ECE 1508: Applied Deep Learning

Chapter 1: Preliminaries

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Function Optimization

The model training always reduces to an optimization problem

$$\mathbf{w}^{\star} = \underset{\mathbf{w}}{\operatorname{argmin}} \, \hat{R}\left(\mathbf{w}\right) = \underset{\mathbf{w}}{\operatorname{argmin}} \, \frac{1}{I} \sum_{i=1}^{I} \mathcal{L}\left(f_{\mathbf{h}}\left(x_{i} \middle| \mathbf{w}\right), \underline{y_{i}}\right) \tag{Training}$$

Let's recall each component of this optimization problem

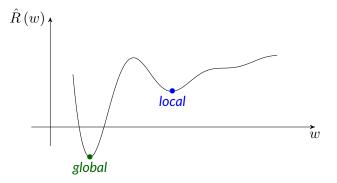
- $f_{\mathbf{h}}\left(\cdot|\mathbf{w}\right)$ model with hyperparameters \mathbf{h} and learnable parameters \mathbf{w}
 - in DL, it is the input-output relation of a neural network whose architecture is specified by **h** and whose weights and biases are collected in **w**
- x_i is a data-point with label y_i , and I is size of dataset
- L is the loss function

No matter what we choose, at the end of the day we need to solve

$$\min_{\mathbf{w}} \hat{R}\left(\mathbf{w}\right)$$

Function Optimization

In general, the empirical risk $\hat{R}(\mathbf{w})$ can have local and global minima Let's take a look at a simple visual case with only one parameter, i.e., $w \in \mathbb{R}$



We are happy if we get the global; but, many times getting to a local is enough!

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Function Optimization

- + Why is it a big problem? We could grid w and search for the grid with smallest empirical risk. We then find it with a good accuracy!
- For only one parameter yes! But, we have seen deep neural networks. They have too many neurons, and hence too many parameters!

Say for an accurate approximation with only one parameter, we need G grids

If we have D parameters, i.e., $\mathbf{w} \in \mathbb{R}^D$, we need

 ${\cal G}^D$ grids

to get an approximation with the same accuracy!

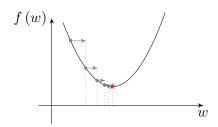
For practical neural networks with $D=10^5$, this is impossible!

we need to have an optimization algorithm with feasible complexity

Optimization Algorithms \equiv Optimizer

We look for an optimization algorithm, or as ML people call it "an optimizer"

- it starts from an initial point and moves for some steps
- in each step, it moves towards where the empirical risk is minimized
- it moves for a feasible number of steps



Optimization Algorithms \equiv Optimizer

Let's clear things up: we are looking for an iterative approach as below

```
1: Initiate at some \mathbf{w}^{(0)} \in \mathbb{R}^D and deviation \Delta = +\infty

2: Choose some small \epsilon

3: \mathbf{while} \ \Delta > \epsilon \ \mathbf{do}

4: Determine a vector \boldsymbol{\mu}^{(t)} \in \mathbb{R}^D based on \hat{R}(\mathbf{w}) \leftarrow we need to figure out

5: Update weights as \mathbf{w}^{(t)} \leftarrow \mathbf{w}^{(t-1)} + \boldsymbol{\mu}^{(t)}

6: Update the deviation \Delta = |\hat{R}(\mathbf{w}^{(t)}) - \hat{R}(\mathbf{w}^{(t-1)})|

7: \mathbf{end} \ \mathbf{while}
```

We would like to have following properties

- → most of the time the empirical risk reduces in line 5

Optimization Algorithms

```
1: Initiate at some \mathbf{w}^{(0)} \in \mathbb{R}^D and deviation \Delta = +\infty
2: Choose some small \epsilon, and set t=1
3: while \Delta > \epsilon do
```

- 3: While $\Delta > \epsilon$ do
- 4: Determine a vector $\boldsymbol{\mu}^{(t)} \in \mathbb{R}^D$ based on $\hat{R}(\mathbf{w}) \leftarrow$ we need to figure out
- 5: Update weights as $\mathbf{w}^{(t)} \leftarrow \mathbf{w}^{(t-1)} + \boldsymbol{\mu}^{(t)}$
- 6: Update the deviation $\Delta = |\hat{R}(\mathbf{w}^{(t)}) \hat{R}(\mathbf{w}^{(t-1)})|$
- 7: end while

We are going to get what we want, if we set

$$\mu^{(t)}$$
 to be proportional to the negative of gradient at $\mathbf{w}^{(t-1)}$

This is what we call the gradient descent algorithm. But, before we start with this algorithm, let's recap some basic notions of calculus!

