Optimization Algorithms

- Optimization algorithms train neural networks faster by adjusting weights and learning rate
- Training deep neural networks on large dataset is slow
- Fast optimization algorithms can speed up training process

Gradient Descent

Gradient descent adjust the parameters iteratively to

minimize cost function

$$w_{new} = w - \alpha \frac{\partial J}{\partial w}$$

$$b_{new} = b - \alpha \frac{\partial J}{\partial b}$$

$$\theta = \begin{bmatrix} w \\ b \end{bmatrix}$$

$$\theta = \theta - \alpha \cdot \nabla_{\theta} J(\theta) \qquad \nabla_{\theta} J(\theta) = \begin{bmatrix} \frac{\partial J}{\partial w} \\ \frac{\partial J}{\partial b} \end{bmatrix}$$



https://ruder.io/optimizing-gradient-descent/

Gradient Descent

If we have dataset with 50,00,000 training examples

$$X = \begin{bmatrix} x^1 \\ x^2 \\ \vdots \\ x^m \end{bmatrix}$$

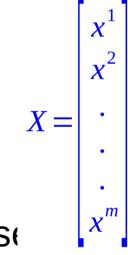
- 1. Will take more time for the convergence
- 2. Large memory to caliculate the gradient
- 3. May trap at local minimum

Stochastic Gradient Descent

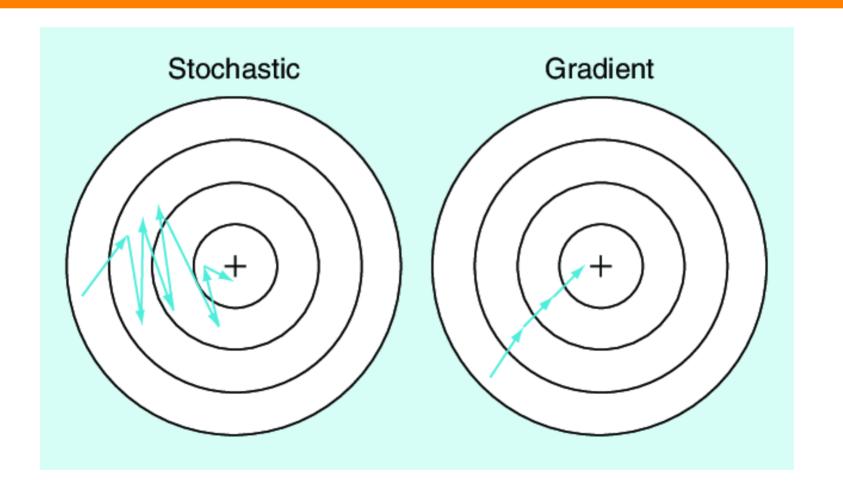
Caliculate gradient based on single training example

$$\theta = \theta - \alpha \cdot \nabla_{\theta} J(\theta; x^{(i)}; y^{(i)})$$

- 1. Less memory to caliculate the gradient
- 2. Faster than Batch Gradient Descent but it lose speed up from vectorization
- 3. May trap at local minimum



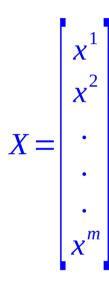
SGD vs GD

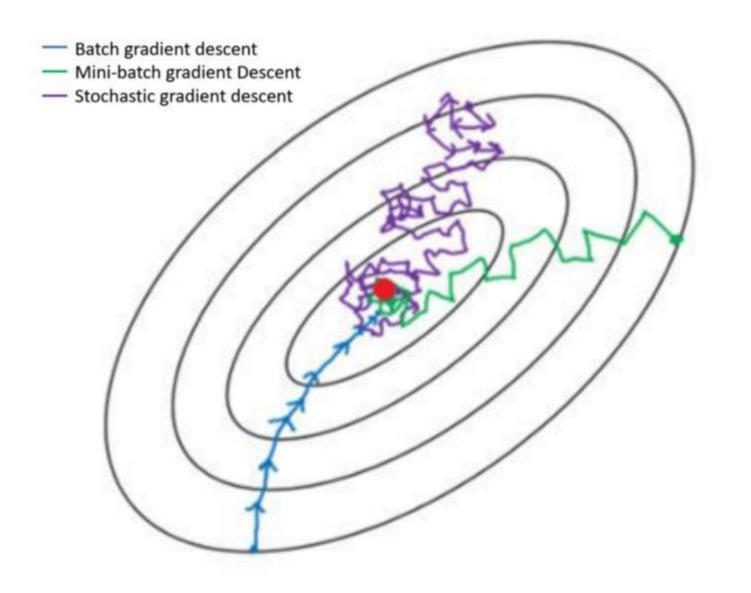


Mini-Batch Gradient Descent

Caliculate gradient based on subset of training example

$$\theta = \theta - \alpha \cdot \nabla_{\theta} J(\theta; x^{(i:i+n)}; y^{(i:i+n)})$$





