

Deep Learning

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- Vanishing Gradient & Activation Functions
- Dropout
- Batch Normalization

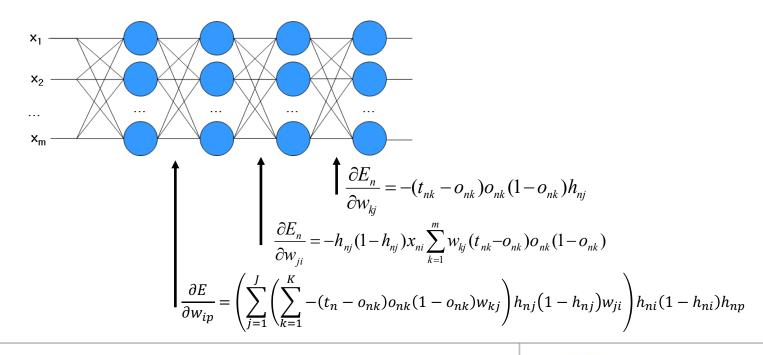


Gradient Vanishing & Activation Functions

Gradient Vanishing & Exploding

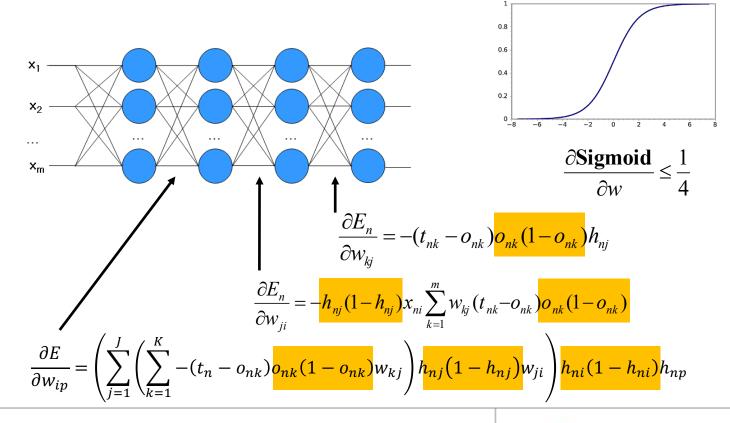
Gradient is easy to vanish or explode

- To many terms are multiplied.
- If some are small numbers, gradient becomes very small.
- If some are large numbers, gradient becomes very large.



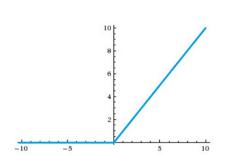
Vanishing Gradient

The major terms are the derivatives of the activation function

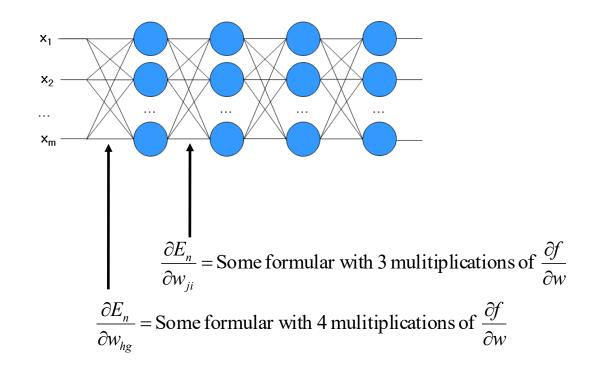


Using another functions instead of sigmoid

Rectified Linear Unit (ReLU)



$$f(x) = \begin{cases} x & \text{if } x > 0\\ 0 & \text{otherwise} \end{cases}$$



Advantage

- No vanishing gradient problems.
 - Deep networks can be trained without pre-training
- Sparse activation
 - In a randomly initialized network, only about 50% of hidden units are activated
- Fast computation:
 - 6 times faster than sigmoid function

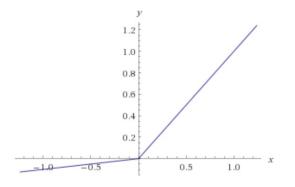
Disadvantage

Knockout Problem

You may use another

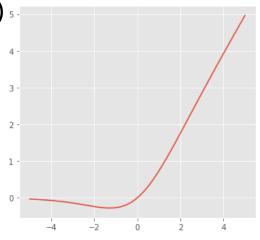
Leaky ReLU

$$f(x) = \begin{cases} x & \text{if } x > 0\\ 0.01x & \text{otherwise} \end{cases}$$

