Lecture 3: Regularization and Optimization

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Image Classification: A core task in Computer Vision



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

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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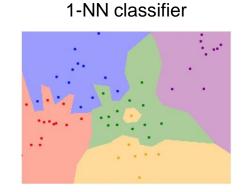
Intraclass Variation

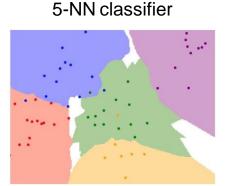


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Recall from last time: data-driven approach, kNN

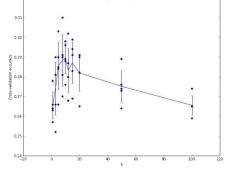






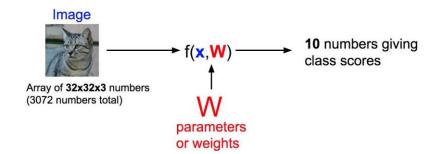


train	validation	test
train	valluation	ເຮວເ





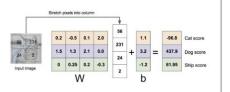
Recall from last time: Linear Classifier



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

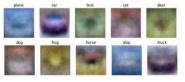
Algebraic Viewpoint

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



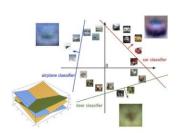
Visual Viewpoint

One template per class



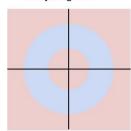
Geometric Viewpoint

Hyperplanes cutting up space



Class 1: 1 <= L2 norm <= 2

Class 2: Everything else



Class 1: Three modes

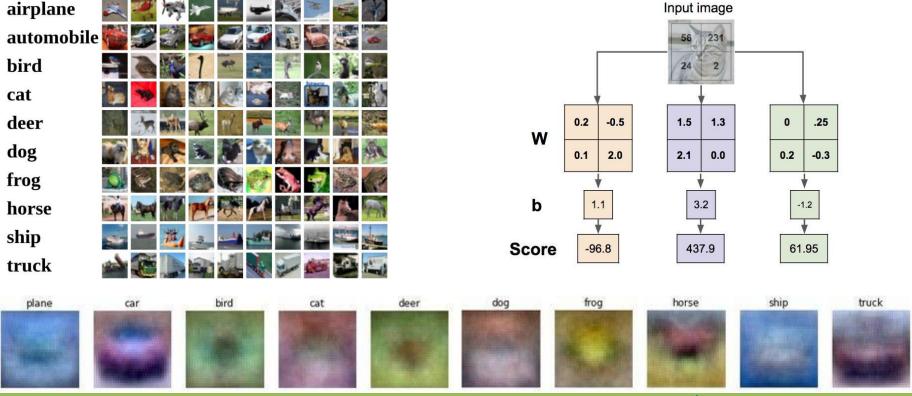
....

Class 2:

Everything else

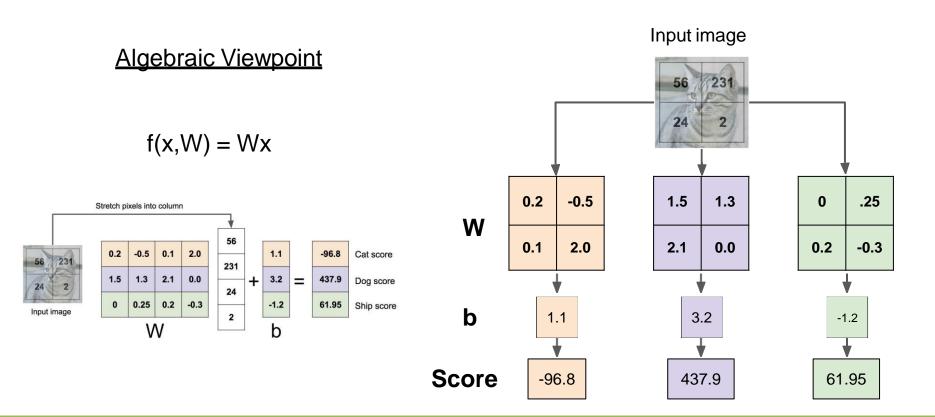


Interpreting a Linear Classifier: Visual Viewpoint



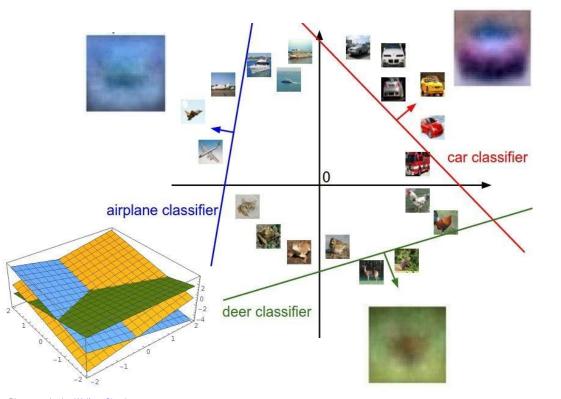


Example with an image with 4 pixels, and 3 classes (cat/dog/ship)





