Representation Learning

Lecture 15 | Part 1

**Backpropagation** 

### **Gradient of a Network**

- We want to compute the gradient ∇<sub>w</sub>H.
   That is, ∂H/∂W<sup>(ℓ)</sup><sub>ii</sub> and ∂H/∂b<sup>(ℓ)</sup><sub>i</sub> for all valid i, j, ℓ.
- A network is a composition of functions.
- We'll make good use of the chain rule.

# **Recall: The Chain Rule**

$$\frac{d}{dx}f(g(x)) = \frac{df}{dg}\frac{dg}{dx}$$
$$= f'(g(x))g'(x)$$

#### **Some Notation**

- We'll consider an arbitrary node in layer l of a neural network.
- Let *g* be the activation function.
- $harpoint n_{\ell}$  denotes the number of nodes in layer  $\ell$ .

# **Arbitrary Node**



#### Claim #1





#### Claim #2





#### Claim #3





# **Exercise** What is $\partial H / \partial b_j^{(l)}$ ?



#### **General Formulas**

For any node in any neural network<sup>1</sup>, we have the following recursive formulas:

$$\frac{\partial H}{\partial a_j^{(\ell)}} = \sum_{k=1}^{n_{\ell+1}} \frac{\partial H}{\partial z_k^{(\ell+1)}} W_{jk}^{(\ell+1)}$$

$$\frac{\partial H}{\partial z_j^{(\ell)}} = \frac{\partial H}{\partial a_j^{(\ell)}} g'(z_j^{\ell})$$

$$\frac{\partial H}{\partial W_{ij}^{(\ell)}} = \frac{\partial H}{\partial z_j^{(\ell)}} a_i^{(\ell-1)}$$

$$\frac{\partial H}{\partial b_i^{(\ell)}} = \frac{\partial H}{\partial z_i^{(\ell)}}$$

<sup>1</sup>Fully-connected, feedforward network

#### Main Idea

The derivatives in layer  $\ell$  depend on derivatives in layer  $\ell + 1$ .

# Backpropagation

- Idea: compute the derivatives in last layers, first.
- That is:

►

- ► Compute derivatives in last layer, *l*; store them.
- Use to compute derivatives in layer l 1.
- ► Use to compute derivatives in layer ℓ 2.

•••

# Backpropagation

Given an input  $\vec{x}$  and a current parameter vector  $\vec{w}$ :

- 1. Evaluate the network to compute  $z_i^{(\ell)}$  and  $a_i^{(\ell)}$  for all nodes.
- 2. For each layer **?** from last to first:

Compute 
$$\frac{\partial H}{\partial a_j^{(\ell)}} = \sum_{k=1}^{n_{\ell+1}} \frac{\partial H}{\partial z_k^{(\ell+1)}} W_{jk}^{(\ell+1)}$$
Compute  $\frac{\partial H}{\partial z_j^{(\ell)}} = \frac{\partial H}{\partial a_j^{(\ell)}} g'(Z_j^{\ell})$ 
Compute  $\frac{\partial H}{\partial W_{ij}^{(\ell)}} = \frac{\partial H}{\partial z_j^{(\ell)}} a_i^{(\ell-1)}$ 
Compute  $\frac{\partial H}{\partial b_j^{(\ell)}} = \frac{\partial H}{\partial z_j^{(\ell)}}$ 

Compute the entries of the gradient given:

 $\frac{\partial H}{\partial a_j^{(\ell)}} = \sum_{k=1}^{n_{\ell+1}} \frac{\partial H}{\partial z_k^{(\ell+1)}} W_{jk}^{(\ell+1)} \qquad \frac{\partial H}{\partial z_j^{(\ell)}} = \frac{\partial H}{\partial a_j^{(\ell)}} g'(z_j^{\ell}) \qquad \frac{\partial H}{\partial W_{ij}^{(\ell)}} = \frac{\partial H}{\partial z_j^{(\ell)}} a_i^{(\ell-1)}$ 



# Summary: Backprop

- Backprop is an algorithm for efficiently computing the gradient of a neural network
- It is not an algorithm you need to carry out by hand: your NN library can do it for you.