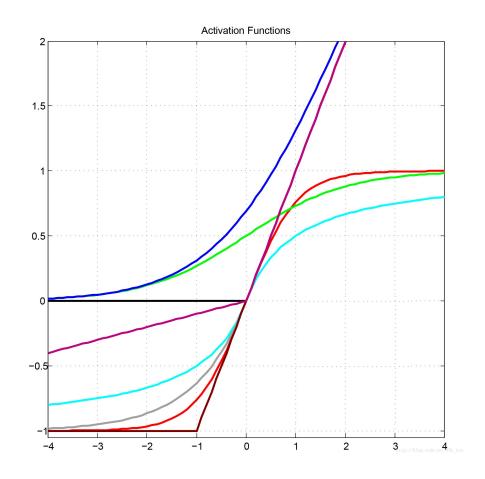


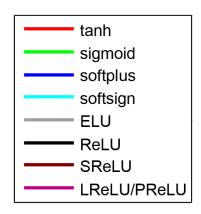
Swish (or SiLU-Sigmoid Linear Unit)

$$f(x) = \frac{x}{1 + e^{-x}}$$



Other Activation Functions





Summary

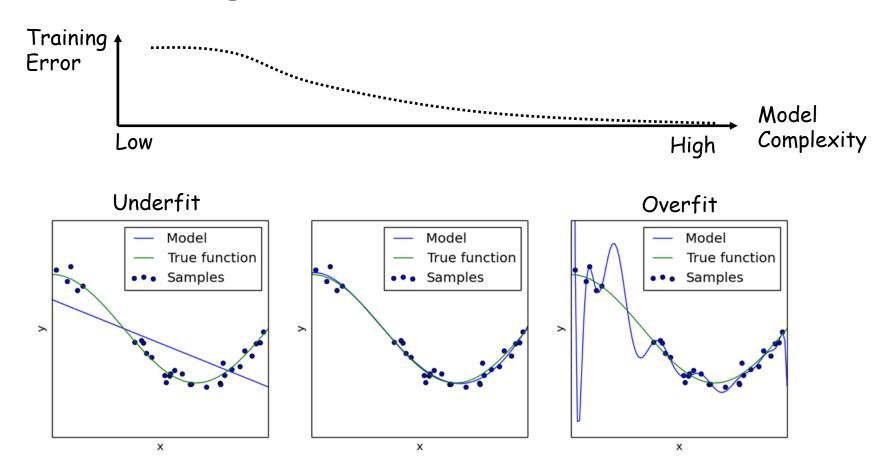
- Sigmoid functions and their combinations generally work better but are sometimes avoided due to the vanishing gradient problem
- ReLU function is a general activation function and is used in most cases these days
- If we encounter a case of dead neurons in our networks the leaky ReLU function is the best choice
- ReLU function is usually used in the hidden layers
- As a rule of thumb, you can begin with using ReLU function and then move over to other activation functions in case ReLU doesn't provide with optimum results



Regularization

Overfitting

Overfitting



Regularization

What is Regularization

Introducing additional information to prevent over-fitting

Approaches

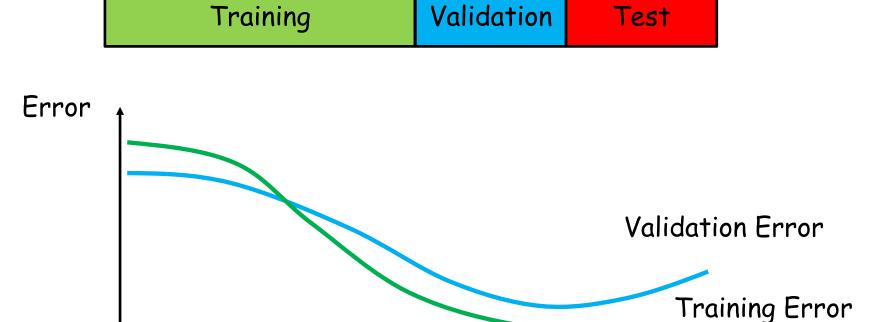
Proper Learning: Early stopping

Proper Structure: Weight decay, Dropout,

DropConnect, Stochastic pooling

Early Stopping

Split data into 3 groups



of updates

L1 Regularization

- Leading most weights very close to zero
- Choosing a small subset of most important inputs
- Resistant to noise in the inputs.

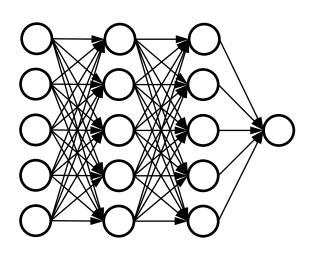
$$\widetilde{E}(\mathbf{w}) = E(\mathbf{w}) + \frac{\lambda}{2} |\mathbf{w}|$$

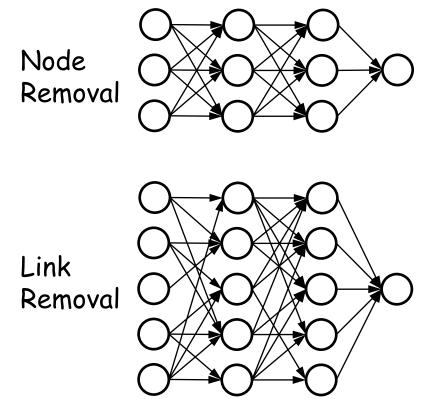
L2 Regularization

- Penalizing peaky weights
- Encouraging to use all of its inputs a little rather than using only some of its inputs a lot.

