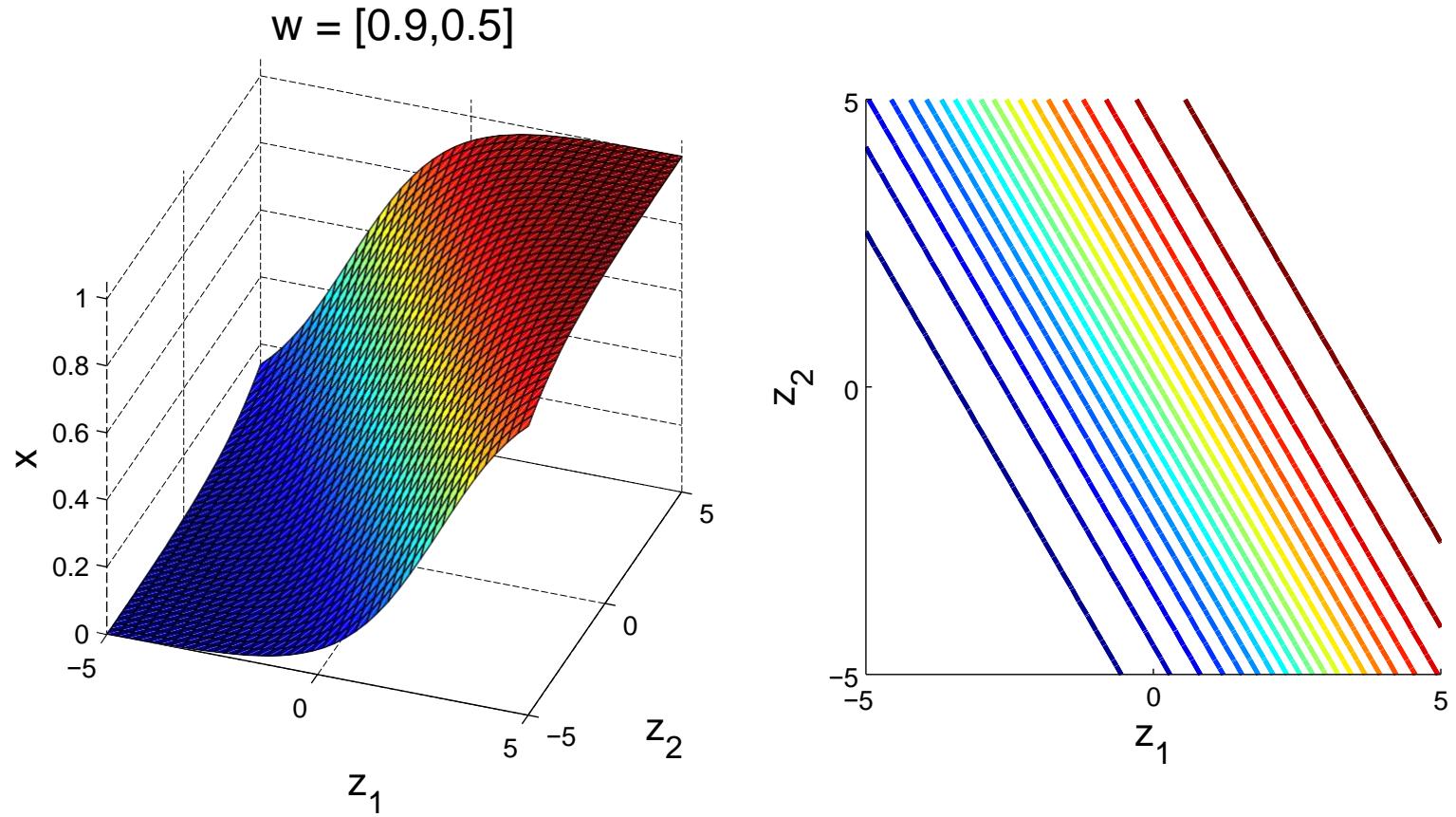
$$x(z_1, z_2) = \frac{1}{1 + \exp(-w_1 z_1 - w_2 z_2)}$$

# Input-Output Function of a Single Neuron



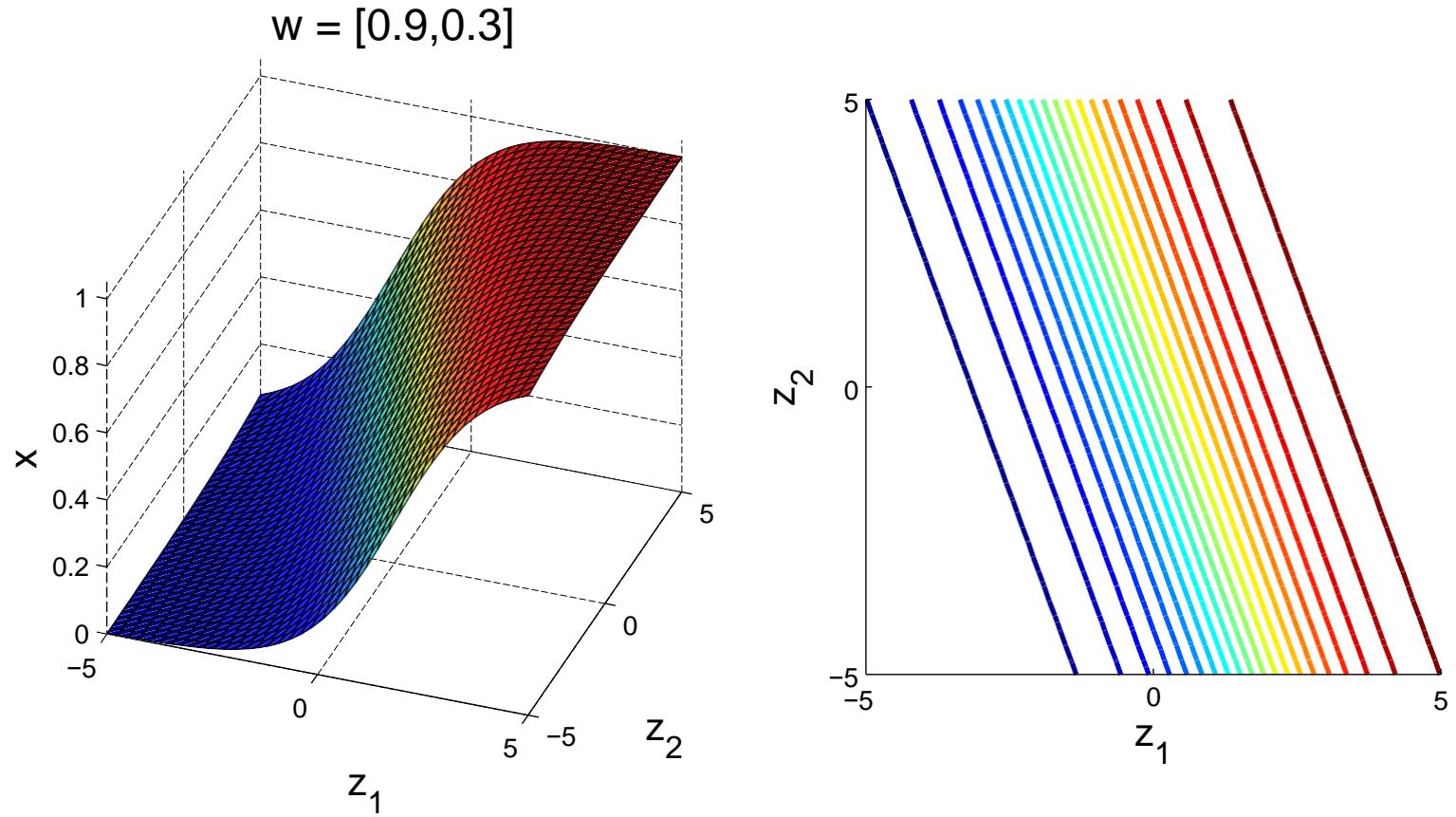
$$x(z_1, z_2) = \frac{1}{1 + \exp(-w_1 z_1 - w_2 z_2)}$$

# Input-Output Function of a Single Neuron



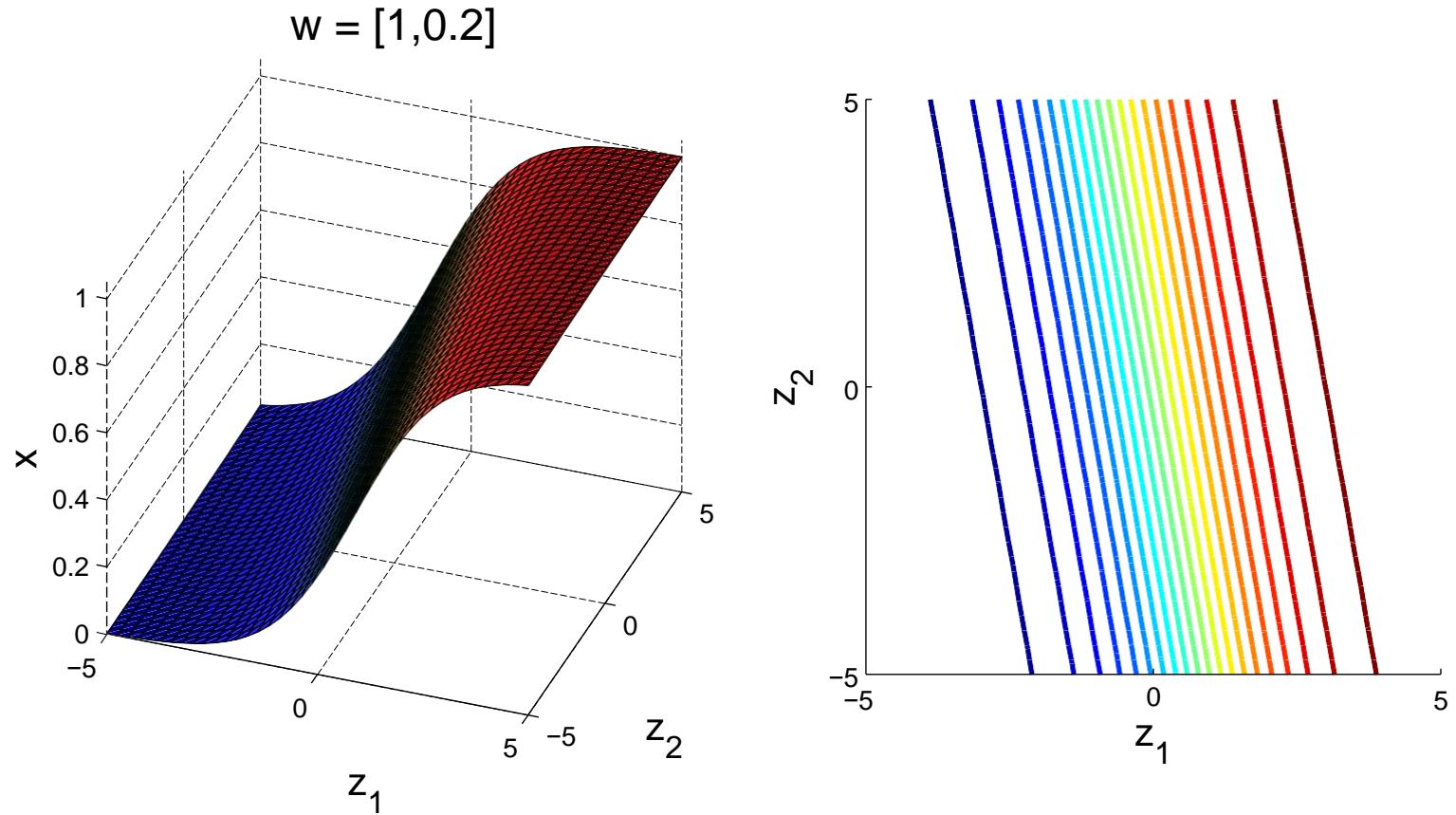
$$x(z_1, z_2) = \frac{1}{1 + \exp(-w_1 z_1 - w_2 z_2)}$$

# Input-Output Function of a Single Neuron



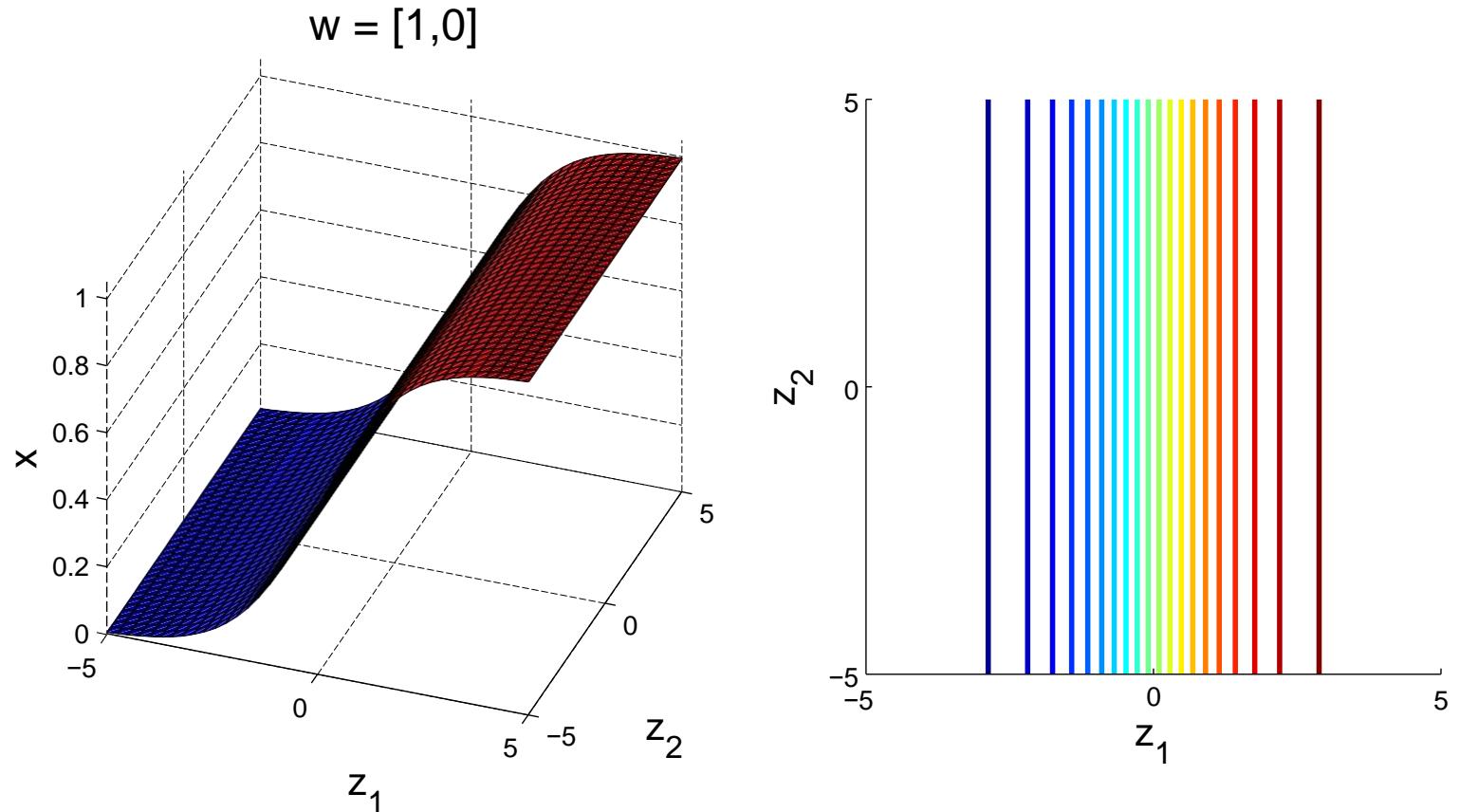
$$x(z_1, z_2) = \frac{1}{1 + \exp(-w_1 z_1 - w_2 z_2)}$$

# Input-Output Function of a Single Neuron



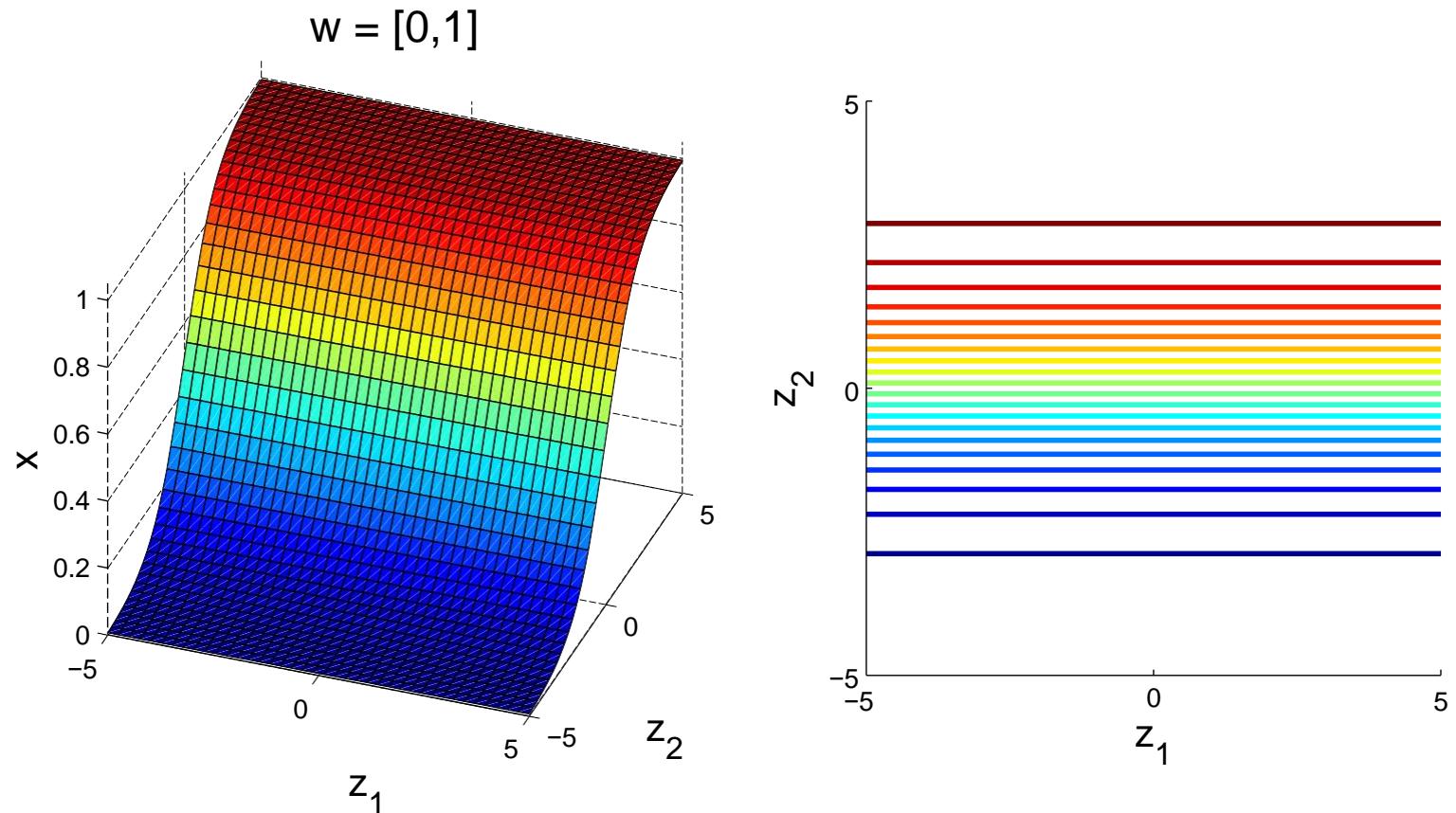
$$x(z_1, z_2) = \frac{1}{1 + \exp(-w_1 z_1 - w_2 z_2)}$$

