Regression and Classification with Neural Networks

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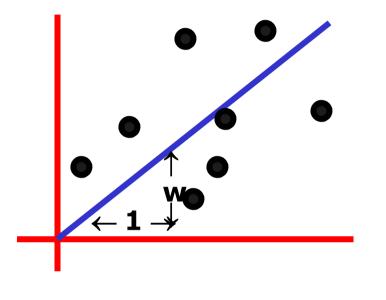
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Sep 25th, 2001

DATASET



inputs	outputs	
$x_1 = 1$	$y_1 = 1$	
$x_2 = 3$	$y_2 = 2.2$	
$x_3 = 2$	$y_3 = 2$	
$x_4 = 1.5$	$y_4 = 1.9$	
$x_5 = 4$	$y_5 = 3.1$	

Linear regression assumes that the expected value of the output given an input, E[y/x], is linear. Simplest case: Out(x) = wx for some unknown w.

Given the data, we can estimate *w*.

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1-parameter linear regression

Assume that the data is formed by

 $y_i = wx_i + noise_i$

where...

- the noise signals are independent
- the noise has a normal distribution with mean 0 and unknown variance σ^2

P(y|w,x) has a normal distribution with

- mean *wx*
- variance σ^2

Bayesian Linear Regression $P(y|w,x) = Normal (mean wx, var \sigma^2)$

We have a set of datapoints $(x_1, y_1) (x_2, y_2) \dots (x_n, y_n)$ which are EVIDENCE about *w*.

We want to infer *w* from the data.

 $\mathsf{P}(W|X_{1'}, X_{2'}, X_{3'}...X_{n'}, Y_{1'}, Y_{2'}...Y_{n})$

•You can use BAYES rule to work out a posterior distribution for *w* given the data.

•Or you could do Maximum Likelihood Estimation

Maximum likelihood estimation of w

Asks the question:

"For which value of *w* is this data most likely to have happened?"

<=>

For what w is

 $P(y_1, y_2...y_n | x_1, x_2, x_3,...x_n, w)$ maximized?

<=>

For what *w* is

$$\prod_{i=1}^{n} P(y_i | w, x_i) \text{ maximized}$$

For what *w* is

$$\prod_{i=1}^{n} P(y_i | w, x_i) \text{ maximized?}$$

For what w is $\prod_{i=1}^{n} \exp(-\frac{1}{2}(\frac{y_i - wx_i}{\sigma})^2) \text{ maximized}?$

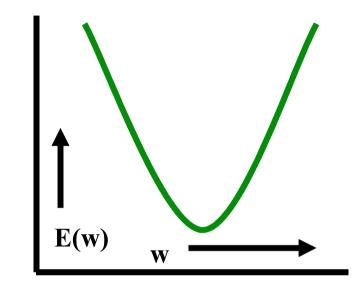
For what *w* is

$$\sum_{i=1}^{n} -\frac{1}{2} \left(\frac{y_i - wx_i}{\sigma} \right)^2 \text{ maximized?}$$

For what *w* is

 $\sum_{i=1}^{n} (y_i - wx_i)^2 \text{ minimized?}$

The maximum likelihood *w* is the one that minimizes sumof-squares of <u>residuals</u>



$$E = \sum_{i} (y_{i} - wx_{i})^{2}$$
$$= \sum_{i} y_{i}^{2} - (2\sum_{i} x_{i}y_{i})w + (\sum_{i} x_{i}^{2})w^{2}$$

We want to minimize a quadratic function of *w*.

Easy to show the sum of squares is minimized

when $w = \frac{\sum x_i y_i}{\sum x_i^2}$

The maximum likelihood model is Out(x) = wx

We can use it for prediction

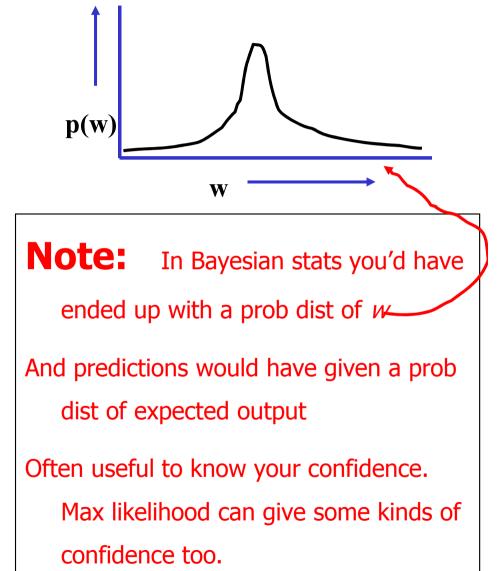
Easy to show the sum of squares is minimized

when

$$w = \frac{\sum x_i y_i}{\sum x_i^2}$$

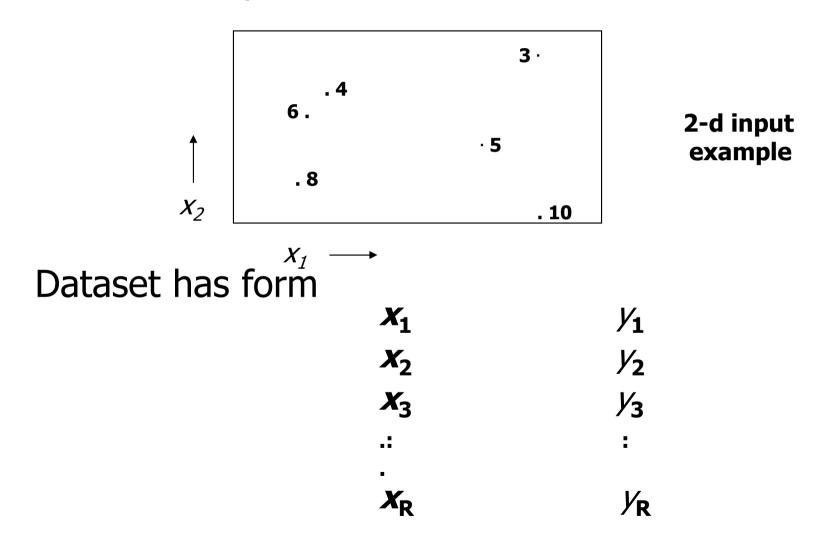
The maximum likelihood model is Out(x) = wx

We can use it for prediction



Multivariate Regression

What if the inputs are vectors?



Multivariate Regression

Write matrix X and Y thus:

$$\mathbf{x} = \begin{bmatrix} \dots & \mathbf{x}_{1} \dots & \mathbf{x}_{1} \\ \dots & \mathbf{x}_{2} \dots & \mathbf{x}_{2} \\ \vdots & \vdots & \vdots \\ \dots & \mathbf{x}_{R} \dots & \mathbf{x}_{R} \end{bmatrix} = \begin{bmatrix} x_{11} & x_{12} & \dots & x_{1m} \\ x_{21} & x_{22} & \dots & x_{2m} \\ \vdots & \vdots & \vdots \\ x_{R1} & x_{R2} & \dots & x_{Rm} \end{bmatrix} \mathbf{y} = \begin{bmatrix} y_{1} \\ y_{2} \\ \vdots \\ y_{R} \end{bmatrix}$$

(there are *R* datapoints. Each input has *m* components) The linear regression model assumes a vector **w** such that $Out(\mathbf{x}) = \mathbf{w}^{T}\mathbf{x} = w_{1}x[1] + w_{2}x[2] + \dots w_{m}x[D]$ The max. likelihood **w** is $\mathbf{w} = (X^{T}X)^{-1}(X^{T}Y)$

