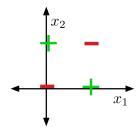
CS 4995 Lecture 2: Multilayer Perceptrons & Backpropagation

Richard Zemel

Richard Zemel

CS 4995 Lecture 2: Multilayer Perceptrons &

- Single neurons (linear classifiers) are very limited in expressive power.
- **XOR** is a classic example of a function that's not linearly separable.



• There's an elegant proof using convexity.

Convex Sets



• A set S is convex if any line segment connecting points in S lies entirely within S. Mathematically,

$$\mathbf{x}_1, \mathbf{x}_2 \in \mathcal{S} \implies \lambda \mathbf{x}_1 + (1 - \lambda) \mathbf{x}_2 \in \mathcal{S} \text{ for } \mathbf{0} \leq \lambda \leq 1.$$

A simple inductive argument shows that for x₁,..., x_N ∈ S, weighted averages, or convex combinations, lie within the set:

$$\lambda_1 \mathbf{x}_1 + \dots + \lambda_N \mathbf{x}_N \in S \quad \text{for } \lambda_i > 0, \ \lambda_1 + \dots + \lambda_N = 1.$$

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Showing that XOR is not linearly separable

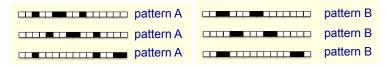
- Half-spaces are obviously convex.
- Suppose there were some feasible hypothesis. If the positive examples are in the positive half-space, then the green line segment must be as well.
- Similarly, the red line segment must line within the negative half-space.



• But the intersection can't lie in both half-spaces. Contradiction!

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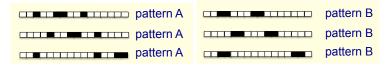
A more troubling example



- These images represent 16-dimensional vectors. White = 0, black = 1.
- Want to distinguish patterns A and B in all possible translations (with wrap-around)
- Translation invariance is commonly desired in vision!

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A more troubling example



- These images represent 16-dimensional vectors. White = 0, black = 1.
- Want to distinguish patterns A and B in all possible translations (with wrap-around)
- Translation invariance is commonly desired in vision!
- Suppose there's a feasible solution. The average of all translations of A is the vector (0.25, 0.25, ..., 0.25). Therefore, this point must be classified as A.
- Similarly, the average of all translations of B is also (0.25, 0.25, ..., 0.25). Therefore, it must be classified as B. Contradiction!

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• Sometimes we can overcome this limitation using feature maps, just like for linear regression. E.g., for **XOR**:

$$\psi(\mathbf{x}) = \begin{pmatrix} x_1 \\ x_2 \\ x_1 x_2 \end{pmatrix}$$

$$x_1 \quad x_2 \quad \phi_1(\mathbf{x}) \quad \phi_2(\mathbf{x}) \quad \phi_3(\mathbf{x}) \quad t$$

$$0 \quad 0 \quad 0 \quad 0 \quad 0$$

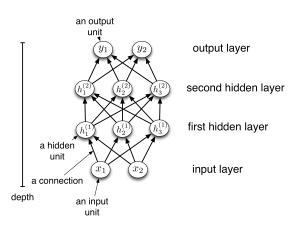
$$0 \quad 1 \quad 0 \quad 1 \quad 0 \quad 1$$

$$1 \quad 0 \quad 1 \quad 0 \quad 0 \quad 1$$

$$1 \quad 1 \quad 1 \quad 1 \quad 0$$

- This is linearly separable. (Try it!)
- Not a general solution: it can be hard to pick good basis functions. Instead, we'll use neural nets to learn nonlinear hypotheses directly.

- We can connect lots of units together into a directed acyclic graph.
- This gives a feed-forward neural network. That's in contrast to recurrent neural networks, which can have cycles. (We'll talk about those later.)
- Typically, units are grouped together into layers.