Representation Learning

#### Lecture 15 | Part 2

#### **Gradient Descent for NN Training**

# **Empirical Risk Minimization**

- 0. Collect a training set,  $\{(\vec{x}^{(i)}, y_i)\}$
- Pick the form of the prediction function, H.
   E.g., a neural network, H.
- 2. Pick a loss function.
- 3. Minimize the empirical risk w.r.t. that loss.

# **Minimizing Risk**

- ► To minimize risk, we often use **vector calculus**.
  - Either set  $\nabla_{\vec{w}} R(\vec{w}) = 0$  and solve...
  - Or use gradient descent: walk in opposite direction of  $\nabla_{\vec{w}} R(\vec{w})$ .

► Recall,  $\nabla_{\vec{w}} R(\vec{w}) = (\partial R / \partial w_0, \partial R / \partial w_1, ..., \partial R / \partial w_d)^T$ 

#### In General

- Let  $\ell$  be the loss function, let  $H(\vec{x}; \vec{w})$  be the prediction function.
- ► The empirical risk:

$$R(\vec{w}) = \frac{1}{n} \sum_{i=1}^{n} \ell(H(\vec{x}^{(i)}; \vec{w}), y_i)$$

Using the chain rule:

$$\nabla_{\vec{w}} R(\vec{w}) = \frac{1}{n} \sum_{i=1}^{n} \frac{\partial \ell}{\partial H} \nabla_{\vec{w}} H(\vec{x}^{(i)}; \vec{w})$$

# **Training Neural Networks**

- For neural networks with nonlinear activations, the risk  $R(\vec{w})$  is typically **complicated**.
- The minimizer cannot be found directly.
- Instead, we use iterative methods, such as gradient descent.

# **Iterative Optimization**

To minimize a function  $f(\vec{x})$ , we may try to compute  $\vec{\nabla} f(\vec{x})$ ; set to 0; solve.

- Often, there is no closed-form solution.
- ► How do we minimize *f*?

• Consider 
$$f(x, y) = e^{x^2 + y^2} + (x - 2)^2 + (y - 3)^2$$
.



Try solving 
$$\vec{\nabla} f(x, y) = 0$$
.

► The gradient is:

$$\vec{\nabla}f(x,y) = \begin{pmatrix} 2xe^{x^2+y^2}+2(x-2)\\ 2ye^{x^2+y^2}+2(y-3) \end{pmatrix}$$

Can we solve the system?

$$2xe^{x^2+y^2} + 2(x-2) = 0$$
$$2ye^{x^2+y^2} + 2(y-3) = 0$$

Try solving 
$$\vec{\nabla} f(x, y) = 0$$
.

► The gradient is:

$$\vec{\nabla}f(x,y) = \begin{pmatrix} 2xe^{x^2+y^2}+2(x-2)\\ 2ye^{x^2+y^2}+2(y-3) \end{pmatrix}$$

Can we solve the system? Not in closed form.

$$2xe^{x^2+y^2} + 2(x-2) = 0$$
$$2ye^{x^2+y^2} + 2(y-3) = 0$$

# Idea

- Apply an iterative approach.
- Start at an arbitrary location.
- "Walk downhill", towards minimum.



# Which way is down?

- Consider a differentiable function f(x, y).
- We are standing at  $P = (x_0, y_0)$ .
- In a small region around P, f looks like a plane.
- Slope of plane in x, y directions:

 $\frac{\partial f}{\partial x}(x_0, y_0) = \frac{\partial f}{\partial y}(x_0, y_0)$ 



#### **The Gradient**

Let  $f : \mathbb{R}^d \to \mathbb{R}$  be differentiable. The gradient of f at  $\vec{x}$  is defined:

$$\vec{\nabla} f(\vec{x}) = \left(\frac{\partial f}{\partial x_1}(\vec{x}), \frac{\partial f}{\partial x_2}(\vec{x}), \dots, \frac{\partial f}{\partial x_d}(\vec{x})\right)^T$$

▶ Note:  $\nabla f(\vec{x})$  is a function mapping  $\mathbb{R}^d \to \mathbb{R}^d$ .

