CS 6120/CS4120: Natural Language Processing

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Outline

- Maximum Entropy
- Feedforward Neural Networks
- Recurrent Neural Networks

Introduction

- So far we've looked at "generative models"
 - Language models, Naive Bayes
- But there is now much use of conditional or discriminative probabilistic models in NLP, Speech, IR (and ML generally)
- Because:
 - They give high accuracy performance
 - They make it easy to incorporate lots of linguistically important features
 - They allow automatic building of language independent, retargetable NLP modules

Joint vs. Conditional Models

- We have some data {(*d*, *c*)} of paired observations *d* and hidden classes *c*.
- Joint (generative) models place probabilities over both observed data and the hidden stuff (generate the observed data from hidden stuff):

P(c,d)

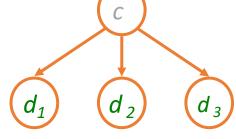
- All the classic statistic NLP models:
 - *n*-gram models, Naive Bayes classifiers, hidden Markov models, probabilistic context-free grammars, IBM machine translation alignment models

Joint vs. Conditional Models

- Discriminative (conditional) models take the data as given, and put a probability over hidden structure given the data: P(c|d)
 - Logistic regression, conditional loglinear or maximum entropy models, conditional random fields
 - Also, SVMs, (averaged) perceptron, etc. are discriminative classifiers (but not directly probabilistic)

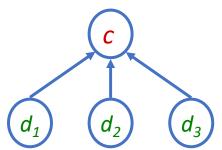
Bayes Network/Graphical Models

- Bayes network diagrams draw circles for random variables, and lines for direct dependencies
- Some variables are observed; some are hidden
- Each node is a little classifier (conditional probability table) based on incoming arcs



Naive Bayes

Generative



Logistic Regression (aka Maximum Entropy)

Discriminative

Conditional vs. Joint Likelihood

- A *joint* model gives probabilities P(*d*,*c*) and tries to maximize this joint likelihood.
 - It turns out to be trivial to choose weights: just relative frequencies.
- A *conditional* model gives probabilities P(c|d). It takes the data as given and models only the conditional probability of the class.
 - We seek to maximize conditional likelihood.
 - Harder to do.
 - More closely related to classification error. (Easy to tune!)

Maximum Entropy (MaxEnt)

• Or logistic regression

Features

- In these slides and most maxent work: *features (or feature functions) f* are elementary pieces of evidence that link aspects of what we observe *d* with a category *c* that we want to predict
- A feature is a function with a **bounded** real value: $f: C \times D \rightarrow \mathbb{R}$

Example Task: Named Entity Type

LOCATION

LOCATION in Arcadia in Québec

DRUG PERSON taking Zantac saw Sue

Example features

- $f_1(c, d) \equiv [c = \text{LOCATION} \land w_{-1} = \text{``in''} \land \text{isCapitalized}(w)]$
- $f_2(c, d) \equiv [c = \text{LOCATION} \land \text{hasAccentedLatinChar}(w)]$
- $f_3(c, d) \equiv [c = DRUG \land ends(w, "c")]$



- Models will assign to each feature a *weight:*
 - A positive weight votes that this configuration is likely correct
 - A negative weight votes that this configuration is likely incorrect

Features

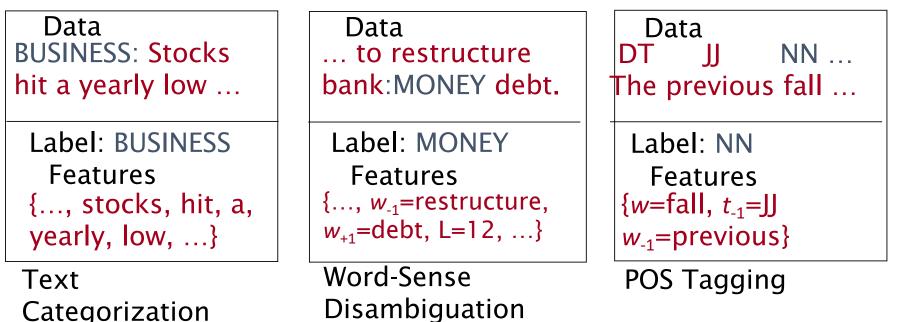
- In NLP uses, usually a feature specifies
 - 1. an indicator function a yes/no boolean matching function of properties of the input and a particular class

$$f_i(c, d) \equiv [\Phi(d) \land c = c_j]$$
 [Value is 0 or 1]

• Each feature picks out a data subset and suggests a label for it

Feature-Based Models

• The decision about a data point is based only on the features active at that point.



Categorization

Other Maxent Classifier Examples

- You can use a maxent classifier whenever you want to assign data points to one of a number of classes:
 - Sentence boundary detection (Mikheev 2000)
 - Is a period end of sentence or abbreviation?
 - Sentiment analysis (Pang and Lee 2002)
 - Word unigrams, bigrams, POS counts, ...
 - Prepositional phrase attachment (Ratnaparkhi 1998)
 - Attach to verb or noun? Features of head noun, preposition, etc.
 - Parsing decisions in general (Ratnaparkhi 1997; Johnson et al. 1999, etc.)

Feature-Based Linear Classifiers

- Linear classifiers at classification time:
 - Linear function from feature sets {*f_i*} to classes {*c*}.
 - Assign a weight λ_i to each feature f_i .
 - We consider each class for sample *d*
 - For a pair (*c*,*d*), features vote with their weights:
 - vote(c) = $\Sigma \lambda_i f_i(c, d)$

| PERSON | LOCATION | DRUG |
|-----------|-----------|-----------|
| in Québec | in Québec | in Québec |

• Choose the class *c* which maximizes $\sum \lambda_i f_i(c,d)$

Feature-Based Linear Classifiers

- $f_1(c, d) \equiv [c = \text{LOCATION} \land w_1 = \text{``in''} \land \text{isCapitalized}(w)] \rightarrow \text{weight 1.8}$
- $f_2(c, d) \equiv [c = \text{LOCATION} \land \text{hasAccentedLatinChar}(w)] \rightarrow \text{weight } -0.6$
- $f_3(c, d) \equiv [c = DRUG \land ends(w, "c")] \rightarrow weight 0.3$

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- Maximum Entropy:
 - Make a probabilistic model from the linear combination $\Sigma \lambda_i f_i(c,d)$

$$P(c \mid d, \lambda) = \frac{\exp \sum_{i} \lambda_{i} f_{i}(c, d)}{\sum_{c'} \exp \sum_{i} \lambda_{i} f_{i}(c', d)} \leftarrow \frac{\text{Makes votes positive}}{\text{Normalizes votes}}$$

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- $P(LOCATION|in Québec) = e^{1.8}e^{-0.6}/(e^{1.8}e^{-0.6} + e^{0.3} + e^{0}) = 0.586$
- $P(DRUG|in Québec) = e^{0.3} / (e^{1.8}e^{-0.6} + e^{0.3} + e^{0}) = 0.238$
- $P(PERSON|in Québec) = e^0 / (e^{1.8}e^{-0.6} + e^{0.3} + e^0) = 0.176$
- The weights are the parameters of the probability model, combined via a "soft max" function

Feature-Based Linear Classifiers

• Exponential models:

- Given this model form, we will choose parameters $\{\lambda_i\}$ that *maximize the conditional likelihood* of the data according to this model.
- We construct not only classifications, but probability distributions over classifications.
 - There are other (good!) ways of discriminating classes SVMs, boosting, even perceptrons – but these methods are not as trivial to interpret as distributions over classes.

