

# **Deep Learning**

#### **Content**

- Vanishing Gradient & Activation Functions
- Dropout
- Batch Normalization

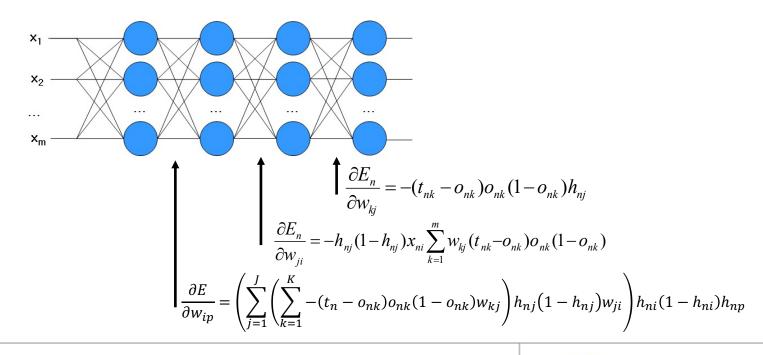


# **Gradient Vanishing & Activation Functions**

# **Gradient Vanishing & Exploding**

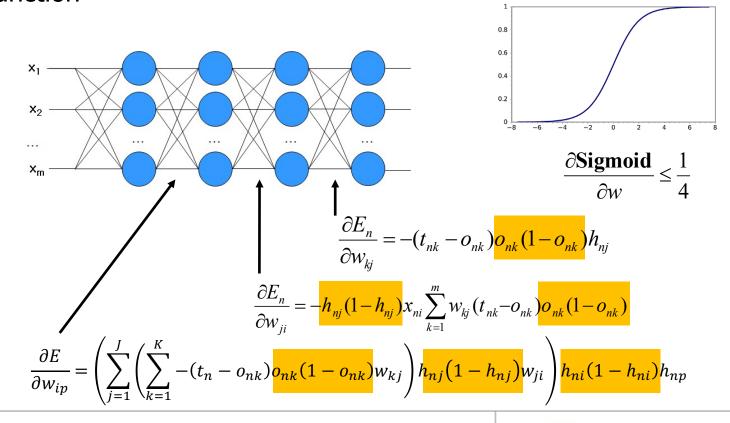
### Gradient is easy to vanish or explode

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



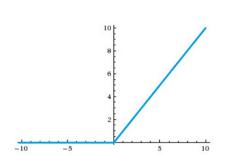
### Vanishing Gradient

The major terms are the derivatives of the activation function

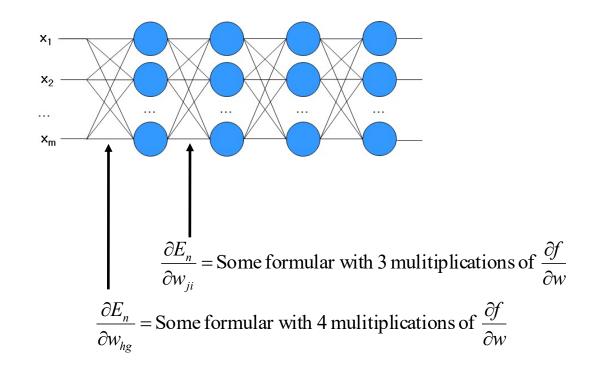


### Using another functions instead of sigmoid

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

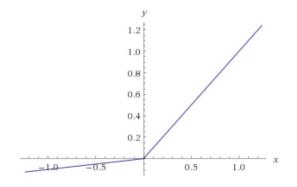
### Disadvantage

Knockout Problem

### You may use another

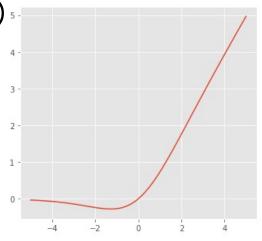
Leaky ReLU

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

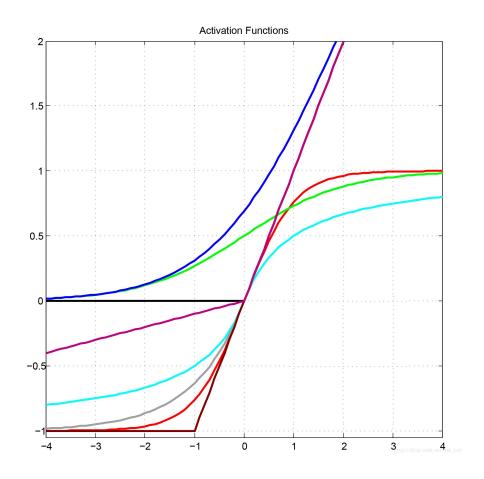


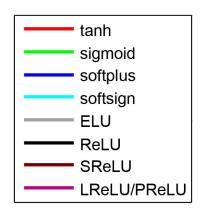
Swish (or SiLU-Sigmoid Linear Unit)

