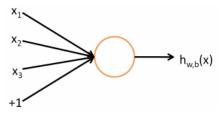
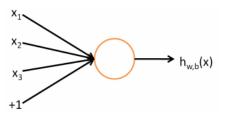


# Multilayer Networks

Natural Language Processing: Jordan Boyd-Graber University of Maryland

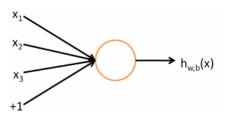




# Input

Vector  $x_1 \dots x_d$ 

inputs encoded as real numbers



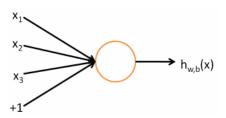
# Output

Input

Vector  $x_1 \dots x_d$ 

$$f\left(\sum_{i}W_{i}X_{i}+b\right)$$

multiply inputs by



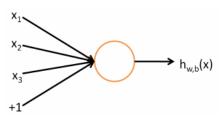
# Output

Input

Vector  $x_1 \dots x_d$ 

$$f\left(\sum_{i}W_{i}x_{i}+b\right)$$

add bias



# Input

Vector  $x_1 \dots x_d$ 

# Output

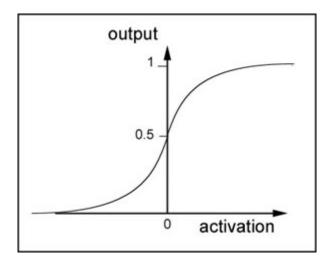
$$f\left(\sum_{i}W_{i}x_{i}+b\right)$$

# Activation

$$f(z) \equiv \frac{1}{1 + \exp(-z)}$$

pass through nonlinear sigmoid

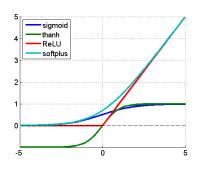
# Why is it called activation?



#### In the shallow end

- This is still logistic regression
- Engineering features *x* is difficult (and requires expertise)
- Can we learn how to represent inputs into final decision?

# Better name: non-linearity



Logistic / Sigmoid

$$f(x) = \frac{1}{1 + e^{-x}} \tag{1}$$

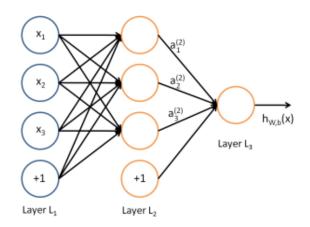
tanh

$$f(x) = \tanh(x) = \frac{2}{1 + e^{-2x}} - 1$$
 (2)

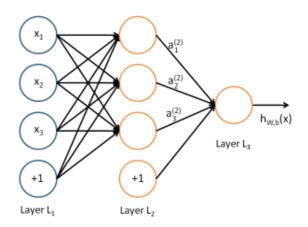
ReLU

$$f(x) = \begin{cases} 0 & \text{for } x < 0 \\ x & \text{for } x \ge 0 \end{cases}$$
 (3)

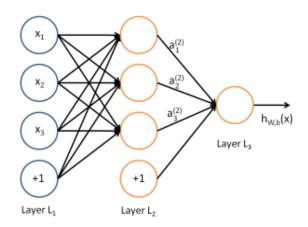
• SoftPlus:  $f(x) = \ln(1 + e^x)$ 



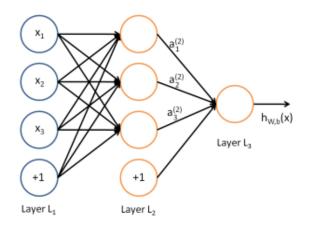
$$a_1^{(2)} = f(W_{11}^{(1)}x_1 + W_{12}^{(1)}x_2 + W_{13}^{(1)}x_3 + b_1^{(1)})$$



$$a_2^{(2)} = f\left(W_{21}^{(1)}x_1 + W_{22}^{(1)}x_2 + W_{23}^{(1)}x_3 + b_2^{(1)}\right)$$



$$a_3^{(2)} = f(W_{31}^{(1)}x_1 + W_{32}^{(1)}x_2 + W_{33}^{(1)}x_3 + b_3^{(1)})$$



$$h_{W,b}(x) = a_1^{(3)} = f\left(W_{11}^{(2)}a_1^{(2)} + W_{12}^{(2)}a_2^{(2)} + W_{13}^{(2)}a_3^{(2)} + b_1^{(2)}\right)$$

• For every example x, y of our supervised training set, we want the label y to match the prediction  $h_{W,b}(x)$ .

$$J(W,b;x,y) \equiv \frac{1}{2} ||h_{W,b}(x) - y||^2$$
 (4)

• For every example x, y of our supervised training set, we want the label y to match the prediction  $h_{W,h}(x)$ .

