CSCI567 Machine Learning (Fall 2018)

Prof. Haipeng Luo

U of Southern California

Sep 12, 2018

Administration

GitHub repos are setup (ask TA Chi Zhang for any issues)

HW 1 is due this Sunday (09/16) 11:59PM

You need to submit a form if you use late days (see course website)

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Effort-based grade for written assignments:

- see the explanation on Piazza
- key: let us know what you have tried and how you thought
- "I spend an hour and came up with nothing" = empty solution

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Outline

Review of Last Lecture

Multiclass Classification

Neural Nets

Outline

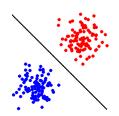
- Review of Last Lecture
- 2 Multiclass Classification
- Neural Nets

Summary

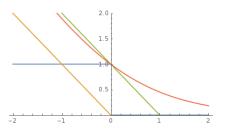
Linear models for binary classification:

Step 1. Model is the set of separating hyperplanes

$$\mathcal{F} = \{ f(\boldsymbol{x}) = \operatorname{sgn}(\boldsymbol{w}^{\mathrm{T}}\boldsymbol{x}) \mid \boldsymbol{w} \in \mathbb{R}^{\mathsf{D}} \}$$



Step 2. Pick the surrogate loss



- perceptron loss $\ell_{perceptron}(z) = \max\{0, -z\}$ (used in Perceptron)
- hinge loss $\ell_{\mathsf{hinge}}(z) = \max\{0, 1-z\}$ (used in SVM and many others)
- \bullet logistic loss $\ell_{\rm logistic}(z) = \log(1 + \exp(-z))$ (used in logistic regression)

Step 3. Find empirical risk minimizer (ERM):

$$\boldsymbol{w}^* = \operatorname*{argmin}_{\boldsymbol{w} \in \mathbb{R}^{\mathsf{D}}} F(\boldsymbol{w}) = \operatorname*{argmin}_{\boldsymbol{w} \in \mathbb{R}^{\mathsf{D}}} \sum_{n=1}^{N} \ell(y_n \boldsymbol{w}^{\mathsf{T}} \boldsymbol{x}_n)$$

using

- GD: $\boldsymbol{w} \leftarrow \boldsymbol{w} \eta \nabla F(\boldsymbol{w})$
- SGD: $\boldsymbol{w} \leftarrow \boldsymbol{w} \eta \tilde{\nabla} F(\boldsymbol{w})$
- Newton: $\boldsymbol{w} \leftarrow \boldsymbol{w} \left(\nabla^2 F(\boldsymbol{w})\right)^{-1} \nabla F(\boldsymbol{w})$

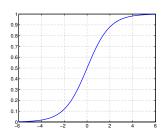
A Probabilistic view of logistic regression

Minimizing logistic loss = MLE for the sigmoid model

$$\boldsymbol{w}^* = \operatorname*{argmin}_{\boldsymbol{w}} \sum_{n=1}^N \ell_{\mathsf{logistic}}(y_n \boldsymbol{w}^{\mathrm{T}} \boldsymbol{x}_n) = \operatorname*{argmax}_{\boldsymbol{w}} \prod_{n=1}^N \mathbb{P}(y_n \mid \boldsymbol{x}_n; \boldsymbol{w})$$

where

$$\mathbb{P}(y \mid \boldsymbol{x}; \boldsymbol{w}) = \sigma(y \boldsymbol{w}^{\mathrm{T}} \boldsymbol{x}) = \frac{1}{1 + e^{-y \boldsymbol{w}^{\mathrm{T}} \boldsymbol{x}}}$$



Outline

- Review of Last Lecture
- Multiclass Classification
 - Multinomial logistic regression
 - Reduction to binary classification
- 3 Neural Nets

Classification

Recall the setup:

- ullet input (feature vector): $oldsymbol{x} \in \mathbb{R}^{\mathsf{D}}$
- output (label): $y \in [C] = \{1, 2, \dots, C\}$
- ullet goal: learn a mapping $f:\mathbb{R}^{\mathsf{D}} o [\mathsf{C}]$

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Examples:

- recognizing digits (C = 10) or letters (C = 26 or 52)
- predicting weather: sunny, cloudy, rainy, etc
- predicting image category: ImageNet dataset (C $\approx 20K$)

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Nearest Neighbor Classifier naturally works for arbitrary C.

