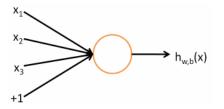
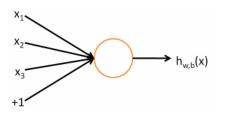
Multilayer Networks

Jordan Boyd-Graber

University of Maryland

Slides adapted from Andrew Ng

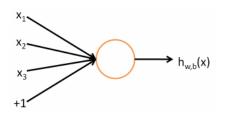




Input

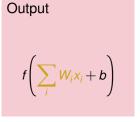
Vector $x_1 \dots x_d$

inputs encoded as real numbers

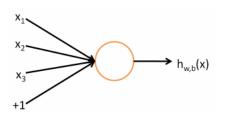


Input

Vector $x_1 \dots x_d$



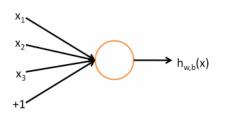
multiply inputs by weights



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

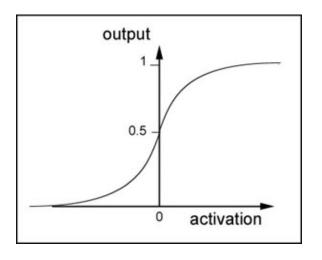
$$f\left(\sum_{i}W_{i}x_{i}+b\right)$$

Activation

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

pass through nonlinear sigmoid

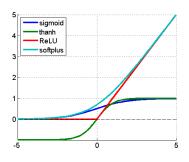
Why is it called activation?



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

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

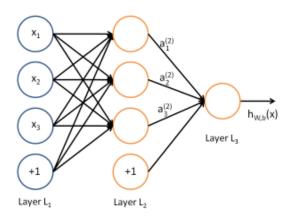
tanh

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

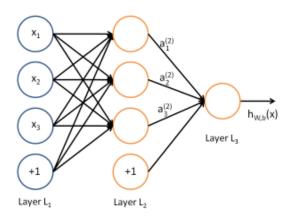
ReLU

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

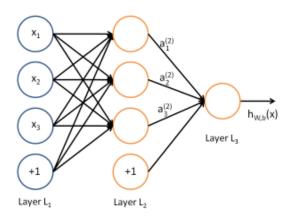
• SoftPlus: $f(x) = \ln(1 + e^x)$



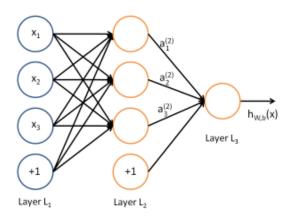
$$a_1^{(2)} = f\left(W_{11}^{(1)}x_1 + W_{12}^{(1)}x_2 + W_{13}^{(1)}x_3 + b_1^{(1)}\right)$$



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



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



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

• 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)

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- · We also want the weights not to be too large

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

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Sum over all layers

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Sum over all destinations

Putting it all together:

$$J(W,b) = \left[\frac{1}{m}\sum_{i=1}^{m}\frac{1}{2}||h_{W,b}(x^{(i)}) - y^{(i)}||^2\right] + \frac{\lambda}{2}\sum_{l=1}^{n_l-1}\sum_{i=1}^{s_l}\sum_{j=1}^{s_{l+1}}\left(W_{ji}^l\right)^2$$
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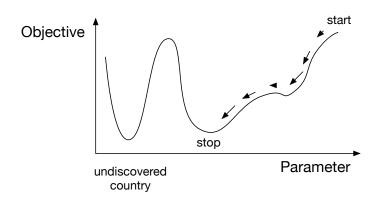
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- Our goal is to minimize J(W, b) as a function of W and b
- Initialize W and b to small random value near zero
- Adjust parameters to optimize J

Gradient Descent

Goal

Optimize J with respect to variables W and b



• For convenience, write the input to sigmoid

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

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

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

$$\delta_{i}^{(l)} = \left(\sum_{j=1}^{s_{t+1}} W_{ji}^{(l+1)} \delta_{j}^{(l+1)}\right) f'(z_{i}^{(l)})$$
(9)

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(chain rule) (9)

Partial Derivatives

• For weights, the partial derivatives are

$$\frac{\partial}{\partial W_{ii}^{(l)}} J(W, b; x, y) = a_j^{(l)} \delta_i^{(l+1)}$$
(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 $\nabla_b J$
 - 2.2 Update weight shifts $U^{(l)} = U^{(l)} + \nabla_{W^{(l)}} J(W, b; x, y)$
 - 2.3 Update bias shifts $V^{(l)} = V^{(l)} + \nabla_{b^{(l)}} J(W, b; x, y)$
- 3. Update the parameters

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

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

4. 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