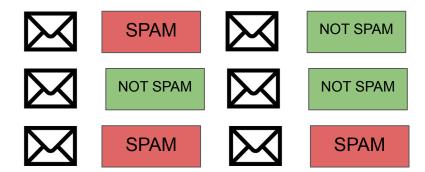


Given a labeled dataset, we want a model that takes as input a text and returns a class label

$$\hat{y} = rg \max_{y} P(y|\mathbf{x}; heta)$$



We represent text as a bag of words, and the model should return the maximum posterior probability given the text

Applying Bayes' Rule and simplifying:

$$\hat{y} = rg \max_{y} P(y|\mathbf{x}; heta) = rg \max_{y} rac{P(\mathbf{x}|y)P(y))}{P(\mathbf{x})} = rg \max_{y} P(\mathbf{x}|y)P(y)$$

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This requires more



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work

Bag of Words

```
P(y \mid \mathbf{x}) = P(SPAM \mid click, here, for, your, new, ...)
= P(y \mid \mathbf{x}) = P(SPAM \mid here, for, click, new, your, ...)
```

= ...

Jurafsky and Martin, Ch. 4

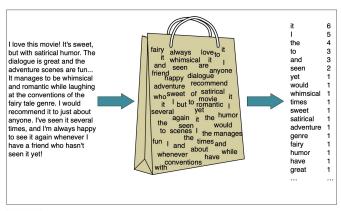


Figure 4.1 Intuition of the multinomial naive Bayes classifier applied to a movie review. The position of the words is ignored (the *bag of words* assumption) and we make use of the frequency of each word.

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Example of this assumption going wrong?

Jurafsky and Martin, Ch. 4

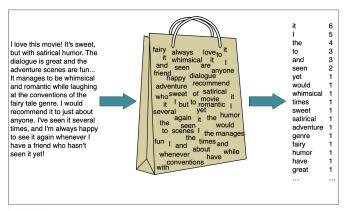


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Example of this assumption going wrong? What issue do we have here with MLE?

Jurafsky and Martin, Ch. 4

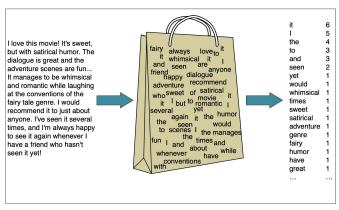


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Multinomial Naive Bayes

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Bag of Words

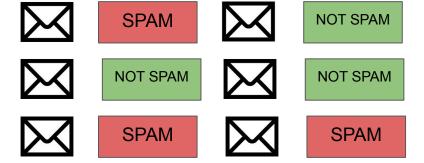
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$$P(| span |) = (# | span | / K) = 3 / 6$$

$$P(NOTSPAM) = (#NOTSPAM) / K) = 3 / 6$$























SPAM

To learn the **likelihoods** P(x|y), the Naive Bayes assumption gives

$$P(\mathbf{x}|y) = \prod_{j} P(x_{j}|y)$$

and the relative frequency for some $P(x_i|y)$ is

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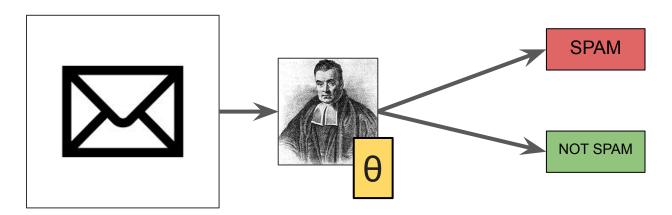
$$P(x_i|y) = count(x_i, y) / count(y)$$

Notice that the product $\prod_i P(x_i|y)$ is zero if any $P(x_i|y)$ is estimated to be zero!

Caveats: Smoothing

We've learned a model but during testing, what if we encounter a word not seen in training?

