Lecture 3: Regularization and Optimization

Administrative: Assignment 1

Released last week, due Fri 4/19 at 11:59pm

Office hours: help with high-level questions only, no code debugging. [No Code Show Policy]

Administrative: Project proposal + Office Hours

Due Mon 4/22

TA expertise + Office Hours are posted on the webpage. Mix of inperson and zoom.

(http://cs231n.stanford.edu/office_hours.html)

Administrative: Ed

Please make sure to check and read all pinned Ed posts.

- <u>Project group</u>: fill in your information in the google form and/or look through existing responses and reach out
- <u>SCPD</u>: if you would like to take the midterm on-campus, send us an email: <u>cs231n-staff-spr24@stanford.edu</u> + cc <u>scpd-exams@stanford.edu</u>

Recap from Last Week

Image Classification: A core task in Computer Vision



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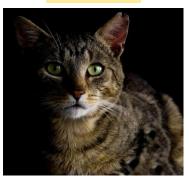
(assume given a set of labels) {dog, cat, truck, plane, ...} cat dog bird deer

truck

Recall from last time: Challenges of recognition

Viewpoint Viewpoint

Illumination



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Deformation



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Occlusion



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Clutter



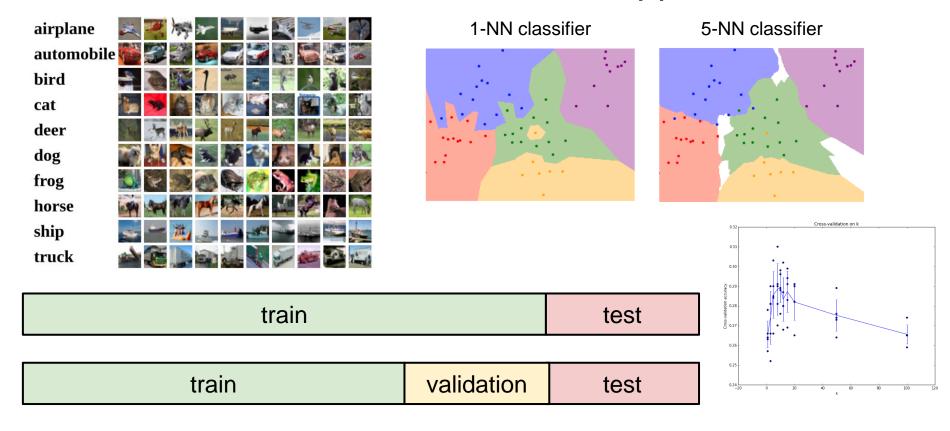
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Intraclass Variation

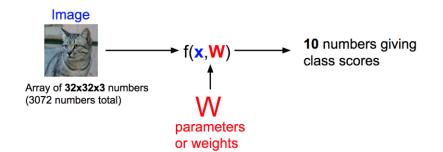


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Recall from last time: data-driven approach, kNN



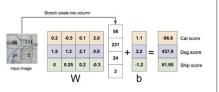
Recall from last time: Linear Classifier



$$f(x,W) = Wx + b$$



$$f(x,W) = Wx$$



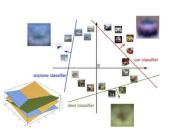


One template per class



Geometric Viewpoint

Hyperplanes cutting up space



Class 1:

Class 2: Everything else

1 <= L2 norm <= 2

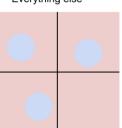


Class 1:

Three modes

Class 2:

Everything else







-1.7 2.0 Given a dataset of examples $\{(x_i, y_i)\}_{i=1}^N$

Where x_i is image and

our current classifier is

 y_i is (integer) label

Loss over the dataset is a average of loss over examples:

A **loss function** tells how good

$$L = \frac{1}{N} \sum_{i} L_i(f(x_i, W), y_i)$$





2.2

3.2

5.1

cat

car

frog

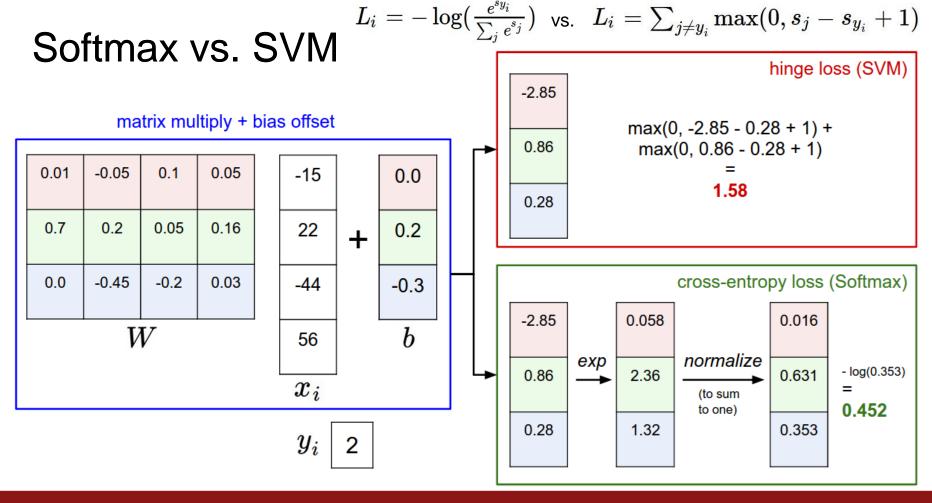
Suppose: 3 training examples, 3 classes.

With some W the scores f(x, W) = Wx are:

1.3

4.9

2.5 -3.1



$$f(x,W)=Wx$$

$$L = rac{1}{N} \sum_{i=1}^{N} \sum_{j
eq y_i} \max(0, f(x_i; W)_j - f(x_i; W)_{y_i} + 1)$$

Q: Suppose that we found a W such that L = 0. Is this W unique?

$$f(x,W) = Wx$$
 $L = rac{1}{N} \sum_{i=1}^N \sum_{j
eq y_i} \max(0, f(x_i; W)_j - f(x_i; W)_{y_i} + 1)$

Q: Suppose that we found a W such that L = 0. Is this W unique?

No! 2W is also has L = 0!

Suppose: 3 training examples, 3 classes. With some W the scores f(x, W) = Wx are:







cat	3.2	1.3	2.2
car	5.1	4.9	2.5
frog	-1.7	2.0	-3.1
Losses:	2.9	0	

$L_i = \sum_{j eq y_i} \max(0, s_j - s_{y_i} + 1)$

Before:

- $= \max(0, 1.3 4.9 + 1)$ $+ \max(0, 2.0 - 4.9 + 1)$
- $= \max(0, -2.6) + \max(0, -1.9)$ = 0 + 0
- = 0

With W twice as large:

- $= \max(0, 2.6 9.8 + 1)$ $+ \max(0, 4.0 - 9.8 + 1)$
- $+\max(0, 4.0 9.8 + 1)$ = $\max(0, -6.2) + \max(0, -4.8)$
- = 0 + 0
- = 0

$$f(x,W) = Wx$$
 $L = rac{1}{N} \sum_{i=1}^N \sum_{j
eq y_i} \max(0, f(x_i; W)_j - f(x_i; W)_{y_i} + 1)$

E.g. Suppose that we found a W such that L = 0. Is this W unique?

No! 2W is also has L = 0! How do we choose between W and 2W?

$$L(W) = \frac{1}{N} \sum_{i=1}^{N} L_i(f(x_i, W), y_i)$$

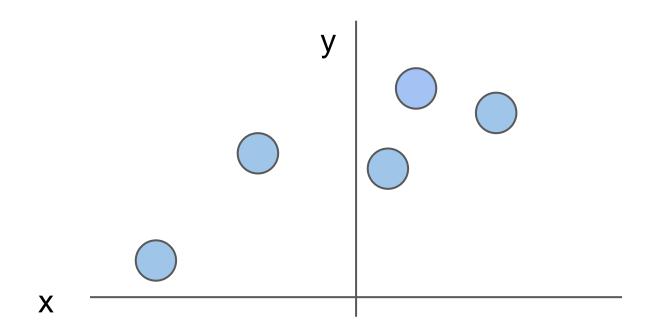
Data loss: Model predictions should match training data

$$L(W) = \underbrace{\frac{1}{N} \sum_{i=1}^{N} L_i(f(x_i, W), y_i) + \lambda R(W)}_{i=1}$$

