CSC 411 Lecture 8: Linear Classification II

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Today's agenda:

- Gradient checking with finite differences
- Learning rates
- Stochastic gradient descent
- Convexity
- Multiclass classification and softmax regression
- Limits of linear classification

- We've derived a lot of gradients so far. How do we know if they're correct?
- Recall the definition of the partial derivative:

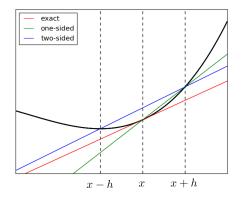
$$\frac{\partial}{\partial x_i}f(x_1,\ldots,x_N) = \lim_{h\to 0}\frac{f(x_1,\ldots,x_i+h,\ldots,x_N)-f(x_1,\ldots,x_i,\ldots,x_N)}{h}$$

• Check your derivatives numerically by plugging in a small value of h, e.g. 10^{-10} . This is known as finite differences.

Gradient Checking

• Even better: the two-sided definition

$$\frac{\partial}{\partial x_i}f(x_1,\ldots,x_N) = \lim_{h\to 0}\frac{f(x_1,\ldots,x_i+h,\ldots,x_N)-f(x_1,\ldots,x_i-h,\ldots,x_N)}{2h}$$



- Run gradient checks on small, randomly chosen inputs
- Use double precision floats (not the default for TensorFlow, PyTorch, etc.!)
- Compute the relative error between derived gradient and finite difference approximation:

$$\frac{|a-b|}{|a|+|b|}$$

• The relative error should be very small, e.g. 10^{-6}

- Gradient checking is really important!
- Learning algorithms often appear to work even if the math is wrong.

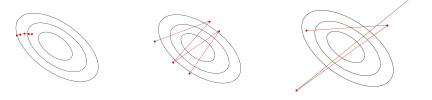
• But:

- They might work much better if the derivatives are correct.
- Wrong derivatives might lead you on a wild goose chase.
- If you implement derivatives by hand, gradient checking is the single most important thing you need to do to get your algorithm to work well.

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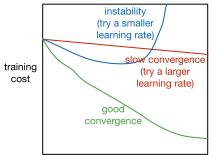
 In gradient descent, the learning rate α is a hyperparameter we need to tune. Here are some things that can go wrong:



- Good values are typically between 0.001 and 0.1. You should do a grid search if you want good performance (i.e. try 0.1, 0.03, 0.01, ...).

Training Curves

• To diagnose optimization problems, it's useful to look at training curves: plot the training cost as a function of iteration.



iteration #

• Warning: it's very hard to tell from the training curves whether an optimizer has converged. They can reveal major problems, but they can't guarantee convergence.

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• So far, the cost function $\mathcal J$ has been the average loss over the training examples:

$$\mathcal{J}(\boldsymbol{\theta}) = \frac{1}{N} \sum_{i=1}^{N} \mathcal{L}^{(i)} = \frac{1}{N} \sum_{i=1}^{N} \mathcal{L}(y(\mathbf{x}^{(i)}, \boldsymbol{\theta}), t^{(i)}).$$

• By linearity,

$$rac{\partial \mathcal{J}}{\partial oldsymbol{ heta}} = rac{1}{N} \sum_{i=1}^{N} rac{\partial \mathcal{L}^{(i)}}{\partial oldsymbol{ heta}}.$$

- Computing the gradient requires summing over *all* of the training examples. This is known as batch training.
- Batch training is impractical if you have a large dataset (e.g. millions of training examples)!

• Stochastic gradient descent (SGD): update the parameters based on the gradient for a single training example, chosen uniformly at random:

$$oldsymbol{ heta} \leftarrow oldsymbol{ heta} - lpha rac{\partial \mathcal{L}^{(i)}}{\partial oldsymbol{ heta}}$$

- SGD can make significant progress before it has even looked at all the data!
- Mathematical justification: if you sample a training example uniformly at random, the stochastic gradient is an unbiased estimate of the batch gradient:

$$\mathbb{E}\left[\frac{\partial \mathcal{L}^{(i)}}{\partial \boldsymbol{\theta}}\right] = \frac{1}{N} \sum_{i=1}^{N} \frac{\partial \mathcal{L}^{(i)}}{\partial \boldsymbol{\theta}} = \frac{\partial \mathcal{J}}{\partial \boldsymbol{\theta}}.$$

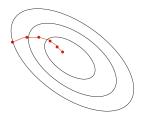
- Problems:
 - Variance in this estimate may be high
 - If we only look at one training example at a time, we can't exploit efficient vectorized operations.

