Machine Intelligence & Deep Learning Workshop

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The Kate Gleason COLLEGE OF ENGINEERING

Introduction to Deep Learning



Raymond Ptucha June 27-29, 2018 Rochester Institute of Technology

www.rit.edu/kgcoe/cqas/machinelearning

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Agenda

Wed, June 27

- 9-10:30am

- 10:30-10:45pm

- 10:45-12:15pm

- 12:15-1:30pm

- 1:30-3:30pm

- 3:30-5pm

• Thur. June 28

9-10:30am10:30-10:45pm

- 10:45-12:15pm

10:45-12:15pm12:15-1:30pm

- 1:30-3:30pm

- 3:30-5.50pii

Fri, June 29

9-10:30am10:30-10:45pm

- 10:45-12:15pm

- 12:15-1:30pm

- 1:30-3:30pm

- 3:30-5pm

Regression and Classification

Break

Boosting and SVM

Lunah

Neural Networks and Dimensionality Reduction

Hands-on Python and Machine Learning

Introduction to deep learning

Break

Convolutional Neural Networks

Lunch

Region and pixel-level convolutions

Hands-on CNNs

Recurrent neural networks

Break

Language and Vision

Lunch

Graph convolutional neural networks; Generative adversarial networks

Hands-on regional CNNs, RNNs

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Machine Learning

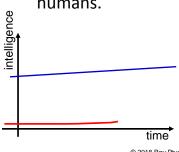
- Machine learning is giving computers the ability to analyze, generalize, think/reason/behave like humans.
- Machine learning is transforming medical research, financial markets, international security, and generally making humans more efficient and improving quality of life.
- Inspired by the mammalian brain, deep learning is machine learning on steroids- bigger, faster, better!

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The point of Singularity

 The point of singularity is when computers become smarter than humans.





Evolution of biology

Advancement of technology

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Unleashing of Intelligence



- Machines will slowly match, then quickly surpass human capabilities.
- Today it is exciting/scary/fun to drive next to an autonomous car.
- Tomorrow it may be considered irresponsible for a human to relinquish control from a car that has faster reaction times, doesn't drink/text/get distracted/tired, and is communicating with surrounding vehicles and objects.

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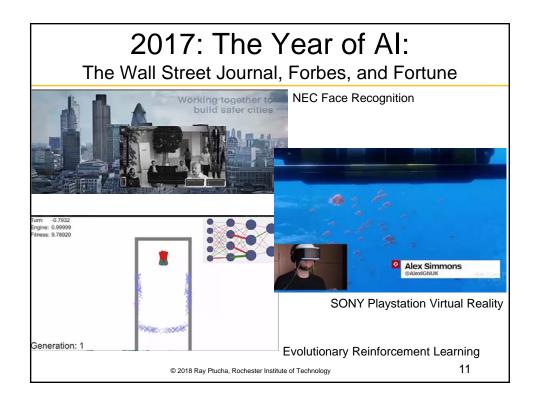
Why is AI (Deep Learning) Just Now Becoming Practical in Many Day-to-Day Situations?

- · Availability of data;
- Sustained advances in hardware capabilities (including GPUs running machine learning workloads);
- Omnipresent connectivity;
- Lower cost and power consumption;
- Sustained advances in algorithmic learning techniques.

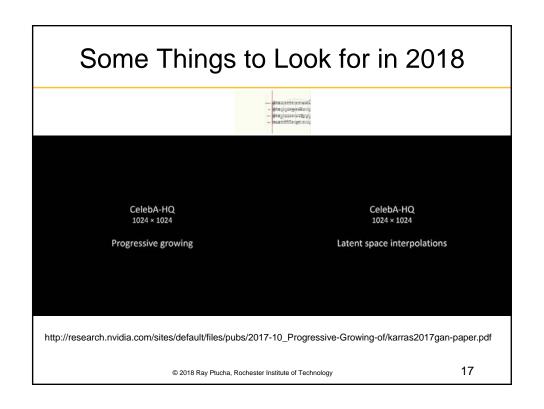
DL with super computers
DL with GPUs
Traditional ML
Amount of data

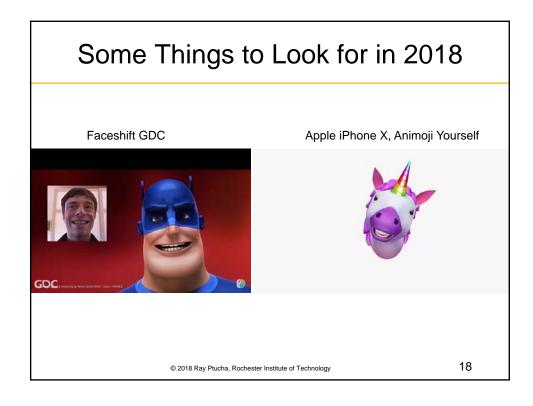
Hot trend:
High performance
architecture experts
teaming up with deep
learning experts

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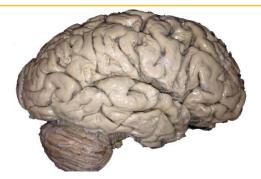
Prerequisites

- Basic probability and statistics
- Some programming experience (we will be using Python)
- Simple linear algebra
- Simple multivariate calculus
- Don't worry if you don't understand all concepts, over time things will make more sense.
- Feel free to stick around from 5-7pm in the evenings where the instructor and TAs can help you setup your own personal laptop, reinforce concepts, or answer personal applications of deep learning for your particular need.