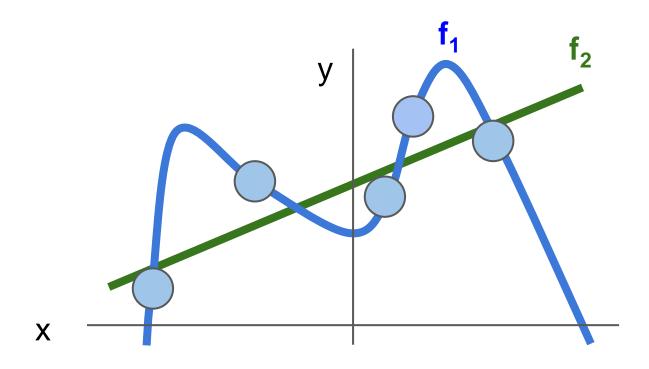
Data loss: Model predictions should match training data

Regularization: Prevent the model from doing *too* well on training data

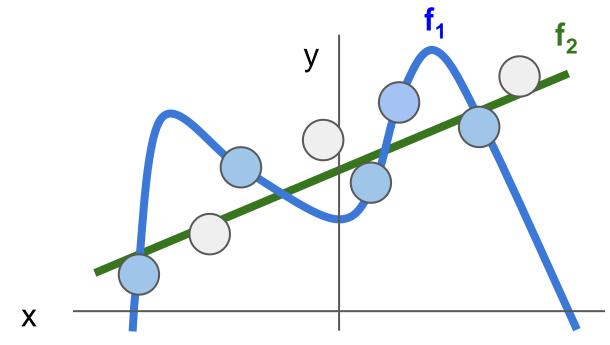
Regularization intuition: toy example training data



Regularization intuition: Prefer Simpler Models



Regularization: Prefer Simpler Models



Regularization pushes against fitting the data too well so we don't fit noise in the data

$$L(W) = \underbrace{\frac{1}{N} \sum_{i=1}^{N} L_i(f(x_i, W), y_i) + \lambda R(W)}_{i=1}$$

Data loss: Model predictions should match training data

Regularization: Prevent the model from doing *too* well on training data

Occam's Razor: Among multiple competing hypotheses, the simplest is the best. William of Ockham 1285-1347

$$\lambda = \frac{\text{regularization strength}}{\text{(hyperparameter)}}$$

$$L(W) = \underbrace{\frac{1}{N} \sum_{i=1}^{N} L_i(f(x_i, W), y_i) + \lambda R(W)}_{i=1}$$

Data loss: Model predictions should match training data

Regularization: Prevent the model from doing *too* well on training data

$$\lambda_{\cdot}$$
 = regularization strength (hyperparameter)

$$L(W) = \frac{1}{N} \sum_{i=1}^{N} L_i(f(x_i, W), y_i) + \lambda R(W)$$

Data loss: Model predictions should match training data

Regularization: Prevent the model from doing *too* well on training data

Simple examples

L2 regularization:
$$R(W) = \sum_k \sum_l W_{k,l}^2$$

L1 regularization:
$$R(W) = \sum_{k} \sum_{l} |W_{k,l}|$$

Elastic net (L1 + L2):
$$R(W) = \sum_k \sum_l \beta W_{k,l}^2 + |W_{k,l}|$$

$$\lambda_{\cdot}$$
 = regularization strength (hyperparameter)

$$L(W) = \underbrace{\frac{1}{N} \sum_{i=1}^{N} L_i(f(x_i, W), y_i) + \lambda R(W)}_{i=1}$$

Data loss: Model predictions should match training data

Regularization: Prevent the model from doing too well on training data

Simple examples

L2 regularization:
$$R(W) = \sum_{k} \sum_{l} W_{k,l}^2$$

L1 regularization: $R(W) = \sum_{k} \sum_{l} |W_{k,l}|$

Dropout

More complex:

Batch normalization

Elastic net (L1 + L2): $R(W) = \sum_{k} \sum_{l} \beta W_{k,l}^{2} + |W_{k,l}|$

April 9, 2024

Stochastic depth, fractional pooling, etc.

$$\lambda_{\cdot}$$
 = regularization strength (hyperparameter)

$$L(W) = \underbrace{\frac{1}{N} \sum_{i=1}^{N} L_i(f(x_i, W), y_i) + \lambda R(W)}_{i=1}$$

Data loss: Model predictions should match training data

Regularization: Prevent the model from doing *too* well on training data

Why regularize?

- Express preferences over weights
- Make the model *simple* so it works on test data
- Improve optimization by adding curvature

Regularization: Expressing Preferences

$$egin{aligned} x &= [1,1,1,1] \ w_1 &= [1,0,0,0] \end{aligned}$$

$$w_2 = [0.25, 0.25, 0.25, 0.25]$$

$$w_1^T x = w_2^T x = 1$$

L2 Regularization

$$R(W) = \sum_k \sum_l W_{k,l}^2$$

Which of w1 or w2 will the L2 regularizer prefer?

Regularization: Expressing Preferences

$$x = [1, 1, 1, 1] \ w_1 = [1, 0, 0, 0]$$

$$w_2 = \left[0.25, 0.25, 0.25, 0.25\right]$$

$$w_1^T x = w_2^T x = 1$$

L2 Regularization

$$R(W) = \sum_{k} \sum_{l} W_{k,l}^2$$

Which of w1 or w2 will the L2 regularizer prefer?

L2 regularization likes to "spread out" the weights

Regularization: Expressing Preferences

$$x = [1, 1, 1, 1]$$

 $w_1 = [1, 0, 0, 0]$

$$w_2 = \left[0.25, 0.25, 0.25, 0.25\right]$$

$$w_1^Tx=w_2^Tx=1$$

L2 Regularization

$$R(W) = \sum_{k} \sum_{l} W_{k,l}^2$$

Which of w1 or w2 will the L2 regularizer prefer?

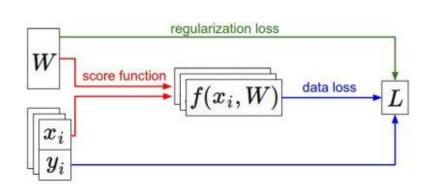
L2 regularization likes to "spread out" the weights

Which one would L1 regularization prefer?

Recap

- We have some dataset of (x,y)
- We have a **score function**: $s = f(x; W) \stackrel{\text{e.g.}}{=} Wx$
- We have a **loss function**:

$$L_i = -\log(rac{e^{sy_i}}{\sum_j e^{s_j}})$$
 SVM $L_i = \sum_{j
eq y_i} \max(0, s_j - s_{y_i} + 1)$ $L = rac{1}{N} \sum_{i=1}^N L_i + R(W)$ Full loss

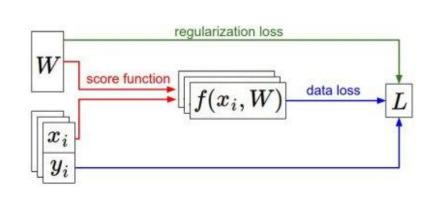


Recap

How do we find the best W?

- We have some dataset of (x,y)
- We have a **score function**: $s = f(x; W) \stackrel{\text{e.g.}}{=} Wx$
- We have a **loss function**:

$$L_i = -\log(rac{e^{sy_i}}{\sum_j e^{s_j}})$$
 SVM $L_i = \sum_{j
eq y_i} \max(0, s_j - s_{y_i} + 1)$ $L = rac{1}{N} \sum_{i=1}^N L_i + R(W)$ Full loss



Optimization



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Walking man image is CC0 1.0 public domain

Strategy #1: A first very bad idea solution: Random search

```
# assume X train is the data where each column is an example (e.g. 3073 x 50,000)
# assume Y train are the labels (e.g. 1D array of 50,000)
# assume the function L evaluates the loss function
bestloss = float("inf") # Python assigns the highest possible float value
for num in xrange(1000):
 W = np.random.randn(10, 3073) * 0.0001 # generate random parameters
 loss = L(X train, Y train, W) # get the loss over the entire training set
 if loss < bestloss: # keep track of the best solution
   bestloss = loss
   bestW = W
 print 'in attempt %d the loss was %f, best %f' % (num, loss, bestloss)
# prints:
# in attempt 0 the loss was 9.401632, best 9.401632
# in attempt 1 the loss was 8.959668, best 8.959668
# in attempt 2 the loss was 9.044034, best 8.959668
# in attempt 3 the loss was 9.278948, best 8.959668
# in attempt 4 the loss was 8.857370, best 8.857370
# in attempt 5 the loss was 8.943151, best 8.857370
# in attempt 6 the loss was 8.605604, best 8.605604
# ... (trunctated: continues for 1000 lines)
```