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Note: a linear model for binary tasks (switching from $\{-1, +1\}$ to $\{1, 2\}$)

$$f(\boldsymbol{x}) = \begin{cases} 1 & \text{if } \boldsymbol{w}^{\mathrm{T}} \boldsymbol{x} \ge 0 \\ 2 & \text{if } \boldsymbol{w}^{\mathrm{T}} \boldsymbol{x} < 0 \end{cases}$$

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can be written as

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for any w_1, w_2 s.t. $w = w_1 - w_2$



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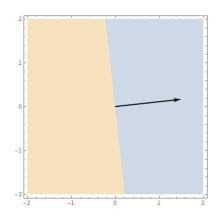
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for any $oldsymbol{w}_1, oldsymbol{w}_2$ s.t. $oldsymbol{w} = oldsymbol{w}_1 - oldsymbol{w}_2$

Think of $\boldsymbol{w}_k^{\mathrm{T}}\boldsymbol{x}$ as a score for class k.



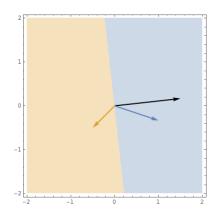
$$\boldsymbol{w} = (\frac{3}{2}, \frac{1}{6})$$

Blue class:

$$\{\boldsymbol{x}: \boldsymbol{w}^{\mathrm{T}} \boldsymbol{x} \geq 0\}$$

• Orange class:

$$\{\boldsymbol{x}: \boldsymbol{w}^{\mathrm{T}}\boldsymbol{x} < 0\}$$



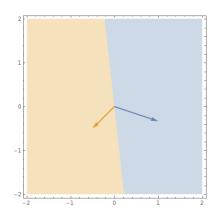
$$w = (\frac{3}{2}, \frac{1}{6}) = w_1 - w_2$$

 $w_1 = (1, -\frac{1}{3})$
 $w_2 = (-\frac{1}{2}, -\frac{1}{2})$

Blue class:

 $\{ \boldsymbol{x} : 1 = \operatorname{argmax}_k \boldsymbol{w}_k^{\mathrm{T}} \boldsymbol{x} \}$

• Orange class: $\{ \boldsymbol{x} : 2 = \operatorname{argmax}_k \boldsymbol{w}_k^{\mathrm{T}} \boldsymbol{x} \}$



$$\mathbf{w}_1 = (1, -\frac{1}{3})$$

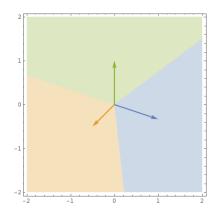
 $\mathbf{w}_2 = (-\frac{1}{2}, -\frac{1}{2})$

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$$\mathbf{w}_1 = (1, -\frac{1}{3})$$

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 $\mathbf{w}_3 = (0, 1)$

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• Green class:

$$\{\boldsymbol{x}: \boldsymbol{3} = \operatorname{argmax}_k \boldsymbol{w}_k^{\mathrm{T}} \boldsymbol{x}\}$$

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How do we generalize perceptron/hinge/logistic loss?

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How do we generalize perceptron/hinge/logistic loss?

This lecture: focus on the more popular logistic loss

Multinomial logistic regression: a probabilistic view

Observe: for binary logistic regression, with $w = w_1 - w_2$:

$$\mathbb{P}(y = 1 \mid \boldsymbol{x}; \boldsymbol{w}) = \sigma(\boldsymbol{w}^{\mathrm{T}} \boldsymbol{x}) = \frac{1}{1 + e^{-\boldsymbol{w}^{\mathrm{T}} \boldsymbol{x}}} = \frac{e^{\boldsymbol{w}_{1}^{\mathrm{T}} \boldsymbol{x}}}{e^{\boldsymbol{w}_{1}^{\mathrm{T}} \boldsymbol{x}} + e^{\boldsymbol{w}_{2}^{\mathrm{T}} \boldsymbol{x}}} \propto e^{\boldsymbol{w}_{1}^{\mathrm{T}} \boldsymbol{x}}$$

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This is called the *softmax function*.

