

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

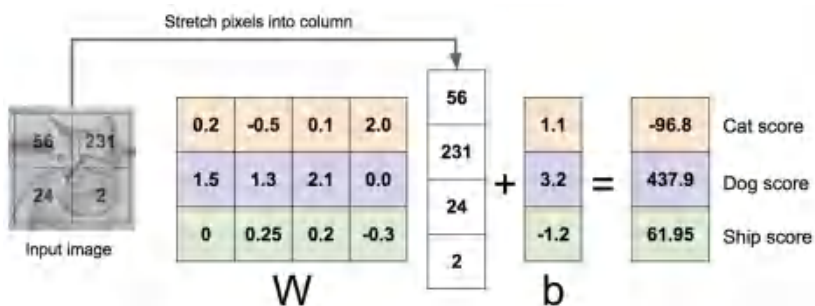
Assignment 2

- 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

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



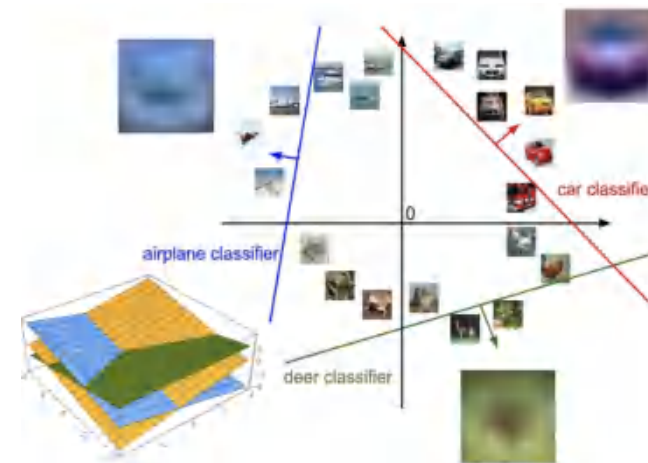
Visual Viewpoint

One template per class



Geometric Viewpoint

Hyperplanes cutting up space



Last Time: Loss Functions quantify preferences

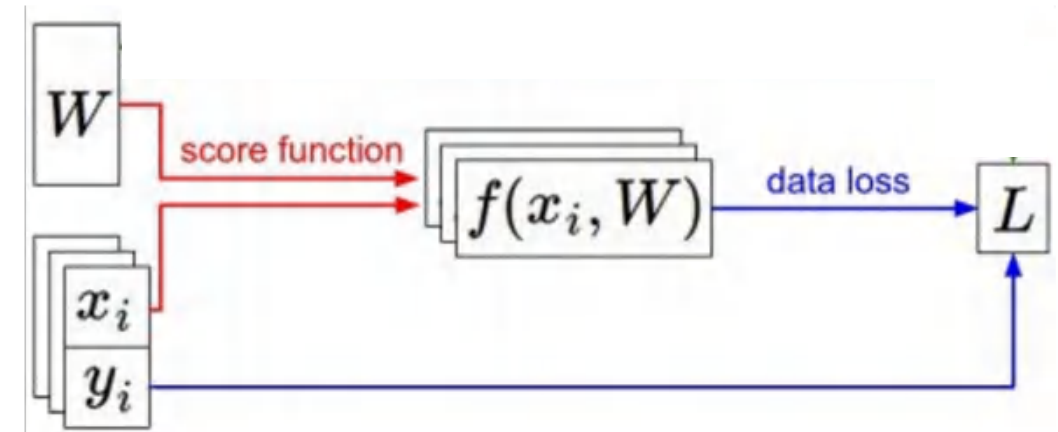
- We have some dataset of (x, y)
- We have a **score function**:
- We have a **loss function**:

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)
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- We have a **loss function**:

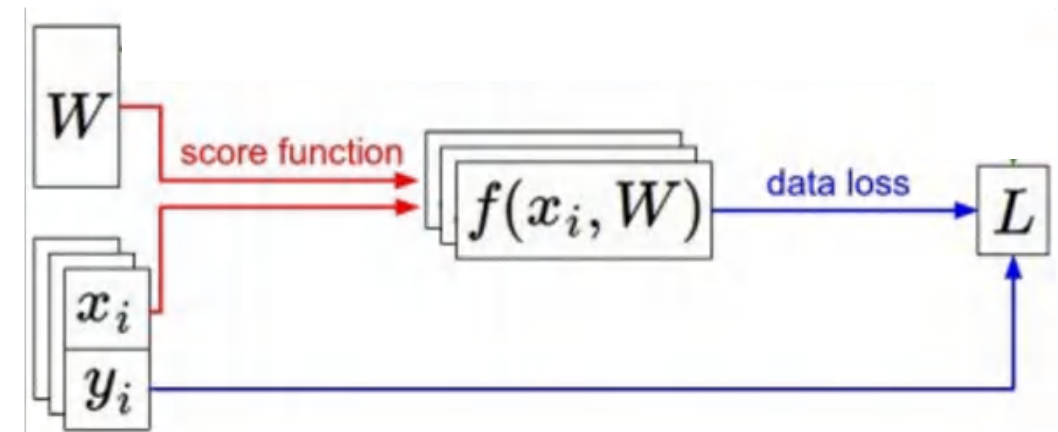
Problem: Loss functions encourage good performance on training data but we really care about test data

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

Linear classifier

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

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



Overfitting

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

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

$$s_i = w_i x + b_i$$

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

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

Overfitting

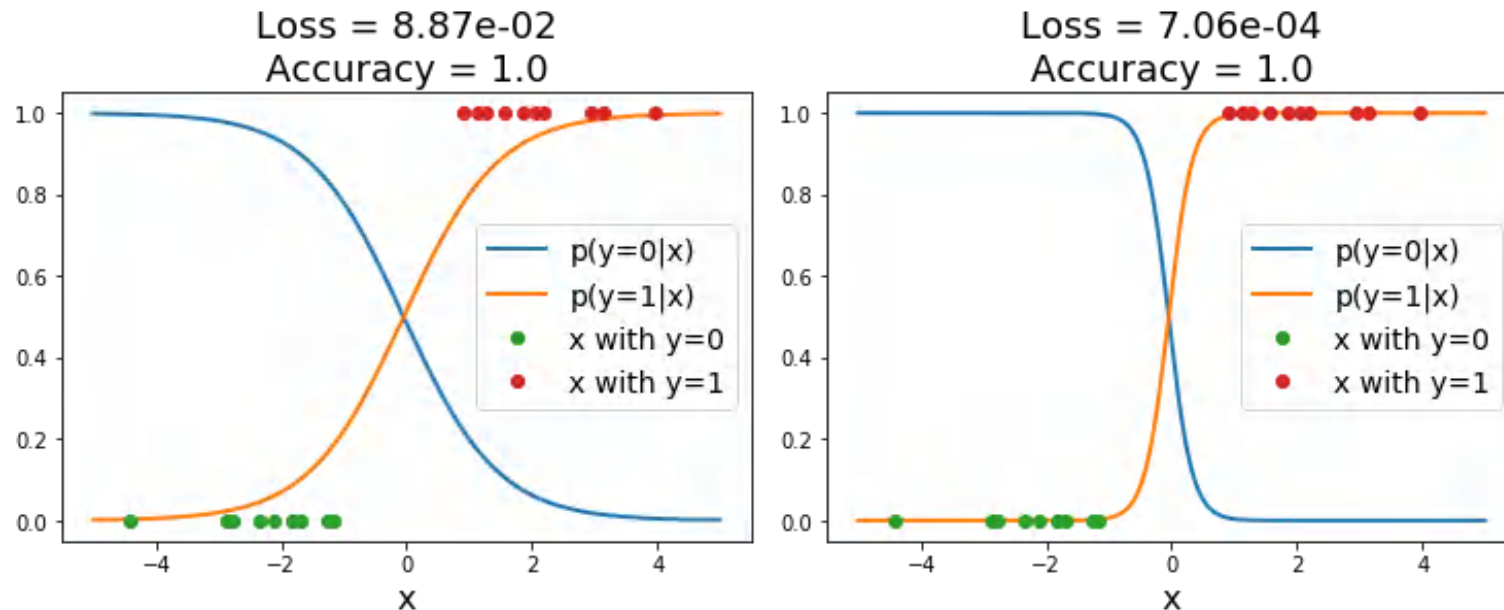
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Both models have perfect accuracy on train data!

Overfitting

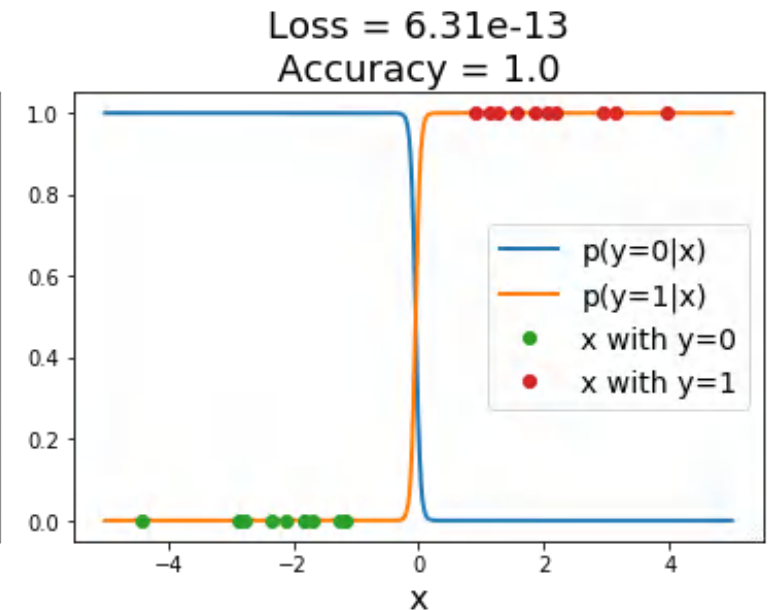
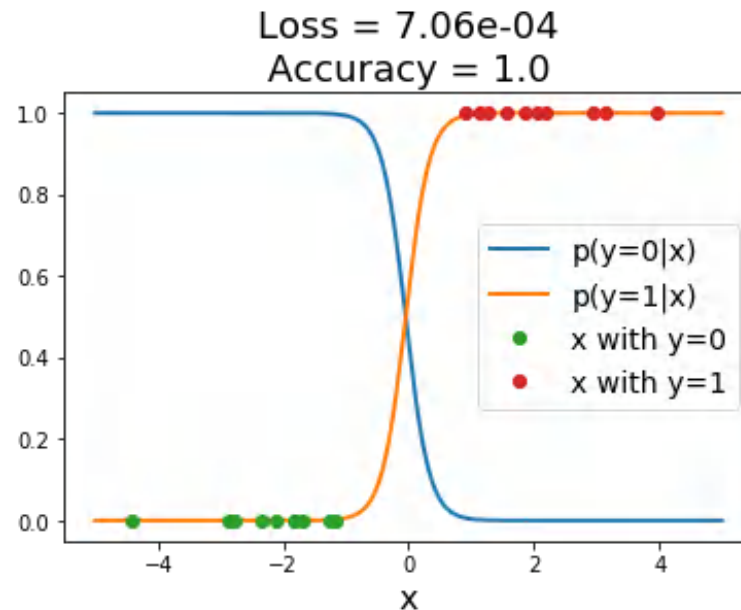
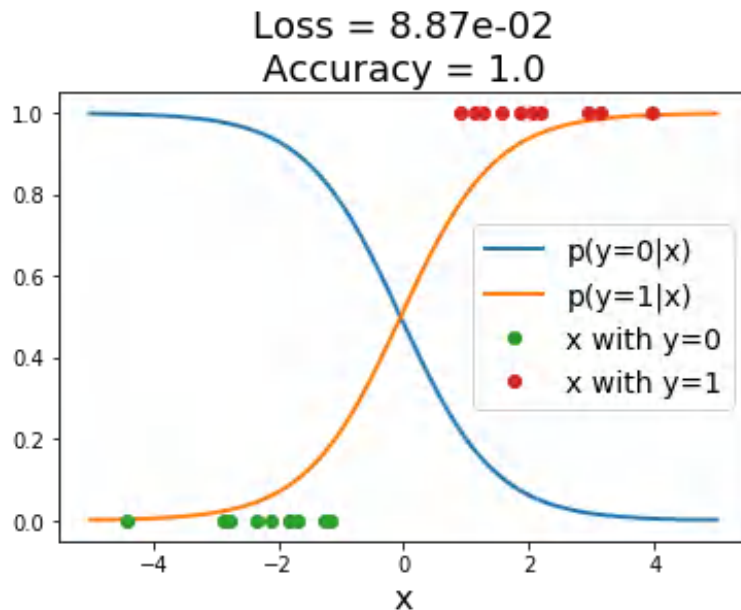
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Both models have perfect accuracy on train data!