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The Human Brain



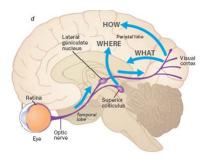
- We've learned more about the brain in the last 5 years than we have learned in the last 5000 years!
- It controls every aspect of our lives, but we still don't understand exactly how it works.

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The Brain on Pattern Recognition

• Airplane, Cat, Car, Dog





http://thebraingeek.blogspot.com/ 2012/08/blindsight.html

STL-10 dataset

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The Brain on Pattern Recognition

Despite Changes in Deformation:









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The Brain on Pattern Recognition

Despite Changes in Occlusion:







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The Brain on Pattern Recognition

Despite Changes in Size, Pose, Angle:













Tardar Sauce "Grumpy Cat"

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The Brain on Pattern Recognition

Despite Changes in Background Clutter:



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The Brain on Pattern Recognition

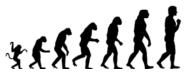
Despite Changes in Class Variation...



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Teaching Computers to See

to develop the marvel of the eye-brain. It took evolution 540M years



- Lets say a child collects a new image every 200msec.
- By age 3, this child has processed ~250M images.



(5 images/sec)(60 sec/min)(60 min/hr)(12 hr/day)(365 days/yr)(3 yrs) = 236 M

Today's computers can do this in a few days...

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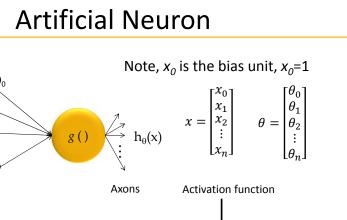
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Neural Nets on Pattern Recognition

- Instead of trying to code simple intuitions/rules on what makes an airplane, car, cat, and dog...
- · We feed neural networks a large number of training samples, and it will automatically learn the rules!
- We will learn the magic behind this today!



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dendrites
$$h_{\theta}(x) = g(x_0\theta_0 + x_1\theta_1 + \dots + x_n\theta_n) = g\left(\sum_{i=0}^n x_i\theta_i\right)$$

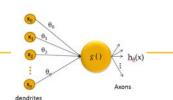
$$h_{\theta}(x) = g(\theta^T x)$$

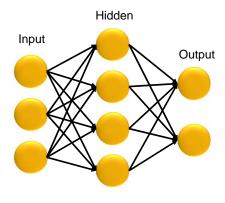
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Artificial Neural Networks

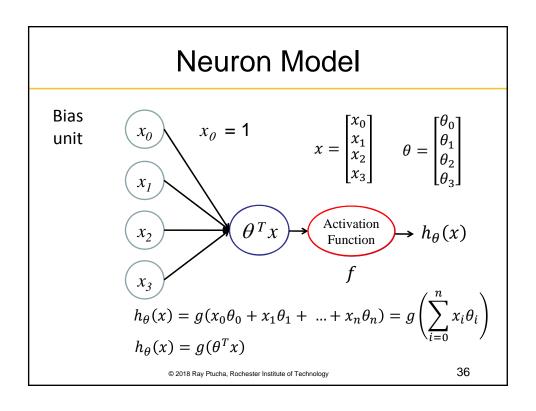
Artificial Neural
 Network (ANN) –
 A network of interconnected nodes that "mimic" the properties of a biological network of neurons



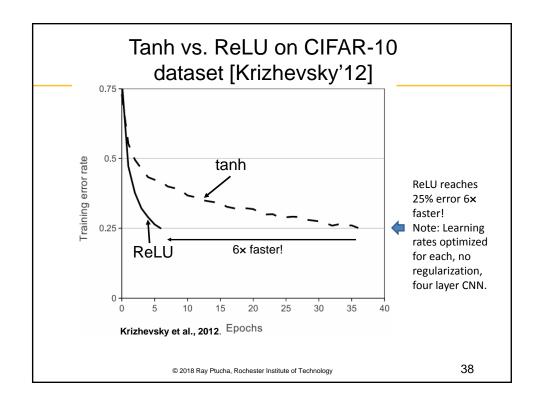


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A-Layer ANN Fully Connected Topology Input Layer Hidden Layer 1 Hidden Layer 2 Output Layer A 20x20 image would a b C nodes, where have 400 input nodes nodes nodes C is the number of classes Backpropagation (~1985) uses $\frac{\partial \Delta}{\partial w}$ for learning Learning happens in the weights- each line is a weight.



Activation Function Comparison Gradient of both Tanh saturates at zero. Sigmoid also non-**Sigmoid** zero centered, so in practice tanh performs better. $h_{\theta}(x) = max(0, x)$ Rectified Linear Units (ReLU) - Better for high dynamic range Faster learning - Overall better result - Neurons can "die" if allowed to grow unconstrained 37 © 2018 Ray Ptucha, Rochester Institute of Technology



Where Do Weights Come From?

