

Gradient and Stochastic Gradient Descent

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Linear Least Squares Regression

Setup

- Input space $\mathcal{X} = \mathbf{R}^d$
 - Output space $\mathcal{Y} = \mathbf{R}$
 - Action space $\mathcal{Y} = \mathbf{R}$
 - Loss: $\ell(\hat{y}, y) = \frac{1}{2} (y - \hat{y})^2$
 - **Hypothesis space:** $\mathcal{F} = \{f : \mathcal{X} \rightarrow \mathcal{Y} \mid f(x) = w^T x\}$
-
- Given data set $\mathcal{D}_n = \{(x_1, y_1), \dots, (x_n, y_n)\}$,
 - Let's find the ERM $\hat{f} \in \mathcal{F}$.

Linear Least Squares Regression

Objective Function: Empirical Risk

The function we want to minimize is the empirical risk:

$$\hat{R}_n(w) = \frac{1}{n} \sum_{i=1}^n (w^T x_i - y_i)^2,$$

where $w \in \mathbf{R}^d$ parameterizes the hypothesis space \mathcal{F} .

Unconstrained Optimization

Setting

Objective function $f : \mathbf{R}^d \rightarrow \mathbf{R}$ is *differentiable*.

Want to find

$$x^* = \arg \min_{x \in \mathbf{R}^d} f(x)$$

The Gradient

Definition

The **gradient** $\nabla_x f(x_0)$ of a differentiable function $f(x)$ at the point x_0 is the direction to move in for the **fastest increase** in $f(x)$, when starting from x_0 .

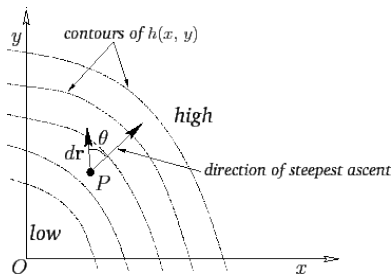


Figure: Figure A.111 from *Newtonian Dynamics*, by Richard Fitzpatrick.

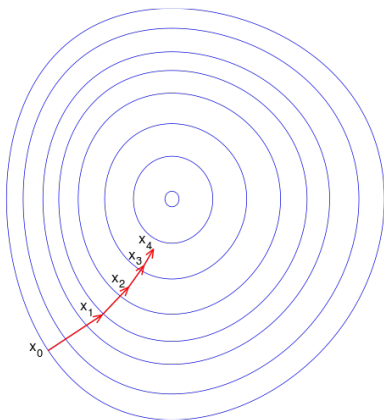
Gradient Descent

Gradient Descent

- Initialize $x = 0$
- repeat
 - $x \leftarrow x - \underbrace{\eta}_{\text{step size}} \nabla f(x)$
- until stopping criterion satisfied

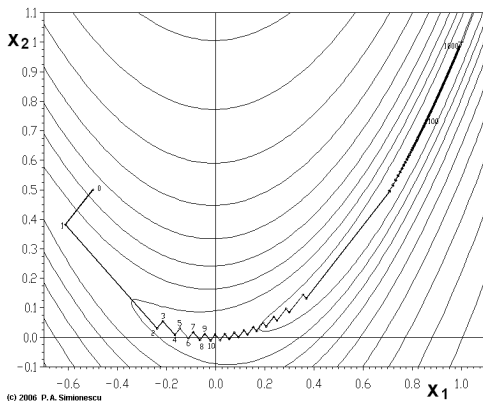
Gradient Descent Path

Gradient Descent for a nice (convex) function



Gradient Descent Path

Gradient Descent Path for the Rosenbrock Function (not convex)



(Figure by P.A. Simionescu from [Wikipedia page on gradient descent](#))

Gradient Descent - Details

Step Size

- Empirically $\eta = 0.1$ often works well
- **Better:** Optimize at every step (e.g. backtracking line search)

Stopping Rule

- Could use a maximum number of steps (e.g. 100)
- Wait until $\|\nabla f(x)\| \leq \varepsilon$.
- Test performance on holdout data (in learning setting)

Gradient Descent for Linear Regression

Gradient of Objective Function:

The gradient of the objective is

$$\begin{aligned}\nabla_w \hat{R}_n(w) &= \nabla_w \left[\frac{1}{n} \sum_{i=1}^n (w^T x_i - y_i)^2 \right] \\ &= \frac{2}{n} \sum_{i=1}^n \underbrace{(w^T x_i - y_i)}_{i\text{th residual}} x_i\end{aligned}$$

Gradient Descent: Does it scale?