Lets see how well this works on the test set...

```
# Assume X_test is [3073 x 10000], Y_test [10000 x 1]
scores = Wbest.dot(Xte_cols) # 10 x 10000, the class scores for all test examples
# find the index with max score in each column (the predicted class)
Yte_predict = np.argmax(scores, axis = 0)
# and calculate accuracy (fraction of predictions that are correct)
np.mean(Yte_predict == Yte)
# returns 0.1555
```

15.5% accuracy! not bad! (SOTA is ~99.7%)

Strategy #2: Follow the slope



Strategy #2: Follow the slope

In 1-dimension, the derivative of a function:

$$rac{df(x)}{dx} = \lim_{h o 0} rac{f(x+h) - f(x)}{h}$$

In multiple dimensions, the **gradient** is the vector of (partial derivatives) along each dimension

The slope in any direction is the **dot product** of the direction with the gradient The direction of steepest descent is the **negative gradient**

[0.34,-1.11, 0.78, 0.12, 0.55, 2.81, -3.1, -1.5, 0.33,...] loss 1.25347 Lecture 3 - 42 April 9, 2024 Fei-Fei Li, Ehsan Adeli, Zane Durante

gradient dW:

current W:

[0.34,[0.34 + 0.0001]-1.11, -1.11, 0.78, 0.78, 0.12, 0.12, 0.55, 0.55, 2.81, 2.81, -3.1, -3.1, -1.5, -1.5, 0.33,...] [0.33,...]?,...] loss 1.25347 loss 1.25322 Fei-Fei Li, Ehsan Adeli, Zane Durante Lecture 3 - 43 April 9, 2024

gradient dW:

W + h (first dim):

current W:

[0.34,[0.34 + 0.0001,**[-2.5**, -1.11, -1.11, 0.78, 0.78, 0.12, 0.12, (1.25322 - 1.25347)/0.00010.55, 0.55, = -2.52.81, 2.81, $\frac{df(x)}{dx} = \lim_{x \to 0} \frac{f(x+h) - f(x)}{f(x+h)}$ -3.1, -3.1, -1.5, -1.5, [0.33,...]0.33,...?,...] loss 1.25347 loss 1.25322

Lecture 3 - 44

gradient dW:

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W + h (first dim):

current W:

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[0.34,[0.34,[-2.5, -1.11, -1.11 + 0.00010.78, 0.78, 0.12, 0.12, 0.55, 0.55, 2.81, 2.81, -3.1, -3.1, -1.5, -1.5, 0.33,...] [0.33,...]?,...] loss 1.25347 loss 1.25353 Fei-Fei Li, Ehsan Adeli, Zane Durante Lecture 3 - 45 April 9, 2024

gradient dW:

W + h (second dim):

current W:

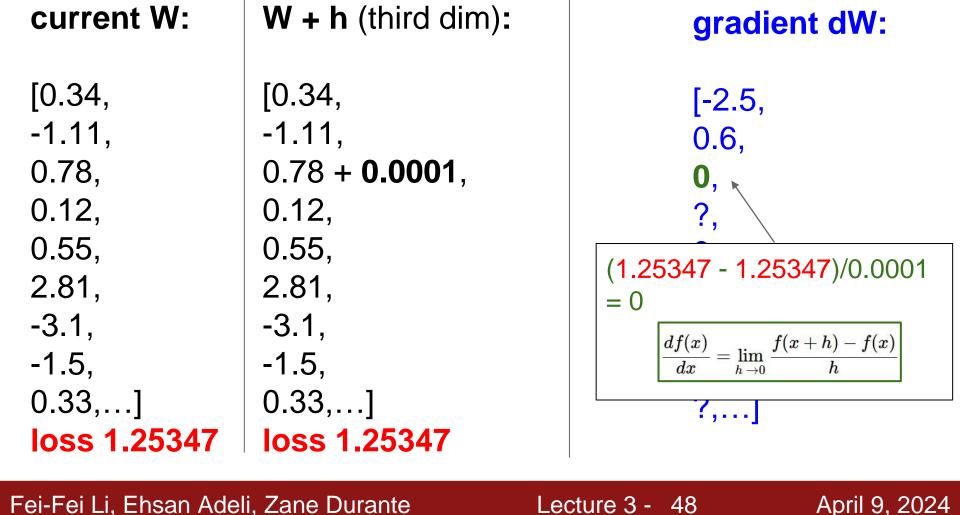
W + h (second dim): current W: gradient dW: [0.34,[0.34,[-2.5, -1.11, -1.11 + 0.0001, **0.6**, k 0.78, 0.78, 0.12, 0.12, 0.55, 0.55, (1.25353 - 1.25347)/0.00012.81, 2.81, = 0.6-3.1, -3.1, -1.5, -1.5, [0.33,...][0.33,...]?,...] loss 1.25347 loss 1.25353

[0.34,[0.34,[-2.5, -1.11, -1.11, 0.6, 0.78, 0.78 + 0.00010.12, 0.12, 0.55, 0.55, 2.81, 2.81, -3.1, -3.1, -1.5, -1.5, 0.33,...] [0.33,...]?,...] loss 1.25347 loss 1.25347 Lecture 3 - 47 Fei-Fei Li, Ehsan Adeli, Zane Durante April 9, 2024

gradient dW:

W + h (third dim):

current W:



W + h (third dim): current W: gradient dW: [0.34,[0.34,[-2.5, -1.11, -1.11, 0.6, 0.78, 0.78 + 0.00010, 0.12, 0.12, 0.55, 0.55, **Numeric Gradient** 2.81, 2.81, Slow! Need to loop over -3.1, -3.1, all dimensions -1.5, -1.5, Approximate 0.33,...[0.33,...]*?*,...| loss 1.25347 loss 1.25347 April 9, 2024 Fei-Fei Li, Ehsan Adeli, Zane Durante Lecture 3 - 49

This is silly. The loss is just a function of W:

$$egin{aligned} L &= rac{1}{N} \sum_{i=1}^{N} L_i + \sum_k W_k^2 \ L_i &= \sum_{j
eq y_i} \max(0, s_j - s_{y_i} + 1) \ s &= f(x; W) = Wx \end{aligned}$$

want $\nabla_W L$

This is silly. The loss is just a function of W:

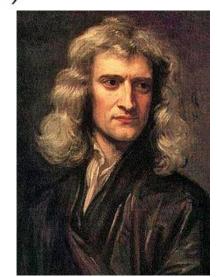
$$L = rac{1}{N} \sum_{i=1}^N L_i + \sum_k W_k^2$$

$$L_i = \sum_{j
eq y_i} \max(0, s_j - s_{y_i} + 1)$$

$$s = f(x; W) = Wx$$

want $\nabla_W L$

Use calculus to compute an analytic gradient





This image is in the public

This image is in the public

[0.34,[-2.5, $dW = \dots$ -1.11, 0.6, (some function 0.78, 0, data and W) 0.12, 0.2, 0.55, 0.7, 2.81, -0.5, -3.1, 1.1, -1.5, 1.3, 0.33,...] -2.1,...] loss 1.25347 Fei-Fei Li, Ehsan Adeli, Zane Durante Lecture 3 - 52 April 9, 2024

gradient dW:

current W:

In summary:

- Numerical gradient: approximate, slow, easy to write
- Analytic gradient: exact, fast, error-prone