Mini-Batch Gradient Descent

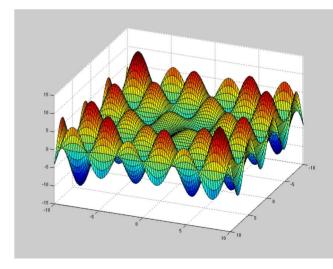
- Pros
 - Takes less noisy steps than SGD
 - Can utilize vectorization speedup
 - Takes less memory than GD
- Cons
 - Takes noisy steps compared to GD
 - Takes longer time compared to full batch GD
 - May stuck at local minimum

Mini-Batch Gradient Descent

Optimizer	Update Speed	Memory Usage
Batch Gradient Descent	Slow	High
Stochastic Gradient Descent	High	Low
Batch Gradient Descent	Medium	Medium

Local Minima

- With non-convex functions, such as neural nets, it is possible to have many local minima.
- The gradients may change dramatically on complex loss functions
- Local minima can be problematic if they have high cost in comparison to the global minimum.



https://ruder.io/optimizing-gradient-descent/

 Plateaus, Saddle Points and Other Flat Regions

 These saddle points are usually surrounded by a plateau of the same error

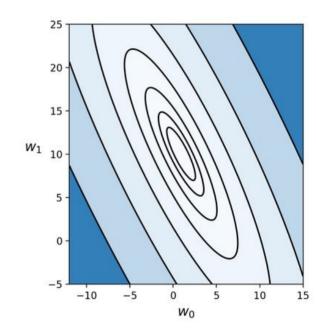
 Plateau is the region where the surface is almost flat and the gradient is close to zero.

 which makes it notoriously hard for GD to escape, as the gradient is close to zero in all dimensions.

Choosing a proper learning rate can be difficult.



- Same learning rate applies to all parameter updates.
- If our data is sparse and our features have very different frequencies, we might not want to update all of them to the same extent.



- Vanilla GD it works based on only Gradients
 - it can't be fast

$$w_{new} = w - \alpha \frac{\partial J}{\partial w}$$

 can't escape from local minima, saddle points and flat regions

 A good algorithm finds the minimum fast and reliably well (i.e. it doesn't get stuck in local minima, saddle points, or plateau regions), but rather goes for the global minimum

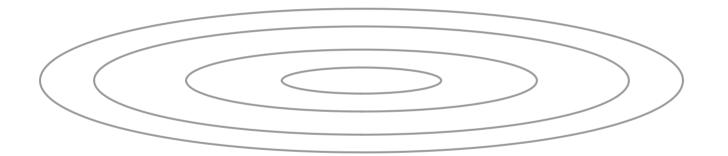
Gradient Descent Optimization Algorithms

- Gradient descent variants
 - Batch gradient descent
 - Stochastic gradient descent
 - Mini-batch gradient descent

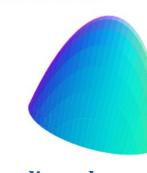
Gradient descent optimization algorithms

- Momentum
- Adagrad
- RMSprop
- Adam

 the surface curves much more steeply in one dimension than in another

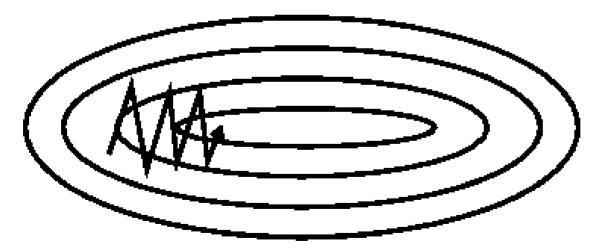


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A contour plot is a graphical technique for representing a 3-dimensional surface by plotting constant z slices, called contours, on a 2-dimensional format.

