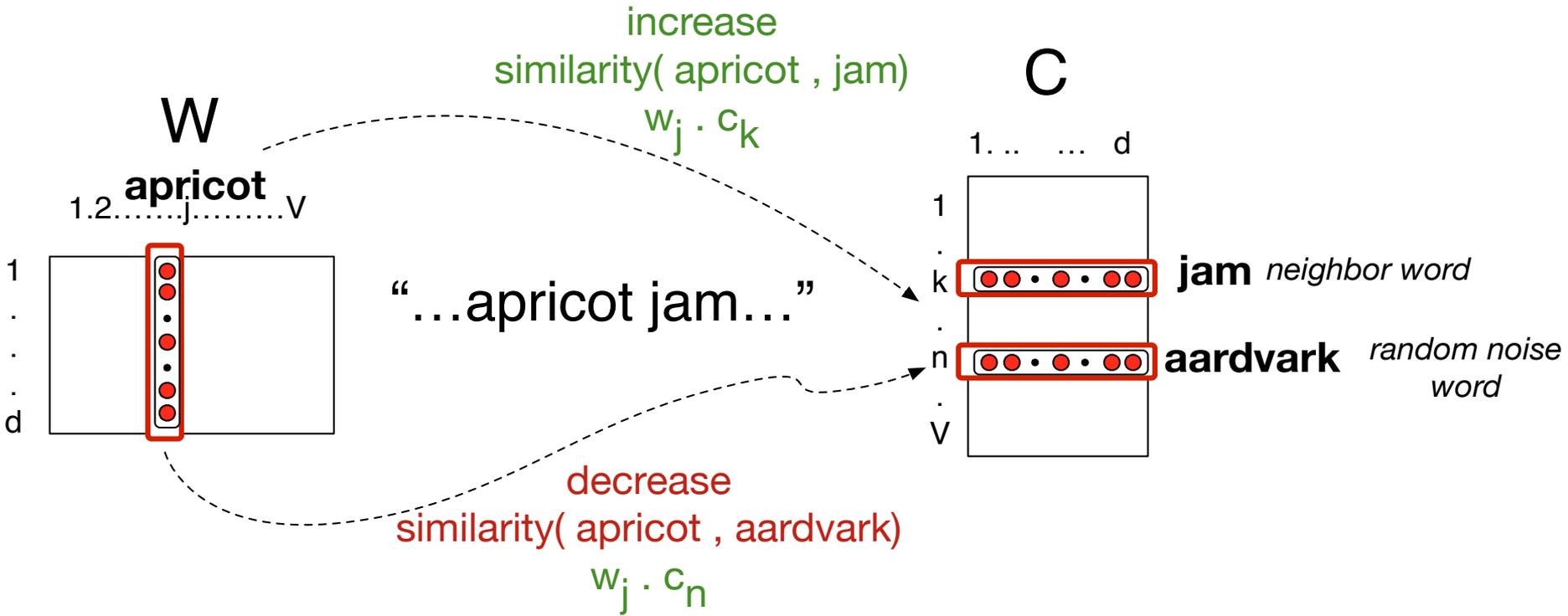


# Logistic Regression

JURAFSKY AND MARTIN CHAPTER 5



# Recap: How to learn word2vec (skip-gram) embeddings

Start with  $V$  random 300-dimensional vectors as initial embeddings

Use **logistic regression** to

- Take a corpus and take pairs of words that co-occur as positive examples
- Take pairs of words that don't co-occur as negative examples
- Train the classifier to distinguish these by slowly adjusting all the embeddings to improve the classifier performance
- Throw away the classifier code and keep the embeddings.

# Classifier components

Machine learning classifiers require a training corpus of  $M$  observations input/output pairs  $(x^{(i)}, y^{(i)})$ .