Low loss, but unnatural “cliff”
between training points

Overfitting

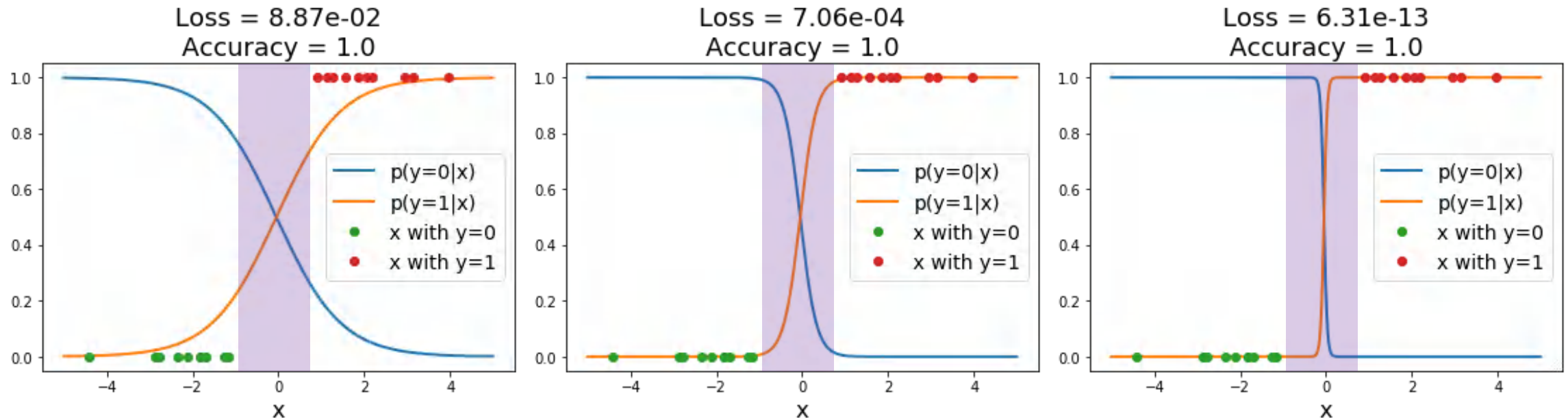
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$$L = -\log(p_y)$$



Overconfidence in regions with no training data could give poor generalization

Regularization: Beyond Training Error

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


Data loss: Model predictions
should match training data

Regularization: Beyond Training Error

$$L(W) = \underbrace{\frac{1}{N} \sum_{i=1}^N L_i(f(x_i, W), y_i)}_{\text{Data loss: Model predictions should match training data}} + \underbrace{\lambda R(W)}_{\text{Regularization: Prevent the model from doing too well on training data}}$$

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

Regularization: Beyond Training Error

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Simple examples

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

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

Regularization: Beyond Training Error

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Simple examples

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

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

More complex:

Dropout

Batch normalization

Cutout, Mixup, Stochastic depth, etc...

Regularization: Prefer Simpler Models

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

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

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

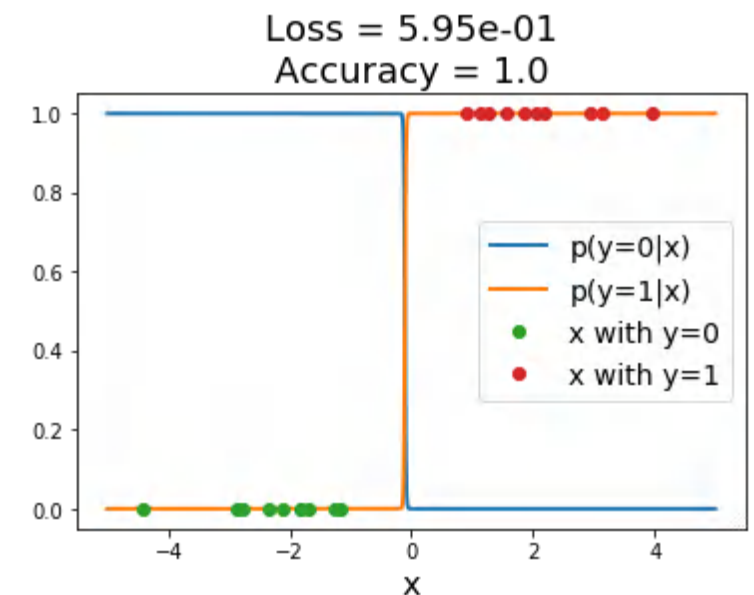
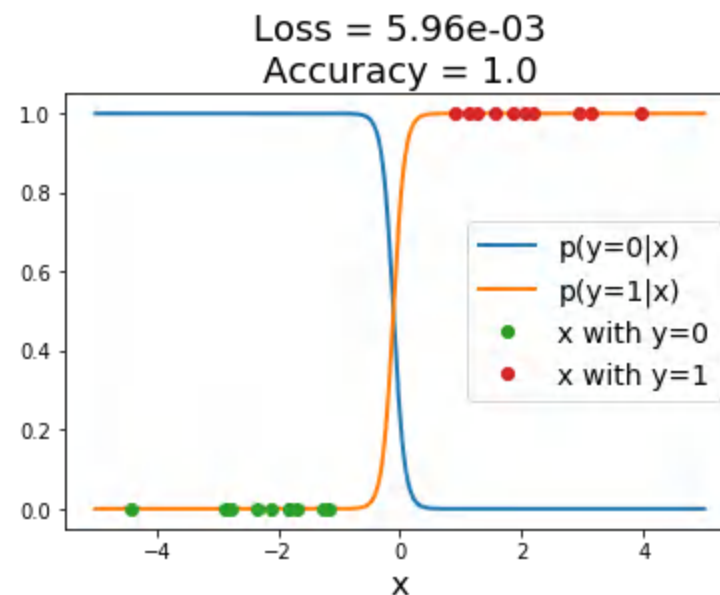
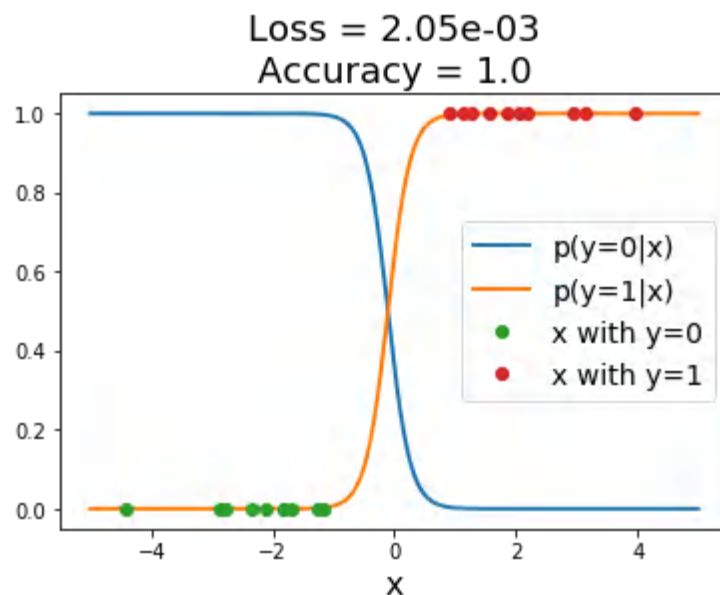
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$$L = -\log(p_y) + \lambda \sum_i w_i^2$$

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



Regularization: Expressing Preferences

L2 Regularization

$$\begin{aligned}x &= [1, 1, 1, 1] \\w_1 &= [1, 0, 0, 0] \\w_2 &= [0.25, 0.25, 0.25, 0.25]\end{aligned}$$

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

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

Same predictions, so data loss will always be the same

Regularization: Expressing Preferences

L2 Regularization

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

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

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

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

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
# ... (truncated: 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:

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

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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**

current W:

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

loss 1.25347

gradient dL/dW :

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

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

gradient dL/dW:

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

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

gradient dL/dW:

[-2.5,
?,
?,
↑

$$(1.25322 - 1.25347)/0.0001 = -2.5$$

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

?,
?,...]

current W:

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

loss 1.25347

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 dL/dW:

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

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 (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 dL/dW:

[-2.5,
0.6,
?,
?,
↑

$$(1.25353 - 1.25347)/0.0001 = 0.6$$

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

?,...]

current W:

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

loss 1.25347

W + h (third dim):

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

loss 1.25347

gradient dL/dW:

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

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 (third dim):

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

loss 1.25347

gradient dL/dW:

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

$$(1.25347 - 1.25347)/0.0001 = 0.0$$

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

current W:

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

loss 1.25347

W + h (third dim):

[0.34,
-1.11,
0.78 + **0.0001**,
0.12,
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2.81,
-3.1,
-1.5,
0.33,...]

loss 1.25347

gradient dL/dW:

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

Numeric Gradient:

- Slow: $O(\text{\#dimensions})$
- 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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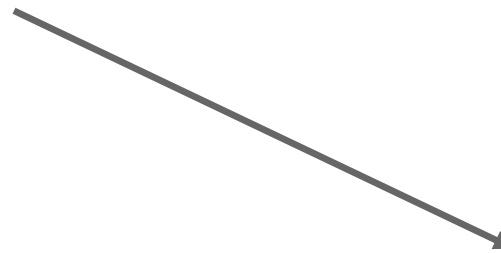
Use calculus to compute an **analytic gradient**

current W:

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

loss 1.25347

$dL/dW = \dots$
(some function
data and W)



gradient dL/dW :

[-2.5,
0.6,
0,
0.2,
0.7,
-0.5,
1.1,
1.3,
-2.1,...]

current W:

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

loss 1.25347

$dL/dW = \dots$
(some function
data and W)

(In practice we will
compute dL/dW using
backpropagation; see
Lecture 6)

gradient dL/dW :

[-2.5,
0.6,
0,
0.2,
0.7,
-0.5,
1.1,
1.3,
-2.1,...]

Computing Gradients

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

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

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In practice: 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.  
    """
```

Computing Gradients

- **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()`.

Computing Gradients

- **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, nondet_tol=0.0)
```

[SOURCE]

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.