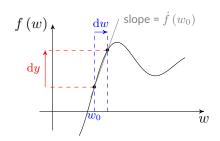
Derivative of one-dimensional function f(w) at point $w = w_0$ is defined as

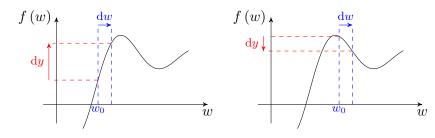
$$\dot{f}(w_0) = \frac{d}{dw} f(w_0) = f'(w_0) = \lim_{\delta \to 0} \frac{f(w_0 + \delta) - f(w_0)}{\delta}$$

This definition is intuitively interpreted as follows:

Let y = f(w). If we vary w around w_0 with a tiny step dw; then,

Variation of
$$y = dy = \dot{f}(w_0) dw$$



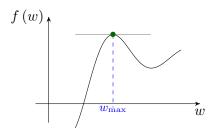


The derivative represents the slope of function

- $\dot{f}(w_0) > 0$ means increasing w will increase y = f(w)
- $\dot{f}\left(w_{0}\right)<0$ means increasing w will decrease $y=f\left(w\right)$

So, we could also say: the derivative shows the moving direction on w-axis towards which the function increases; or alternatively, its negative is the direction that function decreases

When do we have the derivative equal to zero? Either we are at a maximum



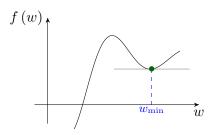
Starting before the maximum,

- The derivative is first positive and gradually reduces to zero
- As we pass the maximum the derivative gets more and more negative

So around the maximum as we increase w, the derivative reduces

$$\ddot{f}(w_{\text{max}}) = \frac{d^2}{dw^2} f(w_{\text{max}}) = f''(w_{\text{max}}) < 0$$

When do we have the derivative equal to zero? Either we are at a minimum



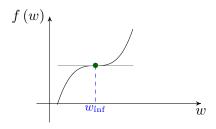
Starting before the minimum,

- The derivative is first negative and gradually increases to zero
- As we pass the minimum the derivative gets more and more positive

So around the maximum as we increase w, the derivative reduces

$$\ddot{f}(w_{\min}) = \frac{d^2}{dw^2} f(w_{\min}) = f''(w_{\min}) > 0$$

When do we have the derivative equal to zero? Either we are at an inflection



Starting before the inflection point,

- The derivative is first positive and gradually decreases to zero
- As we pass the inflection point the derivative gets again positive

So around the inflection point, the second derivative changes sign

$$\ddot{f}(w_{\text{inf}}) = \frac{d^2}{dw^2} f(w_{\text{inf}}) = f''(w_{\text{inf}}) = 0$$

- + What about multi-variable functions, e.g., $f(\mathbf{w})$ for $\mathbf{w} = [w_1, \dots, w_N]$?
- We can take derivative with respect to each variable, i.e.,

$$\dot{f}_n\left(\mathbf{w}_0\right) = \frac{\partial}{\partial w_n} f\left(\mathbf{w}_0\right)$$

This is what we call partial derivative

Partial derivative n represents the same thing: slope in direction of w_n

Let $y = f(\mathbf{w})$. If we vary \mathbf{w} around \mathbf{w}_0 in N-dimensional space with

$$\mathbf{d}\mathbf{w} = [\mathbf{d}w_1, \dots, \mathbf{d}w_N]$$

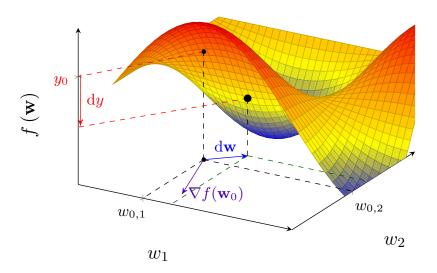
whose entries are very tiny; then, the variation of y is

$$dy = \dot{f}_1(\mathbf{w}_0) dw_1 + \ldots + \dot{f}_N(\mathbf{w}_0) dw_N = \sum_{n=1}^N \dot{f}_n(\mathbf{w}_0) dw_n$$

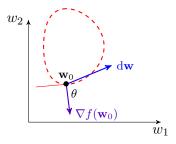
We can use inner-product to represent dy compactly

$$dy = \sum_{n=1}^{N} \dot{f}_n(\mathbf{w}_0) dw_n = \underbrace{\left[\dot{f}_1(\mathbf{w}_0) \dots \dot{f}_N(\mathbf{w}_0)\right]}_{\nabla f(\mathbf{w}_0)^{\mathsf{T}}} \begin{bmatrix} dw_1 \\ \vdots \\ dw_N \end{bmatrix}$$
$$= \nabla f(\mathbf{w}_0)^{\mathsf{T}} d\mathbf{w}$$