1. A **feature representation** of the input. For each input observation  $x^{(i)}$ , this will be a vector of features  $[x_1, x_2, \dots, x_n]$ .
2. A **classification function** that computes the estimated class  $\hat{y}$  via  $p(y|x)$ .
3. An **objective function** for learning, usually involving minimizing error on training examples.
4. An algorithm for **optimizing** the objective function.

# Sentiment classifier

For sentiment classification, consider an input observation  $x$ , represented by a vector of features  $[x_1, x_2, \dots, x_n]$ . The classifier output  $y$  can be 1 (positive sentiment) or 0 (negative sentiment). We want to estimate  $P(y = 1 | x)$ .

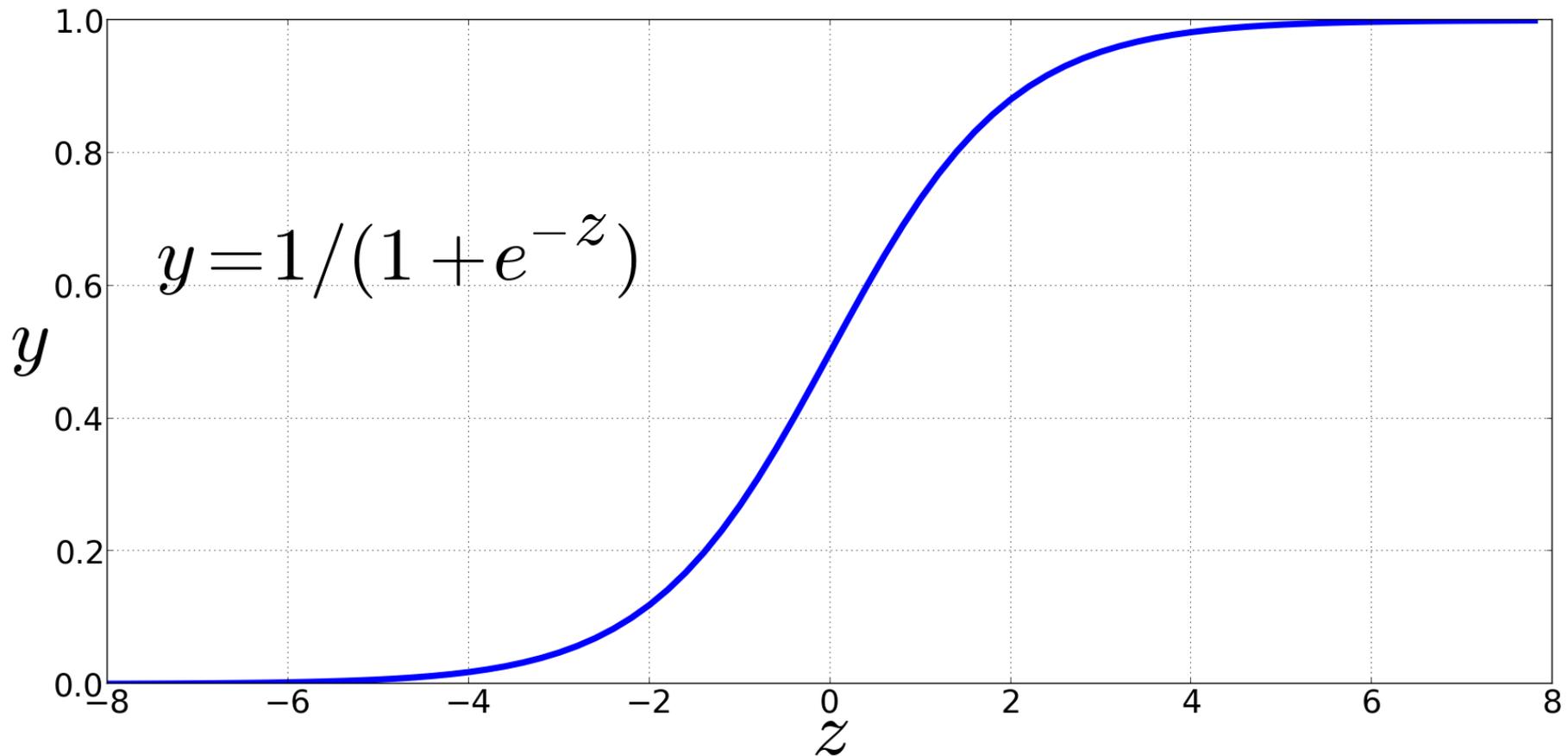
Logistic regression solves this task by learning, from a training set, a vector of **weights** and a **bias term**.