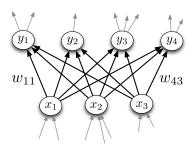


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- Each layer connects N input units to M output units.
- In the simplest case, all input units are connected to all output units. We call this
 a fully connected layer. We'll consider other layer types later.
- Note: the inputs and outputs for a layer are distinct from the inputs and outputs to the network.
- Recall from softmax regression: this means we need an $M \times N$ weight matrix.
- The output units are a function of the input units:

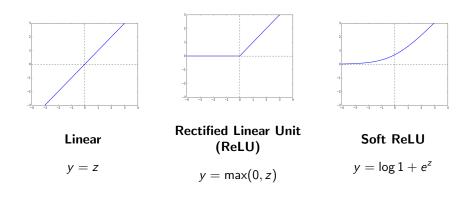
 $\mathbf{y} = f(\mathbf{x}) = \phi \left(\mathbf{W} \mathbf{x} + \mathbf{b} \right)$

• A multilayer network consisting of fully connected layers is called a multilayer perceptron. Despite the name, it has nothing to do with perceptrons!



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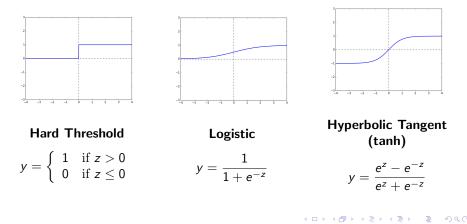
Some activation functions:



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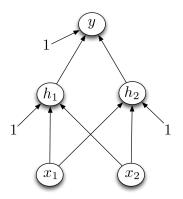
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Some activation functions:

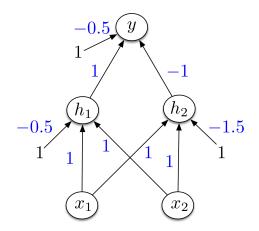


Designing a network to compute XOR:

Assume hard threshold activation function



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• Each layer computes a function, so the network computes a composition of functions:

$$h^{(1)} = f^{(1)}(\mathbf{x})$$

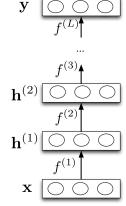
$$h^{(2)} = f^{(2)}(\mathbf{h}^{(1)})$$

$$\vdots$$

$$\mathbf{y} = f^{(L)}(\mathbf{h}^{(L-1)})$$

• Or more simply:

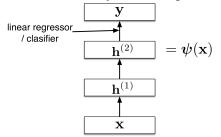
$$\mathbf{y} = f^{(L)} \circ \cdots \circ f^{(1)}(\mathbf{x}).$$



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• Neural nets provide modularity: we can implement each layer's computations as a black box.

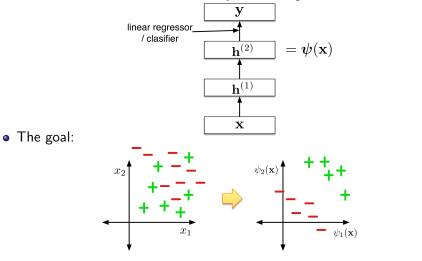
• Neural nets can be viewed as a way of learning features:



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• Neural nets can be viewed as a way of learning features:



Input representation of a digit : 784 dimensional vector.

