Intro to Gradient Descent

Introduction

- We have a loss function associated with how well our function does
 - One such example is MSE: $J(x) = \sum_{i=1}^{m} (f(x_i) y_i)^2$
- How do we minimize the loss?
 - Before, we were able to numerically solve.
 - Can you do this with everything?
 - How can we reliably minimize our loss function?

Introduction

Observation:

- If our function is differentiable, we can find the direction that decreases our function the fastest at any given point
- If we do this over and over again, we can approach the global minimum

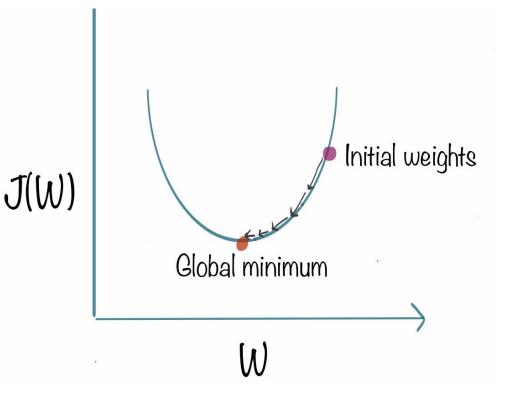
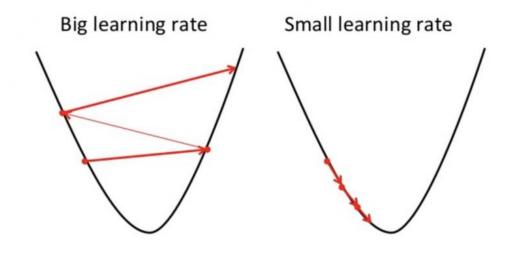


Image credit to KDNuggets

Gradient Descent Rule

$$\overrightarrow{w}_{t+1} = \overrightarrow{w}_t - \alpha \nabla J$$

- w is our parameters. In a linear classifier, it is the coefficients.
- With each iteration, J decreases.
- α is called the learning rate, which dictates how much the algorithm moves at each iteration



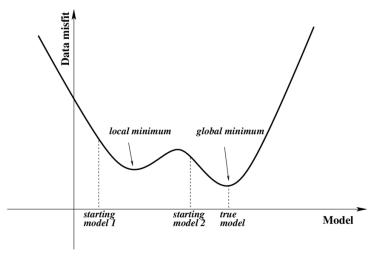
Credit to Towards Data Science for the image

Problems

- Doesn't this only find the local minimum? Won't you get stuck?
 - Yes and yes. This is why people like to use convex loss functions, so the local minimum is the global minimum
 - There are also saddle points, where the gradient is 0, but the point is not a local minimum

• E.g:
$$y = x_1^2 - x_2^2$$

 Techniques to escape from these are outside the scope of the course



Variations

- Batch gradient descent
- Stochastic gradient descent

Stochastic gradient descent

- Classical gradient descent works by generating a gradient for each data point in the dataset, then applying the gradients all at once for each iteration through the dataset
- Stochastic gradient descent instead works by applying the gradient after each individual point, so that the calculation of the loss function uses the updated parameters for the next data point
- In practice, most people use stochastic gradient descent, or a variation called mini-batching, as it seems to converge faster.

Momentum

• See slide 48 of Cottrell slides