We call $\nabla f(\mathbf{w}_0)$ the gradient of $f(\cdot)$ at $\mathbf{w} = \mathbf{w}_0$



Let's get to the w-plane: the gradient is perpendicular to the contour level



The variation of y is the inner product of these two vectors

$$\mathbf{d}y = \nabla f(\mathbf{w}_0)^{\mathsf{T}} \mathbf{d}\mathbf{w} = \|\nabla f(\mathbf{w}_0)\| \|\mathbf{d}\mathbf{w}\| \cos(\theta)$$

where $\|\cdot\|$ is the Euclidean norm, i.e., $\|\mathbf{w}\| = \sqrt{w_1^2 + w_2^2}$

Say we move with a tiny step of fixed size: so we have

$$\|\mathbf{d}\mathbf{w}\| = \epsilon$$

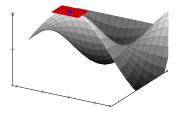
for some small ϵ

- + How can we move, such that y maximally increases?
- Well, we need $\theta = 0$ meaning that we should move in the direction of gradient

Alternatively, the function decreases maximally if $\theta = \pi$ or

we move in the direction of negative gradient

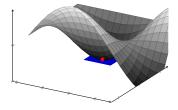
When do we have zero gradient? Either when we are at a maximum



We can again relate it to the second order derivatives of the function

at maximum Hessian matrix is negative definite

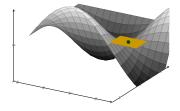
When do we have zero gradient? or when we are at a minimum



We can again relate it to the second order derivatives of the function

at minimum Hessian matrix is positive definite

When do we have zero gradient? or when we are at a saddle point



We can again relate it to the second order derivatives of the function at saddle point Hessian matrix is neither negative nor positive definite

Just as a reminder: Hessian is the matrix of all second order derivatives

$$\nabla^{2} f\left(\mathbf{w}_{0}\right) = \begin{bmatrix} \frac{\partial^{2}}{\partial w_{1}^{2}} f\left(\mathbf{w}_{0}\right) & \frac{\partial^{2}}{\partial w_{1} \partial w_{2}} f\left(\mathbf{w}_{0}\right) & \dots & \frac{\partial^{2}}{\partial w_{1} \partial w_{N}} f\left(\mathbf{w}_{0}\right) \\ \vdots & & & \vdots \\ \frac{\partial^{2}}{\partial w_{N} \partial w_{1}} f\left(\mathbf{w}_{0}\right) & \frac{\partial^{2}}{\partial w_{N} \partial w_{2}} f\left(\mathbf{w}_{0}\right) & \dots & \frac{\partial^{2}}{\partial w_{N}^{2}} f\left(\mathbf{w}_{0}\right) \end{bmatrix}$$

We never use the Hessian matrix in this course

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Moral of Story: Gradient Decent

- + What is the whole motive of this discussions?
- Simple: at any point \mathbf{w}_0 , if we want to move in a direction that the function reduces, the best direction is negative of gradient at \mathbf{w}_0

So, we can complete our optimization algorithm as follows:

```
1: Initiate at some \mathbf{w}^{(0)} \in \mathbb{R}^D and deviation \Delta = +\infty

2: Choose some small \epsilon and \eta, and set t = 1

3: while \Delta > \epsilon do

4: Update weights as \mathbf{w}^{(t)} \leftarrow \mathbf{w}^{(t-1)} - \eta \nabla \hat{R}(\mathbf{w}^{(t-1)})

5: Update the deviation \Delta = |\hat{R}(\mathbf{w}^{(t)}) - \hat{R}(\mathbf{w}^{(t-1)})|

6: end while
```

The scalar η is the step-size we take in each iteration:

we usually call it learning rate

Behavior of Gradient Decent

- + Can we always use gradient descent?
- Pretty much Yes! The problem starts only when the empirical loss is not differentiable

How to handle this problem?