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



### **Other Activation Functions**





#### Summary

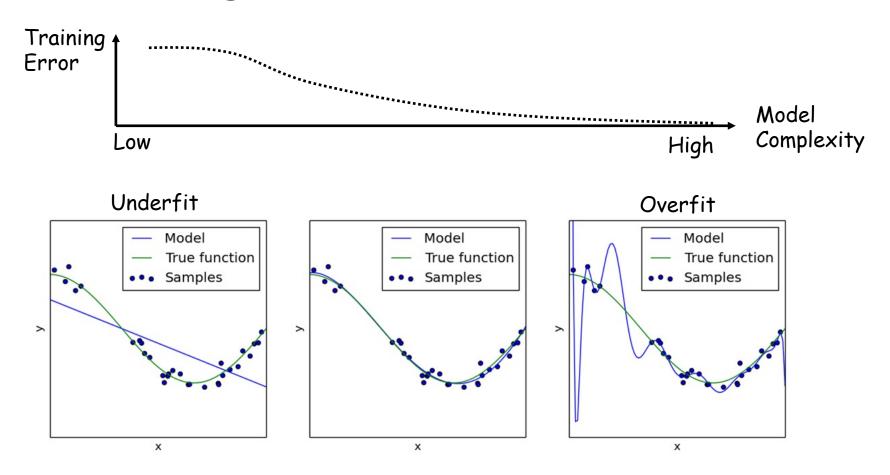
- Sigmoid functions and their combinations generally work better but are sometimes avoided due to the vanishing gradient problem
- ReLU function is a general activation function and is used in most cases these days
- If we encounter a case of dead neurons in our networks the leaky ReLU function is the 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

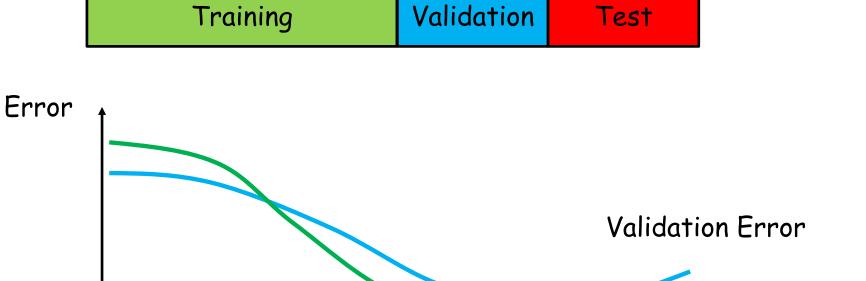
Proper Learning: Early stopping

Proper Structure: Weight decay, Dropout,

DropConnect, Stochastic pooling

# **Early Stopping**

### Split data into 3 groups



# of updates

Training Error

### L1 Regularization

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

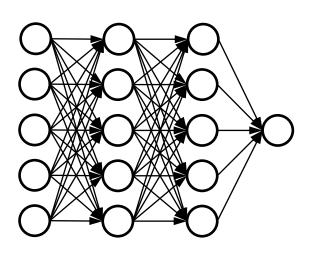
$$\widetilde{E}(\mathbf{w}) = E(\mathbf{w}) + \frac{\lambda}{2} |\mathbf{w}|$$

### L2 Regularization

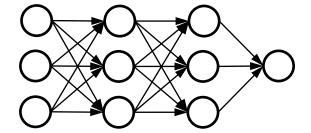
- Penalizing peaky weights
- Encouraging to use all of its inputs a little rather than using only some of its inputs a lot.