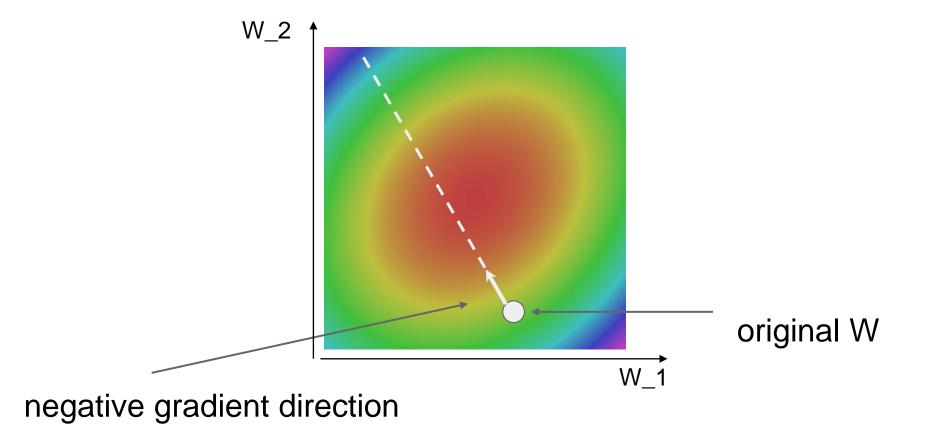
=>

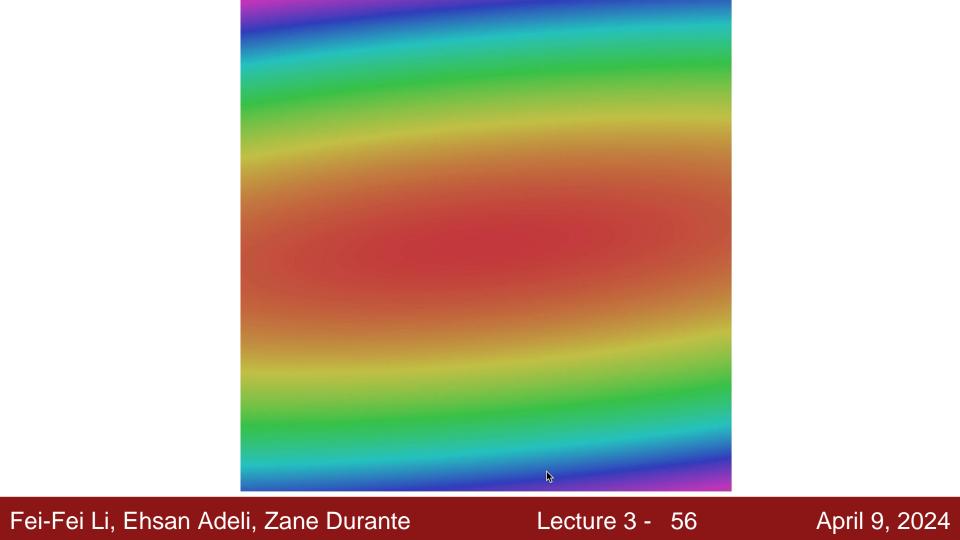
<u>In practice:</u> Always use analytic gradient, but check implementation with numerical gradient. This is called a **gradient check.**

Gradient Descent

```
# Vanilla Gradient Descent

while True:
    weights_grad = evaluate_gradient(loss_fun, data, weights)
    weights += - step_size * weights_grad # perform parameter update
```





Stochastic Gradient Descent (SGD)

$$L(W) = \frac{1}{N} \sum_{i=1}^{N} L_i(x_i, y_i, W) + \lambda R(W)$$

$$\nabla_W L(W) = \frac{1}{N} \sum_{i=1}^{N} \nabla_W L_i(x_i, y_i, W) + \lambda \nabla_W R(W)$$

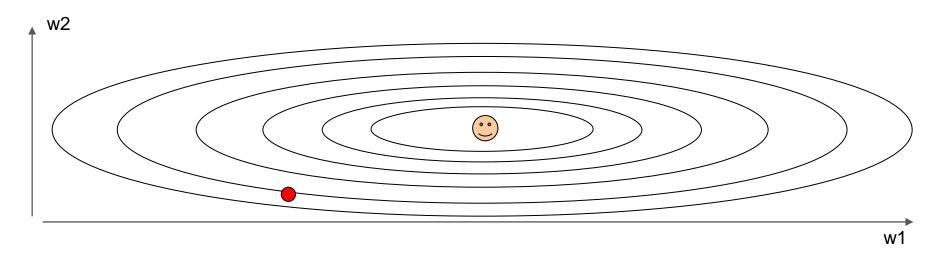
Full sum expensive when N is large!

Approximate sum using a **minibatch** of examples 32 / 64 / 128 common

```
# Vanilla Minibatch Gradient Descent

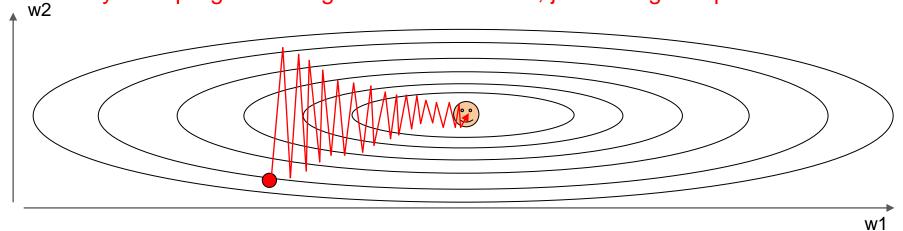
while True:
   data_batch = sample_training_data(data, 256) # sample 256 examples
   weights_grad = evaluate_gradient(loss_fun, data_batch, weights)
   weights += - step_size * weights_grad # perform parameter update
```

What if loss changes quickly in one direction and slowly in another? What does gradient descent do?



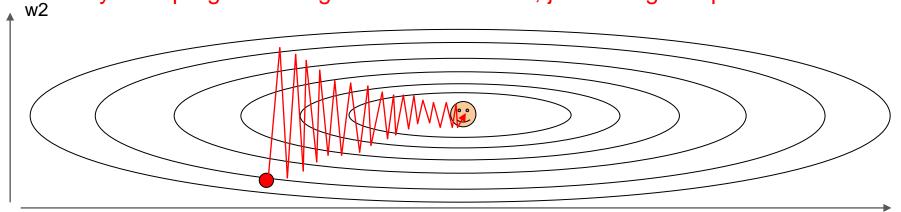
What if loss changes quickly in one direction and slowly in another? What does gradient descent do?

Very slow progress along shallow dimension, jitter along steep direction



What if loss changes quickly in one direction and slowly in another? What does gradient descent do?

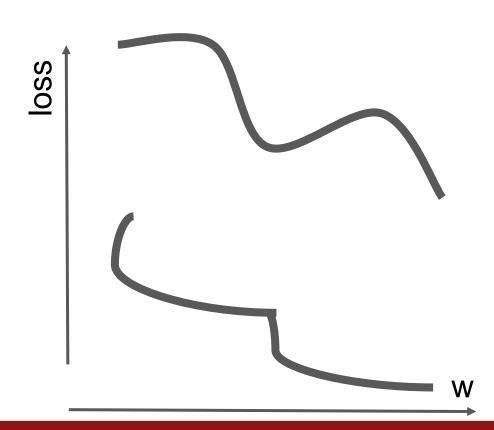
Very slow progress along shallow dimension, jitter along steep direction



Aside: Loss function has high **condition number**: ratio of largest to smallest singular value of the Hessian matrix is large

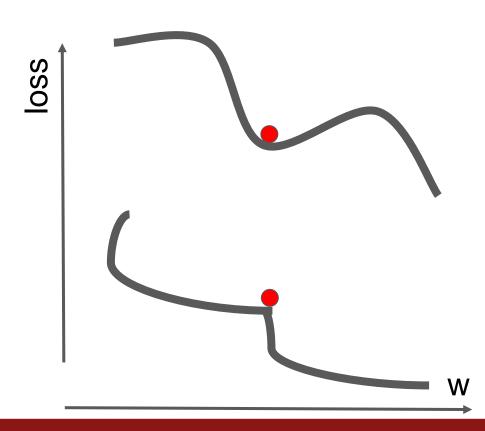
w1

What if the loss function has a local minima or saddle point?



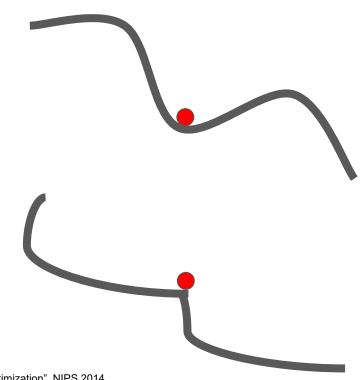
What if the loss function has a local minima or saddle point?

Zero gradient, gradient descent gets stuck



What if the loss function has a local minima or saddle point?