Outline

- Maximum Entropy
- Feedforward Neural Networks
- Recurrent Neural Networks

Neural Network Learning

- Learning approach based on modeling adaptation in biological neural systems.
- Perceptron: Initial algorithm for learning simple neural networks (single layer) developed in the 1950's.
- Backpropagation: More complex algorithm for learning multi-layer neural networks developed in the 1980's.

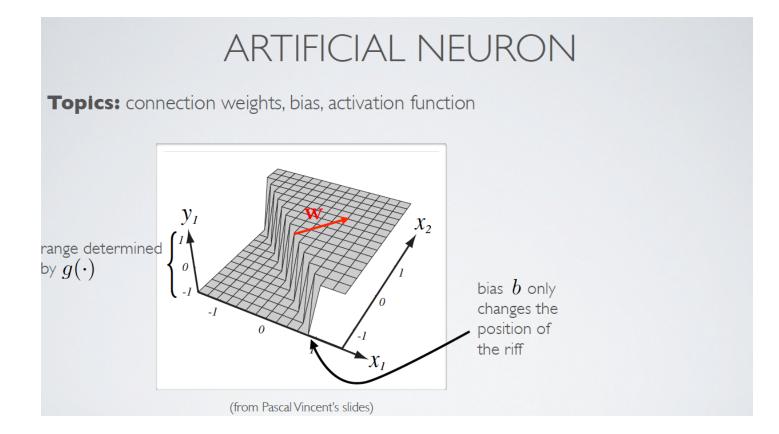
ARTIFICIAL NEURON

Topics: connection weights, bias, activation function

• Neuron pre-activation (or input activation):

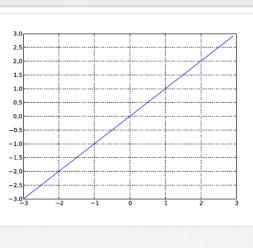
$$a(\mathbf{x}) = b + \sum_{i} w_i x_i = b + \mathbf{w}^\top \mathbf{x}$$

- Neuron (output) activation $h(\mathbf{x}) = g(a(\mathbf{x})) = g(b + \sum_{i} w_{i}x_{i})$ • w are the connection weights
- $\cdot b$ is the neuron bias
- $g(\cdot)$ is called the activation function



Topics: linear activation function

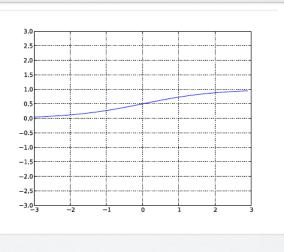
- Performs no input squashing
- Not very interesting...



g(a) = a

Topics: sigmoid activation function

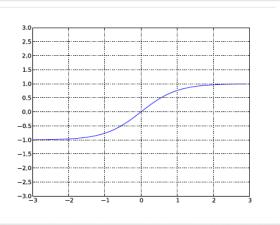
- Squashes the neuron's pre-activation between 0 and 1
- Always positive
- Bounded
- Strictly increasing



$$g(a) = \operatorname{sigm}(a) = \frac{1}{1 + \exp(-a)}$$

Topics: hyperbolic tangent ('tanh'') activation function

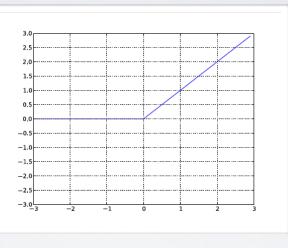
- Squashes the neuron's pre-activation between
 I and I
- Can be positive or negative
- Bounded
- Strictly increasing



$$g(a) = \tanh(a) = \frac{\exp(a) - \exp(-a)}{\exp(a) + \exp(-a)} = \frac{\exp(2a) - 1}{\exp(2a) + 1}$$

Topics: rectified linear activation function

- Bounded below by 0 (always non-negative)
- Not upper bounded
- Strictly increasing
- Tends to give neurons with sparse activities

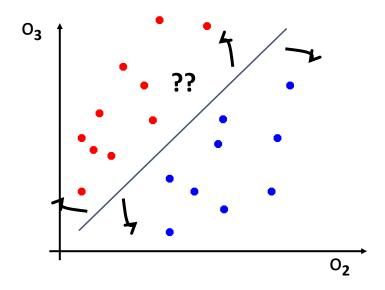


 $g(a) = \operatorname{reclin}(a) = \max(0, a)$

```
class Neuron(object):
    # ...
    def forward(inputs):
        """ assume inputs and weights are 1-D numpy arrays and bias is a number """
        cell_body_sum = np.sum(inputs * self.weights) + self.bias
        firing_rate = 1.0 / (1.0 + math.exp(-cell_body_sum)) # sigmoid activation function
        return firing_rate
```

Linear Separator

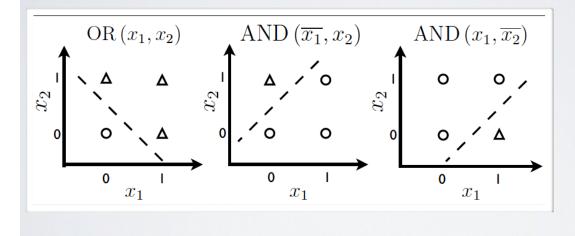
• Since one-layer neuron (aka perceptron) uses linear threshold function, it is searching for a linear separator that discriminates the classes.



ARTIFICIAL NEURON

Topics: capacity of single neuron

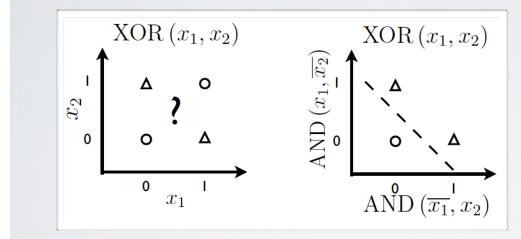
Can solve linearly separable problems



ARTIFICIAL NEURON

Topics: capacity of single neuron

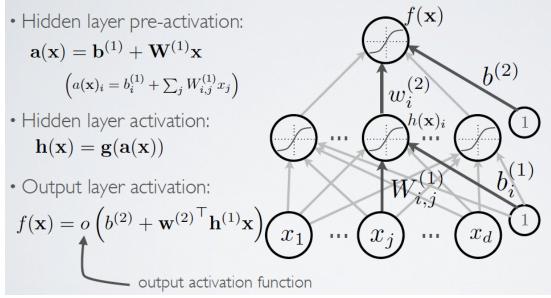
Can't solve non linearly separable problems...