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 We want this value, summed over all of the examples to be as small as possible

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- We want this value, summed over all of the examples to be as small as possible
- We also want the weights not to be too large

$$\frac{\lambda}{2} \sum_{l=1}^{n_{l-1}} \sum_{j=1}^{s_{l}} \sum_{j=1}^{s_{l+1}} \left( W_{ji}^{l} \right)^{2} \tag{5}$$

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Sum over all layers

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Sum over all destinations

Putting it all together:

$$J(W,b) = \left[\frac{1}{m}\sum_{i=1}^{m}\frac{1}{2}||h_{W,b}(x^{(i)}) - y^{(i)}||^2\right] + \frac{\lambda}{2}\sum_{i=1}^{n_{i-1}}\sum_{j=1}^{s_{i}}\sum_{j=1}^{s_{j+1}}\left(W_{ji}^l\right)^2$$
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- Initialize W and b to small random value near zero

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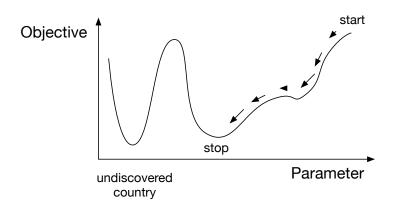
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(6)

- Our goal is to minimize J(W, b) as a function of W and b
- Initialize W and b to small random value near zero
- Adjust parameters to optimize J

#### **Gradient Descent**

# Goal

Optimize J with respect to variables W and b



For convenience, write the input to sigmoid

$$z_i^{(l)} = \sum_{j=1}^n W_{ij}^{(l-1)} x_j + b_i^{(l-1)}$$
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- The gradient is a function of a node's error  $\delta_i^{(l)}$
- For output nodes, the error is obvious:

$$\delta_i^{(n_i)} = \frac{\partial}{\partial z_i^{(n_i)}} ||y - h_{w,b}(x)||^2 = -\left(y_i - a_i^{(n_i)}\right) \cdot f'\left(z_i^{(n_i)}\right) \frac{2}{2}$$
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$$\delta_{i}^{(n_{l})} = \frac{\partial}{\partial z_{i}^{(n_{l})}} ||y - h_{w,b}(x)||^{2} = -\left(y_{i} - a_{i}^{(n_{l})}\right) \cdot f'\left(z_{i}^{(n_{l})}\right) \frac{2}{2}$$
(8)

Other nodes must "backpropagate" downstream error based on connection strength

$$\delta_{i}^{(l)} = \left(\sum_{j=1}^{s_{l+1}} W_{ji}^{(l+1)} \delta_{j}^{(l+1)}\right) f'(z_{i}^{(l)})$$
(9)

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(9)

(chain rule)

#### Partial Derivatives

For weights, the partial derivatives are

$$\frac{\partial}{\partial W_{ij}^{(l)}}J(W,b;x,y)=a_j^{(l)}\delta_j^{(l+1)}$$
(10)

For the bias terms, the partial derivatives are

$$\frac{\partial}{\partial b_i^{(l)}} J(W, b; x, y) = \delta_i^{(l+1)}$$
(11)

But this is just for a single example ...

# **Full Gradient Descent Algorithm**

- 1. Initialize  $U^{(l)}$  and  $V^{(l)}$  as zero
- 2. For each example  $i = 1 \dots m$ 
  - 2.1 Use backpropagation to compute  $\nabla_W J$  and  $\nabla_h J$
  - 2.2 Update weight shifts  $U^{(l)} = U^{(l)} + \nabla_{W^{(l)}} J(W, b; x, y)$
  - **2.3** Update bias shifts  $V^{(l)} = V^{(l)} + \nabla_{b^{(l)}} J(W, b; x, y)$
- Update the parameters

$$W^{(I)} = W^{(I)} - \alpha \left[ \left( \frac{1}{m} U^{(I)} \right) \right]$$
 (12)

$$b^{(l)} = b^{(l)} - \alpha \left[ \frac{1}{m} V^{(l)} \right]$$
 (13)

Repeat until weights stop changing

#### But it is not perfect

- Compare against baselines: randomized features, nearest-neighbors, linear models
- Optimization is hard (alchemy)
- Models are often not interpretable
- Requires specialized hardware and tons of data to scale