Multivariate Regression

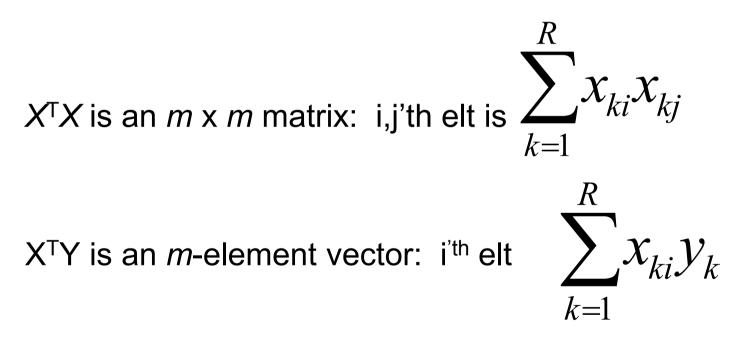
Write matrix X and Y thus:

$$\mathbf{x} = \begin{bmatrix} \dots \mathbf{x}_{1} \dots \mathbf{x}_{2} \dots \\ \dots \mathbf{x}_{2} \dots \\ \vdots \\ \dots \mathbf{x}_{R} \dots \end{bmatrix} = \begin{bmatrix} x_{11} & x_{12} & \dots & x_{1m} \\ x_{21} & x_{22} & \dots & x_{2m} \\ \vdots & \vdots & \vdots \\ x_{R1} & x_{R2} & \dots & x_{Rm} \end{bmatrix} \mathbf{y} = \begin{bmatrix} y_{1} \\ y_{2} \\ \vdots \\ y_{R} \end{bmatrix}$$

(there are *R* datapoints. Each input **IMPORTANT EXERCISE: PROVE IT !!!!!** The linear regression model assumes a vector *w* such that $Out(x) = w^{T}x = w_{1}x[1] + w_{2}x[2] +w_{m}x[D]$ The max. likelihood *w* is $w = (X^{T}X)^{-1}(X^{T}Y)$

Multivariate Regression (con't)

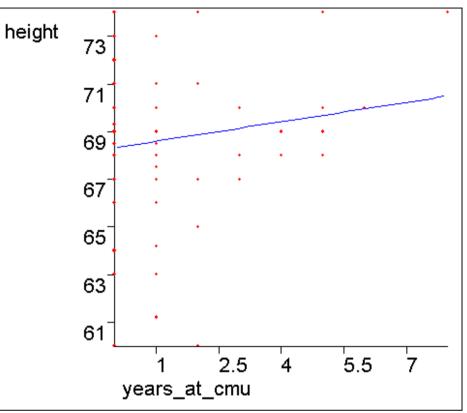
The max. likelihood **w** is $\mathbf{w} = (X^T X)^{-1} (X^T Y)$



What about a constant term?

We may expect linear data that does not go through the origin.

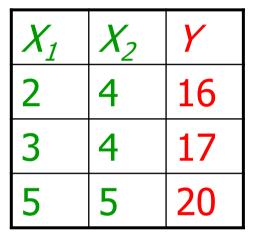
Statisticians and Neural Net Folks all agree on a simple obvious hack.



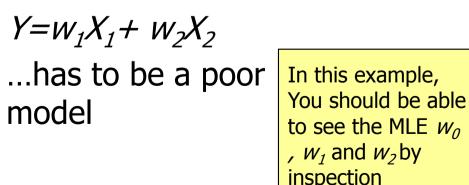
Can you guess??

The constant term

• The trick is to create a fake input " X_0 " that always takes the value 1



Before:



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X ₀	<i>X</i> ₁	<i>X</i> ₂	Y
1	2	4	16
1	3	4	17
1	5	5	20

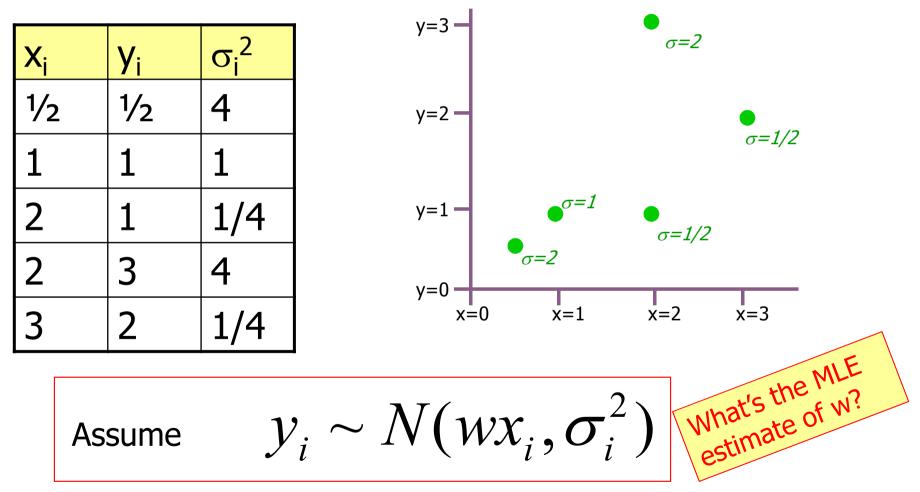
After:

- $Y = w_0 X_0 + w_1 X_1 + w_2 X_2$
- $= W_0 + W_1 X_1 + W_2 X_2$

...has a fine constant term

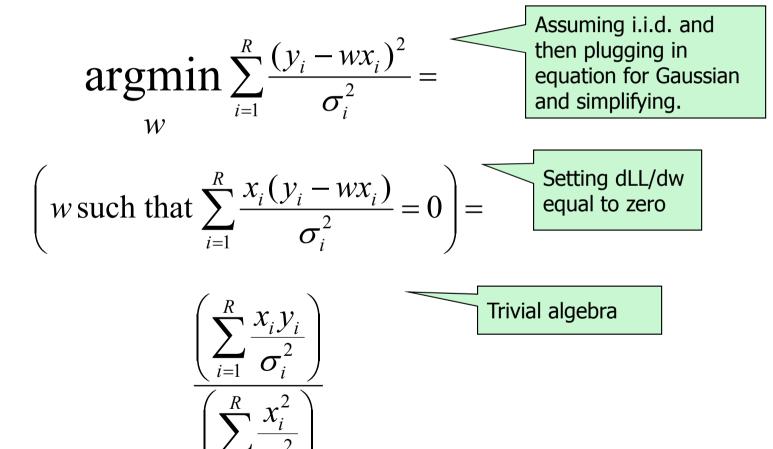
Regression with varying noise

• Suppose you know the variance of the noise that was added to each datapoint.