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

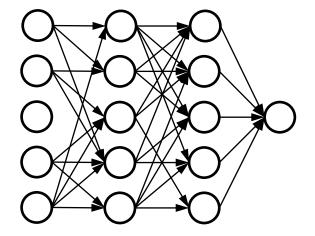
### Complex Structure vs Simple Structure



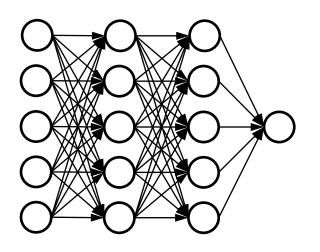
Node Pruning



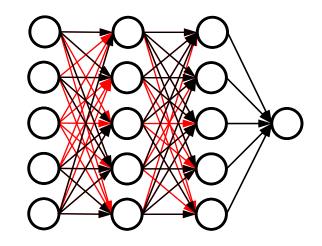
Link Pruning



### Complex Structure vs Simple Structure



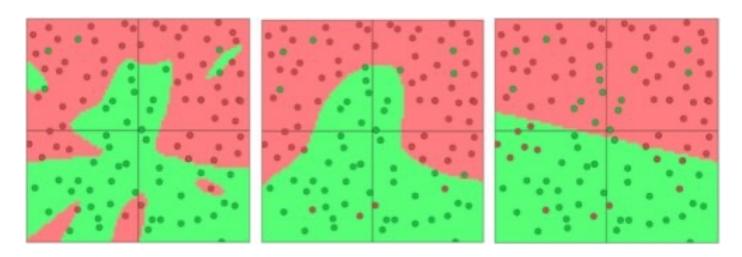
Set many links to zero



|w| is large <-> NN is Complex

|w| is small  $\leftarrow$  NN is Simple

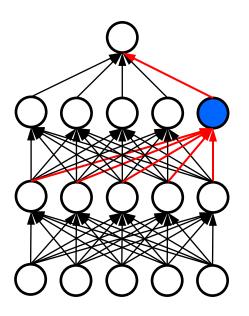
Example: Separating green and red



L2 regularization strengths of 0.01, 0.1, and 1

### In a complex Neural Network

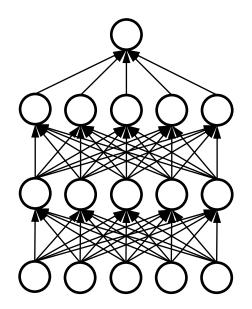
- All nodes do not take the same amount of responsibility
  - While training, some nodes are correlated
- All nodes are not equally trained
  - Some nodes are trained much, but some are not

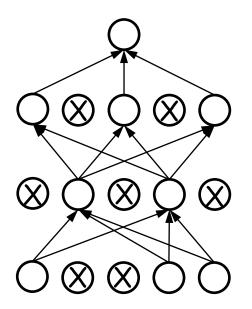


- If the output of the node is bad, the connection weight will decrease.
- If connection weight is close to 0, precedent connection weights are hardly trained.

### How can we reduce the structural complexity?

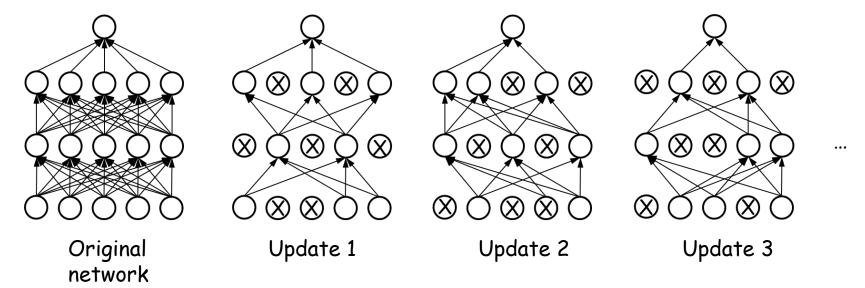
- Let's simply remove some nodes, and
- Train the simplified neural network
- Hmm??





