Lecture 4: Regularization + Optimization Reminder: Assignment 1

Was due on Friday!

If you enrolled late, you can have an extension – but email me / post on Piazza so we can track these

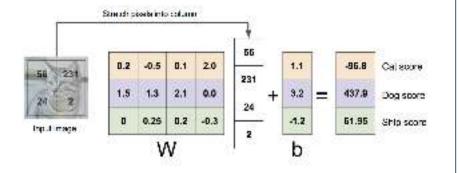
Assignment 2

- Released on Sunday
- Use SGD to train linear classifiers and fully-connected networks
- After today, can do linear classifiers section
- After Lecture 5, can do fully-connected networks
- If you have a hard time computing derivatives, wait for Lecture 6 on backprop
- Due Friday January 28, 11:59pm ET

Last Time: Linear Classifiers

Algebraic Viewpoint

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



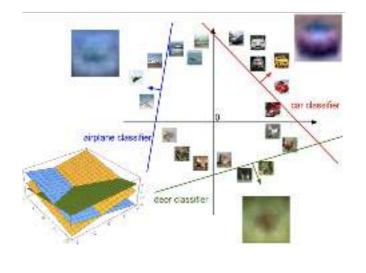
Visual Viewpoint

One template per class



Geometric Viewpoint

Hyperplanes cutting up space



Last Time: Loss Functions quantify preferences

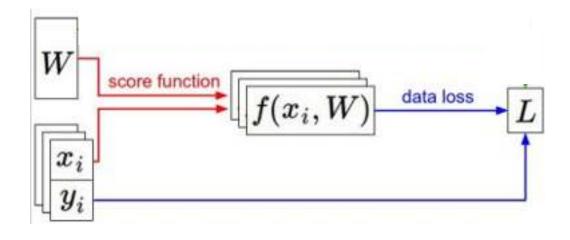
- We have some dataset of (x, y)
- We have a score function:
- We have a loss function:

Softmax:
$$L_i = -\log\left(\frac{\exp(s_{y_i})}{\sum_j \exp(s_j)}\right)$$

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

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

Linear classifier



Last Time: Loss Functions quantify preferences

- We have some dataset of (x, y)
- We have a **score function**:
- We have a loss function:

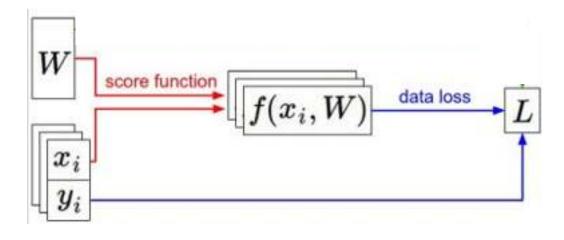
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SVM:
$$L_i = \sum_{j \neq y_i} \max(0, s_j - s_{y_i} + 1)$$

Problem: Loss functions encourage good performance on <u>training</u> data but we really care about <u>test</u> data

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

Linear classifier



A model is **overfit** when it performs too well on the training data, and has poor performance for unseen data

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$$p_i = \frac{\exp(s_i)}{\exp(s_1) + \exp(s_2)}$$

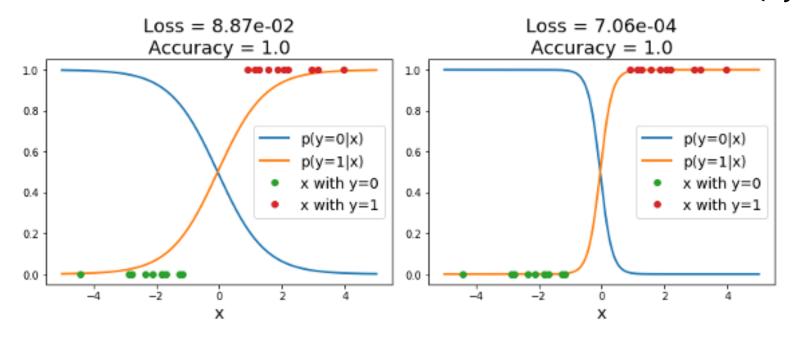
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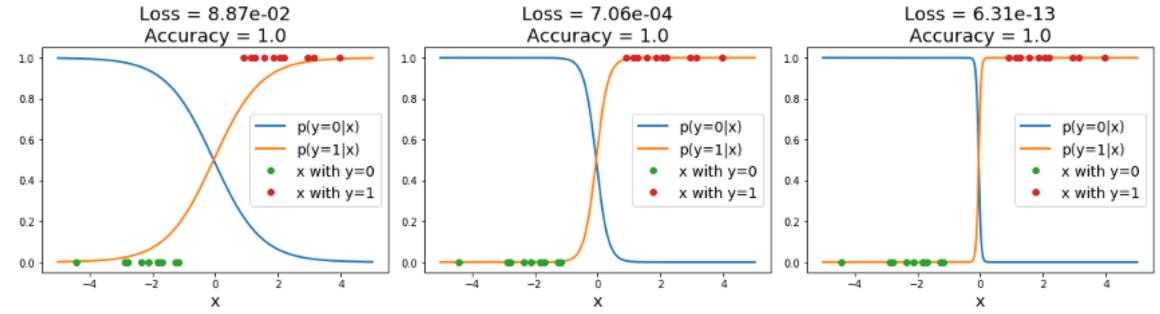
Both models have perfect accuracy on train data!

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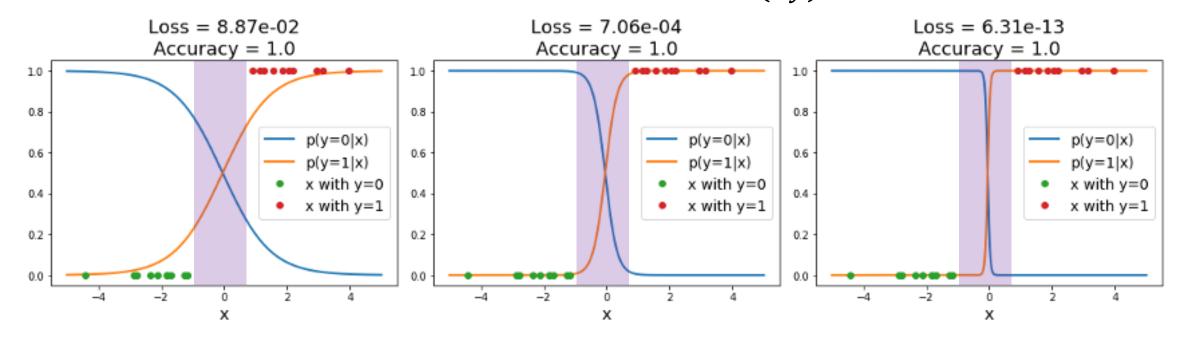
Low loss, but unnatural "cliff" between training points

A model is **overfit** when it performs too well on the training data, and has poor performance for unseen data

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$$p_i = \frac{\exp(s_i)}{\exp(s_1) + \exp(s_2)}$$

$$L = -\log(p_v)$$



Overconfidence in regions with no training data could give poor generalization

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

Data loss: Model predictions should match training data

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

 λ is a hyperparameter giving regularization strength

Data loss: Model predictions should match training data

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

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Data loss: Model predictions should match training data

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

Simple examples

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

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

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Simple examples

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

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

More complex:

Dropout

Batch normalization

Cutout, Mixup, Stochastic depth, etc...

