

# 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



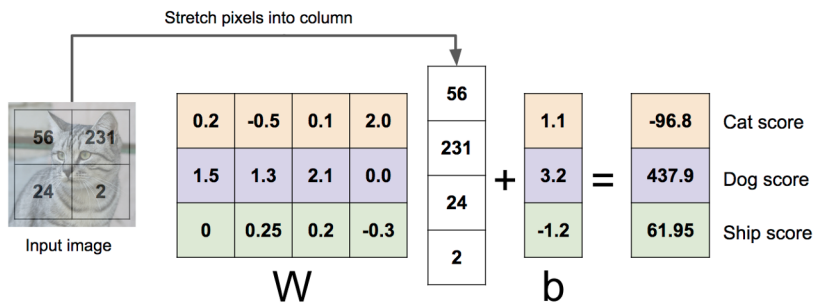
# Assignment 2

- 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

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



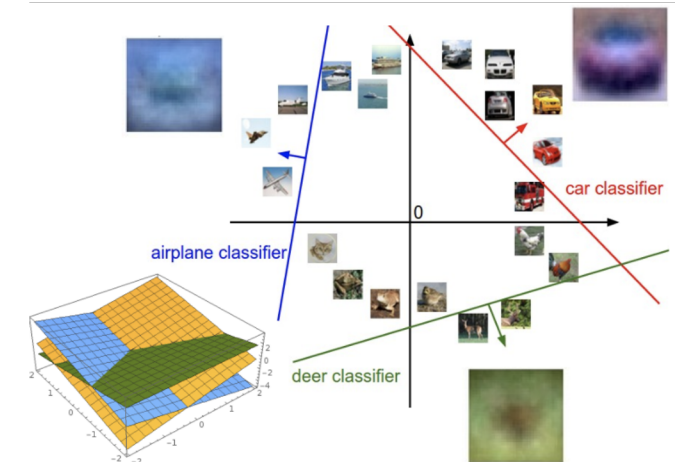
## Visual Viewpoint

One template  
per class



## Geometric Viewpoint

Hyperplanes  
cutting up space



# Last Time: Loss Functions quantify preferences

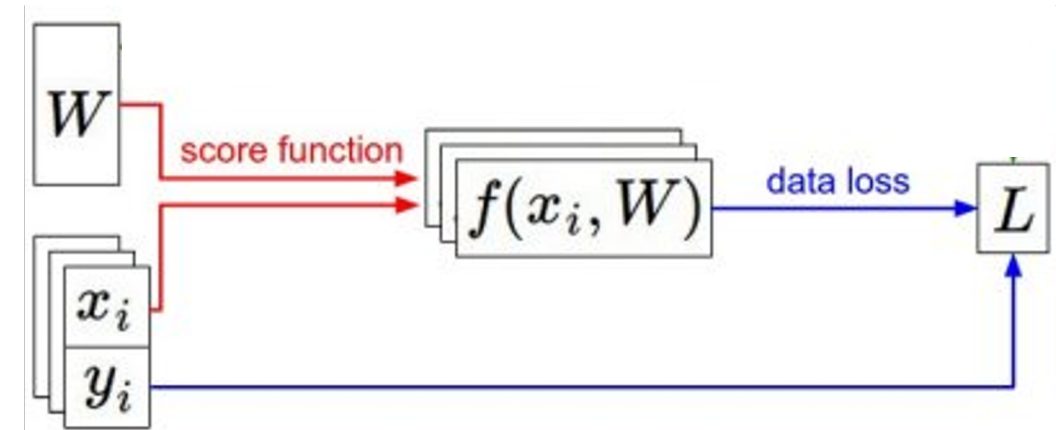
- We have some dataset of  $(x, y)$
- We have a **score function**:
- We have a **loss function**:

**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**:

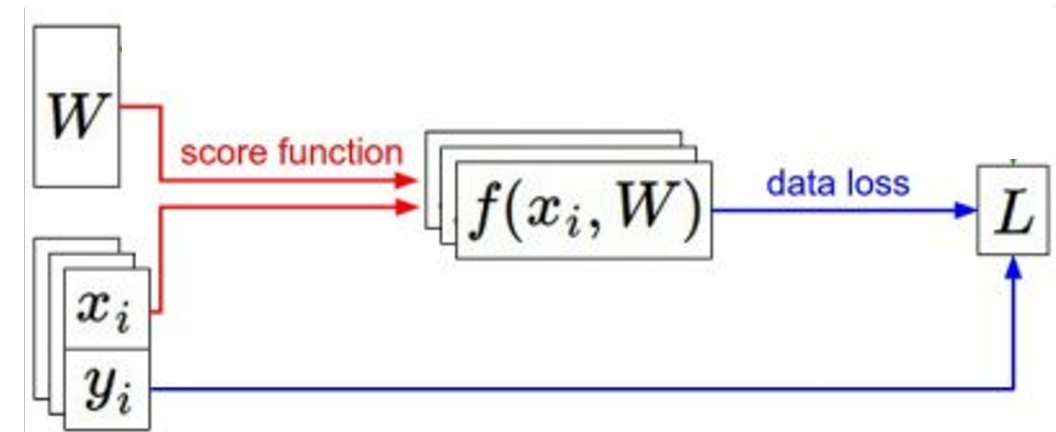
**Problem:** Loss functions encourage good performance on training data but we really care about test data

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

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

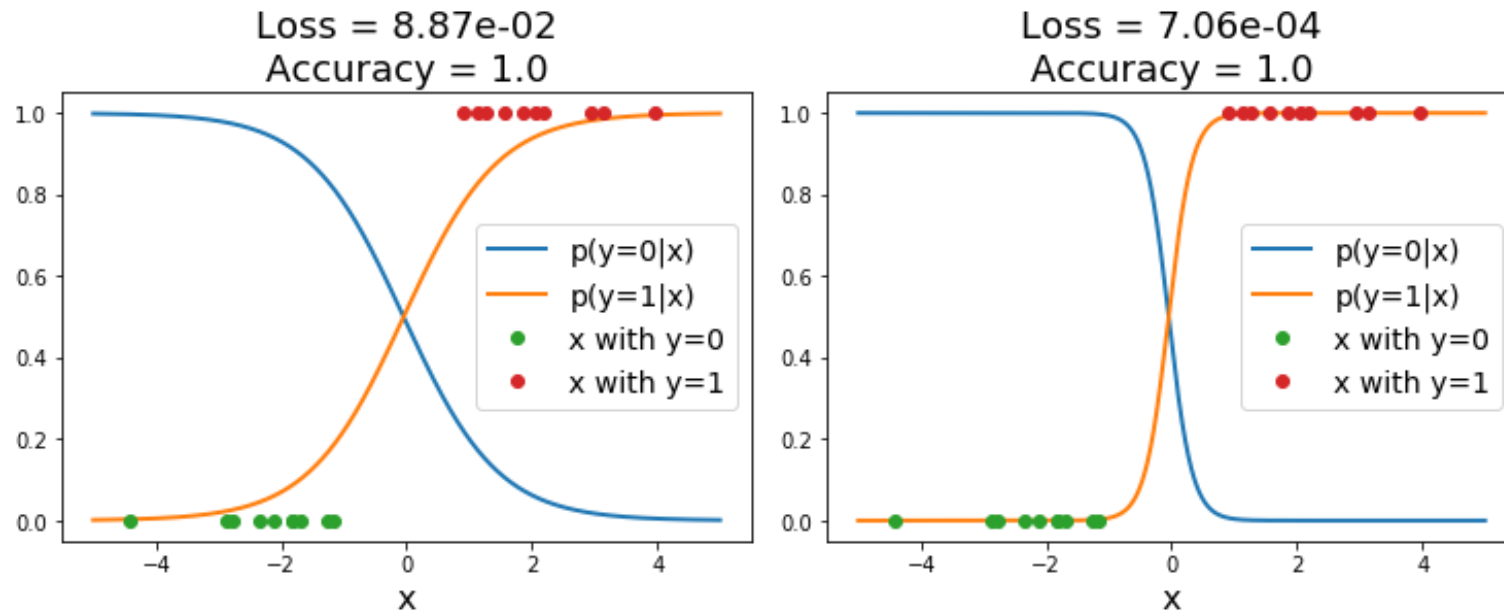
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Both models have perfect accuracy on train data!

# Overfitting

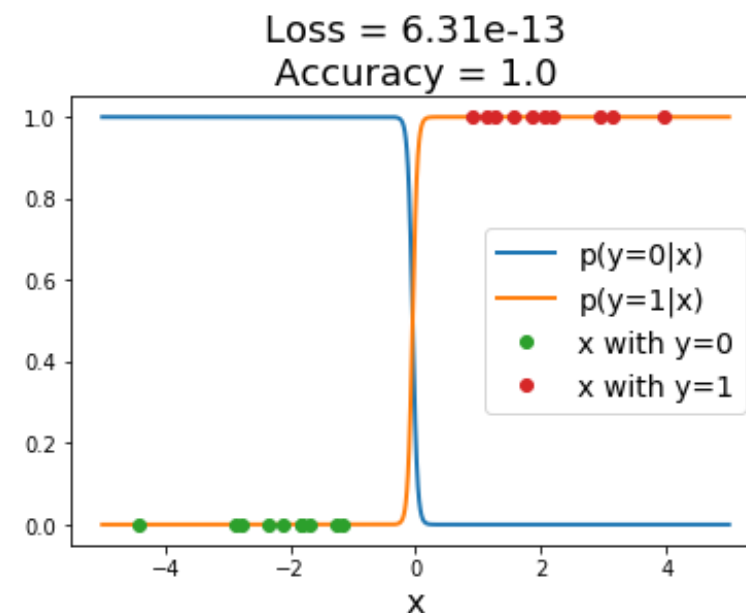
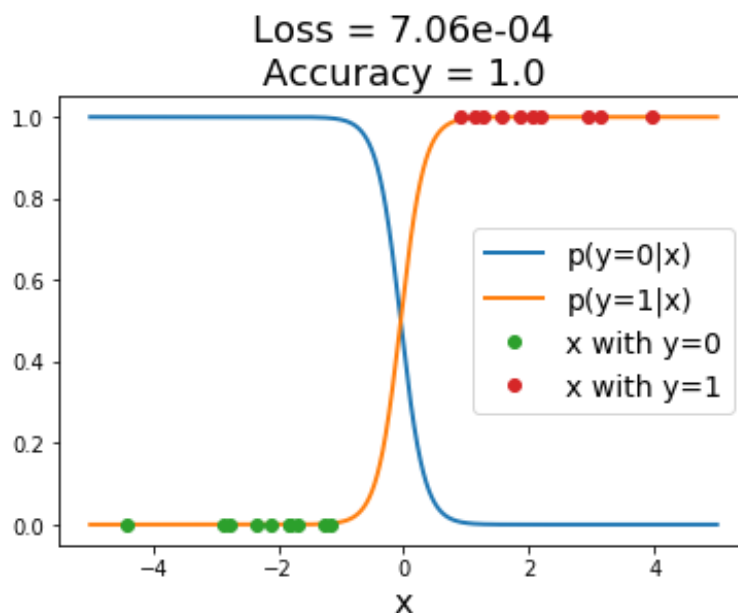
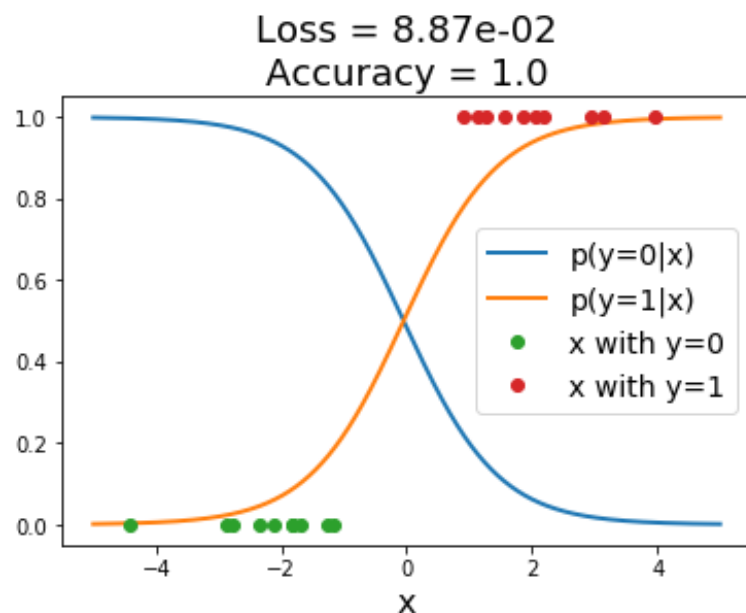
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$$L = -\log(p_y)$$