MLE estimation with varying noise

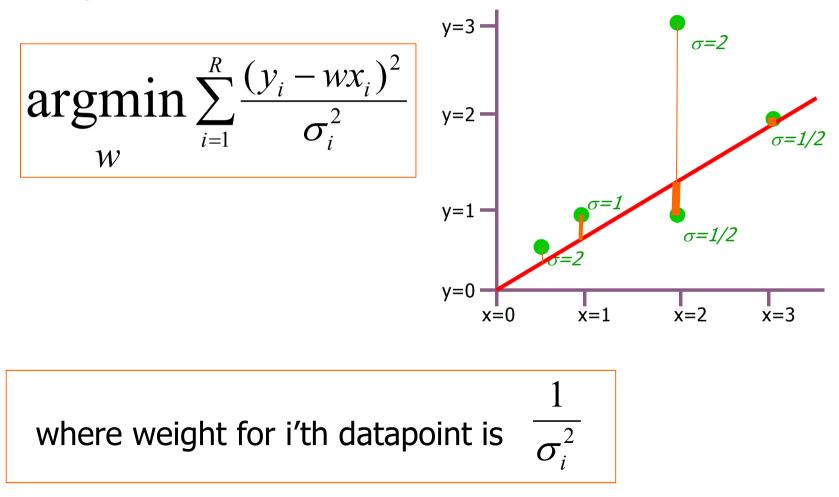
argmax log
$$p(y_1, y_2, ..., y_R | x_1, x_2, ..., x_R, \sigma_1^2, \sigma_2^2, ..., \sigma_R^2, w) =$$



 \mathcal{W}

This is Weighted Regression

• We are asking to minimize the weighted sum of squares

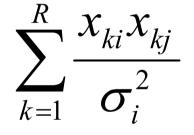


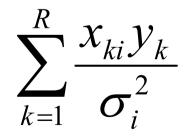
Weighted Multivariate Regression

The max. likelihood \boldsymbol{w} is $\boldsymbol{w} = (WX^TWX)^{-1}(WX^TWY)$

 $(WX^{T}WX)$ is an $m \ge m$ matrix: i,j'th elt is

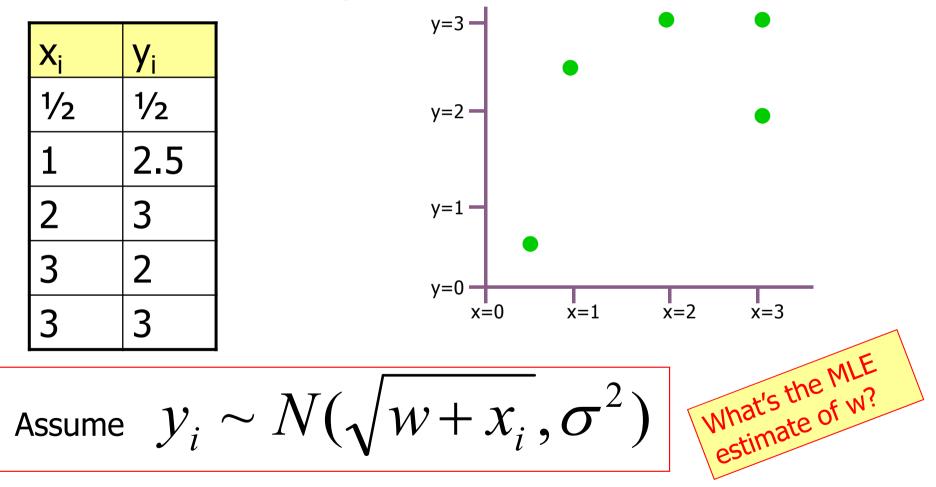
 (WX^TWY) is an *m*-element vector: i'th elt





Non-linear Regression

 Suppose you know that y is related to a function of x in such a way that the predicted values have a non-linear dependence on w, e.g:



Non-linear MLE estimation

$$\operatorname{argmax} \log p(y_1, y_2, ..., y_R | x_1, x_2, ..., x_R, \sigma, w) =$$

$$\operatorname{w}_{W} \operatorname{argmin}_{W} \sum_{i=1}^{R} (y_i - \sqrt{w + x_i})^2 = \operatorname{Assuming i.i.d. and then plugging in equation for Gaussian and simplifying.}$$

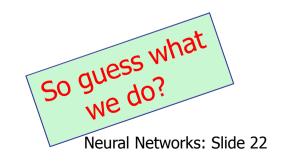
$$\left(w \operatorname{such that} \sum_{i=1}^{R} \frac{y_i - \sqrt{w + x_i}}{\sqrt{w + x_i}} = 0 \right) = \operatorname{Setting dLL/dw}_{equal to zero}$$

Non-linear MLE estimation

argmax log
$$p(y_1, y_2, ..., y_R | x_1, x_2, ..., x_R, \sigma, w) =$$

^W
argmin $\sum_{i=1}^{R} (y_i - \sqrt{w + x_i})^2 =$
^W
Assuming i.i.d. and then plugging in equation for Gaussian and simplifying.
(w such that $\sum_{i=1}^{R} \frac{y_i - \sqrt{w + x_i}}{\sqrt{w + x_i}} = 0$) = Setting dLL/dw equal to zero

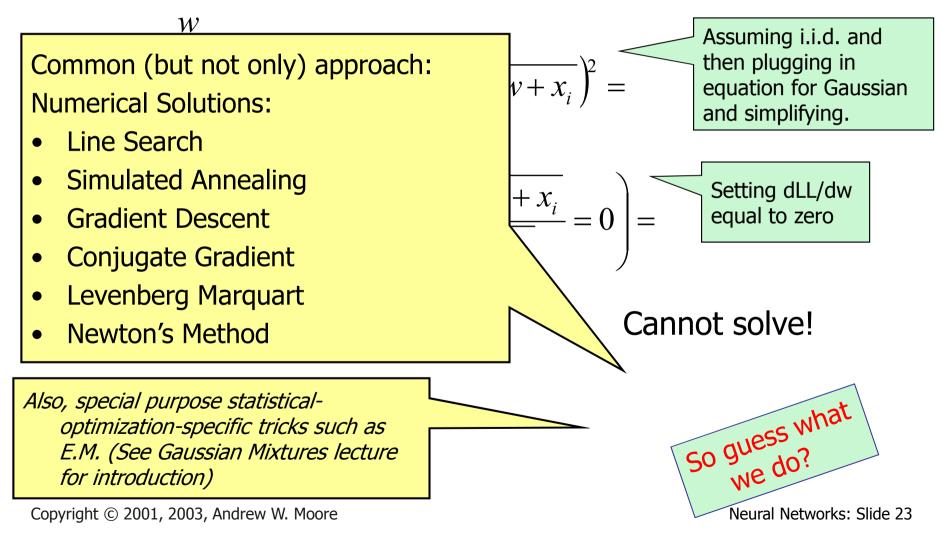
Cannot solve!



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Non-linear MLE estimation

argmax log
$$p(y_1, y_2, ..., y_R | x_1, x_2, ..., x_R, \sigma, w) =$$



GRADIENT DESCENT

Suppose we have a scalar function $f(w): \mathfrak{R} \to \mathfrak{R}$

We want to find a local minimum. Assume our current weight is *w*

GRADIENT DESCENT RULE: $w \leftarrow w - \eta \frac{\partial}{\partial w} f(w)$

 η is called the LEARNING RATE. A small positive number, e.g. $\eta = 0.05$

GRADIENT DESCENT

Suppose we have a scalar function $f(w): \mathfrak{R} \to \mathfrak{R}$

We want to find a local minimum. Assume our current weight is *w*

GRADIENT DESCENT RULE: $w \leftarrow w - \eta \frac{\partial}{\partial w} f(w)$ Recall Andrew's favorite default value for anything η is called the LEARNING FOR A Small positive number, e.g. $\eta = 0.05$ QUESTION: Justify the Gradient Descent Rule

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Gradient Descent in "m" Dimensions

Given $f(\mathbf{w}): \mathfrak{R}^m \to \mathfrak{R}$

 $\nabla f(w) = \begin{pmatrix} \frac{\partial}{\partial w_1} f(w) \\ \vdots \\ \frac{\partial}{\partial w_m} f(w) \end{pmatrix} \text{ points in direction of steepest ascent.}$

 $|\nabla f(w)|$ is the gradient in that direction

GRADIENT DESCENT RULE: $w \leftarrow w - \eta \nabla f(w)$

Equivalently

$$w_j \leftarrow w_j - \eta \frac{\partial}{\partial w_j} f(w)$$

....where w_j is the *j*th weight "just like a linear feedback system"

What's all this got to do with Neural Nets, then, eh??