• ... unless the input is transformed in a better representation

NEURAL NETWORK

Topics: single hidden layer neural network



NEURAL NETWORK

Topics: softmax activation function

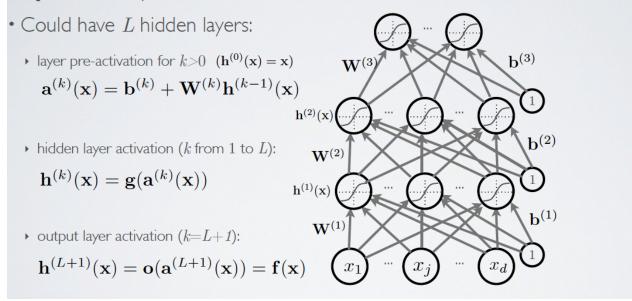
- For multi-class classification:
 - we need multiple outputs (I output per class)
 - ullet we would like to estimate the conditional probability $p(y=c|\mathbf{x})$
- We use the softmax activation function at the output:

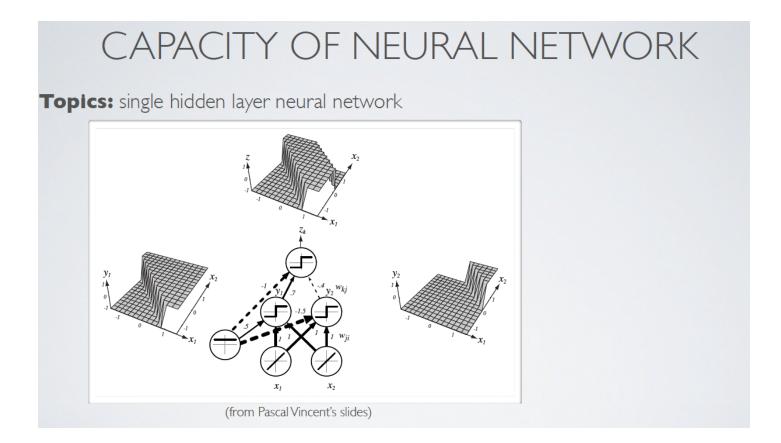
$$\mathbf{o}(\mathbf{a}) = \operatorname{softmax}(\mathbf{a}) = \left[\frac{\exp(a_1)}{\sum_c \exp(a_c)} \dots \frac{\exp(a_C)}{\sum_c \exp(a_c)}\right]^{\mathsf{T}}$$

- strictly positive
- sums to one
- Predicted class is the one with highest estimated probability

NEURAL NETWORK

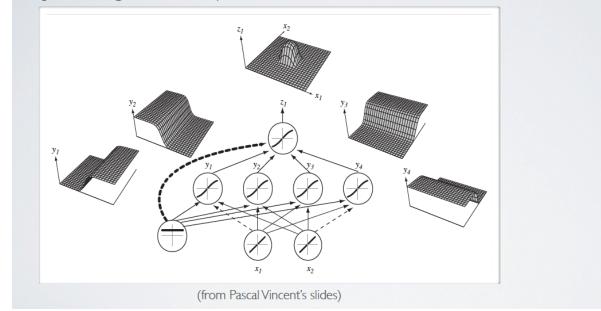
Topics: multilayer neural network





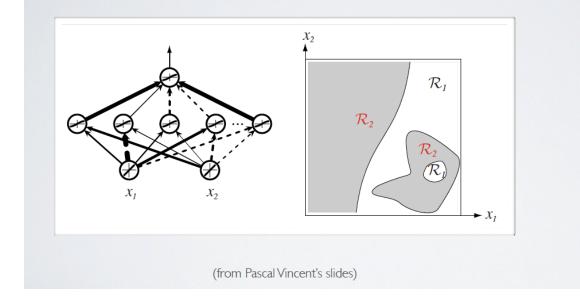
CAPACITY OF NEURAL NETWORK

Topics: single hidden layer neural network



CAPACITY OF NEURAL NETWORK

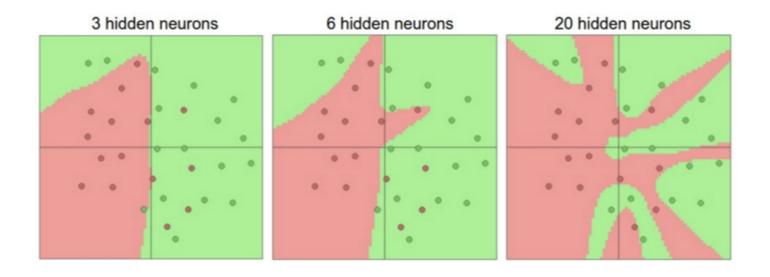
Topics: single hidden layer neural network



CAPACITY OF NEURAL NETWORK

Topics: universal approximation

- Universal approximation theorem (Homik, 1991):
 - "a single hidden layer neural network with a linear output unit can approximate any continuous function arbitrarily well, given enough hidden units"
- The result applies for sigmoid, tanh and many other hidden layer activation functions
- This is a good result, but it doesn't mean there is a learning algorithm that can find the necessary parameter values!



How to train a neural network?

Topics: multilayer neural network

• Could have *L* hidden layers:

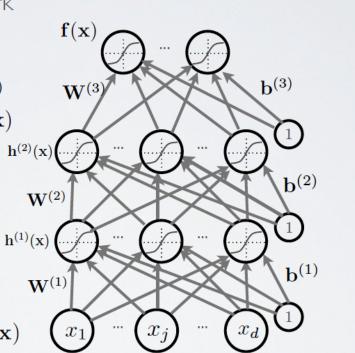
• layer input activation for k>0 $(\mathbf{h}^{(0)}(\mathbf{x}) = \mathbf{x})$ $\mathbf{a}^{(k)}(\mathbf{x}) = \mathbf{b}^{(k)} + \mathbf{W}^{(k)}\mathbf{h}^{(k-1)}(\mathbf{x})$

• hidden layer activation (k from 1 to L):

$$\mathbf{h}^{(k)}(\mathbf{x}) = \mathbf{g}(\mathbf{a}^{(k)}(\mathbf{x}))$$

• output layer activation (k=L+1):

$$\mathbf{h}^{(L+1)}(\mathbf{x}) = \mathbf{o}(\mathbf{a}^{(L+1)}(\mathbf{x})) = \mathbf{f}(\mathbf{x})$$



Empirical Risk Minimization

Topics: empirical risk minimization, regularization

- Empirical risk minimization
 - framework to design learning algorithms

$$\underset{\boldsymbol{\theta}}{\operatorname{arg\,min}} \frac{1}{T} \sum_{t} l(f(\mathbf{x}^{(t)}; \boldsymbol{\theta}), y^{(t)}) + \lambda \Omega(\boldsymbol{\theta})$$

- + $l(f(\mathbf{x}^{(t)}; \boldsymbol{ heta}), y^{(t)})$ is a loss function
- + $\Omega(oldsymbol{ heta})$ is a regularizer (penalizes certain values of $oldsymbol{ heta}$)
- Learning is cast as optimization
 - ideally, we'd optimize classification error, but it's not smooth
 - loss function is a surrogate for what we truly should optimize (e.g. upper bound)

The rest of the slides are for reference only (not covered by lecture or exam)