# Which way is down?

- ▶  $\nabla f(x_0, y_0)$  points in direction of steepest **ascent** at  $(x_0, y_0)$ .
- ►  $-\nabla f(x_0, y_0)$  points in direction of steepest **descent** at  $(x_0, y_0)$ .



# **Gradient Properties**

The gradient is used in the linear approximation of *f*:

$$f(x_0 + \delta_x, y_0 + \delta_y) \approx f(x_0, y_0) + \vec{\delta} \cdot \vec{\nabla} f(x_0, y_0)$$

- Important properties:
  - ▶  $\vec{\nabla} f(\vec{x})$  points in direction of **steepest ascent** at  $\vec{x}$ .
  - ►  $-\vec{\nabla}f(\vec{x})$  points in direction of **steepest descent** at  $\vec{x}$ .
  - ▶ In directions orthogonal to  $\nabla f(\vec{x})$ , f does not change!
  - ▶  $\|\nabla f(\vec{x})\|$  measures steepness of ascent

#### **Gradient Descent**

- Pick arbitrary starting point  $\vec{x}^{(0)}$ , learning rate parameter  $\eta > 0$ .
- Until convergence, repeat:
  - Compute gradient of f at  $\vec{x}^{(i)}$ ; that is, compute  $\vec{\nabla} f(\vec{x}^{(i)})$ .
  - Update  $\vec{x}^{(i+1)} = \vec{x}^{(i)} \eta \vec{\nabla} f(\vec{x}^{(i)})$ .
- When do we stop?
  - When difference between  $\vec{x}^{(i)}$  and  $\vec{x}^{(i+1)}$  is negligible.
  - ► I.e., when  $\|\vec{x}^{(i)} \vec{x}^{(i+1)}\|$  is small.

```
def gradient descent(
         gradient, x, learning rate=.01,
         threshold=.1e-4
):
    while True:
         x \text{ new} = x - \text{learning rate } * \text{gradient}(x)
         if np.linalg.norm(x - x new) < threshold:
             break
         x = x new
    return x
```



# **Backprop Revisited**

- The weights of a neural network can be trained using gradient descent.
- This requires the gradient to be calculated repeatedly; this is where **backprop** enters.
- Sometimes people use "backprop" to mean "backprop + SGD", but this is not strictly correct.

#### **Backprop Revisited**

Consider training a NN using the square loss:

$$\nabla_{\vec{w}} R(\vec{w}) = \frac{1}{n} \sum_{i=1}^{n} \frac{\partial \ell}{\partial H} \nabla_{\vec{w}} H(\vec{x}^{(i)}; \vec{w})$$
$$= \frac{2}{n} \sum_{i=1}^{n} (H(\vec{x}^{(i)}) - y_i) \nabla_{\vec{w}} H(\vec{x}^{(i)}; \vec{w})$$

#### **Backprop Revisited**

Interpretation:

$$\nabla_{\vec{w}} R(\vec{w}) = \frac{2}{n} \sum_{i=1}^{n} \underbrace{(H(\vec{x}^{(i)}) - y_i)}_{\text{Error}} \underbrace{\nabla_{\vec{w}} H(\vec{x}^{(i)}; \vec{w})}_{\text{Blame}}$$

When used in SGD, backprop "propagates error backward" in order to update weights.