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
```

Hyperparameters:

- Weight initialization method
- Number of steps
- Learning rate

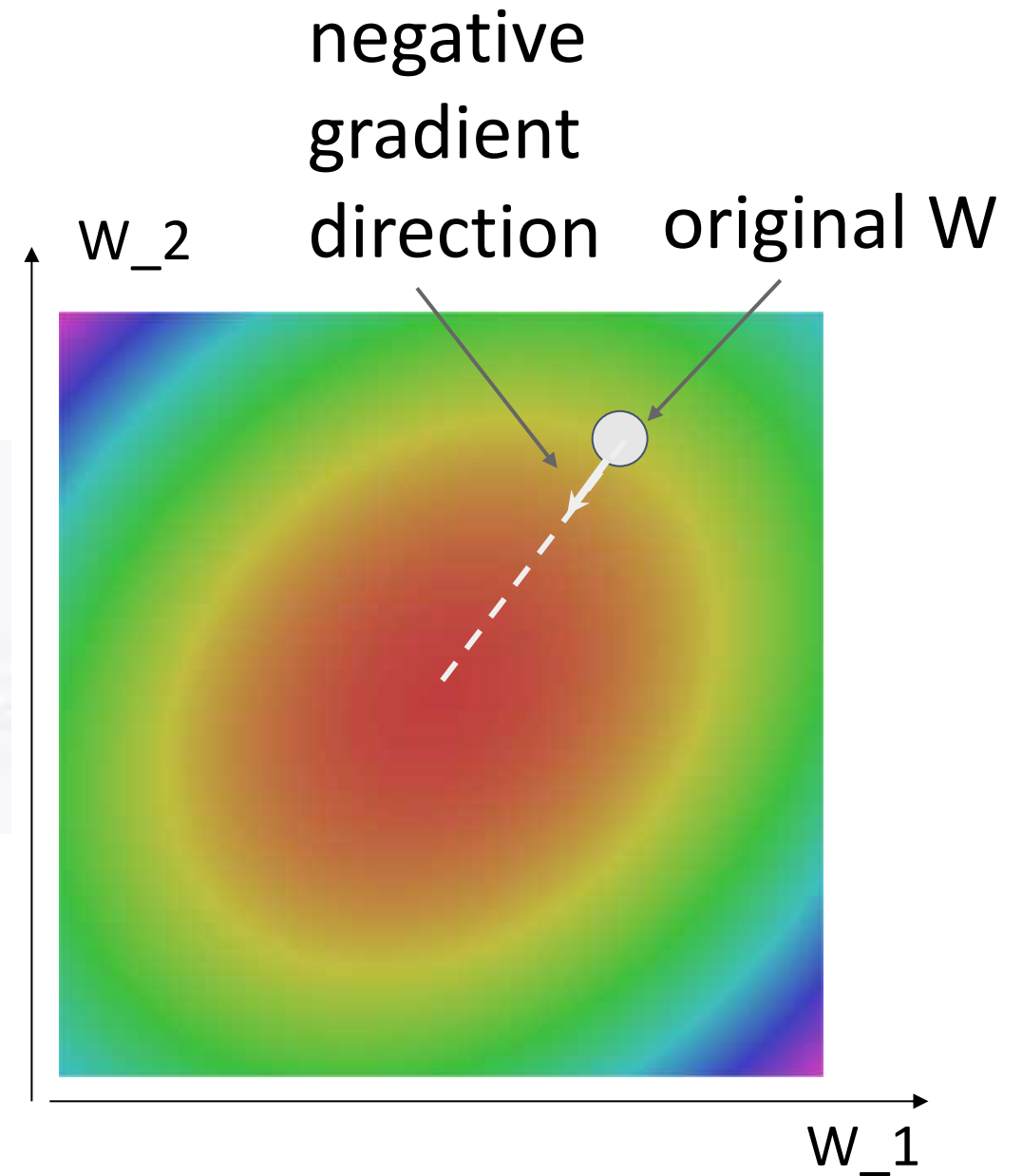
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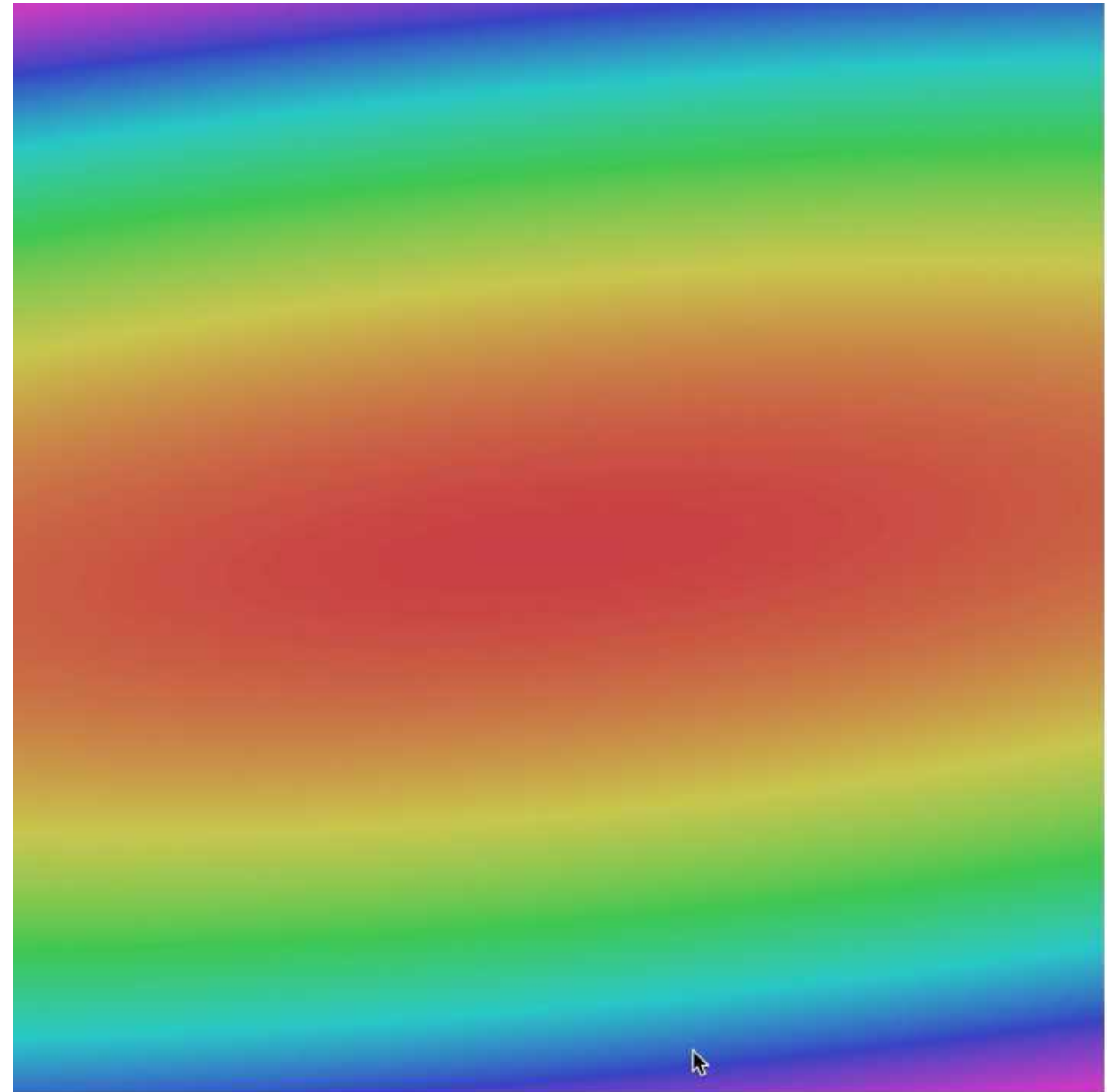
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Hyperparameters:

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- 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
```

Hyperparameters:

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

Stochastic Gradient Descent (SGD)

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

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

Approximate expectation via sampling

Stochastic Gradient Descent (SGD)

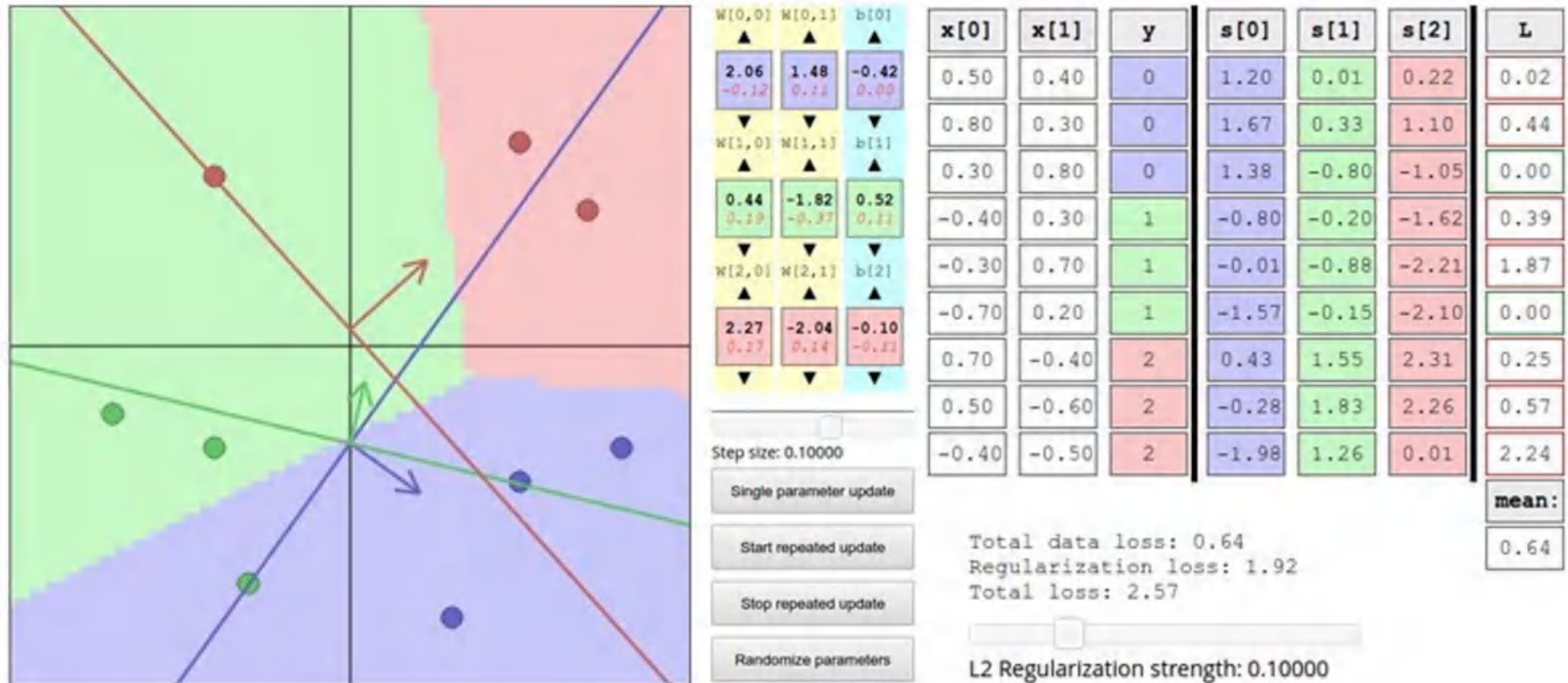
$$\begin{aligned} L(W) &= \mathbb{E}_{(x,y) \sim p_{data}} [L(x, y, W)] + \lambda R(W) \\ &\approx \frac{1}{N} \sum_{i=1}^N L(x_i, y_i, W) + \lambda R(W) \end{aligned}$$

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

Approximate expectation via sampling

$$\begin{aligned} \nabla_W L(W) &= \nabla_W \mathbb{E}_{(x,y) \sim p_{data}} [L(x, y, W)] + \lambda \nabla_W R(W) \\ &\approx \sum_{i=1}^N \nabla_w L_W(x_i, y_i, W) + \nabla_w R(W) \end{aligned}$$