Maximize probability of see labels $y_1,\ldots,y_{\mathsf{N}}$ given ${m x}_1,\ldots,{m x}_{\mathsf{N}}$

$$P(\boldsymbol{W}) = \prod_{n=1}^{\mathsf{N}} \mathbb{P}(y_n \mid \boldsymbol{x}_n; \boldsymbol{W}) = \prod_{n=1}^{\mathsf{N}} \frac{e^{\boldsymbol{w}_{y_n}^{\mathsf{T}} \boldsymbol{x}}}{\sum_{k \in [\mathsf{C}]} e^{\boldsymbol{w}_k^{\mathsf{T}} \boldsymbol{x}}}$$

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By taking **negative log**, this is equivalent to minimizing

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This is the multiclass logistic loss, a.k.a cross-entropy loss.

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When C = 2, this is the same as binary logistic loss.



Optimization

Apply SGD: what is the gradient of

$$g(\boldsymbol{W}) = \ln \left(1 + \sum_{k' \neq y_n} e^{(\boldsymbol{w}_{k'} - \boldsymbol{w}_{y_n})^{\mathrm{T}} \boldsymbol{x}_n} \right) ?$$

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else:

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SGD for multinomial logistic regression

Initialize W = 0 (or randomly). Repeat:

- $oldsymbol{0}$ pick $n \in [\mathsf{N}]$ uniformly at random
- update the parameters

$$m{W} \leftarrow m{W} - \eta \left(egin{array}{ccc} \mathbb{P}(y = 1 \mid m{x}_n; m{W}) & & & \\ & dash & & & \\ \mathbb{P}(y = y_n \mid m{x}_n; m{W}) - 1 & & & \\ & dash & & dash & \\ & \mathbb{P}(y = \mathsf{C} \mid m{x}_n; m{W}) & & \end{array}
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Think about why the algorithm makes sense intuitively.

Having learned $oldsymbol{W}$, we can either

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In either case, (expected) mistake is bounded by logistic loss

Having learned $oldsymbol{W}$, we can either

- ullet make a deterministic prediction $rgmax_{k \in [\mathsf{C}]} ullet w_k^\mathrm{T} oldsymbol{x}$
- ullet make a $\emph{randomized}$ prediction according to $\mathbb{P}(k\mid m{x};m{W}) \propto e^{m{w}_k^{\mathrm{T}}m{x}}$

In either case, (expected) mistake is bounded by logistic loss

deterministic

$$\mathbb{I}[f(\boldsymbol{x}) \neq y] \leq \log_2 \left(1 + \sum_{k \neq y} e^{(\boldsymbol{w}_k - \boldsymbol{w}_y)^{\mathrm{T}} \boldsymbol{x}} \right)$$

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In either case, (expected) mistake is bounded by logistic loss

deterministic

$$\mathbb{I}[f(\boldsymbol{x}) \neq y] \leq \log_2 \left(1 + \sum_{k \neq y} e^{(\boldsymbol{w}_k - \boldsymbol{w}_y)^{\mathrm{T}} \boldsymbol{x}} \right)$$

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Reduce multiclass to binary

Is there an *even more general and simpler approach* to derive multiclass classification algorithms?

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Given a binary classification algorithm (any one, not just linear methods), can we turn it to a multiclass algorithm, in a black-box manner?

Yes, there are in fact many ways to do it.

- one-versus-all (one-versus-rest, one-against-all, etc)
- one-versus-one (all-versus-all, etc)
- Error-Correcting Output Codes (ECOC)
- tree-based reduction

(picture credit: link)

Idea: train C binary classifiers to learn "is class k or not?" for each k.

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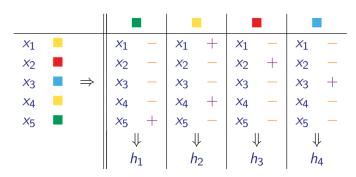
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Issue: will (probably) make a mistake as long as one of h_k errs.