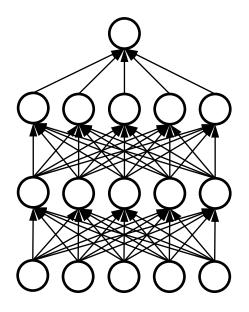
### Do this at every epoch

- 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

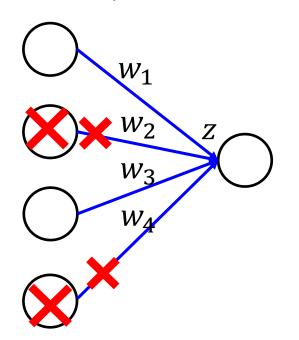
Use all the nodes without dropout



### Testing

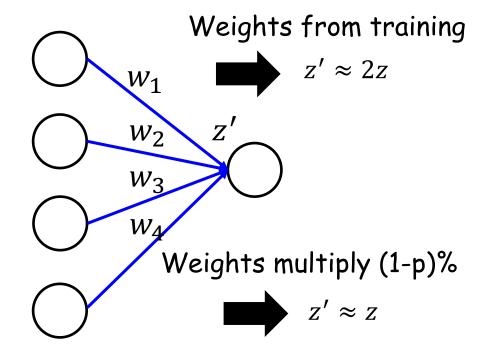
#### Training of Dropout

Assume dropout rate is 50%



### Testing of Dropout

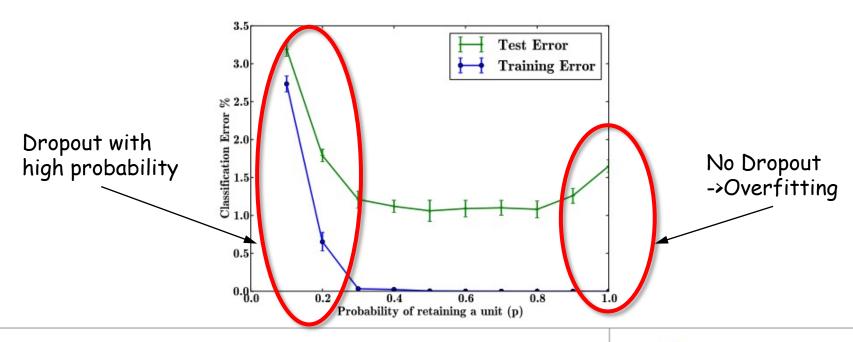
No dropout





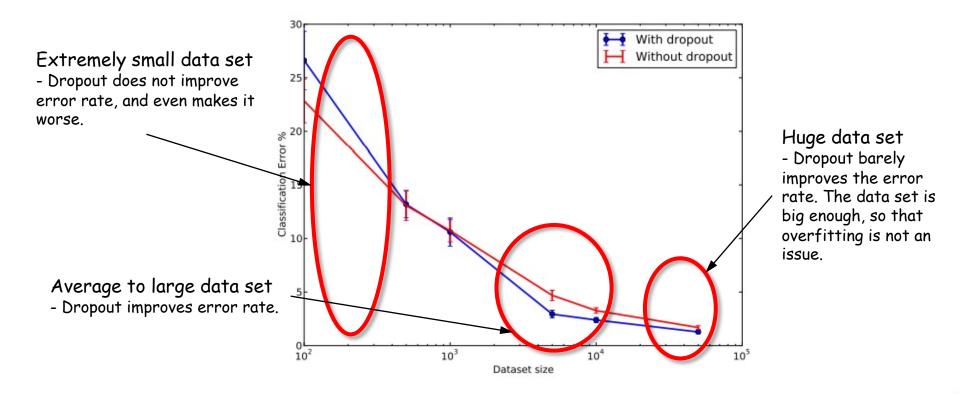
### The effect of the dropout rate p:

- 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.



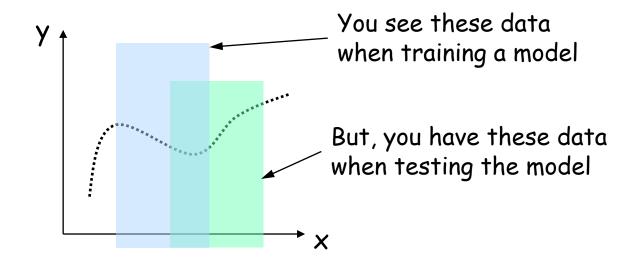
#### Summary

- 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).



#### Covariate Shift

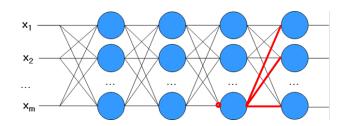
A change in the distribution of a function's domain.

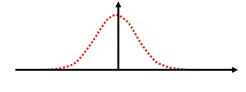


– Can your model work properly?