# **Difficulty of Training NNs**

Gradient descent is guaranteed to find optimum when objective function is convex.<sup>2</sup>



<sup>2</sup>Assuming it is properly initialized

# **Difficulty of Training NNs**

When activations are non-linear, neural network risk is highly non-convex:



# **Non-Convexity**

- When R is non-convex, GD can get "stuck" in local minima.
  - Solution depends on initialization.
- More sophisticated optimizers, using momentum, adaptation, better initialization, etc.
   Adagrad, RMSprop, Adam, etc.

# Difficulty of Training (Deep) NNs

Deep networks can suffer from the problem of vanishing gradients: if w is a weight at the "front" of the network, ∂H/∂w can be very small



# **Vanishing Gradients**

If ∂H/∂w is always close to zero, w is updated very slowly by gradient descent.

- In short: early layers are slower to train.
- One mitigation: use ReLU instead of sigmoid.

# **Vanishing Gradients**



Representation Learning

Lecture 15 | Part 3

**Stochastic Gradient Descent** 

# **Gradient Descent for Minimizing Risk**

In ML, we often want to minimize a risk function:

$$R(\vec{w}) = \frac{1}{n} \sum_{i=1}^{n} \ell(H(\vec{x}^{(i)}; \vec{w}), y_i)$$

#### Observation

The gradient of the risk function is a sum of gradients:

$$\vec{\nabla} R(\vec{w}) = \frac{1}{n} \sum_{i=1}^{n} \vec{\nabla} \ell(H(\vec{x}^{(i)}; \vec{w}), y_i)$$

One term for each point in training data.

#### **Problem**

- In machine learning, the number of training points n can be very large.
- Computing the gradient can be expensive when n is large.
- Therefore, each step of gradient descent can be expensive.

#### Idea

The (full) gradient of the risk uses all of the training data:

$$\nabla R(\vec{w}) = \frac{1}{n} \sum_{i=1}^{n} \nabla \ell(H(\vec{x}^{(i)}; \vec{w}), y_i)$$

- It is an average of n gradients.
- Idea: instead of using all n points, randomly choose << n.</p>

#### **Stochastic Gradient**

Choose a random subset (mini-batch) B of the training data.

Compute a stochastic gradient:

$$\nabla R(\vec{w}) \approx \sum_{i \in B} \vec{\nabla} \ell(H(\vec{x}^{(i)}; \vec{w}), y_i)$$

#### **Stochastic Gradient**

$$\nabla R(\vec{w}) \approx \sum_{i \in B} \vec{\nabla} \ell(H(\vec{x}^{(i)}; \vec{w}), y_i)$$

- **Good:** if  $|B| \ll n$ , this is much faster to compute.
- Bad: it is a (random) approximation of the full gradient, noisy.

#### Stochastic Gradient Descent (SGD) for ERM

- Pick arbitrary starting point  $\vec{x}^{(0)}$ , learning rate parameter  $\eta > 0$ , batch size  $m \ll n$ .
- Until convergence, repeat:
  - Randomly sample a batch B of m training data points (on each iteration).
  - Compute stochastic gradient of f at  $\vec{x}^{(i)}$ :

$$\vec{g} = \sum_{i \in B} \vec{\nabla} \ell(H(\vec{x}^{(i)}; \vec{w}), y_i)$$

• Update 
$$\vec{x}^{(i+1)} = \vec{x}^{(i)} - \eta \vec{g}$$

#### Idea

- In practice, a stochastic gradient often works well enough.
- It is better to take many noisy steps quickly than few exact steps slowly.

### **Batch Size**

- Batch size *m* is a parameter of the algorithm.
- The larger *m*, the more reliable the stochastic gradient, but the more time it takes to compute.
- Extreme case when m = 1 will still work.



# **Usefulness of SGD**

- SGD allows learning on **massive** data sets.
- Useful even when exact solutions available.
   E.g., least squares regression / classification.

# **Training NNs in Practice**

- There are several Python packages for training NNs:
  - PyTorch
  - Tensorflow / Keras
- This week's discussion was a Tensorflow tutorial.