Interpreting a Linear Classifier: Geometric Viewpoint



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



Array of **32x32x3** numbers (3072 numbers total)

Plot created using Wolfram Cloud

Cat image by Nikita is licensed under CC-BY 2.0

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





A **loss function** tells how good our current classifier is

Given a dataset of examples

$$\{(x_i, y_i)\}_{i=1}^N$$

Where x_i is image and y_i is (integer) label

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

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

cat 3.2

1.3

2.2 2.5

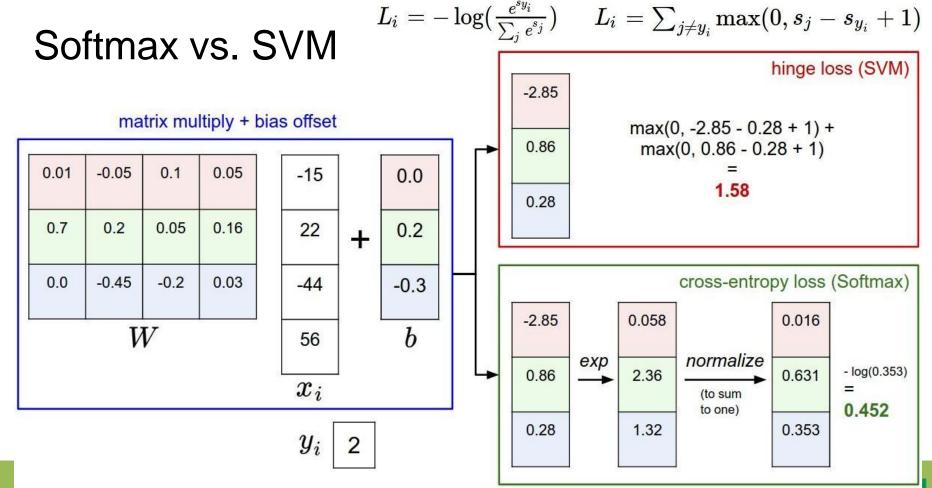
5.1 -1.7 frog

car

4.9 2.0

-3.1

Regularization and Optimization



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

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

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



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

E.g. 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

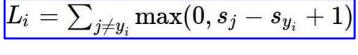
car



-1.7

2.0

2.5



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:

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

$$= \max(0, -6.2) + \max(0, -4.8)$$

$$= 0 + 0$$

$$= 0$$



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

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?

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

Data loss: Model predictions should match training data



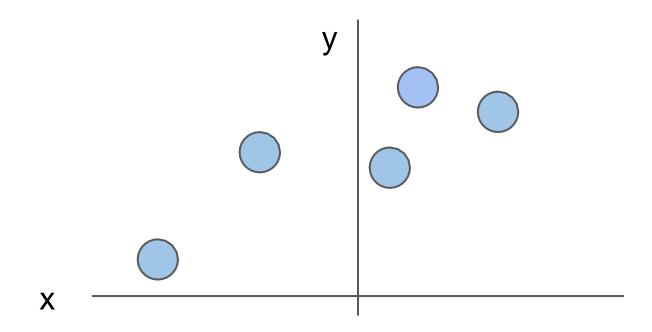
$$L(W) = \underbrace{\frac{1}{N} \sum_{i=1}^{N} L_i(f(x_i, W), y_i) + \lambda R(W)}_{i=1}$$