For supervised learning, neural nets are also models with vectors of **w** parameters in them. They are now called weights.

As before, we want to compute the weights to minimize sumof-squared residuals.

Which turns out, under "Gaussian i.i.d noise" assumption to be max. likelihood.

Instead of explicitly solving for max. likelihood weights, we use **GRADIENT DESCENT** to **SEARCH** for them.

"Why?" you ask, a querulous expression in your eyes. "Aha!!" I reply: "We'll see later."

Linear Perceptrons

They are multivariate linear models:

$$\operatorname{Out}(\mathbf{x}) = \mathbf{w}^{\mathsf{T}}\mathbf{x}$$

And "training" consists of minimizing sum-of-squared residuals by gradient descent.

$$E = \sum_{k} (\text{Out } (\mathbf{x}_{k}) - y_{k})^{2}$$
$$= \sum_{k} (\mathbf{w}^{T} \mathbf{x}_{k} - y_{k})^{2}$$

QUESTION: Derive the perceptron training rule.

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Linear Perceptron Training Rule

$$E = \sum_{k=1}^{R} (y_k - \mathbf{w}^T \mathbf{x}_k)^2$$

Gradient descent tells us we should update **w** thusly if we wish to minimize *E*:

$$w_j \leftarrow w_j - \eta \frac{\partial E}{\partial w_j}$$

So what's $\frac{\partial E}{\partial w_i}$?

Linear Perceptron Training Rule

$$E = \sum_{k=1}^{R} (y_k - \mathbf{w}^T \mathbf{x}_k)^2$$

Gradient descent tells us we should update **w** thusly if we wish to minimize *E*:

$$w_{j} \leftarrow w_{j} - \eta \frac{\partial E}{\partial w_{j}}$$

So what's $\frac{\partial E}{\partial w_{j}}$?

$$\frac{\partial E}{\partial w_j} = \sum_{k=1}^{R} \frac{\partial}{\partial w_j} (y_k - \mathbf{w}^T \mathbf{x}_k)^2$$
$$= \sum_{k=1}^{R} 2(y_k - \mathbf{w}^T \mathbf{x}_k) \frac{\partial}{\partial w_j} (y_k - \mathbf{w}^T \mathbf{x}_k)$$
$$= -2 \sum_{k=1}^{R} \delta_k \frac{\partial}{\partial w_j} \mathbf{w}^T \mathbf{x}_k$$
$$(\dots, where)$$
$$\delta_k = y_k - \mathbf{w}^T \mathbf{x}_k$$
$$= -2 \sum_{k=1}^{R} \delta_k \frac{\partial}{\partial w_j} \sum_{i=1}^{m} w_i x_{ki}$$
$$= -2 \sum_{k=1}^{R} \delta_k \frac{\partial}{\partial w_j} \sum_{i=1}^{m} w_i x_{ki}$$

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Linear Perceptron Training Rule

$$E = \sum_{k=1}^{R} (y_k - \mathbf{w}^T \mathbf{x}_k)^2$$

Gradient descent tells us we should update **w** thusly if we wish to minimize *E*:

$$w_{j} \leftarrow w_{j} - \eta \frac{\partial E}{\partial w_{j}}$$

...where...
$$\frac{\partial E}{\partial w_{j}} = -2\sum_{k=1}^{R} \delta_{k} x_{kj}$$

$$w_{j} \leftarrow w_{j} + 2\eta \sum_{k=1}^{R} \delta_{k} x_{kj}$$

We frequently neglect the 2 (meaning we halve the learning rate)

The "Batch" perceptron algorithm

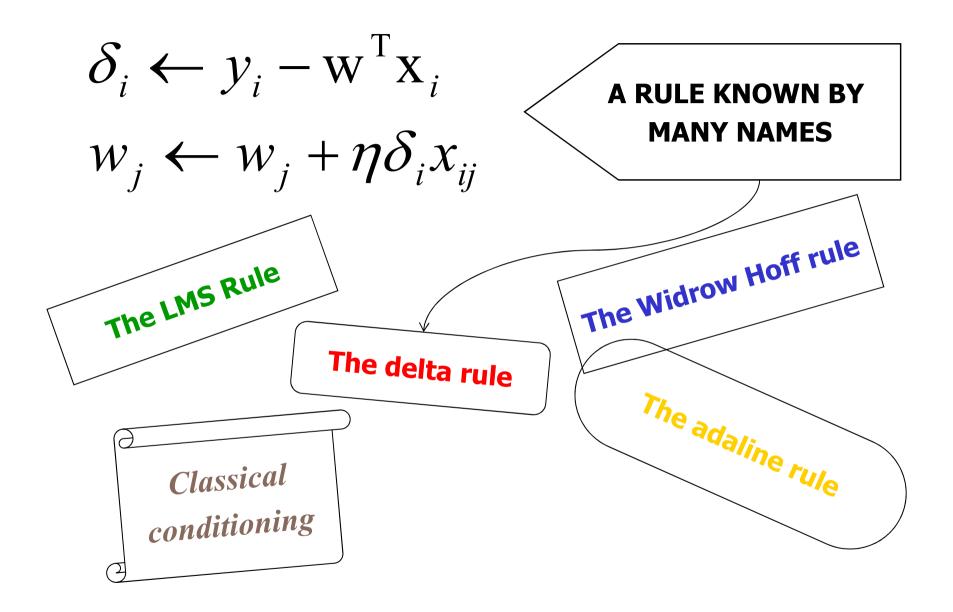
1) Randomly initialize weights w₁ w₂ ... w_m

2) Get your dataset (append 1's to the inputs if you don't want to go through the origin).

3) for
$$i = 1$$
 to R $\delta_i := y_i - \mathbf{W}^T \mathbf{X}_i$

- 4) for j = 1 to m $w_j \leftarrow w_j + \eta \sum_{i=1}^R \delta_i x_{ij}$
- 5) if $\sum \delta_i^2$ stops improving then stop. Else loop back to 3.

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If data is voluminous and arrives fast

Input-output pairs (\mathbf{x}, \mathbf{y}) come streaming in very quickly. THEN

Don't bother remembering old ones. Just keep using new ones.

observe
$$(\mathbf{x}, \mathbf{y})$$

 $\delta \leftarrow \mathbf{y} - \mathbf{w}^{\mathrm{T}} \mathbf{x}$
 $\forall j \ w_{j} \leftarrow w_{j} + \eta \, \delta \, x_{j}$

Gradient Descent vs Matrix Inversion for Linear Perceptrons GD Advantages (MI disadvantages):

- GD Disadvantages (MI advantages):
- •
- •

Gradient Descent vs Matrix Inversion for Linear Perceptrons GD Advantages (MI disadvantages):

- Biologically plausible
- With very very many attributes each iteration costs only O(mR). If fewer than m iterations needed we've beaten Matrix Inversion
- More easily parallelizable (or implementable in wetware)?