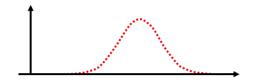
#### Internal Covariate Shift

Input distribution of the red node





- While learning, red connection weights will change based on the input distribution
- After learning, the whole connection weights changes, which cause the change of the input distribution



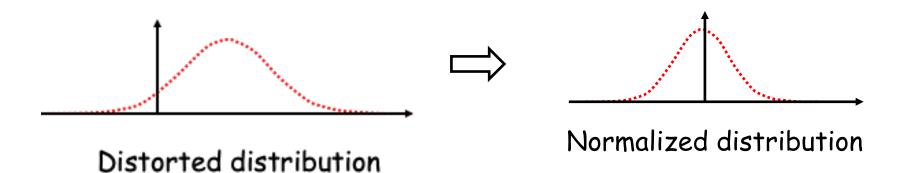
The assumption of the learning is broken

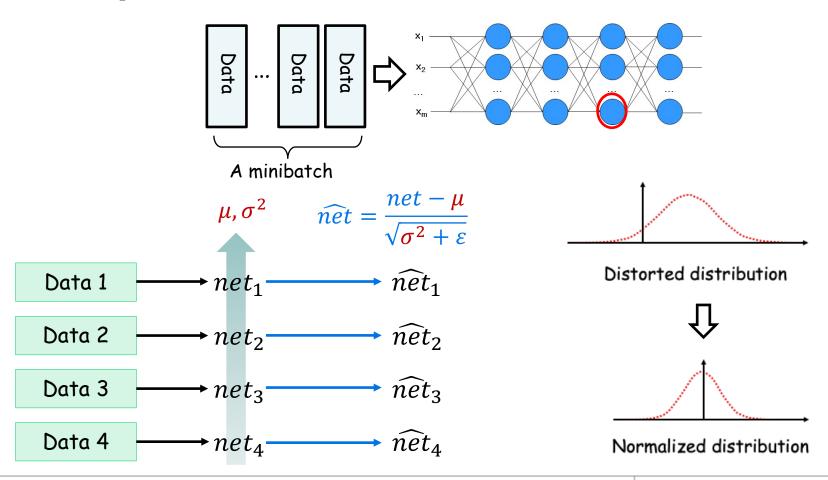
#### Internal Covariate Shift

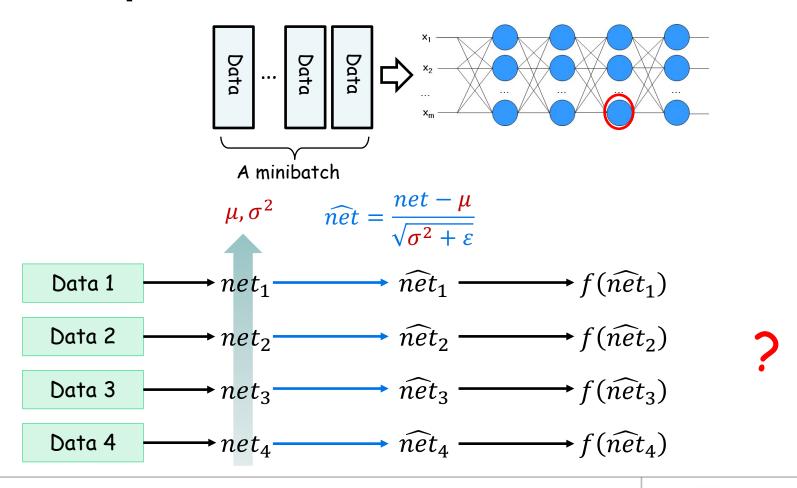
- It disturbs the learning process,
- Learning is getting slow down

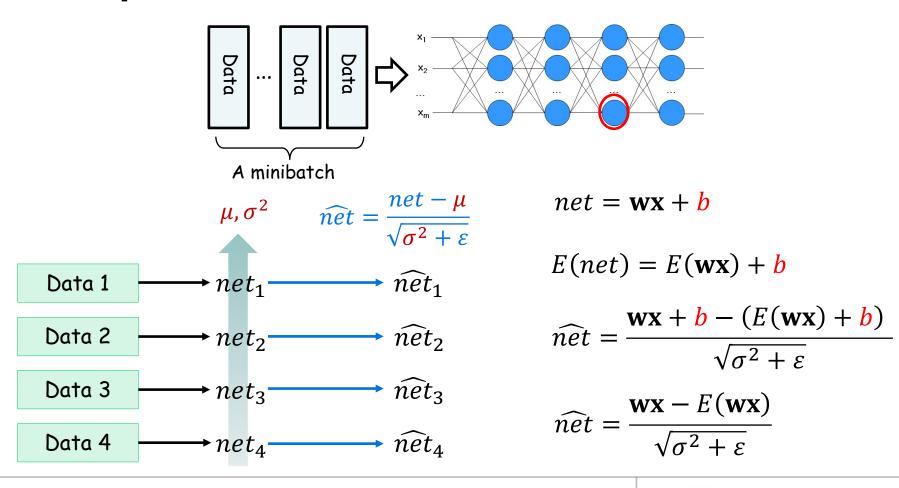
#### What shall we do?