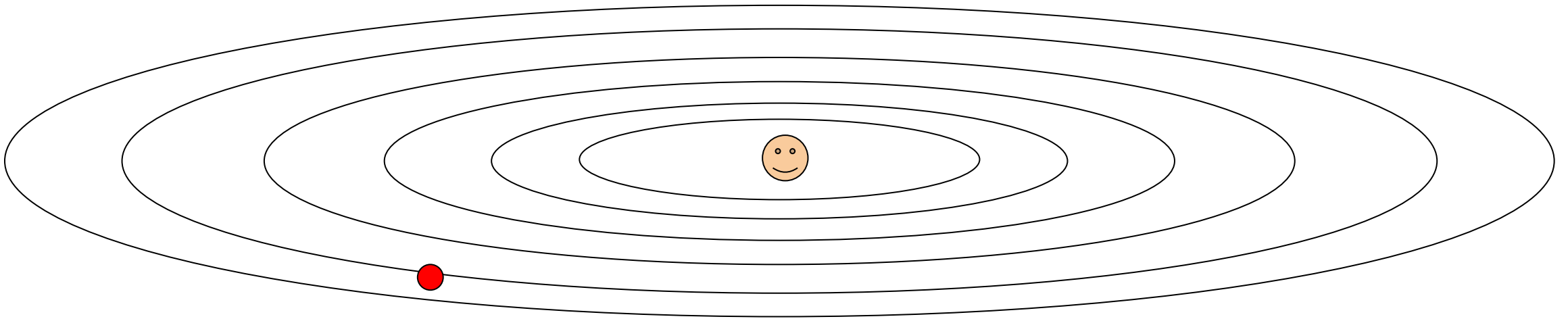
Interactive Web Demo



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

Problems with SGD

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



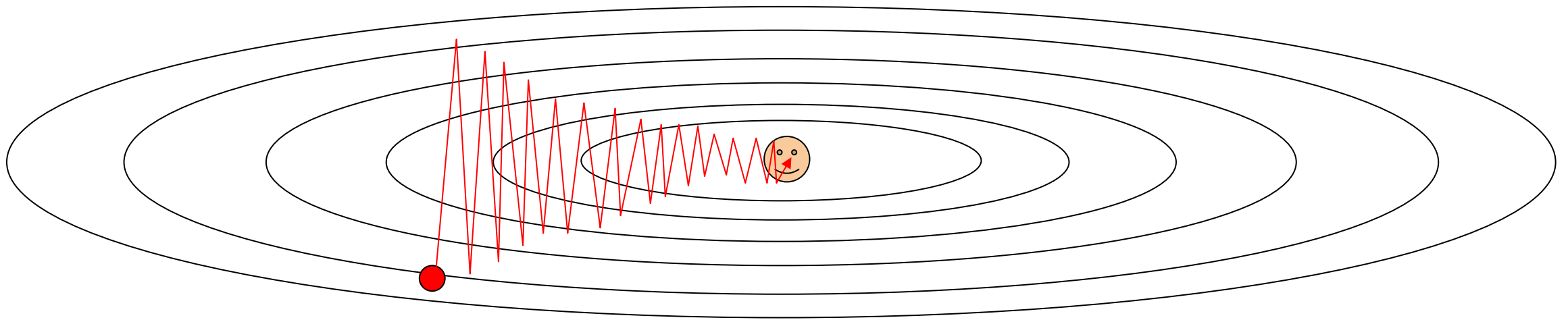
Loss function has high **condition number**: ratio of largest to smallest singular value of the Hessian matrix is large

Problems with SGD

What if loss changes quickly in one direction and slowly in another?

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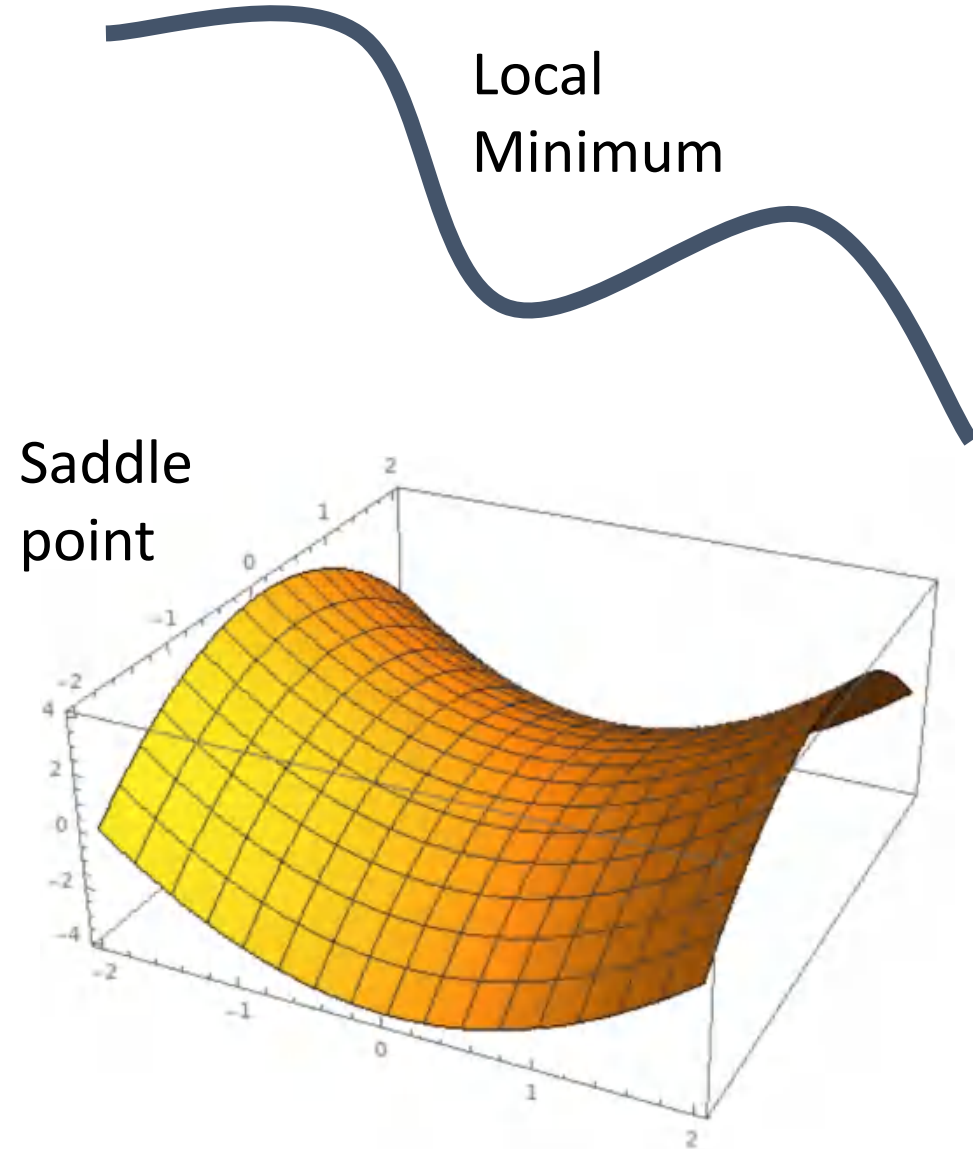
Very slow progress along shallow dimension, jitter along steep direction



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

Problems with SGD

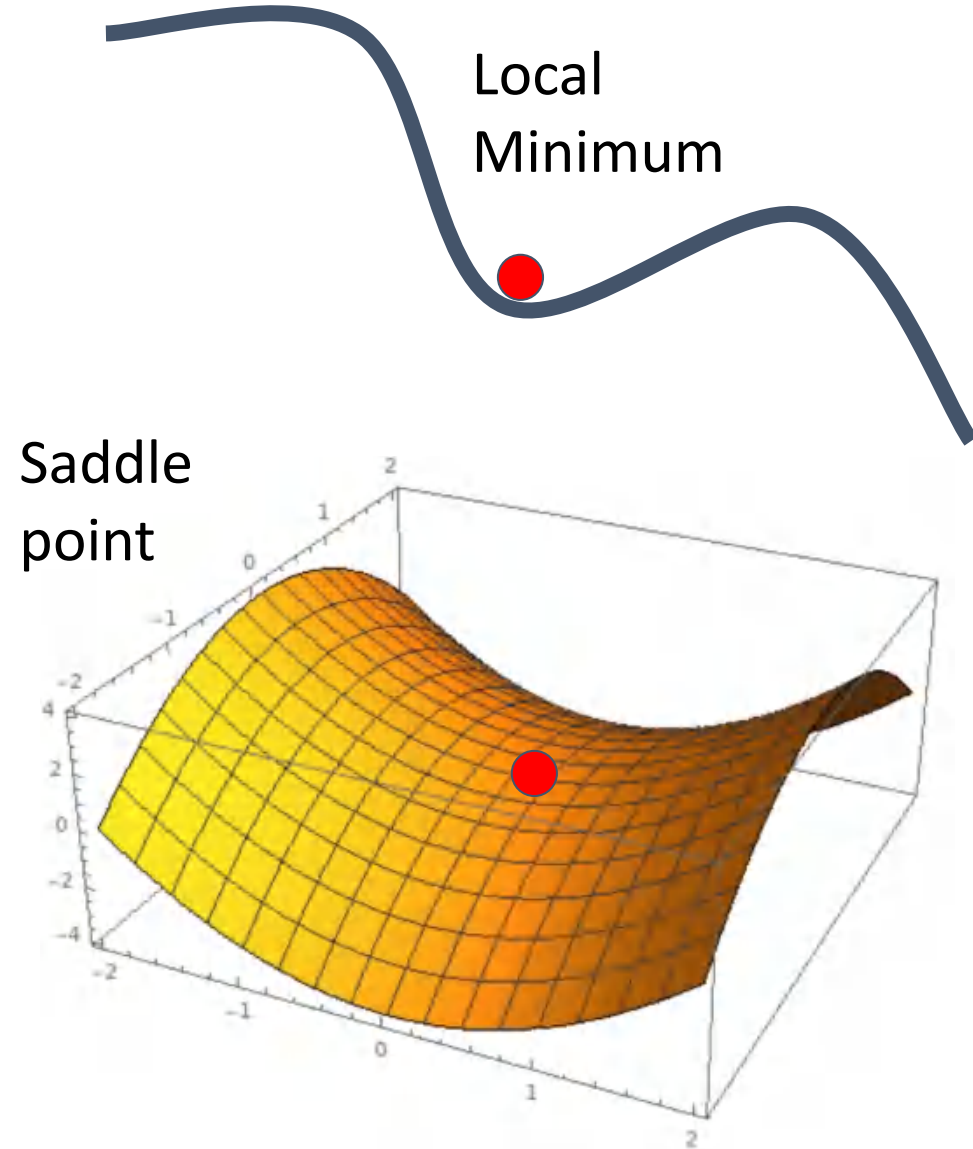
What if the loss function has a **local minimum** or **saddle point**?



