DSC 140B Representation Learning

Lecture 22 | Part 1

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 ℓ 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*?

Example



Example

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?

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

Example

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

Which way is down?

- ▶ $\vec{\nabla} f(x_0, y_0)$ points in direction of steepest **ascent** at (x_0, y_0) .
- ► $-\vec{\nabla}f(x_0, y_0)$ 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:
 - ▶ $\nabla f(\vec{x})$ 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 $\vec{\nabla} f(\vec{x})$, *f* does not change!
 - ▶ $\|\vec{\nabla}f(\vec{x})\|$ 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 new = x - learning rate * 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.¹



¹Assuming 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



 $= \frac{2H}{22^{(0)}} \cdot a^{(0)} = \frac{2}{2a^{(0)}} \cdot g^{(2^{(0)})} \cdot a^{(0)}$ $= \frac{2}{2a^{(0)}} \cdot g^{(2^{(0)})} \cdot a^{(0)}$ $= \frac{2}{2a^{(0)}} \cdot g^{(2^{(0)})} \cdot a^{(0)}$ 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. er plodving goudient 2 20 (1000) gilleting)

Vanishing Gradients



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Lecture 22 Part 2 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>

102 **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 -

nxd)

SGD allows learning on massive data sets.

Useful even when exact solutions available.
E.g., least squares regression / classification.

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Lecture 22 | Part 3

Output Units

Output Units

As with units in hidden layers, can choose different activation functions for the outputs layey.

What activation function?

How many units?

Good choice depends on task:

Regression, binary classification, multiclass, etc.

10.2,

Which loss?

Setting 1: Regression

- Output can be any real number.
- Single output neuron.
- It makes sense to use a linear activation.



Setting 1: Regression

Prediction should not be too high/low.

It makes sense to use the mean squared error.

Setting 1: Regression

- Suppose we use linear activation for output neuron + mean squared error.
- This is very similar to least squares regression...
- But! Features in earlier layers are learned, non-linear.



Special Case: Least Squares



The case of:

- a one layer neural network
- with all linear activations
- trained with square loss

is also called **least squares regression**.

Setting 2: Binary Classification

- Output can be in [0, 1].
- Single output neuron.
- We could use a linear activation, threshold.
- But there is a better way.



Sigmoids for Classification

Natural choice for activation in output layer for binary classification: the sigmoid.



Binary Classification Loss

- We could use square loss for binary classification. There are several reasons not to:
- 1) Square loss penalizes predictions which are "too correct".
- 2) It doesn't work well with the sigmoid due to saturation.

The Cross-Entropy

Instead, we often train deep classifiers using the cross-entropy as loss.

 $\log [1 - f(\vec{x}^{(i)})], \text{ if } y^{(i)}$

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 \mathcal{P} Let $y^{(i)} \in \{0, 1\}$ be true label of *i*th example.

The average cross-entropy loss:

The Cross-Entropy and the Sigmoid

Cross-entropy "undoes" the exponential in the sigmoid, resulting in less saturation.



Summary: Binary Classification

Use sigmoidal activation the output layer + cross-entropy loss.

► This will promote a strong gradient.

Use whatever activation for the hidden layers (e.g., ReLU).

Special Case: Logisitic Regression X₁ The case of: a one layer neural network Wo X₂ with sigmoid activation trained with cross-entropy loss is also called logistic regression. binan Els Wd Хd

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Lecture 22 | Part 4

Convolutions





From Simple to Complex

- Complex shapes are made of simple patterns
- The human visual system uses this fact
- ► Line detector → shape detector → ... → face detector
- Can we replicate this with a deep NN?

Edge Detector

- How do we find vertical edges in an image?
- One solution: convolution with an edge filter.



Vertical Edge Filter



- Take a patch of the image, same size as filter.
- Perform "dot product" between patch and filter.
- If large, this is a (vertical) edge.

image patch:















Convolution

- The result is the (2d) convolution of the filter with the image.
- Output is also 2-dimensional array.
- Called a response map.

Example: Vertical Filter





Example: Horizontal Filter







More About Filters

► Typically 3×3 or 5×5.

Variations: different stride, image padding.

Black and white images are 2-d arrays.

But color images are 3-d arrays:

- a.k.a., tensors
- ► Three color **channels**: red, green, blue.
- height × width × 3

How does convolution work here?

Color Image



The filter must also have three channels:
3 × 3 × 3, 5 × 5 × 3, etc.










Convolution with 3-d Filter

- Filter must have same number of channels as image.
 - ▶ 3 channels if image RGB.
- Result is still a 2-d array.

General Case

- Input "image" has k channels.
- Filter must have k channels as well.
 - ▶ e.g., 3 × 3 × k
- Output is still 2 d