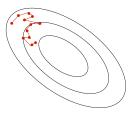
- Compromise approach: compute the gradients on a randomly chosen medium-sized set of training examples *M* ⊂ {1,..., *N*}, called a mini-batch.
- Stochastic gradients computed on larger mini-batches have smaller variance:

$$\operatorname{Var}\left[\frac{1}{|\mathcal{M}|}\sum_{i\in\mathcal{M}}\frac{\partial\mathcal{L}^{(i)}}{\partial\theta_{j}}\right] = \frac{1}{|\mathcal{M}|^{2}}\sum_{i\in\mathcal{M}}\operatorname{Var}\left[\frac{\partial\mathcal{L}^{(i)}}{\partial\theta_{j}}\right] = \frac{1}{|\mathcal{M}|}\operatorname{Var}\left[\frac{\partial\mathcal{L}^{(i)}}{\partial\theta_{j}}\right]$$

- The mini-batch size $|\mathcal{M}|$ is a hyperparameter that needs to be set.
 - Too large: takes more memory to store the activations, and longer to compute each gradient update
 - Too small: can't exploit vectorization
 - A reasonable value might be $|\mathcal{M}| = 100$.

• Batch gradient descent moves directly downhill. SGD takes steps in a noisy direction, but moves downhill on average.



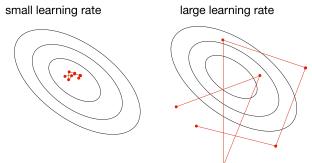


batch gradient descent

stochastic gradient descent

SGD Learning Rate

• In stochastic training, the learning rate also influences the fluctuations due to the stochasticity of the gradients.



- Typical strategy:
 - Use a large learning rate early in training so you can get close to the optimum
 - Gradually decay the learning rate to reduce the fluctuations

• Warning: by reducing the learning rate, you reduce the fluctuations, which can appear to make the loss drop suddenly. But this can come at the expense of long-run performance.



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Convex Sets

Convex Sets



• A set S is convex if any line segment connecting points in S lies entirely within S. Mathematically,

$$\mathbf{x}_1, \mathbf{x}_2 \in \mathcal{S} \implies \lambda \mathbf{x}_1 + (1 - \lambda) \mathbf{x}_2 \in \mathcal{S} \text{ for } \mathbf{0} \leq \lambda \leq 1.$$

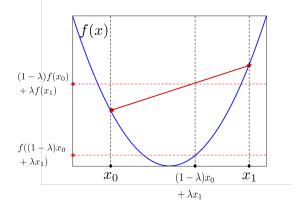
A simple inductive argument shows that for x₁,..., x_N ∈ S, weighted averages, or convex combinations, lie within the set:

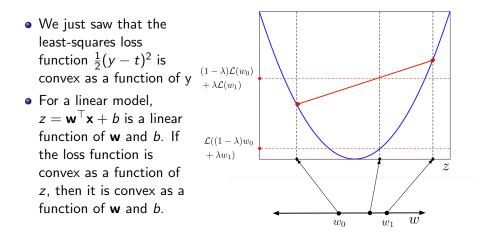
$$\lambda_1 \mathbf{x}_1 + \cdots + \lambda_N \mathbf{x}_N \in S \quad \text{for } \lambda_i > 0, \ \lambda_1 + \cdots + \lambda_N = 1.$$

• A function f is convex if for any $\mathbf{x}_0, \mathbf{x}_1$ in the domain of f,

$$f((1 - \lambda)\mathbf{x}_0 + \lambda \mathbf{x}_1) \leq (1 - \lambda)f(\mathbf{x}_0) + \lambda f(\mathbf{x}_1)$$

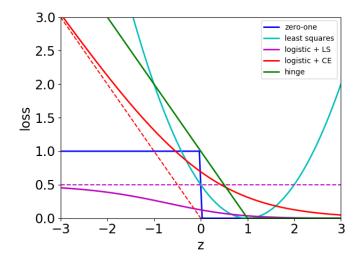
- Equivalently, the set of points lying above the graph of *f* is convex.
- Intuitively: the function is bowl-shaped.





Convex Functions

Which loss functions are convex?