0.0 221.0 115.0 0.0 0.0 0.0 0.0 0.0 0.0 0.0 0.0 0.0 0.0 0.0 0.0 0.0 0.0 0.0 0.0 0.0 0.0 51.0 254.0 145.0 0.0 0.0 0.0 22.0 230.0 134.0 0.0 0.0 0.0 0.0 0.0 0.0 0.0 0.0 0.0 0.0 0.0 0.0 0.0 0.0 0.0 165.0 254.0 115.0 0.0 0.0 24.0 251.0 104.0 0.0 0.0 0.0 0.0 0.0 0.0 0.0 0.0 0.0 0.0 0.0 0.0 0.0 0.0 51.0 234.0 254.0 81.0 0.0 0.0 91.0 253.0 184.0 0.0 0.0 0.0 0.0 0.0 0.0 0.0 0.0 0.0 0.0 0.0 0.0 0.0 0.0 13.0 221.0 254.0 160.0 0.0 0.0 0.0 0.0141.0254.0177.0 0.0 0.0 0.0 0.0 0.0 0.0 0.0 0.0 0.0 0.0 0.0114.0253.0253.076.0 0.0 0.0 0.0 207.0 253.0 93.0 0.0 0.0 0.0 0.0 0.0 0.0 0.0 0.0 0.0 0.0 0.0 0.0 15.0 232.0 253.0 102.0 0.0 0.0 0.0 0.0 0.0 0.0 207.0 253.0 17.0 0.0 0.0 0.0 0.0 0.0 0.0 0.0 0.0 0.0 0.0 0.0 229.0 253.0 202.0 19.0 0.0 0.0 0.0 0.0 0.0 34.0 240.0 253.0 0.0 0.0 0.0 0.0 0.0 0.0 0.0 0.0 0.0 0.0 0.0 7.0 170.0 254.0 254.0 46.0 0.0 0.0 0.0 0.0 0.0 0.0 47.0 254.0 254.0 0.0 0.0 0.0 0.0 0.0 0.0 0.0 0.0 0.0 0.0 0.0 0.0 0.0 0.0 24.0 253.0 254.0 253.0 234.0 163.0 47.0 47.0 26.0 0.0 0.0 130.0 253.0 253.0 0.0 0.0 0.0 0.0 0.0 0.0 0.0 0.0 0.0 0.0 0.0 0.0 21.0 246.0 254.0 253.0 253.0 253.0 254.0 254.0 253.0 232.0 174.0 208.0 232.0 253.0 177.0 0.0 0.0 0.0 0.0 0.0 0.0 0.0 0.0 0.0 0.0 0.0 0.0 0.0 04.0 101.0 211.0 219.0 219.0 254.0 253.0 253.0 253.0 254.0 253.0 244.0 09.0 0.0 0.0 0.0 0.0 0.0 0.0 0.0 0.0 0.0 0.0 0.0 0.0 0.0 0.0 0.0 0.0 0.0 0.0 0.0 93.0 142.0 142.0 93.0 170.0 254.0 230.0 153.0 253.0 213.0 0.0 170.0 253.0 137.0 0.0 220.0 253.0 137.0 0.0 60.0 225.0 254.0 80.0 93.0 254.0 253.0 46.0 0.0 51.0 254.0 215.0 9.8 0.0 0.0 0.0 0.0 0.0 0.0 0.0 0.0 0.0 0.0 0.0 0.0 0.0 0.0 0.0 0.0 0.0 0.0 0.0 105.0 105.0 0.0

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- Each first-layer hidden unit computes $\sigma(\mathbf{w}_i^T \mathbf{x})$
- Here is one of the weight vectors (also called a feature).
- It's reshaped into an image, with gray = 0, white = +, black = -.
- To compute $\mathbf{w}_i^T \mathbf{x}$, multiply the corresponding pixels, and sum the result.



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There are 256 first-level features total. Here are some of them.

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- We've seen that there are some functions that linear classifiers can't represent. Are deep networks any better?
- Any sequence of *linear* layers can be equivalently represented with a single linear layer.

$$\mathbf{y} = \underbrace{\mathbf{W}^{(3)}\mathbf{W}^{(2)}\mathbf{W}^{(1)}}_{\triangleq \mathbf{W}'} \mathbf{x}$$

- Deep linear networks are no more expressive than linear regression!
- Linear layers do have their uses stay tuned!

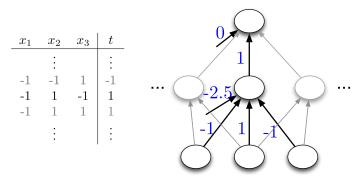
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- Multilayer feed-forward neural nets with *nonlinear* activation functions are <u>universal approximators</u>: they can approximate any function arbitrarily well.
- This has been shown for various activation functions (thresholds, logistic, ReLU, etc.)
 - Even though ReLU is "almost" linear, it's nonlinear enough!