Both models have perfect accuracy on train data!

Low loss, but unnatural “cliff” between training points



# Overfitting

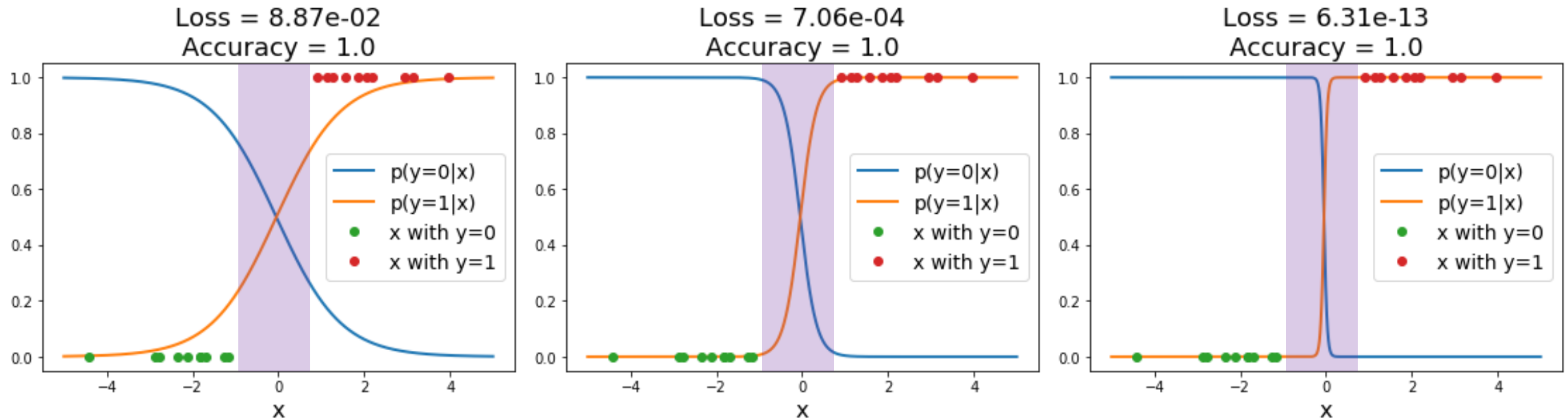
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$$p_i = \frac{\exp(s_i)}{\exp(s_1) + \exp(s_2)}$$

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



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

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

$\lambda$  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

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

$\lambda$  is a hyperparameter giving regularization strength

## 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

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

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

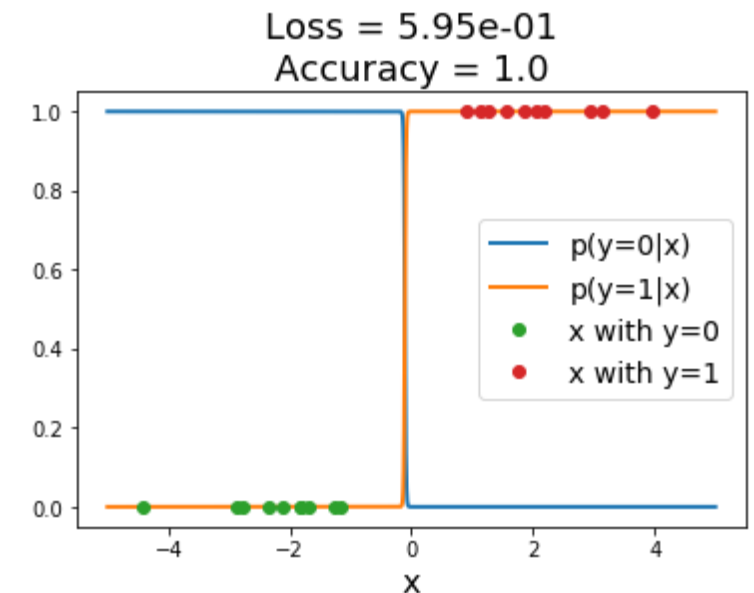
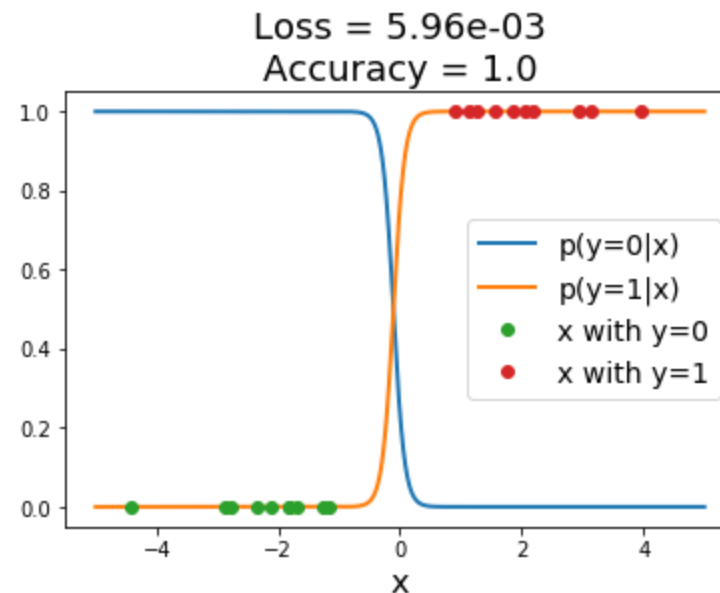
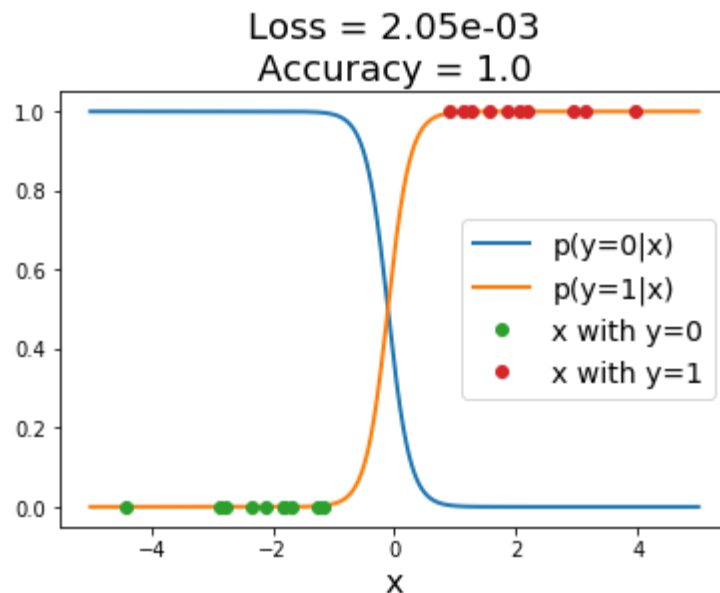
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$$L = -\log(p_y) + \lambda \sum_i w_i^2$$

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



# Regularization: Expressing Preferences

L2 Regularization

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

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

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

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

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

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



## Idea #2: Follow the slope

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 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**

**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,  
0.55,  
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  $W$

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

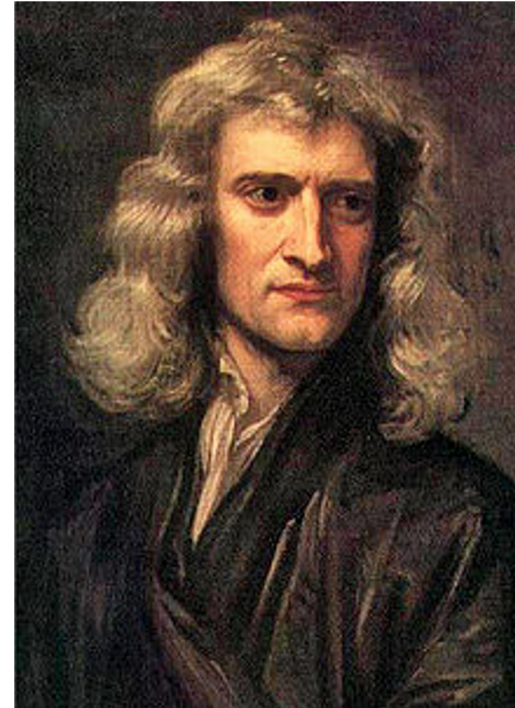
Want  $\nabla_w L$



# 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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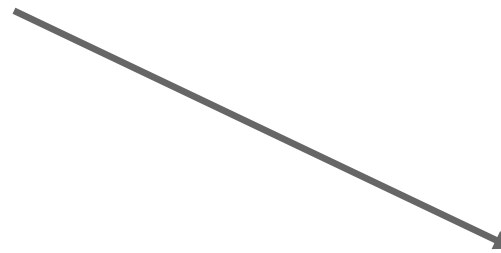
Use calculus to compute an **analytic gradient**