- The weights in a neural network need to be learned such that the errors are minimized.
- Just like logistic regression, we can write a cost function.
- Similar to gradient descent, we can write an iterative procedure to update weights, with each iteration decreasing our cost.
- These iterative methods may be less efficient than a direct analytical solution, but are easier to generalize.

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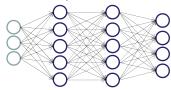
Backpropagation

- We need to solve weights of a network so that the error is minimized.
- Weights can be refined by changing each weight by an amount proportional to the partial derivative of the error with respect to each weight.
- Partial derivatives can be calculated by iteratively changing each weight and measuring the corresponding change in error.
- Hard to do with millions of weights!
- In 1986, a technique called back-propagation was introduced (D. E. Rumelhart, G. E. Hinton, and R. J. Williams "Learning representations by backpropagating errors," J. Nature 323, 533-536, 1986).

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Hyperparameters vs. parameters

- Hyperparameters are the tuning parameters for a nnet- say number of layers, nodes per layer, learning rate, momentum, regularization, etc.
- Parameters are the weights that are being learned. Ignoring bias terms, the below network has 3x5+5x5+5x4 = 60 parameters.
 - If we include bias terms, we have 4x5+6x5+6x4=74 learnable parameters.



Note: deep nets may contain 100M parameters with 100 layers!

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Initialization of Weights

- We initialize weights to small ± values centered on 0.
- If they were all initialized to 0, the network wouldn't learn anything.
 - ie: the output of each node would be equal, therefore the update would update each term identically.
- If they were all initialized to large values (values > +1 or < -1), the inputs to each node would be saturated.
- If they were all initialized to small values around 0, we would be in the linear portion of the activation function.

-W1 = 0.001*rand(length(h1),length(h2))

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Initialization of Weights

- Although the uniform distribution is good, the more inputs to a node, the greater its variance.
- To set output distribution of all nodes equal (this empirically improves convergence). Use h1=# input

Transfer Learning

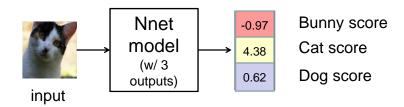
- Other solutions restrict weights: $-\epsilon_{init} < \theta_{ji} < \epsilon_{init}$ $\epsilon_{init} = \sqrt{6}/\sqrt{s_l + s_{l+1}}$ s_l , s_{l+1} are the No. nodes in layers around $\theta^{(l)}$
- For deep ReLU networks, He2015 showed:
 - W1 = 0.001*rand(length(h1),length(h2))./sqrt(2*length(h1))
- Works best and is recommended for them

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nodes

Multiclass Loss Functions



- The input image scores highest against cat, but is also somewhat similar to dog.
- How do we assign a loss function?

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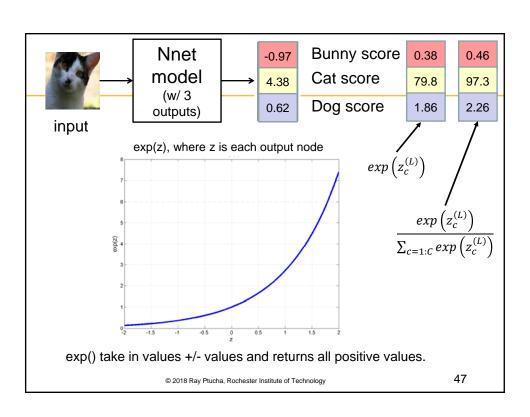
Activation Function of Output Layer

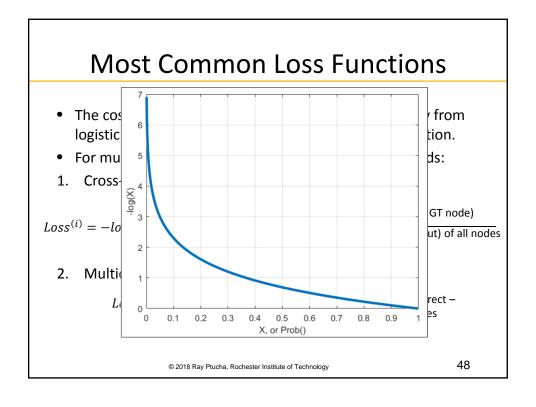
- Sigmoid returns 0 or 1 for each output node.
- What if you wanted a confidence interval?
- Use a linear activation function for regression: $a^{(l)}=z^{(l)}$
- Softmax often used for classification:

$$a_c^{(L)} = h_\theta(x)_c = g\left(z_c^{(L)}\right) = \frac{exp\left(z_c^{(L)}\right)}{\sum_{c=1:C} exp\left(z_c^{(L)}\right)} \underbrace{\qquad exp()}_{c} \text{ of each output nodes}$$

 Note: Only the output layer activation function changes- all hidden layer nodes activation functions would be the sigmoid/tanh/ReLU function.

