# **Gradient Descent**

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What to do if we don't have an optimization method?

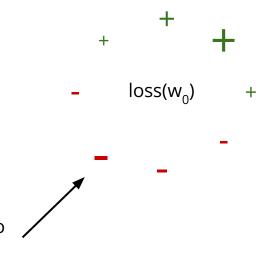
Optimization method when there is no closed form solution to finding the extrema of a function.

**Example**: Logistic Regression

**Goal**: find a sequence of  $w_i$ 's (and b's) that converge toward **a** minimum.

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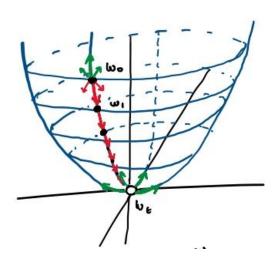
Clearly this is the best nudge to give  $w_0$  to reduce our Loss

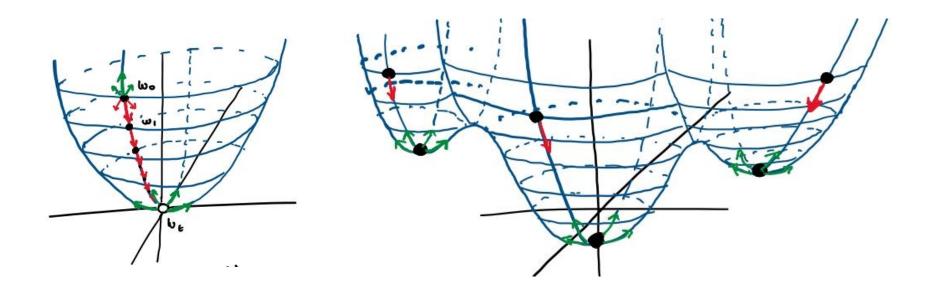
As such we can define the following sequence:

```
w_1 = best nudge to w_0
w_2 = best nudge to w_1
...
```

Until we reach w<sub>+</sub> that looks like this:

At this point we can stop updating w. Why?



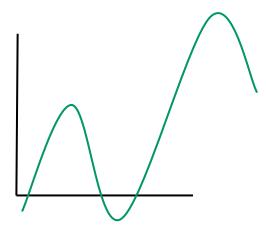


How can we know how much to nudge and in what direction?

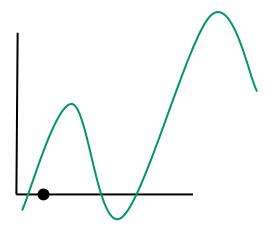
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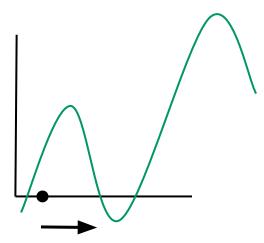
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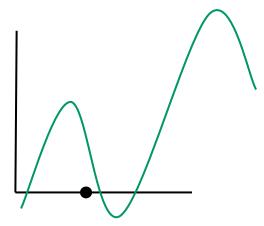
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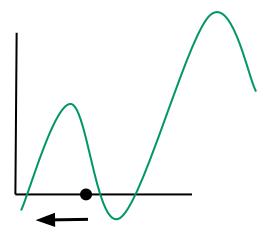
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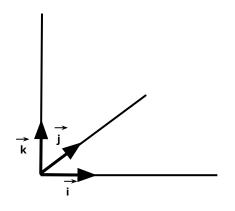


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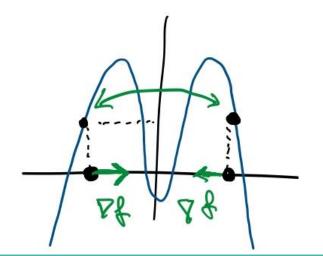


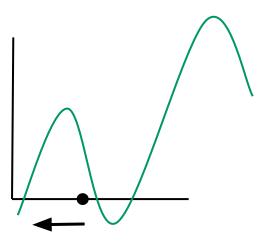
$$\nabla f(x, y, z) = \frac{\partial f}{\partial x}\vec{i} + \frac{\partial f}{\partial y}\vec{j} + \frac{\partial f}{\partial z}\vec{k}$$

# **Example**

However, the gradient expresses the **instantaneous** rate of change. At p,  $\nabla f_p$  is the steepest but the highest value of f will depend on how many units we step in that direction. If we step too many units away, the instantaneous change in f is no longer representative of what values f will take.

Example:





#### **Gradient Descent**

Given a "smooth" function f for which there exists no closed form solution for finding its **maximum**, we can find a local maximum through the following steps:

- 1. Define a step size  $\alpha$  (tuning parameter)
- 2. Initialize p to be random
- 3.  $p_{\text{new}} = \alpha \nabla f_p + p$
- 4.  $p \square p_{new}$
- 5. Repeat 3 & 4 until p  $\sim$  p<sub>new</sub>

To find a local **minimum**, just use  $-\nabla f_{D}$ 

#### **Gradient Descent**

#### Notes about $\alpha$ :

- If  $\alpha$  is too large, GD may overshoot the maximum, take a long time to or never be able to converge
- If  $\alpha$  is too small, GD may take too long to converge

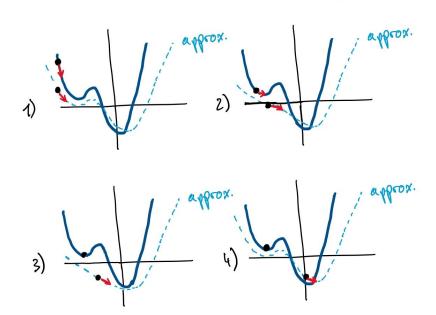
#### **Stochastic Gradient Descent**

Recall the Cost is computed for the entire dataset. This has some limitations:

- 1. It's expensive to run
- 2. The result we get depends only on the initial starting point

#### **Stochastic Gradient Descent**

**Goal**: Approximate the gradient of the Cost using a sample of the data (batch)



### Note

The magnitude of  $\nabla f_p$  depends on p. A p gets closer to the min / max, the size of  $\nabla f_p$  decreases.

This also means that points p that contain more "information" have larger gradients. So the order with which this process is exposed to examples matters.