













Lecture 4 **Regularization + Optimization**







- Instructions and code available on the website • Here: <u>deeprob.org/projects/project1/</u>
- Uses Python, PyTorch and Google Colab
- Implement KNN, linear SVM, and linear softmax classifiers
- Autograder is online
- Due Thursday, January 26th 11:59 PM EST



Project 1 – Reminder



Progress Robot Object Perception Samples Dataset



Chen et al., "ProgressLabeller: Visual Data Stream Annotation for Training Object-Centric 3D Perception", IROS, 2022.



Project 1 – Dataset

10 classes 32x32 RGB images **50k** training images (5k per class) **10k** test images (1k per class)



DR Discussion 2-How was this dataset created?

Xiaotong Chen Huijie Zhang Zeren Yu Stanley Lewis Odest Chadwicke Jenkins

Rough Pose Estimates from Pretrained Model

6D pose annotation through interactive interface









Human Annotator



ProgressLabeller: Visual Data Stream Annotation for Training Object-Centric 3D Perception

Fine-tuned Pose Estimates

Pose-based Robot Grasping





Idea:

- 1. Record video of scene
- 2. Human labels object pose in selected frames
- 3. Pose labels propagate to (large number of) remaining frames



Gradescope Quizzes

- Let course staff know if you have issues accessing
- Quiz links available through gradescope course 480760
- Time limit of 15 min once quiz is opened
- Each available to take from 7:00AM-3:00PM EST on quiz days
- Covers material from previous lectures and graded projects
- Today only: quiz 1 available until 6:00PM EST





Recap—Linear Classifiers

Algebraic Viewpoint

f(x,W) = Wx





Plot created using Wolfram Clou

master chef can



fish can











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



Recap—Loss Functions Quantify Preferences

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





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

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



Recap—Loss Functions Quantify Preferences

Q: How do we find the best W,b?

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

Linear classifier





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



Recap—Loss Functions Quantify Preferences

Problem: Loss functions encourage good performance on training data but we care about test data

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

Linear classifier





Regularization + Optimization





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





Overfitting

Both models have perfect accuracy on the training data!





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

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

$$s_i = w_i x + b_i$$

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

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



Overfitting

Both models have perfect accuracy on the training data!





1.0

0.8

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

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

$$s_{i} = w_{i}x + b_{i}$$

$$p_{i} = \frac{exp(s_{i})}{exp(s_{1}) + exp(s_{2})}$$

$$L = -\log(p_{y})$$

$$0.6$$

$$0.4$$

$$0.4$$

$$0.2$$

$$0.0$$



Overfitting





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







Overfitting

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



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





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

Data loss: Model predictions should match training data



 Hyperparameter giving regularization strength

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



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

Data loss: Model predictions should match training data

Simple examples:

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



 Hyperparameter giving regularization strength

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



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

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Simple examples:

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



 Hyperparameter giving regularization strength

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

More complex: Dropout Batch normalization Cutout, Mixup, Stochastic depth, etc...



Regularization: Prefer Simpler Models

Example: Linear classifier with 1D inputs, 2 classes, and 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: Prefer Simpler Models



$$w_{i}x + b_{i}$$

$$exp(s_{i})$$

$$exp(s_{1}) + exp(s_{2})$$

$$- log(p_{y}) + \lambda \sum_{i} w_{i}^{2}$$

$$ss = 5.96e-03$$

$$curacy = 1.0$$

$$p(y=0|x)$$

$$p(y=1|x)$$

$$x \text{ with } y=0$$

$$x \text{ with } y=1$$

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







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

Same predictions, so data loss will always be the same



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 $R(W) = \sum_{k,l} W_{k,l}^{2}$ L2 Regularization prefers weights to be

L2 Regularization prefers weights to be "spread out"

Same predictions, so data loss will always be the same



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

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



Finding a good W

$$f(x_i, W), y_i) + \lambda R(W)$$







DR

Optimization

$w^* = \arg\min L(w)$ W







The valley image and the walking man image are in CC0 1.0 public domain







The valley image and the walking man image are in CC0 1.0 public domain



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)</pre>

prints:

#	in	attempt	0	the	loss	Was	9.401632,	1
#	in	attempt	1	the	loss	was	8.959668,	1
#	in	attempt	2	the	loss	was	9.044034,	1
#	in	attempt	3	the	loss	was	9.278948,	1
#	in	attempt	4	the	loss	was	8.857370,	1
#	in	attempt	5	the	loss	was	8.943151,	1
#	in	attempt	6	the	loss	was	8.605604,	1
#	<pre># (trunctated:</pre>				continues for 1000			



best 9.401632 best 8.959668 best 8.959668 best 8.857370 best 8.857370 best 8.605604 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 on CIFAR-10! not bad!