GD Disadvantages (MI advantages):

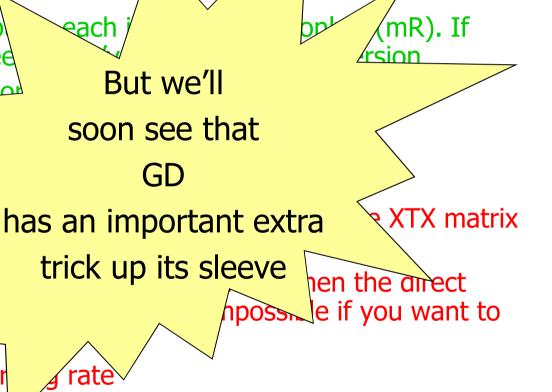
- It's moronic
- It's essentially a slow implementation of a way to build the XTX matrix and then solve a set of linear equations
- If m is small it's especially outageous. If m is large then the direct matrix inversion method gets fiddly but not impossible if you want to be efficient.
- Hard to choose a good learning rate
- Matrix inversion takes predictable time. You can't be sure when gradient descent will stop.

Gradient Descent vs Matrix Inversion for Linear Perceptrons GD Advantages (MI disadvantage):

- Biologically plausible
- With very very many attrib fewer than m iterations nee
- More easily parallelizable (or

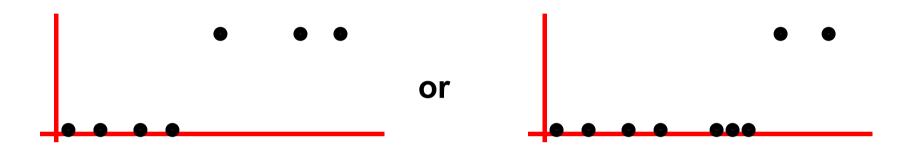
GD Disadvanta

- It's moronic
- It's essentially and then solve a second
- If m is small it's especimatrix inversion met be efficient.
- Hard to choose a good lear
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Perceptrons for Classification

What if all outputs are 0's or 1's ?



We can do a linear fit.

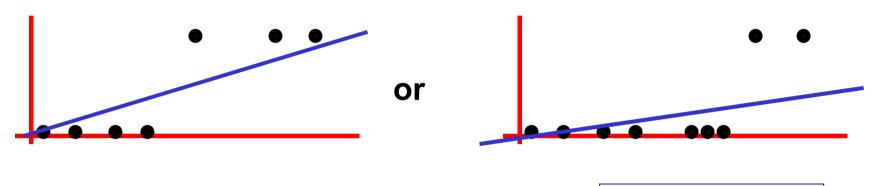
Our prediction is 0 if $out(\mathbf{x}) \le 1/2$

1 if out(*x*)>1/2

WHAT'S THE BIG PROBLEM WITH THIS???

Perceptrons for Classification

What if all outputs are 0's or 1's ?



We can do a linear fit.

Blue = Out(x)

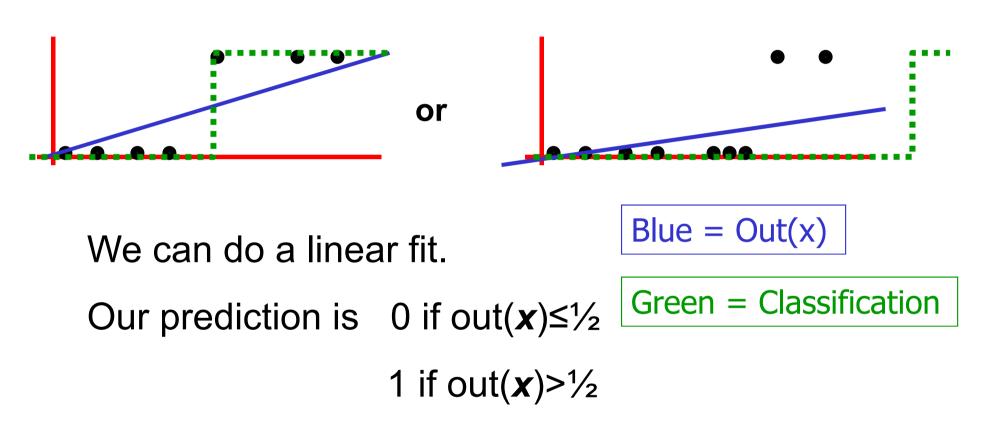
Our prediction is 0 if $out(\mathbf{x}) \le \frac{1}{2}$

1 if out(**x**)>¹/₂

WHAT'S THE BIG PROBLEM WITH THIS???

Perceptrons for Classification

What if all outputs are 0's or 1's ?



Classification with Perceptrons I

Don't minimize

$$\sum \left(y_i - \mathbf{W}^{\mathrm{T}} \mathbf{X}_i \right)^2.$$

Minimize number of misclassifications instead. [Assume outputs are

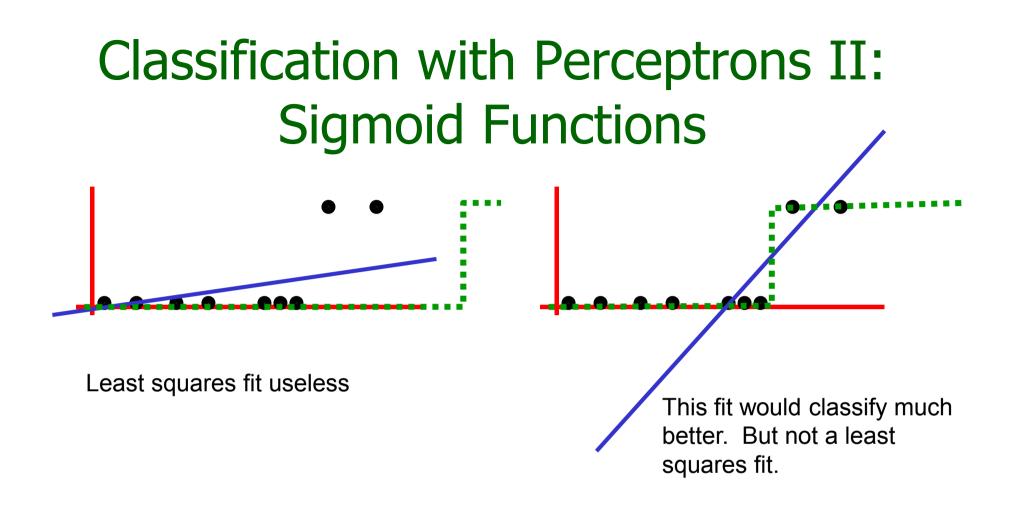
+1 & -1, not +1 & 0]
where Round(x) = -1 if x<0
1 if x \ge 0
$$\sum (y_i - \text{Round}(w^T x_i))$$
NOTE: CUTE & NON OBVIOUS WHY
THIS WORKS!!