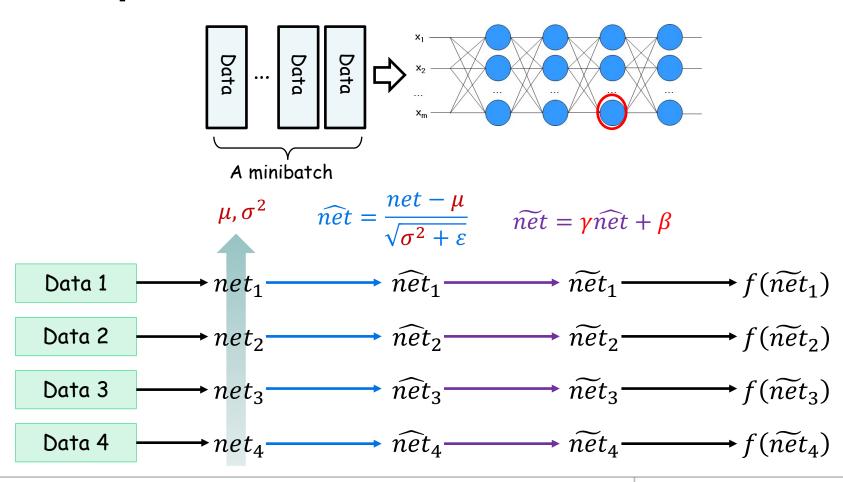
Why don't we normalize the distribution of inputs