Regularization: Prefer Simpler Models

$$s_i = w_i x + b_i \qquad p_i = \frac{\exp(s_i)}{\exp(s_1) + \exp(s_2)}$$

$$L = -\log(p_y) + \lambda \sum_i w_i^2$$

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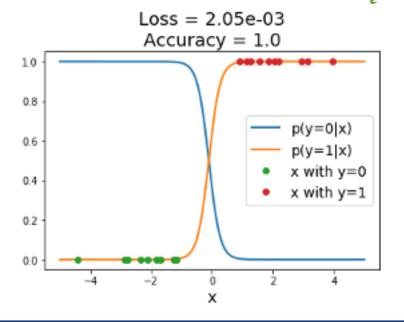
Regularization: Prefer Simpler Models

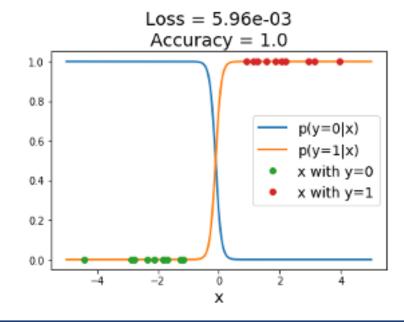
Example: Linear classifier with 1D inputs, 2 classes, softmax loss

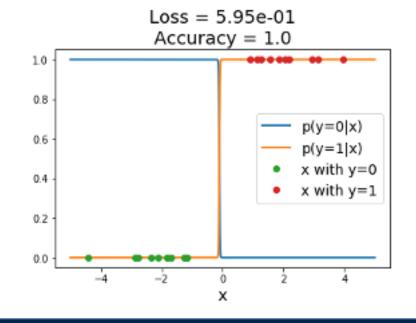
$$s_i = w_i x + b_i$$
 $p_i = \frac{\exp(s_i)}{\exp(s_1) + \exp(s_2)}$

$$L = -\log(p_y) + \lambda \sum_{i} w_i^2$$

Regularization term causes loss to **increase** for model with sharp cliff







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Regularization: Expressing Preferences

L2 Regularization

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

 $w_1 = [1, 0, 0, 0]$
 $w_2 = [0.25, 0.25, 0.25, 0.25]$

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

$$w_1^T x = w_2^T x = 1$$
 Same predictions, so data loss will always be the same

Regularization: Expressing Preferences

L2 Regularization

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

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

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

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

L2 regularization prefers weights to be "spread out"

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

Same predictions, so data loss will always be the same

Finding a good W

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

Loss function consists of data loss to fit the training data and regularization to prevent overfitting

Optimization

$$w^* = \arg\min_{w} L(w)$$



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Idea #1: Random Search (bad idea!)

```
# 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)
```

Idea #1: Random Search (bad idea!)

```
# 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!

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```

15.5% accuracy! not bad! (SOTA is ~95%)

Idea #2: Follow the slope



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In 1-dimension, the derivative of a function gives the slope:

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

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In 1-dimension, the derivative of a function gives the slope:

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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

W + h (first dim):

```
[0.34,
-1.11,
0.78,
0.12,
0.55,
2.81,
-3.1,
-1.5,
[0.33,...]
loss 1.25347
```

[?,
?,
?,
?,
?,
?,
?,
?,
?,]

W + h (first dim):

```
[-2.5,
(1.25322 - 1.25347)/0.0001
= -2.5
         \frac{df(x)}{df(x)} = \lim_{h \to \infty} \frac{f(x+h) - f(x)}{f(x+h) - f(x)}
               ?,...]
```

W + h (second dim):

```
[0.34,
-1.11,
0.78,
0.12,
0.55,
2.81,
-3.1,
-1.5,
0.33,...]
loss 1.25347
```

```
[-2.5,
?,...]
```

W + h (second dim):

```
[0.34,
-1.11,
0.78,
0.12,
0.55,
2.81,
-3.1,
-1.5,
[0.33,...]
loss 1.25347
```

```
[0.34,
-1.11 + 0.0001
0.78,
0.12,
0.55,
2.81,
-3.1,
-1.5,
0.33,...]
loss 1.25353
```

```
[-2.5,
              0.6,
(1.25353 - 1.25347)/0.0001
= 0.6
         \frac{df(x)}{df(x)} = \lim_{x \to \infty} \frac{f(x+h) - f(x)}{f(x+h)}
```

W + h (third dim):

gradient dL/dW:

[0.34,
-1.11,
0.78,
0.12,
0.55,
2.81,
-3.1,
-1.5,
0.33,]
loss 1.25347

```
[0.34,
-1.11,
0.78 + 0.0001,
0.12,
0.55,
2.81,
-3.1,
-1.5,
0.33,...]
loss 1.25347
```

[-2.5,0.6, ?,...]

W + h (third dim):

```
[0.34,
-1.11,
0.78,
0.12,
0.55,
2.81,
-3.1,
-1.5,
[0.33,...]
loss 1.25347
```

```
[-2.5,
        0.6,
        0.0,
(1.25347 - 1.25347)/0.0001
= 0.0
```

current W:

W + h (third dim):

gradient dL/dW:

```
[0.34,
-1.11,
0.78,
0.12,
0.55,
2.81,
-3.1,
-1.5,
[0.33,...]
loss 1.25347
```

[-2.5, 0.6, **0.0**, ?, ?,

Numeric Gradient:

- Slow: O(#dimensions)
- Approximate

Loss is a function of W

$$L = \frac{1}{2} \sum_{i=1}^{N} L_i + \sum_{k} W_k^2$$

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

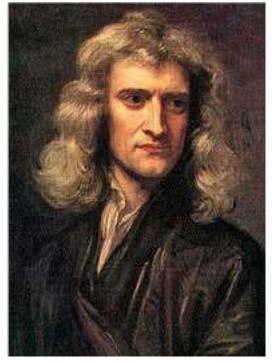
$$s = f(x, W) = Wx$$
Want $\nabla_W L$

Loss is a function of W: Analytic Gradient

$$L = \frac{1}{2} \sum_{i=1}^{N} L_i + \sum_{k} W_k^2$$

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

$$s = f(x, W) = Wx$$
Want $\nabla_W L$





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Use calculus to compute an analytic gradient

current W:

gradient dL/dW:

```
[0.34,
-1.11,
0.78,
0.12,
0.55,
2.81,
-3.1,
-1.5,
0.33,...]
loss 1.25347
```

current W:

gradient dL/dW:

```
[0.34,
-1.11,
0.78,
0.12,
0.55,
2.81,
-3.1,
-1.5,
[0.33,...]
loss 1.25347
```

```
[-2.5,
 dL/dW = ...
                                 0.6,
 (some function
                                 0,
 data and W)
                                0.2,
                                0.7,
                                 -0.5,
(In practice we will
                                 1.1,
compute dL/dW using
                                 1.3,
backpropagation; see
                                -2.1,...]
Lecture 6)
```

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

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

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

```
def grad_check_sparse(f, x, analytic_grad, num_checks=10, h=1e-7):
    sample a few random elements and only return numerical
    in this dimensions.
    """
```

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

```
torch.autograd.gradcheck(func, inputs, eps=1e-06, atol=1e-05, rtol=0.001, raise_exception=True, check_sparse_nnz=False, nondet_tol=0.0) [SOURCE] &
```

Check gradients computed via small finite differences against analytical gradients w.r.t. tensors in inputs that are of floating point type and with requires_grad=True.

The check between numerical and analytical gradients uses allclose().

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