# Input-Output Function of a Single Neuron



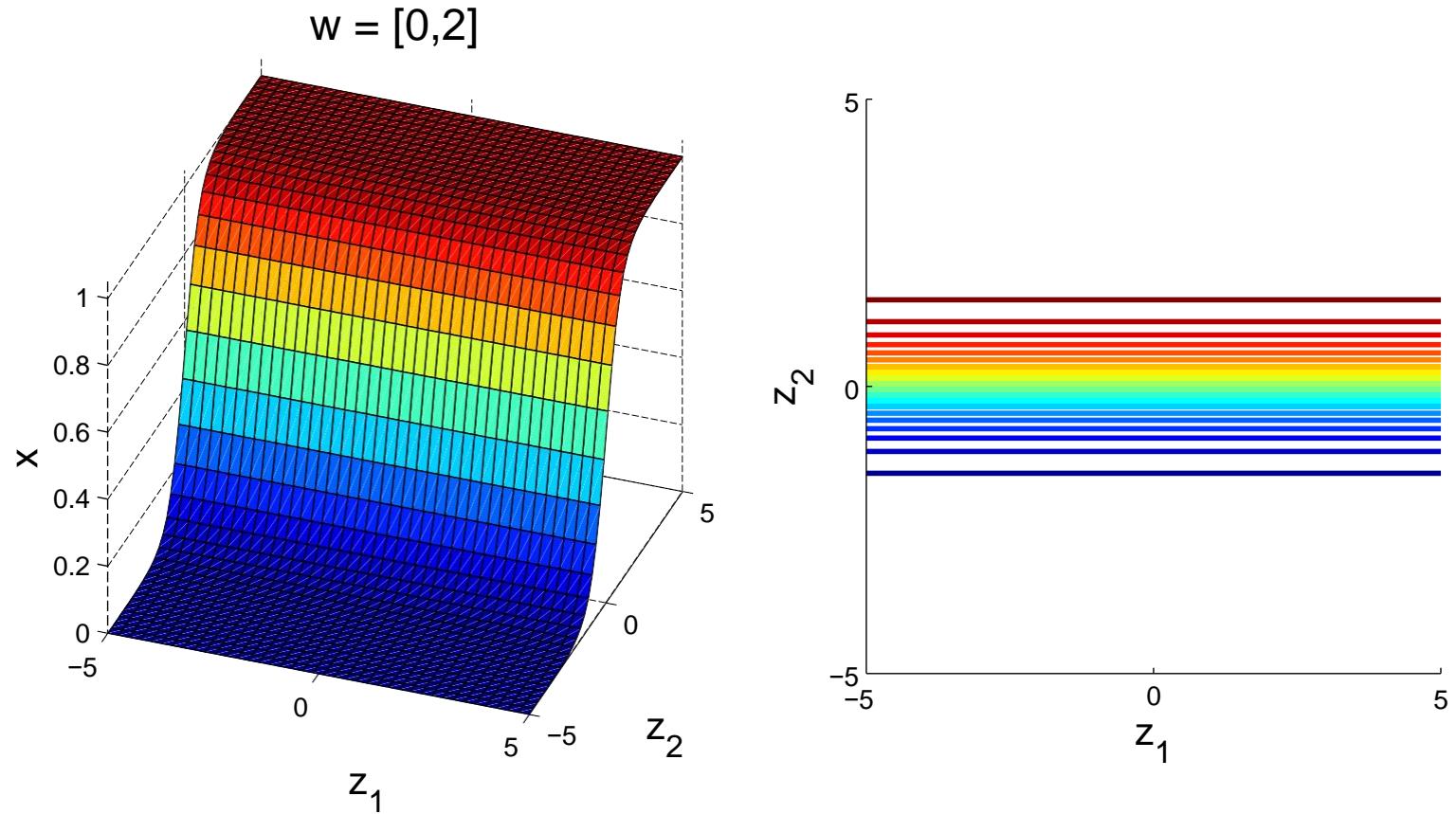
$$x(z_1, z_2) = \frac{1}{1 + \exp(-w_1 z_1 - w_2 z_2)}$$

# Input-Output Function of a Single Neuron (cont'd)



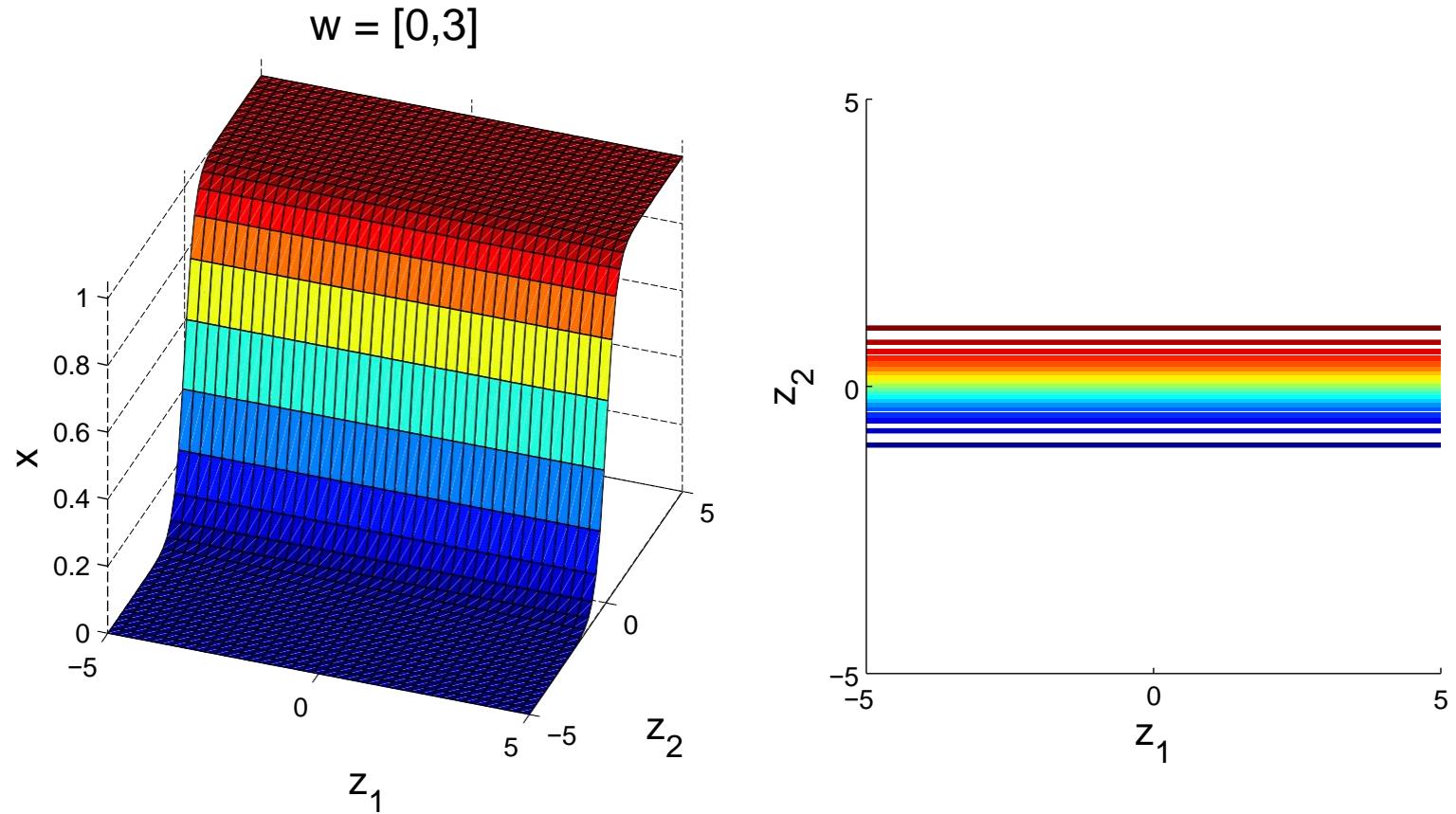
$$x(z_1, z_2) = \frac{1}{1 + \exp(-w_1 z_1 - w_2 z_2)}$$

# Input-Output Function of a Single Neuron (cont'd)



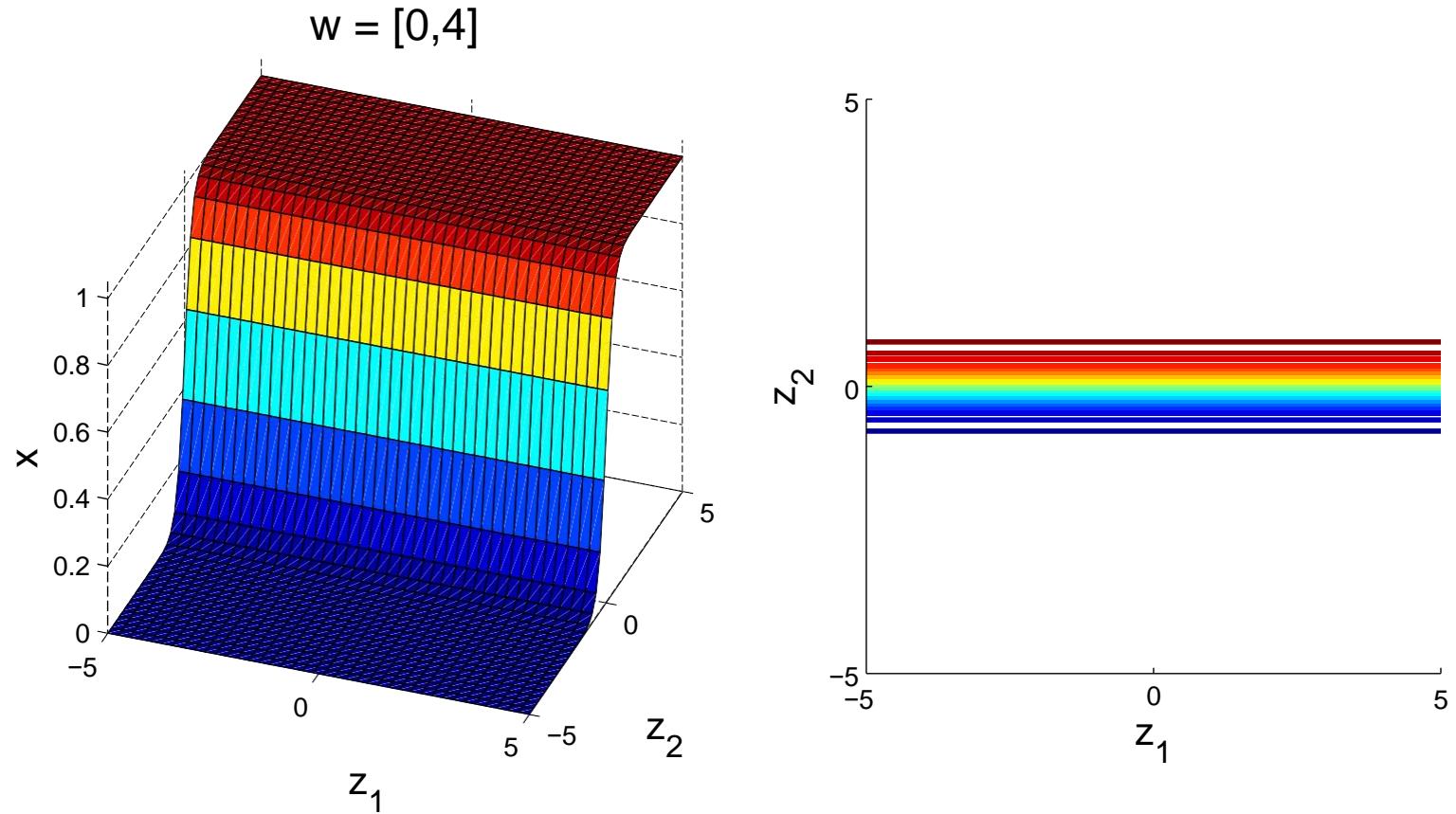
$$x(z_1, z_2) = \frac{1}{1 + \exp(-w_1 z_1 - w_2 z_2)}$$

# Input-Output Function of a Single Neuron (cont'd)



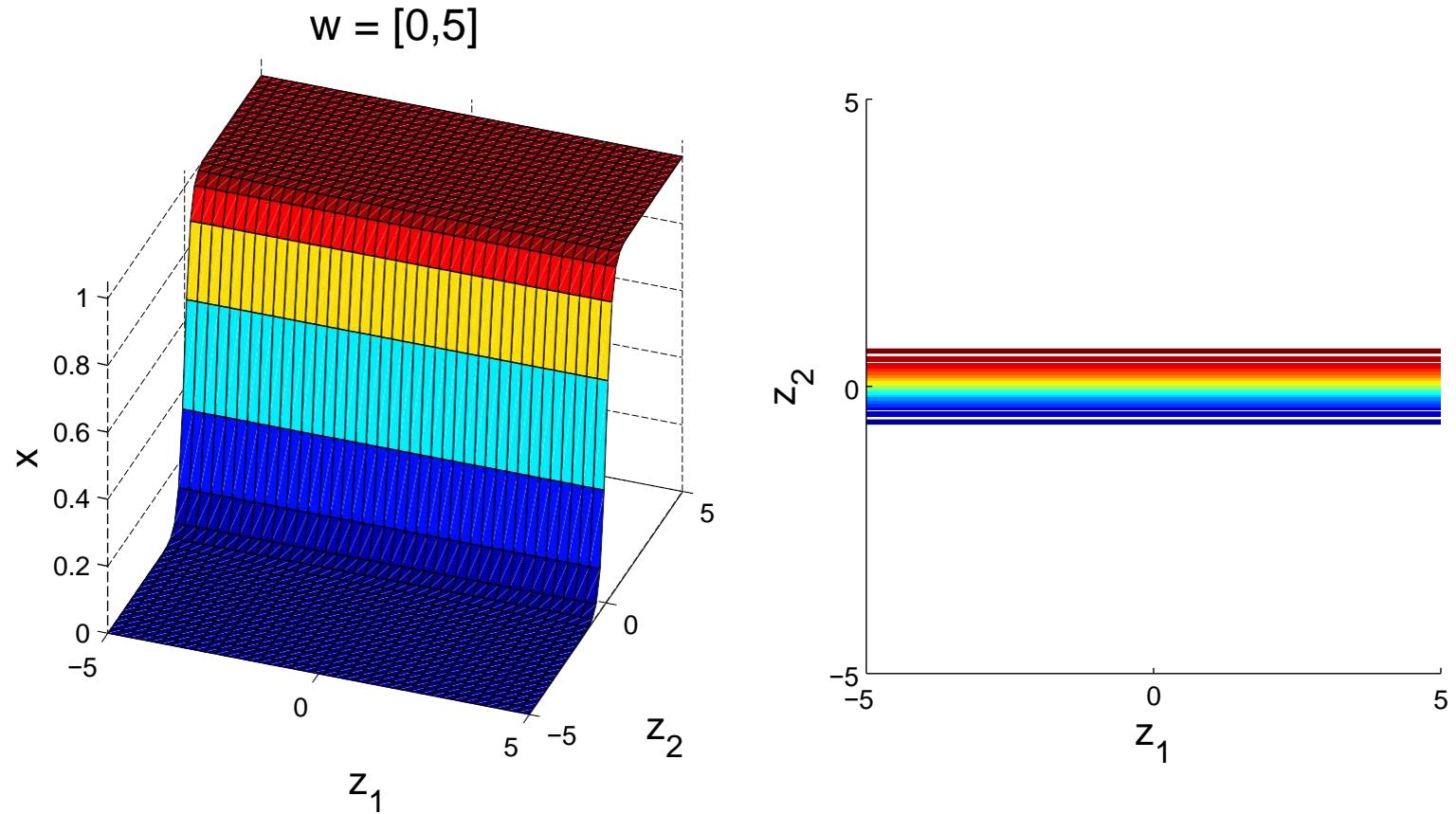
$$x(z_1, z_2) = \frac{1}{1+\exp(-w_1 z_1 - w_2 z_2)}$$

# Input-Output Function of a Single Neuron (cont'd)



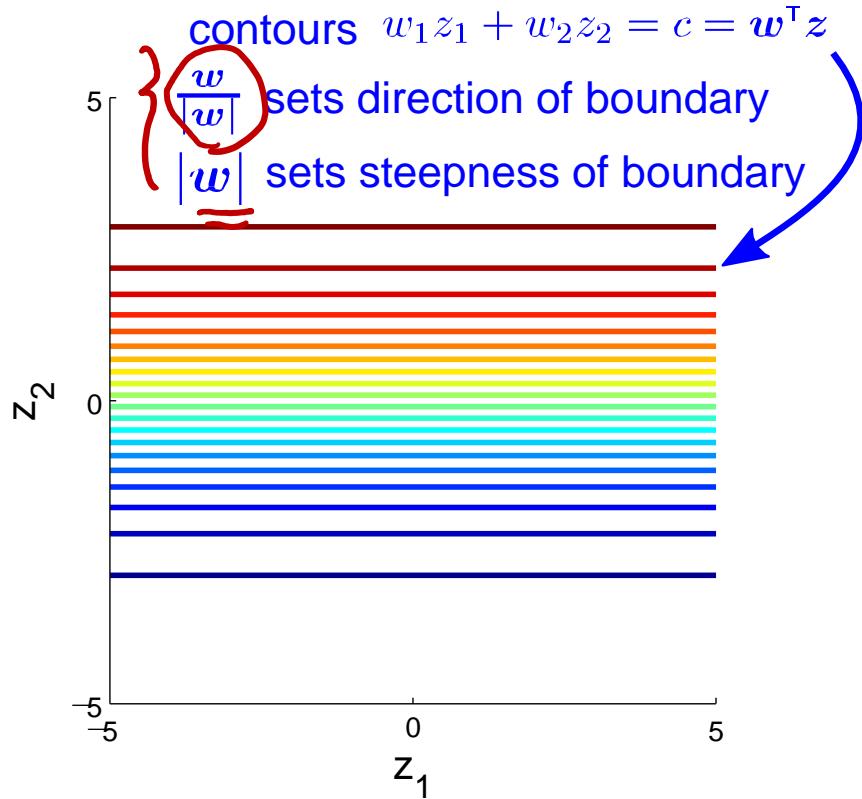
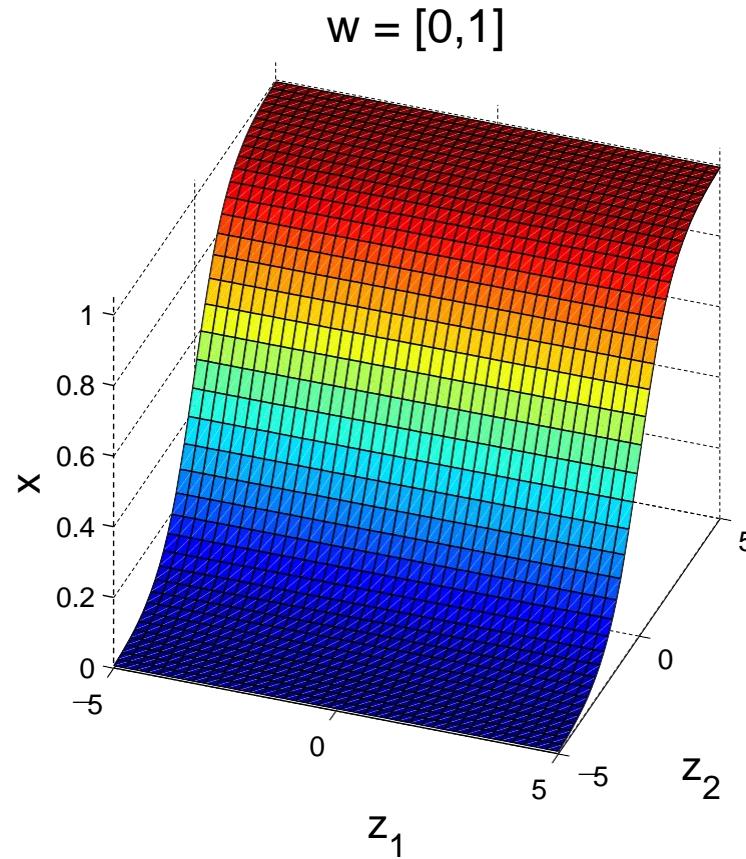
$$x(z_1, z_2) = \frac{1}{1+\exp(-w_1 z_1 - w_2 z_2)}$$