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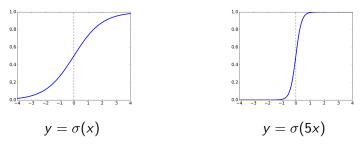
Universality for binary inputs and targets:

- Hard threshold hidden units, linear output
- Strategy: 2^D hidden units, each of which responds to one particular input configuration



• Only requires one hidden layer, though it needs to be extremely wide!

- What about the logistic activation function?
- You can approximate a hard threshold by scaling up the weights and biases:



• This is good: logistic units are differentiable, so we can tune them with gradient descent. (Stay tuned!)

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• Limits of universality

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- Limits of universality
 - You may need to represent an exponentially large network.
 - If you can learn any function, you'll just overfit.
 - Really, we desire a *compact* representation!

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- Limits of universality
 - You may need to represent an exponentially large network.
 - If you can learn any function, you'll just overfit.
 - Really, we desire a *compact* representation!
- We've derived units which compute the functions AND, OR, and NOT. Therefore, any Boolean circuit can be translated into a feed-forward neural net.
 - This suggests you might be able to learn *compact* representations of some complicated functions

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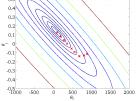
Overview

- We've seen that multilayer neural networks are powerful. But how can we actually learn them?
- Backpropagation is the central algorithm in this course.
 - It's is an algorithm for computing gradients.
 - Really it's an instance of reverse mode automatic differentiation, which is much more broadly applicable than just neural nets.
 - This is "just" a clever and efficient use of the Chain Rule for derivatives.
 - We'll see how to implement an automatic differentiation system next week.

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Recap: Gradient Descent

• **Recall:** gradient descent moves opposite the gradient (the direction of steepest descent)



- Weight space for a multilayer neural net: one coordinate for each weight or bias of the network, in *all* the layers
- Conceptually, not any different from what we've seen so far just higher dimensional and harder to visualize!
- We want to compute the cost gradient ${\rm d}\mathcal{J}/{\rm d}\bm{w},$ which is the vector of partial derivatives.
 - This is the average of $d\mathcal{L}/d\mathbf{w}$ over all the training examples, so in this lecture we focus on computing $d\mathcal{L}/d\mathbf{w}$.

- We've already been using the univariate Chain Rule.
- Recall: if f(x) and x(t) are univariate functions, then

$$\frac{\mathrm{d}}{\mathrm{d}t}f(x(t))=\frac{\mathrm{d}f}{\mathrm{d}x}\frac{\mathrm{d}x}{\mathrm{d}t}.$$

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Recall: Univariate logistic least squares model

$$z = wx + b$$

$$y = \sigma(z)$$

$$\mathcal{L} = \frac{1}{2}(y - t)^{2}$$

Let's compute the loss derivatives.

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How you would have done it in calculus class

$$\mathcal{L} = \frac{1}{2} (\sigma(wx+b)-t)^2$$

$$\frac{\partial \mathcal{L}}{\partial w} = \frac{\partial}{\partial w} \left[\frac{1}{2} (\sigma(wx+b)-t)^2 \right]$$

$$= \frac{1}{2} \frac{\partial}{\partial w} (\sigma(wx+b)-t)^2$$

$$= (\sigma(wx+b)-t) \frac{\partial}{\partial w} (\sigma(wx+b)-t)$$

$$= (\sigma(wx+b)-t) \sigma'(wx+b) \frac{\partial}{\partial w} (wx+b)$$

$$= (\sigma(wx+b)-t) \sigma'(wx+b) \frac{\partial}{\partial w} (wx+b)$$

$$= (\sigma(wx+b)-t) \sigma'(wx+b) \frac{\partial}{\partial w} (wx+b)$$

$$= (\sigma(wx+b)-t) \sigma'(wx+b) x$$

$$\frac{\partial \mathcal{L}}{\partial b} = \frac{\partial}{\partial b} \left[\frac{1}{2} (\sigma(wx+b)-t)^2 \right]$$