Saddle points much more common in high dimension



Dauphin et al, "Identifying and attacking the saddle point problem in high-dimensional non-convex optimization", NIPS 2014

saddle point in two dimension

$$f(x,y) = x^2 - y^2$$

$$rac{\partial}{\partial x}(x^2-y^2)=2x
ightarrow 2(0)=0$$

$$rac{\partial}{\partial oldsymbol{y}}(x^2-oldsymbol{y}^2)=-2y
ightarrow -2({\color{red}0})=0$$

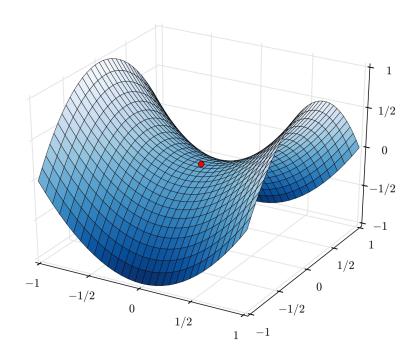
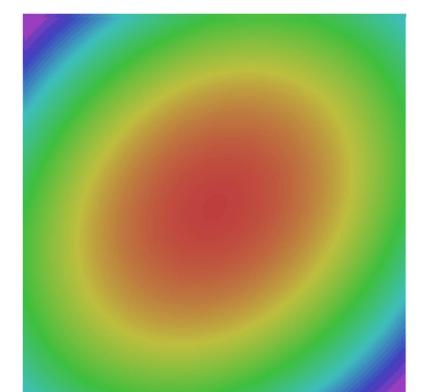


Image source: https://en.wikipedia.org/wiki/Saddle_point

Our gradients come from minibatches so they can be noisy!

$$L(W) = \frac{1}{N} \sum_{i=1}^{N} L_i(x_i, y_i, W)$$

$$\nabla_W L(W) = \frac{1}{N} \sum_{i=1}^{N} \nabla_W L_i(x_i, y_i, W)$$

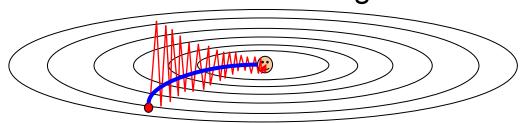


SGD + Momentum

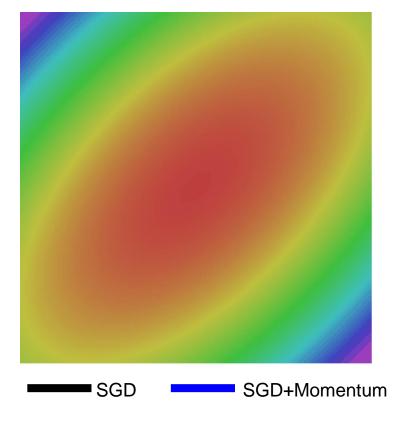
Local Minima Saddle points



Poor Conditioning



Gradient Noise



SGD: the simple two line update code

SGD

```
x_{t+1} = x_t - \alpha \nabla f(x_t)
```

```
while True:
    dx = compute_gradient(x)
    x -= learning_rate * dx
```

SGD + Momentum:

continue moving in the general direction as the previous iterations

SGD

$$x_{t+1} = x_t - \alpha \nabla f(x_t)$$

```
while True:
    dx = compute_gradient(x)
    x -= learning_rate * dx
```

SGD+Momentum

```
v_{t+1} = \rho v_t + \nabla f(x_t)x_{t+1} = x_t - \alpha v_{t+1}
```

- Build up "velocity" as a running mean of gradients
- Rho gives "friction"; typically rho=0.9 or 0.99

Sutskever et al, "On the importance of initialization and momentum in deep learning", ICML 2013

SGD + Momentum:

continue moving in the general direction as the previous iterations

SGD

```
x_{t+1} = x_t - \alpha \nabla f(x_t)
```

```
while True:
    dx = compute_gradient(x)
    x -= learning_rate * dx
```

SGD+Momentum

 $v_{t+1} = \rho v_t + \nabla f(x_t)$

```
x_{t+1} = x_t - \alpha v_{t+1}

vx = 0

while True:
 dx = compute_gradient(x)
```

x -= learning_rate * vx

vx = rho * vx + dx

- Build up "velocity" as a running mean of gradients
- Rho gives "friction"; typically rho=0.9 or 0.99

Sutskever et al, "On the importance of initialization and momentum in deep learning", ICML 2013

SGD + Momentum:

alternative equivalent formulation

SGD+Momentum

```
v_{t+1} = \rho v_t - \alpha \nabla f(x_t)x_{t+1} = x_t + v_{t+1}
```

```
vx = 0
while True:
    dx = compute_gradient(x)
    vx = rho * vx - learning_rate * dx
    x += vx
```

SGD+Momentum

```
v_{t+1} = \rho v_t + \nabla f(x_t)
x_{t+1} = x_t - \alpha v_{t+1}
v_t = 0
```

```
while True:
    dx = compute_gradient(x)
    vx = rho * vx + dx
    x -= learning_rate * vx
```

You may see SGD+Momentum formulated different ways, but they are equivalent - give same sequence of x

Sutskever et al, "On the importance of initialization and momentum in deep learning", ICML 2013

More Complex Optimizers: RMSProp

```
SGD +
Momentum
```

```
vx = 0
while True:
    dx = compute_gradient(x)
    vx = rho * vx + dx
    x -= learning_rate * vx
```

Adds element-wise scaling of the gradient based on the historical sum of squares in each dimension (with decay)

RMSProp

```
grad_squared = 0
while True:
    dx = compute_gradient(x)

    grad_squared = decay_rate * grad_squared + (1 - decay_rate) * dx * dx
    x -= learning_rate * dx / (np.sqrt(grad_squared) + 1e-7)
```

More Complex Optimizers: RMSProp

SGD + Momentum

```
vx = 0
while True:
    dx = compute_gradient(x)
    vx = rho * vx + dx
    x -= learning_rate * vx
```

"Per-parameter learning rates" or "adaptive learning rates"

RMSProp

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grad_squared = 0
while True:
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    grad_squared = decay_rate * grad_squared + (1 - decay_rate) * dx * dx
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RMSProp

RMSProp

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    grad_squared = decay_rate * grad_squared + (1 - decay_rate) * dx * dx
    x -= learning_rate * dx / (np.sqrt(grad_squared) + 1e-7)
```

Q: What happens with RMSProp?

RMSProp

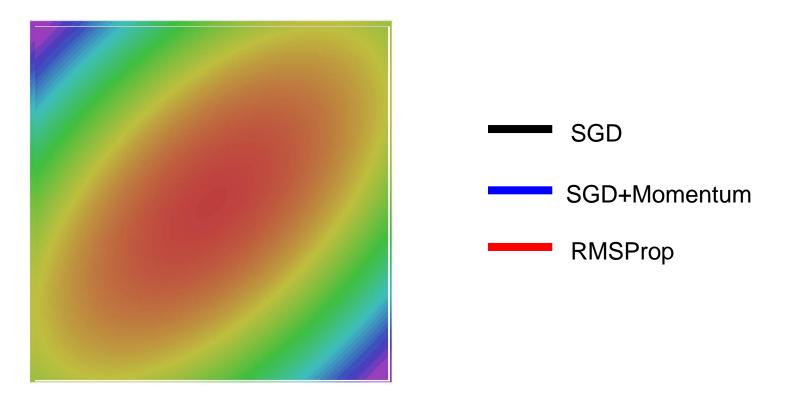
RMSProp

```
grad_squared = 0
while True:
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    x -= learning_rate * dx / (np.sqrt(grad_squared) + 1e-7)
```

Q: What happens with RMSProp?

Progress along "steep" directions is damped; progress along "flat" directions is accelerated

RMSProp



Optimizers: Adam (almost)

```
first_moment = 0
second_moment = 0
while True:
    dx = compute_gradient(x)
    first_moment = beta1 * first_moment + (1 - beta1) * dx
    second_moment = beta2 * second_moment + (1 - beta2) * dx * dx
    x -= learning_rate * first_moment / (np.sqrt(second_moment) + 1e-7))
```

Kingma and Ba, "Adam: A method for stochastic optimization", ICLR 2015

Adam (almost)

```
first_moment = 0
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while True:
    dx = compute_gradient(x)
    first_moment = beta1 * first_moment + (1 - beta1) * dx
    second_moment = beta2 * second_moment + (1 - beta2) * dx * dx
    x -= learning_rate * first_moment / (np.sqrt(second_moment) + 1e-7))
Momentum
RMSProp
```

Sort of like RMSProp with momentum

Q: What happens at first timestep?

Kingma and Ba, "Adam: A method for stochastic optimization", ICLR 2015

Adam (full form)

```
first_moment = 0
second_moment = 0
for t in range(1, num_iterations):
    dx = compute_gradient(x)
    first_moment = beta1 * first_moment + (1 - beta1) * dx
    second_moment = beta2 * second_moment + (1 - beta2) * dx * dx

first_unbias = first_moment / (1 - beta1 ** t)
    second_unbias = second_moment / (1 - beta2 ** t)

x -= learning_rate * first_unbias / (np.sqrt(second_unbias) + 1e-7))