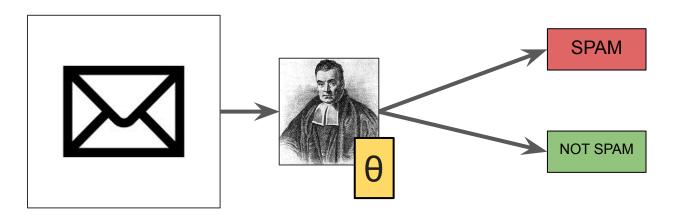
Laplace smoothing: $P(x_j|y) = count(x_j, y) + 1 / count(y) + V + 1$



Caveats: Smoothing

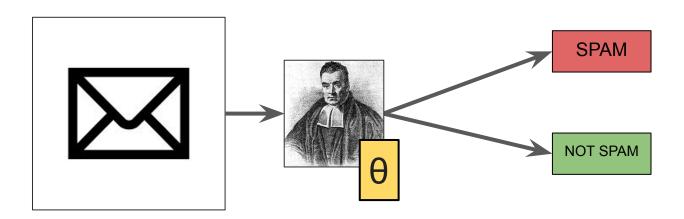
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Lidstone smoothing: $P(x_i|y) = count(x_i, y) + \alpha / count(y) + \alpha(V + 1)$



Caveats: Underflow

We store log probabilities to avoid numerical underflow



Gradient Descent Optimization

The chicken and the egg

Training neural networks involves two basic ingredients

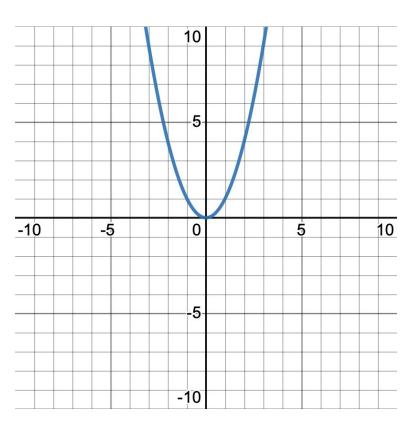
- Optimizing the parameters of a network with an iterative approach like gradient descent
- Computing the gradient of the loss function efficiently using backpropagation

We've previously talked about the central ingredients to a machine learning problem: data, a model with an associated choice of loss function, and an optimization algorithm

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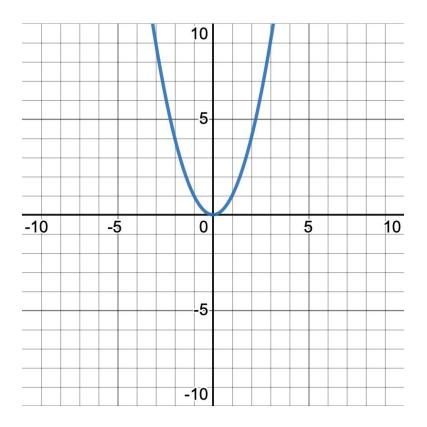
- In supervised learning, we saw that the learning process unfolds using labeled data
- We select a **model** that specifies inductive biases that are hopefully relevant to the task (c.f., BOW, Naive Bayes assumptions)
- The **loss function** measures how well the model is predicting the labels
- The optimization algorithm describes how the model parameters are adjusted

What do we do when we want to determine the optima of a function?



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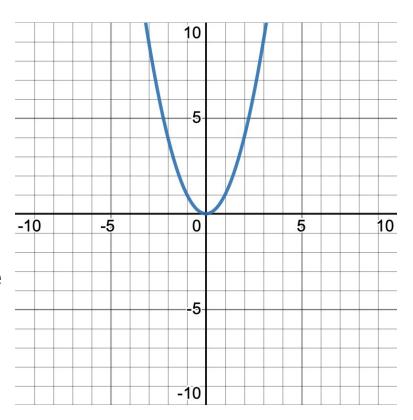
How do we find the minimum of $f(x) = x^2$? Exercise: find the minimum.