$$z = \left( \sum_{i=1}^n w_i x_i \right) + b$$

We can also write this as a dot product:

$$z = w \cdot x + b$$

# Sigmoid function



# Probabilities

$$\begin{aligned} P(y = 1) &= \sigma(w \cdot x + b) \\ &= \frac{1}{1 + e^{-(w \cdot x + b)}} \end{aligned}$$

# Decision boundary

Now we have an algorithm that given an instance  $x$  computes the probability  $P(y = 1 | x)$ . How do we make a decision?

$$\hat{y} = \begin{cases} 1 & \text{if } P(y = 1 | x) > 0.5 \\ 0 & \text{otherwise} \end{cases}$$

For a test instance  $x$ , we say yes if the probability  $P(y = 1 | x)$  is more than .5, and no otherwise. We call .5 the decision boundary

# Extracting Features

It's hokey. There are virtually no surprises , and the writing is second-rate . So why was it so enjoyable? For one thing , the cast is great . Another nice touch is the music . I was overcome with the urge to get off the couch and start dancing . It sucked me in , and it'll do the same to you .

Var	Definition	Value
$x_1$	Count of positive lexicon words	
$x_2$	Count of negative lexicon words	
$x_3$	Does no appear? (binary feature)	
$x_4$	Number of 1 <sup>st</sup> and 2 <sup>nd</sup> person pronouns	
$x_5$	Does ! appear? (binary feature)	
$x_6$	Log of the word count for the document	

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Var	Definition	Value
$x_1$	Count of positive lexicon words	3
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$x_3$	Does no appear? (binary feature)	1
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Word count = 64,  $\ln(64) = 4.15$

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$x_3$	Does no appear? (binary feature)	1
$x_4$	Number of 1 <sup>st</sup> and 2 <sup>nd</sup> person pronouns	3
$x_5$	Does ! appear? (binary feature)	0
$x_6$	Log of the word count for the document	4.15

Var	Definition	Value	Weight	Product
$x_1$	Count of positive lexicon words	3	2.5	
$x_2$	Count of negative lexicon words	2	-5.0	
$x_3$	Does no appear? (binary feature)	1	-1.2	
$x_4$	Num 1 <sup>st</sup> and 2nd person pronouns	3	0.5	
$x_5$	Does ! appear? (binary feature)	0	2.0	
$x_6$	Log of the word count for the doc	4.15	0.7	
$b$	bias	1	0.1	

$$z = \left( \sum_{i=1}^n w_i x_i \right) + b$$

# Computing Z

Var	Definition	Value	Weight	Product
$x_1$	Count of positive lexicon words	3	2.5	7.5
$x_2$	Count of negative lexicon words	2	-5.0	-10
$x_3$	Does no appear? (binary feature)	1	-1.2	-1.2
$x_4$	Num 1 <sup>st</sup> and 2nd person pronouns	3	0.5	1.5
$x_5$	Does ! appear? (binary feature)	0	2.0	0
$x_6$	Log of the word count for the doc	4.15	0.7	2.905
b	bias	1	0.1	.1

$$z = \left( \sum_{i=1}^n w_i x_i \right) + b$$

$$Z=0.805$$

# Sigmoid(Z)

Var	Definition	Value	Weight	Product
$x_1$	Count of positive lexicon words	3	2.5	7.5
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$x_3$	Does no appear? (binary feature)	1	-1.2	-1.2
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$x_5$	Does ! appear? (binary feature)	0	2.0	0
$x_6$	Log of the	0.7	4.2	2.905
b	bias	0.1	0.1	.1



$$\sigma(0.805) = 0.69$$

# Learning in logistic regression

How do we get the weights of the model? We learn the parameters (weights + bias) via learning. This requires 2 components:

1. An objective function or **loss function** that tells us *distance* between the system output and the gold output. We will use **cross-entropy loss**.
2. An algorithm for optimizing the objective function. We will use stochastic gradient descent to **minimize** the **loss function**.