AdaGrad / RMSProp
```

Bias correction for the fact that first and second moment estimates start at zero

Kingma and Ba, "Adam: A method for stochastic optimization", ICLR 2015

Adam (full form)

```
first_moment = 0
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for t in range(1, num_iterations):
    dx = compute_gradient(x)
    first_moment = beta1 * first_moment + (1 - beta1) * dx

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    x -= learning_rate * first_unbias / (np.sqrt(second_unbias) + 1e-7))

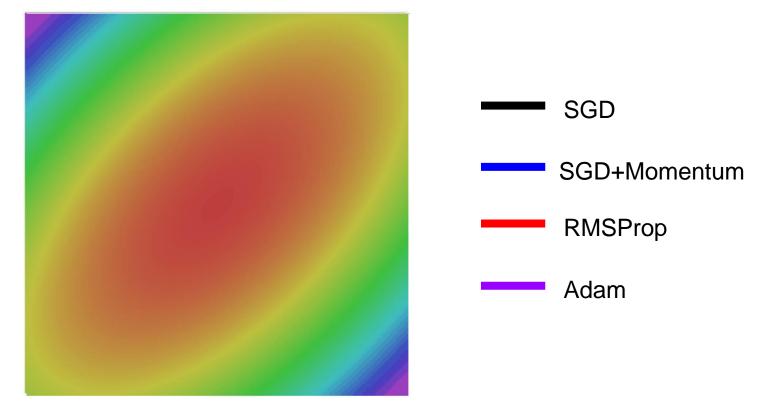
AdaGrad / RMSProp
```

Bias correction for the fact that first and second moment estimates start at zero

Adam with beta1 = 0.9, beta2 = 0.999, and learning_rate = 1e-3 or 5e-4 is a great starting point for many models!

Kingma and Ba, "Adam: A method for stochastic optimization", ICLR 2015

Adam



Q: How does regularization interact with the optimizer? (e.g., L2)

```
first_moment = 0
second_moment = 0
for t in range(1, num_iterations):
    dx = compute_gradient(x)
    first_moment = beta1 * first_moment + (1 - beta1) * dx
    second_moment = beta2 * second_moment + (1 - beta2) * dx * dx
    first_unbias = first_moment / (1 - beta1 ** t)
    second_unbias = second_moment / (1 - beta2 ** t)
    x -= learning_rate * first_unbias / (np.sqrt(second_unbias) + 1e-7))
```

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    second_unbias = second_moment / (1 - beta2 ** t)
    x -= learning_rate * first_unbias / (np.sqrt(second_unbias) + 1e-7))
```

A: It depends!

Q: How does regularization interact with the optimizer? (e.g., L2)

```
first_moment = 0
second_moment = 0
for t in range(1, num_iterations):
    dx = compute_gradient(x)
    first_moment = beta1 * first_moment + (1 - beta1) * dx
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    first_unbias = first_moment / (1 - beta1 ** t)
    second_unbias = second_moment / (1 - beta2 ** t)
    x -= learning_rate * first_unbias / (np.sqrt(second_unbias) + 1e-7))
```

Used during moment calculations!

Q: How does regularization interact with the optimizer? (e.g., L2)

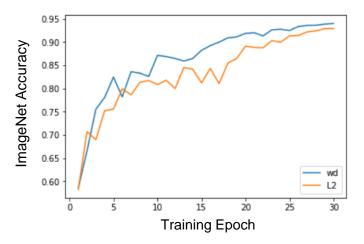
```
first_moment = 0
second moment = 0
for t in range(1, num_iterations):
  dx = compute\_gradient(x)
  first_moment = beta1 * first_moment + (1 - beta1) * dx
  second_moment = beta2 * second_moment + (1 - beta2) * dx * dx
  first_unbias = first_moment / (1 - beta1 ** t)
  second_unbias = second_moment / (1 - beta2 ** t)
  x -= learning_rate * first_unbias / (np.sgrt(second_unbias) + 1e-7))
```

AdamW (Weight Decay) adds term here

Computed after the moments!

Q: How does regularization interact with the optimizer? (e.g., L2)

```
Standard Adam computes L2 here
first_moment = 0
second moment = 0
for t in range(1, num_iterations):
  dx = compute_gradient(x) 
  first_moment = beta1 * first_moment + (1 - beta1) * dx
  second_moment = beta2 * second_moment + (1 - beta2) * dx * dx
  first_unbias = first_moment / (1 - beta1 ** t)
  second_unbias = second_moment / (1 - beta2 ** t)
  x -= learning_rate * first_unbias / (np.sqrt(second_unbias) + 1e-7)),
                        AdamW (Weight Decay) adds term here
```



Source: https://www.fast.ai/posts/2018-07-02adam-weight-decay.html

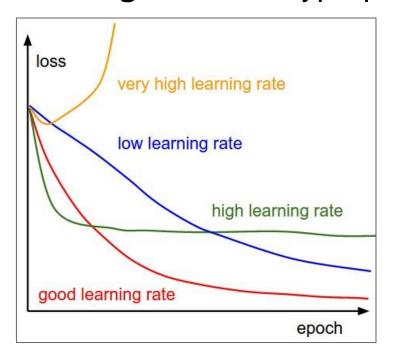
Learning rate schedules

```
# Vanilla Gradient Descent

while True:
    weights_grad = evaluate_gradient(loss_fun, data, weights)
    weights += - step_size * weights_grad # perform parameter update

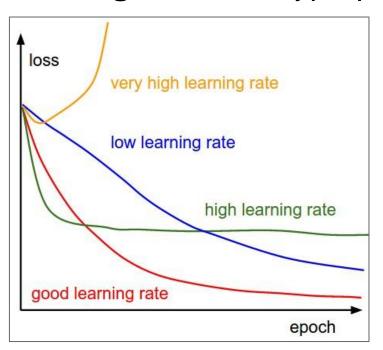
    Learning rate
```

SGD, SGD+Momentum, RMSProp, Adam, AdamW all have learning rate as a hyperparameter.



Q: Which one of these learning rates is best to use?

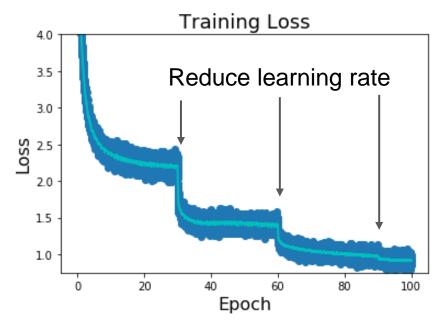
SGD, SGD+Momentum, RMSProp, Adam, AdamW all have learning rate as a hyperparameter.



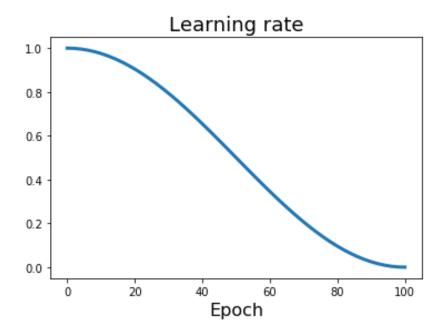
Q: Which one of these learning rates is best to use?

A: In reality, all of these could be good learning rates.

Learning rate decays over time



Step: Reduce learning rate at a few fixed points. E.g. for ResNets, multiply LR by 0.1 after epochs 30, 60, and 90.



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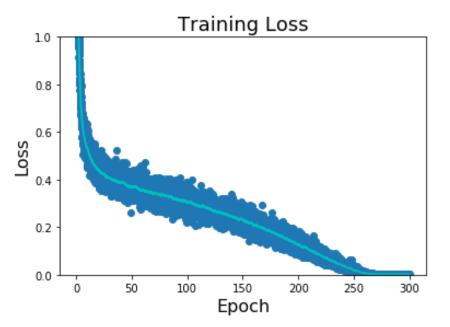
Cosine:
$$\alpha_t = \frac{1}{2}\alpha_0 \left(1 + \cos(t\pi/T)\right)$$

Loshchilov and Hutter, "SGDR: Stochastic Gradient Descent with Warm Restarts", ICLR 2017 Radford et al, "Improving Language Understanding by Generative Pre-Training", 2018 Feichtenhofer et al, "SlowFast Networks for Video Recognition", arXiv 2018 Child at al, "Generating Long Sequences with Sparse Transformers", arXiv 2019

 $lpha_0$: Initial learning rate

 $lpha_t$: Learning rate at epoch t

T: Total number of epochs



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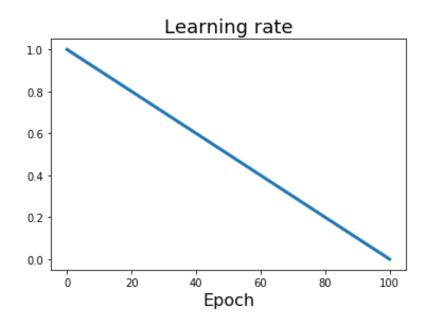
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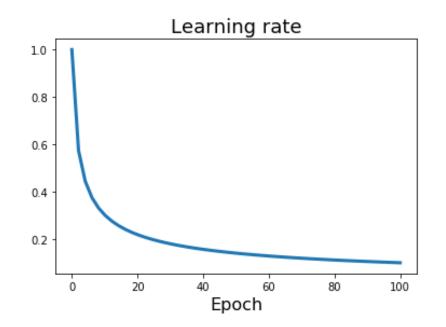
Linear:
$$\alpha_t = \alpha_0(1 - t/T)$$

 $lpha_0$: Initial learning rate

 $lpha_t$: Learning rate at epoch t

T: Total number of epochs

Devlin et al, "BERT: Pre-training of Deep Bidirectional Transformers for Language Understanding", 2018



Step: Reduce learning rate at a few fixed points. E.g. for ResNets, multiply LR by 0.1 after epochs 30, 60, and 90.