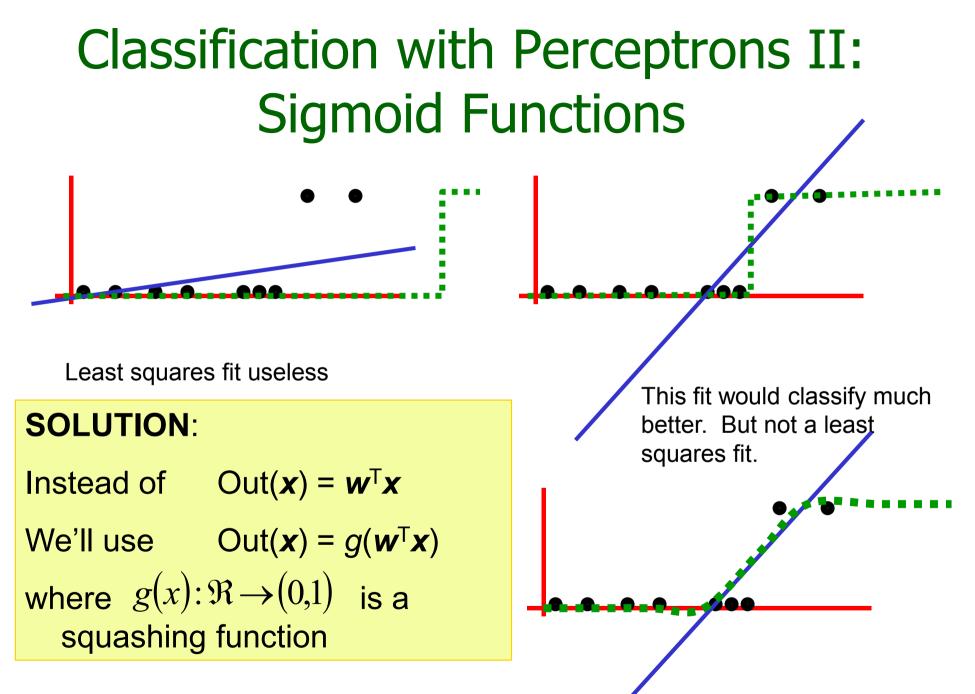
The gradient descent rule can be changed to:

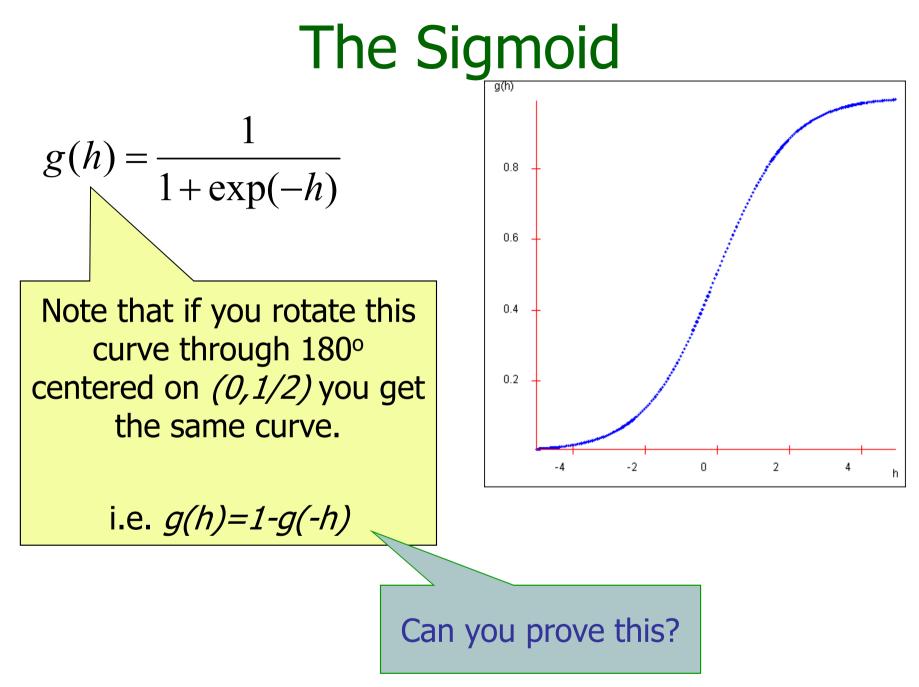
if $(\mathbf{x}_i, \mathbf{y}_i)$ correctly classed, don't change

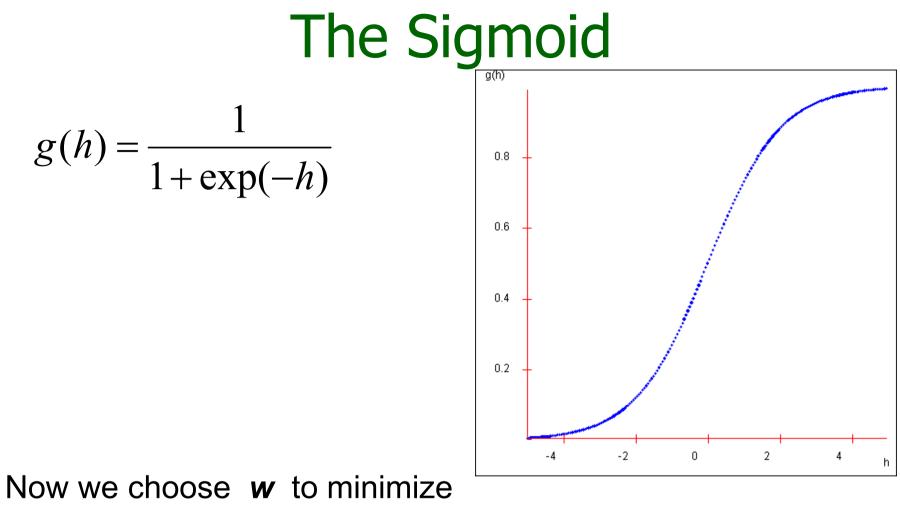
if wrongly predicted as 1 $w \leftarrow w - x_i$

if wrongly predicted as -1 $w \leftarrow w + x_i$





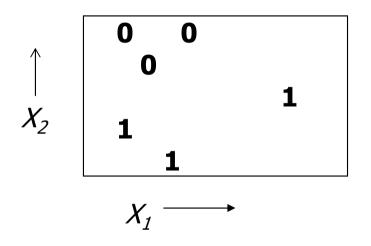




$$\sum_{i=1}^{R} [y_i - \text{Out}(\mathbf{x}_i)]^2 = \sum_{i=1}^{R} [y_i - g(\mathbf{w}^{\mathsf{T}}\mathbf{x}_i)]^2$$

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Linear Perceptron Classification Regions



We'll use the model $\operatorname{Out}(\mathbf{x}) = g(\mathbf{w}^{\mathsf{T}}(\mathbf{x},1))$ = $g(w_1x_1 + w_2x_2 + w_0)$

Which region of above diagram classified with +1, and which with 0 ??

Gradient descent with sigmoid on a perceptron

First, notice
$$g'(x) = g(x)(1-g(x))$$

Because: $g(x) = \frac{1}{1+e^{-x}}$ so $g'(x) = \frac{-e^{-x}}{(1+e^{-x})^2}$
 $= \frac{1-1-e^{-x}}{(1+e^{-x})^2} = \frac{1}{(1+e^{-x})^2} - \frac{1}{1+e^{-x}} = \frac{-1}{1+e^{-x}} \left(1 - \frac{1}{1+e^{-x}}\right) = -g(x)(1-g(x))$
Out $(x) = g\left(\sum_{k} w_k x_k\right)$
 $E = \sum_{i} \left(y_i - g\left(\sum_{k} w_k x_k\right)\right)^2$
 $\frac{\partial E}{\partial w_j} = \sum_{i} 2\left(y_i - g\left(\sum_{k} w_k x_k\right)\right) \left(-\frac{\partial}{\partial w_j} g\left(\sum_{k} w_k x_{ik}\right)\right)$
 $= \sum_{i} -2\left(y_i - g\left(\sum_{k} w_k x_k\right)\right)g'\left(\sum_{k} w_k x_{ik}\right)\frac{\partial}{\partial w_j}\sum_{k} w_k x_{ik}$
 $= \sum_{i} -2\delta_i g(\operatorname{net}_i)(1-g(\operatorname{net}_i))x_{ij}$
where $\delta_i = y_i - \operatorname{Out}(x_i)$ net $i = \sum w_k x_k$

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 $(g_i)x_{ij}$

Other Things about Perceptrons

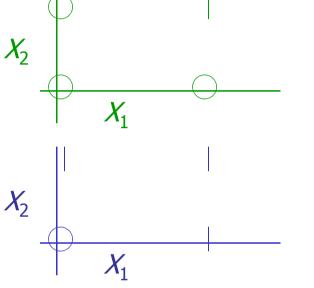
- Invented and popularized by Rosenblatt (1962)
- Even with sigmoid nonlinearity, correct convergence is guaranteed
- Stable behavior for overconstrained and underconstrained problems

Perceptrons and Boolean Functions

If inputs are all 0's and 1's and outputs are all 0's and 1's...