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Most Common Loss Functions

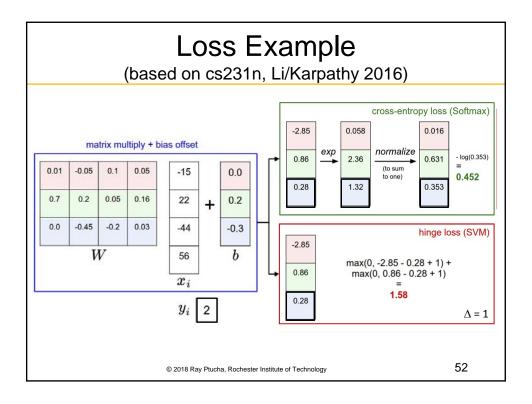
- The cost function we previously used was a direct copy from logistic regression and works great for binary classification.
- For multi-class, there are two popular data loss methods:
- 1. Cross-entropy loss, which uses softmax:

$$Loss^{(i)} = -log\left(\frac{exp\left(out_{yi}^{(i)}\right)}{\sum_{c=1:C} exp\left(out_{c}^{(i)}\right)}\right) \qquad \text{Loss for sample } i = \frac{\exp(\text{output of GT node})}{\text{Sum of exp(output) of all nodes}}$$

2. Multiclass SVM Loss (Weston Watkins formulation):

$$Loss^{(i)} = \sum_{j \neq yi} max(0, out_j - out_{yi} + \Delta)$$
 Sum of incorrect – correct classes

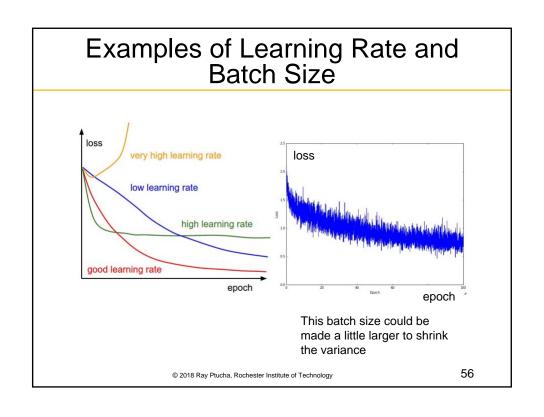
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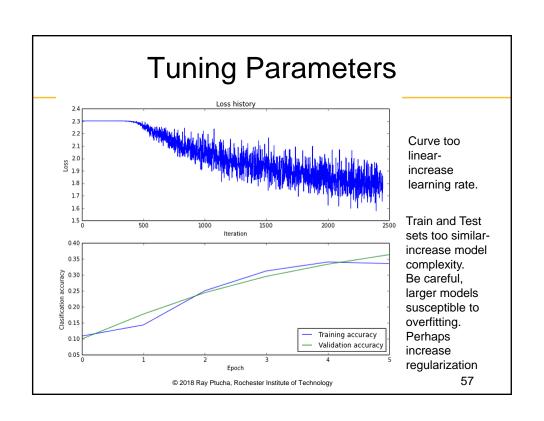


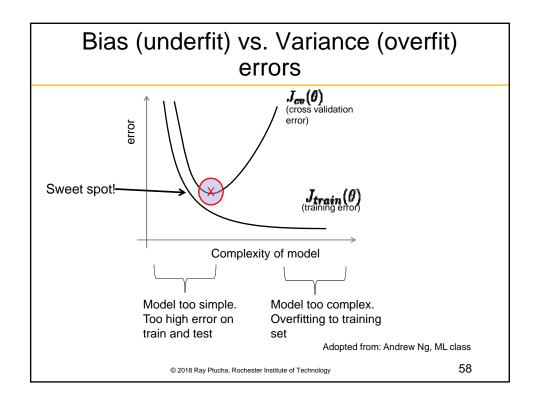
Sequential vs. Batch Training

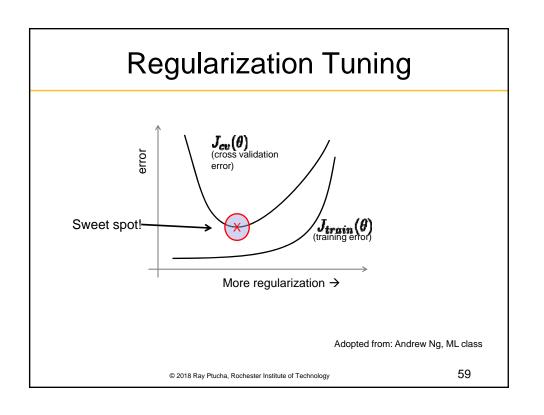
- Back propagation can be done:
 - Sequential: one sample at a time,
 - Weights are shifted back and forth quite a bit
 - Group (minibatch): a group of samples at a time, or
 - Batch: all training samples at once
 - Weights are shifted in direction that makes most input samples better
 - Generally converges the fastest
- Recommended to use largest minibatch possible and stay within memory constraints of hardware.