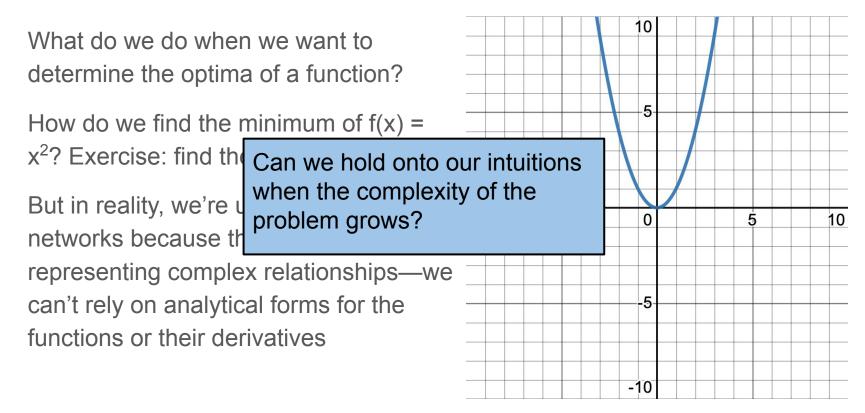


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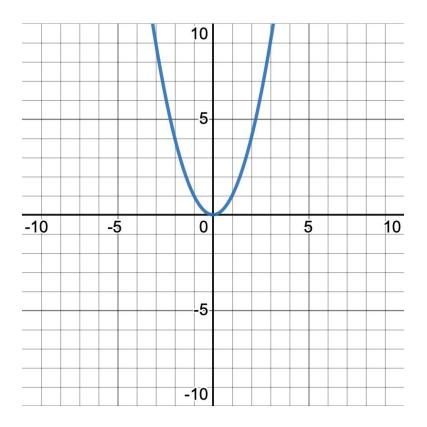
How do we find the minimum of $f(x) = x^2$? Exercise: find the minimum.

But in reality, we're using neural networks because they're capable of representing complex relationships—we can't rely on analytical forms for the functions or their derivatives



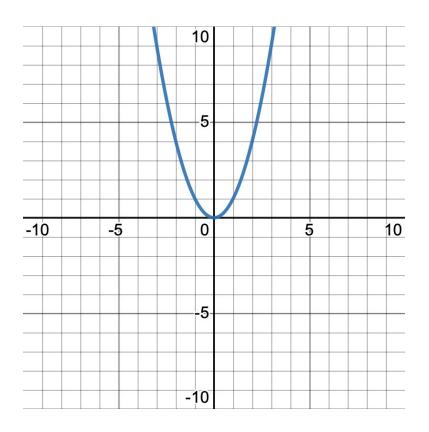


If we can still compute outputs and derivatives of the function for particular inputs, then we can still rely on the same intuition as before, that derivatives provide us with useful optimization information



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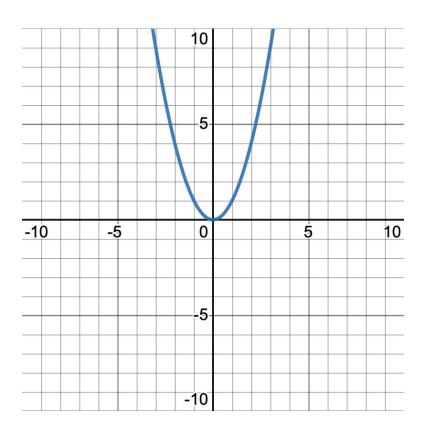
The non-linearity of neural networks means that the loss functions can be complex (non-convex)



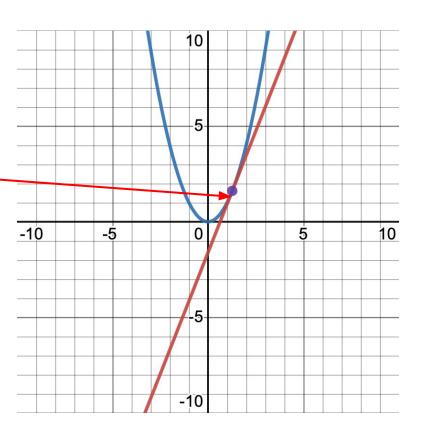
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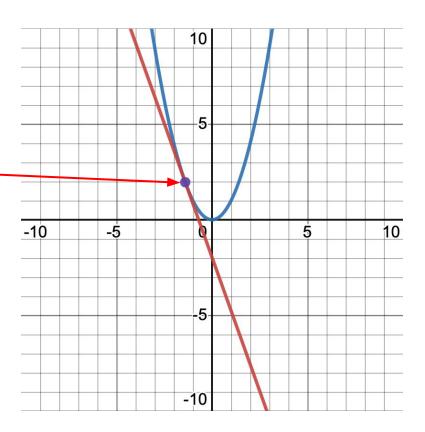
So, we will use **iterative methods** that slowly decrement the loss