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

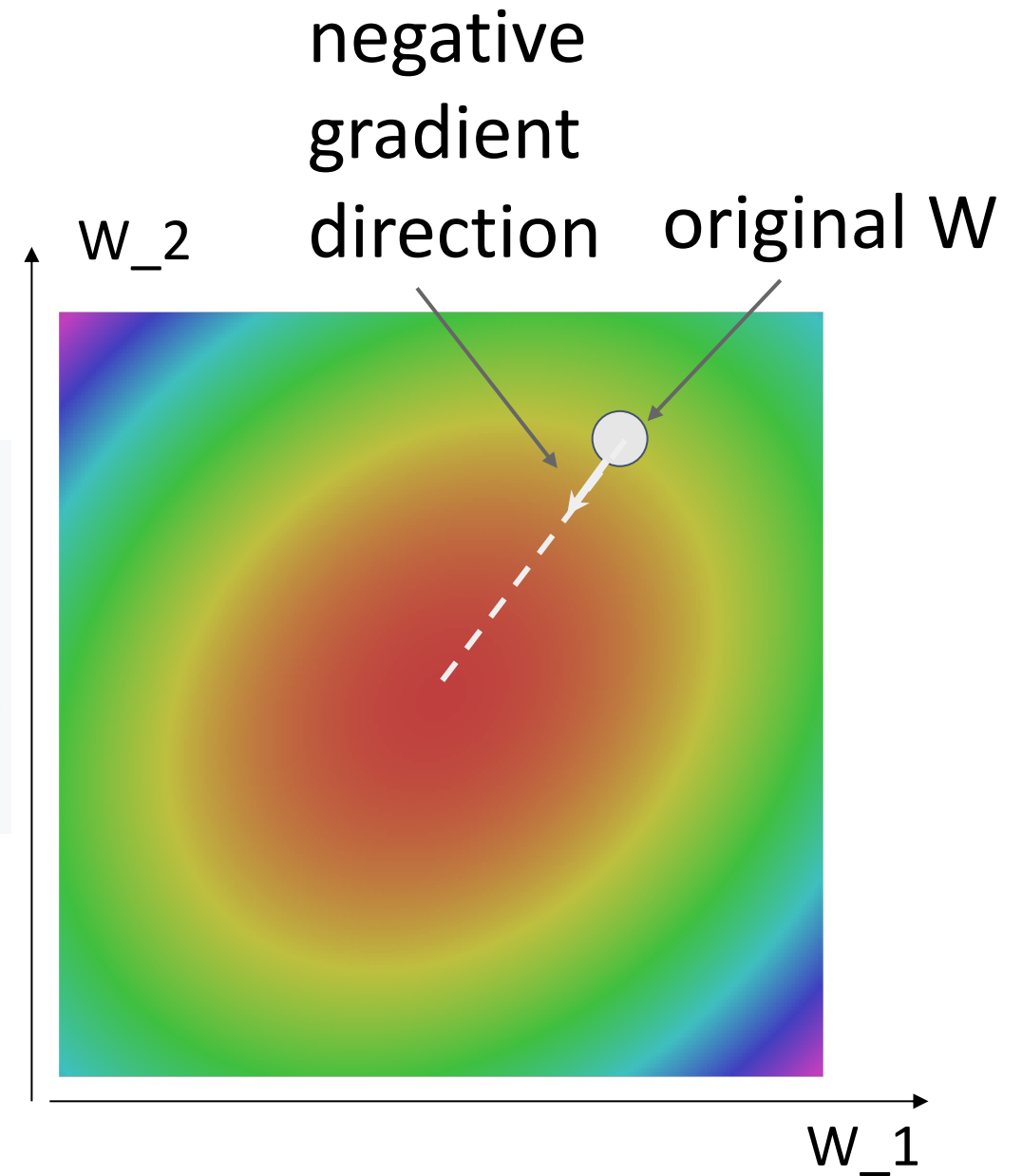
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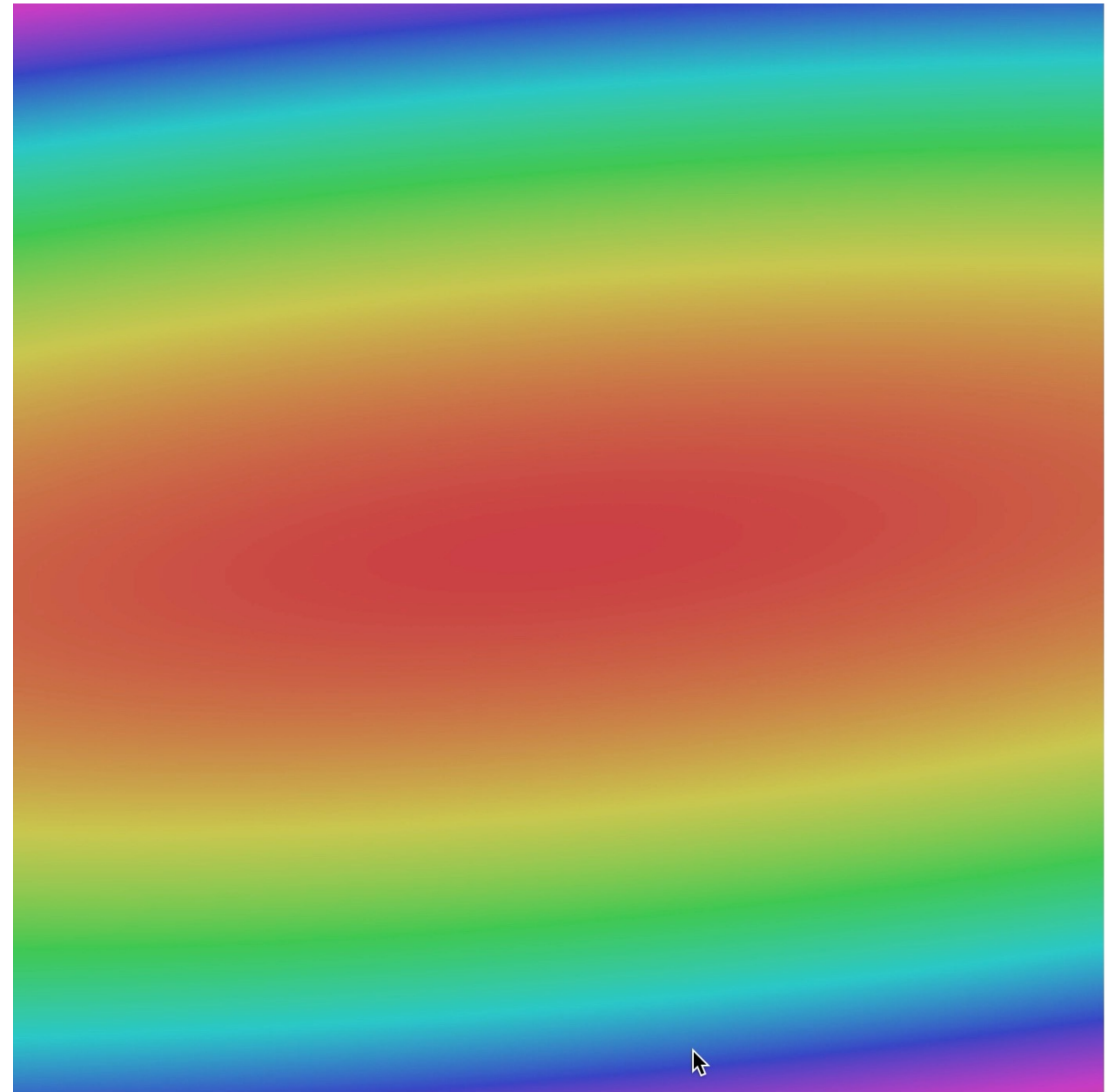
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```