# Input-Output Function of a Single Neuron (cont'd)



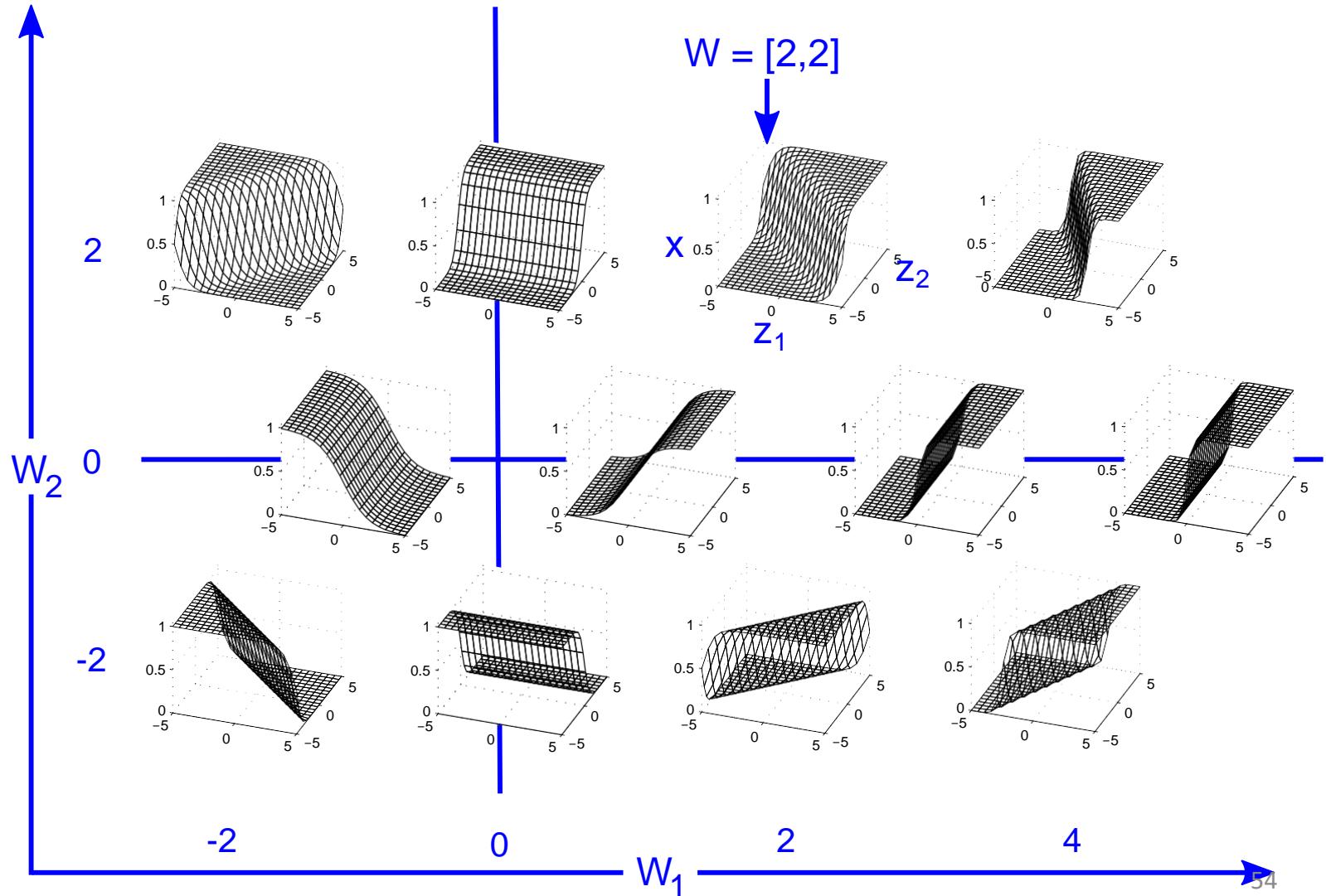
$$x(z_1, z_2) = \frac{1}{1+\exp(-w_1 z_1 - w_2 z_2)}$$

# Input-Output Function of a Single Neuron (cont'd)

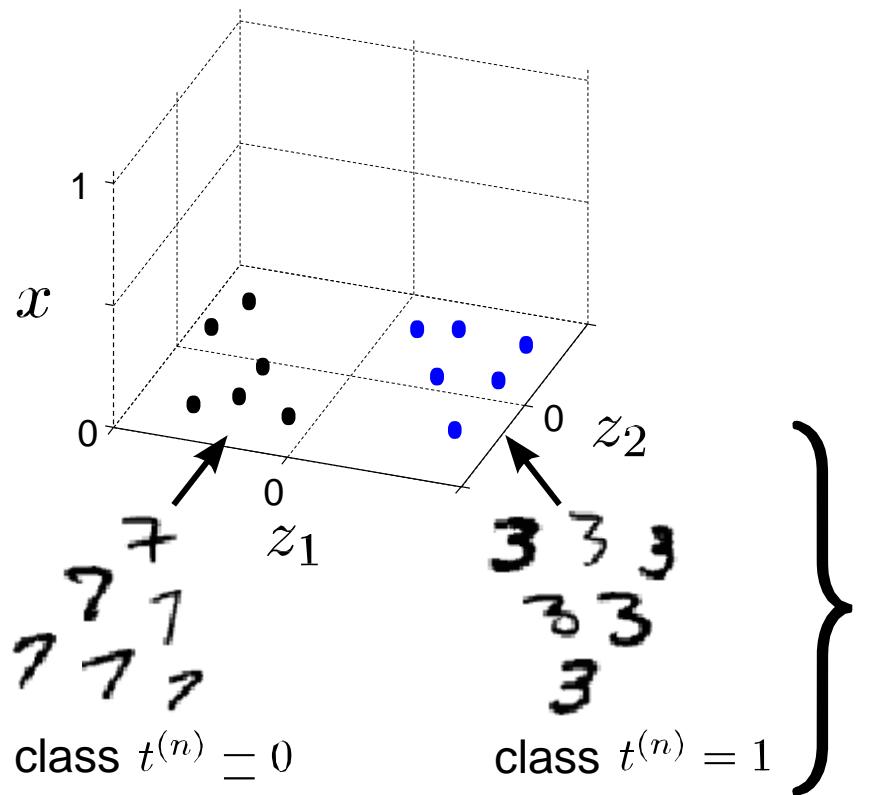


$$x(z_1, z_2) = \frac{1}{1 + \exp(-w_1 z_1 - w_2 z_2)}$$

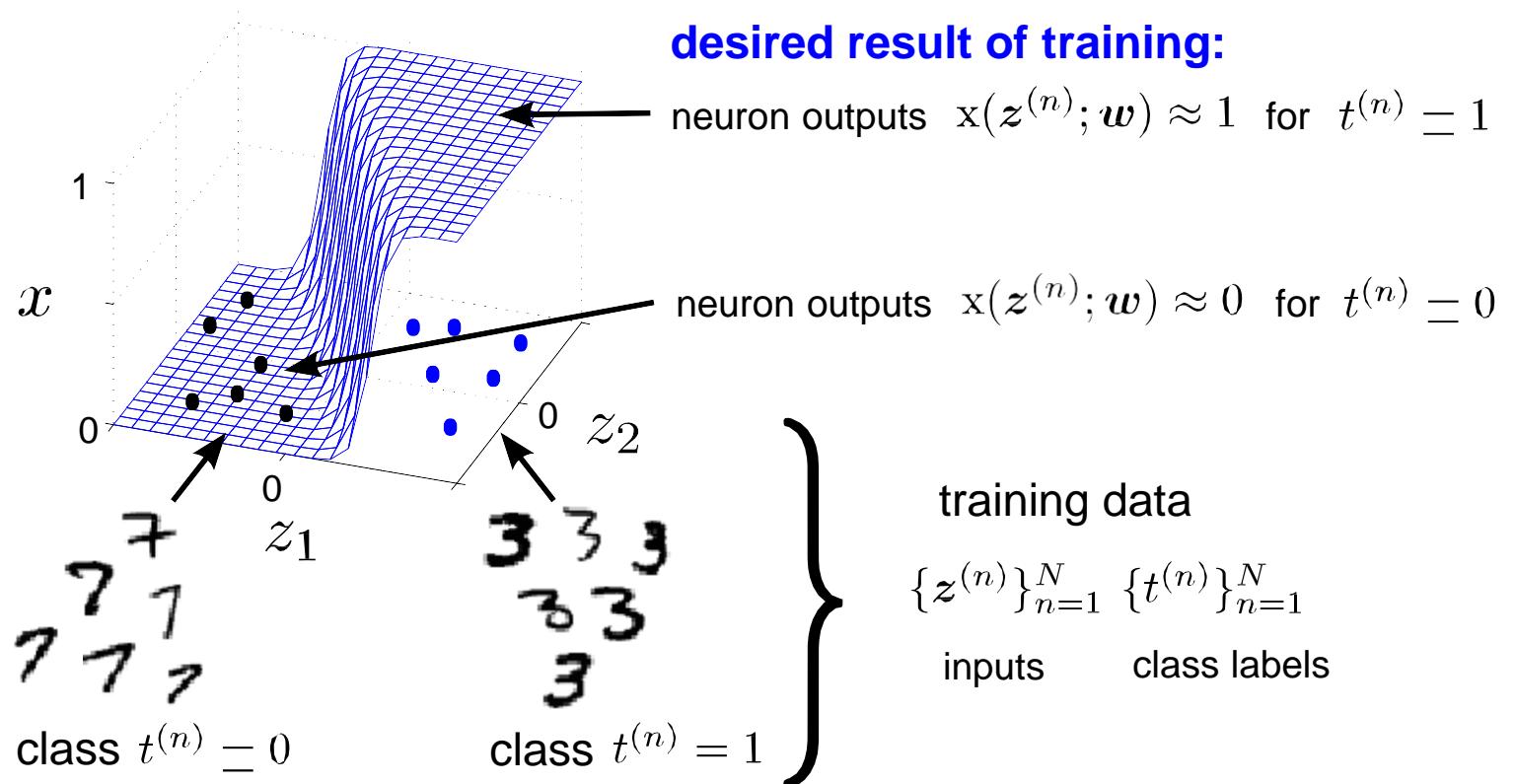
# Weight Space of a Single Neuron



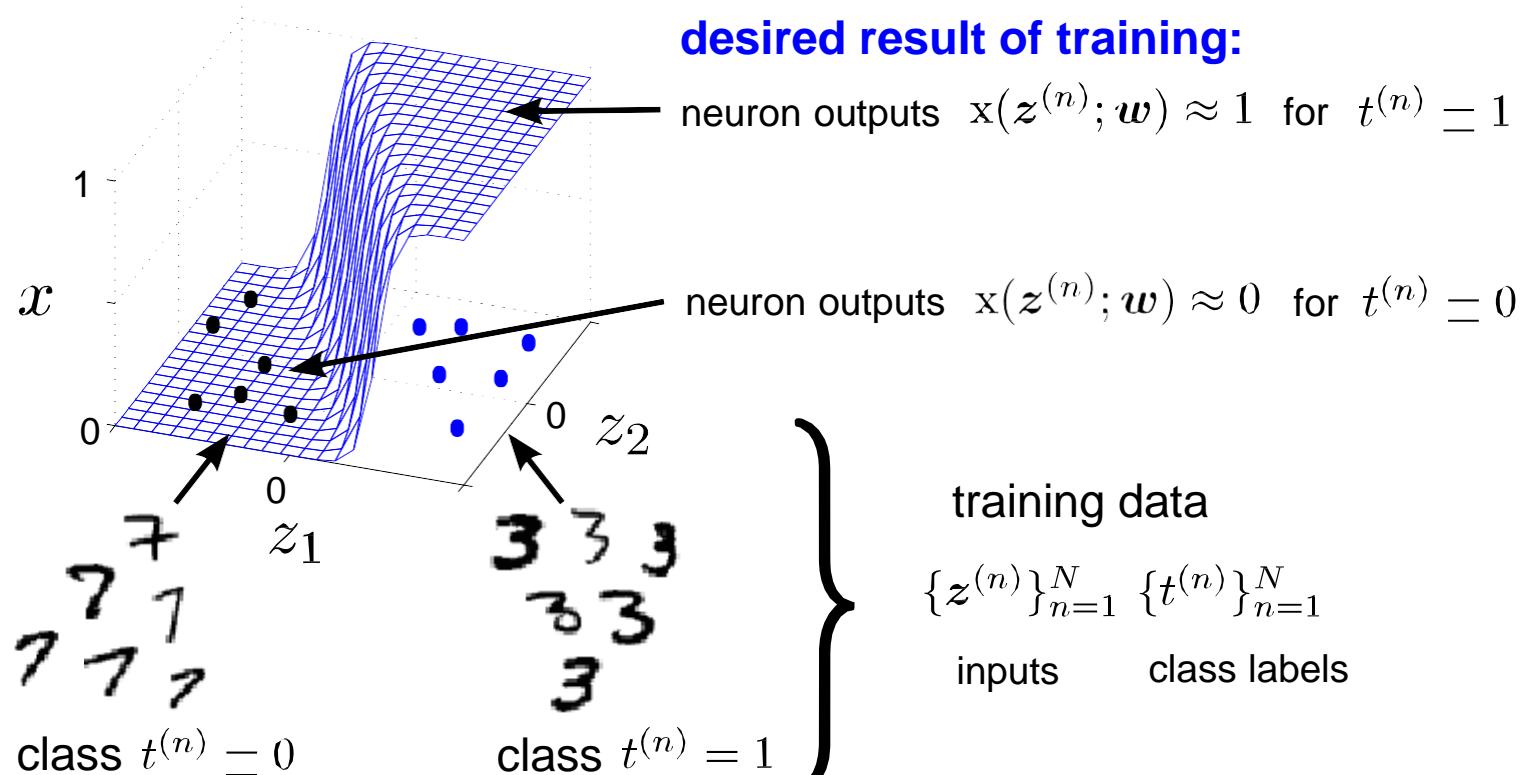
# Training a Single Neuron



# Training a Single Neuron



# Training a Single Neuron

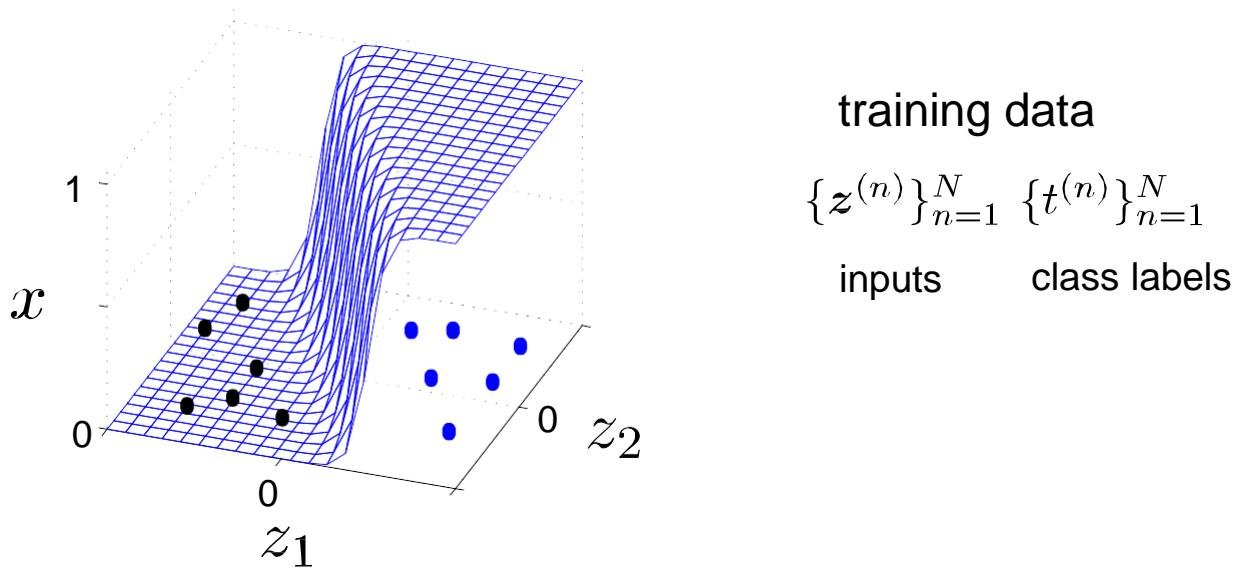


## objective function:

$$G(\mathbf{w}) = - \sum_n [t^{(n)} \log x(z^{(n)}; \mathbf{w}) + (1 - t^{(n)}) \log (1 - x(z^{(n)}; \mathbf{w}))] \geq 0$$

surprise  $- \log p(\text{outcome})$  when observing  $t^{(n)}$  } encourages neuron output  
 relative entropy between  $x(z^{(n)}; \mathbf{w})$  and  $t^{(n)}$  } to match training data