# Loss functions

We need to determine for some observation  $x$  how close the classifier output ( $\hat{y} = \sigma(w \cdot x + b)$ ) is to the correct output ( $y$ , which is 0 or 1).

$L(\hat{y}, y)$  = how much  $\hat{y}$  differs from the true  $y$

One example is mean squared error

$$L_{MSE}(\hat{y}, y) = \frac{1}{2} (\hat{y} - y)^2$$

# Loss functions for probabilistic classification

We use a loss function that prefers the correct class labels of the training example to be more likely.

Conditional maximum likelihood estimation: Choose parameters  $w, b$  that maximize the (log) probabilities of the true labels in the training data.

The resulting loss function is the negative log likelihood loss, more commonly called the **cross entropy loss**.

# Loss functions for probabilistic classification

For one observation  $x$ , let's **maximize** the probability of the correct label  $p(y|x)$ .

$$p(y|x) = \hat{y}^y (1 - \hat{y})^{1-y}$$

If  $y = 1$ , then  $p(y|x) = \hat{y}$ .

If  $y = 0$ , then  $p(y|x) = 1 - \hat{y}$ .

# Loss functions for probabilistic classification

Change to logs (still maximizing)

$$\begin{aligned}\log p(y|x) &= \log[\hat{y}^y (1 - \hat{y})^{1-y}] \\ &= y \log \hat{y} + (1 - y) \log(1 - \hat{y})\end{aligned}$$

This tells us what log likelihood should be maximized. But for loss functions, we want to minimize things, so we'll flip the sign.

# Cross-entropy loss

The result is cross-entropy loss:

$$L_{CE}(\hat{y}, y) = -\log p(y|x) = -[y \log \hat{y} + (1 - y) \log(1 - \hat{y})]$$

Finally, plug in the definition for  $\hat{y} = \sigma(w \cdot x + b)$

$$L_{CE}(\hat{y}, y) = -[y \log \sigma(w \cdot x + b) + (1 - y) \log(1 - \sigma(w \cdot x + b))]$$

# Cross-entropy loss

Why does minimizing this negative log probability do what we want?

A perfect classifier would assign probability 1 to the correct outcome ( $y=1$  or  $y=0$ ) and probability 0 to the incorrect outcome.

That means the higher  $\hat{y}$  (the closer it is to 1), the better the classifier; the lower  $\hat{y}$  is (the closer it is to 0), the worse the classifier.

The negative log of this probability is a convenient loss metric since it goes from 0 (negative log of 1, no loss) to infinity (negative log of 0, infinite loss).

# Loss on all training examples

$$\begin{aligned}\log p(\text{training labels}) &= \log \prod_{i=1}^m p(y^{(i)} | x^{(i)}) \\ &= \sum_{i=1}^m \log p(y^{(i)} | x^{(i)}) \\ &= - \sum_{i=1}^m L_{CE}(\hat{y}^{(i)}, y^{(i)})\end{aligned}$$

# Average Loss

$$\begin{aligned} \text{Cost}(w, b) &= \frac{1}{m} \sum_{i=1}^m L_{CE}(\hat{y}^{(i)}, y^{(i)}) \\ &= -\frac{1}{m} \sum_{i=1}^m y^{(i)} \log \sigma(w \cdot x^{(i)} + b) + (1 - y^{(i)}) \log (1 - \sigma(w \cdot x^{(i)} + b)) \end{aligned}$$