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 on CIFAR-10! not bad! (SOTA is ~95%)





Idea #2: Follow the slope





The valley image and the walking man image are in CC0 1.0 public domain



Idea #2: Follow the slope

 $\frac{df}{dx} = \lim_{h \to 0} \frac{f}{dx}$



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

$$\frac{f(x+h)-f(x)}{h}$$



Idea #2: Follow the slope

- In 1-dimension, the derivative of a function gives the slope:
 - $\frac{df}{dx} = \lim_{h \to 0} \frac{f}{dx}$
- 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**.



$$\frac{f(x+h)-f(x)}{h}$$



Current W:

[0.34, -1.11,

0.78,

0.12,

0.55,

2.81,

- -3.1,
- -1.5,

0.33, ...]

loss 1.25347



 $\frac{dL}{dW}$ Gradient [?, ?, ?, ?, ?, ?, ?, ?, ?, ...]



Current W:

[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 + **0.0001**, -1.11, 0.78, 0.12, 0.55, 2.81, -3.1, -1.5, 0.33, …] loss 1.25322







Current W:

[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 + **0.0001**, -1.11, 0.78, 0.12, 0.55, 2.81, -3.1, -1.5, 0.33, …] loss 1.25322





?, ...]


[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

W + h (second dim):



 $\frac{dL}{dW}$ Gradient [-2.5, ?, ?, ?, ?, ?, ?, ?, ?, ...]



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



W + h (second dim): [0.34, -1.11 + **0.0001**, 0.78, 0.12, 0.55, 2.81, -3.1, -1.5, 0.33, …] loss 1.25353

Gradient
$$\frac{dL}{dW}$$

[-2.5,
0.6,
?,
?,
?,
?,
(1.25353 - 1.25347)/
0.0001
= 0.6
 $\frac{df}{dx} = \lim_{h \to 0} \frac{f(x+h) - f(x)}{h}$





[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 1.25353

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



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

W + h (third dim):

 $\frac{dL}{dW}$ Gradient [-2.5, 0.6, 0.0, ?, ?,

Numeric Gradient: - Slow: O(#dimensions) - Approximate





Loss is a function of W

 $L = \frac{1}{N} \sum_{i=1}^{N} L_i + \sum_{k=1}^{N} W_k^2$ $L_i = \sum \max(0, s_i - s_{y_i} + 1)$ $j \neq y_i$ s = f(x, W) = Wx

Want $\nabla_w L$





Loss is a function of W

 $L = \frac{1}{N} \sum_{i=1}^{N} L_i + \sum_{k=1}^{N} W_k^2$ $L_i = \sum \max(0, s_i - s_{v_i} + 1)$ $j \neq y_i$ s = f(x, W) = Wx

Want $\nabla_w L$



Use calculus to compute an analytic gradient



[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, 0.12, 0.55, 2.81, -3.1, -1.5, 0.33, ...] loss 1.25347

dW

In practice we will compute using back propagation; see Lecture 6







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



e, slow, easy to write error-prone



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

with numerical gradient. This is called a gradient check.



In practice: Always use analytic gradient, but check implementation



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

with numerical gradient. This is called a gradient check.

11 11 11 sample a few random elements and only return numerical in this dimensions.



- In practice: Always use analytic gradient, but check implementation
- def grad_check_sparse(f, x, analytic_grad, num checks=10, h=1e-7):



- 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)

that are of floating point type and with requires_grad=True.