- At every iteration, we compute the gradient at current w :

$$\nabla_w \hat{R}_n(w) = \frac{2}{n} \sum_{i=1}^n \underbrace{(w^T x_i - y_i)}_{i\text{th residual}} x_i$$

- We have to touch all n training points to take a single step. $[O(n)]$
 - Called a **batch optimization** method
- Can we make progress without looking at all the data?

Gradient Descent on the Risk

- Real goal is to minimize the risk (expected loss):

$$\arg \min_{f \in \mathcal{F}} \mathbb{E}[\ell(f(X), Y)]$$

- For linear regression, that's

$$\arg \min_w \mathbb{E} (w^T X - Y)^2$$

- Gradient descent on this?

$$\nabla_w \mathbb{E} (w^T X - Y)^2 = \mathbb{E} [2 (w^T X - Y) X]$$

Gradient Descent on the Risk [approximately]

- Want to find gradient of the risk:

$$\nabla R(w) = \mathbb{E} [2 (w^T X - Y) X]$$

- Can estimate expectation with a sample:

$$\widehat{\nabla R(w)} = \frac{1}{n} \sum_{i=1}^n \left[2 \left(\underbrace{w^T x_i - y_i}_{\text{i'th residual}} \right) x_i \right]$$

- Let's return to the general case...

Gradient Descent on the Risk: General Case

Gradient of Risk:

- Say hypothesis space \mathcal{F} is parameterized by $w \in \mathbf{R}^d$.
- Switching ∇_w and \mathbb{E} we can write the gradient of risk as

$$\text{Gradient(Risk)} = \nabla_w \mathbb{E}[\ell(f(X), Y)] = \mathbb{E}[\nabla_w \ell(f(X), Y)]$$

Unbiased estimator for Gradient(Risk):

$$\frac{1}{n} \sum_{i=1}^n [\nabla_w \ell(f_w(x_i), y_i)] \approx \underbrace{\mathbb{E}[\nabla_w \ell(f(X), Y)]}_{\text{Gradient(Risk)}}$$

Gradient Descent on the Risk: General Case

- We want $\text{Gradient}(\text{Risk})$
- Estimate it using sample of size n
- Bigger $n \implies$ Better estimate
- Bigger $n \implies$ Touching more data (slower!)
- But how big an n do we need?

Gradient Descent on the Risk [approximately]

- Gradient descent takes a bunch of steps whether we use
 - the perfect step direction $\nabla R(w)$,
 - an empirical estimate using all training data $\nabla \hat{R}_n(w)$, or
 - an empirical estimate using a random subset of data $\nabla \hat{R}_N(w)$ ($N \ll n$)
- What about $N = 1$?
- Even with a sample of size 1, the estimate

$$\nabla_w \ell(f_w(x_i), y_i)$$

is still **unbiased for gradient(Risk)**.

Stochastic Gradient Descent (SGD)

Stochastic Gradient Descent

- initialize $w = 0$
- repeat
 - randomly choose training point $(x_i, y_i) \in \mathcal{D}_n$
 - $w \leftarrow w - \eta \underbrace{\nabla_w \ell(f_w(x_i), y_i)}_{\text{Grad(Loss on i'th example)}}$
- until stopping criteria met

SGD: Step Size

- Let η_t be the step size at the t 'th step.
- How should η_t 's decrease with each step?

Robbins-Monro Conditions

Many classical convergence results depend on the following two conditions:

$$\sum_{t=1}^{\infty} \eta_t^2 < \infty \quad \sum_{t=1}^{\infty} \eta_t = \infty$$

- As fast as $\eta_t = O\left(\frac{1}{t}\right)$ would satisfy this... but should be faster than $O\left(\frac{1}{\sqrt{t}}\right)$.
- A useful reference for practical techniques: Leon Bottou's "Tricks":
<http://research.microsoft.com/pubs/192769/tricks-2012.pdf>