## Hyperparameters:

- 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
```

## 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)

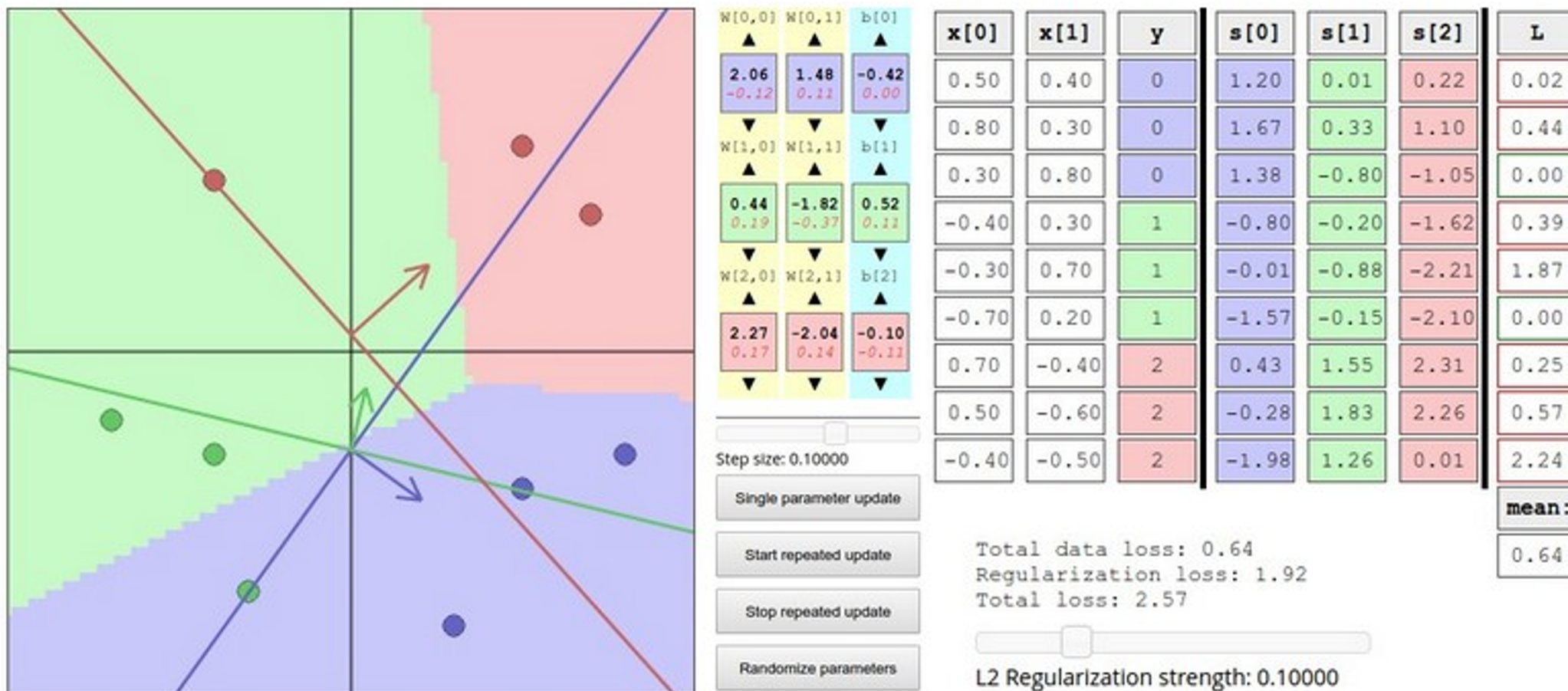
$$\begin{aligned} L(W) &= \mathbb{E}_{(x,y) \sim p_{data}} [L(x, y, W)] + \lambda R(W) \\ &\approx \frac{1}{N} \sum_{i=1}^N L(x_i, y_i, W) + \lambda R(W) \end{aligned}$$

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

Approximate expectation via sampling

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

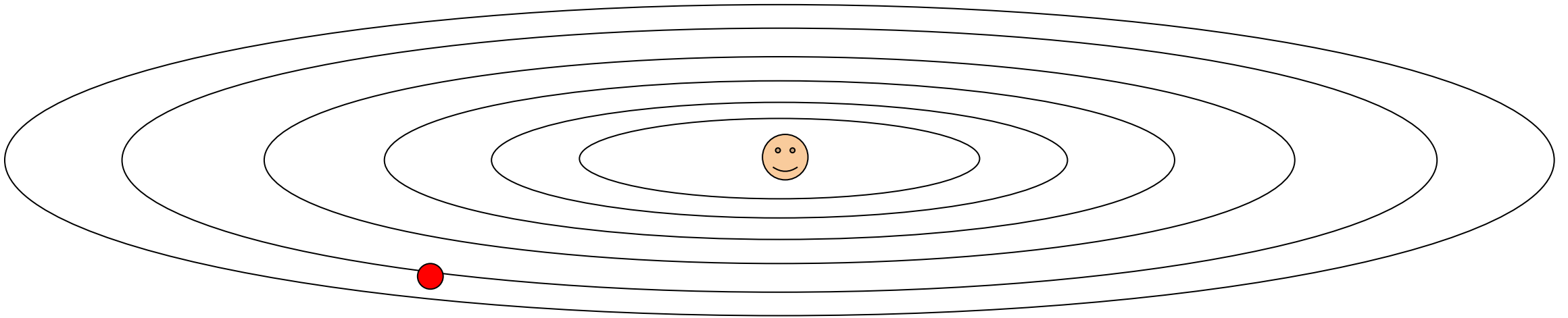
# Interactive Web Demo



<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?



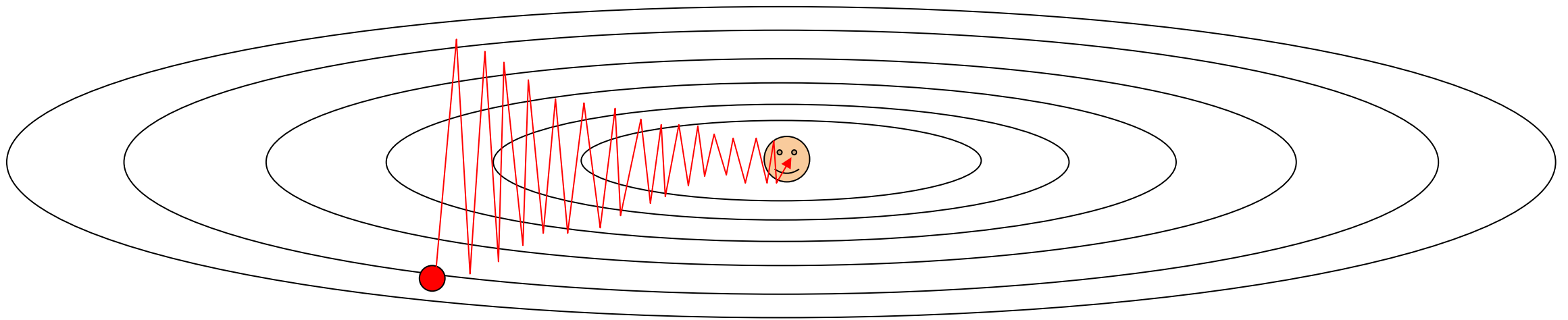
Loss function has high **condition number**: ratio of largest to smallest singular value of the Hessian matrix is large

# Problems with SGD

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

What does gradient descent do?

Very slow progress along shallow dimension, jitter along steep direction

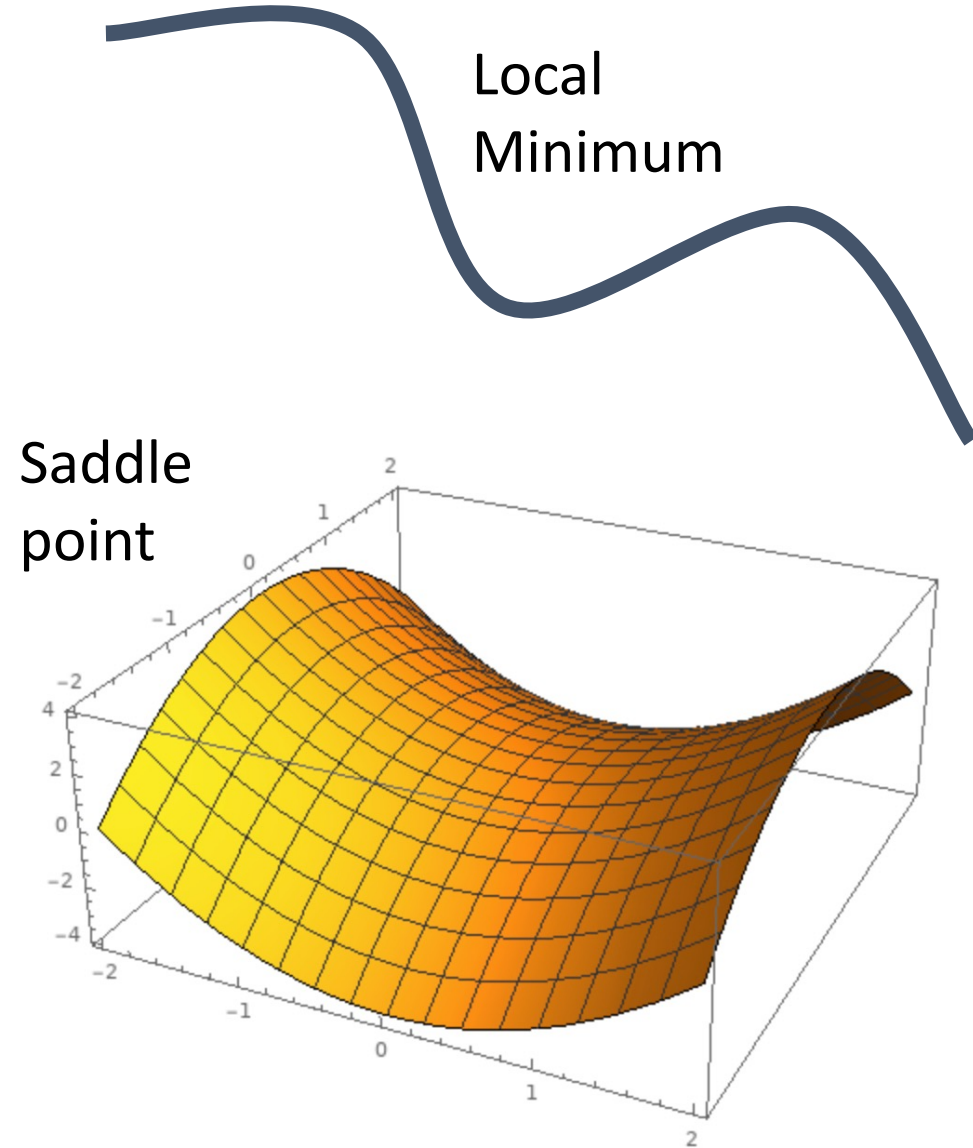


Loss function has high **condition number**: ratio of largest to smallest singular value of the Hessian matrix is large



# Problems with SGD

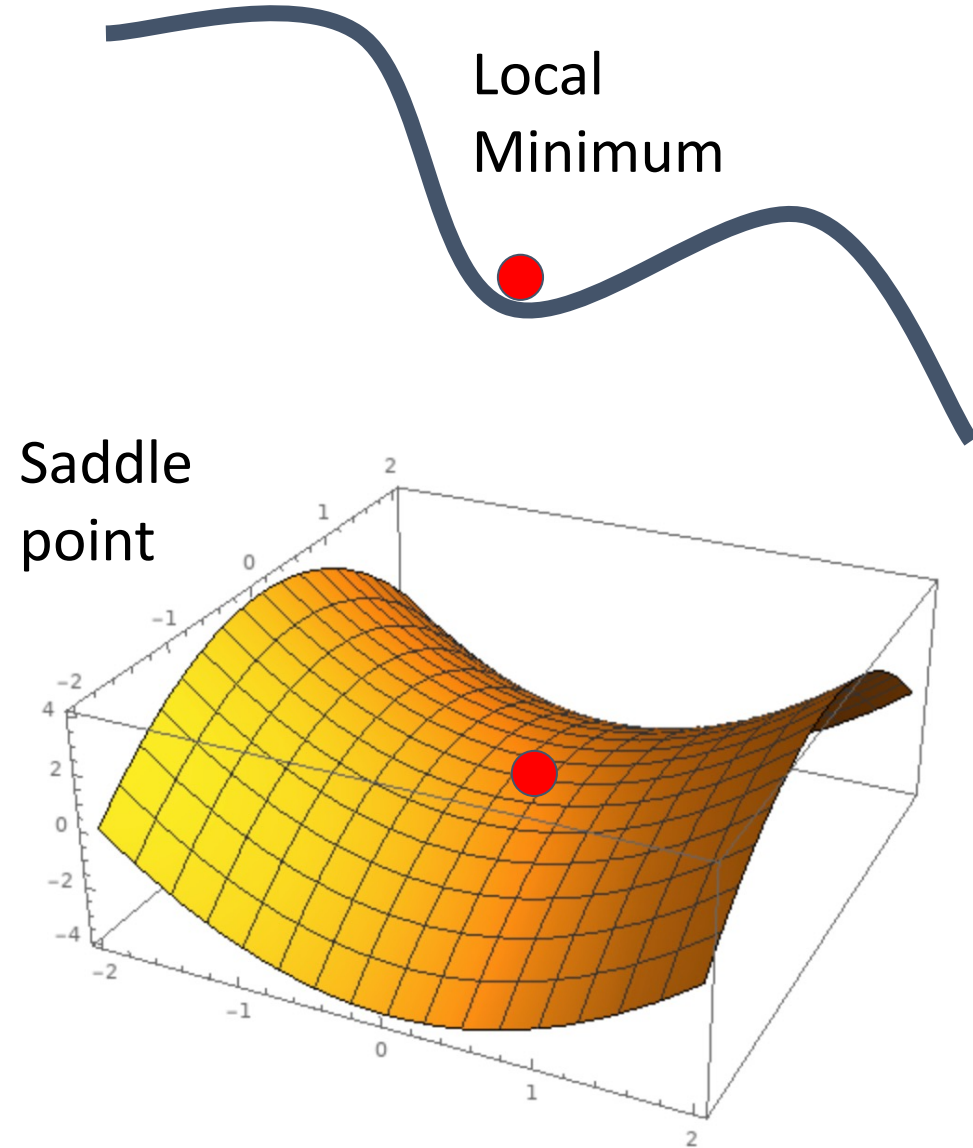
What if the loss function has a **local minimum** or **saddle point**?



# 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

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

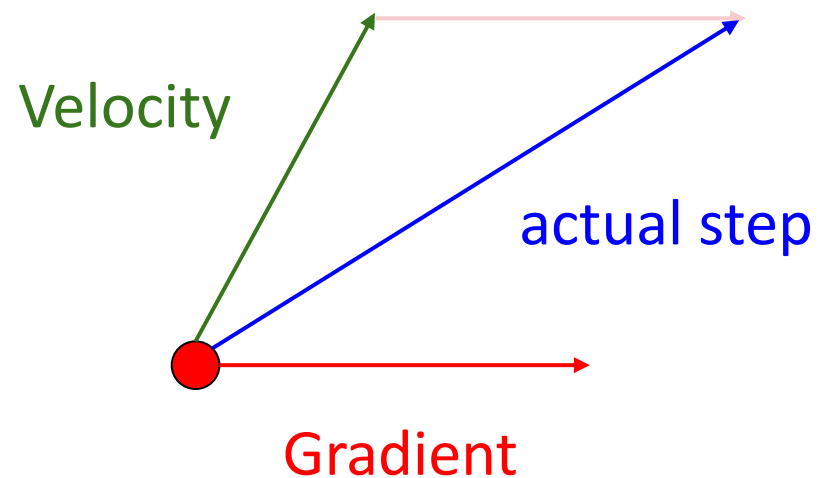
$$x_{t+1} = x_t - \alpha v_{t+1}$$