• Can learn the function $x_1 \wedge x_2$

Can learn the function $x_1 \vee x_2$.



• Can learn any conjunction of literals, e.g.

 $X_1 \wedge \gamma X_2 \wedge \gamma X_3 \wedge X_4 \wedge X_5$

QUESTION: WHY?

ullet

Perceptrons and Boolean Functions

• Can learn any disjunction of literals

e.g. $x_1 \wedge \neg x_2 \wedge \neg x_3 \wedge x_4 \wedge x_5$

• Can learn majority function

$$f(x_1, x_2 \dots x_n) = \begin{cases} 1 \text{ if } n/2 x_i \text{'s or more are} = 1 \\ 0 \text{ if less than } n/2 x_i \text{'s are} = 1 \end{cases}$$

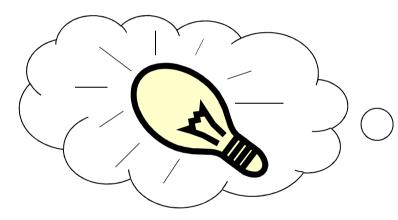
• What about the exclusive or function?

$$f(x_1, x_2) = x_1 \forall x_2 = (x_1 \land \neg x_2) \lor (\neg x_1 \land x_2)$$

Multilayer Networks

The class of functions representable by perceptrons is limited 1

$$\operatorname{Out}(\mathbf{x}) = g\left(\mathbf{w}^{\mathrm{T}}\mathbf{x}\right) = g\left(\sum_{j} w_{j} x_{j}\right)$$



^{OUse} a wider representation !

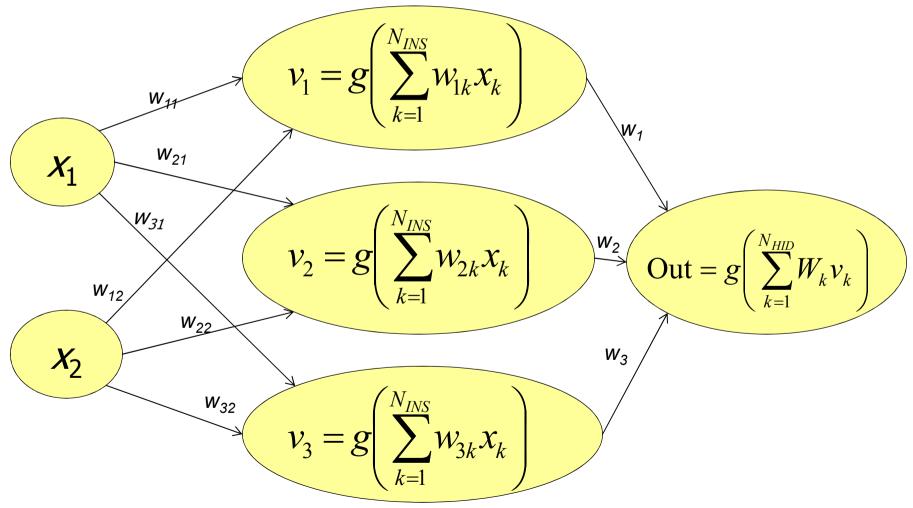
$$\operatorname{Out}(\mathbf{x}) = g\left(\sum_{j} W_{j}g\left(\sum_{k} W_{jk} x_{jk}\right)\right)$$

This is a nonlinear function Of a linear combination Of non linear functions Of linear combinations of inputs

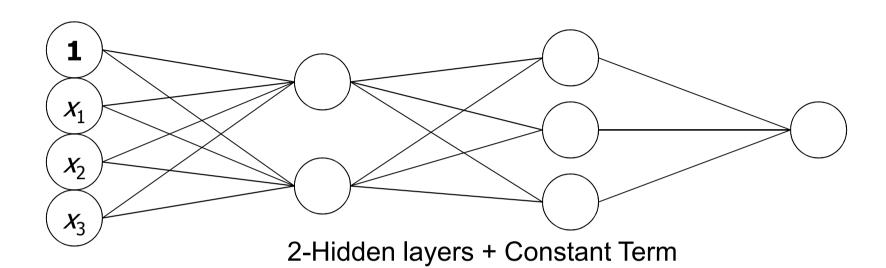
A 1-HIDDEN LAYER NET



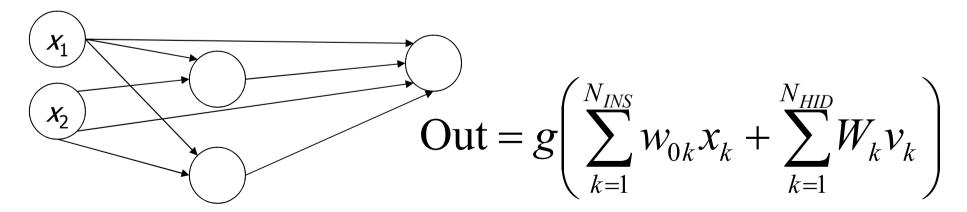
 $N_{HIDDEN} = 3$



OTHER NEURAL NETS



"JUMP" CONNECTIONS



Neural Networks: Slide 53

Backpropagation

$$\operatorname{Out}(\mathbf{x}) = g\left(\sum_{j} W_{j}g\left(\sum_{k} W_{jk}x_{k}\right)\right)$$

Find a set of weights $\{W_j\}, \{w_{jk}\}$

to minimize

$$\sum_{i} (y_i - \operatorname{Out}(\mathbf{x}_i))^2$$

by gradient descent.



Backpropagation Convergence

Convergence to a global minimum is <u>not</u> guaranteed.

•In practice, this is not a problem, apparently.

Tweaking to find the right number of hidden units, or a useful learning rate η , is more hassle, apparently.

IMPLEMENTING BACKPROP: \supseteq Differentiate Monster sum-square residual Write down the Gradient Descent Rule It turns out to be easier & computationally efficient to use lots of local variables with names like $h_j o_k v_j net_i$ etc...

Choosing the learning rate

- This is a subtle art.
- Too small: can take days instead of minutes to converge
- Too large: diverges (MSE gets larger and larger while the weights increase and usually oscillate)
- Sometimes the "just right" value is hard to find.