Problems with SGD

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

Zero gradient, gradient descent gets stuck

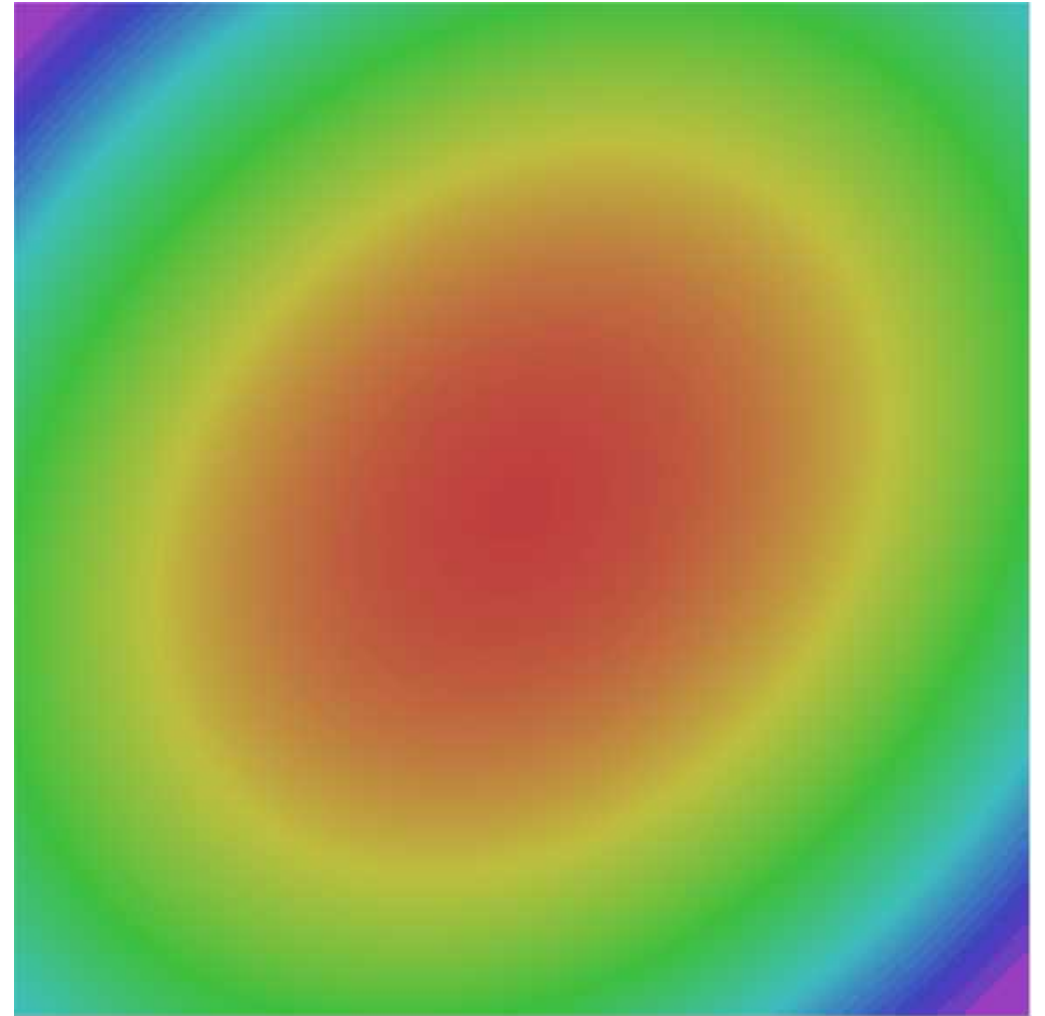


Problems with SGD

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 + Momentum

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

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

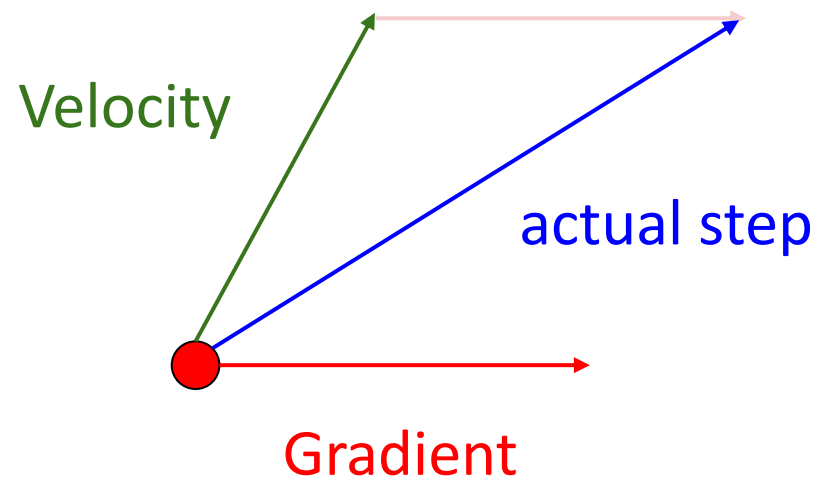
$$x_{t+1} = x_t - \alpha v_{t+1}$$

```
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

SGD + Momentum

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

SGD + Momentum

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

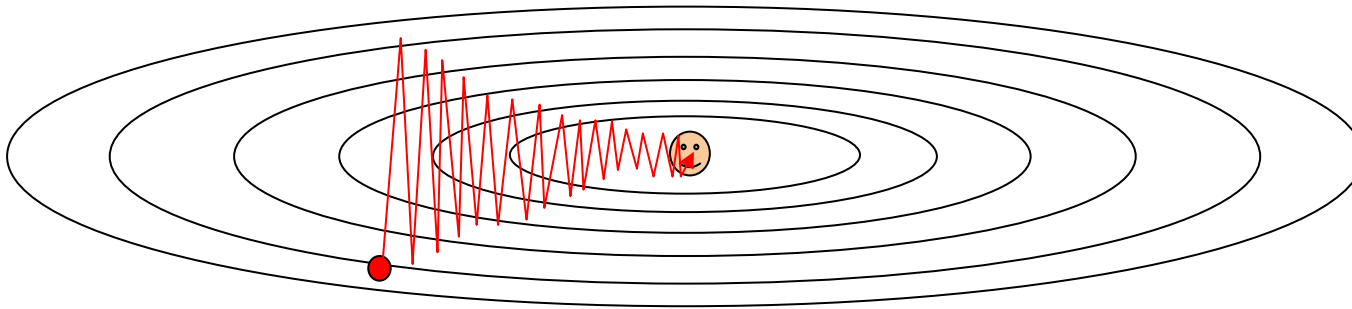
SGD + Momentum

Local Minima

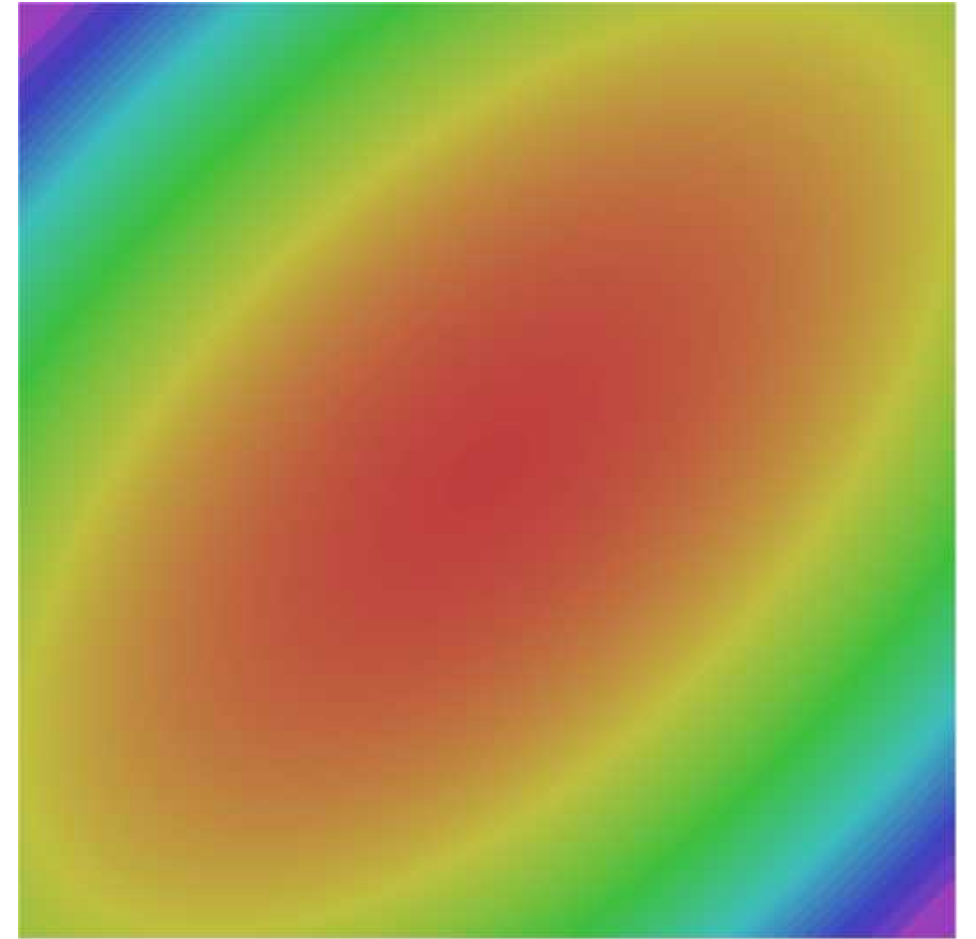
Saddle points



Poor Conditioning



Gradient Noise

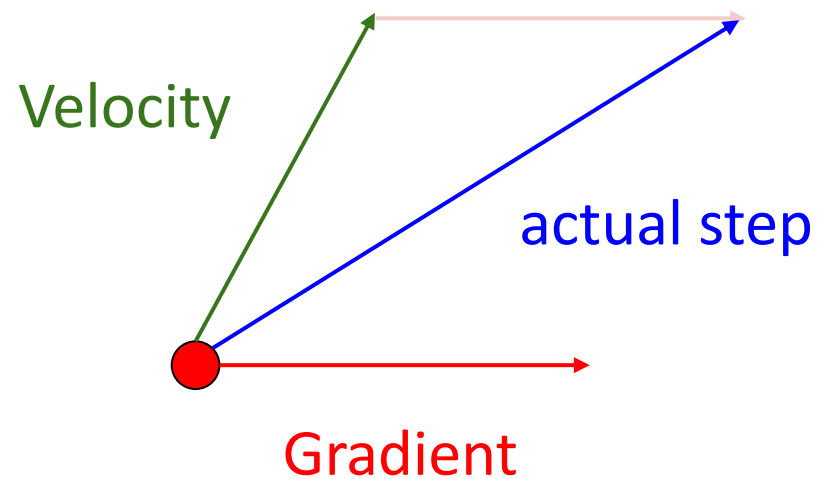


— SGD — SGD+Momentum

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

SGD + Momentum

Momentum update:



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

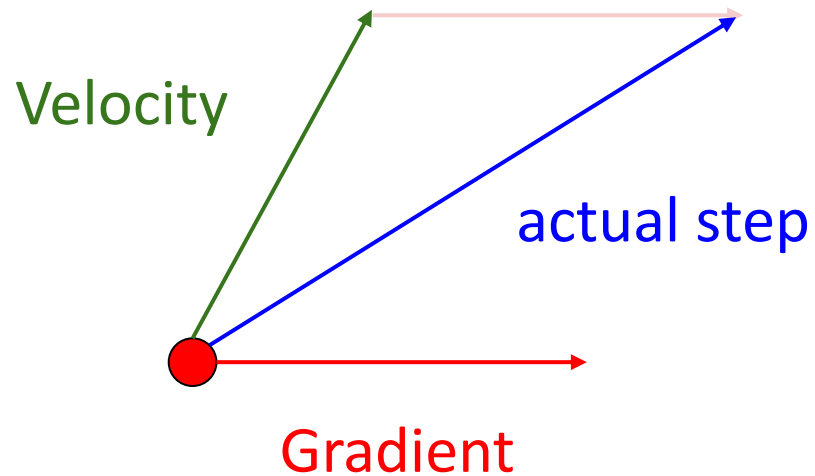
Nesterov, "A method of solving a convex programming problem with convergence rate $O(1/k^2)$ ", 1983

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

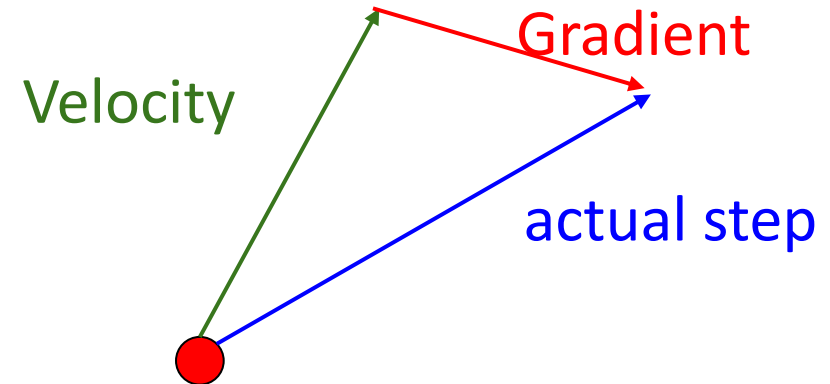
Nesterov Momentum

Momentum update:



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

Nesterov Momentum



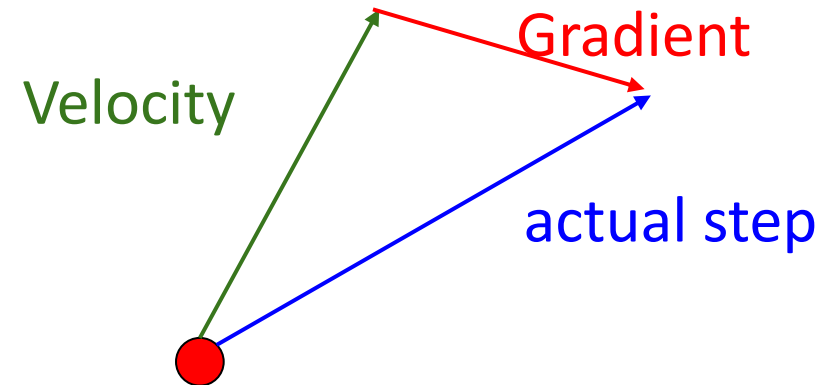
“Look ahead” to the point where updating using velocity would take us; compute gradient there and mix it with velocity to get actual update direction

Nesterov, “A method of solving a convex programming problem with convergence rate $O(1/k^2)$ ”, 1983
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Nesterov Momentum

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

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



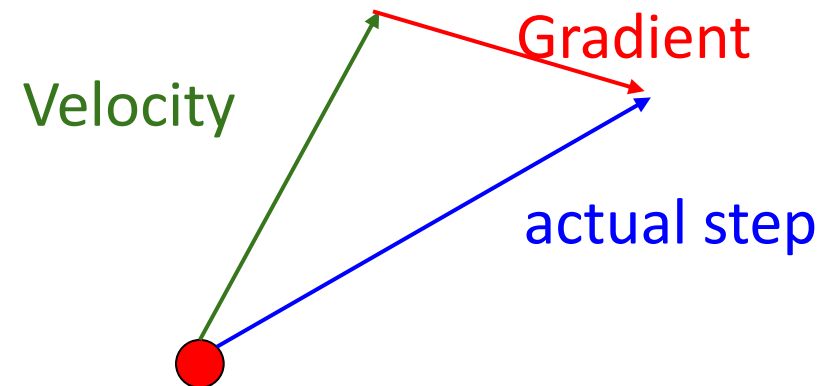
“Look ahead” to the point where updating using velocity would take us; compute gradient there and mix it with velocity to get actual update direction

Nesterov Momentum

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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

Nesterov Momentum

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

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

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

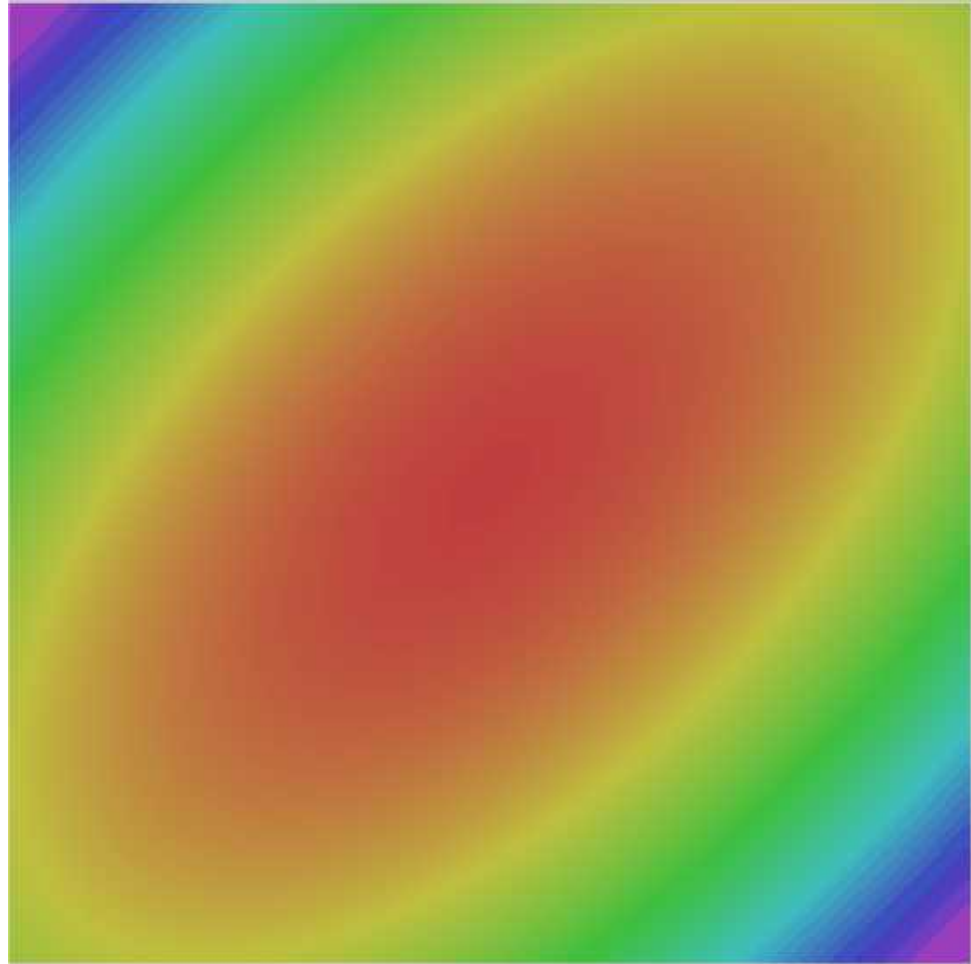
$$v_{t+1} = \rho v_t - \alpha \nabla f(\tilde{x}_t)$$

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

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