# Training a Single Neuron



**objective function:**

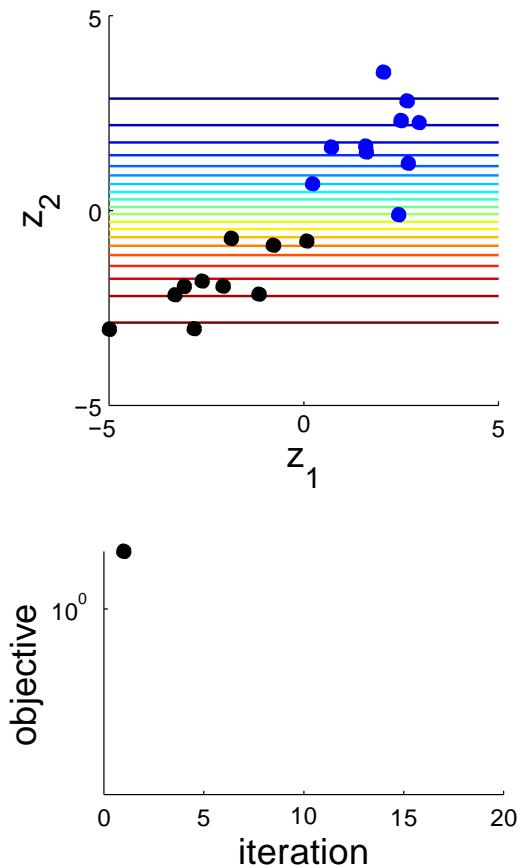
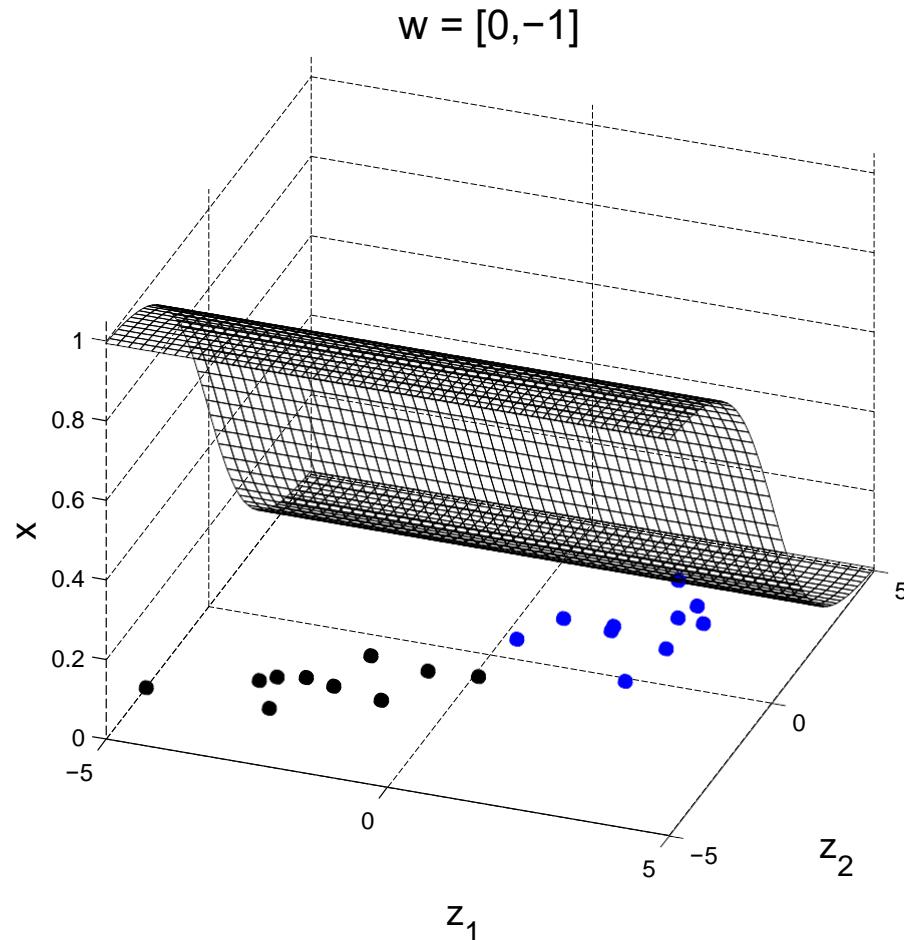
$$G(\mathbf{w}) = - \sum_n [t^{(n)} \log x(\mathbf{z}^{(n)}; \mathbf{w}) + (1 - t^{(n)}) \log (1 - x(\mathbf{z}^{(n)}; \mathbf{w}))] \geq 0$$

$\mathbf{w}^* = \arg \min_{\mathbf{w}} G(\mathbf{w})$  choose the weights that minimise the network's surprise about the training data

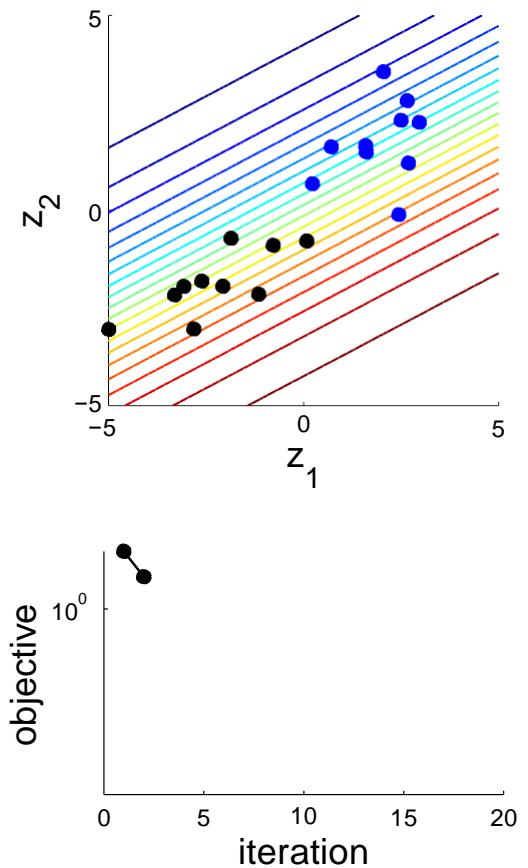
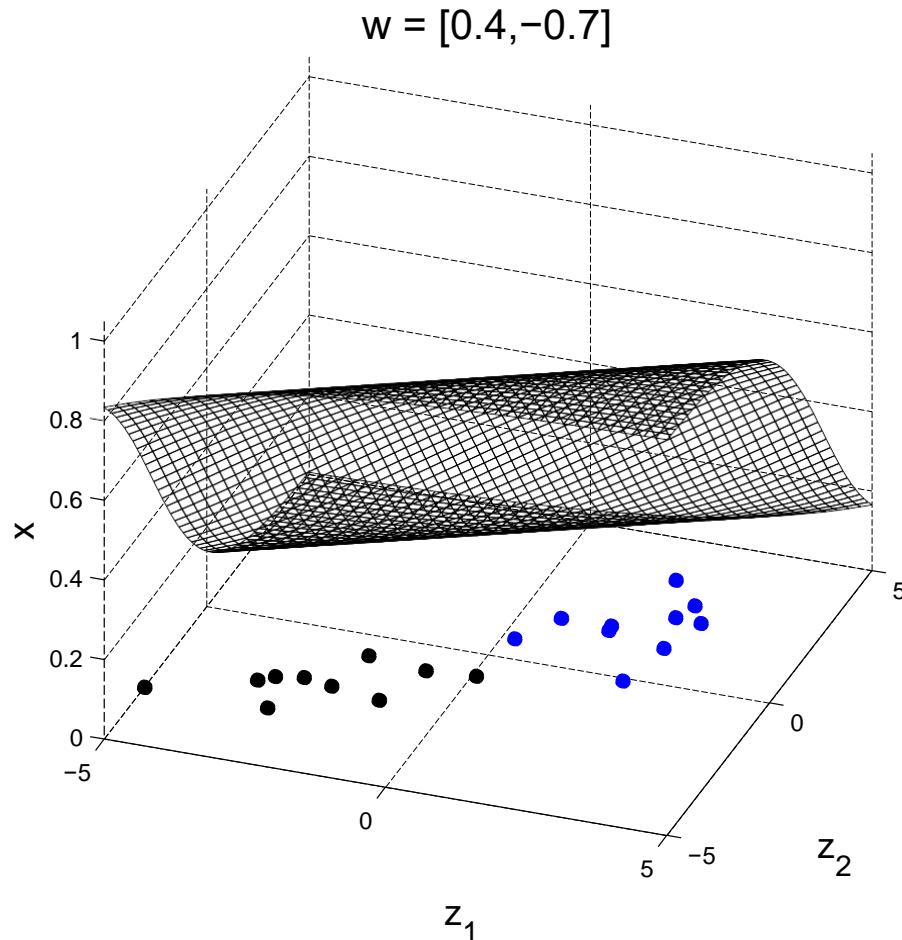
$$\frac{d}{d\mathbf{w}} G(\mathbf{w}) = \sum_n \frac{dG(\mathbf{w})}{dx^{(n)}} \frac{dx^{(n)}}{d\mathbf{w}} = - \sum_n (t^{(n)} - x^{(n)}) \mathbf{z}^{(n)} = \text{prediction error} \times \text{feature}$$

$\mathbf{w} \leftarrow \mathbf{w} - \eta \frac{d}{d\mathbf{w}} G(\mathbf{w})$  iteratively step down the objective (gradient points up hill)

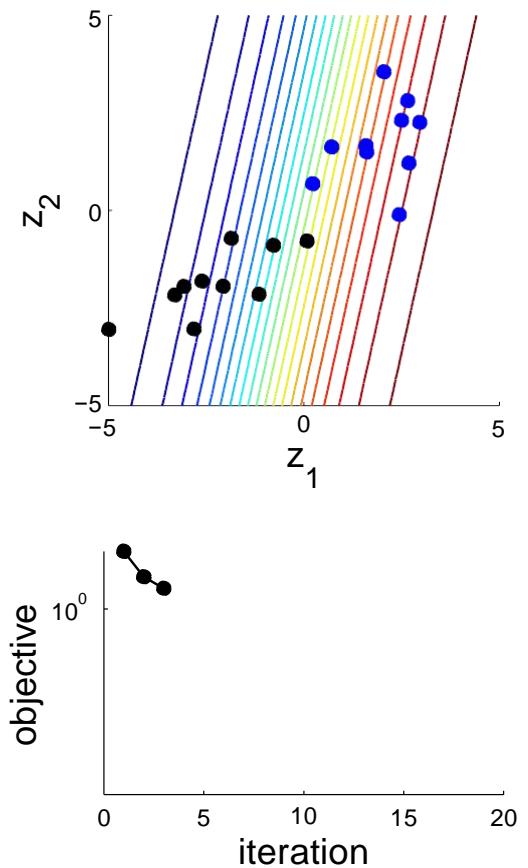
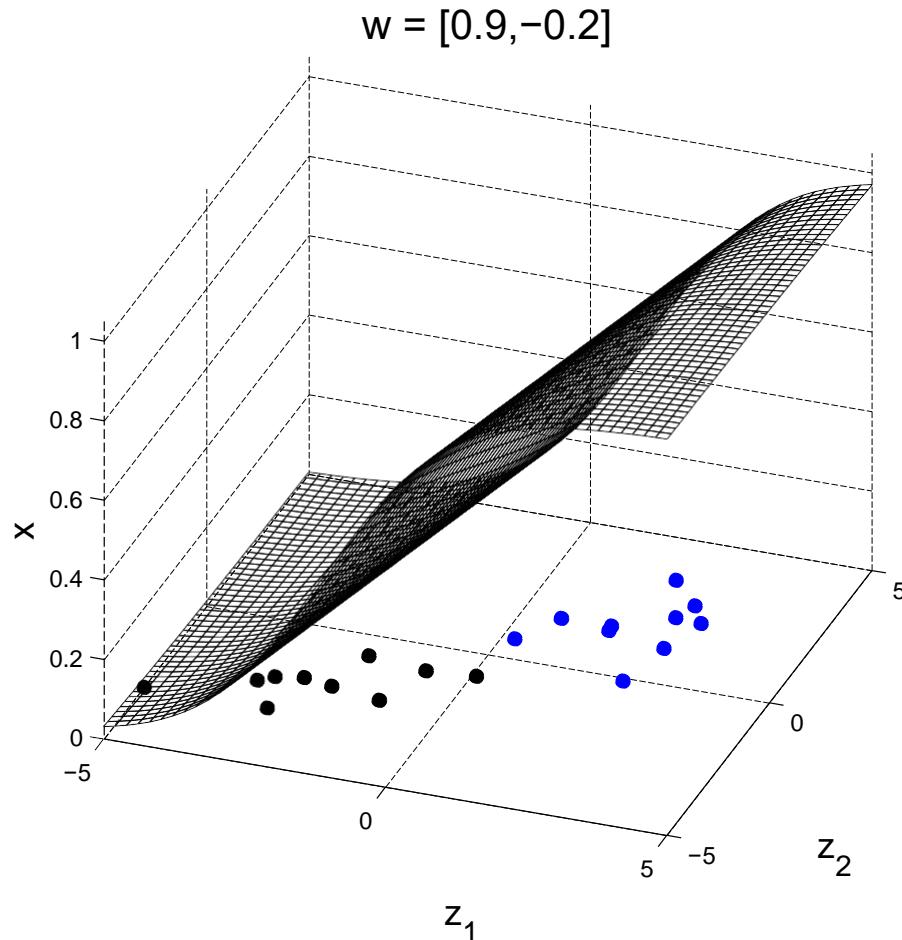
# Training a Single Neuron



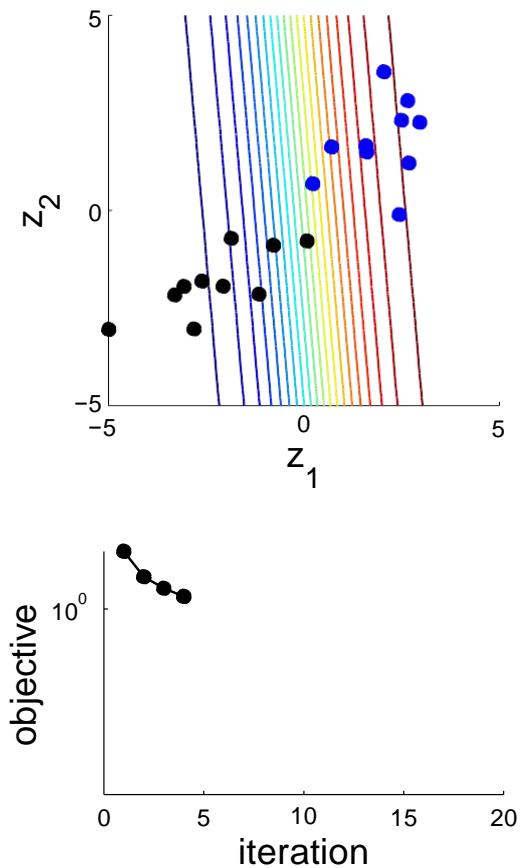
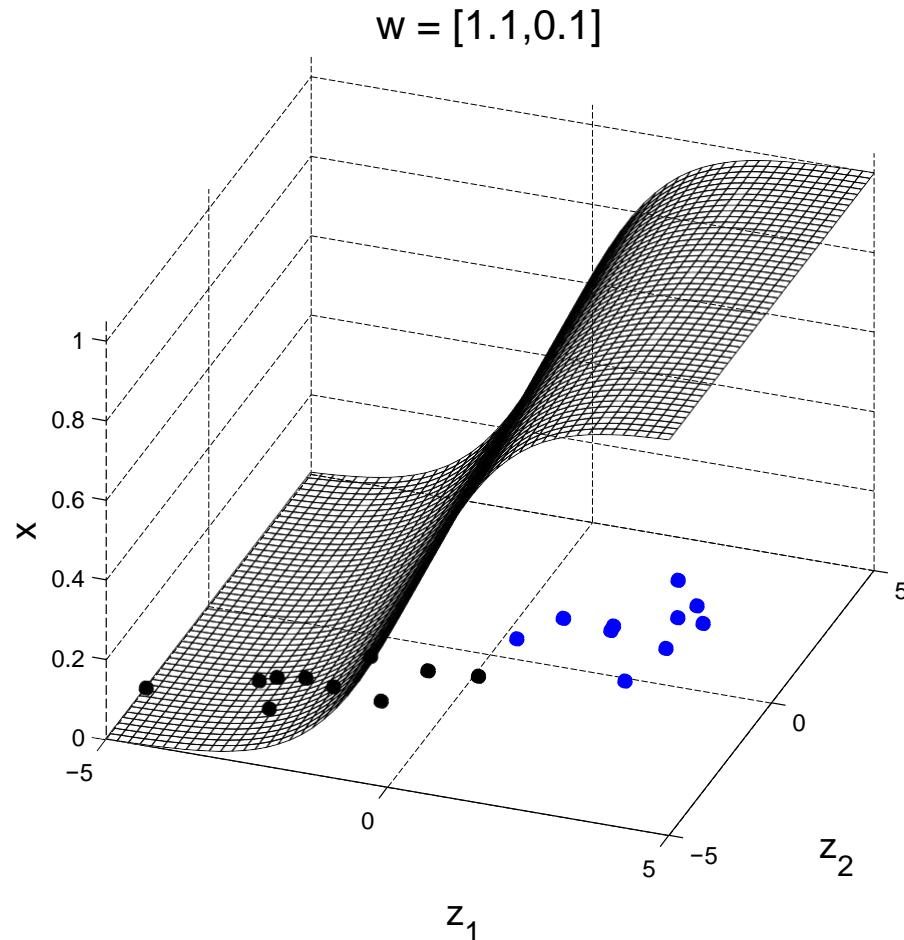
# Training a Single Neuron