Cosine:
$$\alpha_t = \frac{1}{2}\alpha_0 \left(1 + \cos(t\pi/T)\right)$$

Linear:
$$\alpha_t = \alpha_0(1 - t/T)$$

Inverse sqrt:
$$\alpha_t = \alpha_0/\sqrt{t}$$

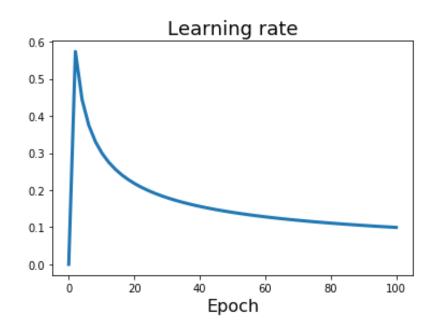
 $lpha_0$: Initial learning rate

 $lpha_t$: Learning rate at epoch t

T: Total number of epochs

Vaswani et al, "Attention is all you need", NIPS 2017

Learning Rate Decay: Linear Warmup

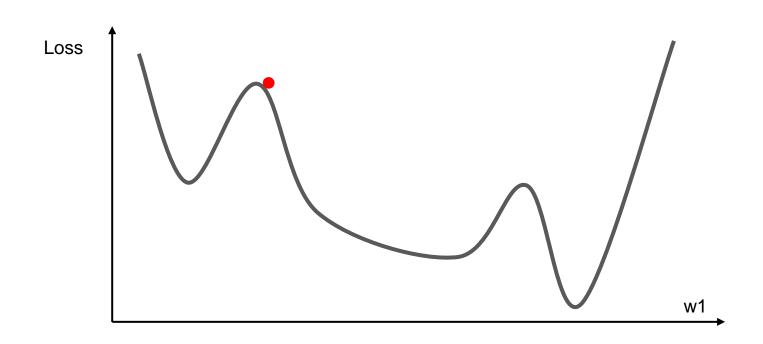


High initial learning rates can make loss explode; linearly increasing learning rate from 0 over the first ~5,000 iterations can prevent this.

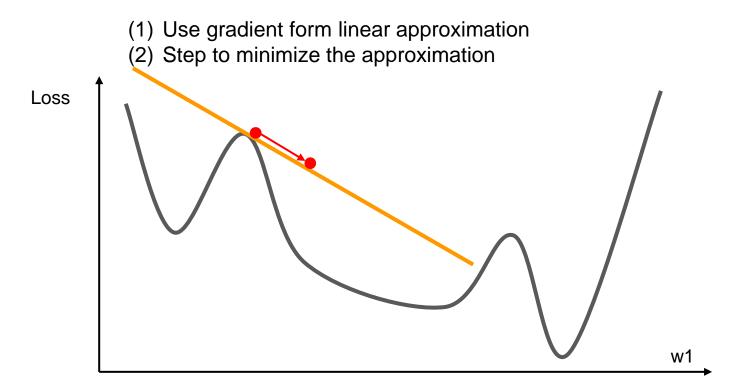
Empirical rule of thumb: If you increase the batch size by N, also scale the initial learning rate by N

Goyal et al, "Accurate, Large Minibatch SGD: Training ImageNet in 1 Hour", arXiv 2017

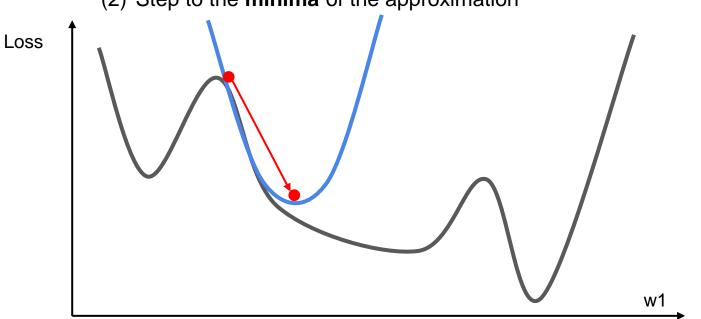
First-Order Optimization



First-Order Optimization



- (1) Use gradient and Hessian to form quadratic approximation
- (2) Step to the **minima** of the approximation



second-order Taylor expansion:

$$J(\boldsymbol{\theta}) \approx J(\boldsymbol{\theta}_0) + (\boldsymbol{\theta} - \boldsymbol{\theta}_0)^{\top} \nabla_{\boldsymbol{\theta}} J(\boldsymbol{\theta}_0) + \frac{1}{2} (\boldsymbol{\theta} - \boldsymbol{\theta}_0)^{\top} \boldsymbol{H} (\boldsymbol{\theta} - \boldsymbol{\theta}_0)$$

Solving for the critical point we obtain the Newton parameter update:

$$\boldsymbol{\theta}^* = \boldsymbol{\theta}_0 - \boldsymbol{H}^{-1} \nabla_{\boldsymbol{\theta}} J(\boldsymbol{\theta}_0)$$

Q: Why is this bad for deep learning?

second-order Taylor expansion:

$$J(\boldsymbol{\theta}) \approx J(\boldsymbol{\theta}_0) + (\boldsymbol{\theta} - \boldsymbol{\theta}_0)^{\top} \nabla_{\boldsymbol{\theta}} J(\boldsymbol{\theta}_0) + \frac{1}{2} (\boldsymbol{\theta} - \boldsymbol{\theta}_0)^{\top} \boldsymbol{H} (\boldsymbol{\theta} - \boldsymbol{\theta}_0)$$

Solving for the critical point we obtain the Newton parameter update:

$$\boldsymbol{\theta}^* = \boldsymbol{\theta}_0 - \boldsymbol{H}^{-1} \nabla_{\boldsymbol{\theta}} J(\boldsymbol{\theta}_0)$$

Hessian has O(N^2) elements Inverting takes O(N^3) N = (Tens or Hundreds of) Millions

Q: Why is this bad for deep learning?

In practice:

- Adam(W) is a good default choice in many cases; it often works ok even with constant learning rate
- SGD+Momentum can outperform Adam but may require more tuning of LR and schedule
- If you can afford to do full batch updates then look beyond 1st order optimization (2nd order and beyond)

Looking Ahead: How to optimize more complex functions?

(Currently) Linear score function: f=Wx

$$x \in \mathbb{R}^D, W \in \mathbb{R}^{C \times D}$$

Neural networks: 2 layers

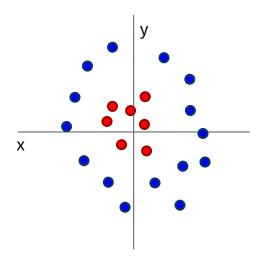
(Currently) Linear score function: f=Wx

(Next Class) 2-layer Neural Network $\,f = W_2 \max(0, W_1 x)\,$

$$x \in \mathbb{R}^D, W_1 \in \mathbb{R}^{H \times D}, W_2 \in \mathbb{R}^{C \times H}$$

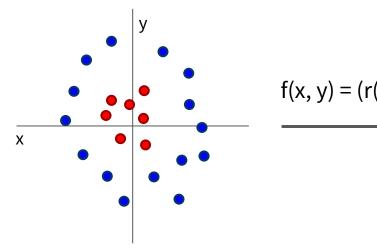
(In practice we will usually add a learnable bias at each layer as well)

Why do we want non-linearity?