```
torch.autograd.gradgradcheck(func, inputs, grad_outputs=None, eps=1e-06, atol=1e-
05, rtol=0.001, gen_non_contig_grad_outputs=False, raise_exception=True, [SOURCE]
nondet_tol=0.0)
```

Check gradients of gradients computed via small finite differences against analytical gradients w.r.t. tensors in inputs and grad_outputs that are of floating point type and with requires_grad=True.

This function checks that backpropagating through the gradients computed to the given <code>grad_outputs</code> are correct.

Gradient Descent

Iteratively step in the direction of the negative gradient (direction of local steepest descent)

```
# Vanilla gradient descent
w = initialize_weights()
for t in range(num_steps):
   dw = compute_gradient(loss_fn, data, w)
   w -= learning_rate * dw
```

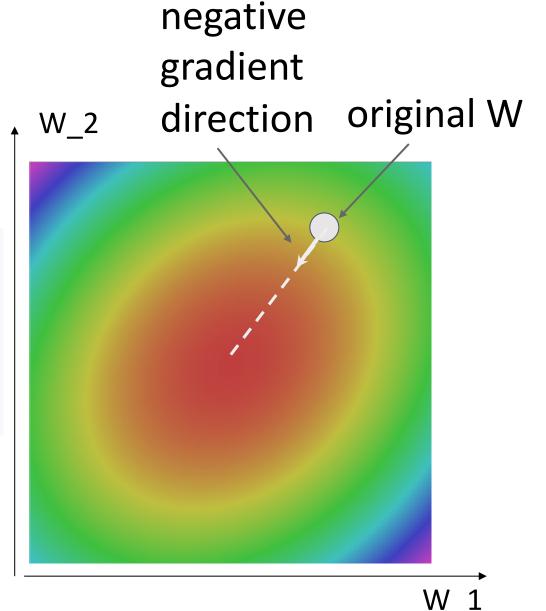
- Weight initialization method
- Number of steps
- Learning rate

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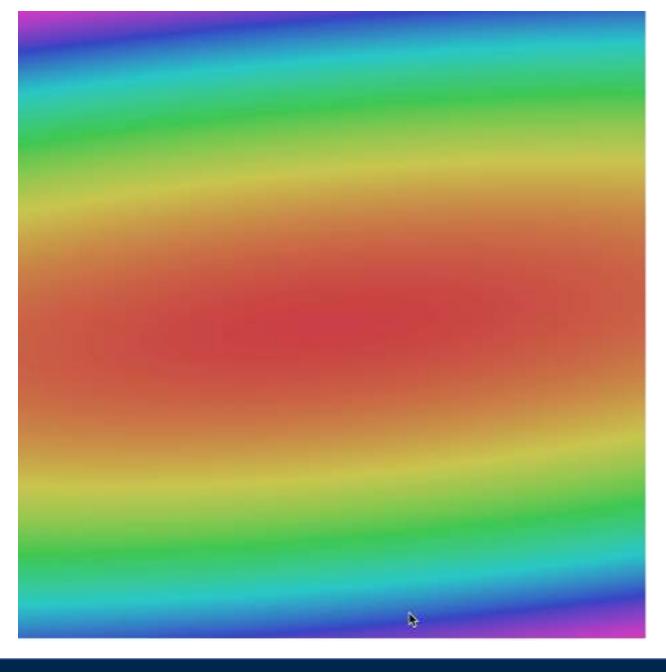


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```

- Weight initialization method
- Number of steps
- Learning rate



Batch Gradient Descent

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

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

Full sum expensive when N is large!

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