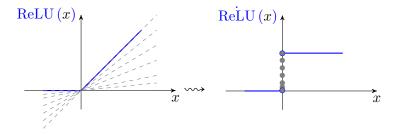
There are two sources for being non-differentiable

- 1 a function that is continuous but not differentiable
- 2 a discontinuous function

Let's look at each case separately

Behavior of Gradient Decent: Non-differentiable Elements

An example of a non-differentiable continuous function is ReLU

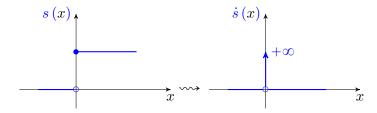


In this case, we define a sub-gradient and use it instead of gradient

take the slope of a line that lies below the curve at the point

Behavior of Gradient Decent: Discontinuous Elements

An example of a discontinuous function is the step function



The gradient is somehow infinite! We can only rely on the sign of variation

we always avoid such elements in our model and loss

Bingo! You may recall that we discouraged the choice of activation and loss function in the perceptron machine

- + Now, let's assume that we've handled differentiability. Does gradient decent always end up at the minimum point?
- This brings up the concept of *convergence*

Let's look at the algorithm again

```
1: Initiate at some \mathbf{w}^{(0)} \in \mathbb{R}^D and deviation \Delta = +\infty

2: Choose some small \epsilon and \eta, and set t=1

3: while \Delta > \epsilon do

4: Update weights as \mathbf{w}^{(t)} \leftarrow \mathbf{w}^{(t-1)} - \eta \nabla \hat{R}(\mathbf{w}^{(t-1)})

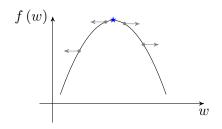
5: Update the deviation \Delta = |\hat{R}(\mathbf{w}^{(t)}) - \hat{R}(\mathbf{w}^{(t-1)})|

6: end while
```

Intuitively, if we set ϵ very small: the algorithm stops when the gradient is close to zero, i.e., when we are at a maximum, minimum or an inflection/saddle-point

Let's see how the algorithm behaves when we get close to such point

When we are around a maximum



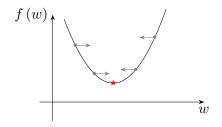
If we are exactly at a maximum; then, the algorithm stops. But, in reality

we land somewhere around it

at such points, the algorithm always pushes us outwards

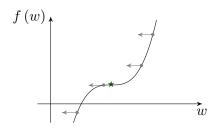
Gradient descent practically does not get into a maximum

When we are around a minimum



Around minima, the algorithm always pushes us towards the minimum

When we are around an inflection point



If we are exactly at an inflection; then, the algorithm stops. But, in reality

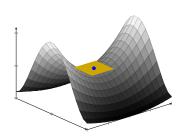
we land somewhere around it

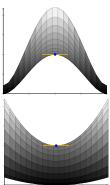
at such points, the algorithm always pushes us somewhere else

Gradient descent practically does not get into an inflection

- + Can we extend this conclusion to saddle-points?
- Yes, but with a bit of caution!

At saddle points, function is maximized in a direction and minimized in another





So, for a saddle-point we can conclude: if we are exactly at a saddle-point; then, the algorithm stops. But, in reality

we land somewhere around it

If at that point, the gradient has a component in the direction that the function is maximized; then, the algorithm pushes us outwards.

- + Can it happen that we do not land at such point?
- Thinking with an engineer's mind: Not really!

So we could say

Gradient descent almost never gets trapped at a saddle-point

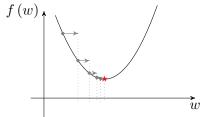
Moral of Story

Gradient descent almost never gets trapped at a point that is not minimum

- + Nice! But, does it always converge?
- Well! If we choose the *learning rate* properly; then, Yes!

With small learning rates, the algorithm converges; how small? η <

f(w)

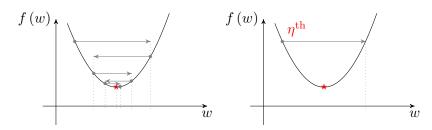


Moral of Story

Gradient descent almost never gets trapped at a point that is not minimum

- + Nice! But, does it always converge?
- Well! If we choose the learning rate properly; then, Yes!

With larger learning rates, the algorithm starts oscillating: $|\eta^{\star} < \eta < \eta^{ ext{th}}|$

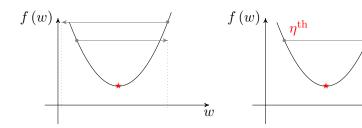


Moral of Story

Gradient descent almost never gets trapped at a point that is not minimum

- + Nice! But, does it always converge?
- Well! If we choose the *learning rate* properly; then, Yes!