```
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
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    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  $x$



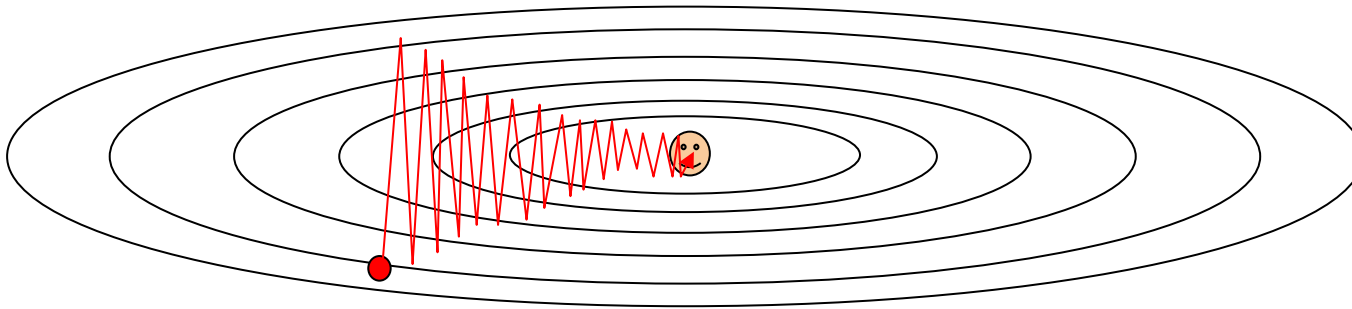
# SGD + Momentum

Local Minima

Saddle points



Poor Conditioning



Gradient Noise

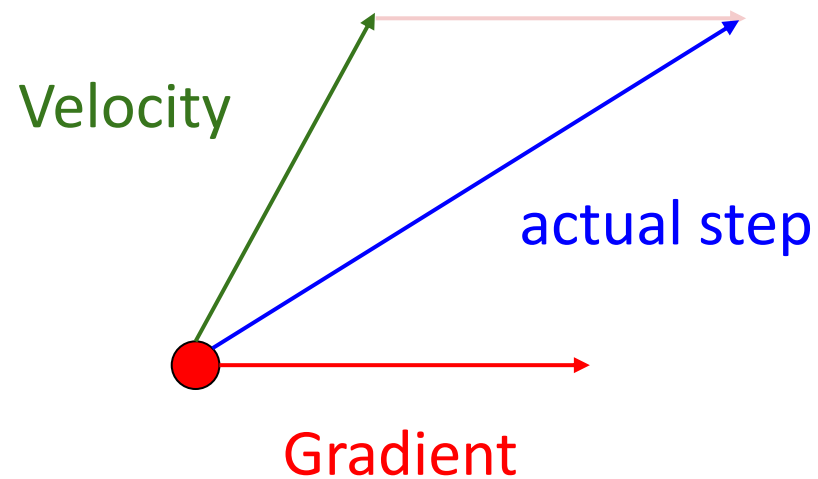


— SGD — SGD+Momentum

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

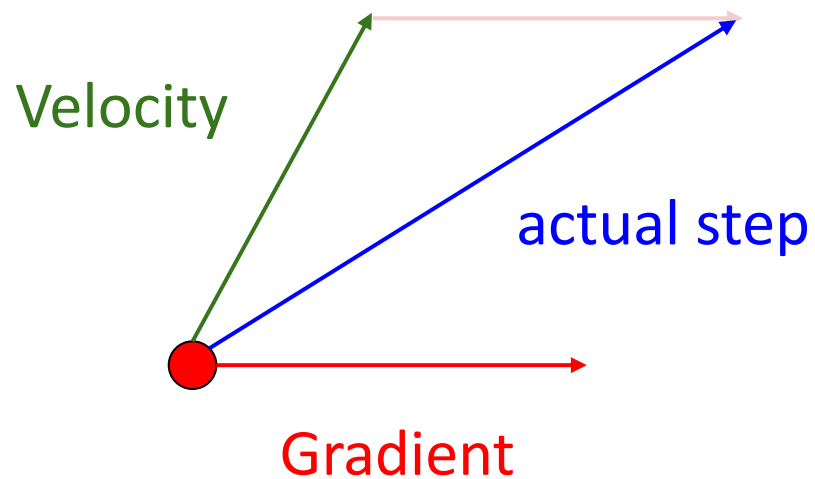
Nesterov, "A method of solving a convex programming problem with convergence rate  $O(1/k^2)$ ", 1983

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

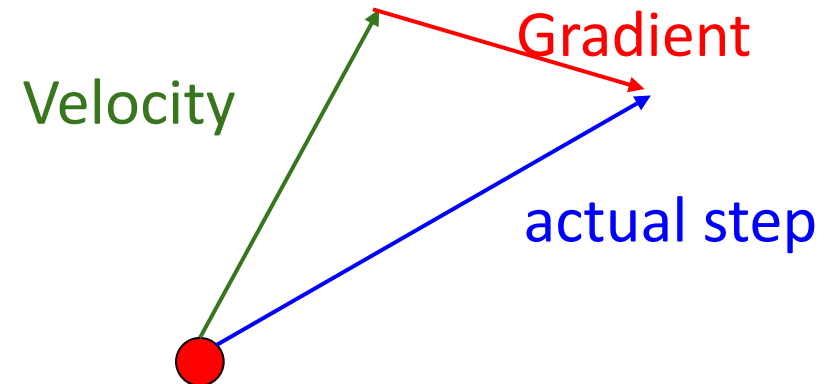
# Nesterov Momentum

Momentum update:



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

Nesterov Momentum



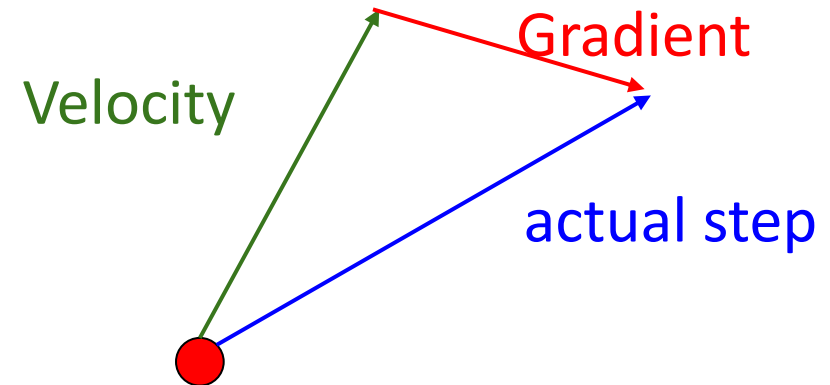
“Look ahead” to the point where updating using velocity would take us; compute gradient there and mix it with velocity to get actual update direction

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Nesterov, “Introductory lectures on convex optimization: a basic course”, 2004  
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# Nesterov Momentum

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

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



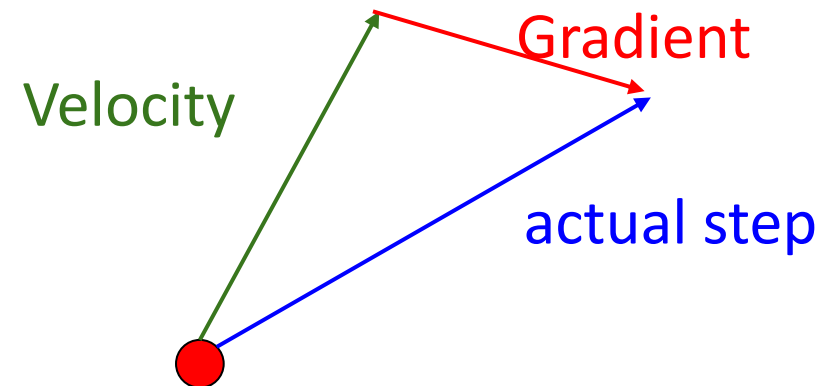
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# Nesterov Momentum