$$= \frac{1}{2} \frac{\partial}{\partial b} (\sigma(wx+b)-t)^2$$

$$= (\sigma(wx+b)-t) \frac{\partial}{\partial b} (\sigma(wx+b)-t)$$

$$= (\sigma(wx+b)-t) \sigma'(wx+b) \frac{\partial}{\partial b} (wx+b)$$

$$= (\sigma(wx+b)-t) \sigma'(wx+b) x$$

What are the disadvantages of this approach?

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A more structured way to do it

Computing the derivatives:

Computing the loss: z = wx + b $y = \sigma(z)$ $\mathcal{L} = \frac{1}{2}(y - t)^{2}$ $\frac{d\mathcal{L}}{dy} = y - t$ $\frac{d\mathcal{L}}{dz} = \frac{d\mathcal{L}}{dy} \sigma'(z)$ $\frac{\partial\mathcal{L}}{\partial w} = \frac{d\mathcal{L}}{dz} \times$ $\frac{\partial\mathcal{L}}{\partial b} = \frac{d\mathcal{L}}{dz}$

Remember, the goal isn't to obtain closed-form solutions, but to be able to write a program that efficiently computes the derivatives.

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- We can diagram out the computations using a computation graph.
- The nodes represent all the inputs and computed quantities, and the edges represent which nodes are computed directly as a function of which other nodes.

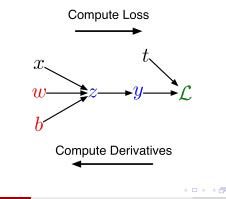


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A slightly more convenient notation:

- Use \overline{y} to denote the derivative $d\mathcal{L}/dy$, sometimes called the error signal.
- This emphasizes that the error signals are just values our program is computing (rather than a mathematical operation).
- This is not a standard notation, but I couldn't find another one that I liked.

Computing the loss:

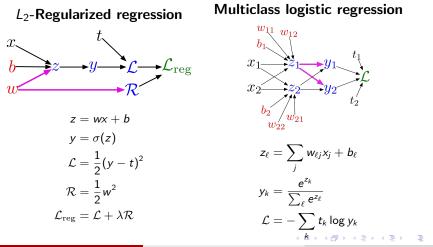
Computing the derivatives:

 $z = wx + b \qquad \qquad \overline{y} = y - t$ $y = \sigma(z) \qquad \qquad \overline{z} = \overline{y} \sigma'(z)$ $\mathcal{L} = \frac{1}{2}(y - t)^2 \qquad \qquad \overline{w} = \overline{z} x$ $\overline{b} = \overline{z}$

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Multivariate Chain Rule

Problem: what if the computation graph has fan-out > 1? This requires the multivariate Chain Rule!



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Multivariate Chain Rule

• Suppose we have a function f(x, y) and functions x(t) and y(t). (All the variables here are scalar-valued.) Then

$$\frac{\mathrm{d}}{\mathrm{d}t}f(x(t),y(t)) = \frac{\partial f}{\partial x}\frac{\mathrm{d}x}{\mathrm{d}t} + \frac{\partial f}{\partial y}\frac{\mathrm{d}y}{\mathrm{d}t}$$



• Example:

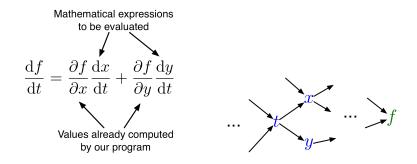
$$f(x, y) = y + e^{xy}$$
$$x(t) = \cos t$$
$$y(t) = t^{2}$$

• Plug in to Chain Rule:

$$\frac{\mathrm{d}f}{\mathrm{d}t} = \frac{\partial f}{\partial x}\frac{\mathrm{d}x}{\mathrm{d}t} + \frac{\partial f}{\partial y}\frac{\mathrm{d}y}{\mathrm{d}t}$$
$$= (ye^{xy}) \cdot (-\sin t) + (1 + xe^{xy}) \cdot 2t$$