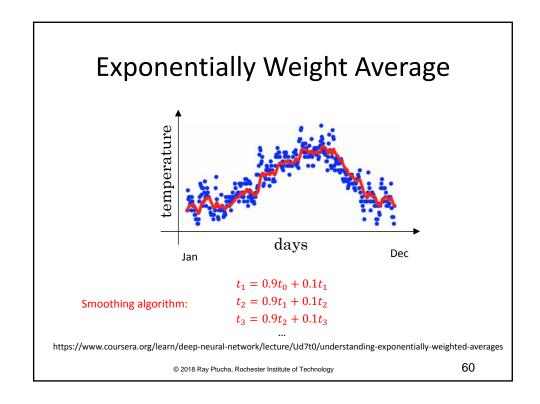
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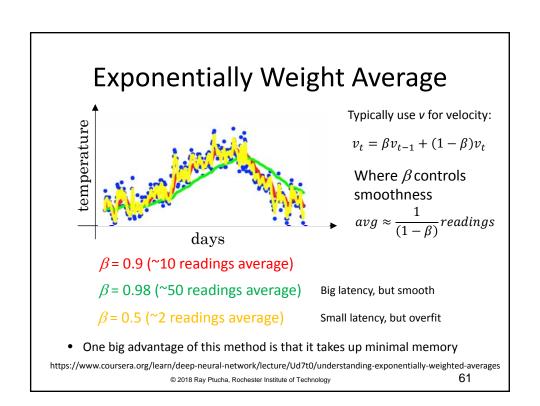










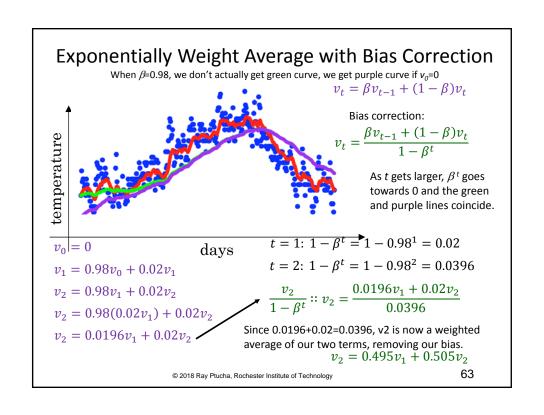


Exponentially Weight Average with Bias Correction When
$$\beta$$
=0.98, we don't actually get green curve, we get purple curve if v_0 =0
$$v_t = \beta v_{t-1} + (1-\beta)v_t$$
 Bias correction:
$$v_t = \frac{\beta v_{t-1} + (1-\beta)v_t}{1-\beta t}$$

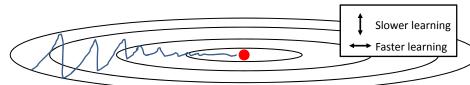
$$v_t = \frac{\beta v_{t-1} + (1-\beta)v_t}{1-\beta t}$$

$$v_t = \frac{\beta v_{t-1} + (1-\beta)v_t}{1-\beta t}$$

$$v_t = \frac{0.98v_0 + 0.02v_1}{0.02} = v_1$$
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Gradient Descent with Momentum



- Typical gradient descent- oscillations slow down progress
- Further, if learning rate too large, initial step would cause significant overshooting, and perhaps divergence.

For each iteration, t:

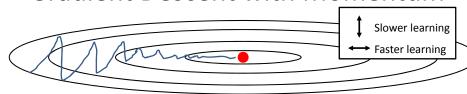
Compute dW, db on current mini-batch

$$\begin{array}{l} V_{dW} = \beta V_{dW} + (1-\beta)dW \\ V_{db} = \beta V_{db} + (1-\beta)db \end{array} \right] \mbox{Replace each partial derivative} \\ W = W - \eta V_{dW} \\ b = b - \eta V_{db} \end{array} \right] \mbox{Update using smoothed derivatives}$$

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Gradient Descent with Momentum



For each iteration, t:

Compute dW, db on current mini-batch

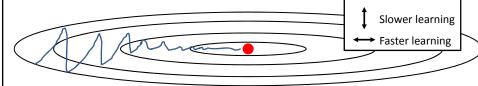
$$\begin{array}{l} V_{dW} = \beta V_{dW} + (1-\beta)dW \\ V_{db} = \beta V_{db} + (1-\beta)db \end{array} \right] \text{ Replace each partial derivative } \\ W = W - \eta V_{dW} \\ b = b - \eta V_{db} \end{array} \right] \text{ Update using smoothed derivatives}$$

We end up with two hyperparameters, η and β (learning rate and momentum) Very common to set β =0.9

Because moving average moves up quickly, bias correction not commonly used

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Gradient Descent with Momentum



For each iteration, t:

Compute dW, db on current mini-batch

$$\begin{aligned} V_{dW} &= \beta V_{dW} + (1-\beta)dW \\ V_{db} &= \beta V_{db} + (1-\beta)db \end{aligned} \quad \begin{array}{l} \text{Sometimes} \\ \text{approximated} \\ \text{with:} \end{aligned} \quad \begin{array}{l} V_{dW} &= \beta V_{dW} + dW \\ V_{db} &= \beta V_{db} + db \end{aligned}$$

$$W &= W - \eta V_{dW}$$

$$b &= b - \eta V_{db}$$

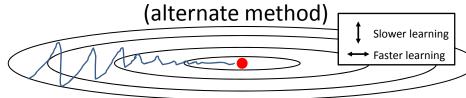
We end up with two hyperparameters, η and β (learning rate and momentum) Very common to set $\beta\!\!=\!\!0.9$