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

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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

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

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

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

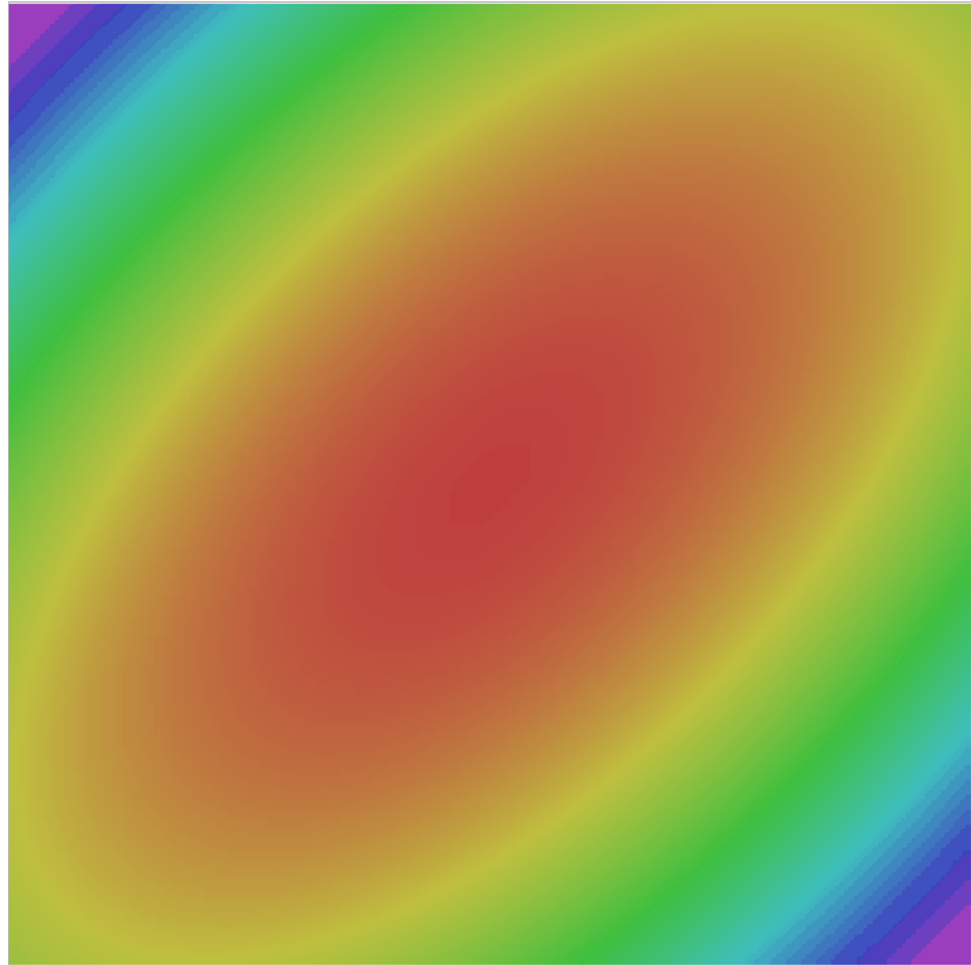
Change of variables  $\tilde{x}_t = x_t + \rho v_t$   
and rearrange:

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

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

```
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

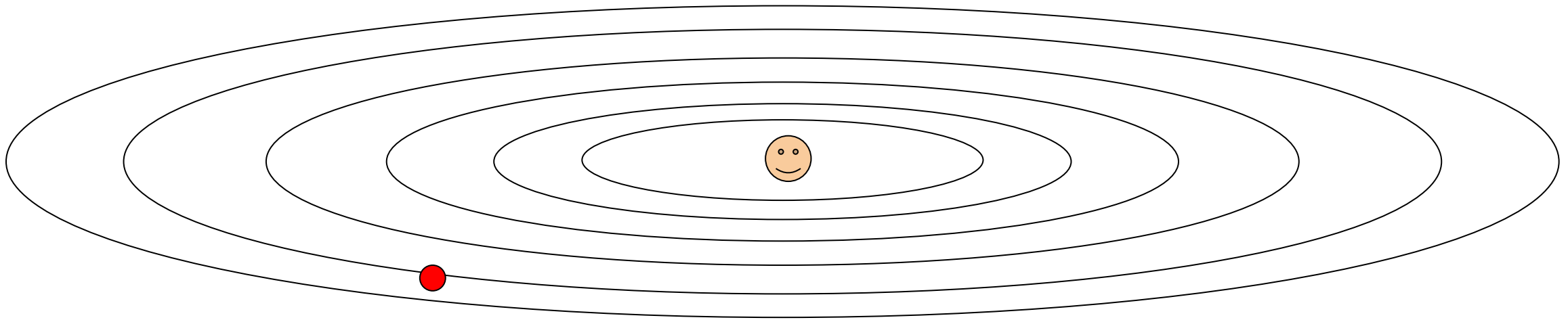
```
grad_squared = 0
for t in range(num_steps):
    dw = compute_gradient(w)
    grad_squared += dw * dw
    w -= learning_rate * dw / (grad_squared.sqrt() + 1e-7)
```

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”

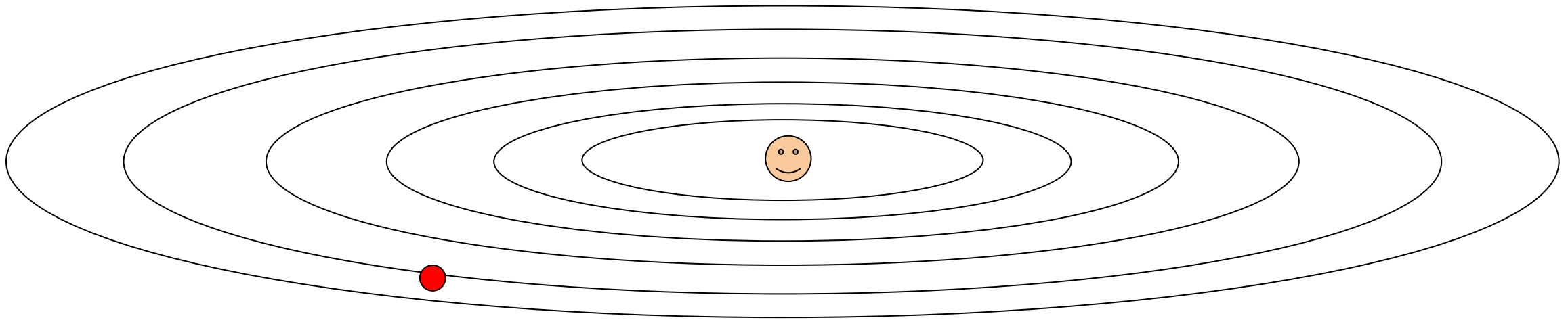
# AdaGrad

```
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

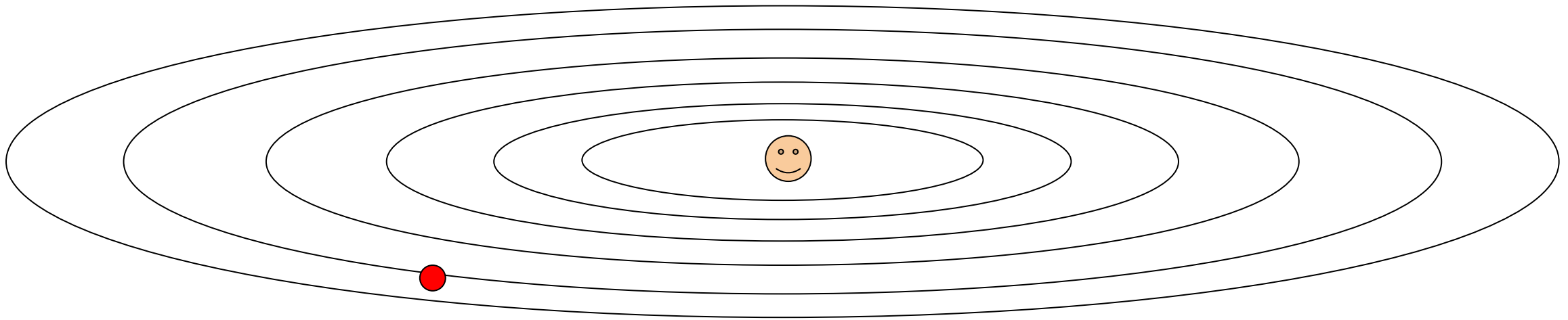
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for t in range(num_steps):
    dw = compute_gradient(w)
    grad_squared += dw * dw
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```



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)
```

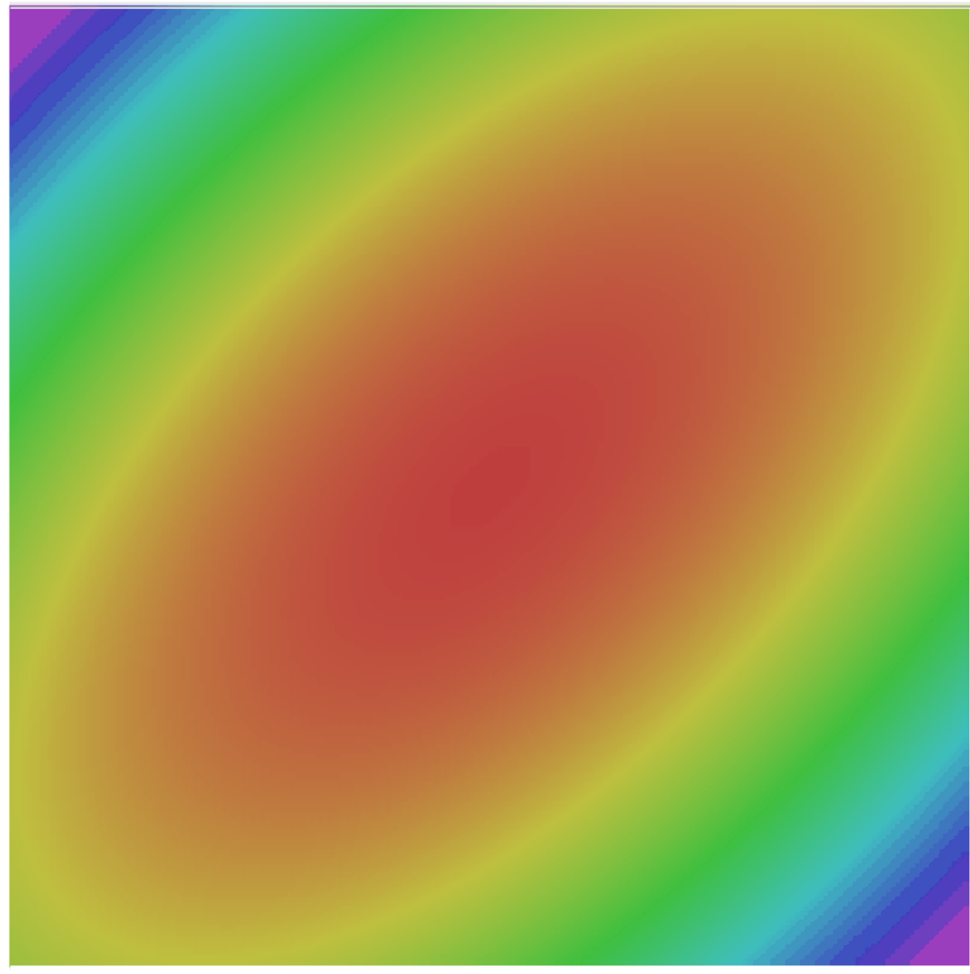


AdaGrad