Multivariable Chain Rule

• In the context of backpropagation:



In our notation:

$$\overline{t} = \overline{x} \, \frac{\mathrm{d}x}{\mathrm{d}t} + \overline{y} \, \frac{\mathrm{d}y}{\mathrm{d}t}$$

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Backpropagation

Full backpropagation algorithm:

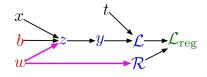
Let v_1, \ldots, v_N be a topological ordering of the computation graph (i.e. parents come before children.)

 v_N denotes the variable we're trying to compute derivatives of (e.g. loss).

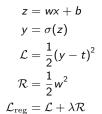
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Backpropagation

Example: univariate logistic least squares regression



Forward pass:



Backward pass:

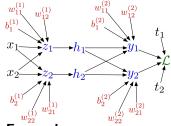
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Backpropagation

Multilayer Perceptron (multiple outputs):



Forward pass:

$$egin{aligned} &z_i = \sum_j w_{ij}^{(1)} x_j + b_i^{(1)} \ &h_i = \sigma(z_i) \ &y_k = \sum_i w_{ki}^{(2)} h_i + b_k^{(2)} \ &\mathcal{L} = rac{1}{2} \sum_k (y_k - t_k)^2 \end{aligned}$$

Backward pass:

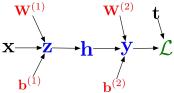
$$\begin{split} \overline{\mathcal{L}} &= 1\\ \overline{y_k} &= \overline{\mathcal{L}} \left(y_k - t_k \right)\\ \overline{w_{ki}^{(2)}} &= \overline{y_k} h_i\\ \overline{b_k^{(2)}} &= \overline{y_k}\\ \overline{h_i} &= \sum_k \overline{y_k} w_{ki}^{(2)}\\ \overline{z_i} &= \overline{h_i} \sigma'(z_i)\\ \overline{w_{ij}^{(1)}} &= \overline{z_i} x_j\\ \overline{b_i^{(1)}} &= \overline{z_i} \end{split}$$

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Richard Zemel

CS 4995 Lecture 2: Multilayer Perceptrons &

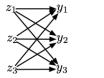
- Computation graphs showing individual units are cumbersome.
- As you might have guessed, we typically draw graphs over the vectorized variables.



• We pass messages back analogous to the ones for scalar-valued nodes.

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• Consider this computation graph:



• Backprop rules:

$$\overline{z_j} = \sum_k \overline{y_k} \frac{\partial y_k}{\partial z_j} \qquad \overline{z} = \frac{\partial \mathbf{y}}{\partial \mathbf{z}}^\top \overline{\mathbf{y}},$$

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where $\partial \mathbf{y} / \partial \mathbf{z}$ is the Jacobian matrix:

$$\frac{\partial \mathbf{y}}{\partial \mathbf{z}} = \begin{pmatrix} \frac{\partial y_1}{\partial z_1} & \cdots & \frac{\partial y_1}{\partial z_n} \\ \vdots & \ddots & \vdots \\ \frac{\partial y_m}{\partial z_1} & \cdots & \frac{\partial y_m}{\partial z_n} \end{pmatrix}$$

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Examples

Matrix-vector product

$$z = Wx$$
 $\frac{\partial z}{\partial x} = W$ $\overline{x} = W^{\top}\overline{z}$

• Elementwise operations

$$\mathbf{y} = \exp(\mathbf{z})$$
 $\frac{\partial \mathbf{y}}{\partial \mathbf{z}} = \begin{pmatrix} \exp(z_1) & 0 \\ & \ddots & \\ 0 & \exp(z_D) \end{pmatrix}$ $\overline{\mathbf{z}} = \exp(\mathbf{z}) \circ \overline{\mathbf{y}}$

• Note: we never explicitly construct the Jacobian. It's usually simpler and more efficient to compute the VJP directly.