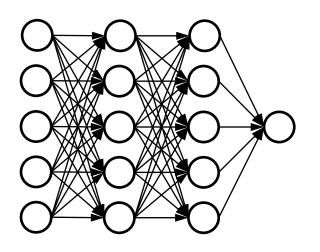
$$\widetilde{E}(\mathbf{w}) = E(\mathbf{w}) + \frac{\lambda}{2} \mathbf{w}^T \mathbf{w}$$

Complex Structure vs Simple Structure

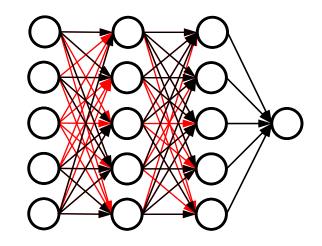




Complex Structure vs Simple Structure



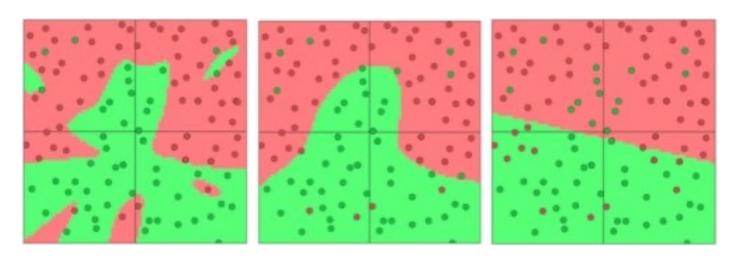
Set many links to zero



|w| is large <-> NN is Complex

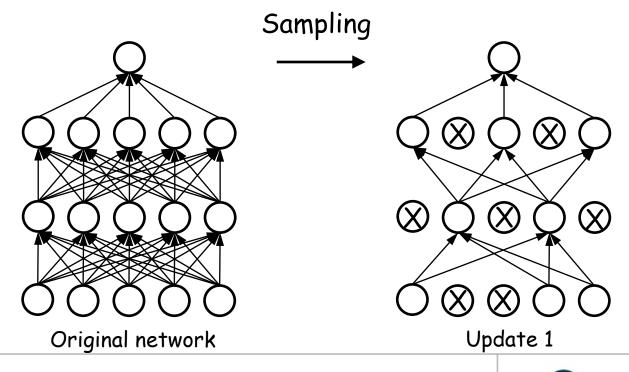
|w| is small \leftarrow NN is Simple

Example: Separating green and red

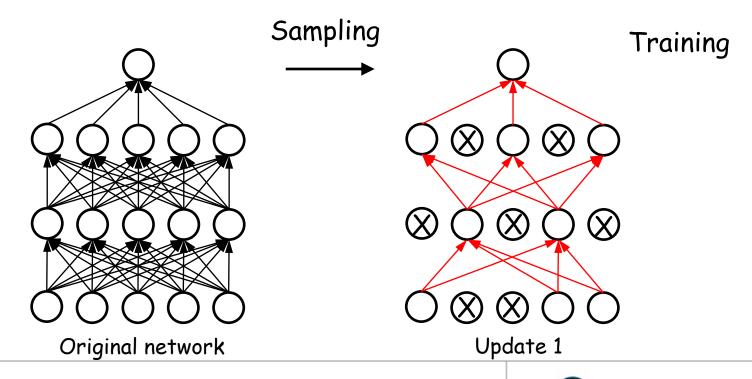


L2 regularization strengths of 0.01, 0.1, and 1

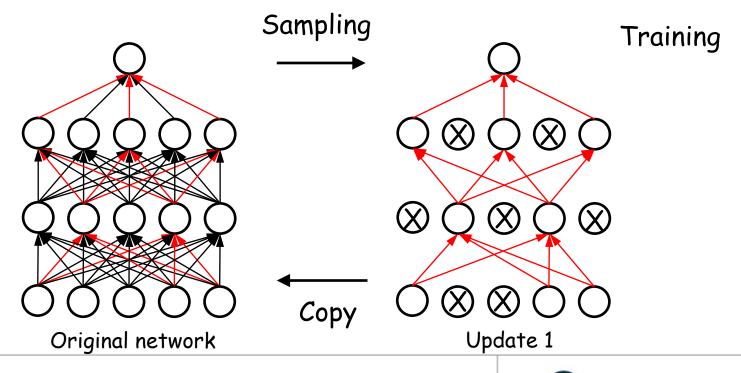
- How can we reduce the structural complexity without removing nodes?
 - Hmm??