```
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
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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

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

SGD+Momentum

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```

Adam

Momentum

AdaGrad / RMSProp

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

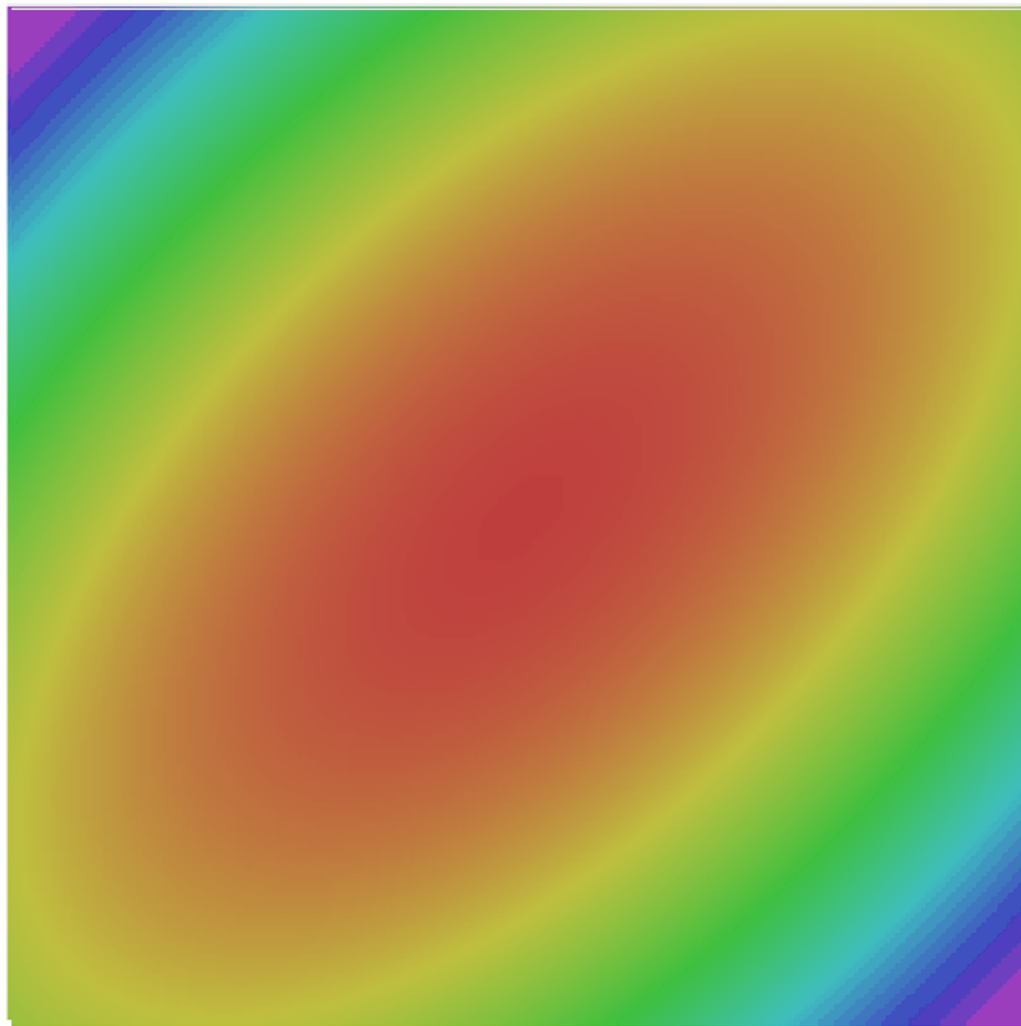
Zhu, Kaplan, Johnson, and Fei-Fei, ECCV 2018

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  $\text{beta1} = 0.9$ ,  
 $\text{beta2} = 0.999$ , and  $\text{learning\_rate} = 1\text{e-}3, 5\text{e-}4, 1\text{e-}4$   
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	<b>X</b>	<b>X</b>	<b>X</b>	<b>X</b>
SGD+Momentum	✓	<b>X</b>	<b>X</b>	<b>X</b>
Nesterov	✓	<b>X</b>	<b>X</b>	<b>X</b>
AdaGrad	<b>X</b>	✓	<b>X</b>	<b>X</b>
RMSProp	<b>X</b>	✓	✓	<b>X</b>
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

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

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

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## L2 Regularization

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

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

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

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## Weight Decay

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

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

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

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

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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

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

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

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## 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)

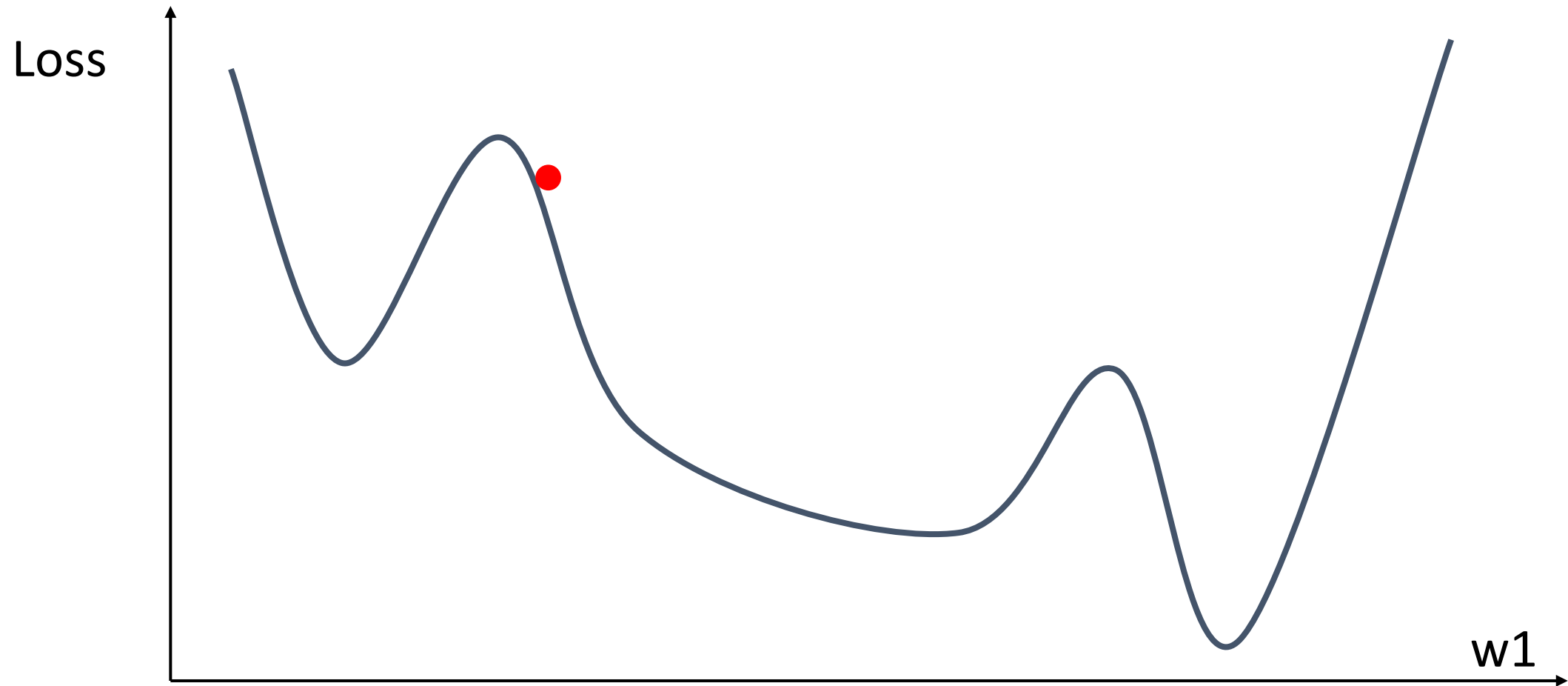
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- 1: **given**  $\alpha = 0.001, \beta_1 = 0.9, \beta_2 = 0.999, \epsilon = 10^{-8}, \lambda \in \mathbb{R}$
- 2: **initialize** time step  $t \leftarrow 0$ , parameter vector  $\theta_{t=0} \in \mathbb{R}^n$ , first moment vector  $m_{t=0} \leftarrow \mathbf{0}$ , 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