The Learning Algorithm Topics: stochastic gradient descent (SGD) • Algorithm that performs updates after each example • initialize θ ($\theta \equiv \{W^{(1)}, b^{(1)}, \dots, W^{(L+1)}, b^{(L+1)}\}$) • for N iterations • for each training example $(\mathbf{x}^{(t)}, y^{(t)})$ $\prec \Delta = -\nabla_{\theta} l(f(\mathbf{x}^{(t)}; \theta), y^{(t)}) - \lambda \nabla_{\theta} \Omega(\theta)$ $\prec \theta \leftarrow \theta + \alpha \Delta$ • To apply this algorithm to neural network training, we need • the loss function $l(\mathbf{f}(\mathbf{x}^{(t)}; \theta), y^{(t)})$

- a procedure to compute the parameter gradients $abla_{m{ heta}} l(\mathbf{f}(\mathbf{x}^{(t)};m{ heta}),y^{(t)})$
- the regularizer $\Omega({m heta})$ (and the gradient $abla_{{m heta}}\Omega({m heta})$)
- initialization method

LOSS FUNCTION

Topics: loss function for classification

- Neural network estimates $f(\mathbf{x})_c = p(y = c | \mathbf{x})$
 - ightarrow we could maximize the probabilities of $y^{(t)}$ given $\mathbf{x}^{(t)}$ in the training set
- To frame as minimization, we minimize the negative log-likelihood natural log (In)

$$l(\mathbf{f}(\mathbf{x}), y) = -\sum_{c} 1_{(y=c)} \log f(\mathbf{x})_{c} = -\log f(\mathbf{x})_{y}$$

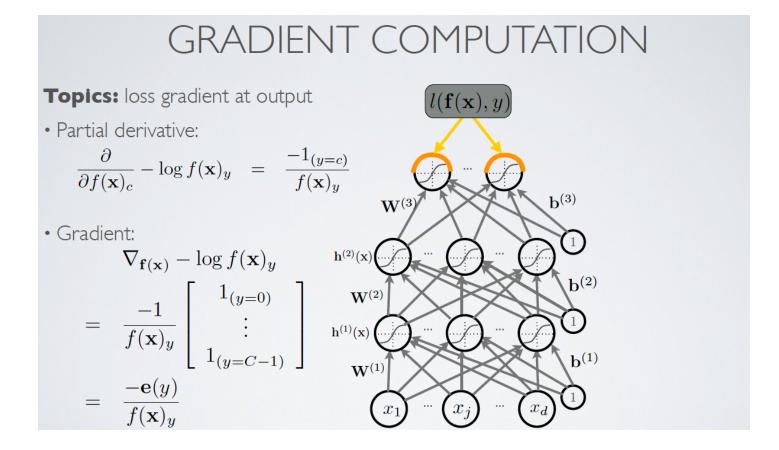
- we take the log to simplify for numerical stability and math simplicity
- sometimes referred to as cross-entropy

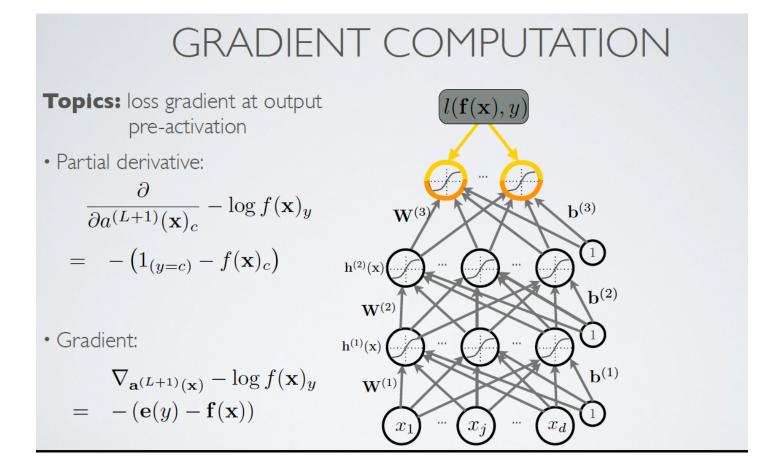
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$$rac{\partial}{\partial a^{(L+1)}(\mathbf{x})_c} - \log f(\mathbf{x})_y$$

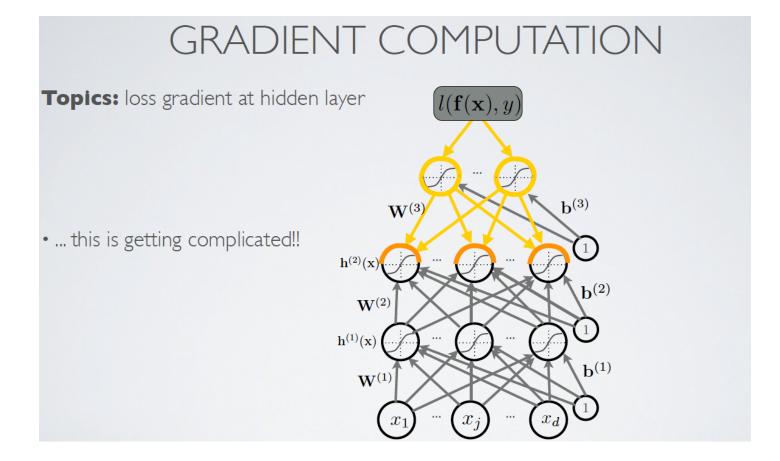
$$= \frac{\frac{\partial}{\partial a^{(L+1)}(\mathbf{x})_c} - \log f(\mathbf{x})_y}{\frac{\partial}{\partial a^{(L+1)}(\mathbf{x})_c}} f(\mathbf{x})_y$$

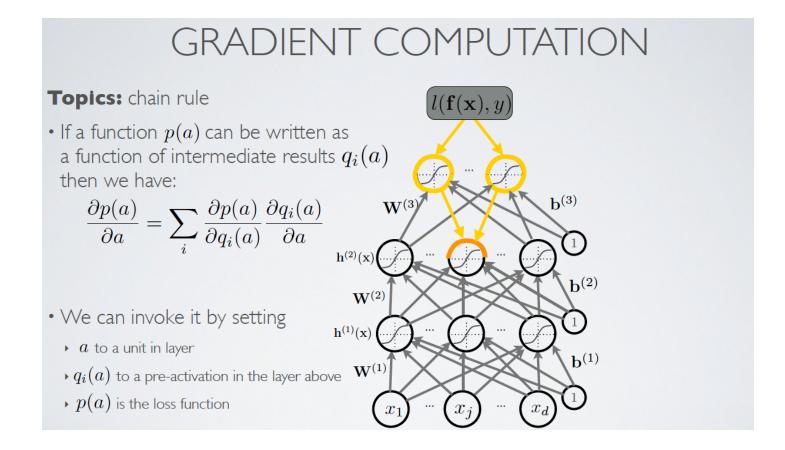
$$\begin{aligned} & \frac{\partial}{\partial a^{(L+1)}(\mathbf{x})_c} - \log f(\mathbf{x})_y \\ &= \frac{-1}{f(\mathbf{x})_y} \frac{\partial}{\partial a^{(L+1)}(\mathbf{x})_c} f(\mathbf{x})_y \\ &= \frac{-1}{f(\mathbf{x})_y} \frac{\partial}{\partial a^{(L+1)}(\mathbf{x})_c} \operatorname{softmax}(\mathbf{a}^{(L+1)}(\mathbf{x}))_y \end{aligned}$$