```
# Stochastic gradient descent
w = initialize_weights()
for t in range(num_steps):
   minibatch = sample_data(data, batch_size)
   dw = compute_gradient(loss_fn, minibatch, w)
   w -= learning_rate * dw
```

- Weight initialization
- Number of steps
- Learning rate
- Batch size
- Data sampling

Stochastic Gradient Descent (SGD)

$$L(W) = \mathbb{E}_{(x,y)\sim p_{data}}[L(x,y,W)] + \lambda R(W)$$

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

Think of loss as an expectation over the full data distribution p_{data}

Approximate expectation via sampling

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Stochastic Gradient Descent (SGD)

$$L(W) = \mathbb{E}_{(x,y) \sim p_{data}} [L(x,y,W)] + \lambda R(W)$$

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

Think of loss as an expectation over the full data distribution p_{data}

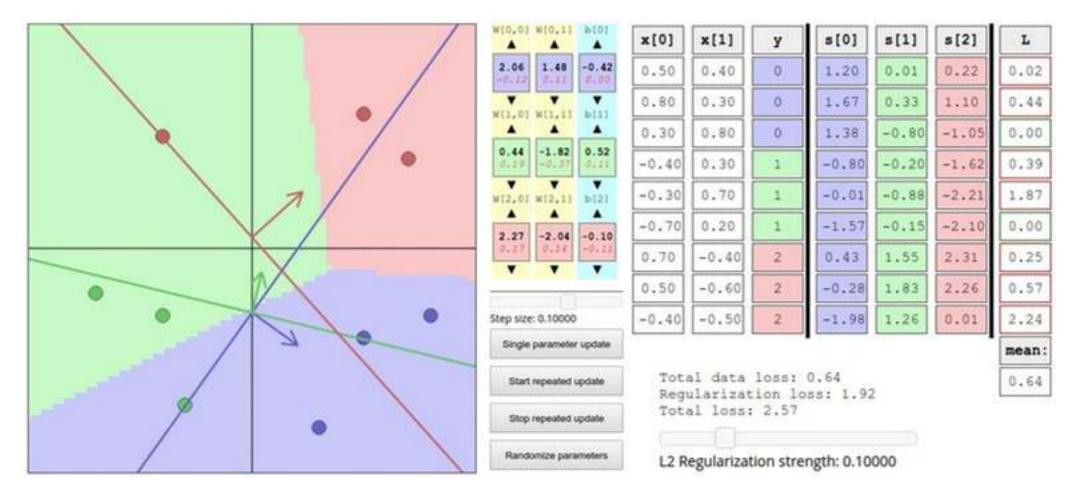
Approximate expectation via sampling

$$\nabla_{W} L(W) = \nabla_{W} \mathbb{E}_{(x,y) \sim p_{data}} [L(x,y,W)] + \lambda \nabla_{W} R(W)$$

$$\approx \sum_{i=1}^{N} \nabla_{w} L_{W}(x_{i},y_{i},W) + \nabla_{w} R(W)$$

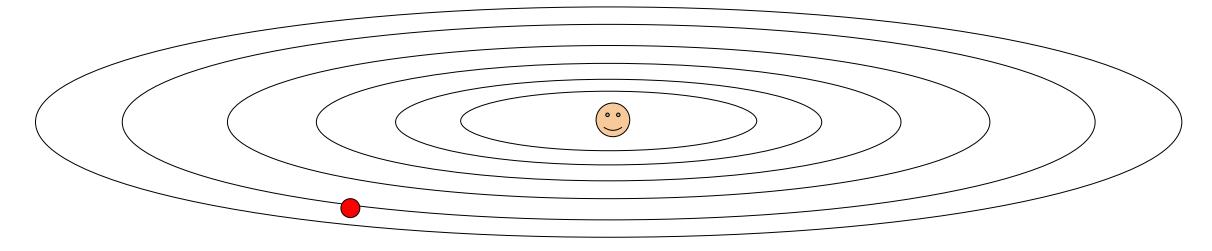
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Interactive Web Demo



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

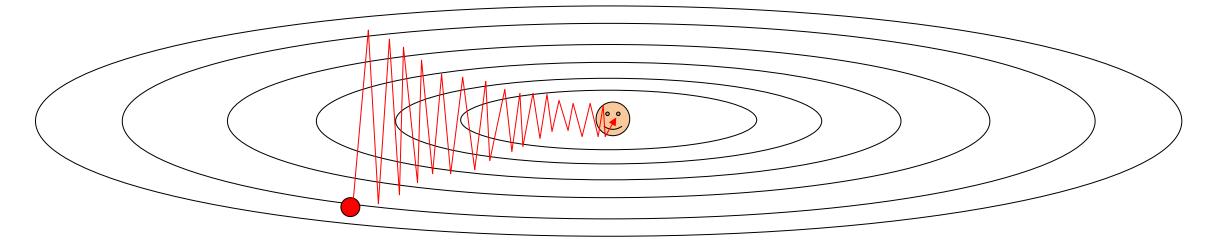
What if loss changes quickly in one direction and slowly in another? What does gradient descent do?



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

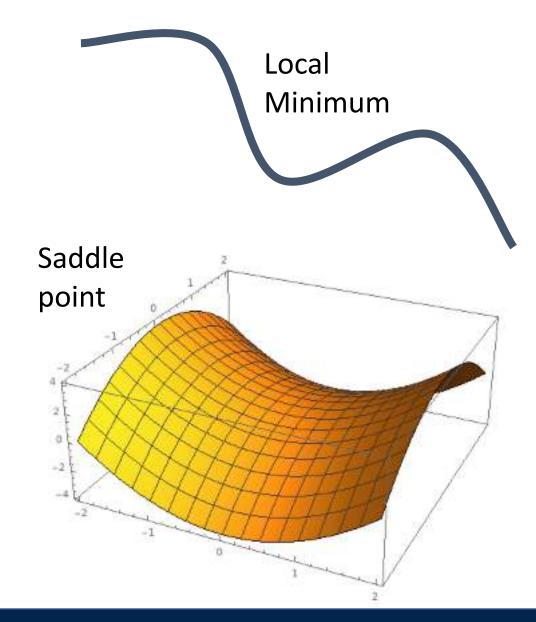
What if loss changes quickly in one direction and slowly in another? What does gradient descent do?

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

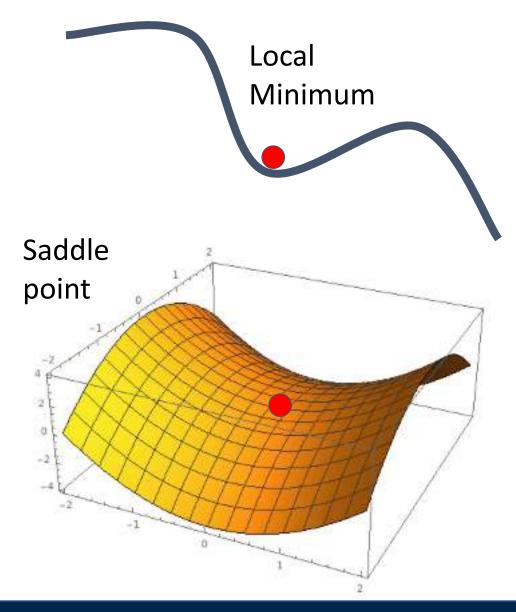
What if the loss function has a **local minimum** or **saddle point**?



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

Zero gradient, gradient descent gets stuck

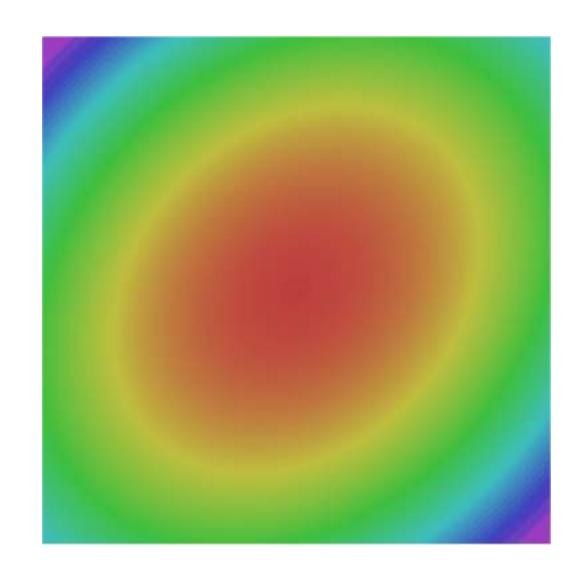


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Our gradients come from minibatches so they can be noisy!

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

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



SGD

SGD

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

```
for t in range(num_steps):
   dw = compute_gradient(w)
   w -= learning_rate * dw
```

SGD

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

```
for t in range(num_steps):
   dw = compute_gradient(w)
   w -= learning_rate * dw
```

SGD+Momentum

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

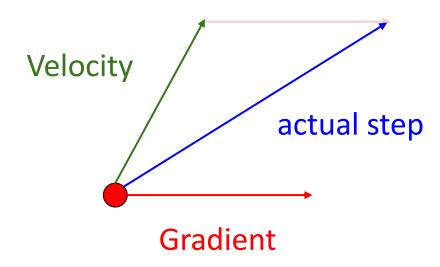
$$x_{t+1} = x_t - \alpha v_{t+1}$$

```
v = 0
for t in range(num_steps):
   dw = compute_gradient(w)
   v = rho * v + dw
   w -= learning_rate * v
```

- 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

Momentum update:



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

SGD+Momentum

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

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

```
v = 0
for t in range(num_steps):
   dw = compute_gradient(w)
   v = rho * v + dw
   w -= learning_rate * v
```

- 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

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

$$x_{t+1} = x_t + v_{t+1}$$

```
v = 0
for t in range(num_steps):
  dw = compute_gradient(w)
  v = rho * v - learning_rate * dw
  w += v
```

SGD+Momentum

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

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