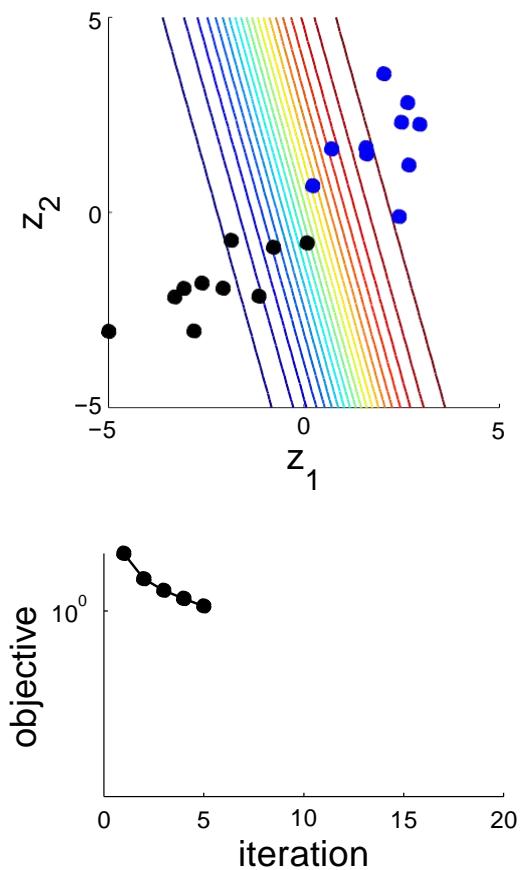
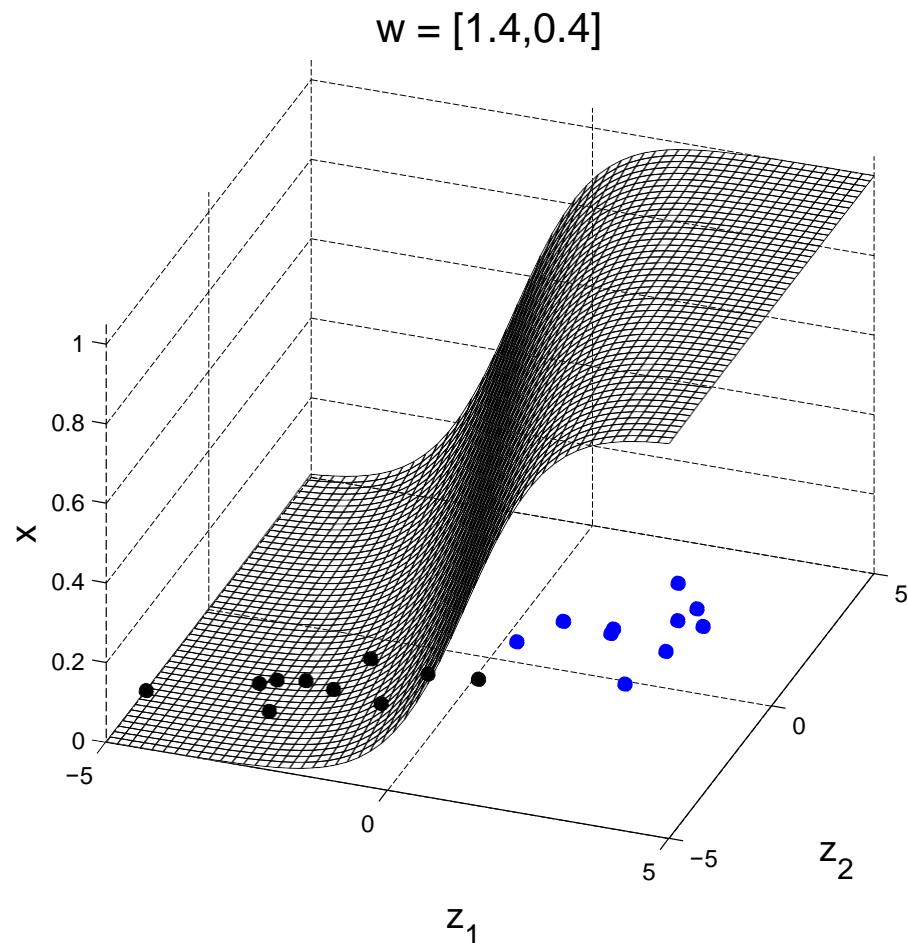
# Training a Single Neuron



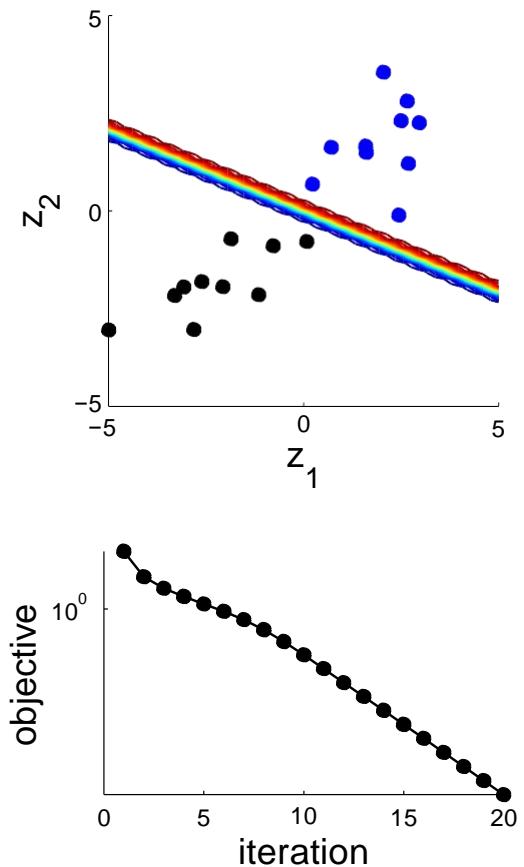
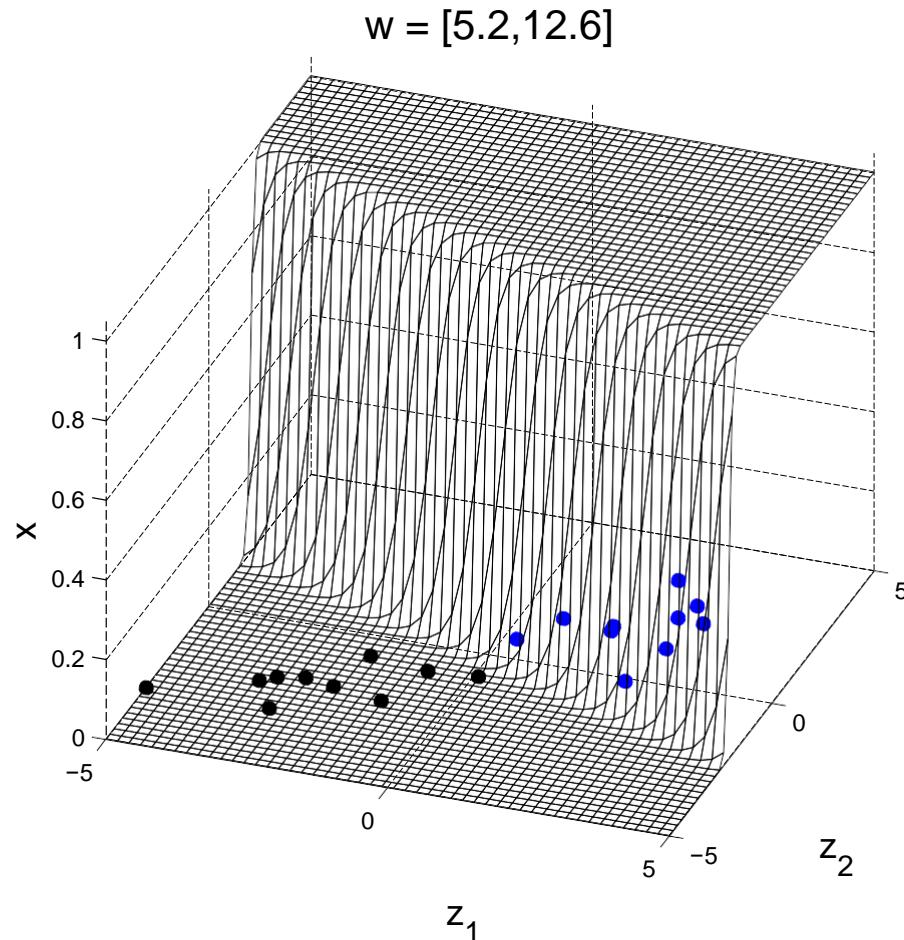
# Training a Single Neuron



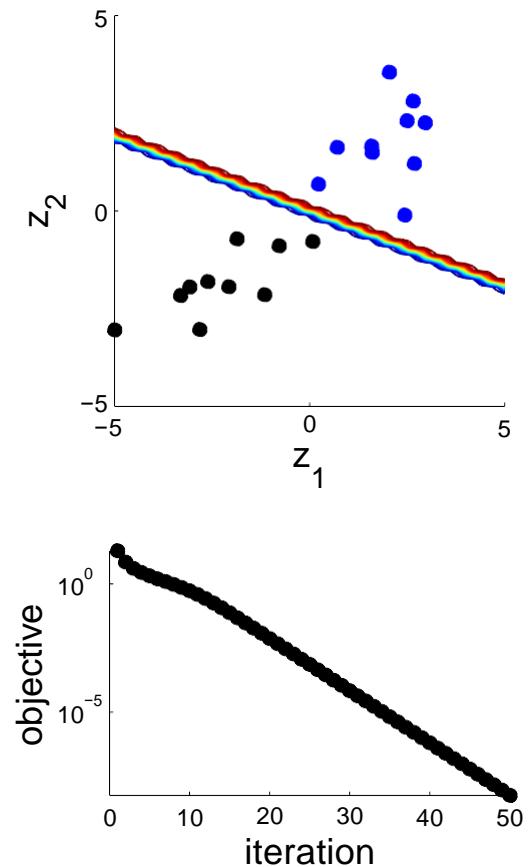
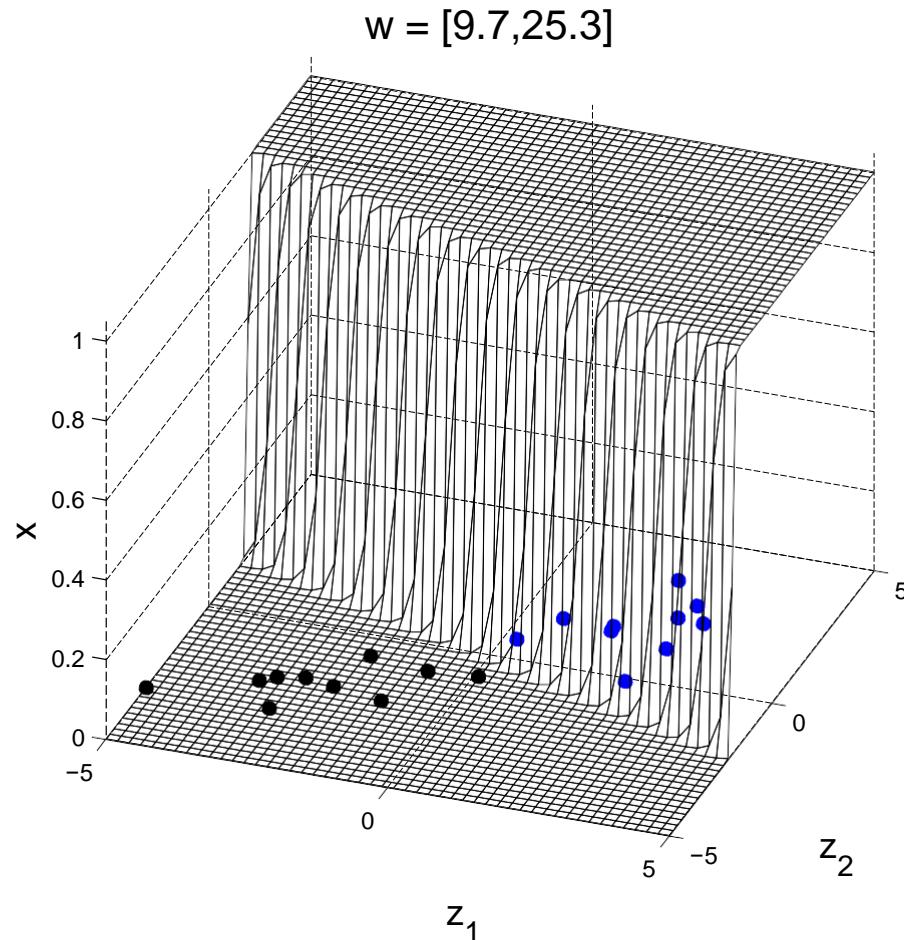
# Training a Single Neuron



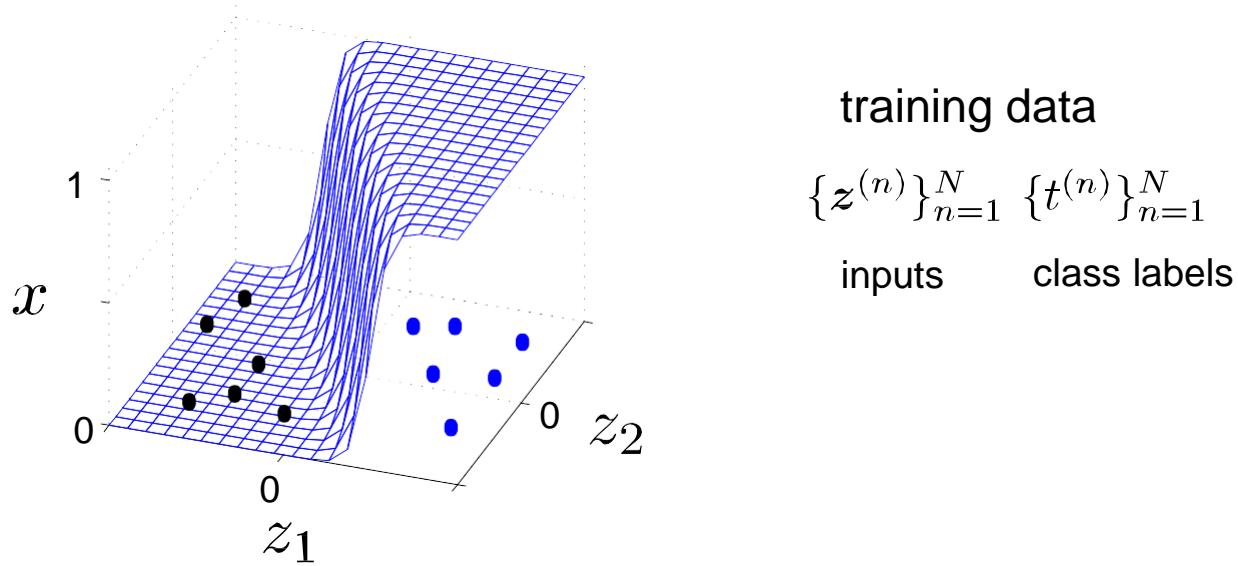
# Training a Single Neuron



# Training a Single Neuron



# Overfitting and Weight Decay



**objective function:**

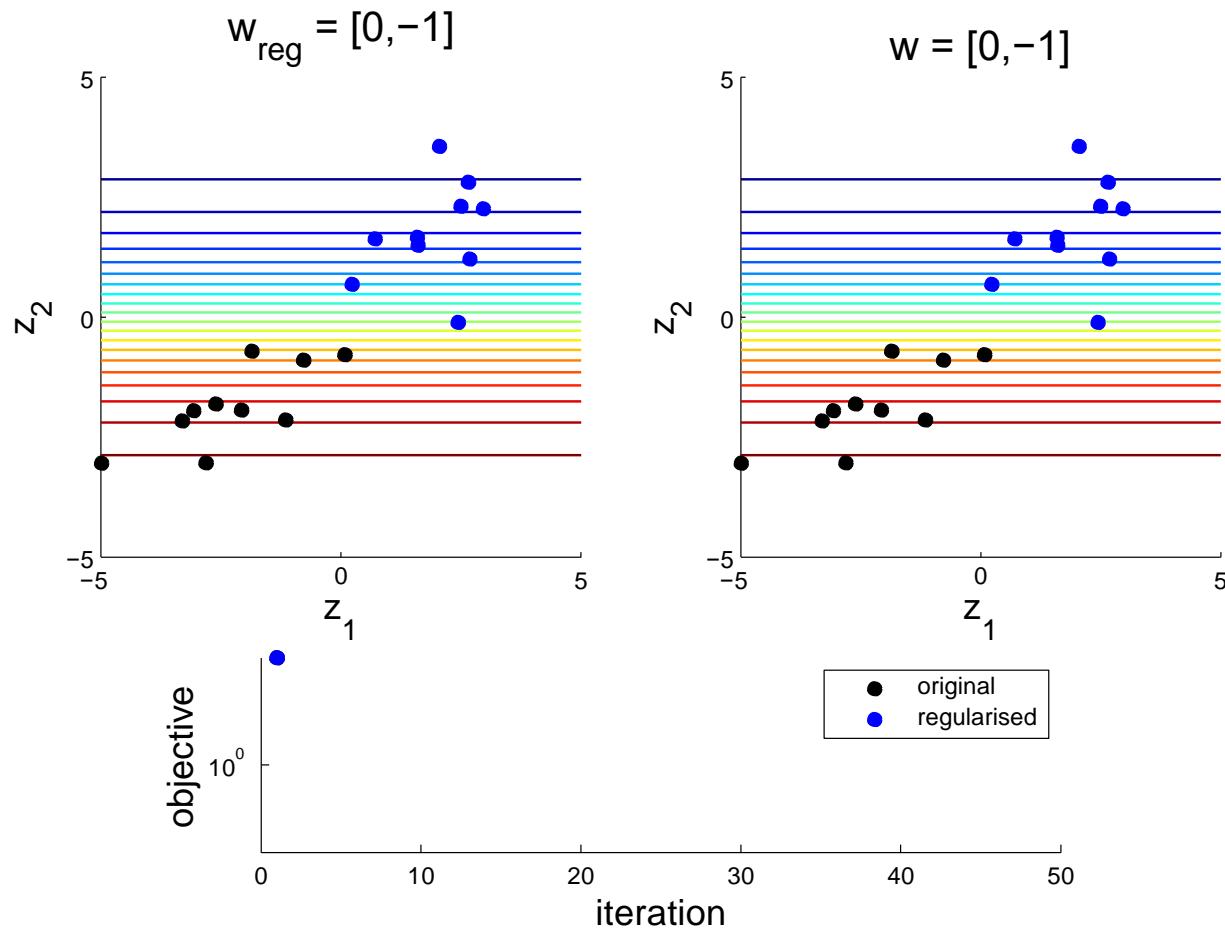
$$G(\mathbf{w}) = - \sum_n [t^{(n)} \log x(\mathbf{z}^{(n)}; \mathbf{w}) + (1 - t^{(n)}) \log (1 - x(\mathbf{z}^{(n)}; \mathbf{w}))]$$

$$E(\mathbf{w}) = \frac{1}{2} \sum_i w_i^2 \quad \text{regulariser discourages the network using extreme weights}$$