Data loss: Model predictions should match training data

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

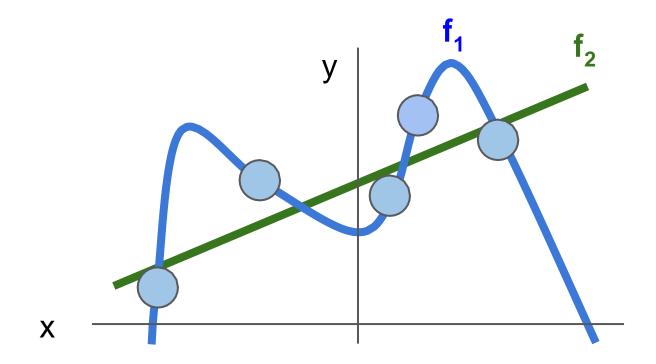


Regularization intuition: toy example training data



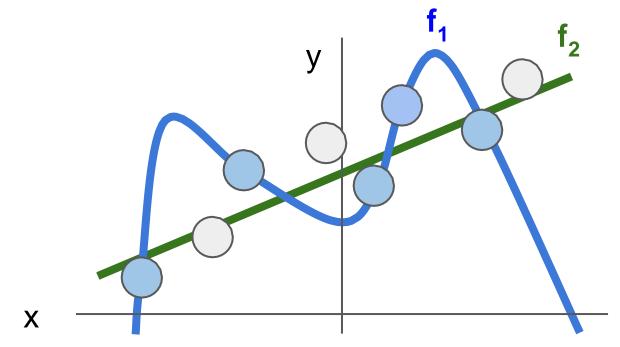


Regularization intuition: Prefer Simpler Models





Regularization: Prefer Simpler Models



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

Regularization and Optimization

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

Data loss: Model predictions should match training data

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

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



 λ = 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



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



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

More complex:

Dropout

Batch normalization

Stochastic depth, fractional pooling, etc





$$\lambda$$
 = 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

Why regularize?

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



Regularization: Expressing Preferences

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

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

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

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

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

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

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

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

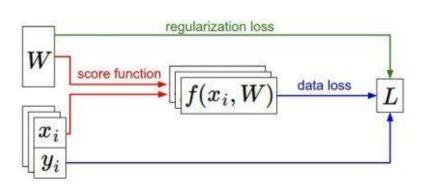
Which one would L1 regularization prefer?



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

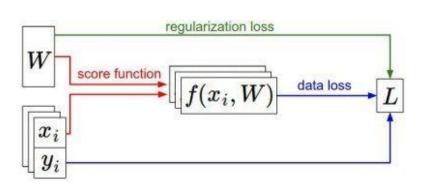


Recap

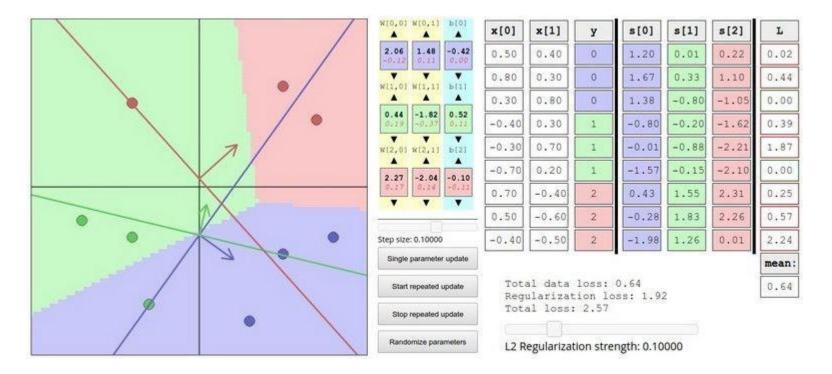
How do we find the best W?

- 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



Interactive Web Demo



http://vision.stanford.edu/teaching/cs231n-demos/linear-classify/



Optimization











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**

Regularization and Optimization

-1.11, 0.78, 0.12, 0.55, 2.81, -3.1, -1.5, 0.33,...] loss 1.25347 Regularization and Optimization

gradient dW:

current W:

[0.34,

[0.34 + 0.0001][0.34,-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,...?,...]

W + h (first dim):



gradient dW:

loss 1.25322

current W:

loss 1.25347

-1.11, -1.11, 0.78, 0.78, 0.12, 0.12, (1.25322 - 1.25347)/0.0001 0.55, 0.55, = -2.52.81, 2.81, $\frac{df(x)}{dx} = \lim \frac{f(x+h) - f(x)}{dx}$ -3.1, -3.1, -1.5, -1.5, [0.33,...]0.33,...?,...] loss 1.25347 loss 1.25322 Regularization and Optimization

gradient dW:

[-2.5,

W + h (first dim):

[0.34 + 0.0001,

current W:

[0.34,

[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,...?,...]

W + h (second dim):



gradient dW:

loss 1.25353

current W:

loss 1.25347

[0.34, [0.34,[-2.5, -1.11 + 0.0001-1.11, 0.6, 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

gradient dW:

W + h (second dim):

current W:

Regularization and Optimization

[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,...?,...]

W + h (third dim):



gradient dW:

loss 1.25347

current W:

loss 1.25347

W + h (third dim): gradient dW: [0.34,[0.34,[-2.5, -1.11, -1.11, 0.6, 0.78 + 0.00010.78, 0.12, 0.12, 0.55, 0.55, (1.25347 - 1.25347)/0.00012.81, 2.81, = 0-3.1, -3.1, $\frac{df(x)}{dx} = \lim_{x \to 0} \frac{f(x+h) - f(x)}{f(x+h)}$ -1.5, -1.5, [0.33,...]0.33,...loss 1.25347 loss 1.25347

current W:

Regularization and Optimization

[-2.5, -1.11, -1.11, 0.6, 0.78 + 0.00010.78, 0, 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 Regularization and Optimization

gradient dW:

W + h (third dim):

[0.34,

current W:

[0.34,

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 $\nabla_W L$



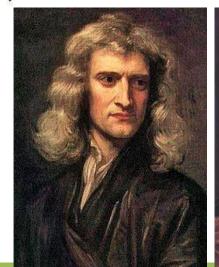
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eq y_i} \max(0, s_j - s_{y_i} + 1) \end{aligned}$$

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

want $\nabla_W L$

Use calculus to compute an analytic gradient



is in the public domain



-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 Regularization and Optimization

dW = ...

gradient dW:

[-2.5,

current W:

[0.34,

In summary:

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

=>

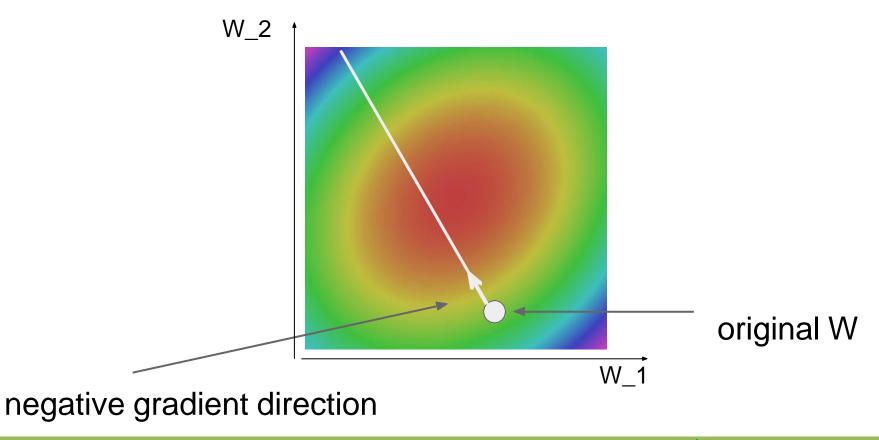
<u>In practice:</u> Always use analytic gradient, but check implementation with numerical gradient. This is called a **gradient check.**



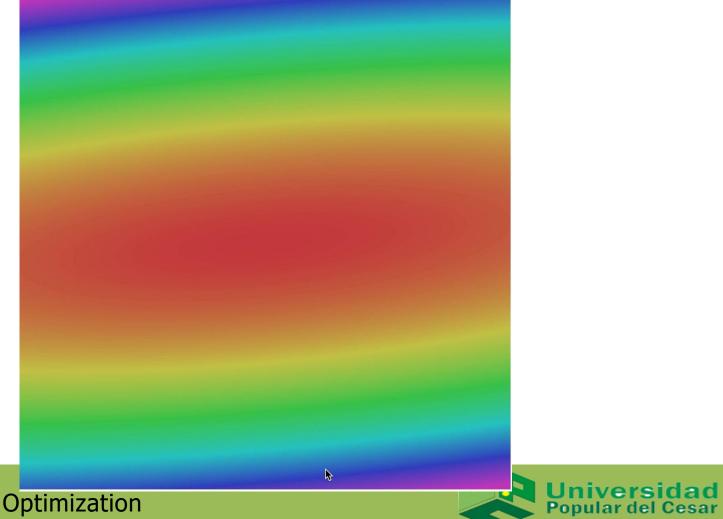
Gradient Descent

```
# 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!

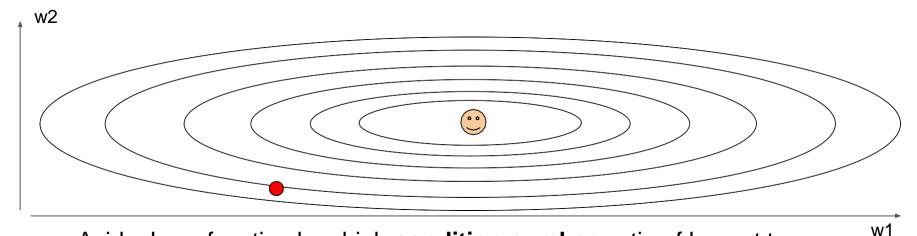
Approximate sum using a **minibatch** of examples 32 / 64 / 128 common

```
# 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?