```
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
```

Nesterov Momentum



— SGD

— SGD+Momentum

— Nesterov

AdaGrad

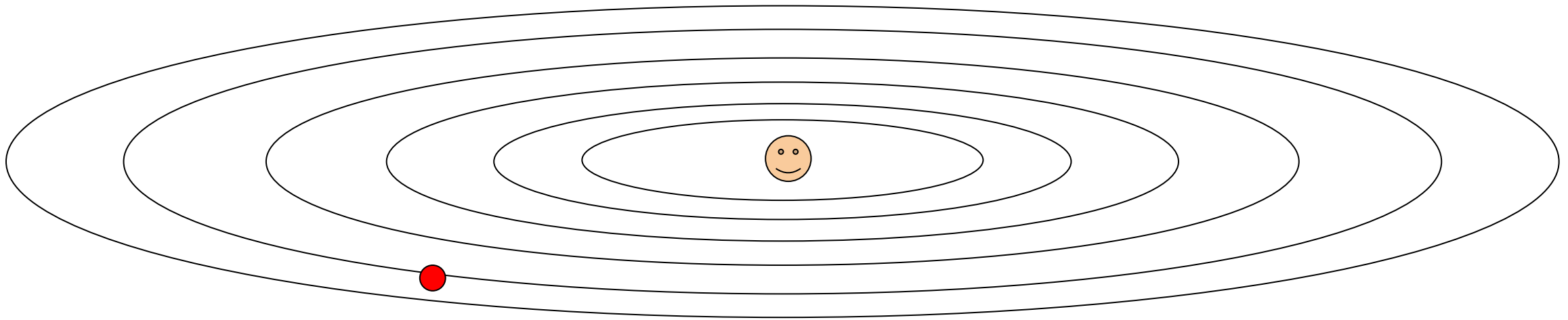
```
grad_squared = 0
for t in range(num_steps):
    dw = compute_gradient(w)
    grad_squared += dw * dw
    w -= learning_rate * dw / (grad_squared.sqrt() + 1e-7)
```

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”

AdaGrad

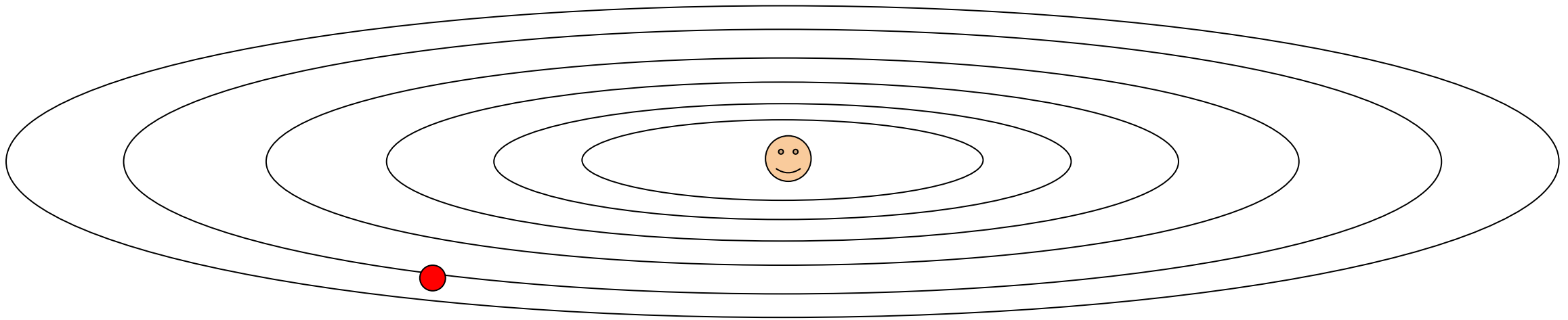
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    w -= learning_rate * dw / (grad_squared.sqrt() + 1e-7)
```



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

AdaGrad

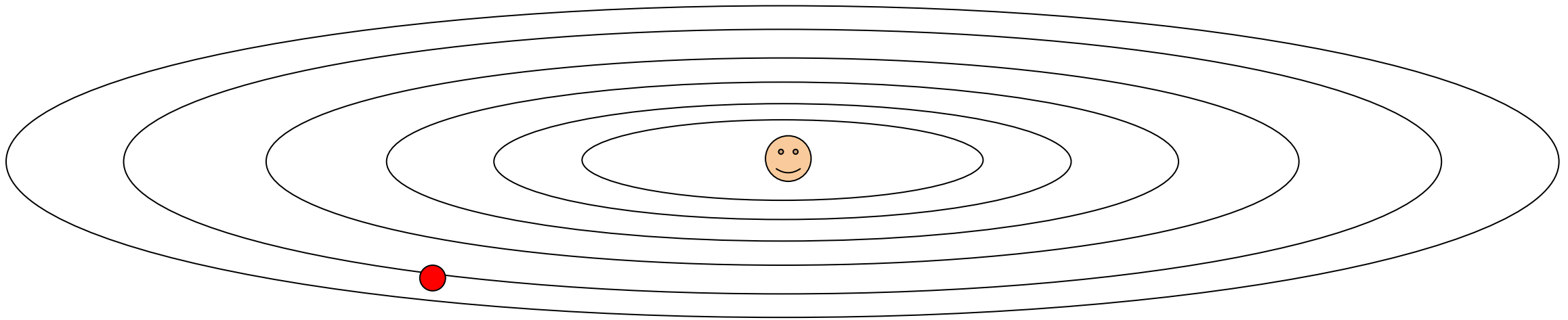
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for t in range(num_steps):
    dw = compute_gradient(w)
    grad_squared += dw * dw
    w -= learning_rate * dw / (grad_squared.sqrt() + 1e-7)
```



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

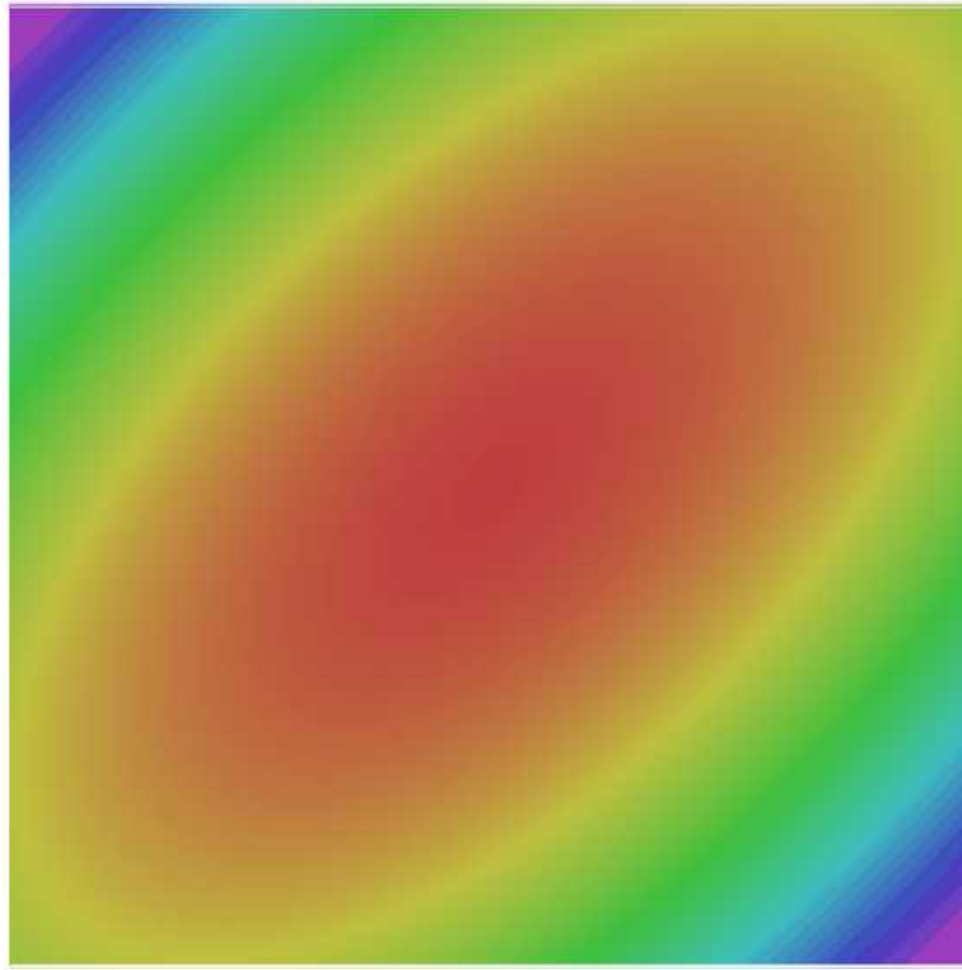
AdaGrad

↓