```
v = 0
for t in range(num_steps):
   dw = compute_gradient(w)
   v = rho * v + dw
   w -= learning_rate * v
```

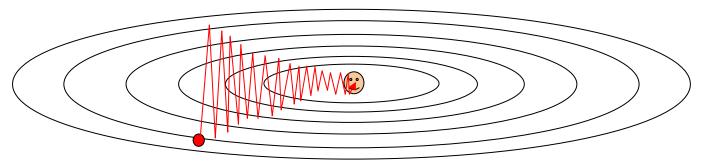
You may see SGD+Momentum formulated different ways, but they are equivalent - give same sequence of x

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

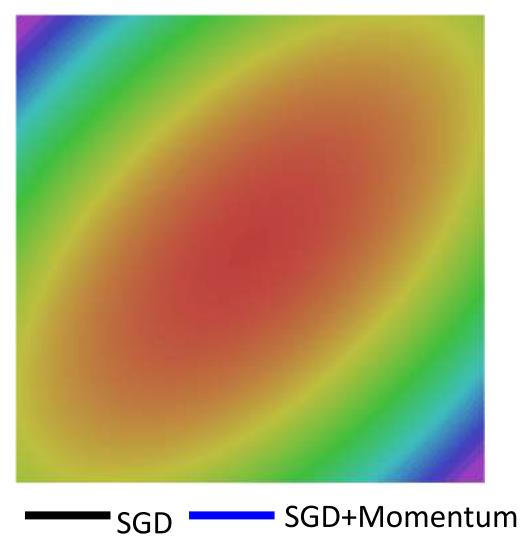
Local Minima Saddle points



Poor Conditioning



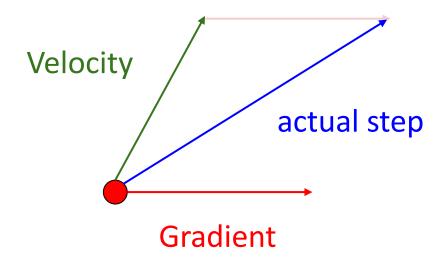
Gradient Noise



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

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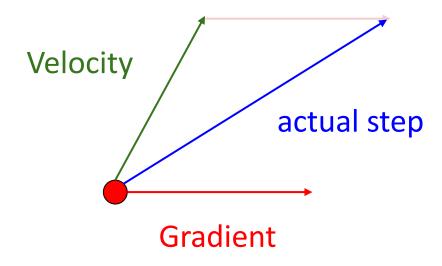
Momentum update:



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

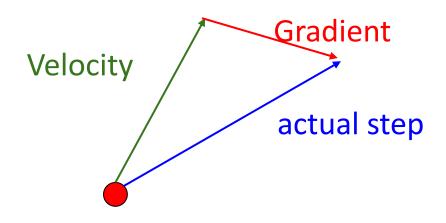
Momentum update:



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

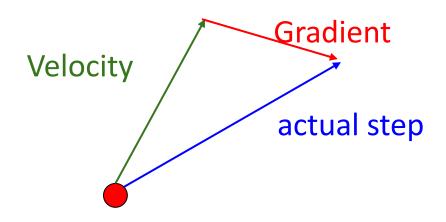
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Nesterov Momentum



"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

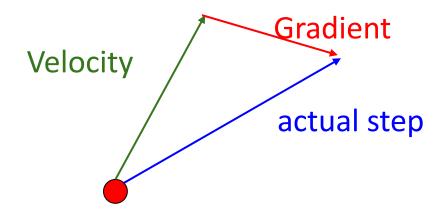
$$v_{t+1} = \rho v_t - \alpha \nabla f(x_t + \rho v_t)$$
$$x_{t+1} = x_t + v_{t+1}$$



"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

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

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



"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

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

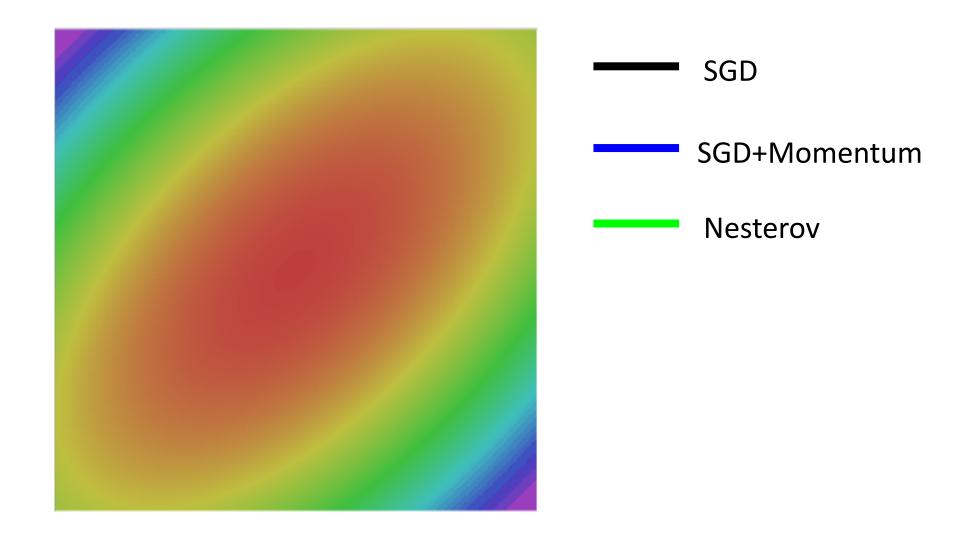
Change of variables $\tilde{x}_t = x_t + \rho v_t$ and rearrange:

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

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

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

```
v = 0
for t in range(num_steps):
   dw = compute_gradient(w)
   old_v = v
   v = rho * v - learning_rate * dw
   w -= rho * old_v - (1 + rho) * v
```



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AdaGrad

```
grad_squared = 0
for t in range(num_steps):
   dw = compute_gradient(w)
   grad_squared += dw * dw
w -= learning_rate * dw / (grad_squared.sqrt() + 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

AdaGrad

```
grad_squared = 0
for t in range(num_steps):
  dw = compute_gradient(w)
  grad_squared += dw * dw
  w -= learning_rate * dw / (grad_squared.sqrt() + 1e-7)
```

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

AdaGrad

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grad_squared = 0
for t in range(num_steps):
  dw = compute_gradient(w)
  grad_squared += dw * dw
  w -= learning_rate * dw / (grad_squared.sqrt() + 1e-7)
```

Q: What happens with AdaGrad?

AdaGrad

```
grad_squared = 0
for t in range(num_steps):
  dw = compute_gradient(w)
 grad_squared += dw * dw
 w -= learning_rate * dw / (grad_squared.sqrt() + 1e-7)
```

Q: What happens with AdaGrad?

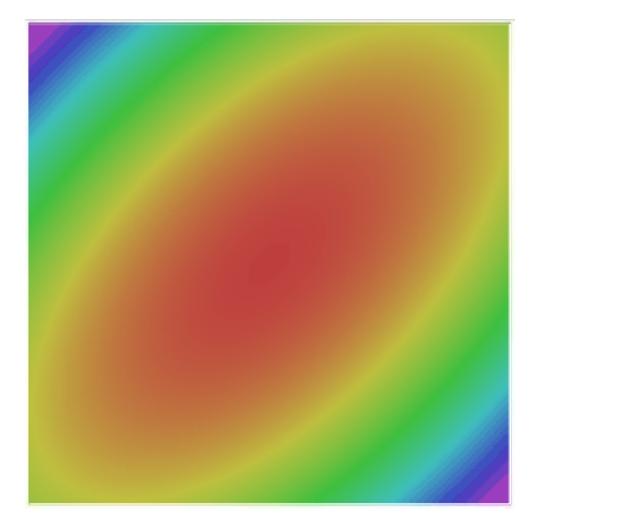
Progress along "steep" directions is damped; progress along "flat" directions is accelerated

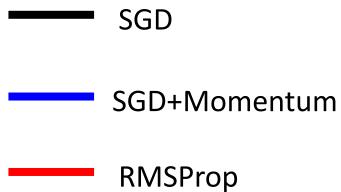
RMSProp: "Leaky Adagrad"

```
grad_squared = 0
for t in range(num_steps):
                                                             AdaGrad
  dw = compute_gradient(w)
  grad_squared += dw * dw
  w -= learning_rate * dw / (grad_squared.sqrt() + 1e-7)
grad_squared = 0
for t in range(num_steps):
 dw = compute_gradient(w)
                                                                       RMSProp
 grad_squared = decay_rate * grad_squared + (1 - decay_rate) * dw * dw
 w -= learning_rate * dw / (grad_squared.sqrt() + 1e-7)
```

Tieleman and Hinton, 2012

RMSProp