Because moving average moves up quickly, bias correction not commonly used

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Gradient Descent with Momentum (alternate method)



For each iteration, t:

Compute dW, db on current mini-batch

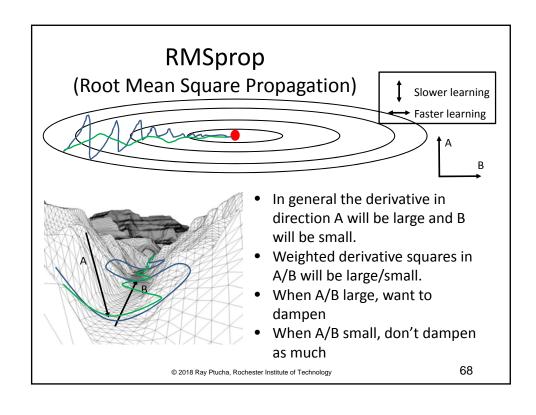
$$V_{dW} = \beta V_{dW} - \eta dW$$
 Smoothed version of dW, but include learning rate to increase momentum more.

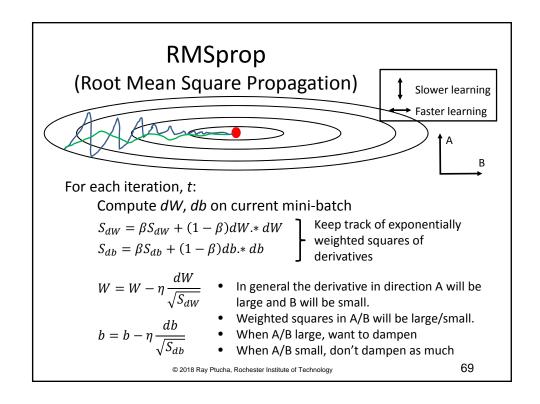
$$W = W + V_{dW}$$
 Since we were learning '-' dW above, we add, not subtract here.

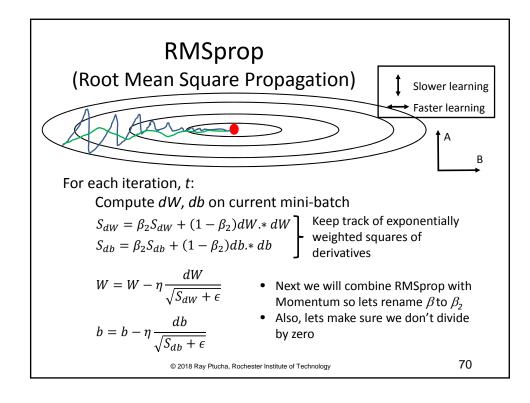
We end up with two hyperparameters, η and β (learning rate and momentum) Very common to set β =0.9

Because moving average moves up quickly, bias correction not commonly used

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Adam Optimization

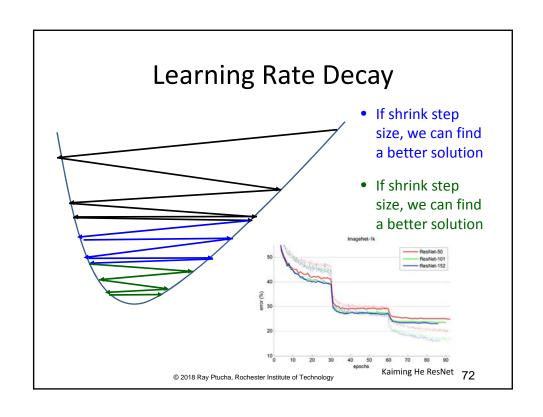
Combine momentum with RMSprop with Bias Correction

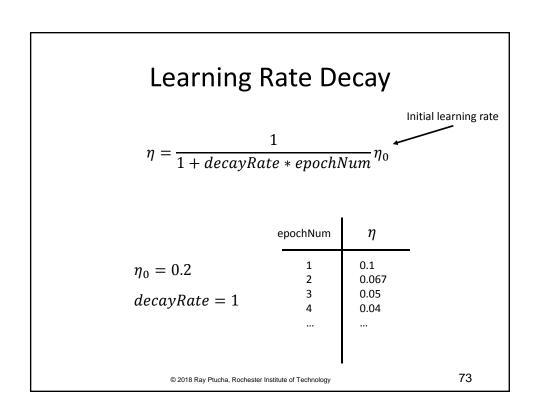
 V_{dW} =0, S_{dW} =0, V_{db} =0, S_{db} =0 For each iteration, t:

Compute dW, db on current mini-batch

momentum
$$\begin{cases} V_{dW} = \beta_1 V_{dW} + (1-\beta_1) dW & \text{Hyperparameters to tune:} \\ V_{db} = \beta_1 V_{db} + (1-\beta_1) db & \beta_2 : 0.99 \\ S_{dW} = \beta_2 S_{dW} + (1-\beta_2) dW.* dW & \epsilon : 10^{-8} \\ S_{db} = \beta_2 S_{db} + (1-\beta_2) db.* db & \pi: \text{need to solve} \end{cases}$$
Bias Correction
$$\begin{cases} V_{dW}^{Corrected} = V_{dW}/(1-\beta_1^t); & V_{db}^{Corrected} = V_{db}/(1-\beta_1^t) \\ S_{dW}^{Corrected} = S_{dW}/(1-\beta_2^t); & S_{db}^{Corrected} = S_{db}/(1-\beta_2^t) \end{cases}$$

$$W = W - \eta \left(V_{dW}^{Corrected} / \sqrt{S_{dW}^{Corrected} + \epsilon} \right); & b = b - \eta \left(V_{db}^{Corrected} / \sqrt{S_{db}^{Corrected} + \epsilon} \right)$$