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



[SOURCE]

- Check gradients computed via small finite differences against analytical gradients w.r.t. tensors in inputs





- 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=1e05, 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 grad_outputs are correct.



e, slow, easy to write error-prone





- 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

Hyperparameters:

- Weight initialization method
- Number of steps
- Learning rate



Iteratively step in the direction of the negative gradient (direction of local



steepest descent)





Iteratively step in the direction of the negative gradient (direction of local



steepest descent)





• Iteratively step in the direction of the negative gradient (direction of local



steepest descent)





• Iteratively step in the direction of the negative gradient (direction of local





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)

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

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



 $\nabla_W R(W)$

Full sum expensive when N is large!

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

Hyperparameters:

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



Stochastic Gradient Descent (SGD)

 $L(W) = \mathbb{E}_{(x,y) \sim p_{data}}[L(x, y, W)] + \lambda R(W)]$ $\approx \frac{1}{N} \sum_{i=1}^{N} L(x_i, y_i, W) + \lambda R(W)$

 $\approx \sum N \nabla_{w} L(x_{i}, y_{i}, W) + \nabla_{w} \lambda R(W)$ i=1



Think of loss as an expectation over the full data distribution Pdata

Approximate expectation via sampling

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





Interactive Web Demo



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



0,0] ▲	W[0,1]	b[0]	x [0]	x[1]	У	s[0]	s[1]	s[2]	L
. 30	2.00	0.00	0.50	0.40	0	1.95	-0.10	0.60	0.00
V	V	V	0.80	0.30	0	2.44	0.90	1.60	0.16
⊥,0] ▲	w[1,1]	b[1]	0.30	0.80	0	2.29	-2.10	-0.40	0.00
.00 .42	- 4 .00 -0.61	0.50 -0.22	-0.40	0.30	1	-0.32	-1.50	-2.00	2.68
▼ 2,0]	▼ W[2,1]	▼ b[2]	-0.30	0.70	1	0.71	-2.90	-2.10	6.41
A	A	<u>۸</u>	-0.70	0.20	1	-1.21	-1.70	-2.80	1.49
.12	0.32	-0.11	0.70	-0.40	2	0.81	3.50	2.00	2.50
V	•	•	0.50	-0.60	2	-0.05	3.90	1.60	3.30
p size: 0.09976			-0.40	-0.50	2	-1.92	1.70	-1.20	4.18
Single	parameter	rupdate							mean:
Start	repeated u	update	Tota Regi	Total data loss: 2.30 Regularization loss: 3.93					
Stop	repeated u	update	Tota	Total loss: 6.23					

Randomize parameters

L2 Regularization strength: 0.10000



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



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





What does gradient decent do?



the Hessian matrix is large



- What if loss changes quickly in one direction and slowly in another?
- Very slow progress along shallow dimension, jitter along steep direction

Loss function has high condition number: ratio of largest to smallest singular value of



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







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

Zero gradient, gradient descent gets stuck







Our gradients come from mini batches so they can be noisy!

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







Our gradients come from mini batches so they can be noisy!

$$L(W) = \frac{1}{N} \bigvee_{i=1}^{N} \forall anilla \text{ gradient descent} \\ w = initialize_weights() \\ \text{for t in range(num_steps):} \\ dw = compute_gradient(loss_w -= learning_rate * dw \\ N \swarrow_{i=1}^{M}$$









Our gradients come from mini batches so they can be noisy!

$$L(W) = \frac{1}{N} \bigvee_{i=1}^{N} \forall anilla \text{ gradient descent} \\ w = initialize_weights() \\ \text{for t in range(num_steps):} \\ dw = compute_gradient(loss_w -= learning_rate * dw \\ N \swarrow_{i=1}^{M}$$







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

Batched gradient descent always computes same gradients

SGD computes noisy gradients, may help to escape saddle points









SGD

$$w_{t+1} = w_t - \alpha \nabla L(w_t)$$

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







SGD

$$w_{t+1} = w_t - \alpha \nabla L(w_t)$$

for t in range(num steps): for t in range(num_steps): dw = $dw = compute_gradient(w)$ W -= w -= learning_rate * dw





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





Momentum update:



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



SGD + Momentum

$$v_{t+1} = \rho v_t + \nabla L(w_t)$$

$$w_{t+1} = w_t - \alpha v_{t+1}$$
):

$$v = 0$$
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





SGD + Momentum

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

 $w_{t+1} = w_t + v_{t+1}$
 $v = 0$
for t in range(num_steps):
dw = compute_gradient(w)
v = rho * v - learning_rate * dw
w += v