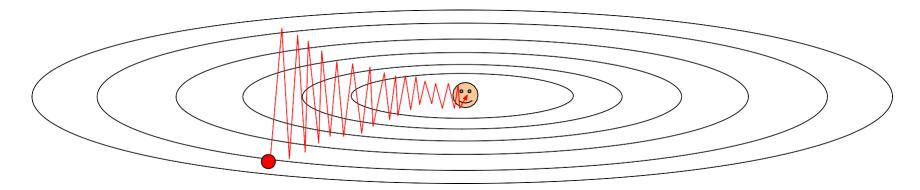


Aside: Loss function has high **condition number**: ratio of largest to smallest singular value of the Hessian matrix is large



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

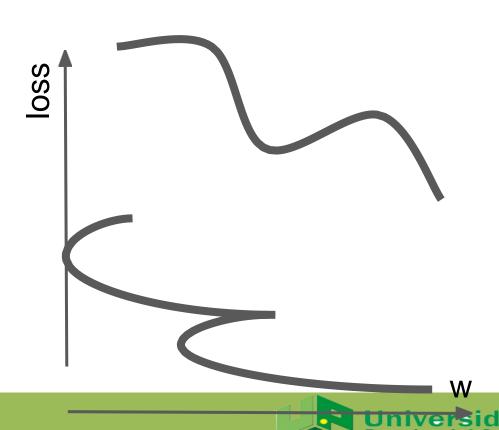
Very slow progress along shallow dimension, jitter along steep direction



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

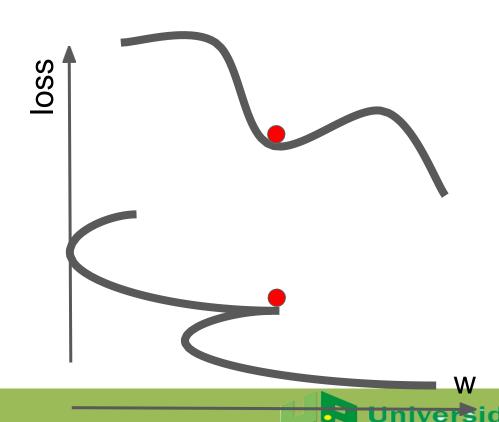


What if the loss function has a local minima or saddle point?



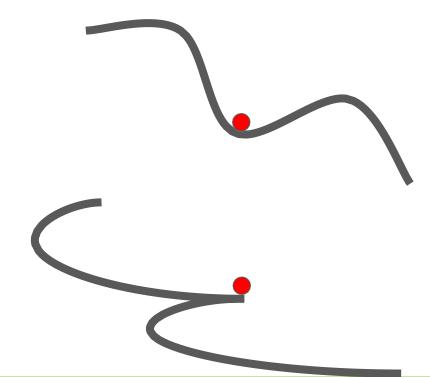
What if the loss function has a local minima or saddle point?

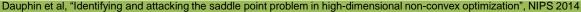
Zero gradient, gradient descent gets stuck



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

Saddle points much more common in high dimension





saddle point in two dimension

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

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

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

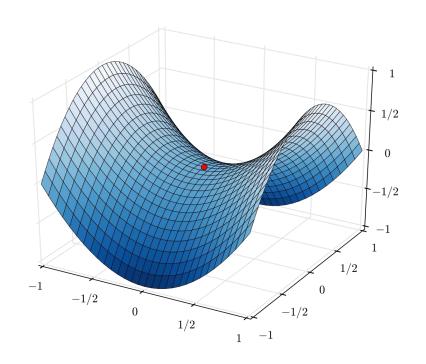


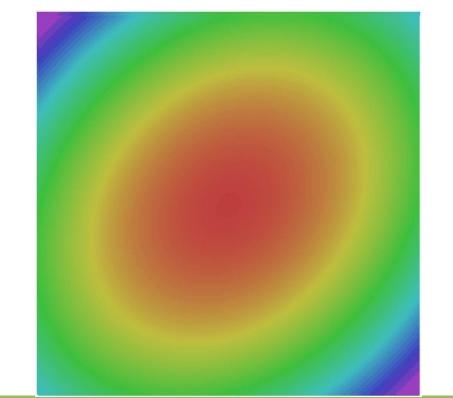
Image source: https://en.wikipedia.org/wiki/Saddle_point



Our gradients come from minibatches so they can be noisy!

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

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



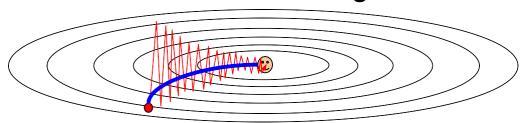


SGD + Momentum

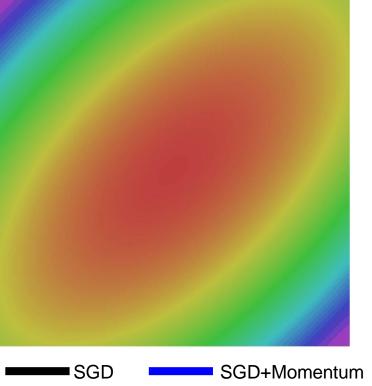
Local Minima Saddle points



Poor Conditioning



Gradient Noise



3GD+Momentum



SGD: the simple two line update code

SGD

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

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

SGD + Momentum:

continue moving in the general direction as the previous iterations SGD+Momentum

```
x_{t+1} = x_t - \alpha \nabla f(x_t) while True:  dx = compute\_gradient(x)   x -= learning\_rate * dx
```