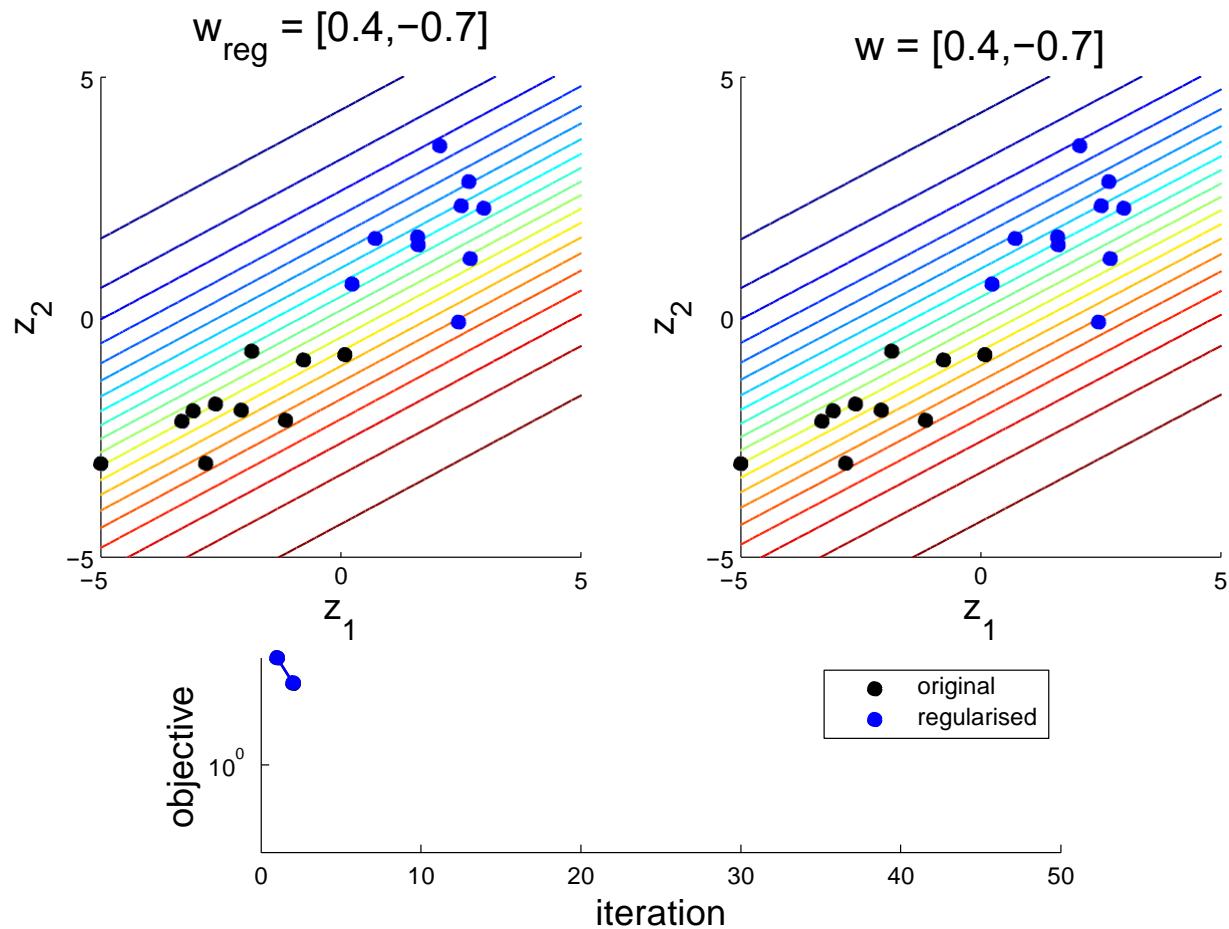
$$\mathbf{w}^* = \arg \min_{\mathbf{w}} M(\mathbf{w}) = \arg \min_{\mathbf{w}} [G(\mathbf{w}) + \alpha E(\mathbf{w})]$$

$$\frac{d}{d\mathbf{w}} M(\mathbf{w}) = - \sum_n (t^{(n)} - x^{(n)}) \mathbf{z}^{(n)} + \alpha \mathbf{w} \quad \text{weight decay - shrinks weights towards zero}$$

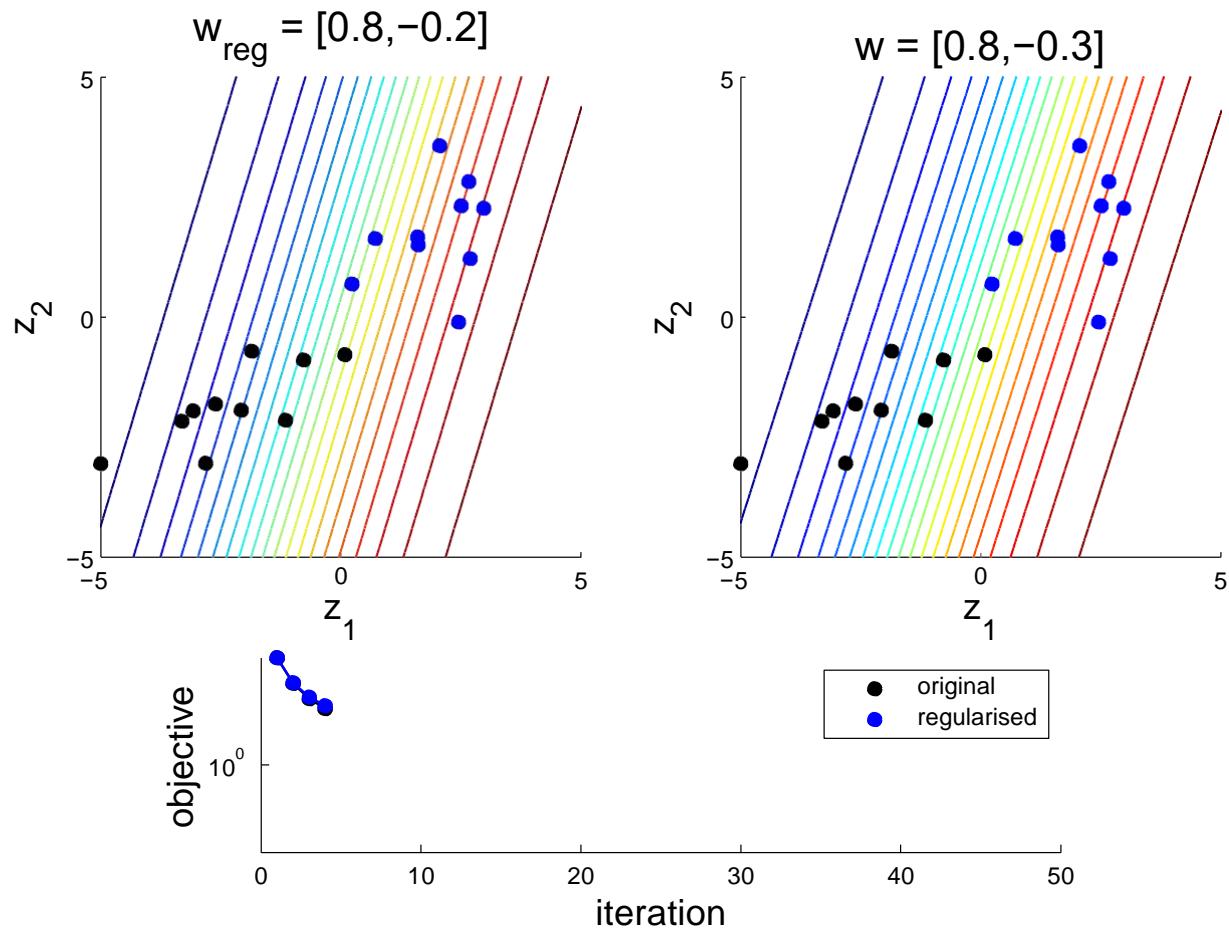
# Training a Single Neuron (cont'd)



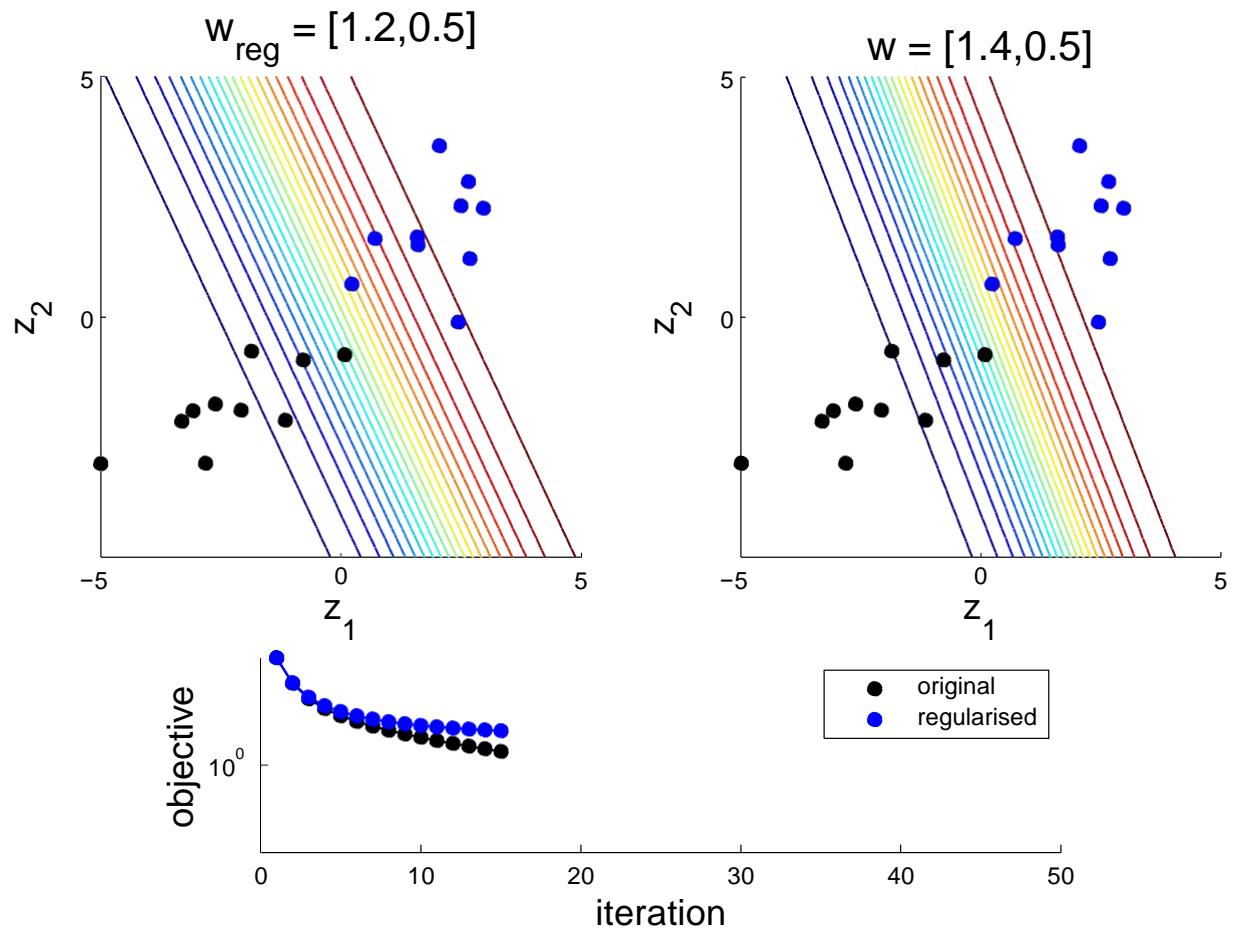
# Training a Single Neuron (cont'd)



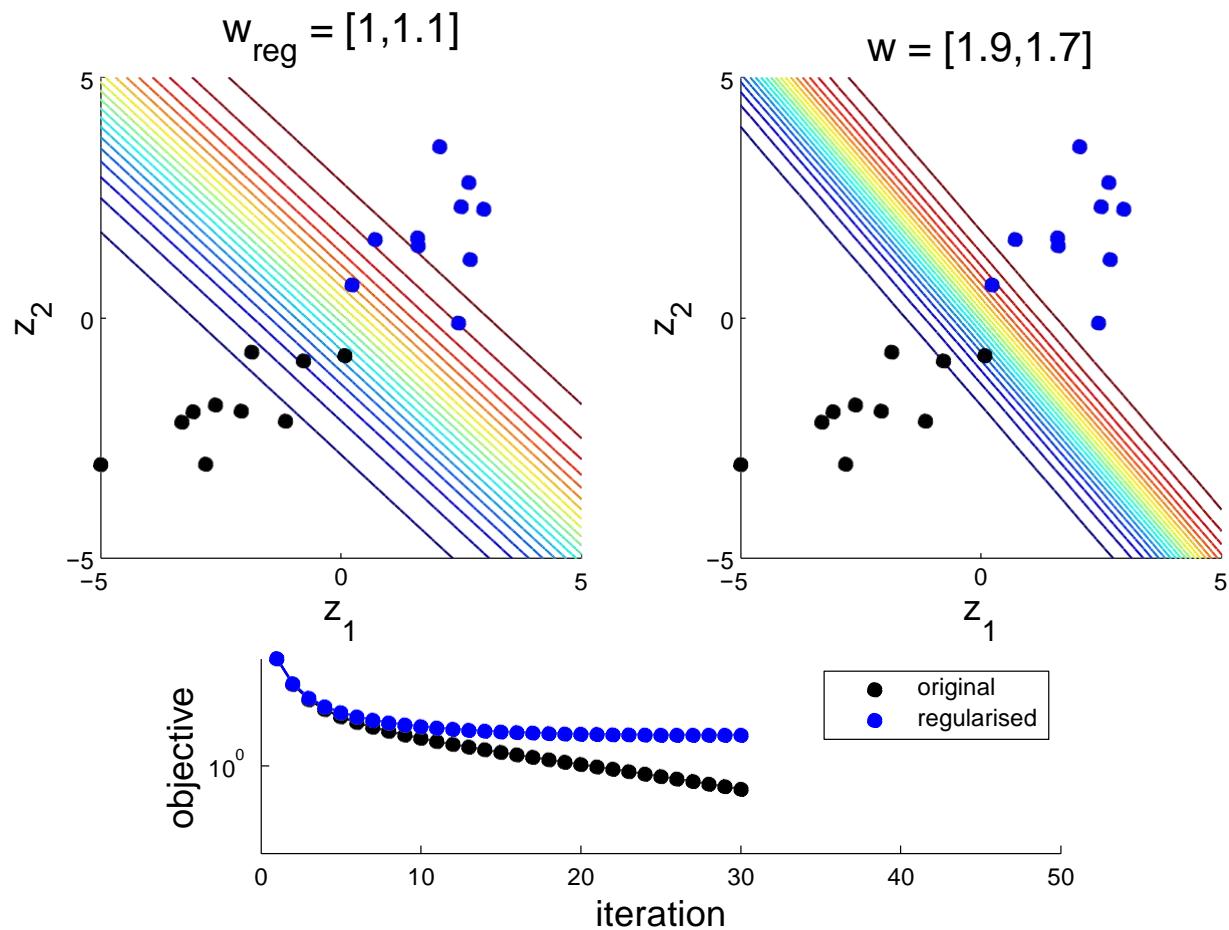
# Training a Single Neuron (cont'd)



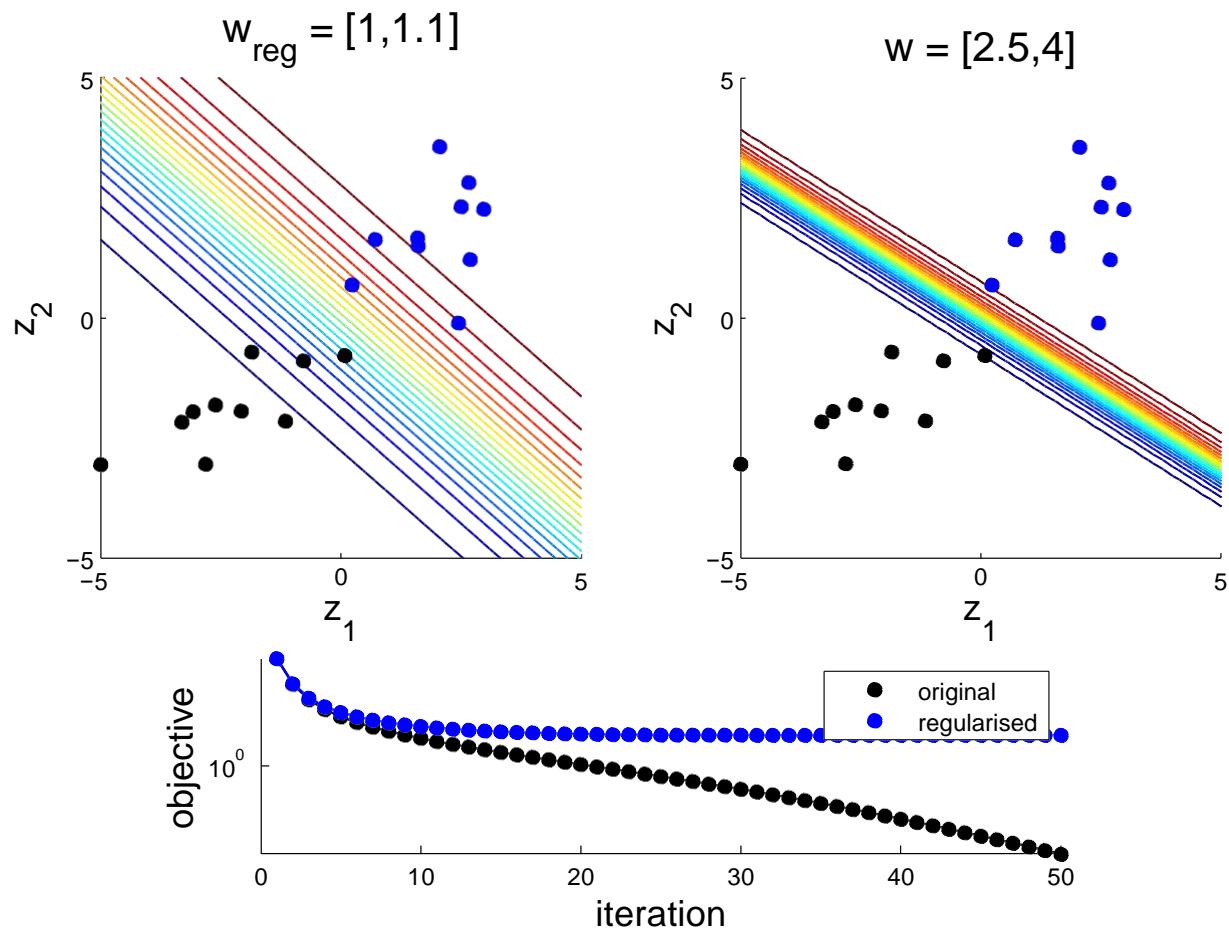
# Training a Single Neuron (cont'd)



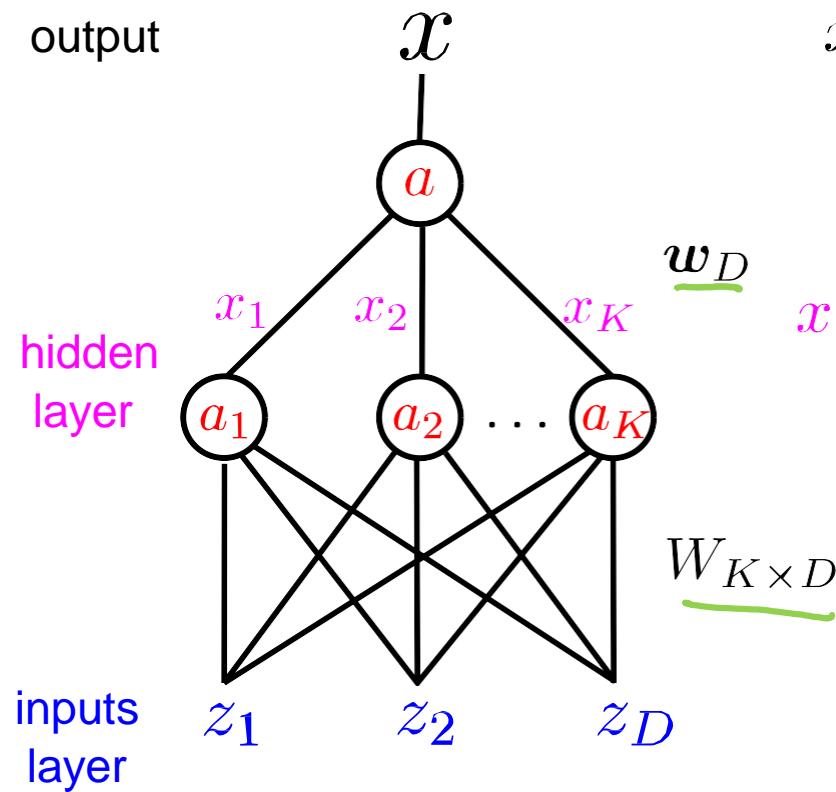
# Training a Single Neuron (cont'd)