```
moment1 = 0
moment2 = 0
for t in range(1, num_steps + 1):  # Start at t = 1
  dw = compute_gradient(w)
  moment1 = beta1 * moment1 + (1 - beta1) * dw
  moment2 = beta2 * moment2 + (1 - beta2) * dw * dw
  w -= learning_rate * moment1 / (moment2.sqrt() + 1e-7)
```

```
moment1 = 0
moment2 = 0
for t in range(1, num_steps + 1): # Start at t = 1
  dw = compute_gradient(w)

moment1 = beta1 * moment1 + (1 - beta1) * dw

moment2 = beta2 * moment2 + (1 - beta2) * dw * dw

w -= learning_rate * moment1 / (moment2.sqrt() + 1e-7)
```

Adam

Momentum

```
v = 0
for t in range(num_steps):
   dw = compute_gradient(w)
   v = rho * v + dw
   w -= learning_rate * v
```

SGD+Momentum

```
moment1 = 0
moment2 = 0
for t in range(1, num_steps + 1): # Start at t = 1
   dw = compute_gradient(w)

moment1 = beta1 * moment1 + (1 - beta1) * dw

moment2 = beta2 * moment2 + (1 - beta2) * dw * dw

w -= learning_rate * moment1 / (moment2.sqrt() + 1e-7)
```

Adam

Momentum

AdaGrad / RMSProp

```
grad_squared = 0
for t in range(num_steps):
   dw = compute_gradient(w)
   grad_squared = decay_rate * grad_squared + (1 - decay_rate) * dw * dw
   w -= learning_rate * dw / (grad_squared.sqrt() + 1e-7)
```

RMSProp

```
moment1 = 0
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for t in range(1, num_steps + 1):  # Start at t = 1
  dw = compute_gradient(w)
  moment1 = beta1 * moment1 + (1 - beta1) * dw
  moment2 = beta2 * moment2 + (1 - beta2) * dw * dw
  w -= learning_rate * moment1 / (moment2.sqrt() + 1e-7)
```

Adam

Momentum

AdaGrad / RMSProp

Bias correction

Q: What happens at t=1? (Assume beta2 = 0.999)

```
moment1 = 0
moment2 = 0
for t in range(1, num_steps + 1): # Start at t = 1
    dw = compute_gradient(w)
    moment1 = beta1 * moment1 + (1 - beta1) * dw
    moment2 = beta2 * moment2 + (1 - beta2) * dw * dw
    moment1_unbias = moment1 / (1 - beta1 ** t)
    moment2_unbias = moment2 / (1 - beta2 ** t)
    w -= learning_rate * moment1_unbias / (moment2_unbias.sqrt() + 1e-7)
```

Momentum

AdaGrad / RMSProp

Bias correction

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

```
moment1 = 0
moment2 = 0
for t in range(1, num_steps + 1):  # Start at t = 1
  dw = compute_gradient(w)
  moment1 = beta1 * moment1 + (1 - beta1) * dw
  moment2 = beta2 * moment2 + (1 - beta2) * dw * dw
  moment1_unbias = moment1 / (1 - beta1 ** t)
  moment2_unbias = moment2 / (1 - beta2 ** t)
  w -= learning_rate * moment1_unbias / (moment2_unbias.sqrt() + 1e-7)
```

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, 5e-4, 1e-4 is a great starting point for many models!

Adam: Very Common in Practice!

for input to the CNN; each colored pixel in the image yields a 7D one-hot vector. Following common practice, the network is trained end-to-end using stochastic gradient descent with the Adam optimizer [22]. We anneal the learning rate to 0 using a half cosine schedule without restarts [28].

Bakhtin, van der Maaten, Johnson, Gustafson, and Girshick, NeurIPS 2019

We train all models using Adam [23] with learning rate 10^{-4} and batch size 32 for 1 million iterations; training takes about 3 days on a single Tesla P100. For each minibatch we first update f, then update D_{img} and D_{obj} .

Johnson, Gupta, and Fei-Fei, CVPR 2018

ganized into three residual blocks. We train for 25 epochs using Adam [27] with learning rate 10^{-4} and 32 images per batch on 8 Tesla V100 GPUs. We set the cubify thresh-

Gkioxari, Malik, and Johnson, ICCV 2019

sampled with each bit drawn uniformly at random. For gradient descent, we use Adam [29] with a learning rate of 10^{-3} and default hyperparameters. All models are trained with batch size 12. Models are trained for 200 epochs, or 400 epochs if being trained on multiple noise layers.

Zhu, Kaplan, Johnson, and Fei-Fei, ECCV 2018

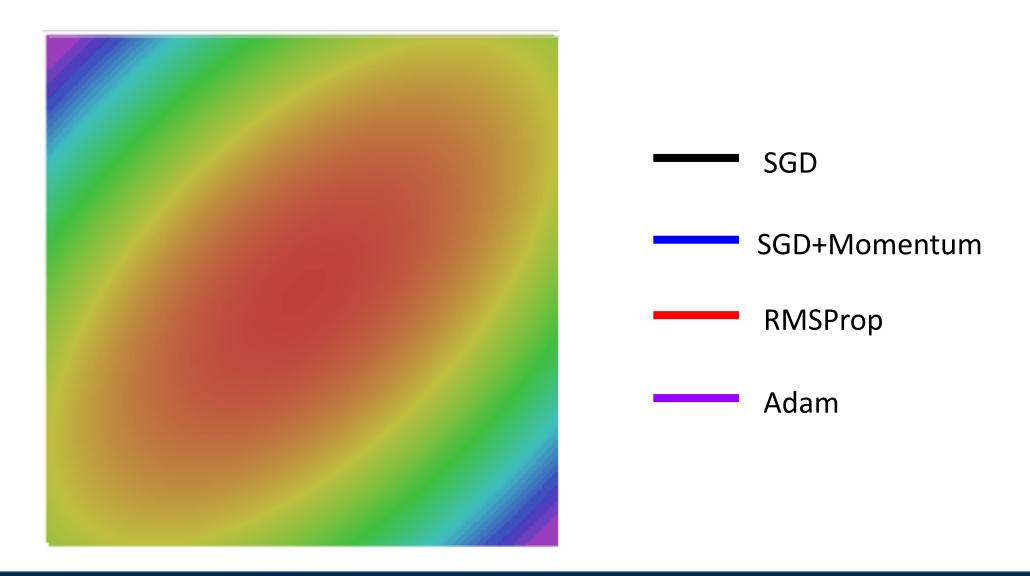
16 dimensional vectors. We iteratively train the Generator and Discriminator with a batch size of 64 for 200 epochs using Adam [22] with an initial learning rate of 0.001.

Gupta, Johnson, et al, CVPR 2018

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

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Adam



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Optimization Algorithm Comparison

Algorithm	Tracks first moments (Momentum)	Tracks second moments (Adaptive learning rates)	Leaky second moments	Bias correction for moment estimates
SGD	X	X	X	X
SGD+Momentum	✓	X	X	X
Nesterov	✓	X	X	X
AdaGrad	X	✓	X	X
RMSProp	X	✓	✓	X
Adam	✓	✓	✓	✓

Optimization Algorithm

$$L(w) = L_{data}(w) + L_{reg}(w)$$

$$g_t = \nabla L(w_t)$$

$$s_t = optimizer(g_t)$$

$$w_{t+1} = w_t - \alpha s_t$$

Optimization Algorithm

$$L(w) = L_{data}(w) + L_{reg}(w)$$

$$g_t = \nabla L(w_t)$$

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$$w_{t+1} = w_t - \alpha s_t$$

L2 Regularization

$$L(w) = L_{data}(w) + \lambda |w|^{2}$$

$$g_{t} = \nabla L(w_{t}) = \nabla L_{data}(w_{t}) + 2\lambda w_{t}$$

$$s_{t} = optimizer(g_{t})$$

$$w_{t+1} = w_{t} - \alpha s_{t}$$

Optimization Algorithm

$$L(w) = L_{data}(w) + L_{reg}(w)$$

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Weight Decay

$$L(w) = L_{data}(w)$$

$$g_t = \nabla L_{data}(w_t)$$

$$s_t = optimizer(g_t) + 2\lambda w_t$$

$$w_{t+1} = w_t - \alpha s_t$$

Optimization Algorithm

$$L(w) = L_{data}(w) + L_{reg}(w)$$

$$g_t = \nabla L(w_t)$$

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L2 Regularization and Weight Decay are equivalent for SGD, SGD+Momentum so people often use the terms interchangeably!

L2 Regularization

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Optimization Algorithm

$$L(w) = L_{data}(w) + L_{reg}(w)$$

$$g_t = \nabla L(w_t)$$

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$$w_{t+1} = w_t - \alpha s_t$$

L2 Regularization and Weight Decay are equivalent for SGD, SGD+Momentum so people often use the terms interchangeably!

But they are not the same for adaptive methods (AdaGrad, RMSProp, Adam, etc)

L2 Regularization

$$L(w) = L_{data}(w) + \lambda |w|^{2}$$

$$g_{t} = \nabla L(w_{t}) = \nabla L_{data}(w_{t}) + 2\lambda w_{t}$$

$$s_{t} = optimizer(g_{t})$$

$$w_{t+1} = w_{t} - \alpha s_{t}$$

Weight Decay

$$L(w) = L_{data}(w)$$

$$g_t = \nabla L_{data}(w_t)$$

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$$w_{t+1} = w_t - \alpha s_t$$

AdamW: Decoupled Weight Decay

Algorithm 2 Adam with L₂ regularization and Adam with decoupled weight decay (AdamW)