 In these scenarios, GD oscillates across the slopes of the ravine while only making hesitant progress along the bottom towards the local optimum as shown in image



- Momentum is a method that helps accelerate GD in the relevant direction and dampens oscillations.
- Adding a fraction y of the update vector of the past time step to the current update vector
- The momentum term γ is usually set to 0.9 or a similar value.

$$v_t = \gamma v_{t-1} + \alpha \cdot \nabla_{\theta} J(\theta)$$

$$\theta = \theta - v_t$$



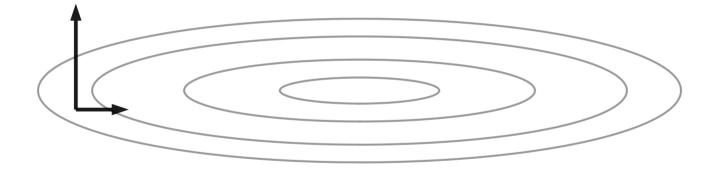
- The momentum term increases for dimensions whose gradients point in the same directions and reduces updates for dimensions whose gradients change directions.
- As a result, we gain faster convergence and reduced oscillation.
- When we push a ball down a hill. The ball accumulates momentum as it rolls downhill, becoming faster and faster on the way. The same thing happens to our parameter updates:

Problem with Momentum

- Learning rate should set manually.
- Same learning rate will be used for all parameters(in all dimentions).
- We may require small learning rate in some dimention and large learning rate in other dimention

Motivation for Adagrad

Consider two dimentional optimization problem



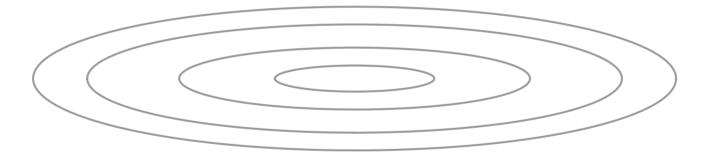
Motivation for Adagrad

Consider two dimentional optimization problem



Adagrad

• Adagrad uses a different learning rate for every parameter θ_i at every time step t



- It adapts lower learning rate for the parameters having larger gradients
- high learning rates for parameters having smaller gradients.

Adagrad

 At time t, it maintain the squared sum of past gradients.

$$S_t = S_{t-1} + g_t^2$$

Update rule

$$\theta_{t+1} = \theta_t - \frac{\alpha_t}{\sqrt{s_t + \epsilon}} \odot g_t \qquad \theta = \theta - \alpha \cdot \nabla_{\theta} J(\theta) \\ \theta = \theta - \alpha \cdot g$$

 Scale factor of a parameter is inversely proportional to the square root of squared sum of past gradients

Adagrad

- One of Adagrad's main benefits is that it eliminates the need to manually tune the learning rate. Most implementations use a default value of 0.01.
- Adagrad's main weakness is its accumulation of the squared gradients in the denominator:
- Since every added term is positive, the accumulated sum keeps growing during training.
- This in turn causes the learning rate to shrink and eventually become infinitesimally small, at which point the algorithm is no longer able to acquire additional knowledge.

RMSprop(Root Mean Squared Propagation)

- In RMSprop is also the learning rate is adapted for each of the parameters.
- RMSprop have been developed to resolve Adagrad's radically diminishing learning rates.
- RMSprop divides the learning rate by an exponentially decaying average of squared gradients.

$$s_t = \gamma s_{t-1} + (1 - \gamma) g_t^2$$
 $s_t = s_{t-1} + g_t^2$

 Replace the sum of the squares of the past gradients with the decaying average over past squared gradients

$$\theta_{t+1} = \theta_t - \frac{\alpha}{\sqrt{s_t + \epsilon}} g_t$$

- Adaptive Moment Estimation (Adam) is also computes adaptive learning rates for each parameter.
- In addition to storing an exponentially decaying average of past squared gradients like RMSprop,
- Adam also keeps an exponentially decaying average of past gradients, similar to momentum.

- We compute the decaying averages of past gradients v_t and and also past squared gradients s_t as follows
- $v_t = \beta_1 v_{t-1} + (1 \beta_1) g_t$
- $s_t = \beta_2 s_{t-1} + (1 \beta_2) g_t^2$
- As v_t and s_t are initialized as vectors of 0's,
- the authors of Adam observe that they are biased towards zero, especially during the initial time steps, and especially when the decay rates are small (i.e. β_1 and β_2 are close to 1).

 They reduce these biases by computing biascorrected first and second moment estimates:

$$\hat{\mathbf{v}}_t = \frac{\mathbf{v}_t}{1 - \beta_1^t} \qquad \hat{\mathbf{s}}_t = \frac{\mathbf{s}_t}{1 - \beta_2^t}$$

 They then use these to update the parameters just as we have seen in RMSprop, which yields the Adam update rule:

$$\theta_{t+1} = \theta_t - \frac{\alpha}{\sqrt{\hat{S}_t + \epsilon}} \hat{v}_t$$

- It combines best features of adagrad and RMSprop
- It adds momentum and adaptive learning rate to the basic gradient descent.

Adam is the best among adaptive optimizers in most of the cases