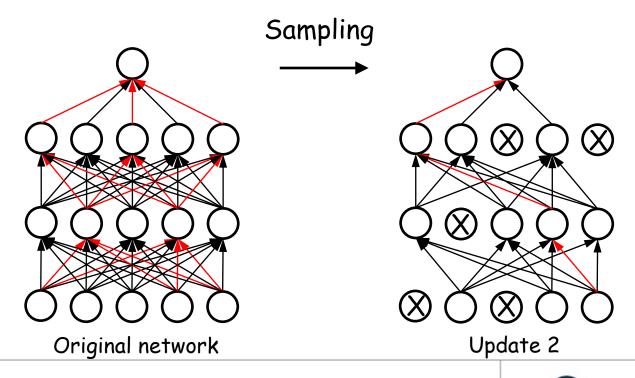
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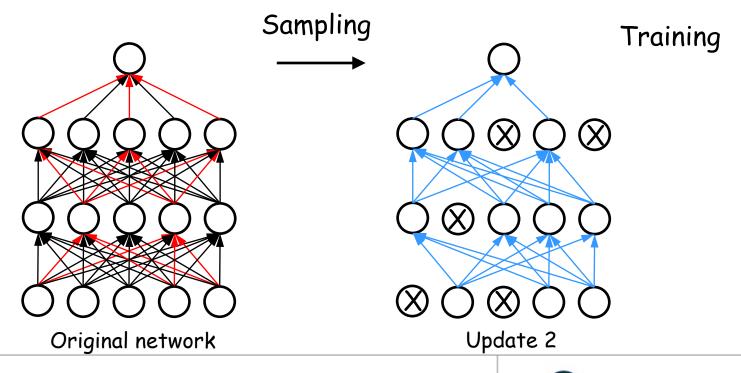
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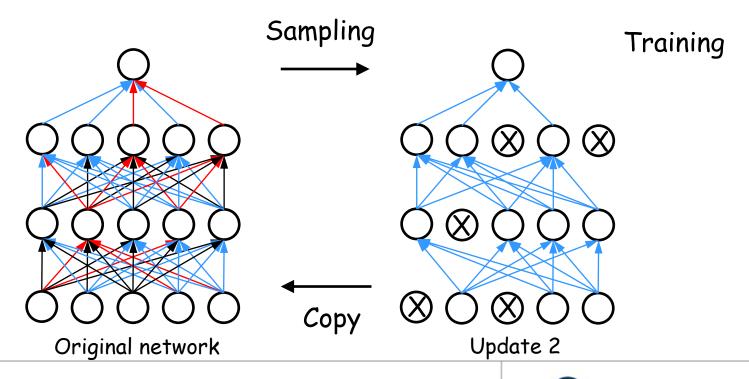
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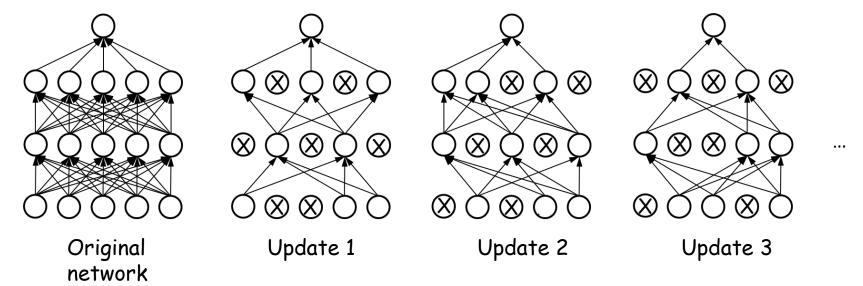


- How can we reduce the structural complexity without removing nodes?
 - Hmm??



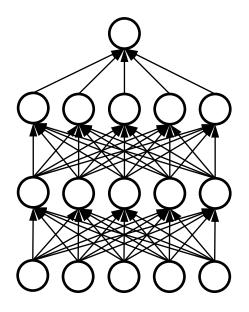
Do this at every epoch

- Randomly choose nodes with a probability of p
 - Usually p = 0.5
- Train the simplified neural network
 - At every epoch, we train different neural network which share connection weight each other