```
1: given \alpha = 0.001, \beta_1 = 0.9, \beta_2 = 0.999, \epsilon = 10^{-8}, \lambda \in \mathbb{R}
  2: initialize time step t \leftarrow 0, parameter vector \boldsymbol{\theta}_{t=0} \in \mathbb{R}^n, first moment vector \boldsymbol{m}_{t=0} \leftarrow \boldsymbol{\theta}, second moment
        vector \mathbf{v}_{t=0} \leftarrow \mathbf{0}, schedule multiplier \eta_{t=0} \in \mathbb{R}
 3: repeat
 4: t \leftarrow t+1
 5: \nabla f_t(\boldsymbol{\theta}_{t-1}) \leftarrow \text{SelectBatch}(\boldsymbol{\theta}_{t-1})

▷ select batch and return the corresponding gradient

 6: \mathbf{g}_t \leftarrow \nabla f_t(\boldsymbol{\theta}_{t-1}) + \lambda \boldsymbol{\theta}_{t-1}
 7: \boldsymbol{m}_t \leftarrow \beta_1 \boldsymbol{m}_{t-1} + (1 - \beta_1) \boldsymbol{g}_t
                                                                                                                ▶ here and below all operations are element-wise
 8: \mathbf{v}_t \leftarrow \beta_2 \mathbf{v}_{t-1} + (1 - \beta_2) \mathbf{g}_t^2