(picture credit: link)

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		■ vs. ■		■ vs. ■		■ vs. ■		■ vs. ■		■ vs. ■		■ vs. ■	
x_1		<i>x</i> ₁	_					<i>x</i> ₁	_			<i>x</i> ₁	_
<i>x</i> ₂				<i>x</i> ₂	_	<i>x</i> ₂	+					<i>x</i> ₂	+
<i>X</i> 3	\Rightarrow					<i>X</i> 3	_	<i>X</i> 3	+	<i>X</i> 3	_		
<i>X</i> ₄		<i>X</i> ₄	_					<i>X</i> 4	_			<i>X</i> 4	_
<i>X</i> 5		<i>X</i> 5	+	<i>X</i> 5	+					<i>X</i> 5	+		
		↓		↓ ↓		↓		↓		\		↓	
		$h_{(1,2)}$		$h_{(1,3)}$		$h_{(3,4)}$		$h_{(4,2)}$		$h_{(1,4)}$		$h_{(3,2)}$	

Prediction: for a new example x

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More robust than one-versus-all, but slower in prediction.

(picture credit: link)

Idea: based on a code $M \in \{-1,+1\}^{\mathsf{C} \times \mathsf{L}}$, train L binary classifiers to learn "is bit b on or off".

(picture credit: link)

Idea: based on a code $M \in \{-1, +1\}^{\mathsf{C} \times \mathsf{L}}$, train L binary classifiers to learn "is bit b on or off".

Training: for each bit $b \in [\mathsf{L}]$

- ullet relabel example x_n as $M_{y_n,b}$
- train a binary classifier h_b using this new dataset.

М	1	2	3	4	5
	+	_	+	_	+
	_	- + +	+	+	+
	+	+	_	_	_
	+	+	+	+	_

		1	1		2		3		4		5	
<i>x</i> ₁		<i>x</i> ₁	_	<i>x</i> ₁	_		+	<i>x</i> ₁	+	<i>x</i> ₁	+	
<i>x</i> ₂		<i>x</i> ₂	+	<i>x</i> ₂	+				_	<i>x</i> ₂	_	
<i>X</i> 3	\Rightarrow	<i>X</i> 3	+					<i>X</i> 3	+	<i>X</i> 3	_	
<i>X</i> ₄		<i>X</i> ₄	_	<i>X</i> ₄	_	<i>X</i> ₄				<i>X</i> ₄	+	
<i>X</i> 5		<i>X</i> 5		<i>X</i> 5	_	<i>X</i> 5			_	<i>X</i> 5	+	
		↓	\Downarrow		↓		↓					
		h_1		h_2		h_3		h_4		h_5		

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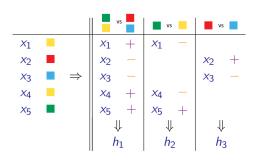
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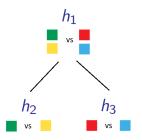
- the more dissimilar the codes between different classes are, the better
- random code is a good choice, but might create hard training sets

Idea: train \approx C binary classifiers to learn "belongs to which half?".

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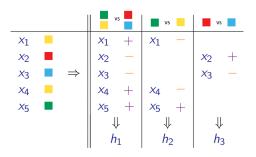
Training: see pictures

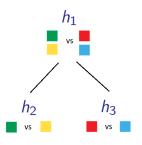




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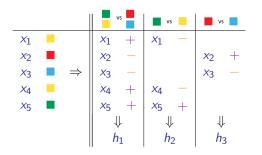


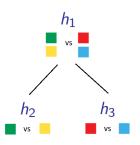


Prediction is also natural,

Idea: train \approx C binary classifiers to learn "belongs to which half?".