**This is what we want to minimize!!**

# Finding good parameters

We use gradient descent to find good settings for our weights and bias by minimizing the loss function.

$$\hat{\theta} = \underset{\theta}{\operatorname{argmin}} \frac{1}{m} \sum_{i=1}^m L_{CE}(y^{(i)}, x^{(i)}; \theta)$$

Gradient descent is a method that finds a minimum of a function by figuring out in which direction (in the space of the parameters  $\theta$ ) the function's slope is rising the most steeply, and moving in the opposite direction.

# Gradient descent



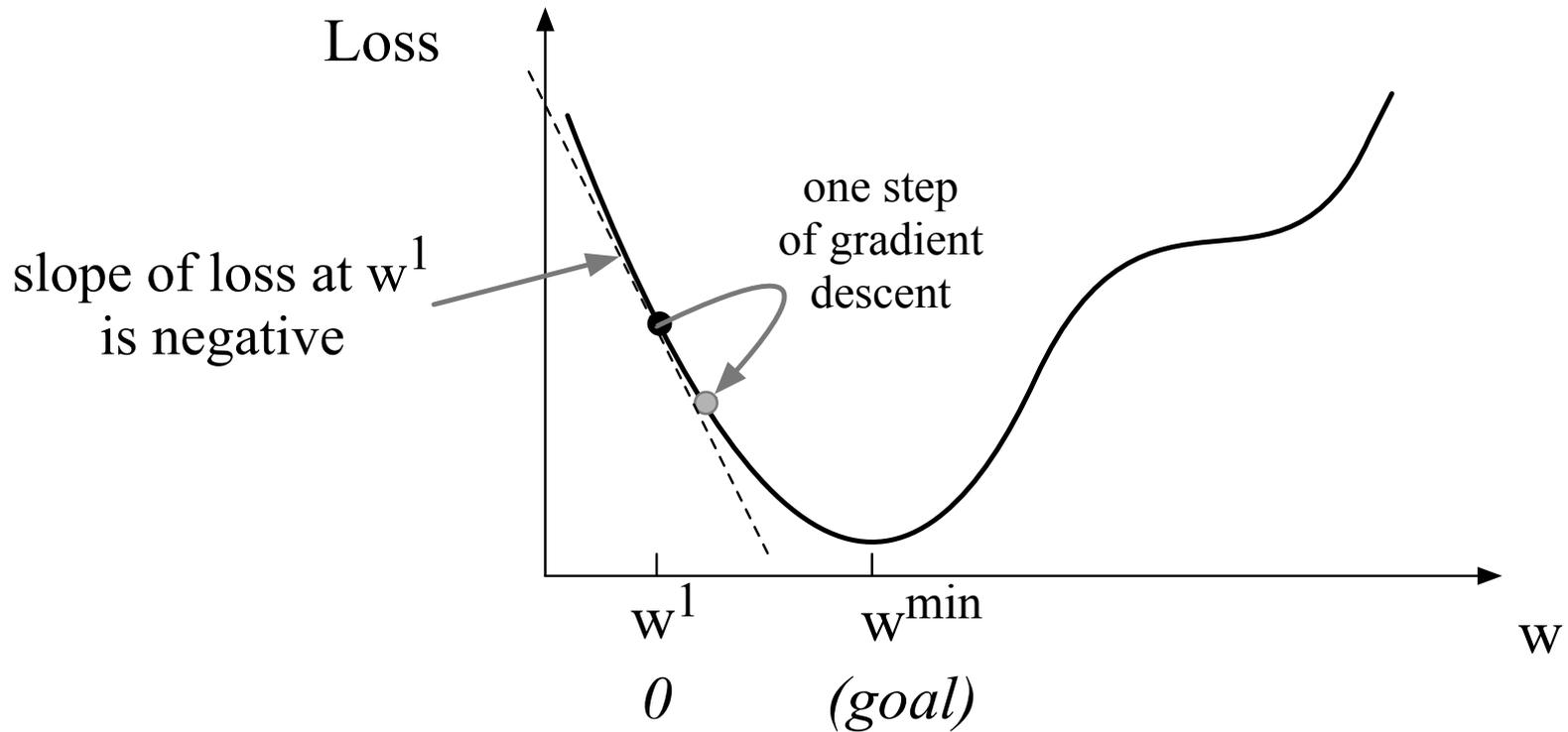
# Global v. Local Minimums

For logistic regression, this loss function is conveniently convex.

A convex function has just one minimum, so there are no local minima to get stuck in.

So gradient descent starting from any point is guaranteed to find the minimum.

# Iteratively find minimum



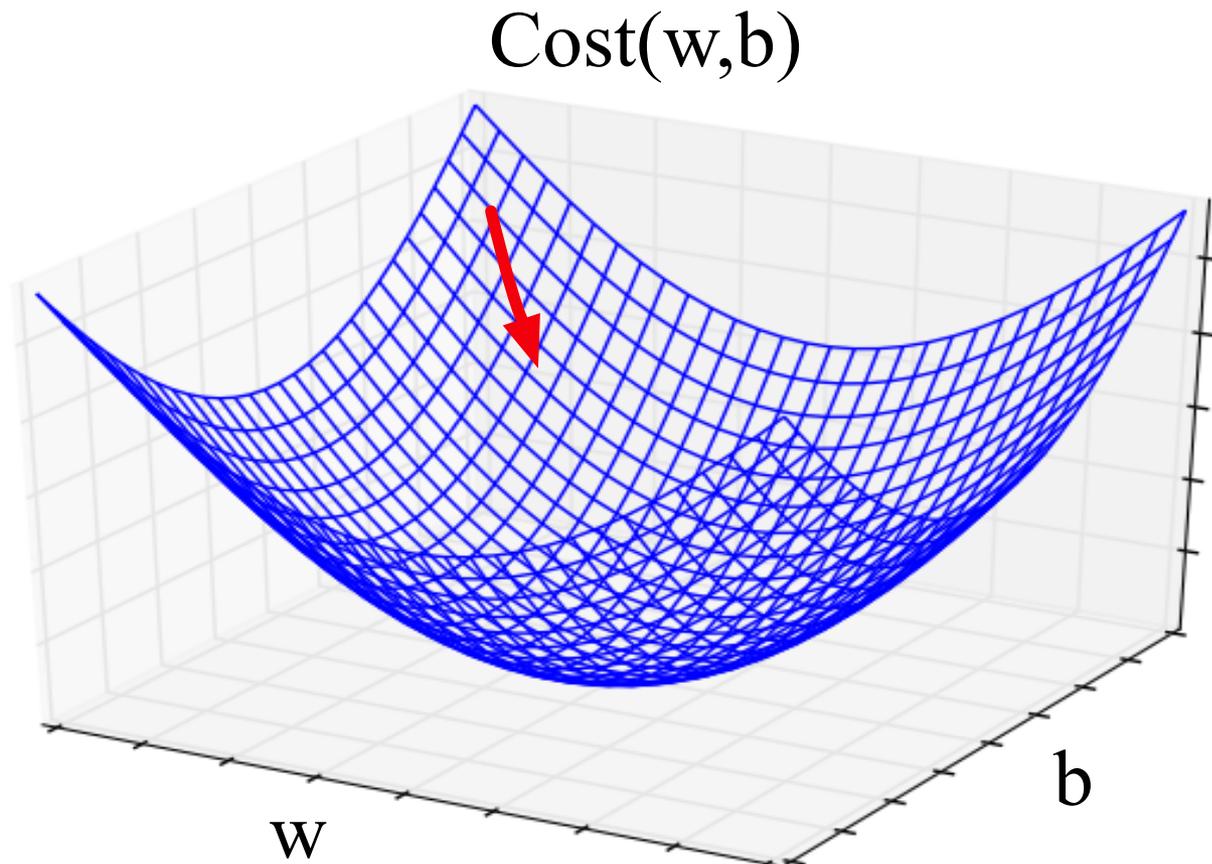
# How much should we update the parameter by?