Learning-rate problems

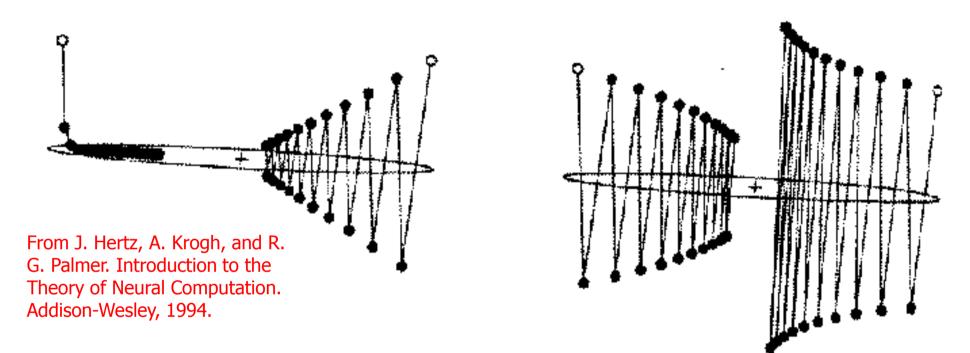


FIGURE 5.10 Gradient descent on a simple quadratic surface (the left and right parts are copies of the same surface). Four trajectories are shown, each for 20 steps from the open circle. The minimum is at the + and the ellipse shows a constant error contour. The only significant difference between the trajectories is the value of η , which was 0.02, 0.0476, 0.049, and 0.0505 from left to right.

Improving Simple Gradient Descent

Momentum

Don't just change weights according to the current datapoint. Re-use changes from earlier iterations.

Let $\Delta \mathbf{w}(t)$ = weight changes at time *t*.

Let $-\eta \frac{\partial E}{\partial W}$ be the change we would make with regular gradient descent.

Instead we use

$$\Delta \mathbf{w}(t+1) = -\eta \frac{\partial \mathbf{E}}{\partial \mathbf{w}} + \alpha \Delta \mathbf{w}(t)$$
$$\mathbf{w}(t+1) = \mathbf{w}(t) + \Delta \mathbf{w}(t)$$
Momentum damps oscillations.
A hack? Well, maybe.

Momentum illustration

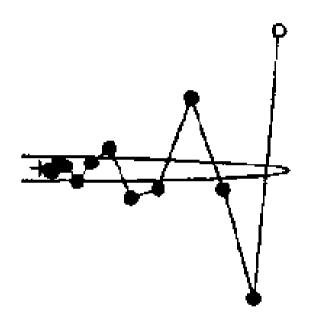


FIGURE 6.3 Gradient descent on the simple quadratic surface of Fig. 5.10. Both trajectories are for 12 steps with $\eta = 0.0476$, the best value in the absence of momentum. On the left there is no momentum ($\alpha = 0$), while $\alpha = 0.5$ on the right. Improving Simple Gradient Descent Newton's method

$$E(\mathbf{w} + \mathbf{h}) = E(\mathbf{w}) + \mathbf{h}^T \frac{\partial E}{\partial \mathbf{w}} + \frac{1}{2} \mathbf{h}^T \frac{\partial^2 E}{\partial \mathbf{w}^2} \mathbf{h} + O(|\mathbf{h}|^3)$$

If we neglect the $O(h^3)$ terms, this is a *quadratic form*

Quadratic form fun facts:

If
$$y = c + b^T x - 1/2 x^T A x$$

And if **A** is SPD

Then

 $\mathbf{x}^{opt} = \mathbf{A}^{-1}\mathbf{b}$ is the value of \mathbf{x} that maximizes \mathbf{y}

Improving Simple Gradient Descent Newton's method

$$E(\mathbf{w} + \mathbf{h}) = E(\mathbf{w}) + \mathbf{h}^T \frac{\partial E}{\partial \mathbf{w}} + \frac{1}{2} \mathbf{h}^T \frac{\partial^2 E}{\partial \mathbf{w}^2} \mathbf{h} + O(|\mathbf{h}|^3)$$

If we neglect the $O(h^3)$ terms, this is a *quadratic form*

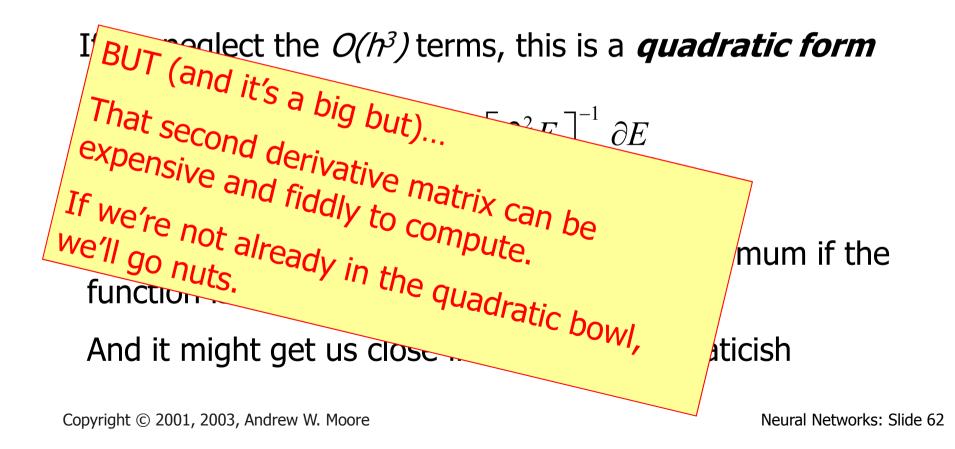
$$\mathbf{w} \leftarrow \mathbf{w} - \left[\frac{\partial^2 E}{\partial \mathbf{w}^2}\right]^{-1} \frac{\partial E}{\partial \mathbf{w}}$$

This should send us directly to the global minimum if the function is truly quadratic.

And it might get us close if it's locally quadraticish

Improving Simple Gradient Descent Newton's method

$$E(\mathbf{w} + \mathbf{h}) = E(\mathbf{w}) + \mathbf{h}^T \frac{\partial E}{\partial \mathbf{w}} + \frac{1}{2} \mathbf{h}^T \frac{\partial^2 E}{\partial \mathbf{w}^2} \mathbf{h} + O(|\mathbf{h}|^3)$$



Improving Simple Gradient Descent Conjugate Gradient

Another method which attempts to exploit the "local quadratic bowl" assumption

But does so while only needing to use ∂E

∂w

and not
$$\frac{\partial^2 E}{\partial \mathbf{w}^2}$$

It is also more stable than Newton's method if the local quadratic bowl assumption is violated.

It's complicated, outside our scope, but it often works well. More details in Numerical Recipes in C.

BEST GENERALIZATION

Intuitively, you want to use the smallest, simplest net that seems to fit the data.



- 1. Don't. Just use intuition
- 2. Bayesian Methods Get it Right
- 3. Statistical Analysis explains what's going on
- 4. Cross-validation Discussed in the next lecture

What You Should Know

- How to implement multivariate Leastsquares linear regression.
- Derivation of least squares as max. likelihood estimator of linear coefficients
- The general gradient descent rule

What You Should Know

- Perceptrons
 - \rightarrow Linear output, least squares
 - \rightarrow Sigmoid output, least squares
- Multilayer nets
 - \rightarrow The idea behind back prop
 - \rightarrow Awareness of better minimization methods
- Generalization. What it means.

APPLICATIONS

To Discuss:

- What can non-linear regression be useful for?
- What can neural nets (used as non-linear regressors) be useful for?
- What are the advantages of N. Nets for nonlinear regression?
- What are the disadvantages?

Other Uses of Neural Nets...

- Time series with recurrent nets
- Unsupervised learning (clustering principal components and non-linear versions thereof)
- Combinatorial optimization with Hopfield nets, Boltzmann Machines
- Evaluation function learning (in reinforcement learning)