Training: see pictures





Prediction is also natural, *but is very fast!* (think ImageNet where $C \approx 20K$)

Reduction	#training points	test time	remark
OvA			
OvO			
ECOC			
Tree			

Reduction	#training points	test time	remark
OvA	CN		
OvO			
ECOC			
Tree			

Reduction	#training points	test time	remark
OvA	CN	С	
OvO			
ECOC			
Tree			

Reduction	#training points	test time	remark
OvA	CN	С	not robust
OvO			
ECOC			
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Tree			

Reduction	#training points	test time	remark
OvA	CN	С	not robust
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ECOC			
Tree			

Reduction	#training points	test time	remark
OvA	CN	С	not robust
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Tree			

Reduction	#training points	test time	remark
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ECOC	LN		
Tree			

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ECOC	LN	L	
Tree			

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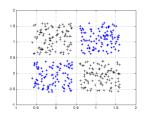
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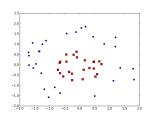
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Tree	$(\log_2C)N$	\log_2C	good for "extreme classification"

Outline

- Review of Last Lecture
- 2 Multiclass Classification
- Neural Nets
 - Definition
 - Backpropagation
 - Preventing overfitting

Linear models are not always adequate

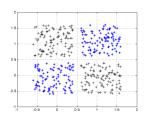


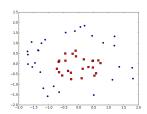


We can use a nonlinear mapping as discussed:

$$\phi(x): x \in \mathbb{R}^\mathsf{D}
ightarrow z \in \mathbb{R}^\mathsf{M}$$

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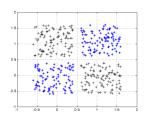


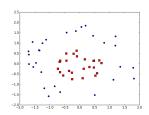
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But what kind of nonlinear mapping ϕ should be used? Can we actually learn this nonlinear mapping?

Linear models are not always adequate





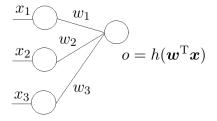
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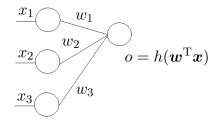
THE most popular nonlinear models nowadays: neural nets

Linear model as a one-layer neural net



h(a) = a for linear model

Linear model as a one-layer neural net



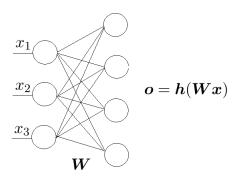
h(a) = a for linear model

To create non-linearity, can use

- Rectified Linear Unit (ReLU): $h(a) = \max\{0, a\}$
- sigmoid function: $h(a) = \frac{1}{1+e^{-a}}$
- TanH: $h(a) = \frac{e^a e^{-a}}{e^a + e^{-a}}$
- many more



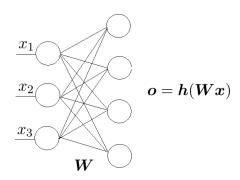
More output nodes



$$m{W} \in \mathbb{R}^{4 imes 3}$$
, $m{h}: \mathbb{R}^4 o \mathbb{R}^4$ so $m{h}(m{a}) = (h_1(a_1), h_2(a_2), h_3(a_3), h_4(a_4))$



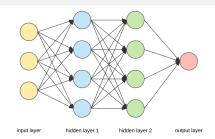
More output nodes



$$W \in \mathbb{R}^{4 \times 3}$$
, $h : \mathbb{R}^4 \to \mathbb{R}^4$ so $h(a) = (h_1(a_1), h_2(a_2), h_3(a_3), h_4(a_4))$

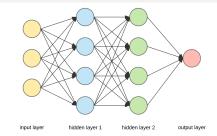
Can think of this as a nonlinear basis: $\Phi(m{x}) = m{h}(m{W}m{x})$

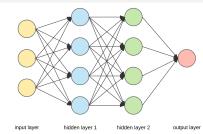




Becomes a network:

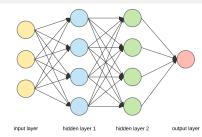
• each node is called a neuron



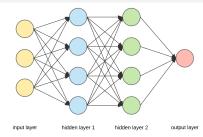


- each node is called a neuron
- h is called the activation function
 - can use h(a) = 1 for one neuron in each layer to *incorporate bias term*
 - output neuron can use h(a) = a

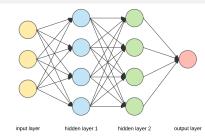




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- deep neural nets can have many layers and millions of parameters
- this is a **feedforward**, **fully connected** neural net, there are many variants

How powerful are neural nets?

Universal approximation theorem (Cybenko, 89; Hornik, 91):

A feedforward neural net with a single hidden layer can approximate any continuous functions.