# Training a Single Neuron (cont'd)



# Single Hidden Layer Neural Networks

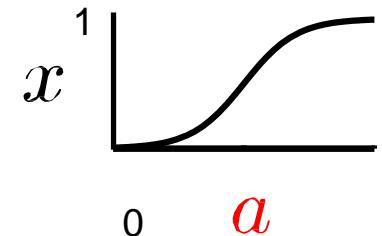


$$x(\textcolor{red}{a}) = \frac{1}{1 + \exp(-\textcolor{red}{a})}$$

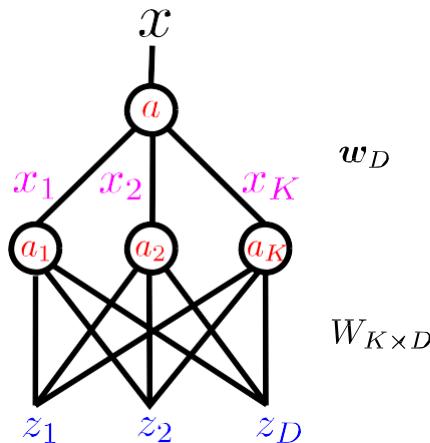
$$\textcolor{red}{a} = \sum_{k=1}^K w_k \textcolor{magenta}{x}_k$$

$$x(\textcolor{red}{a}_k) = \frac{1}{1 + \exp(-\textcolor{red}{a}_k)}$$

$$a_k = \sum_{d=1}^D W_{k,d} \textcolor{blue}{z}_d$$



# Training a Neural Network with a Single Hidden Layer



$$x(\textcolor{red}{a}) = \frac{1}{1 + \exp(-\textcolor{red}{a})}$$

$$\textcolor{red}{a} = \sum_{k=1}^K w_k \textcolor{violet}{x}_k$$

$$\textcolor{violet}{x}(\textcolor{red}{a}_k) = \frac{1}{1 + \exp(-\textcolor{red}{a}_k)}$$

$$a_k = \sum_{d=1}^D W_{k,d} \textcolor{blue}{z}_d$$

**objective function:**

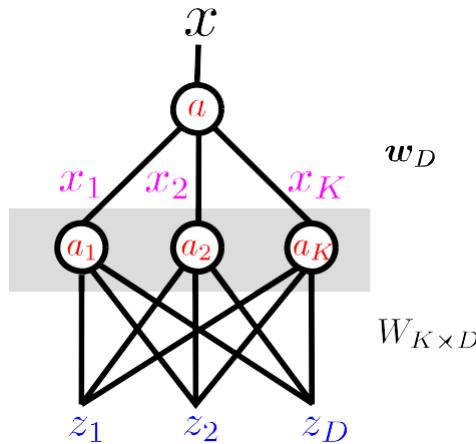
$$G(W, \mathbf{w}) = - \sum_n [t^{(n)} \log x^{(n)} + (1 - t^{(n)}) \log (1 - x^{(n)})] \quad \text{likelihood same as before}$$

$$E(W, \mathbf{w}) = \frac{1}{2} \sum_i w_i^2 + \frac{1}{2} \sum_{ij} W_{ij}^2 \quad \text{regulariser discourages extreme weights}$$

$$\{W, \mathbf{w}^*\} = \arg \min_{W, \mathbf{w}} M(W, \mathbf{w}) = \arg \min_{W, \mathbf{w}} [G(W, \mathbf{w}) + \alpha E(W, \mathbf{w})]$$

# Training a Neural Network with a Single Hidden Layer

Networks with hidden layers can be fit using gradient descent using an algorithm called **back-propagation**.



$$x(\mathbf{a}) = \frac{1}{1 + \exp(-\mathbf{a})}$$

$$\mathbf{a} = \sum_{k=1}^K w_k \mathbf{x}_k$$

$$\mathbf{x}(\mathbf{a}_k) = \frac{1}{1 + \exp(-\mathbf{a}_k)}$$

$$a_k = \sum_{d=1}^D W_{k,d} z_d$$

## objective function:

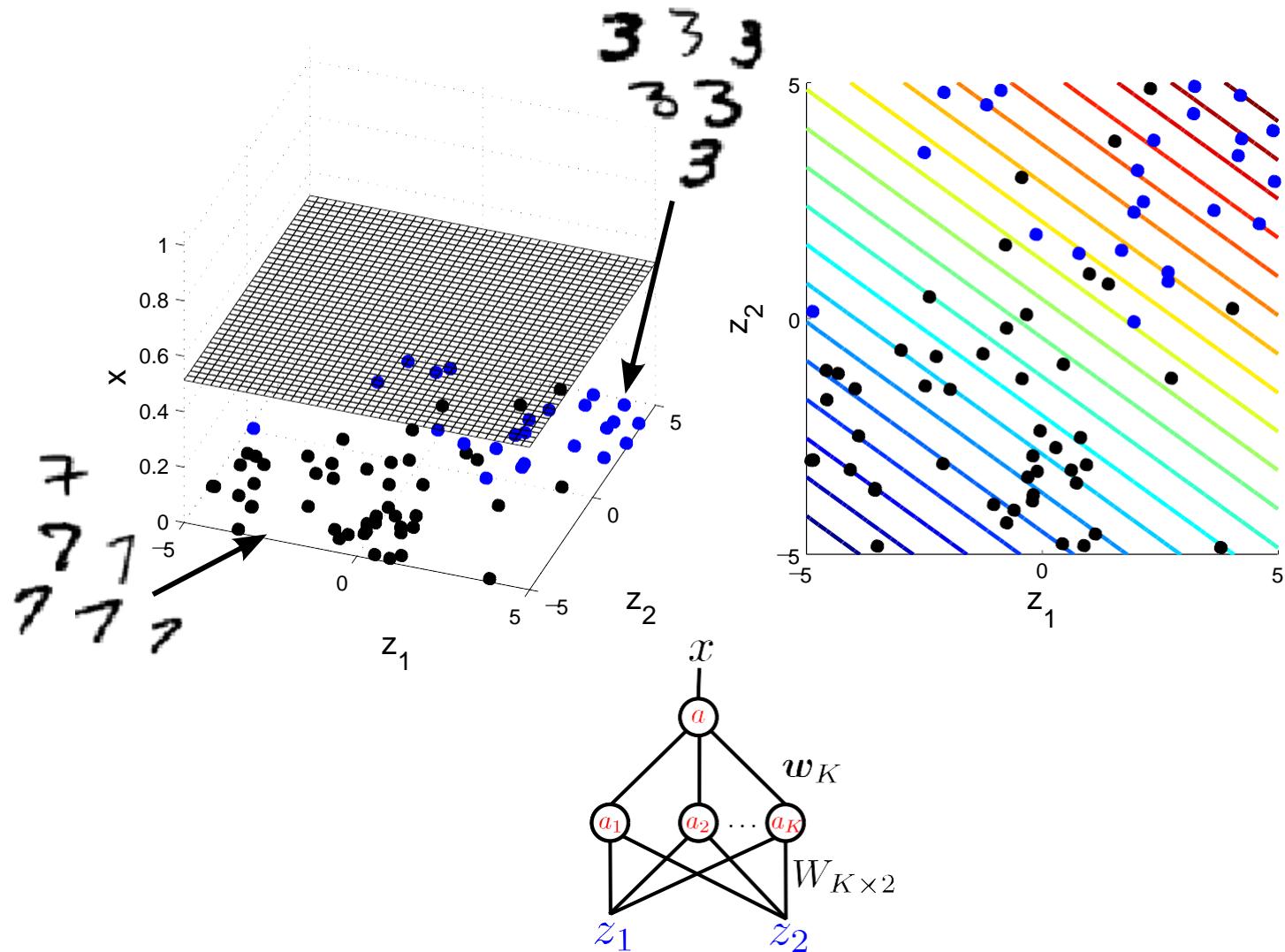
$$G(W, \mathbf{w}) = - \sum_n [t^{(n)} \log \mathbf{x}^{(n)} + (1 - t^{(n)}) \log (1 - \mathbf{x}^{(n)})] \quad \text{likelihood same as before}$$

$$E(W, \mathbf{w}) = \frac{1}{2} \sum_i w_i^2 + \frac{1}{2} \sum_{ij} W_{ij}^2 \quad \text{regulariser discourages extreme weights}$$

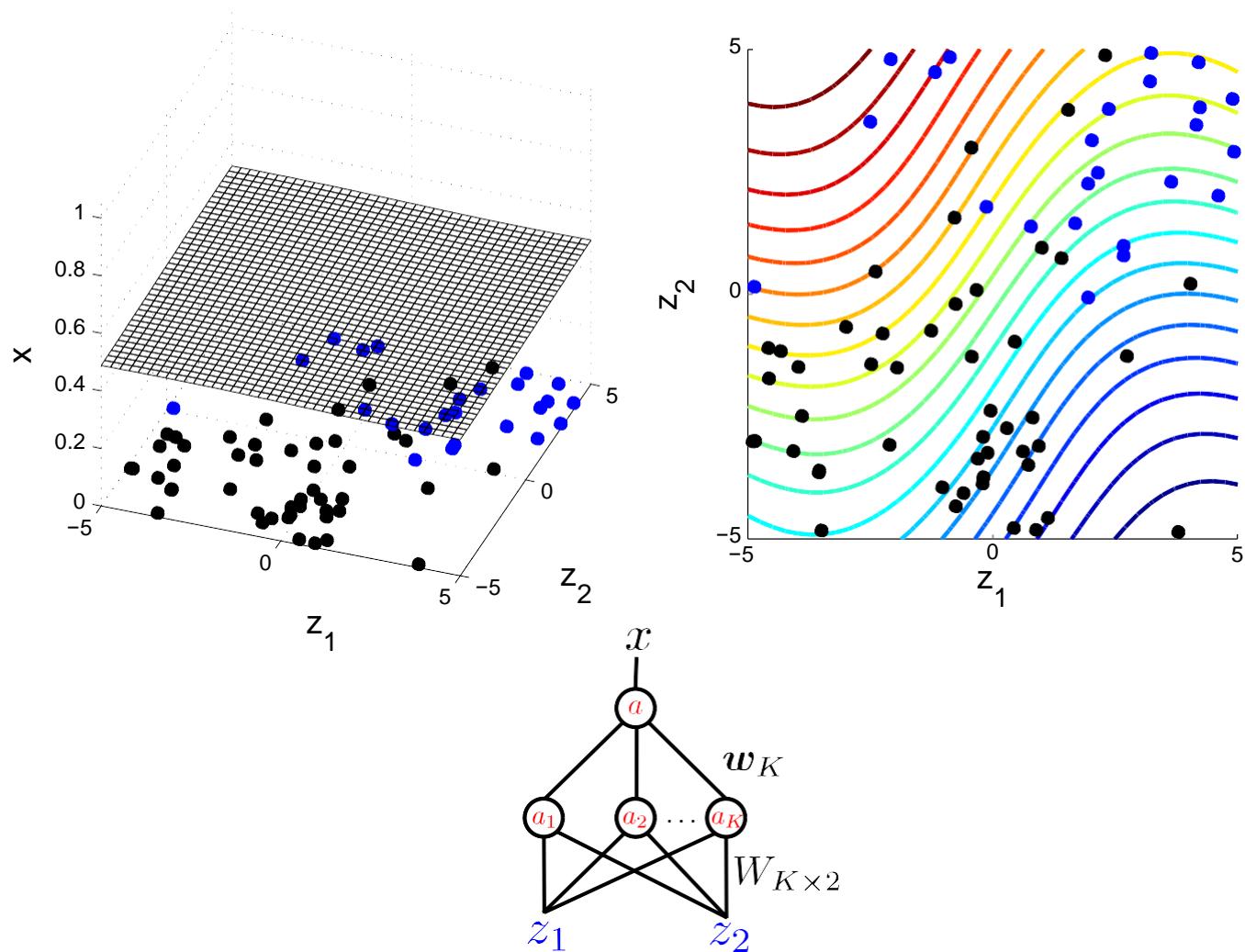
$$\{W, \mathbf{w}^*\} = \arg \min_{W, \mathbf{w}} M(W, \mathbf{w}) = \arg \min_{W, \mathbf{w}} [G(W, \mathbf{w}) + \alpha E(W, \mathbf{w})]$$

$$\begin{aligned} \frac{dG(W, \mathbf{w})}{dW_{ij}} &= \sum_n \frac{dG(W, \mathbf{w})}{dx^{(n)}} \frac{dx^{(n)}}{dW_{ij}} = \sum_n \frac{dG(W, \mathbf{w})}{dx^{(n)}} \frac{dx^{(n)}}{da^{(n)}} \frac{da^{(n)}}{dW_{ij}} \\ &= \sum_n \frac{dG(W, \mathbf{w})}{dx^{(n)}} \frac{dx^{(n)}}{da^{(n)}} \frac{da^{(n)}}{dx_i^{(n)}} \frac{dx_i^{(n)}}{dW_{ij}} = \sum_n \frac{dG(W, \mathbf{w})}{dx^{(n)}} \frac{dx^{(n)}}{da^{(n)}} \frac{da^{(n)}}{dx_i^{(n)}} \frac{dx_i^{(n)}}{da_i^{(n)}} \frac{da_i^{(n)}}{dW_{ij}} \end{aligned}$$

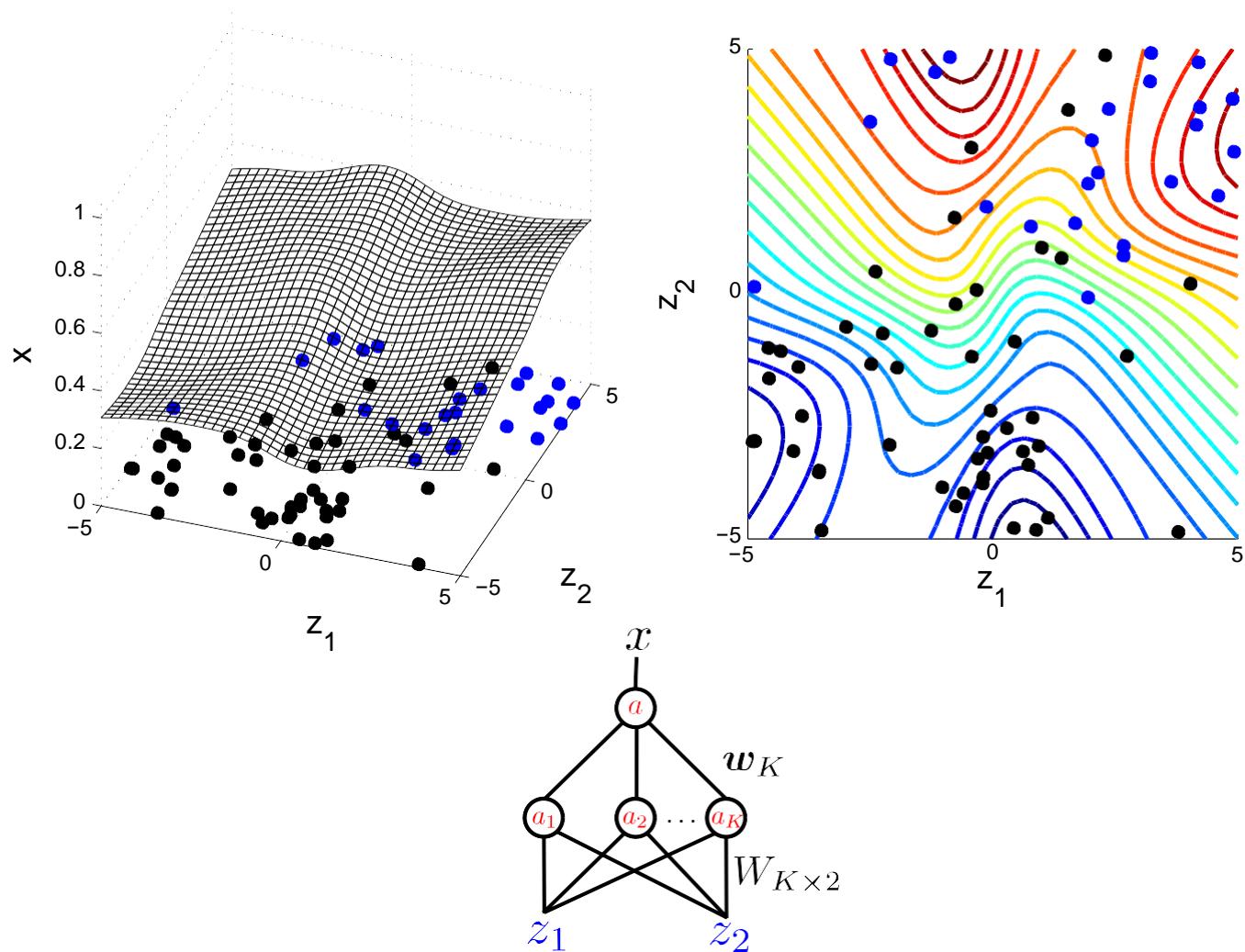
# Training a Neural Network with a Single Hidden Layer