If we compute the derivative here, how we should we adjust our estimate of the minimum x?



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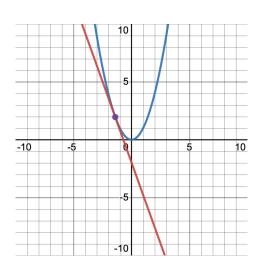


Procedure:

- Randomly select x
- Evaluate gradient
- Move in the direction of the negative gradient (i.e., opposite the slope in each dimension)

$$\hat{g} \leftarrow rac{1}{N}
abla_{ heta} \sum_{i} L\left(f(x^{(i)}; heta), y^{(i)}
ight)$$

$$\theta \leftarrow \theta - \alpha \hat{g}$$



The key issue with vanilla/batch gradient descent is that the computation of the gradient is based on the entire dataset

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Mini-batch gradient descent computes gradients on small random sets of instances called *mini-batches*

$$\hat{g} \leftarrow rac{1}{m}
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Often, authors refer to mini-batch gradient descent and stochastic gradient descent interchangeably

Algorithm 8.1 Stochastic gradient descent (SGD) update

```
Require: Learning rate schedule \epsilon_1, \epsilon_2, \ldots
Require: Initial parameter \theta
   k \leftarrow 1
    while stopping criterion not met do
       Sample a minibatch of m examples from the training set \{x^{(1)}, \dots, x^{(m)}\} with
       corresponding targets y^{(i)}.
       Compute gradient estimate: \hat{\boldsymbol{g}} \leftarrow \frac{1}{m} \nabla_{\boldsymbol{\theta}} \sum_{i} L(f(\boldsymbol{x}^{(i)}; \boldsymbol{\theta}), \boldsymbol{y}^{(i)})
       Apply update: \boldsymbol{\theta} \leftarrow \boldsymbol{\theta} - \epsilon_k \hat{\boldsymbol{q}}
       k \leftarrow k + 1
    end while
```

Recall that the softmax function produces a probability distribution from a numeric vector When working with neural networks, we call these raw outputs **logits**. They are the inputs to the softmax.

The network outputs a categorical distribution, and through learning, the parameters are adjusted so the model distribution matches the target distribution codified by the labeled data

$$\sigma(\mathbf{z})_i = rac{e^{z_i}}{\sum_{j=1}^K e^{z_j}}$$

We use the **cross-entropy loss**, which is the negative log likelihood:

$$egin{aligned} \hat{ heta} = rgmin_{ heta} - \mathbb{E}_{\mathbf{x}, \mathbf{y} \sim P_D}[\log P(\mathbf{y} | \mathbf{x}; heta)] \ \sigma(\mathbf{z})_i = rac{e^{z_i}}{\sum_{j=1}^K e^{z_j}} \end{aligned}$$

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$$\hat{ heta} = \operatorname*{argmin}_{ heta} - \mathbb{E}_{\mathbf{x},\mathbf{y}\sim P_D}[\log P(\mathbf{y}|\mathbf{x}; heta)]$$

Note that Naive Bayes is **generative** (since it classifies according to the joint probability of **x** and y), while logistic regression (K=2) and softmax regression (K>3) are **discriminative** (since they are trained to directly classify **x**)

$$\sigma(\mathbf{z})_i = rac{e^{z_i}}{\sum_{j=1}^K e^{z_j}}$$

Softmax Regression Demo