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

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



SGD + Momentum:

continue moving in the general direction as the previous iterations SGD SGD+Momentum

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

```
while True:
  dx = compute\_gradient(x)
  x -= learning_rate * dx
```

```
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)
  vx = rho * vx + dx
  x -= learning_rate * vx
```

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



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}
```

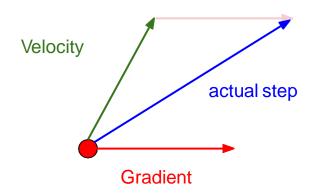
```
vx = 0
while True:
    dx = compute_gradient(x)
    vx = rho * vx + dx
    x -= learning_rate * vx
```

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



SGD+Momentum

Momentum update:

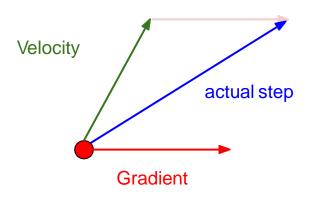


Combine gradient at current point with velocity to get step used to update weights

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



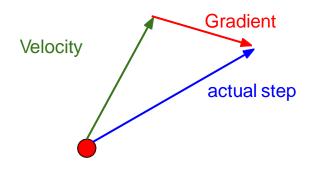
Momentum update:



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

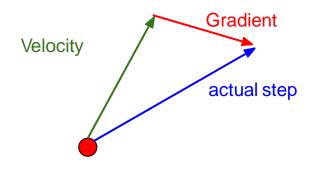
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Nesterov Momentum





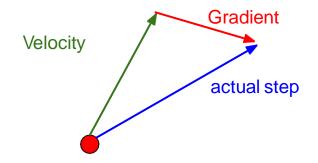
$$v_{t+1} = \rho v_t - \alpha \nabla f(x_t + \rho v_t)$$
$$x_{t+1} = x_t + v_{t+1}$$





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

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

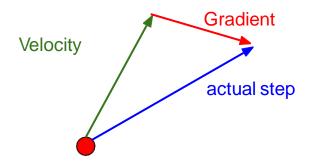




$$v_{t+1} = \rho v_t - \alpha \nabla f(x_t + \rho v_t)$$
$$x_{t+1} = x_t + v_{t+1}$$

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

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





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

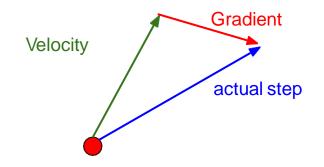
Change of variables $\tilde{x}_t = x_t + \rho v_t$ and rearrange:

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

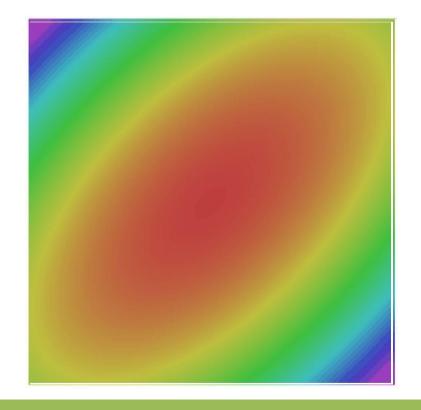
$$\tilde{x}_{t+1} = \tilde{x}_t - \rho v_t + (1+\rho)v_{t+1}$$

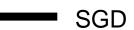
$$= \tilde{x}_t + v_{t+1} + \rho(v_{t+1} - v_t)$$

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









SGD+Momentum

Nesterov

```
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"

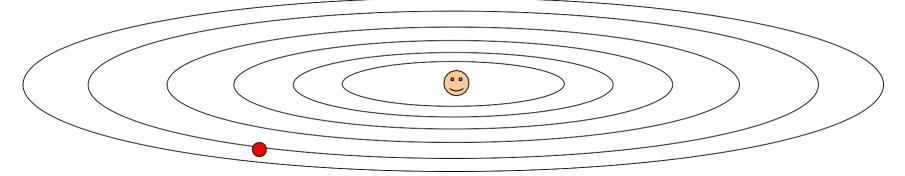


```
grad_squared = 0
while True:
  dx = compute\_gradient(x)
  grad_squared += dx * dx
 x -= learning_rate * dx / (np.sqrt(grad_squared) + 1e-7)
                                         ( 0 0 )
```

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

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```
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)
                                       00
```

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)
```