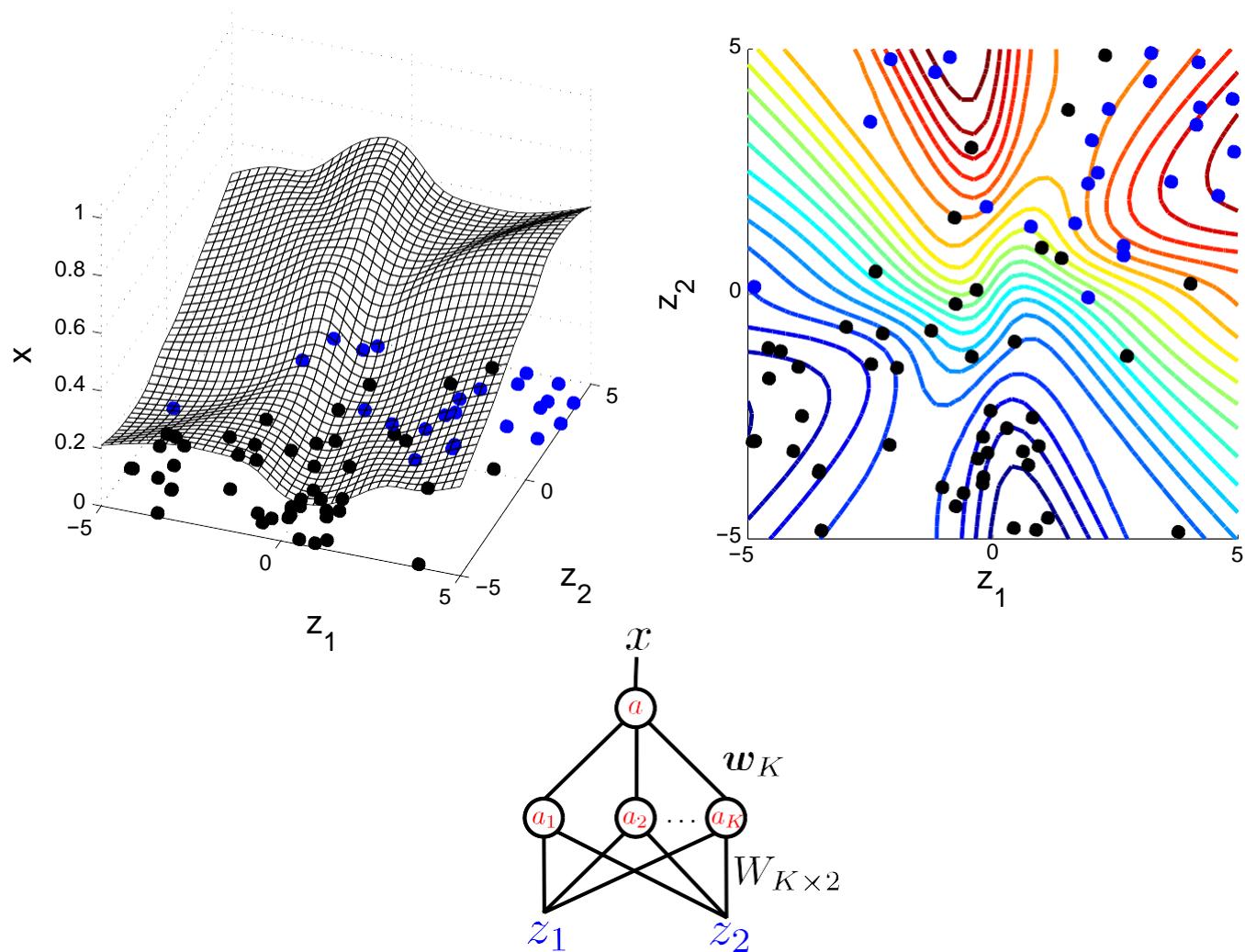
# Training a Neural Network with a Single Hidden Layer



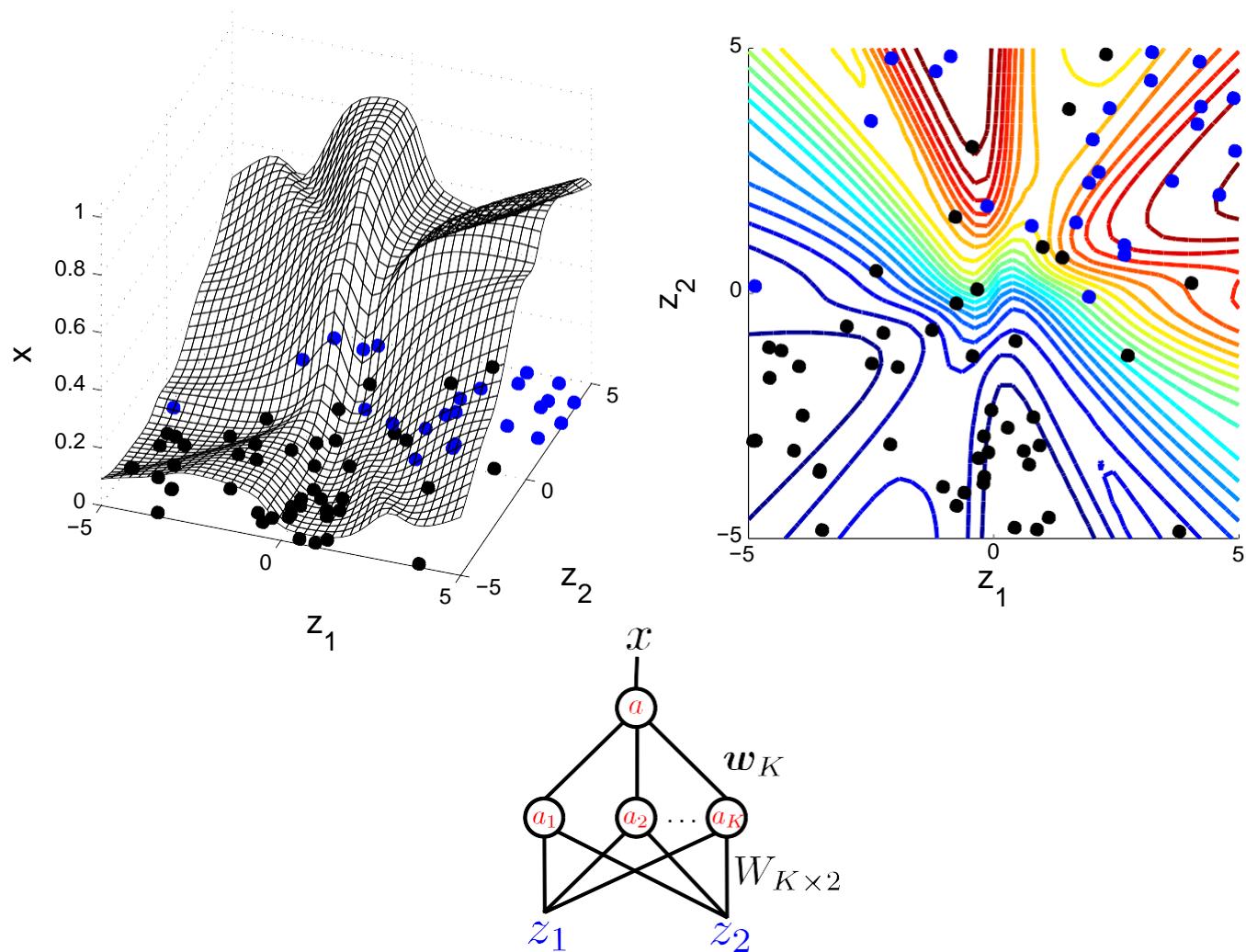
# Training a Neural Network with a Single Hidden Layer



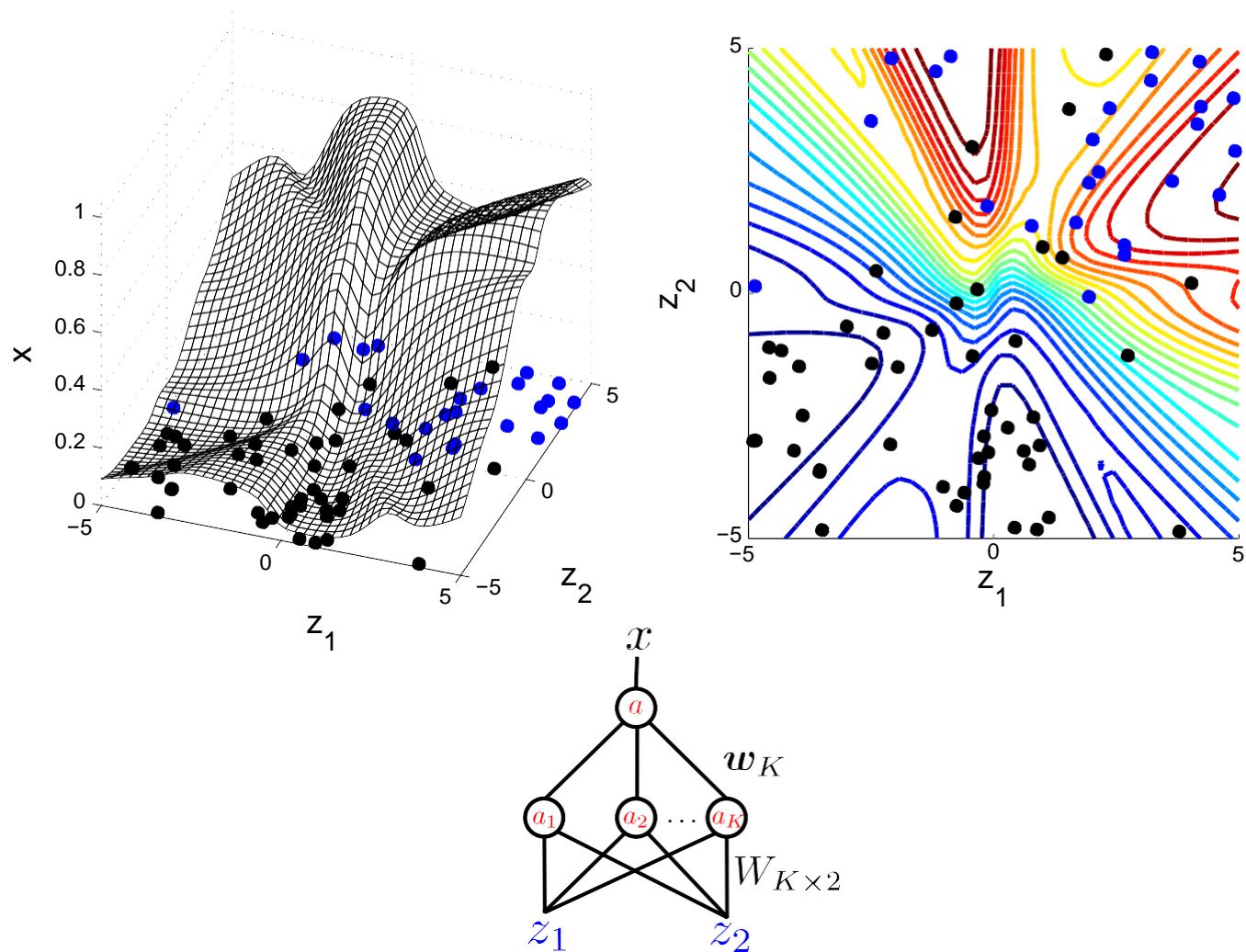
# Training a Neural Network with a Single Hidden Layer



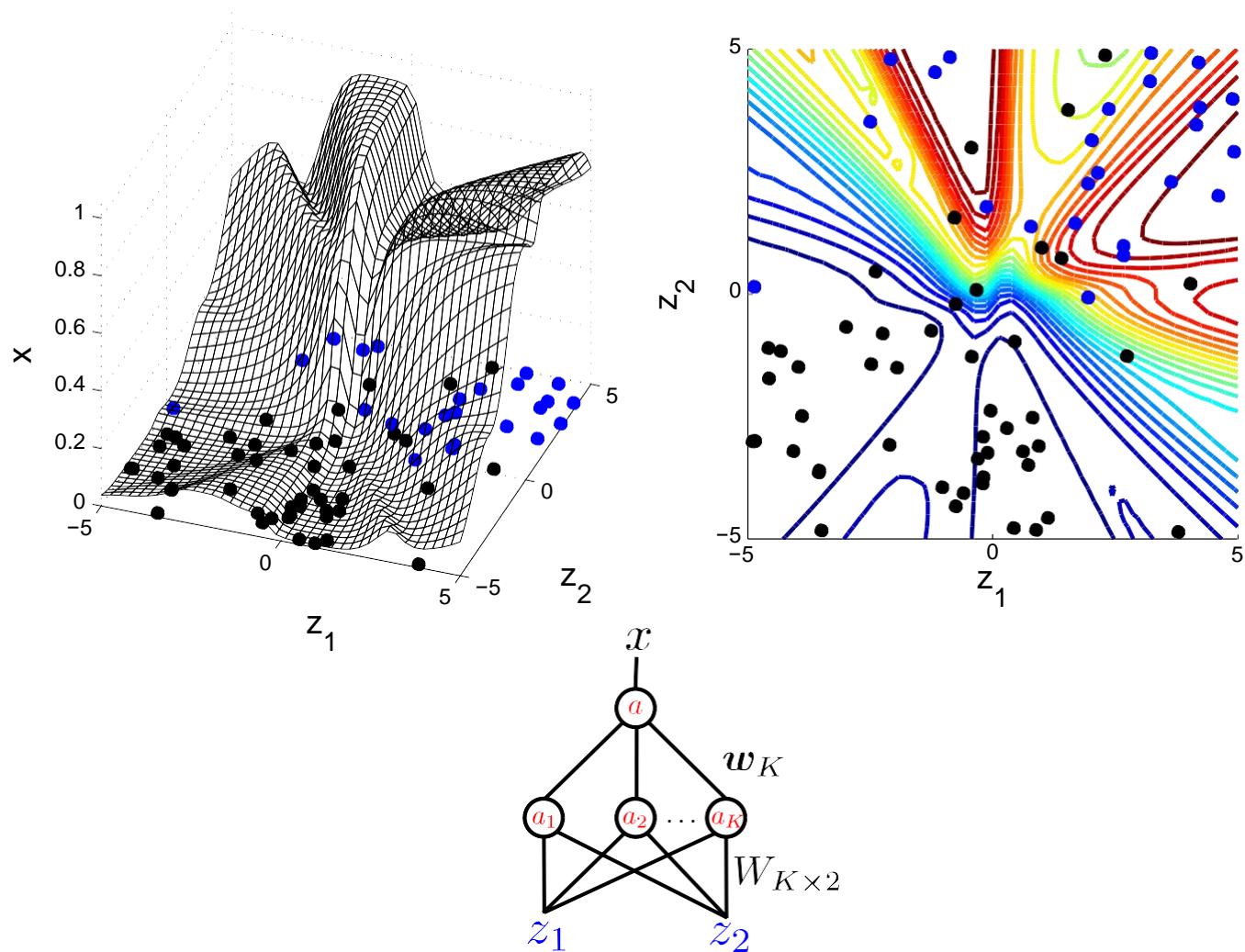
# Training a Neural Network with a Single Hidden Layer



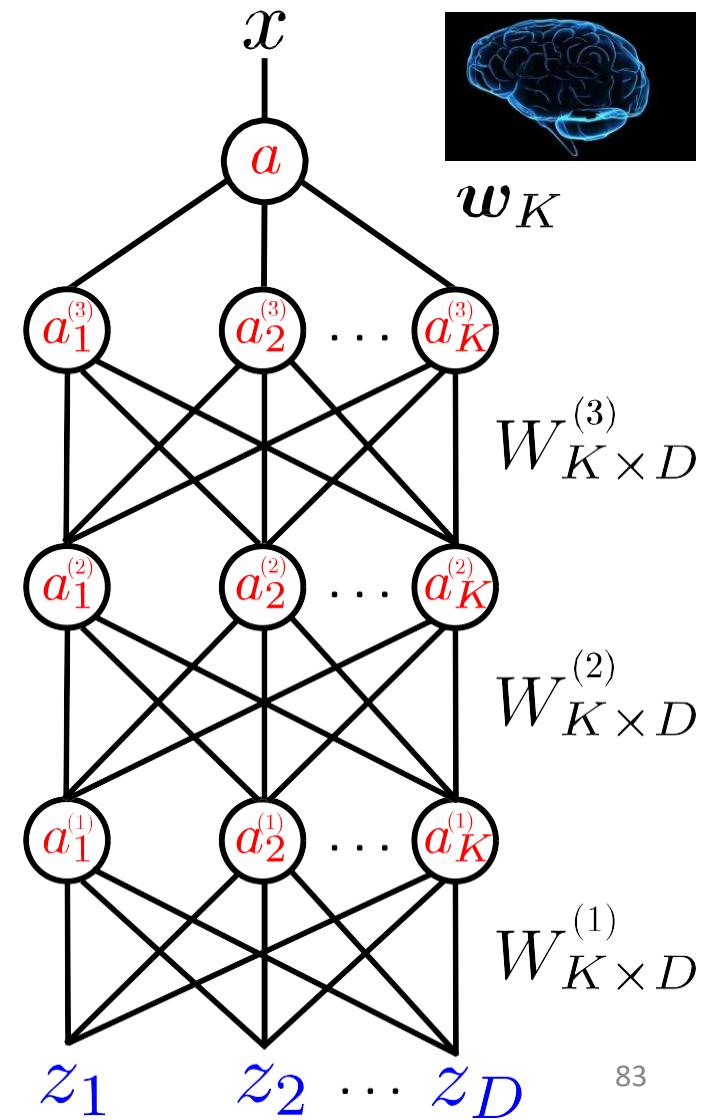
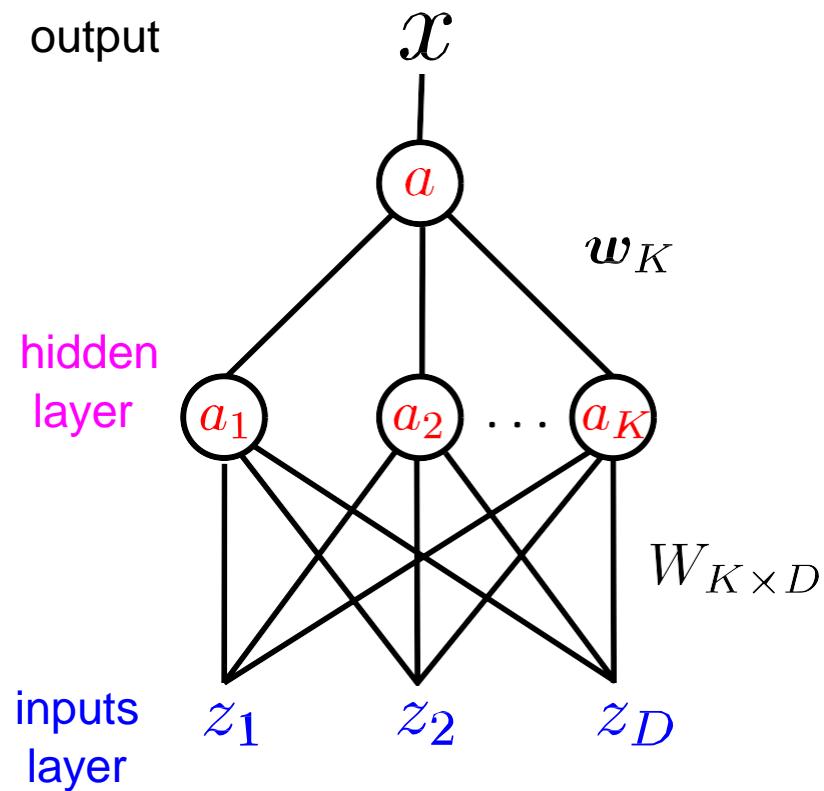
# Training a Neural Network with a Single Hidden Layer



# Training a Neural Network with a Single Hidden Layer



# Hierarchical Models with Many Layers



# What We've Covered Today...

- Unsupervised vs. Supervised Learning
  - Clustering & Dimension Reduction
  - Training, testing, & validation
  - Linear Classification
  - From Linear Classifier to Neural Nets