Why we care about convexity

- All critical points are minima
- Gradient descent finds the optimal solution (more on this in a later lecture)

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• What about classification tasks with more than two categories?





- Targets form a discrete set $\{1, \ldots, K\}$.
- It's often more convenient to represent them as one-hot vectors, or a one-of-K encoding:

$$\mathbf{t} = \underbrace{(0, \dots, 0, 1, 0, \dots, 0)}_{\text{entry } k \text{ is } 1}$$

- Now there are D input dimensions and K output dimensions, so we need K × D weights, which we arrange as a weight matrix W.
- Also, we have a K-dimensional vector **b** of biases.
- Linear predictions:

$$z_k = \sum_j w_{kj} x_j + b_k$$

• Vectorized:

 $\mathbf{z} = \mathbf{W}\mathbf{x} + \mathbf{b}$

• A natural activation function to use is the softmax function, a multivariable generalization of the logistic function:

$$y_k = ext{softmax}(z_1, \dots, z_K)_k = rac{e^{z_k}}{\sum_{k'} e^{z_{k'}}}$$

- The inputs z_k are called the logits.
- Properties:
 - Outputs are positive and sum to 1 (so they can be interpreted as probabilities)
 - If one of the z_k 's is much larger than the others, $\operatorname{softmax}(z)$ is approximately the argmax. (So really it's more like "soft-argmax".)
 - **Exercise:** how does the case of K = 2 relate to the logistic function?
- Note: sometimes σ(z) is used to denote the softmax function; in this class, it will denote the logistic function applied elementwise.

• If a model outputs a vector of class probabilities, we can use cross-entropy as the loss function:

$$egin{aligned} \mathcal{L}_{ ext{CE}}(\mathbf{y},\mathbf{t}) &= -\sum_{k=1}^{K} t_k \log y_k \ &= -\mathbf{t}^{ op}(\log \mathbf{y}), \end{aligned}$$

where the log is applied elementwise.

• Just like with logistic regression, we typically combine the softmax and cross-entropy into a softmax-cross-entropy function.

• Softmax regression:

$$\begin{split} \mathbf{z} &= \mathbf{W}\mathbf{x} + \mathbf{b} \\ \mathbf{y} &= \operatorname{softmax}(\mathbf{z}) \\ \mathcal{L}_{\operatorname{CE}} &= -\mathbf{t}^{\top}(\log \mathbf{y}) \end{split}$$

• Gradient descent updates are derived in the readings:

$$\frac{\partial \mathcal{L}_{\rm CE}}{\partial \textbf{z}} = \textbf{y} - \textbf{t}$$

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• Visually, it's obvious that **XOR** is not linearly separable. But how to show this?



Showing that XOR is not linearly separable

- Half-spaces are obviously convex.
- Suppose there were some feasible hypothesis. If the positive examples are in the positive half-space, then the green line segment must be as well.
- Similarly, the red line segment must line within the negative half-space.



• But the intersection can't lie in both half-spaces. Contradiction!

Limits of Linear Classification

A more troubling example



- These images represent 16-dimensional vectors. White = 0, black = 1.
- Want to distinguish patterns A and B in all possible translations (with wrap-around)
- Translation invariance is commonly desired in vision!

Limits of Linear Classification

A more troubling example



- These images represent 16-dimensional vectors. White = 0, black = 1.
- Want to distinguish patterns A and B in all possible translations (with wrap-around)
- Translation invariance is commonly desired in vision!
- Suppose there's a feasible solution. The average of all translations of A is the vector (0.25, 0.25, ..., 0.25). Therefore, this point must be classified as A.
- Similarly, the average of all translations of B is also (0.25, 0.25, ..., 0.25). Therefore, it must be classified as B. Contradiction!

Limits of Linear Classification

• Sometimes we can overcome this limitation using feature maps, just like for linear regression. E.g., for **XOR**:

$$\psi(\mathbf{x}) = egin{pmatrix} x_1 \ x_2 \ x_1 x_2 \end{pmatrix}$$

x_1	<i>x</i> ₂	$\psi_1(\mathbf{x})$	$\psi_2(\mathbf{x})$	$\psi_{3}(\mathbf{x})$	t
0	0	0	0	0	0
0	1	0	1	0	1
1	0	1	0	0	1
1	1	1	1	1	0

- This is linearly separable. (Try it!)
- Not a general solution: it can be hard to pick good basis functions. Instead, we'll use neural nets to learn nonlinear hypotheses directly.