Testing

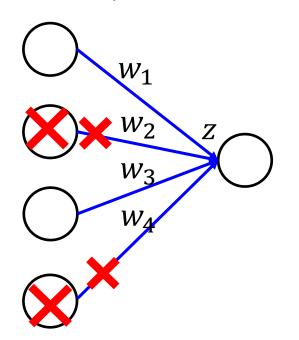
Use all the nodes without dropout



Testing

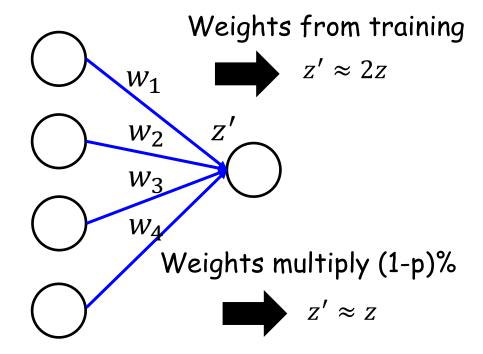
Training of Dropout

Assume dropout rate is 50%



Testing of Dropout

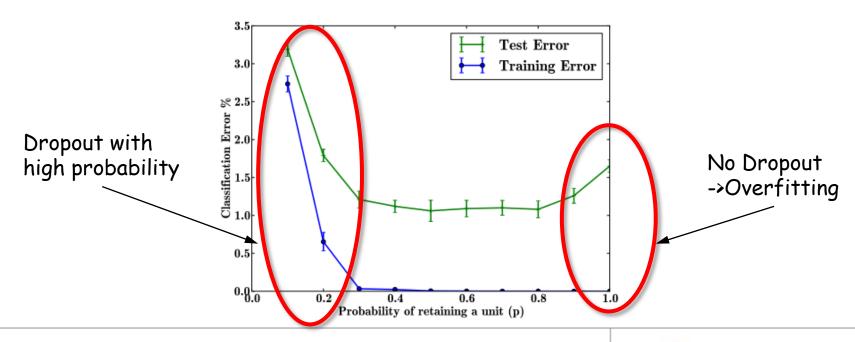
No dropout





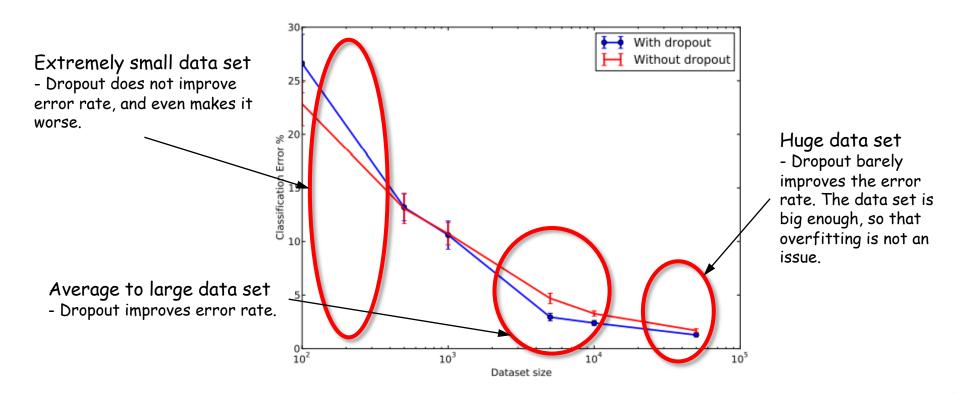
The effect of the dropout rate p:

- An architecture of 784-2048-2048-2048-10 is used on the MNIST dataset.
- The dropout rate p is changed from small numbers (most units are dropped out) to 1.0 (no dropout).



The effect of data set size:

 An architecture of 784-1024-1024-2048-10 is used on the MNIST dataset.



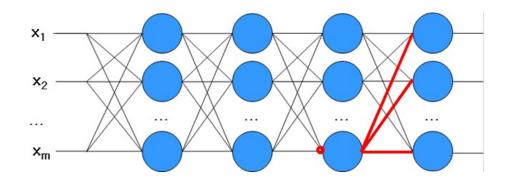
Summary

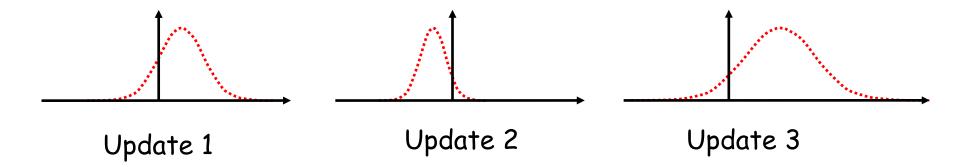
- Dropout is a very good and fast regularization method.
- Dropout is a bit slow to train (2-3 times slower than without dropout).
- If the amount of data is average-large dropout excels.
 When data is big enough, dropout does not help much.
- Dropout achieves better results than former used regularization methods (Weight Decay).