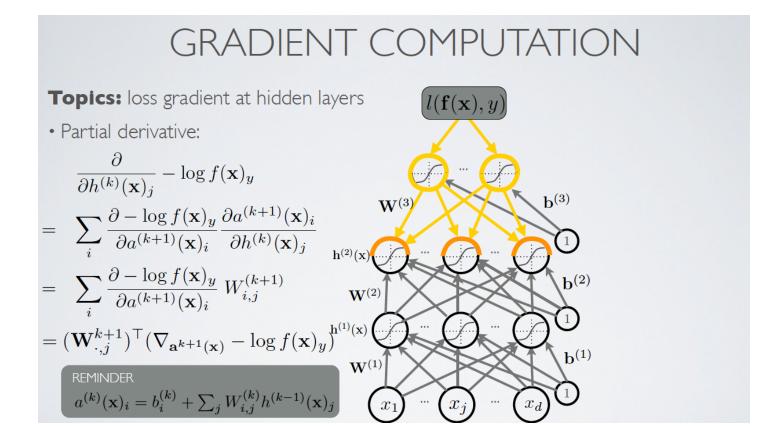
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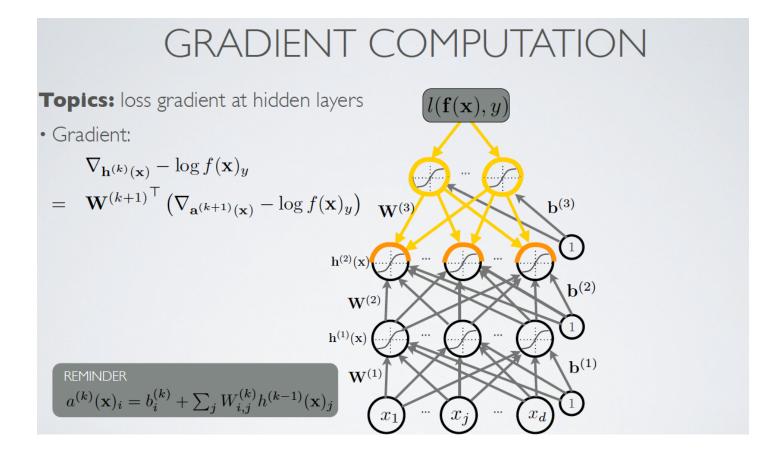
$$\begin{split} &\frac{\partial}{\partial a^{(L+1)}(\mathbf{x})_{c}} - \log f(\mathbf{x})_{y} \\ &= \frac{-1}{f(\mathbf{x})_{y}} \frac{\partial}{\partial a^{(L+1)}(\mathbf{x})_{c}} f(\mathbf{x})_{y} \\ &= \frac{-1}{f(\mathbf{x})_{y}} \frac{\partial}{\partial a^{(L+1)}(\mathbf{x})_{c}} \operatorname{softmax}(\mathbf{a}^{(L+1)}(\mathbf{x}))_{y} \end{split} \\ \begin{split} & \boxed{\frac{\partial \frac{g(x)}{h(x)}}{\partial x} = \frac{\partial g(x)}{\partial x} \frac{1}{h(x)} - \frac{g(x)}{h(x)^{2}} \frac{\partial h(x)}{\partial x}}{\partial x} \\ &= \frac{-1}{f(\mathbf{x})_{y}} \frac{\partial}{\partial a^{(L+1)}(\mathbf{x})_{c}} \operatorname{softmax}(\mathbf{a}^{(L+1)}(\mathbf{x}))_{y} \\ &= \frac{-1}{f(\mathbf{x})_{y}} \frac{\partial}{\partial a^{(L+1)}(\mathbf{x})_{c}} \frac{\exp(a^{(L+1)}(\mathbf{x})_{y})}{\sum_{c'} \exp(a^{(L+1)}(\mathbf{x})_{c'})} - \frac{\exp(a^{(L+1)}(\mathbf{x})_{y}) \left(\frac{\partial}{\partial a^{(L+1)}(\mathbf{x})_{c}} \sum_{c'} \exp(a^{(L+1)}(\mathbf{x})_{c'})\right)}{\left(\sum_{c'} \exp(a^{(L+1)}(\mathbf{x})_{c'}\right)^{2}} \end{split} \\ &= \frac{-1}{f(\mathbf{x})_{y}} \left(\frac{1_{(y=c)} \exp(a^{(L+1)}(\mathbf{x})_{y})}{\sum_{c'} \exp(a^{(L+1)}(\mathbf{x})_{c'})} - \frac{\exp(a^{(L+1)}(\mathbf{x})_{y})}{\sum_{c'} \exp(a^{(L+1)}(\mathbf{x})_{c'})} \frac{\exp(a^{(L+1)}(\mathbf{x})_{c'})}{\sum_{c'} \exp(a^{(L+1)}(\mathbf{x})_{c'})} \right) \\ &= \frac{-1}{f(\mathbf{x})_{y}} \left(\frac{1_{(y=c)} \exp(a^{(L+1)}(\mathbf{x})_{y}}{1_{(y=c)} \operatorname{softmax}(\mathbf{a}^{(L+1)}(\mathbf{x}))_{y}} - \operatorname{softmax}(\mathbf{a}^{(L+1)}(\mathbf{x}))_{y} \operatorname{softmax}(\mathbf{a}^{(L+1)}(\mathbf{x}))_{c} \right) \\ &= \frac{-1}{f(\mathbf{x})_{y}} \left(1_{(y=c)} \operatorname{softmax}(\mathbf{a}^{(L+1)}(\mathbf{x}))_{y} - \operatorname{softmax}(\mathbf{a}^{(L+1)}(\mathbf{x}))_{y} \operatorname{softmax}(\mathbf{a}^{(L+1)}(\mathbf{x}))_{c} \right) \end{aligned}$$

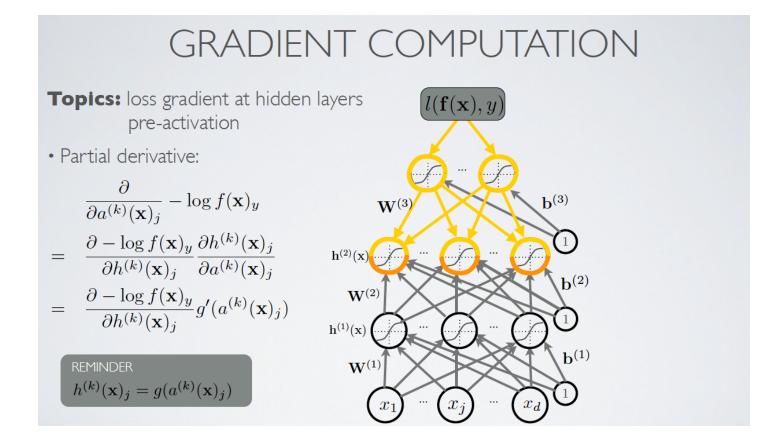
- Output layer gradient (o)
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- Activation function gradient (a)
- Parameter gradient (W, b)