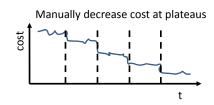


Learning Rate Decay- Other formulas

Some hyperparameter <1
$$\eta = 0.95^{epochNum} \cdot \eta_0 \qquad \text{(Exponential decay)}$$

$$\eta = \frac{k}{\sqrt{epochNum}} \cdot \eta_0 \quad \text{\tiny --OR--} \quad \eta = \frac{k}{\sqrt{t}} \cdot \eta_0 \quad \text{\tiny t can be custom} \quad \text{to problem}$$

Discrete staircase: Divide learning rate in half every 100 iterations



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Local Optima in Neural Networks

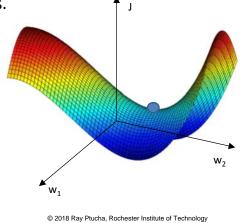
Local Optima
Global Optima

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Tocal Optima

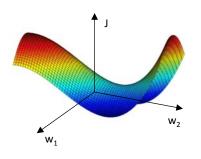
Local Optima in Neural Networks

 As luck would have it, most points with gradient zero are not like previous slide, but occur on saddle points.

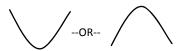


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Local Optima in Neural Networks



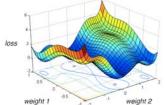
For a gradient to be zero, in each direction we either have a concave or a convex shape.



Now...in a very high dimensional space, say 10,000 dimensions, what are the odds all will be concave?

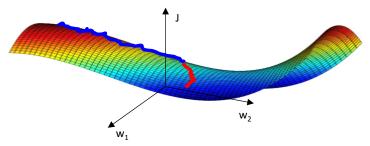
Pretty small!

We are much more likely to have some directions run up, and some run down, like the saddle point on the upper left or chaos below.



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Problem of Plateaus



- Given a surface like this, we traverse the cost function for a long time making little progress.
- Eventually, or perhaps by luck, we step to the side and finally begin to learn more aggressively again.
- But...thankfully in high dimensional spaces, unlikely to get stuck in local optima

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Batch Normalization: Accelerating Deep Network Training by Reducing Internal Covariate Shift

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Abstract

Training Deep Neural Networks is complicated by the fact that the distribution of each layer's inputs changes during training, as the parameters of the previous layers change. This slows down the training by requiring lower learning rates and careful parameter initialization, and makes it notoriously hard to train models with saturating nonlinearities. We refer to this phenomenon as internal covariate shift, and address the problem by normalizing layer inputs. Our method draws its strength from making normalization a part of the model architecture and performing the normalization for each training

minimize the loss

$$\Theta = \arg \min_{\Theta} \frac{1}{N} \sum_{i=1}^{N} \ell(\mathbf{x}_i, \Theta)$$

where $x_{1...N}$ is the training data set. With SGD, the training proceeds in steps, at each step considering a *minibatch* $x_{1...m}$ of size m. Using mini-batches of examples, as opposed to one example at a time, is helpful in several ways. First, the gradient of the loss over a mini-batch $\frac{1}{m}\sum_{i=1}^{m}\frac{\partial \ell(x_i,\Theta)}{\partial \Theta}$ is an estimate of the gradient over the training set, whose quality improves as the batch size increases. Second, computation over a mini-batch can be more efficient than m computations for individual examples on modern computing platforms.

http://jmlr.org/proceedings/papers/v37/ioffe15.pdf

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Batch Normalization

- Makes hyperparameter search easier
- Makes learning more robust to the choice of hyperparameters
- Makes learning work with a larger range of hyperparameters
- Enables easier and often faster training of very deep networks

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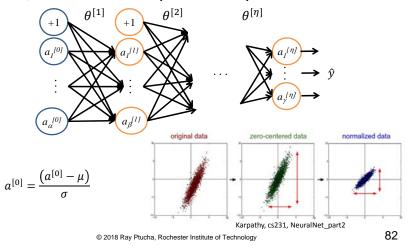
Normalization of Features

- We have seen how it is beneficial to subtract the mean and divide by the standard deviation for our input vector X (note, each dimension has its own μ and σ).
- In a deeper net, it would make the training of $\theta^{[l]}$ more efficient if the input to the layer, $a^{[l-1]}$ was also normalized.
- Note, in batch normalization, we will actually normalize z, not a. There is some disagreement on this, but most normalize z.