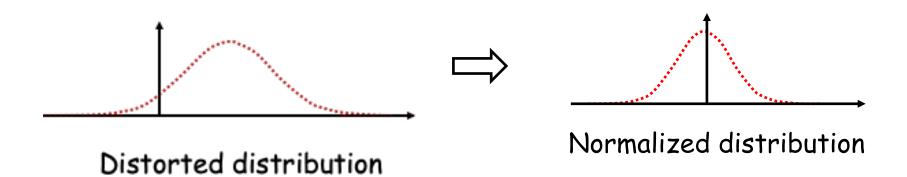
Distribution Shift

Output distribution of the red node

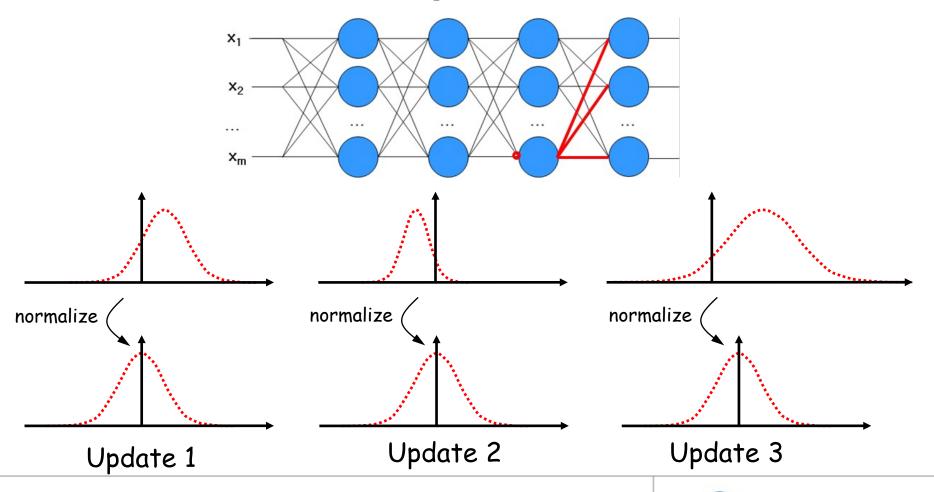




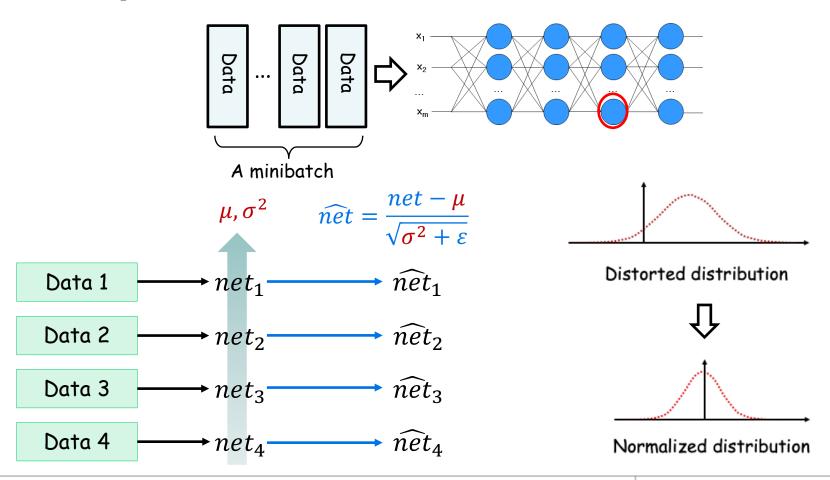
- Distribution Shift
 - It disturbs the learning process,
 - Learning is getting slow down
- Why don't we normalize the distribution of inputs