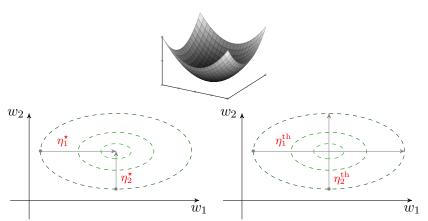
With extremely large *learning rates*, the algorithm *diverges*: $|\eta>\eta^{
m th}$



11)

We can extend this idea to the multi-dimensional functions; however,

 η^* and η^{th} are different on each axis



One may suggest that we use a vector of learning rates, i.e.,

$$w_n^{(t)} \leftarrow w_n^{(t-1)} - \frac{\partial}{\partial w_n} \hat{R}(\mathbf{w}^{(t-1)})$$

for each n = 1, ..., N. This is however not easy; the easier way is to focus on

$$\eta^{\star} = \min_{n} \, \eta_{n}^{\star} \quad \text{and} \quad \eta^{\text{th}} = \min_{n} \, \eta_{n}^{\text{th}}$$

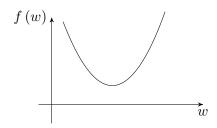
Clearly, there is always a trade-off

- We can choose a large learning rate
 - + Gradient descent converges faster: high convergence speed
 - The chance of divergence however increases: high divergence rate
- We can choose a small learning rate
 - Gradient descent converges slowly: low convergence speed
 - + The chance of divergence is now very low: low divergence rate
- + Well, say we are patient! Then, choosing a small learning rate is safe! Right?!
- Well! If the empirical risk is convex; then, Yes! But, with non-convex risks not always!

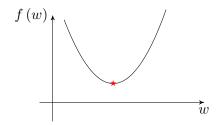
Recap: Convex Function

You really don't need to know the definition of a convex function for this course; however, just in case you're interested, here it goes:

$$f\left(\cdot\right):\mathbb{R}^{N}\mapsto\mathbb{R}$$
 is convex, if for any two points $x_{1},x_{2}\in\mathbb{R}^{N}$, we have
$$f(\lambda x_{1}+\left(1-\lambda\right)x_{2})\leqslant\lambda f(x_{1})+\left(1-\lambda\right)f(x_{2})$$
 for all $0\leqslant\lambda\leqslant1$



In convex functions, we don't have disjoint local minima



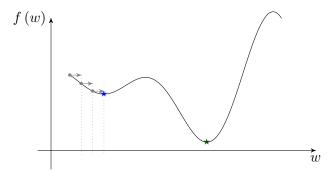
So, if we choose a small learning rate

we surely converge to the global minimum

But, most empirical losses in deep learning are non-convex:

we have multiple disjoint local minima

Gradient descent converges to one of them, but not necessarily the global one

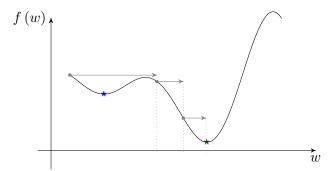


Too small learning rate can leave us in a bad local minimum!

But, most empirical losses in deep learning are non-convex:

we have multiple disjoint local minima

Gradient descent converges to one of them, but not necessarily the global one



Initial larger learning rates can take us out of bad local minima!

We can hence conclude a more general trade-off

- We can choose a large learning rate
 - + Gradient descent converges faster: high convergence speed
 - + Gradient descent may fall out of a local minimum: lower risk
 - The chance of divergence however increases: high divergence rate
- We can choose a small learning rate
 - Gradient descent converges slowly: low convergence speed
 - Gradient descent traps in a local minimum: high risk
 - + The chance of divergence is now very low: low divergence rate
- + How do we do it in practice?
- In practice, we start with large learning rates and reduce it gradually as we get close to the minimum

We will talk about this more once we start training practical neural networks!

Gradient Decent: Summary

```
1: Initiate at some \mathbf{w}^{(0)} \in \mathbb{R}^D and deviation \Delta = +\infty
```

- 2: Choose some small ϵ and η , and set t=1
- 3: while $\Delta > \epsilon$ do
- Update weights as $\mathbf{w}^{(t)} \leftarrow \mathbf{w}^{(t-1)} \eta \nabla \hat{R}(\mathbf{w}^{(t-1)})$
- Update the deviation $\Delta = |\hat{R}(\mathbf{w}^{(t)}) \hat{R}(\mathbf{w}^{(t-1)})|$
- 6: end while

Gradient descent converges almost always to a local minimum

- With convex empirical risks, this is global minimum
- With non-convex empirical risks, this is not necessarily global minimum
- Learning rate is a crucial parameter that tunes the convergence

There are other optimization algorithms that work based on gradient: we will talk about them later!