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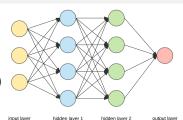
Designing network architecture is important and very complicated

• for feedforward network, need to decide number of hidden layers, number of neurons at each layer, activation functions, etc.

Math formulation

An L-layer neural net can be written as

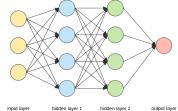
$$oldsymbol{f}(oldsymbol{x}) = oldsymbol{h}_{\mathsf{L}} \left(oldsymbol{W}_{L} oldsymbol{h}_{\mathsf{L}-1} \left(oldsymbol{W}_{L-1} \cdots oldsymbol{h}_{1} \left(oldsymbol{W}_{1} oldsymbol{x}
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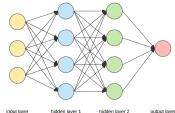
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where

- $oldsymbol{W}_\ell \in \mathbb{R}^{\mathsf{D}_\ell imes \mathsf{D}_{\ell-1}}$ is the weights for layer ℓ
- ullet $D_0 = D, D_1, \dots, D_L$ are numbers of neurons at each layer
- $oldsymbol{a}_\ell \in \mathbb{R}^{\mathsf{D}_\ell}$ is input to layer ℓ
- $oldsymbol{o}_\ell \in \mathbb{R}^{\mathsf{D}_\ell}$ is output to layer ℓ
- $m{h}:\mathbb{R}^{\mathsf{D}_\ell} o\mathbb{R}^{\mathsf{D}_\ell}$ is activation functions at layer ℓ



Learning the model

No matter how complicated the model is, our goal is the same: minimize

$$\mathcal{E}(\boldsymbol{W}_1,\ldots,\boldsymbol{W}_{\mathsf{L}}) = \sum_{n=1}^{\mathsf{N}} \mathcal{E}_n(\boldsymbol{W}_1,\ldots,\boldsymbol{W}_{\mathsf{L}})$$

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where

$$\mathcal{E}_n(\boldsymbol{W}_1,\dots,\boldsymbol{W}_{\mathsf{L}}) = \begin{cases} \|\boldsymbol{f}(\boldsymbol{x}_n) - \boldsymbol{y}_n\|_2^2 & \text{for regression} \\ \ln\left(1 + \sum_{k \neq y_n} e^{f(\boldsymbol{x}_n)_k - f(\boldsymbol{x}_n)_{y_n}}\right) & \text{for classification} \end{cases}$$

Same thing: apply **SGD**! even if the model is *nonconvex*.

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What is the gradient of this complicated function?

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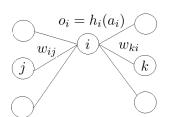
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the simplest example $f(g_1(w), g_2(w)) = g_1(w)g_2(w)$

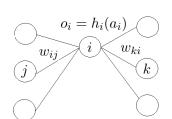


Drop the subscript ℓ for layer for simplicity.

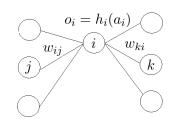


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$$\frac{\partial \mathcal{E}_n}{\partial w_{ij}} = \frac{\partial \mathcal{E}_n}{\partial a_i} \frac{\partial a_i}{\partial w_{ij}}$$

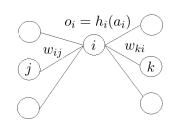


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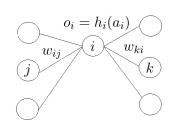
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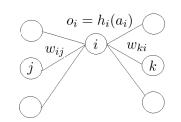
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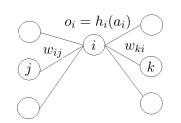


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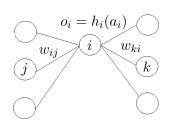
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Adding the subscript for layer:

$$\frac{\partial \mathcal{E}_n}{\partial w_{\ell,ij}} = \frac{\partial \mathcal{E}_n}{\partial a_{\ell,i}} o_{\ell-1,j}$$

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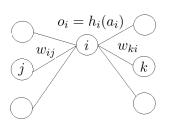
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For the last layer, for square loss

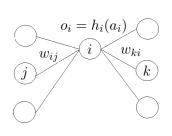
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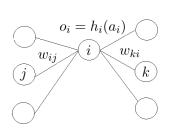
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Exercise: try to do it for logistic loss yourself.