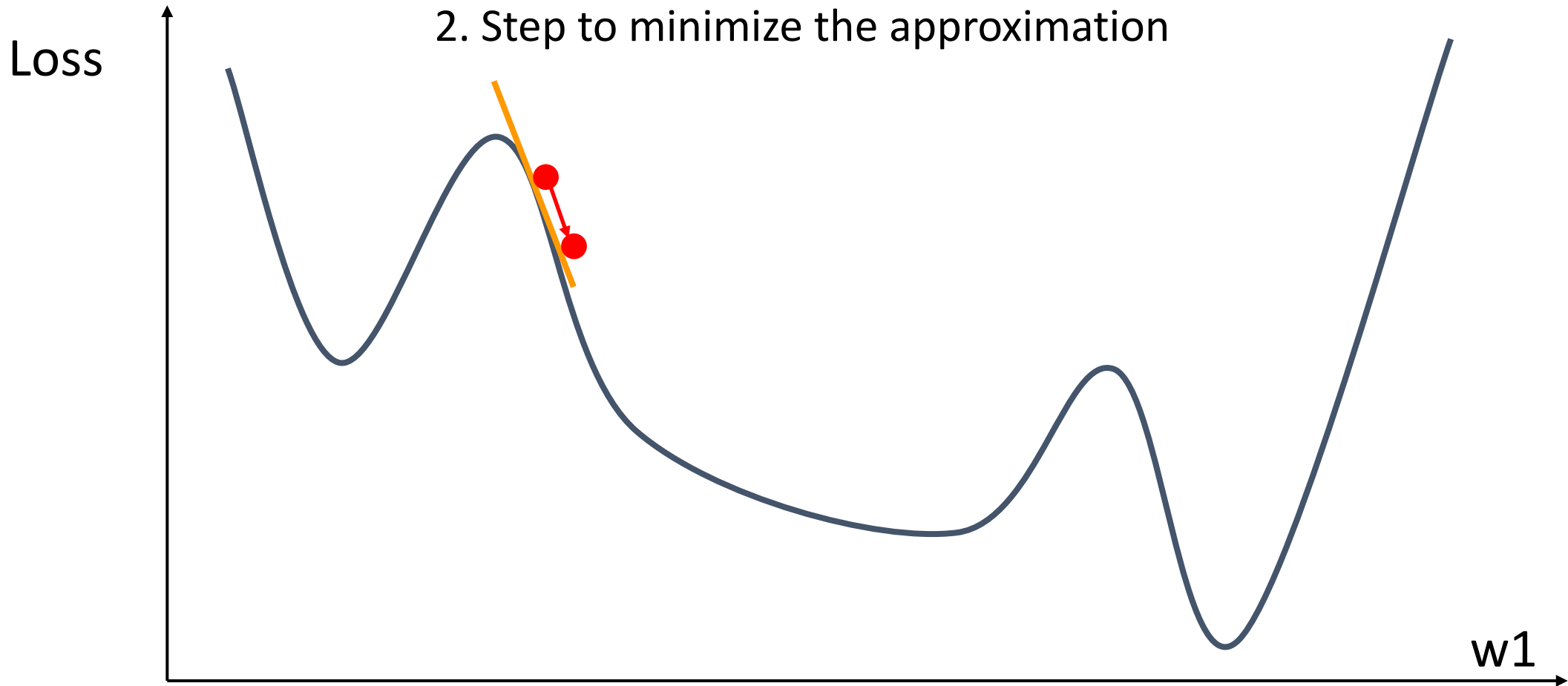
- 11:  $\eta_t \leftarrow \text{SetScheduleMultiplier}(t)$  ▷ can be fixed, decay, or also be used for warm restarts
  - 12:  $\theta_t \leftarrow \theta_{t-1} - \eta_t \left( \alpha \hat{m}_t / (\sqrt{\hat{v}_t} + \epsilon) + \lambda \theta_{t-1} \right)$
  - 13: **until** *stopping criterion is met*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

# So far: First-Order Optimization



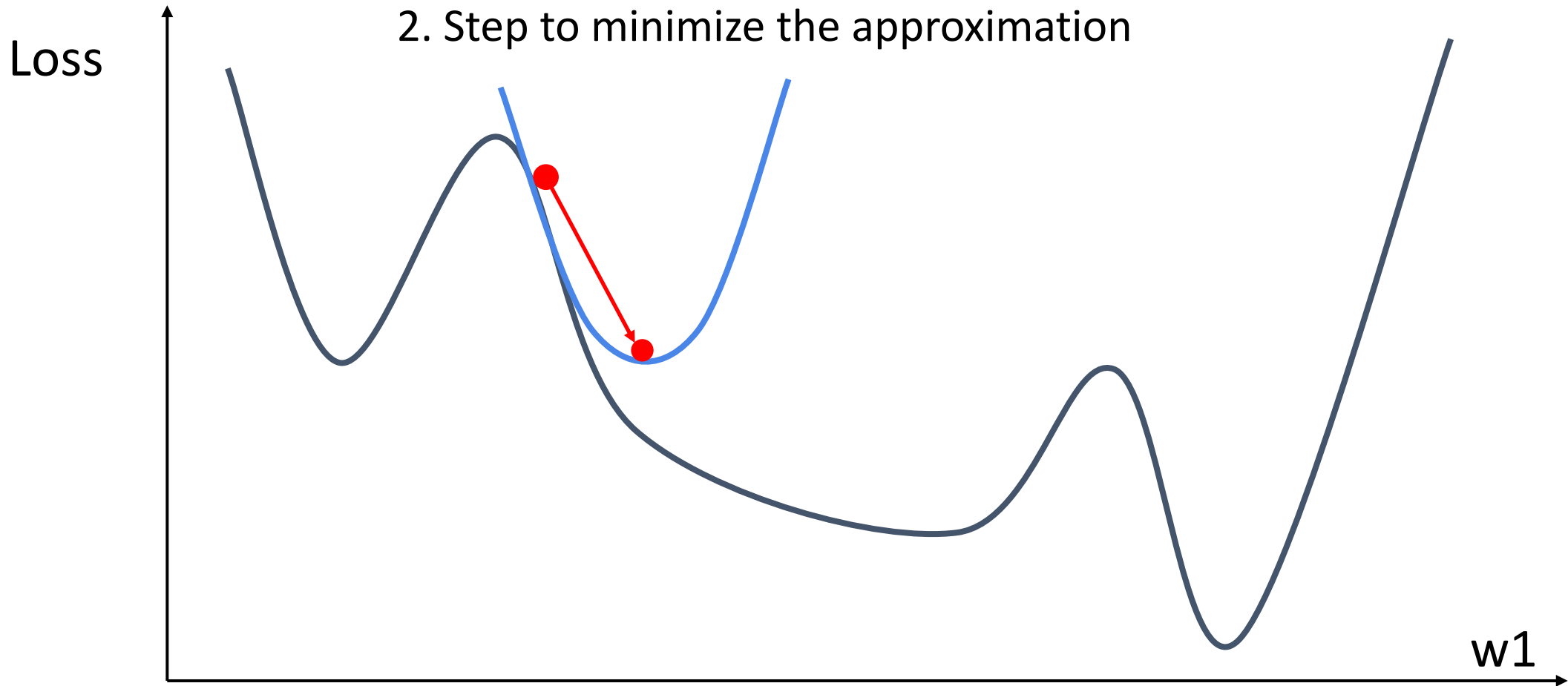
# So far: First-Order Optimization

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



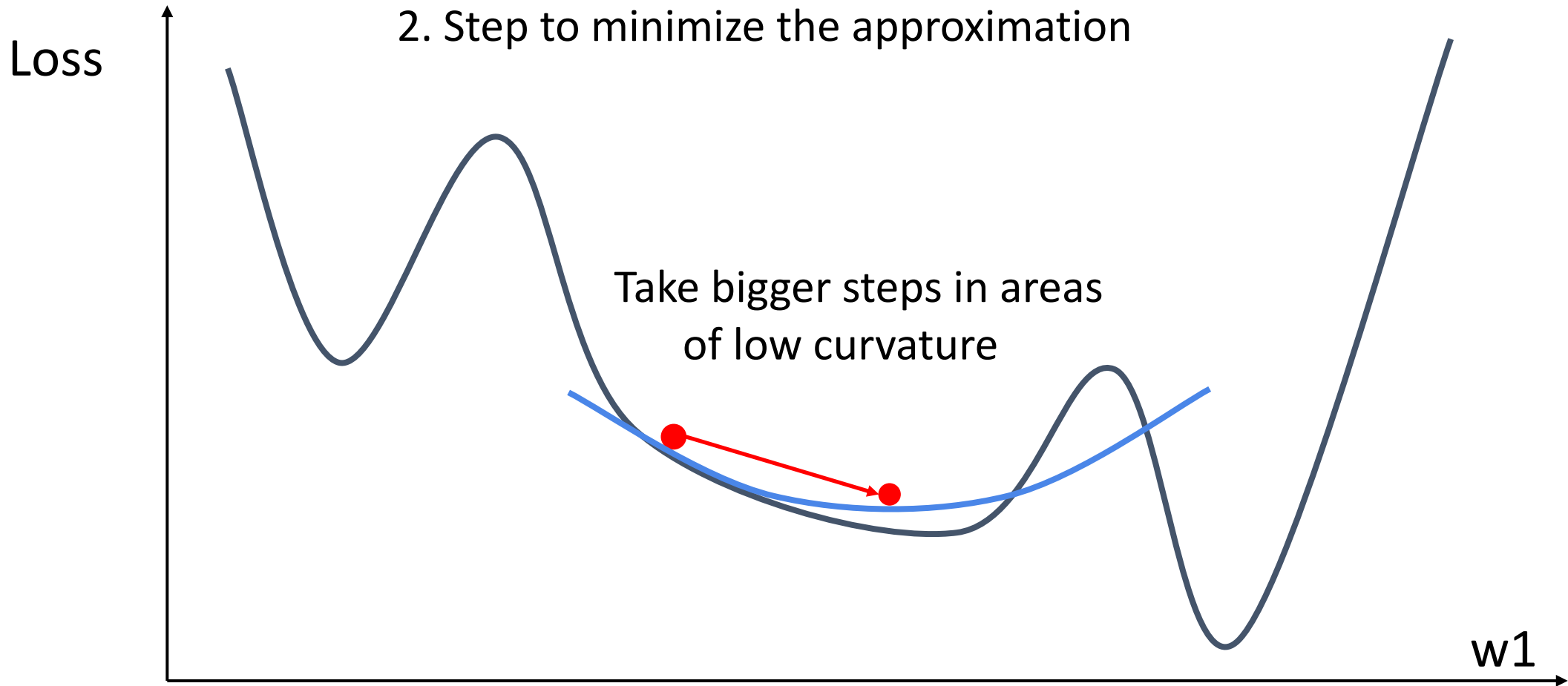
# Second-Order Optimization

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



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1. Use gradient and Hessian to make quadratic approximation
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# 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:

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

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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

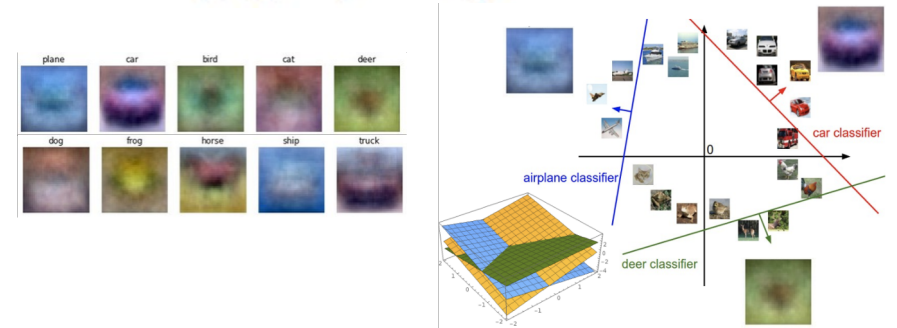
In practice:

- **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