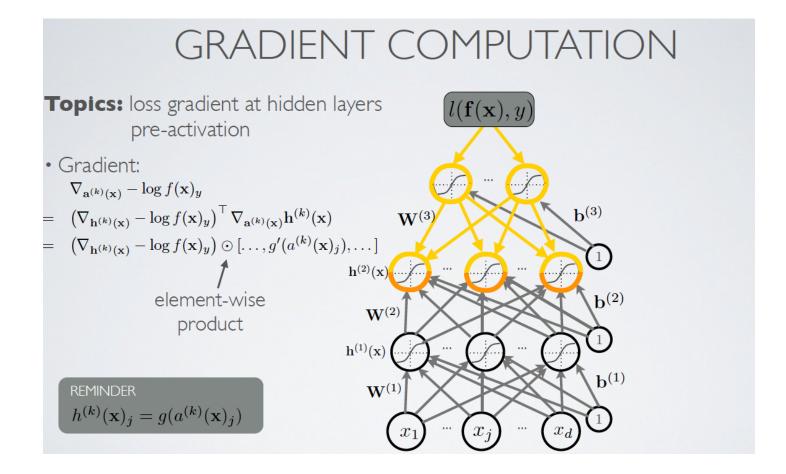








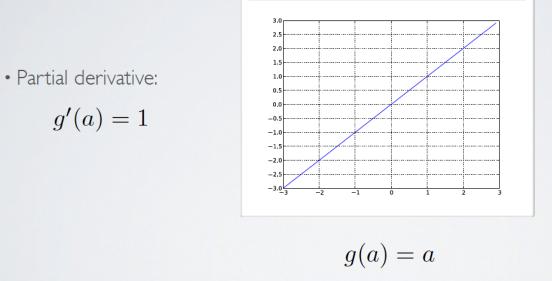




- Output layer gradient (o)
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ACTIVATION FUNCTION

Topics: linear activation function gradient

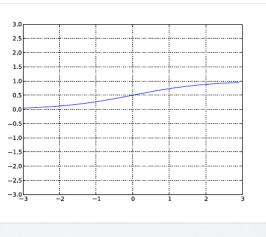


ACTIVATION FUNCTION

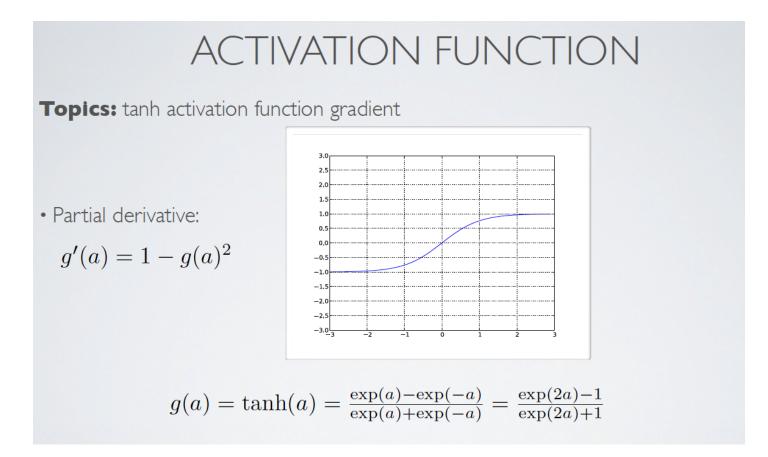
Topics: sigmoid activation function gradient

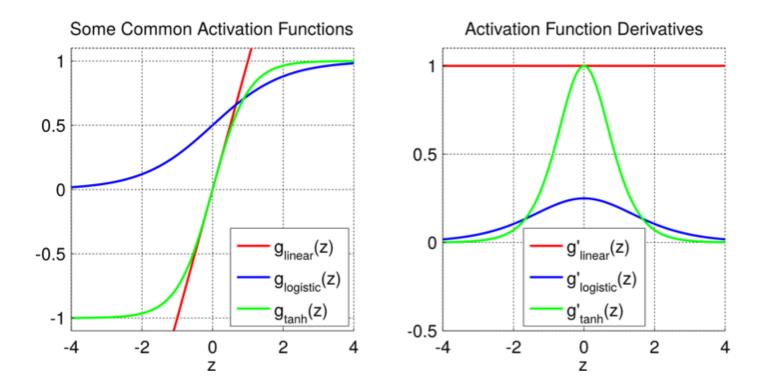
• Partial derivative:

g'(a) = g(a)(1 - g(a))

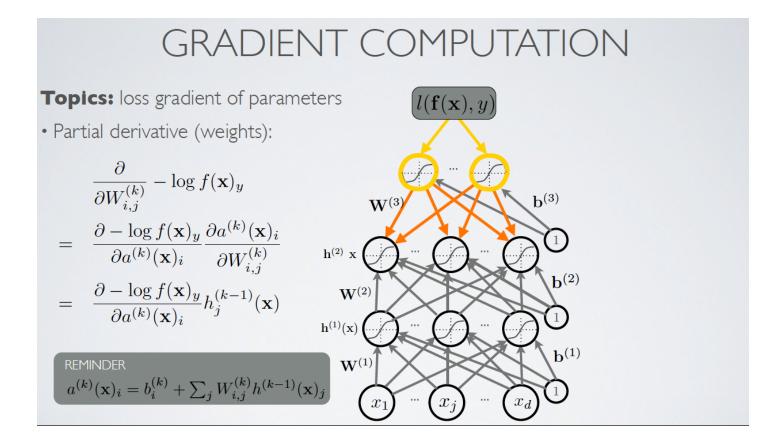


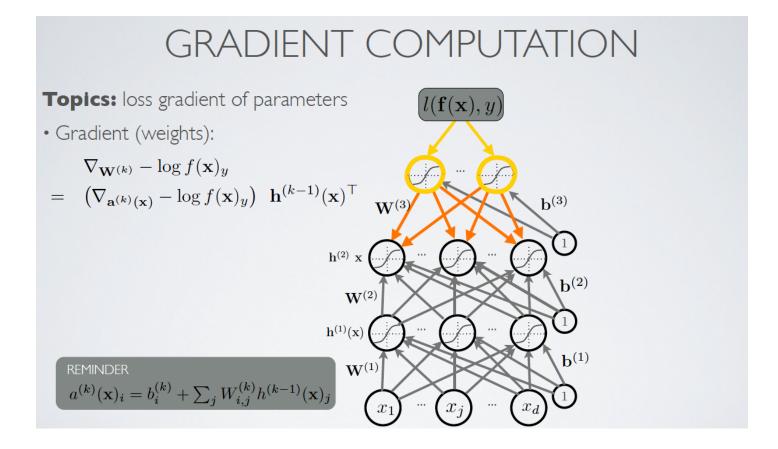
$$g(a) = \operatorname{sigm}(a) = \frac{1}{1 + \exp(-a)}$$