```
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

RMSProp



- SGD
- SGD+Momentum
- RMSProp

Adam (almost): RMSProp + 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 (almost): RMSProp + Momentum

```
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

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

SGD+Momentum

Adam (almost): RMSProp + Momentum

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    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
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RMSProp

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    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 $\beta_2 = 0.999$)

Adam (almost): RMSProp + 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
    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

Adam (almost): RMSProp + 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
    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 $\beta_1 = 0.9$,
 $\beta_2 = 0.999$, and $\text{learning_rate} = 1e-3, 5e-4, 1e-4$
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 mini-batch 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.

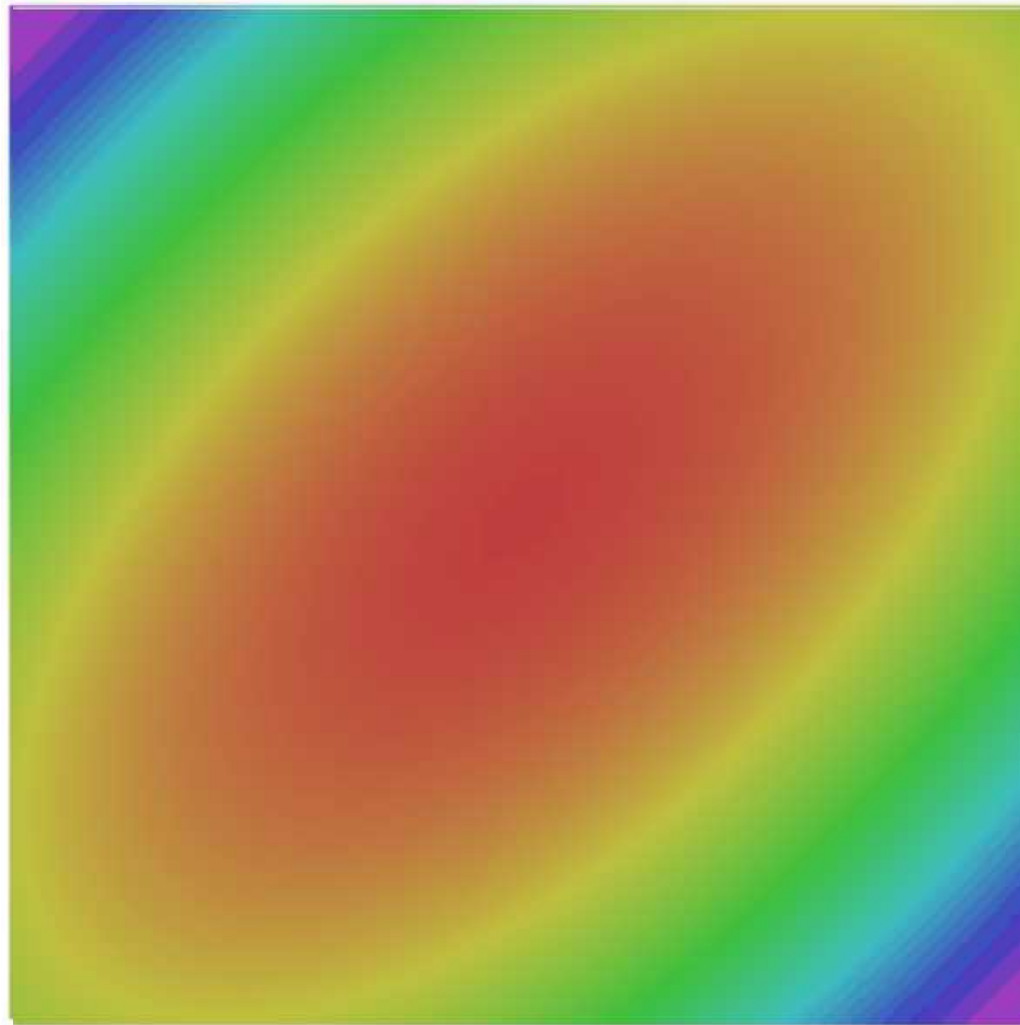
Zhu, Kaplan, Johnson, and Fei-Fei, ECCV 2018

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 $\beta_1 = 0.9$,
 $\beta_2 = 0.999$, and $\text{learning_rate} = 1e-3, 5e-4, 1e-4$
is a great starting point for many models!

Adam



- SGD
- SGD+Momentum
- RMSProp
- Adam

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	✓	✓	✓	✓

L2 Regularization vs Weight Decay

Optimization Algorithm

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

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

$$s_t = \text{optimizer}(g_t)$$

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

L2 Regularization vs Weight Decay

Optimization Algorithm

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

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L2 Regularization

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

$$g_t = \nabla L(w_t) = \nabla L_{data}(w_t) + 2\lambda w_t$$

$$s_t = \text{optimizer}(g_t)$$

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

L2 Regularization vs Weight Decay

Optimization Algorithm

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

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Weight Decay

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

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

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

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

L2 Regularization vs Weight Decay

Optimization Algorithm

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

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

L2 Regularization

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

$$g_t = \nabla L(w_t) = \nabla L_{data}(w_t) + 2\lambda w_t$$

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

Weight Decay

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

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

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L2 Regularization vs Weight Decay

Optimization Algorithm

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

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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

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

$$g_t = \nabla L(w_t) = \nabla L_{data}(w_t) + 2\lambda w_t$$

$$s_t = \text{optimizer}(g_t)$$

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

Weight Decay

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

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

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

AdamW: Decoupled Weight Decay

Algorithm 2 Adam with L_2 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  $\theta_{t=0} \in \mathbb{R}^n$ , first moment vector  $\mathbf{m}_{t=0} \leftarrow \mathbf{0}$ , 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:    $\nabla f_t(\theta_{t-1}) \leftarrow \text{SelectBatch}(\theta_{t-1})$  ▷ select batch and return the corresponding gradient
6:    $\mathbf{g}_t \leftarrow \nabla f_t(\theta_{t-1}) + \lambda \theta_{t-1}$ 
7:    $\mathbf{m}_t \leftarrow \beta_1 \mathbf{m}_{t-1} + (1 - \beta_1) \mathbf{g}_t$  ▷ 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:    $\hat{\mathbf{m}}_t \leftarrow \mathbf{m}_t / (1 - \beta_1^t)$  ▷  $\beta_1$  is taken to the power of  $t$ 
10:   $\hat{\mathbf{v}}_t \leftarrow \mathbf{v}_t / (1 - \beta_2^t)$  ▷  $\beta_2$  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
12:   $\theta_t \leftarrow \theta_{t-1} - \eta_t \left( \alpha \hat{\mathbf{m}}_t / (\sqrt{\hat{\mathbf{v}}_t} + \epsilon) + \lambda \theta_{t-1} \right)$ 
13: until stopping criterion is met
14: return optimized parameters  $\theta_t$ 
```

AdamW: Decoupled Weight Decay

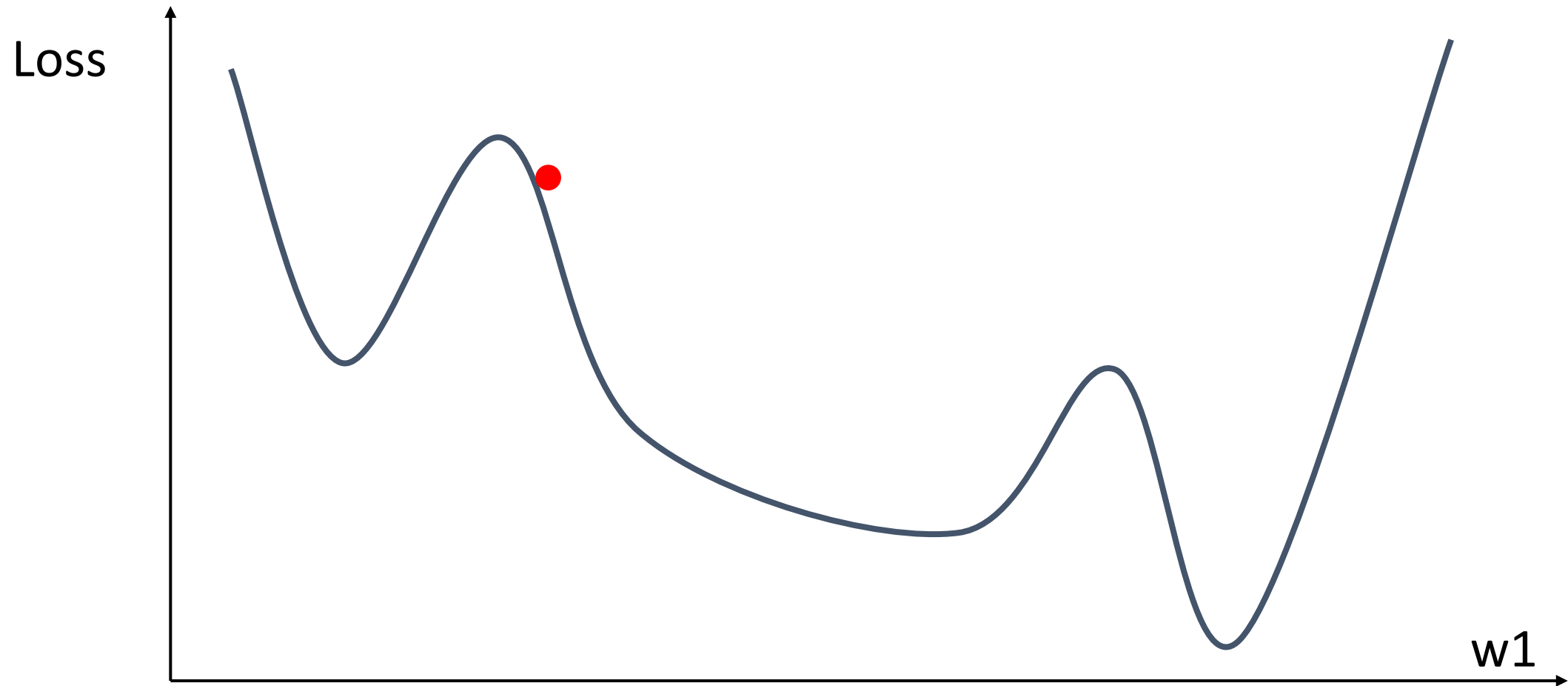
Algorithm 2 Adam with L_2 regularization and Adam with decoupled weight decay (AdamW)

- 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 $\theta_{t=0} \in \mathbb{R}^n$, first moment vector $m_{t=0} \leftarrow \mathbf{0}$, second moment vector $v_{t=0} \leftarrow \mathbf{0}$, schedule multiplier $\eta_{t=0} \in \mathbb{R}$