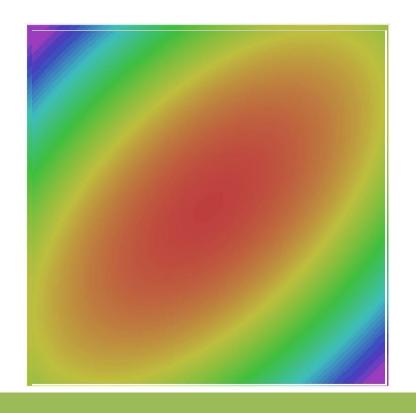


RMSProp

```
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)
```

RMSProp





SGD+Momentum

RMSProp

AdaGrad
(stuck due to decaying lr)



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))
```

Adam (almost)

```
first moment = 0
second moment = 0
while True:
 dx = compute\_gradient(x)
                                                                        Momentum
 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))
```

AdaGrad / RMSProp

Sort of like RMSProp with momentum

Q: What happens at first timestep?



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))
```

Momentum

Bias correction

AdaGrad / RMSProp

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



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))
```

Momentum

Bias correction

AdaGrad / RMSProp

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

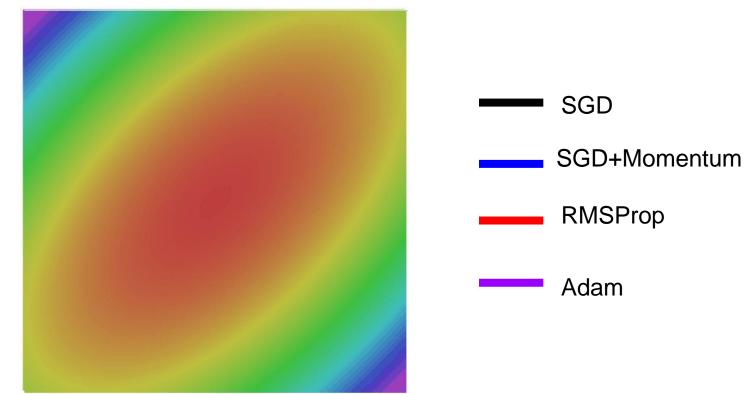
Adam with beta1 = 0.9, beta2 = 0.999, and learning_rate = 1e-3 or 5e-4 is a great starting point for many models!

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





Adam





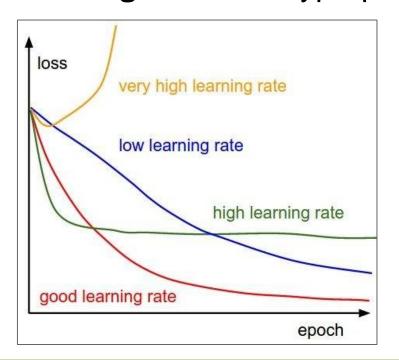
Learning rate schedules