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Full backpropagation algorithm (vector form):

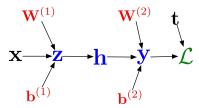
Let $\mathbf{v}_1, \ldots, \mathbf{v}_N$ be a topological ordering of the computation graph (i.e. parents come before children.)

 \mathbf{v}_N denotes the variable we're trying to compute derivatives of (e.g. loss). It's a scalar, which we can treat as a 1-D vector.

forward pass
$$\begin{bmatrix} & \text{For } i = 1, \dots, N \\ & \text{Compute } \mathbf{v}_i \text{ as a function of } Pa(\mathbf{v}_i \\ & \overline{\mathbf{v}_N} = 1 \\ & \text{For } i = N - 1, \dots, 1 \\ & \overline{\mathbf{v}_i} = \sum_{j \in Ch(\mathbf{v}_i)} \frac{\partial \mathbf{v}_j}{\partial \mathbf{v}_i}^\top \overline{\mathbf{v}_j} \end{bmatrix}$$

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MLP example in vectorized form:



Forward pass:

$$\begin{aligned} \mathbf{z} &= \mathbf{W}^{(1)}\mathbf{x} + \mathbf{b}^{(1)} \\ \mathbf{h} &= \sigma(\mathbf{z}) \\ \mathbf{y} &= \mathbf{W}^{(2)}\mathbf{h} + \mathbf{b}^{(2)} \\ \mathcal{L} &= \frac{1}{2}\|\mathbf{t} - \mathbf{y}\|^2 \end{aligned}$$

Backward pass:

$$\begin{split} \overline{\mathcal{L}} &= 1\\ \overline{\mathbf{y}} = \overline{\mathcal{L}} \left(\mathbf{y} - \mathbf{t} \right)\\ \overline{\mathbf{W}^{(2)}} &= \overline{\mathbf{y}} \mathbf{h}^{\top}\\ \overline{\mathbf{b}^{(2)}} &= \overline{\mathbf{y}}\\ \overline{\mathbf{h}} &= \mathbf{W}^{(2)\top} \overline{\mathbf{y}}\\ \overline{\mathbf{z}} &= \overline{\mathbf{h}} \circ \sigma'(\mathbf{z})\\ \overline{\mathbf{W}^{(1)}} &= \overline{\mathbf{z}} \mathbf{x}^{\top}\\ \overline{\mathbf{b}^{(1)}} &= \overline{\mathbf{z}} \end{split}$$

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Computational Cost

 Computational cost of forward pass: one add-multiply operation per weight

$$z_i = \sum_j w_{ij}^{(1)} x_j + b_i^{(1)}$$

 Computational cost of backward pass: two add-multiply operations per weight

$$\overline{v_{ki}^{(2)}} = \overline{y_k} h_i$$
$$\overline{h_i} = \sum_k \overline{y_k} w_{ki}^{(2)}$$

 Rule of thumb: the backward pass is about as expensive as two forward passes.

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• For a multilayer perceptron, this means the cost is linear in the number of layers, quadratic in the number of units per layer.

Closing Thoughts

- Backprop is used to train the overwhelming majority of neural nets today.
 - Even optimization algorithms much fancier than gradient descent (e.g. second-order methods) use backprop to compute the gradients.
- Despite its practical success, backprop is believed to be neurally implausible.
 - No evidence for biological signals analogous to error derivatives.
 - All the biologically plausible alternatives we know about learn much more slowly (on computers).
 - So how on earth does the brain learn?

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The psychological profiling [of a programmer] is mostly the ability to shift levels of abstraction, from low level to high level. To see something in the small and to see something in the large.

– Don Knuth

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- By now, we've seen three different ways of looking at gradients:
 - Geometric: visualization of gradient in weight space
 - Algebraic: mechanics of computing the derivatives
 - Implementational: efficient implementation on the computer
- When thinking about neural nets, it's important to be able to shift between these different perspectives!