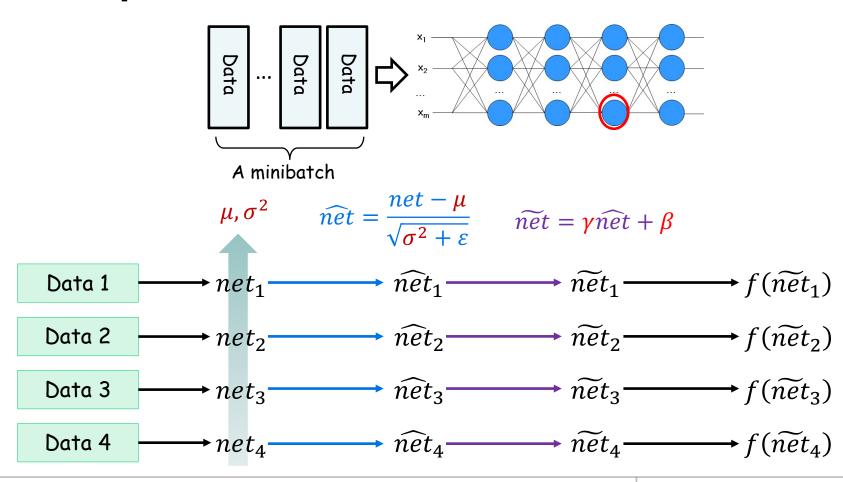




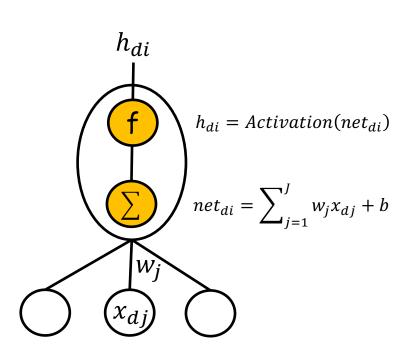


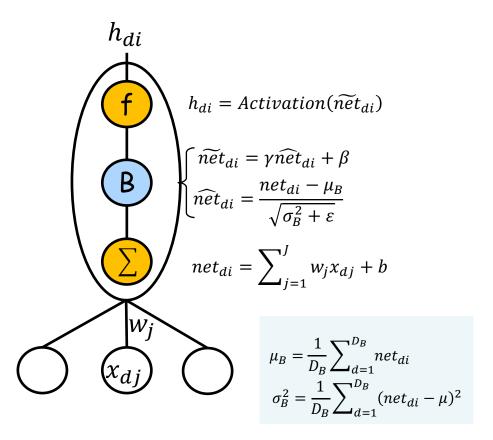






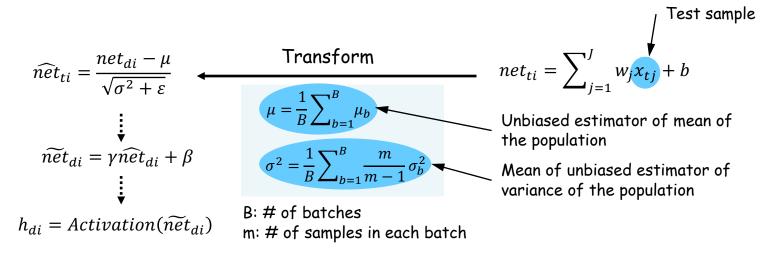
### For a Single Node





#### Testing

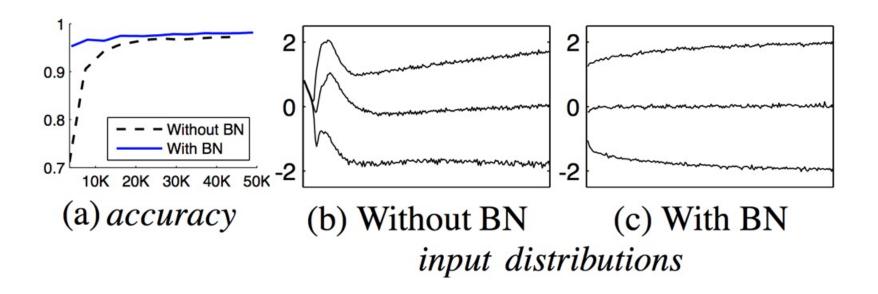
- For Training, the mean and variance of each batch are used for normalization
- For Testing, of which data the mean and variance will be used?
  - Estimated with those of batches in the training



### 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

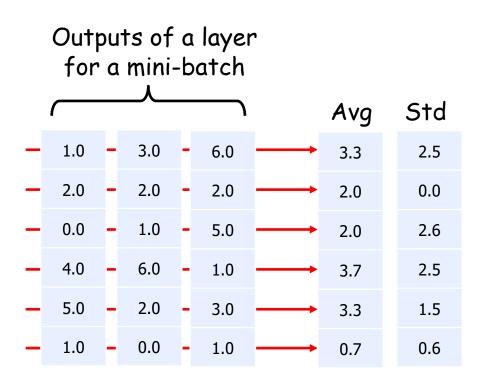
#### Batch normalization

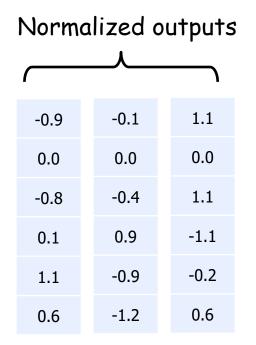
$$\mu_{j} = \frac{1}{m} \sum_{i=1}^{m} x_{ij} 
\sigma_{j}^{2} = \frac{1}{m} \sum_{i=1}^{m} (x_{ij} - \mu_{j})^{2} 
\hat{x}_{ij} = \frac{x_{ij} - \mu_{j}}{\sqrt{\sigma_{j}^{2} + \epsilon}}$$

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

### Recap: Batch Normalization

Normalization of each node output





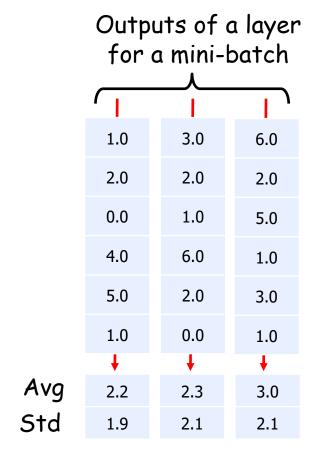
### Proposed as an alternative to Batch Normalization

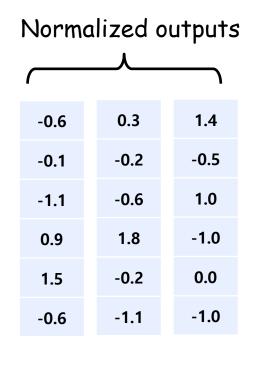
- Works regardless of batch size (batch size = 1)
- Performs well with RNNs

#### 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