AdamW should probably be your
“default” optimizer for new problems

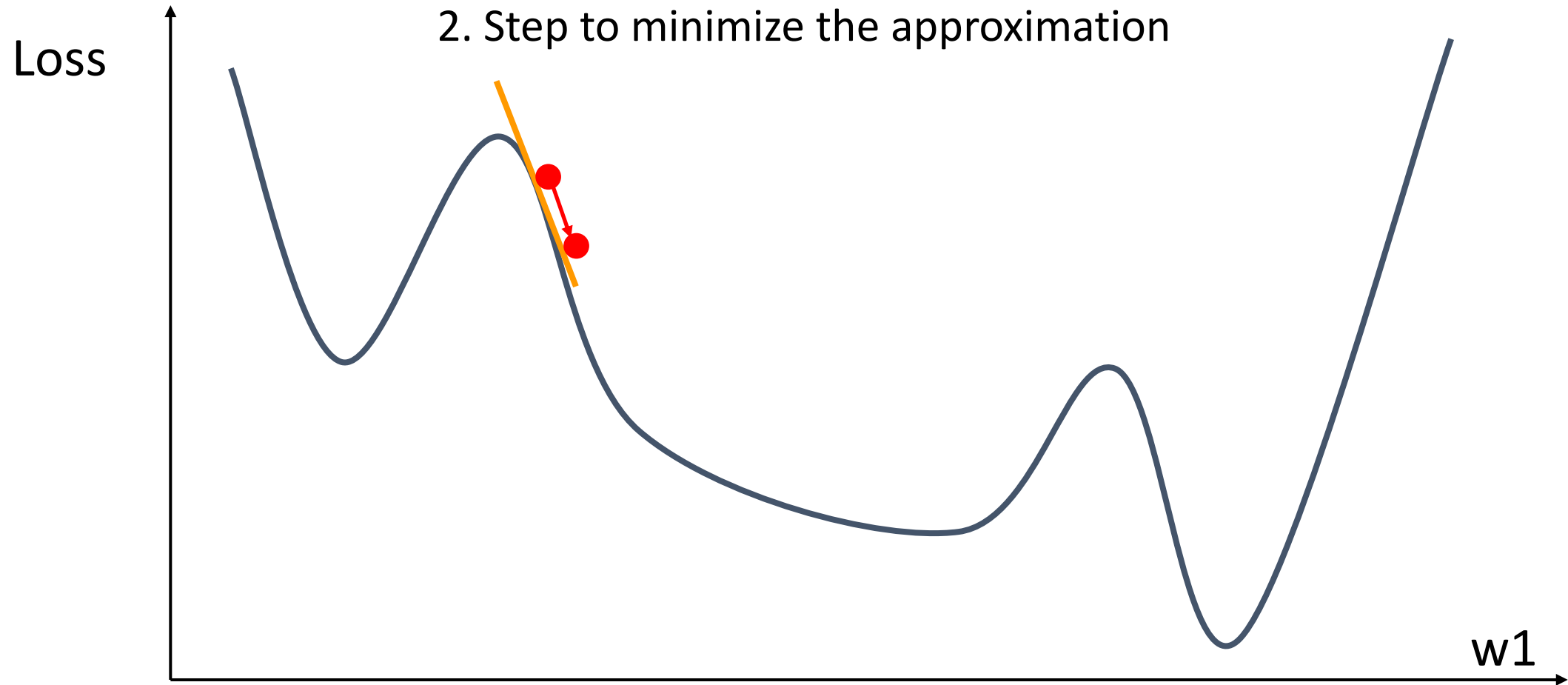
- 11: $\eta_t \leftarrow \text{SetScheduleMultiplier}(t)$ \triangleright can be fixed, decay, or also be used for warm restarts
- 12: $\theta_t \leftarrow \theta_{t-1} - \eta_t \left(\alpha \hat{m}_t / (\sqrt{\hat{v}_t} + \epsilon) + \lambda \theta_{t-1} \right)$
- 13: **until** *stopping criterion is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

So far: First-Order Optimization



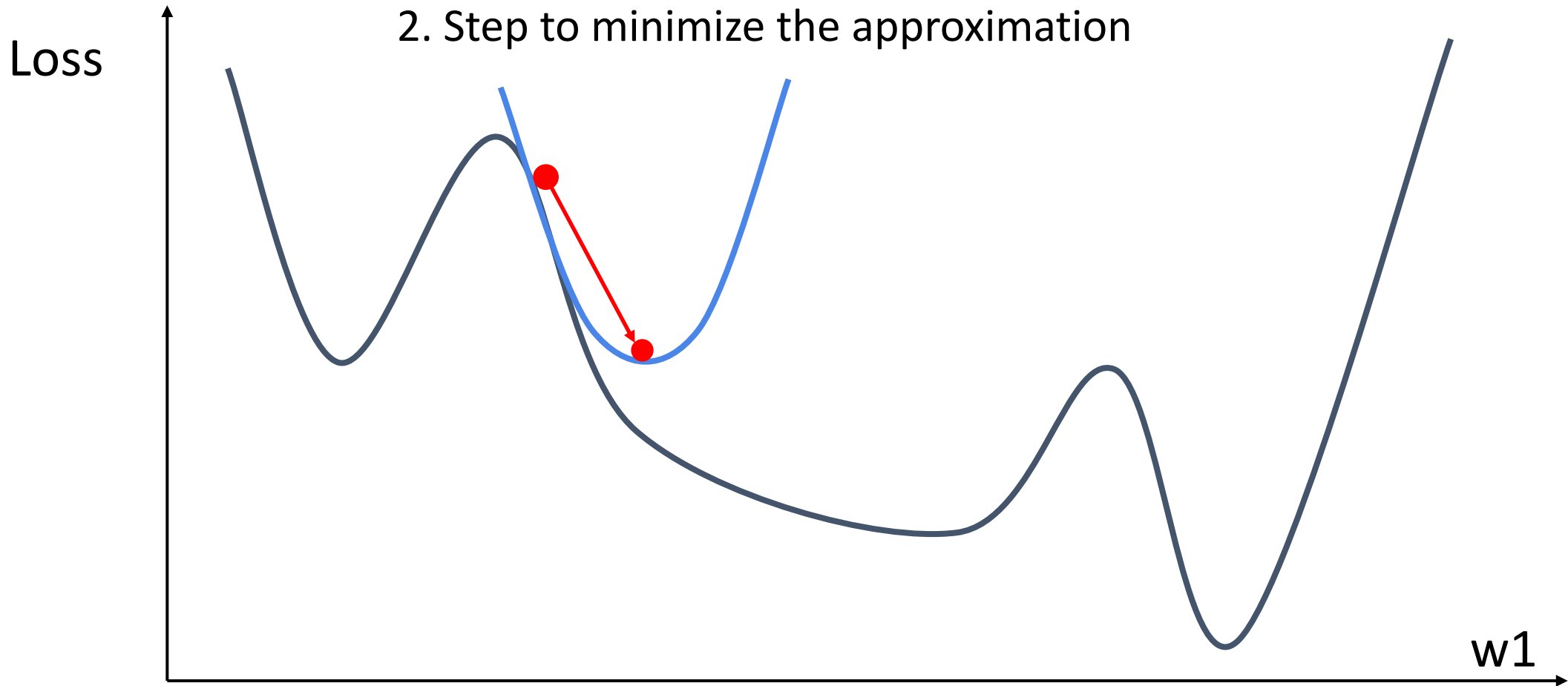
So far: First-Order Optimization

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



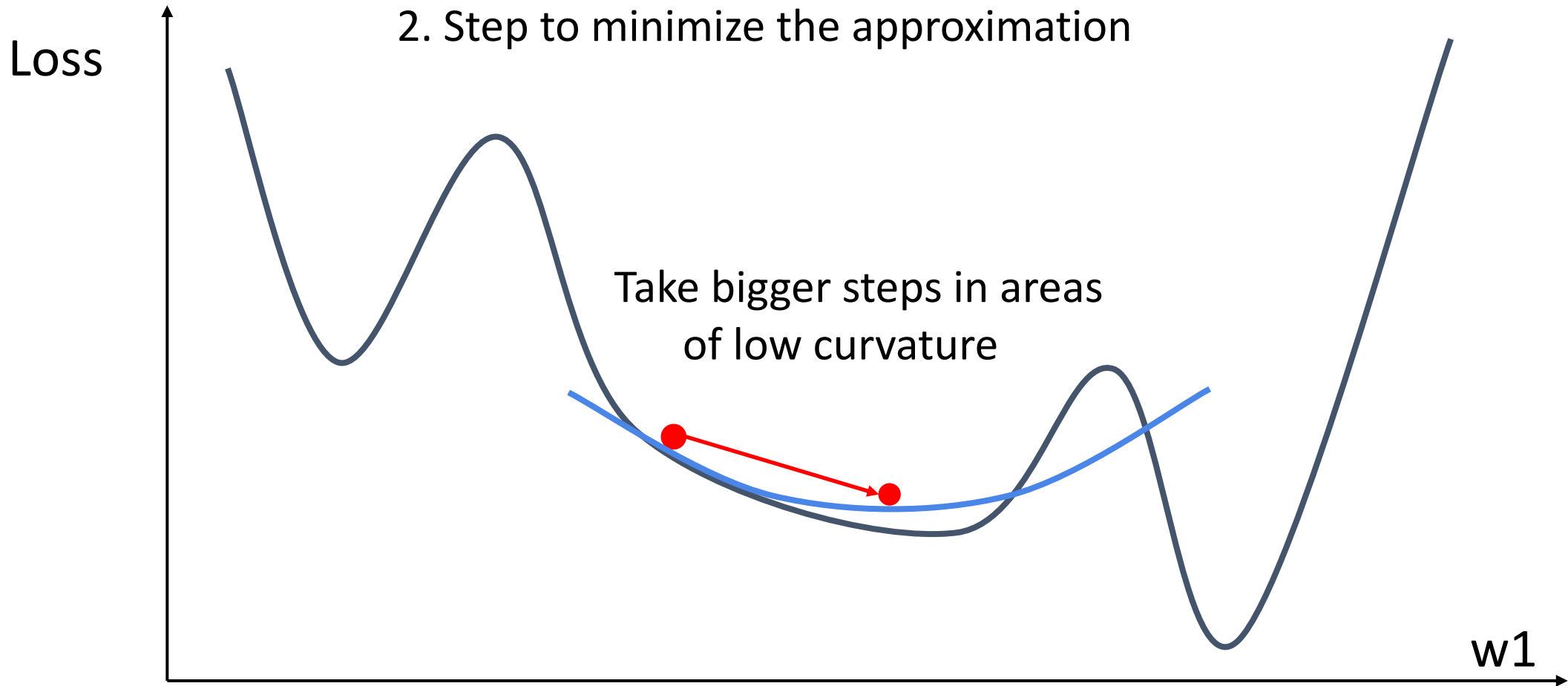
Second-Order Optimization

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



Second-Order Optimization

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



Second-Order Optimization

Second-Order Taylor Expansion:

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

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 Optimization

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Q: Why is this impractical?

Second-Order Optimization

Second-Order Taylor Expansion:

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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

Second-Order Optimization

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

- 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

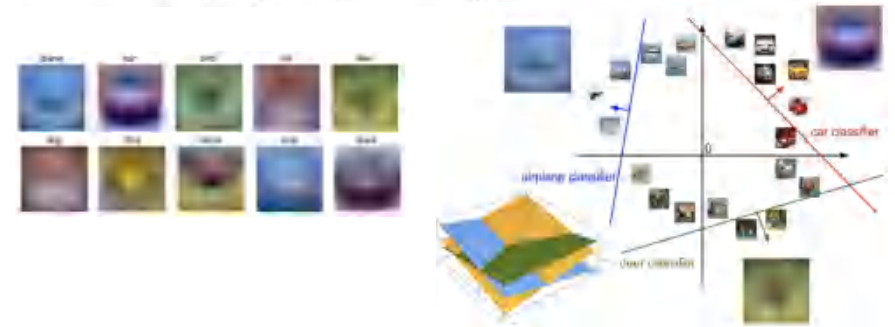
In practice:

- **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
4. Use **Stochastic Gradient Descent** to minimize our loss functions and train the model

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

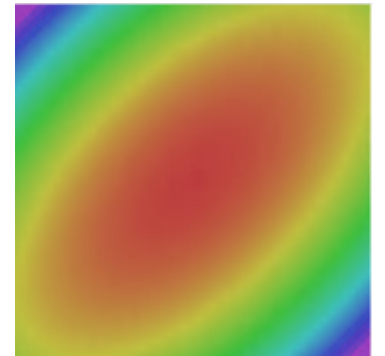


$$L_i = -\log\left(\frac{e^{s_{y_i}}}{\sum_j e^{s_j}}\right) \quad \text{Softmax} \quad \text{SVM}$$

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

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

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



Next time:
Neural Networks