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



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





Local Minima







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

SGD + Momentum

Gradient Noise



Local Minima







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

SGD + Momentum

Gradient Noise



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

SGD + Momentum

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


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

Nesterov Momentum

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



$v_{t+1} = \rho v_t - \alpha \nabla L (w_t + \rho v_t)$ $w_{t+1} = w_t + v_{t+1}$



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



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

$$w_{t+1} = w_t + v_{t+1}$$

Change of variables
and rearrange:

$$\tilde{w}_t = w_t + \rho v_t$$

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

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

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



Nesterov Momentum

Annoying, usually we want to update in terms of W_t , $\nabla L(W_t)$







AdaGrad

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 sub gradient methods for online learning and stochastic optimization," JMLR 2011

squared.sqrt() + 1e-7)



AdaGrad

00





Q: What happens with AdaGrad?



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



AdaGrad





Q: What happens with AdaGrad?



Problem: AdaGrad will slow over many iterations

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







_ 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) w -= learning_rate * dw / (grad_squared.sqrt() + 1e-7)







moment1 = 0moment2 = 0for t in range(1, num_steps + 1): # Start at t = 1 dw = compute_gradient(w) moment1 = beta1 * moment1 + (1 - beta1) * dwmoment2 = beta2 * moment2 + (1 - beta2) * dw * dww -= learning_rate * moment1 / (moment2.sqrt() + 1e-7)





moment1 = 0moment2 = 0for t in range(1, num_steps + 1): # St dw = compute_gradient(w) moment1 = beta1 * moment1 + (1 - beta1) * dwmoment2 = beta2 * moment2 + (1 - beta2) * dw * dww -= learning_rate * moment1 / (moment2.sqrt() + 1e-7)

ge(num_steps): te_gradient(w) ing_rate * dw

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



art at t =
$$1$$

Adam

Momentum

```
-= learning_rate * v
```

SGD+Momentum



moment1 = 0moment1 = 0moment2 = 0for t in range(1, num_steps + 1): # Sta dw = compute_gradient(w) moment1 = beta1 * moment1 + (1 - beta moment2 = beta2 * moment2 + (1 - beta w -= learning_rate * moment1 / (moment

grad_squared = 0 RMSProp 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)







moment1 = 0moment2 = 0for t in range(1, num_steps + 1): # Start at t = 1 dw = compute_gradient(w) moment1 = beta1 * moment1 + (1 - beta1) * dwmoment2 = beta2 * moment2 + (1 - beta2) * dw * dww -= learning_rate * moment1 / (moment2.sqrt() + 1e-7)

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



Adam Momentum AdaGrad / RMSProp **Bias correction**







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



Adam with beta 1 = 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

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 threshsampled 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.

Gkioxari, Malik, and Johnson, ICCV 2019

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



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

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

Adam with beta 1 = 0.9,

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





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	\checkmark	X	X	X
Nesterov	\checkmark	X	X	X
AdaGrad	X	\checkmark	X	X
RMSProp	X	\checkmark	\checkmark	X
Adam	\checkmark	\checkmark	\checkmark	\checkmark





L2 Regularization vs Weight Decay

Optimization Algorithm

 $L(w) = L_{data}(w) + L_{reg}(w)$ $g_t = \nabla L(w_t)$ $s_t = optimizer(g_t)$ $W_{t+1} = W_t - \alpha S_t$

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

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

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

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

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

Optimization Algorithm $L(w) = L_{data}(w)$ $g_t = \nabla L_{data}(w_t)$ $s_t = optimizer(g_t) + 2\lambda w_t$ $W_{t+1} = W_t - \alpha S_t$

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



AdamW: Decouple Weight Decay

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

- 1: given $\alpha = 0.001, \beta_1 = 0.9, \beta_2 = 0.999, \epsilon = 10^{-8}, \lambda \in \mathbb{R}$
- vector $\mathbf{v}_{t=0} \leftarrow \mathbf{0}$, schedule multiplier $\eta_{t=0} \in \mathbb{R}$