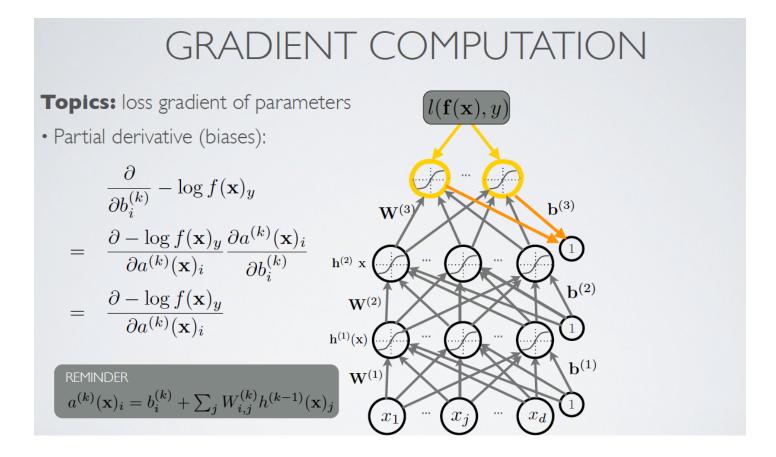


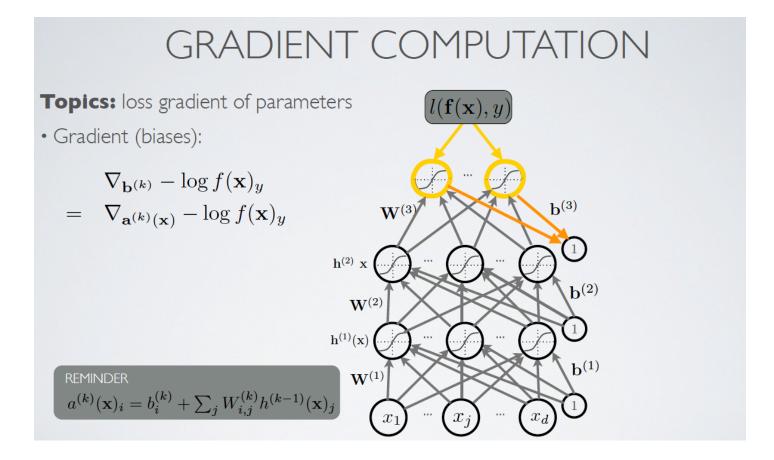


- Output layer gradient (o)
- Hidden layer gradient (h)
- Activation function gradient (a)
- Parameter gradient (W, b)

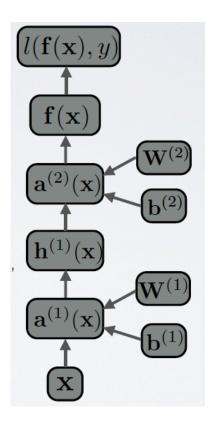








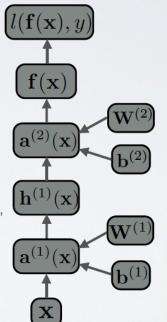
Backpropagation



FLOW GRAPH

Topics: flow graph

- Forward propagation can be represented as an acyclic flow graph
- It's a nice way of implementing forward propagation in a modular way
 - each box could be an object with an fprop method, that computes the value of the box given its children
 - calling the fprop method of each box in the right order yield forward propagation



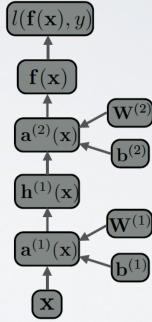
forward-pass of a 3-layer neural network:

f = lambda x: 1.0/(1.0 + np.exp(-x)) # activation function (use sigmoid)
x = np.random.randn(3, 1) # random input vector of three numbers (3x1)
h1 = f(np.dot(W1, x) + b1) # calculate first hidden layer activations (4x1)
h2 = f(np.dot(W2, h1) + b2) # calculate second hidden layer activations (4x1)
out = np.dot(W3, h2) + b3 # output neuron (1x1)

FLOW GRAPH

Topics: automatic differentiation

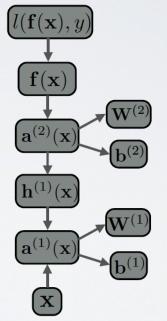
- Each object also has a bprop method
 - it computes the gradient of the loss with respect to each children
 - fprop depends on the fprop of a box's children, while bprop depends the bprop of a box's parents
- By calling bprop in the reverse order, we get backpropagation
 - only need to reach the parameters



FLOW GRAPH

Topics: automatic differentiation

- Each object also has a bprop method
 - it computes the gradient of the loss with respect to each children
 - fprop depends on the fprop of a box's children, while bprop depends the bprop of a box's parents
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BACKPROPAGATION

Topics: backpropagation algorithm

- This assumes a forward propagation has been made before
 - compute output gradient (before activation)

$$\nabla_{\mathbf{a}^{(L+1)}(\mathbf{x})} - \log f(\mathbf{x})_y \iff -(\mathbf{e}(y) - \mathbf{f}(\mathbf{x}))$$

- for k from L+1 to 1
 - compute gradients of hidden layer parameter

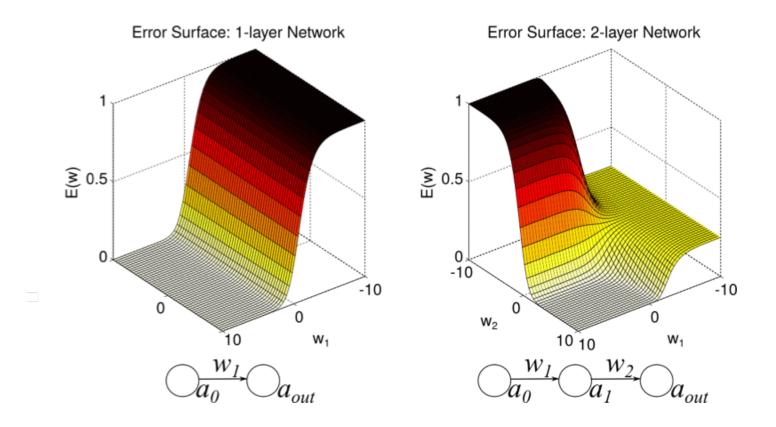
$$\nabla_{\mathbf{W}^{(k)}} - \log f(\mathbf{x})_y \iff \left(\nabla_{\mathbf{a}^{(k)}(\mathbf{x})} - \log f(\mathbf{x})_y\right) \quad \mathbf{h}^{(k-1)}(\mathbf{x})^\top$$
$$\nabla_{\mathbf{b}^{(k)}} - \log f(\mathbf{x})_y \iff \nabla_{\mathbf{a}^{(k)}(\mathbf{x})} - \log f(\mathbf{x})_y$$