The magnitude of the amount to move in gradient descent is the value of the slope weighted by a learning rate  $\eta$ .

A higher/faster learning rate means that we should move  $w$  more on each step.

$$w^{t+1} = w^t - \eta \frac{d}{dw} f(x; w)$$

# Many dimensions



# Updating each dimension $w_i$

$$\nabla_{\theta} L(f(x; \theta), y) = \begin{bmatrix} \frac{\partial}{\partial w_1} L(f(x; \theta), y) \\ \frac{\partial}{\partial w_2} L(f(x; \theta), y) \\ \vdots \\ \frac{\partial}{\partial w_n} L(f(x; \theta), y) \end{bmatrix}$$

The final equation for updating  $\theta$  based on the gradient is

$$\theta_{t+1} = \theta_t - \eta \nabla L(f(x; \theta), y)$$

# The Gradient

To update  $\theta$ , we need a definition for the gradient  $\nabla L(f(x; \theta), y)$ .

For logistic regression the cross-entropy loss function is:

$$L_{CE}(w, b) = -[y \log \sigma(w \cdot x + b) + (1 - y) \log (1 - \sigma(w \cdot x + b))]$$

The derivative of this function for one observation vector  $x$  for a single weight  $w^j$  is

$$\frac{\partial L_{CE}(w, b)}{\partial w_j} = [\sigma(w \cdot x + b) - y]x_j$$

The gradient is a very intuitive value: the difference between the true  $y$  and our estimate for  $x$ , multiplied by the corresponding input value  $x_j$ .

# The Gradient

The loss for a batch of data or an entire dataset is just the average loss over the  $m$  examples

$$\text{Cost}(w, b) = -\frac{1}{m} \sum_{i=1}^m y^{(i)} \log \sigma(w \cdot x^{(i)} + b) + (1 - y^{(i)}) \log (1 - \sigma(w \cdot x^{(i)} + b))$$

The gradient for multiple data points is the sum of the individual gradients:

$$\frac{\partial \text{Cost}(w, b)}{\partial w_j} = \sum_{i=1}^m \left[ \sigma(w \cdot x^{(i)} + b) - y^{(i)} \right] x_j^{(i)}$$

# Stochastic gradient descent

## algorithm

**function** STOCHASTIC GRADIENT DESCENT( $L()$ ,  $f()$ ,  $x$ ,  $y$ ) **returns**  $\theta$

# where:  $L$  is the loss function

#  $f$  is a function parameterized by  $\theta$

#  $x$  is the set of training inputs  $x^{(1)}, x^{(2)}, \dots, x^{(n)}$

#  $y$  is the set of training outputs (labels)  $y^{(1)}, y^{(2)}, \dots, y^{(n)}$

$\theta \leftarrow 0$

**repeat**  $T$  times

For each training tuple  $(x^{(i)}, y^{(i)})$  (in random order)

Compute  $\hat{y}^{(i)} = f(x^{(i)}; \theta)$  # What is our estimated output  $\hat{y}$ ?

Compute the loss  $L(\hat{y}^{(i)}, y^{(i)})$  # How far off is  $\hat{y}^{(i)}$  from the true output  $y^{(i)}$ ?

$g \leftarrow \nabla_{\theta} L(f(x^{(i)}; \theta), y^{(i)})$  # How should we move  $\theta$  to maximize loss ?

$\theta \leftarrow \theta - \eta g$  # go the other way instead

**return**  $\theta$

# Worked example

# Neural Networks

The building block of a neural network is a single computational unit. A unit takes a set of real valued numbers as input, performs some computation.