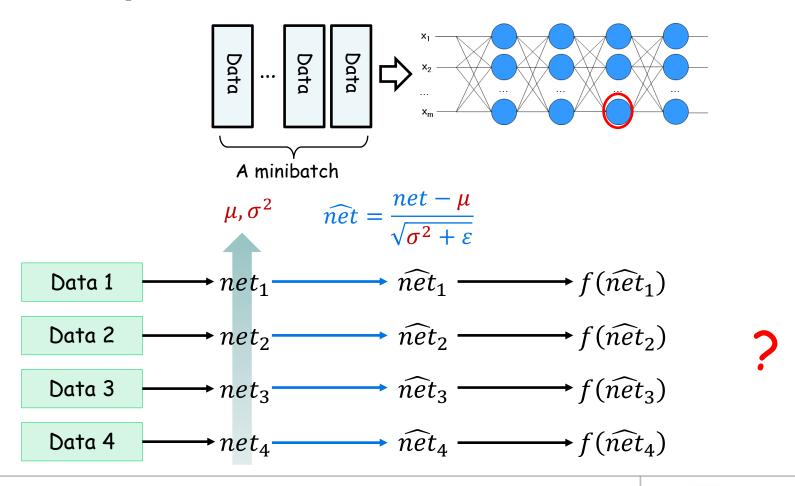
Normalization of outputs



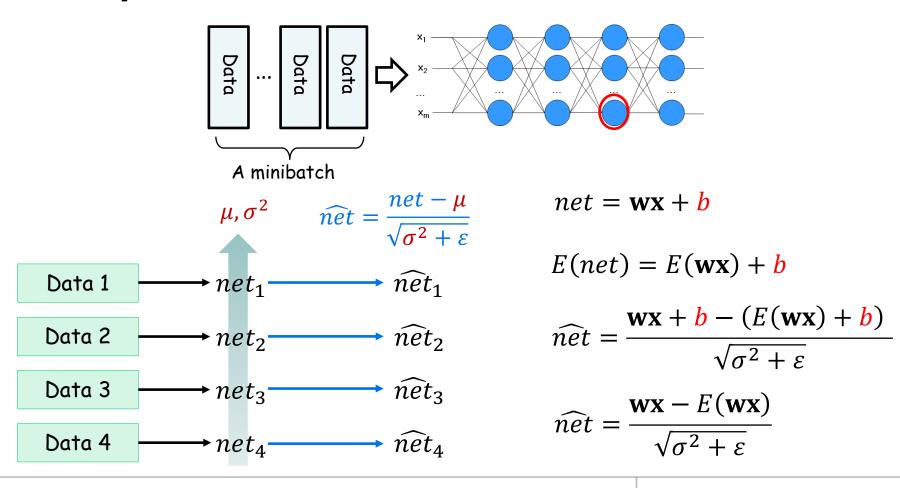
Input Normalization



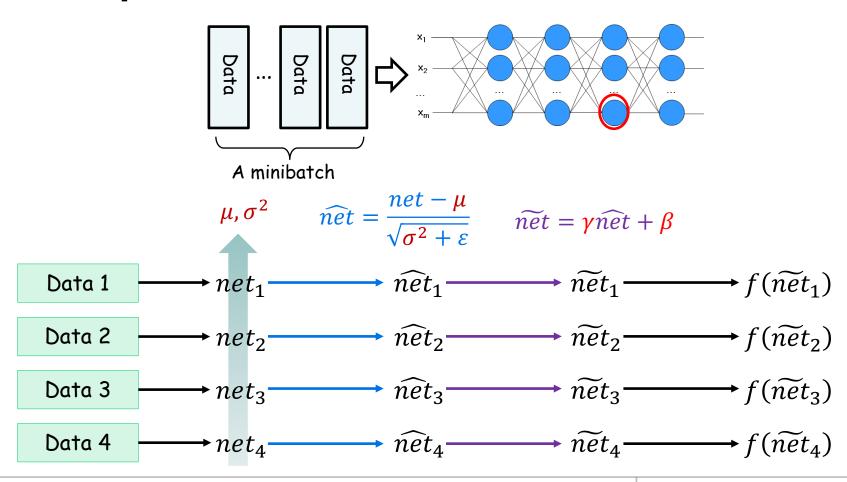
Input Normalization



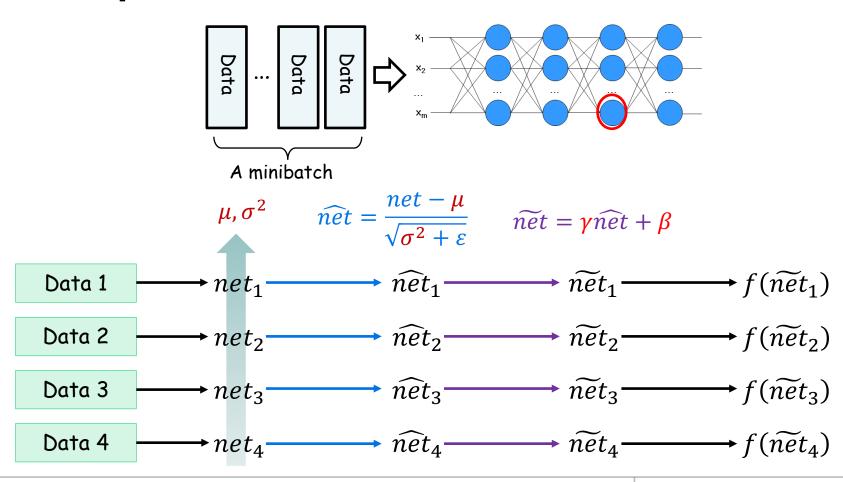
Input Normalization



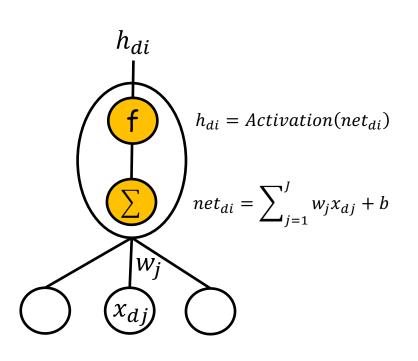
Input Normalization

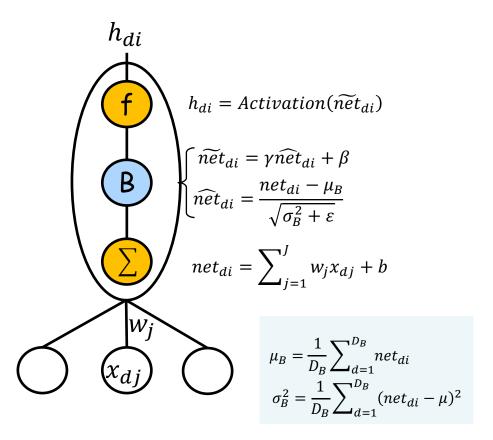


Input Normalization



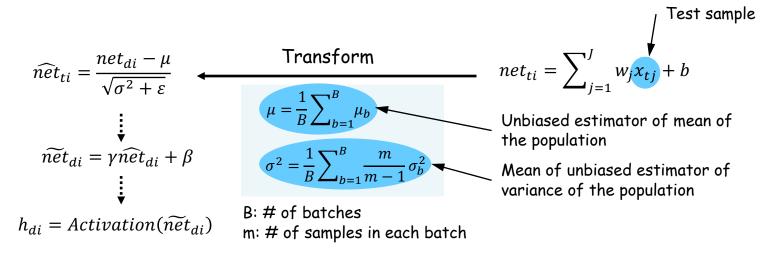
For a Single Node





Testing

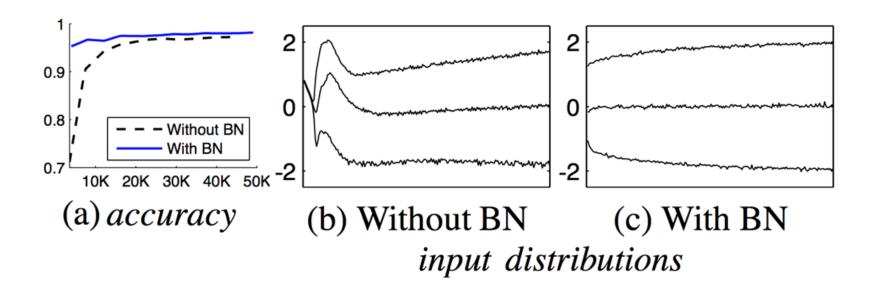
- For Training, the mean and variance of each batch are used for normalization
- For Testing, of which data the mean and variance will be used?
 - Estimated with those of batches in the training



Advantage

- Reduces internal covariant shift.
- Reduces the dependence of gradients on the scale of the connection weights.
- Regularizes the model and reduces the need for regularization techniques.
 - It adds some stochastic noise to the activations as a result of using noisy estimates computed on the mini-batches. This has a regularization effect in some applications,

Performance with BN



Disadvantage

- Expensive: Memory and time
 - Must keep interim results of all instances in a batch
 - Especially in CNN, usually an image is large
- Hard to apply when the batch size is small