Cannot separate red and blue points with linear classifier

Why do we want non-linearity?



 $f(x, y) = (r(x, y), \theta(x, y))$

After applying feature

Cannot separate red and blue points with linear classifier

After applying feature transform, points can be separated by linear classifier

Neural networks: also called fully connected network

(Currently) Linear score function: f=Wx

(Next Class) 2-layer Neural Network $f = W_2 \max(0, W_1 x)$

$$x \in \mathbb{R}^D, W_1 \in \mathbb{R}^{H \times D}, W_2 \in \mathbb{R}^{C \times H}$$

"Neural Network" is a very broad term; these are more accurately called "fully-connected networks" or sometimes "multi-layer perceptrons" (MLP)

(In practice we will usually add a learnable bias at each layer as well)

Next time:

Introduction to neural networks

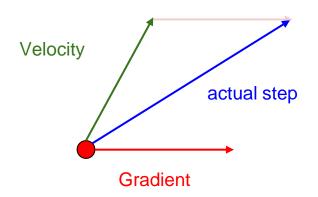
Backpropagation (How do you calculate dx for neural nets?)

Appendix Material

Extraneous content that may be of interest

SGD+Momentum

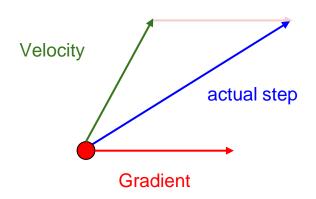
Momentum update:



Combine gradient at current point with velocity to get step used to update weights

Nesterov, "A method of solving a convex programming problem with convergence rate O(1/k^2)", 1983 Nesterov, "Introductory lectures on convex optimization: a basic course", 2004 Sutskever et al, "On the importance of initialization and momentum in deep learning", ICML 2013

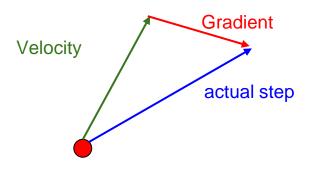
Momentum update:



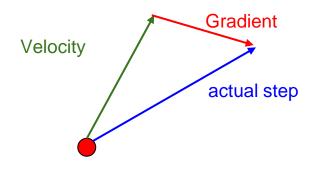
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Nesterov Momentum

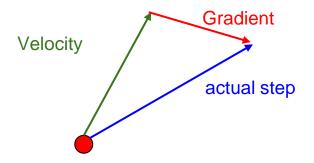


$$v_{t+1} = \rho v_t - \alpha \nabla f(x_t + \rho v_t)$$
$$x_{t+1} = x_t + v_{t+1}$$



$$v_{t+1} = \rho v_t - \alpha \nabla f(x_t + \rho v_t)$$
$$x_{t+1} = x_t + v_{t+1}$$

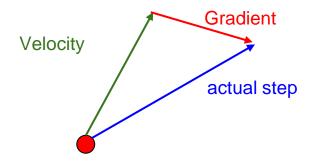
Annoying, usually we want update in terms of $x_t, \nabla f(x_t)$



$$v_{t+1} = \rho v_t - \alpha \nabla f(x_t + \rho v_t)$$
$$x_{t+1} = x_t + v_{t+1}$$

Change of variables $\tilde{x}_t = x_t + \rho v_t$ and rearrange:

Annoying, usually we want update in terms of $x_t, \nabla f(x_t)$



$$v_{t+1} = \rho v_t - \alpha \nabla f(x_t + \rho v_t)$$
$$x_{t+1} = x_t + v_{t+1}$$

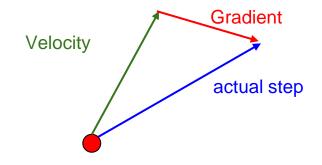
Change of variables $\tilde{x}_t = x_t + \rho v_t$ and rearrange:

$$v_{t+1} = \rho v_t - \alpha \nabla f(\tilde{x}_t)$$

$$\tilde{x}_{t+1} = \tilde{x}_t - \rho v_t + (1+\rho)v_{t+1}$$

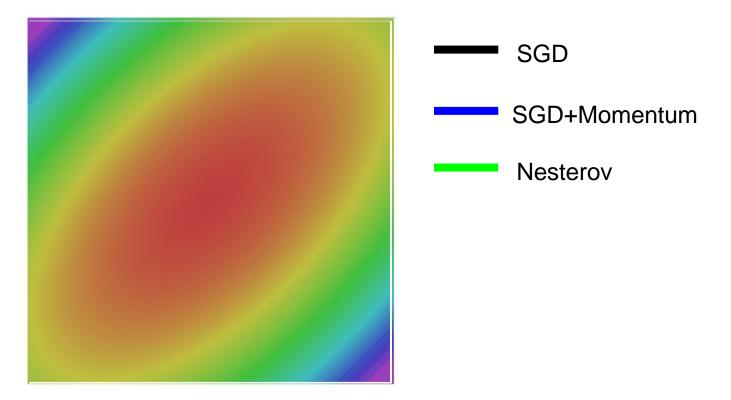
$$= \tilde{x}_t + v_{t+1} + \rho(v_{t+1} - v_t)$$

Annoying, usually we want update in terms of $x_t, \nabla f(x_t)$



"Look ahead" to the point where updating using velocity would take us; compute gradient there and mix it with velocity to get actual update direction

https://cs231n.github.io/neural-networks-3/



```
grad_squared = 0
while True:
    dx = compute_gradient(x)
    grad_squared += dx * dx
x -= learning_rate * dx / (np.sqrt(grad_squared) + 1e-7)
```

Added element-wise scaling of the gradient based on the historical sum of squares in each dimension

"Per-parameter learning rates" or "adaptive learning rates"

Duchi et al, "Adaptive subgradient methods for online learning and stochastic optimization", JMLR 2011

```
grad_squared = 0
while True:
  dx = compute\_gradient(x)
  grad_squared += dx * dx
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```

Q: What happens with AdaGrad?

```
grad_squared = 0
while True:
  dx = compute\_gradient(x)
  grad_squared += dx * dx
 x -= learning_rate * dx / (np.sqrt(grad_squared) + 1e-7)
```

Q: What happens with AdaGrad?

Progress along "steep" directions is damped; progress along "flat" directions is accelerated

```
grad_squared = 0
while True:
  dx = compute\_gradient(x)
  grad_squared += dx * dx
 x -= learning_rate * dx / (np.sqrt(grad_squared) + 1e-7)
```

Q2: What happens to the step size over long time?

```
grad_squared = 0
while True:
  dx = compute\_gradient(x)
  grad_squared += dx * dx
 x -= learning_rate * dx / (np.sqrt(grad_squared) + 1e-7)
```

Q2: What happens to the step size over long time? Decays to zero

RMSProp: "Leaky AdaGrad"

AdaGrad

```
grad_squared = 0
while True:
    dx = compute_gradient(x)
    grad_squared += dx * dx
    x -= learning_rate * dx / (np.sqrt(grad_squared) + 1e-7)
```

RMSProp

```
grad_squared = 0
while True:
    dx = compute_gradient(x)
    grad_squared = decay_rate * grad_squared + (1 - decay_rate) * dx * dx
    x -= learning_rate * dx / (np.sqrt(grad_squared) + 1e-7)
```

Tieleman and Hinton, 2012

$$\boldsymbol{\theta}^* = \boldsymbol{\theta}_0 - \boldsymbol{H}^{-1} \nabla_{\boldsymbol{\theta}} J(\boldsymbol{\theta}_0)$$

- Quasi-Newton methods (BGFS most popular):
 instead of inverting the Hessian (O(n^3)), approximate
 inverse Hessian with rank 1 updates over time (O(n^2)
 each).
- L-BFGS (Limited memory BFGS):
 Does not form/store the full inverse Hessian.

L-BFGS

- Usually works very well in full batch, deterministic mode
 i.e. if you have a single, deterministic f(x) then L-BFGS will
 probably work very nicely
- **Does not transfer very well to mini-batch setting**. Gives bad results. Adapting second-order methods to large-scale, stochastic setting is an active area of research.

Le et al, "On optimization methods for deep learning, ICML 2011"

Ba et al, "Distributed second-order optimization using Kronecker-factored approximations", ICLR 2017

In practice:

- Adam is a good default choice in many cases; it often works ok even with constant learning rate
- **SGD+Momentum** can outperform Adam but may require more tuning of LR and schedule
- If you can afford to do full batch updates then try out **L-BFGS** (and don't forget to disable all sources of noise)