9: \hat{\mathbf{m}}_t \leftarrow \mathbf{m}_t / (1 - \beta_1^t)
                                                                                                                                                    \triangleright \beta_1 is taken to the power of t
10: \hat{\mathbf{v}}_t \leftarrow \mathbf{v}_t/(1-\beta_2^t)
                                                                                                                                                    \triangleright \beta_2 is taken to the power of t
11: \eta_t \leftarrow \text{SetScheduleMultiplier}(t)
                                                                                                     > can be fixed, decay, or also be used for warm restarts
         \boldsymbol{\theta}_t \leftarrow \boldsymbol{\theta}_{t-1} - \eta_t \left( \alpha \hat{\boldsymbol{m}}_t / (\sqrt{\hat{\boldsymbol{v}}_t} + \epsilon) + \lambda \boldsymbol{\theta}_{t-1} \right)
13: until stopping criterion is met
```

Loshchilov and Hutter, "Decoupled Weight Decay Regularization", ICLR 2019

14: **return** optimized parameters θ_t

AdamW: Decoupled Weight Decay

Algorithm 2 Adam with L₂ regularization and Adam with decoupled weight decay (AdamW)

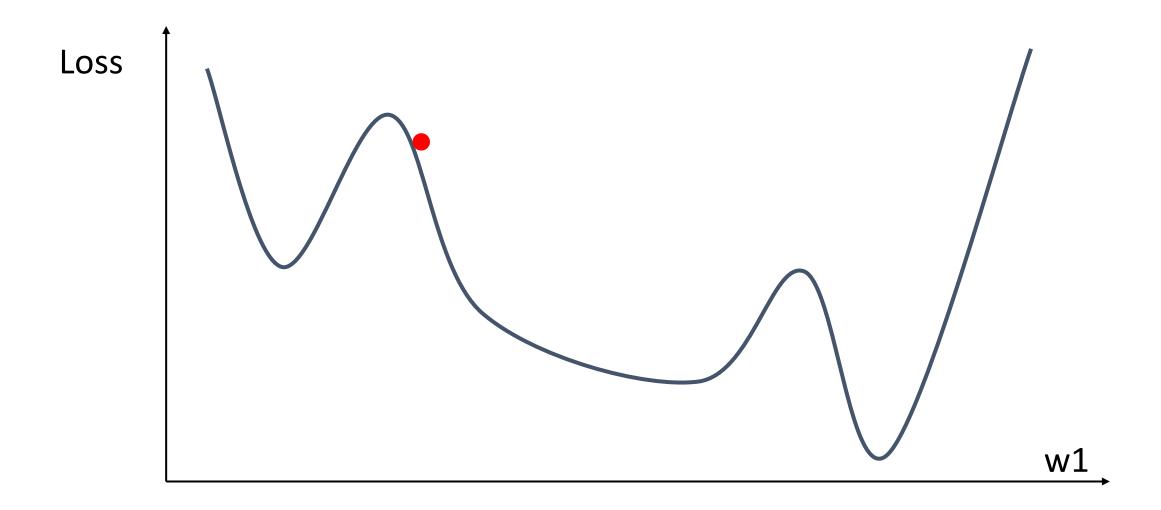
- 1: given $\alpha = 0.001, \beta_1 = 0.9, \beta_2 = 0.999, \epsilon = 10^{-8}, \lambda \in \mathbb{R}$
- 2: **initialize** time step $t \leftarrow 0$, parameter vector $\boldsymbol{\theta}_{t=0} \in \mathbb{R}^n$, first moment vector $\boldsymbol{m}_{t=0} \leftarrow \boldsymbol{\theta}$, second moment vector $\boldsymbol{v}_{t=0} \leftarrow \boldsymbol{\theta}$, schedule multiplier $n_{t=0} \in \mathbb{R}$.

AdamW should probably be your "default" optimizer for new problems

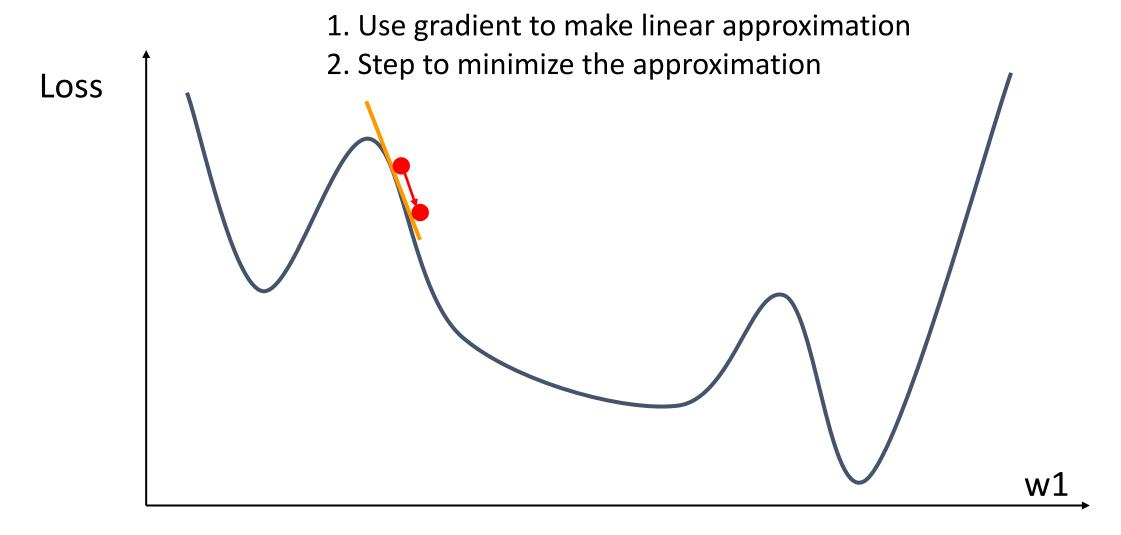
11: $\eta_t \leftarrow \text{SetScheduleMultiplier}(t)$

- > can be fixed, decay, or also be used for warm restarts
- 12: $\boldsymbol{\theta}_t \leftarrow \boldsymbol{\theta}_{t-1} \eta_t \left(\alpha \hat{\boldsymbol{m}}_t / (\sqrt{\hat{\boldsymbol{v}}_t} + \epsilon) + \lambda \boldsymbol{\theta}_{t-1} \right)$
- 13: until stopping criterion is met
- 14: **return** optimized parameters θ_t

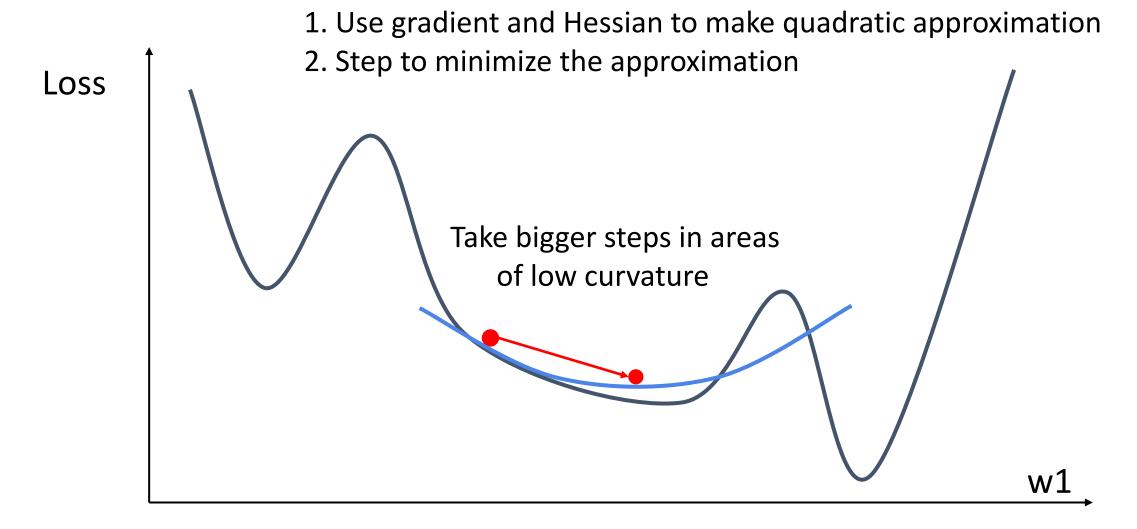
So far: First-Order Optimization



So far: First-Order Optimization



1. Use gradient and Hessian to make quadratic approximation 2. Step to minimize the approximation Loss w1



Second-Order Taylor Expansion:

$$L(w) \approx L(w_0) + (w - w_0)^{\mathsf{T}} \nabla_w L(w_0) + \frac{1}{2} (w - w_0)^{\mathsf{T}} \mathbf{H}_w L(w_0) (w - w_0)$$

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

$$w^* = w_0 - \mathbf{H}_w L(w_0)^{-1} \nabla_w L(w_0)$$

Second-Order Taylor Expansion:

$$L(w) \approx L(w_0) + (w - w_0)^{\mathsf{T}} \nabla_w L(w_0) + \frac{1}{2} (w - w_0)^{\mathsf{T}} \mathbf{H}_w L(w_0) (w - w_0)$$

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Q: Why is this impractical?

Second-Order Taylor Expansion:

$$L(w) \approx L(w_0) + (w - w_0)^{\mathsf{T}} \nabla_w L(w_0) + \frac{1}{2} (w - w_0)^{\mathsf{T}} \mathbf{H}_w L(w_0) (w - w_0)$$

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

$$w^* = w_0 - \mathbf{H}_w L(w_0)^{-1} \nabla_w L(w_0)$$

Q: Why is this impractical?

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

$$w^* = w_0 - \mathbf{H}_w L(w_0)^{-1} \nabla_w L(w_0)$$

- 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.

Second-Order Optimization: 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.

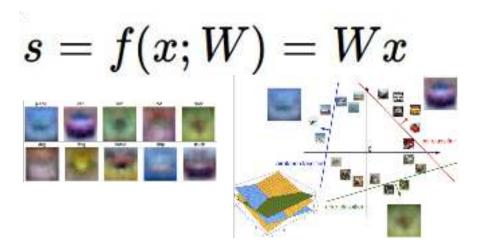
Le et al, "On optimization methods for deep learning, ICML 2011"
Ba et al, "Distributed second-order optimization using Kronecker-factored approximations", ICLR 2017

In practice:

- Adam is a good default choice in many cases
 SGD+Momentum can outperform Adam but may require more tuning
- If you can afford to do full batch updates then try out **L-BFGS** (and don't forget to disable all sources of noise)

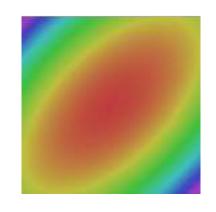
Summary

- 1. Use **Linear Models** for image classification problems
- 2. Use **Loss Functions** to express preferences over different choices of weights
- 3. Use **Regularization** to prevent overfitting to training data
- Use Stochastic Gradient
 Descent to minimize our loss functions and train the model



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

```
v = 0
for t in range(num_steps):
   dw = compute_gradient(w)
   v = rho * v + dw
   w -= learning_rate * v
```



Next time: Neural Networks