3: repeat

- 4: $t \leftarrow t+1$ 5: $\nabla f_t(\boldsymbol{\theta}_{t-1}) \leftarrow \text{SelectBatch}(\boldsymbol{\theta}_{t-1})$
- 6: $\boldsymbol{g}_t \leftarrow \nabla f_t(\boldsymbol{\theta}_{t-1}) + \lambda \boldsymbol{\theta}_{t-1}$
- 7: $\boldsymbol{m}_t \leftarrow \beta_1 \boldsymbol{m}_{t-1} + \overline{(1-\beta_1)\boldsymbol{g}}_t$ 8: $\mathbf{v}_t \leftarrow \beta_2 \mathbf{v}_{t-1} + (1 - \beta_2) \mathbf{g}_t^2$
- 9: $\hat{\boldsymbol{m}}_t \leftarrow \boldsymbol{m}_t / (1 \beta_1^t)$
- 10: $\hat{\mathbf{v}}_t \leftarrow \mathbf{v}_t / (1 \beta_2^t)$
- 11: $\eta_t \leftarrow \text{SetScheduleMultiplier}(t)$
- 12: $\boldsymbol{\theta}_t \leftarrow \boldsymbol{\theta}_{t-1} \eta_t \left(\alpha \hat{\boldsymbol{w}}_t / (\sqrt{\hat{\boldsymbol{v}}_t} + \epsilon) + \lambda \boldsymbol{\theta}_{t-1} \right)$
- 13: **until** stopping criterion is met
- 14: return optimized parameters θ_t



2: initialize time step $t \leftarrow 0$, parameter vector $\theta_{t=0} \in \mathbb{R}^n$, first moment vector $m_{t=0} \leftarrow 0$, second moment

▷ select batch and return the corresponding gradient

▷ here and below all operations are element-wise

 $\triangleright \beta_1$ is taken to the power of t

 $\triangleright \beta_2$ is taken to the power of t

▷ can be fixed, decay, or also be used for warm restarts



AdamW: Decouple Weight Decay

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

- 1: given $\alpha = 0.001, \beta_1 = 0.9, \beta_2 = 0.999, \epsilon = 10^{-8}, \lambda \in \mathbb{R}$
- vector $\mathbf{v}_{t=0} \leftarrow \mathbf{0}$, schedule multiplier $\eta_{t=0} \in \mathbb{R}$

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

12: $\boldsymbol{\theta}_t \leftarrow \boldsymbol{\theta}_{t-1} - \eta_t \left(\alpha \hat{\boldsymbol{w}}_t / (\sqrt{\hat{\boldsymbol{v}}_t} + \epsilon) + \lambda \boldsymbol{\theta}_{t-1} \right)$

13: **until** stopping criterion is met 14: return optimized parameters θ_t



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

2: initialize time step $t \leftarrow 0$, parameter vector $\theta_{t=0} \in \mathbb{R}^n$, first moment vector $m_{t=0} \leftarrow 0$, second moment









DR

So far: First-order Optimization





So far: First-order Optimization

1. Use gradient to make linear approximation 2. Step to minimize the approximation w1















DR

Second-order Optimization

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

Take bigger steps in areas of low curvature

w1





Second-order Taylor Expansion:

$$L(w) \approx L(w_0) + (w - w_0)^T \nabla_w L(w_0) + \frac{1}{2} (w - w_0)^T H_w L(w_0) (w - w_0)$$

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:

$$L(w) \approx L(w_0) + (w - w_0)^T \nabla_w L(w_0) + \frac{1}{2} (w - w_0)^T H_w L(w_0) (w - w_0)$$

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



- time ($O(n^2)$ each).
- Hessian



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

- Quasi-Newton methods (BGFS most popular): instead of inverting the Hessian ((O(n^3)), approximate inverse Hessian with rank 1 updates over

- L-BFGS (Limited memory BFGS): Does not form/store the full inverse



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.





In practice:

- Adam is a good default choice in many cases more tuning.
- (and don't forget to disable all sources of noise)



SGD+Momentum can outperform Adam but may require

If you can afford to do full batch updates then try out L-BFGS



- Use Linear Models for image classification problems.
- Use Loss Functions to express preferences over different choices of weights.
- Use **Regularization** to prevent overfitting to training data.
- for t in range(num_steps): scent t(dw = compute_gradient(w) and t w -= learning_rate * dw



Summary







Next time: Neural Networks

















Lecture 4 **Regularization + Optimization**