 - If batches are small, the means and variances cannot approximate the global ones.
- Hard to apply to recurrent networks
 - It doesn't match to structure of recurrent networks
 - Hard to implement with recurrent networks

Recap: Batch Normalization

Normalization of each node output

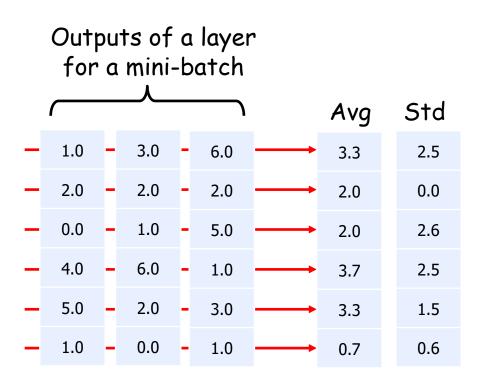
Batch normalization

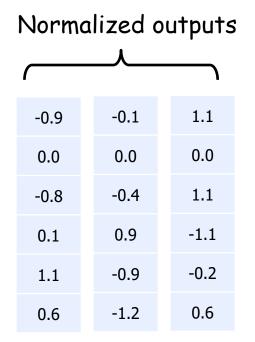
$$\mu_{j} = \frac{1}{m} \sum_{i=1}^{m} x_{ij}
\sigma_{j}^{2} = \frac{1}{m} \sum_{i=1}^{m} (x_{ij} - \mu_{j})^{2}
\hat{x}_{ij} = \frac{x_{ij} - \mu_{j}}{\sqrt{\sigma_{j}^{2} + \epsilon}}$$

i, j: index of the batch and the node of hidden layers

Recap: Batch Normalization

Normalization of each node output





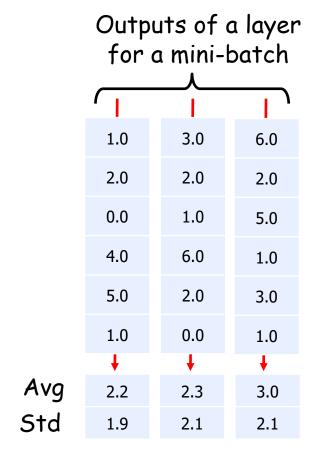
Proposed as an alternative to Batch Normalization

- Works regardless of batch size (batch size = 1)
- Performs well with RNNs

Layer normalization:

$$\mu_{i} = \frac{1}{m} \sum_{j=1}^{m} x_{ij}
\sigma_{i}^{2} = \frac{1}{m} \sum_{j=1}^{m} (x_{ij} - \mu_{i})^{2}
\hat{x}_{ij} = \frac{x_{ij} - \mu_{i}}{\sqrt{\sigma_{i}^{2} + \epsilon}}$$

i, j: index of the batch and the node of hidden layers





Group Normalization shows consistent accuracy with smaller batches

Tested on ImageNet (1000 Classes, 1.28M training, 50K validation), ResNet-50