Using matrix notation greatly simplifies presentation and implementation:

$$\frac{\partial \mathcal{E}_n}{\partial \boldsymbol{W}_{\ell}} = \frac{\partial \mathcal{E}_n}{\partial \boldsymbol{a}_{\ell}} \boldsymbol{o}_{\ell-1}^{\mathrm{T}}$$

$$\frac{\partial \mathcal{E}_n}{\partial \boldsymbol{a}_{\ell}} = \begin{cases} \left(\boldsymbol{W}_{\ell+1}^{\mathrm{T}} \frac{\partial \mathcal{E}_n}{\partial \boldsymbol{a}_{\ell+1}}\right) \circ \boldsymbol{h}'_{\ell}(\boldsymbol{a}_{\ell}) & \text{if } \ell < \mathsf{L} \\ 2(\boldsymbol{h}_{\mathsf{L}}(\boldsymbol{a}_{\mathsf{L}}) - \boldsymbol{y}_n) \circ \boldsymbol{h}'_{\mathsf{L}}(\boldsymbol{a}_{\mathsf{L}}) & \text{else} \end{cases}$$

where $v_1 \circ v_2 = (v_{11}v_{21}, \cdots, v_{1D}v_{2D})$ is the element-wise product (a.k.a. Hadamard product).

Verify yourself!



The **backpropagation** algorithm (**Backprop**)

Initialize W_1, \ldots, W_L (all 0 or randomly). Repeat:

 $\textbf{ 1} \text{ randomly pick one data point } n \in [\mathsf{N}]$

The backpropagation algorithm (Backprop)

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 - ullet compute $oldsymbol{a}_\ell = oldsymbol{W}_\ell oldsymbol{o}_{\ell-1}$ and $oldsymbol{o}_\ell = oldsymbol{h}_\ell (oldsymbol{a}_\ell)$

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update weights

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Think about how to do the last two steps properly!

More tricks to optimize neural nets

Many variants based on backprop

- SGD with minibatch: randomly sample a batch of examples to form a stochastic gradient
- SGD with momentum
- o . . .

SGD with momentum

Initialize $oldsymbol{w}_0$ and $oldsymbol{ ext{velocity}} oldsymbol{v} = oldsymbol{0}$

For t = 1, 2, ...

- ullet form a stochastic gradient $oldsymbol{g}_t$
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Updates for first few rounds:

- $w_1 = w_0 \eta g_1$
- $\bullet \ \boldsymbol{w}_2 = \boldsymbol{w}_1 \alpha \eta \boldsymbol{g}_1 \eta \boldsymbol{g}_2$
- $\mathbf{w}_3 = \mathbf{w}_2 \alpha^2 \eta \mathbf{g}_1 \alpha \eta \mathbf{g}_2 \eta \mathbf{g}_3$
-



Overfitting

Overfitting is very likely since the models are too powerful.

Methods to overcome overfitting:

- data augmentation
- regularization
- dropout
- early stopping
-

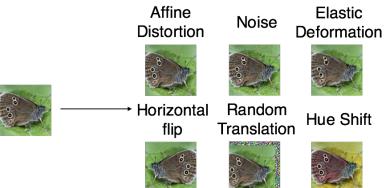
Data augmentation

Data: the more the better. How do we get more data?

Data augmentation

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Exploit prior knowledge to add more training data



Regularization

L2 regularization: minimize

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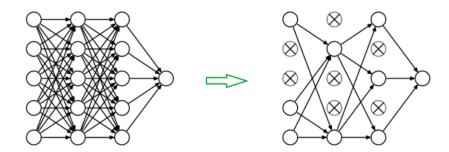
Simple change to the gradient:

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Introduce weight decaying effect

Dropout

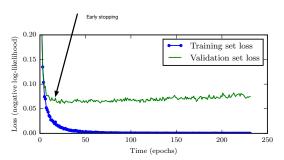
Randomly delete neurons during training



Very effective, makes training faster as well

Early stopping

Stop training when the performance on validation set stops improving



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