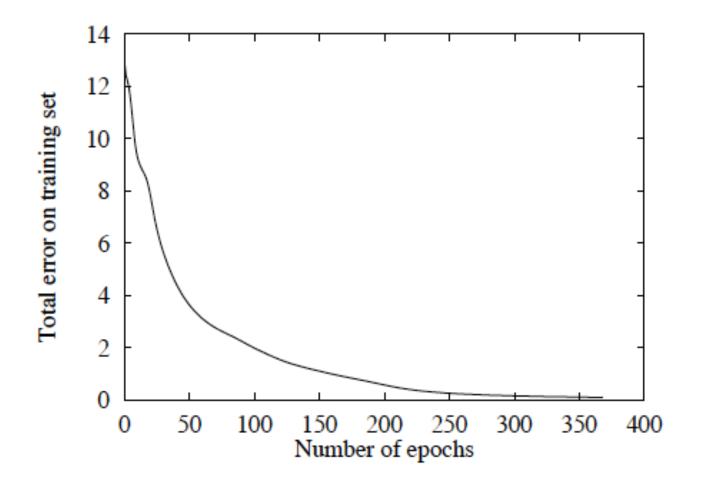
- compute gradient of hidden layer below

$$7_{\mathbf{h}^{(k-1)}(\mathbf{x})} - \log f(\mathbf{x})_y \iff \mathbf{W}^{(k)^{\top}} \left(\nabla_{\mathbf{a}^{(k)}(\mathbf{x})} - \log f(\mathbf{x})_y \right)$$

- compute gradient of hidden layer below (before activation)

$$\nabla_{\mathbf{a}^{(k-1)}(\mathbf{x})} - \log f(\mathbf{x})_y \iff \left(\nabla_{\mathbf{h}^{(k-1)}(\mathbf{x})} - \log f(\mathbf{x})_y\right) \odot \left[\dots, g'(a^{(k-1)}(\mathbf{x})_j), \dots\right]$$





[figure from Greg Mori's slides]

The Learning Algorithm Topics: stochastic gradient descent (SGD) • Algorithm that performs updates after each example • initialize θ ($\theta \equiv \{W^{(1)}, b^{(1)}, \dots, W^{(L+1)}, b^{(L+1)}\}$) • for N iterations • for each training example $(\mathbf{x}^{(t)}, y^{(t)})$ $< \Delta = -\nabla_{\theta} l(f(\mathbf{x}^{(t)}; \theta), y^{(t)}) - \lambda \nabla_{\theta} \Omega(\theta)$ $< \theta \leftarrow \theta + \alpha \Delta$ • To apply this algorithm to neural network training, we need • the loss function $l(\mathbf{f}(\mathbf{x}^{(t)}; \theta), y^{(t)})$

- a procedure to compute the parameter gradients $abla_{m{ heta}} l(\mathbf{f}(\mathbf{x}^{(t)};m{ heta}),y^{(t)})$
- + the regularizer $\Omega({m heta})$ (and the gradient $abla_{{m heta}}\Omega({m heta})$)
- initialization method

REGULARIZATION

Topics: L2 regularization

$$\Omega(\boldsymbol{\theta}) = \sum_{k} \sum_{i} \sum_{j} \left(W_{i,j}^{(k)} \right)^2 = \sum_{k} ||\mathbf{W}^{(k)}||_F^2$$

• Gradient:
$$abla_{\mathbf{W}^{(k)}}\Omega(oldsymbol{ heta}) = 2\mathbf{W}^{(k)}$$

- Only applied on weights, not on biases (weight decay)
- Can be interpreted as having a Gaussian prior over the weights

REGULARIZATION

Topics: LI regularization

$$\Omega(\boldsymbol{\theta}) = \sum_{k} \sum_{i} \sum_{j} |W_{i,j}^{(k)}|$$

• Gradient:
$$abla_{\mathbf{W}^{(k)}}\Omega(oldsymbol{ heta}) = \mathrm{sign}(\mathbf{W}^{(k)})$$

• where
$$\operatorname{sign}(\mathbf{W}^{(k)})_{i,j} = 1_{\mathbf{W}^{(k)}_{i,j} > 0} - 1_{\mathbf{W}^{(k)}_{i,j} < 0}$$

- Also only applied on weights
- Unlike L2, L1 will push certain weights to be exactly 0
- Can be interpreted as having a Laplacian prior over the weights

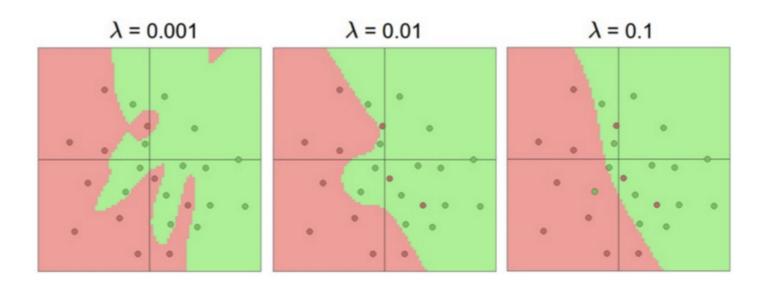
Empirical Risk Minimization

Topics: empirical risk minimization, regularization

- Empirical risk minimization
 - framework to design learning algorithms

$$\underset{\boldsymbol{\theta}}{\operatorname{arg\,min}} \frac{1}{T} \sum_{t} l(f(\mathbf{x}^{(t)}; \boldsymbol{\theta}), y^{(t)}) + \lambda \Omega(\boldsymbol{\theta})$$

- + $l(f(\mathbf{x}^{(t)}; \boldsymbol{ heta}), y^{(t)})$ is a loss function
- + $\Omega(oldsymbol{ heta})$ is a regularizer (penalizes certain values of $oldsymbol{ heta}$)
- Learning is cast as optimization
 - ideally, we'd optimize classification error, but it's not smooth
 - loss function is a surrogate for what we truly should optimize (e.g. upper bound)



[http://cs231n.github.io/neural-networks-1/]

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INITIALIZATION

size of $\mathbf{h}^{(k)}(\mathbf{x})$

Topics: initialization

- For biases
 - initialize all to 0
- For weights
 - · Can't initialize weights to 0 with tanh activation
 - we can show that all gradients would then be 0 (saddle point)
 - Can't initialize all weights to the same value
 - we can show that all hidden units in a layer will always behave the same
 - need to break symmetry
 - Recipe: sample $\mathbf{W}_{i,j}^{(k)}$ from $U\left[-b,b
 ight]$ where $b=\frac{\sqrt{6}}{\sqrt{H_k+H_{k-1}}}$
 - the idea is to sample around 0 but break symmetry
 - other values of *b* could work well (not an exact science) (see Glorot & Bengio, 2010)

The Learning Algorithm Topics: stochastic gradient descent (SGD) • Algorithm that performs updates after each example • initialize θ ($\theta \equiv \{W^{(1)}, b^{(1)}, \dots, W^{(L+1)}, b^{(L+1)}\}$) • for N iterations • for each training example $(\mathbf{x}^{(t)}, y^{(t)})$ $\prec \Delta = -\nabla_{\theta} l(f(\mathbf{x}^{(t)}; \theta), y^{(t)}) - \lambda \nabla_{\theta} \Omega(\theta)$ $\prec \theta \leftarrow \theta + \alpha \Delta$ • To apply this algorithm to neural network training, we need • the loss function $l(\mathbf{f}(\mathbf{x}^{(t)}; \theta), y^{(t)})$

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Toolkits

- TensorFlow
 - <u>https://www.tensorflow.org/</u>
- Theano (not maintained any more)
 - <u>http://deeplearning.net/software/theano/</u>
- PyTorch
 - http://pytorch.org/