```
# Vanilla Gradient Descent

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

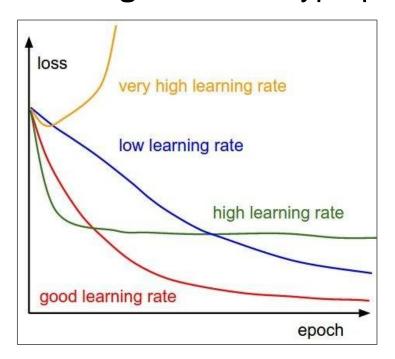
    Learning rate
```

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



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

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

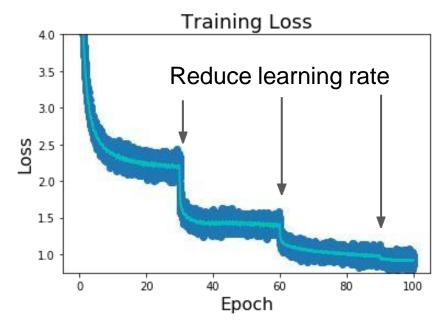


Q: Which one of these learning rates is best to use?

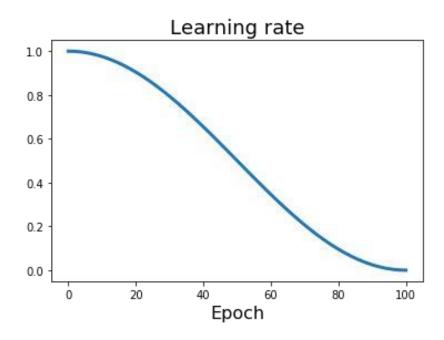
A: In reality, all of these are good learning rates.



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

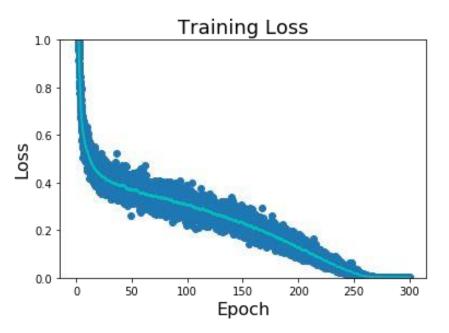
Regularization and Optimization

 $lpha_0$: Initial learning rate

 $lpha_t$: Learning rate at epoch t

T: Total number of epochs





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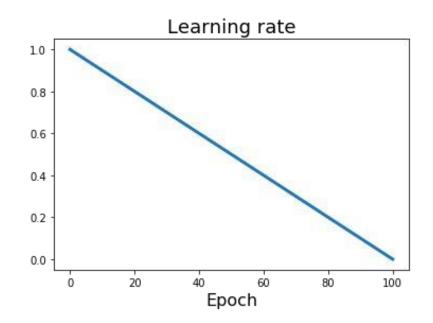
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Linear:
$$\alpha_t = \alpha_0(1 - t/T)$$

 $lpha_0$: Initial learning rate

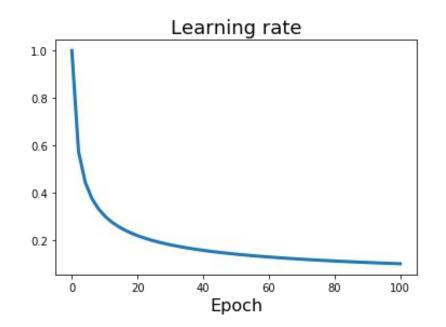
 $lpha_t$: Learning rate at epoch t

T : Total number of epochs

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

Regularization and Optimization





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

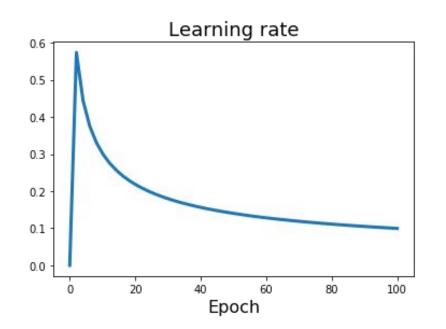
 $lpha_0$: Initial learning rate

 $lpha_t$: Learning rate at epoch t T : Total number of epochs

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



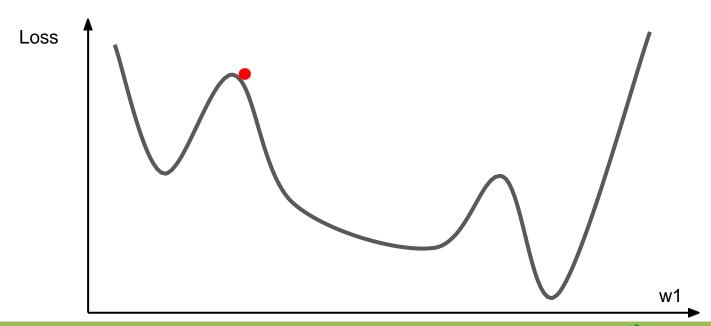
Learning Rate Decay: Linear Warmup



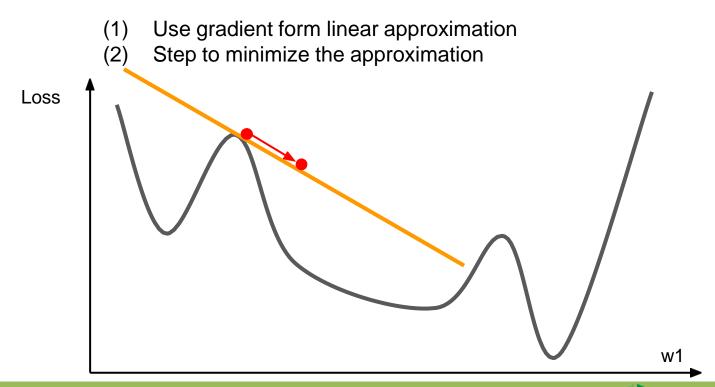
High initial learning rates can make loss explode; linearly increasing learning rate from 0 over the first ~5,000 iterations can prevent this.

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

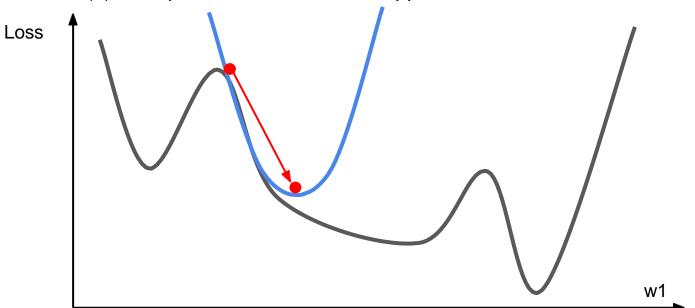
First-Order Optimization



First-Order Optimization



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





second-order Taylor expansion:

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

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:

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

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?



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

- 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.

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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)

Next time:

Introduction to neural networks

Backpropagation