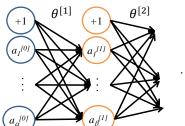
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Implementing Batch Normalization

• Analogous to the way we normalized the input to $\theta^{[1]}$, normalize the input to all layers:

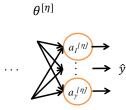


Implementing Batch Normalization



 As we then pass training samples through the net, replace z^[i] with z_{norm}^[l]:

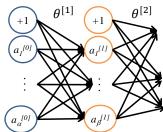
$$z_{norm}^{[l]} = \frac{\left(z^{[l]} - \mu^{[l]}\right)}{\sqrt{\sigma^{[l]2} + \epsilon}}$$

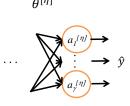


- Normalize z, not a
- Calculate μ and σ^2 from minibatch training samples:
- $\mu^{[l]} = \frac{1}{n} \sum_{i} z^{[l](i)}$
- $\left(\sigma^{[l]}\right)^{2} = \frac{1}{n} \sum_{i} (z^{[l](i)} \mu^{[l]})^{2}$

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Implementing Batch Normalization





$$z_{norm}^{[l]} = \frac{\left(z^{[l]} - \mu^{[l]}\right)}{\sqrt{\sigma^{[l]2} + \epsilon}}$$

$$\tilde{z}^{[l]} = \gamma z_{norm}^{[l]} + \beta$$

Then go through activation function

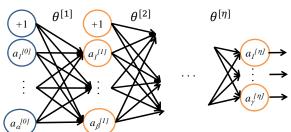
 But, do we don't always want our distribution to have mean=0 and unit variance.

So we learn two new parameters, γ and β , these will be learned along with our $\theta^{[l]'}$ s

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Implementing Batch Normalization



$$z_{norm}^{[l]} = \frac{\left(z^{[l]} - \mu^{[l]}\right)}{\sqrt{\sigma^{[l]2} + \epsilon}}$$

$$\tilde{z}^{[l]} = \gamma^{[l]} z_{norm}^{[l]} + \beta^{[l]}$$

Then go through activation function

Note $\gamma^{[l]}$ and $\beta^{[l]}$ are vectors for each layer, and get applied independently to each dimension.

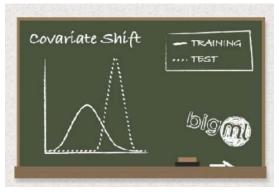
• If using batch normalization with $\beta^{[l]}$, no need to use bias values as they both do the same job.

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Batch Normalization Intuition

 When training set statistics differ from test set statistics, we call this covariate shift.



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Batch Normalization Intuition

- When training set statistics differ from test set statistics, we call this covariate shift.
- Typically we would have to relearn the model if we have a covariate shift, but batch norm minimizes covariate shift.
- In effect, it forces the input to each layer to have a predictable distribution, μ =0, σ =1, or whatever distribution the γ and β determine works best for that layer.
- This predictable distribution makes later layers somewhat independent of earlier layers, enabling robust behavior even in deep networks.

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Batch Normalization at Test Time

- Using mini-batch training samples, we calculate μ and σ^2 for each layer:
- $\mu^{[l]} = \frac{1}{n} \sum_{i} z^{[l](i)}$
- $\left(\sigma^{[l]}\right)^2 = \frac{1}{n} \sum_{i} (z^{[l](i)} \mu^{[l]})^2$

$$z_{norm}^{[l]} = \frac{\left(z^{[l]} - \mu^{[l]}\right)}{\sqrt{\sigma^{[l]^2} + \epsilon}}$$

$$\tilde{z}^{[l]} = \gamma z_{norm}^{[l]} + \beta$$

- During test, if we were testing a mini-batch at once, we could use the test mini-batch to calculate μ and σ.
- Alternately, if only have one or a few samples you need an estimate for μ and σ .
 - Can compute during training using a full epoch.
 - Can use an exponentially weighted average for each:

$$v_t = \beta v_{t-1} + (1 - \beta)v_t$$

Note: this β is from exponentially weighted average and not related to batch norm β !

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Andrew Ng, 2017 https://www.deeplearning.ai/

Deep Learning Specialization, Five courses:

- 1. Neural Networks and Deep Learning
- 2. Improving Deep Neural Networks
- 3. Structured Machine Learning Projects
- 4. Convolutional Neural Networks
- 5. Sequence Models

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CS231n: Convolutional Neural Networks for Visual Recognition







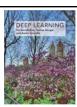


Li, Johnson, Yeung 2017 http://cs231n.stanford.edu/ I

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Books on Deep Learning



- Grokking Deep Learning by Andrew Trask. Use Udacity discount code traskud17 for 40% off. This provides a very gentle introduction to Deep Learning and covers the intuition more than the theory.
 - https://www.manning.com/books/grokking-deep-learning (\$49)
- Neural Networks And Deep Learning by Michael Nielsen. This book is more rigorous than Grokking Deep Learning and includes a lot of fun, interactive visualizations to play with.
 - http://neuralnetworksanddeeplearning.com/ (free!)
- The Deep Learning Textbook from Ian Goodfellow, Yoshua Bengio, and Aaron Courville. This online book contains a lot of material and is the most rigorous of the three books suggested.
 - http://www.deeplearningbook.org/ (html- free!)
 - https://www.amazon.com/Deep-Learning-Adaptive-Computation-Machine/dp/0262035618/ (hard copy \$50)

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