

Network Regularization

Weight initialization, dropout, batch normalization

Machine Learning and Data Mining, 2023

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National Research University Higher School of Economics



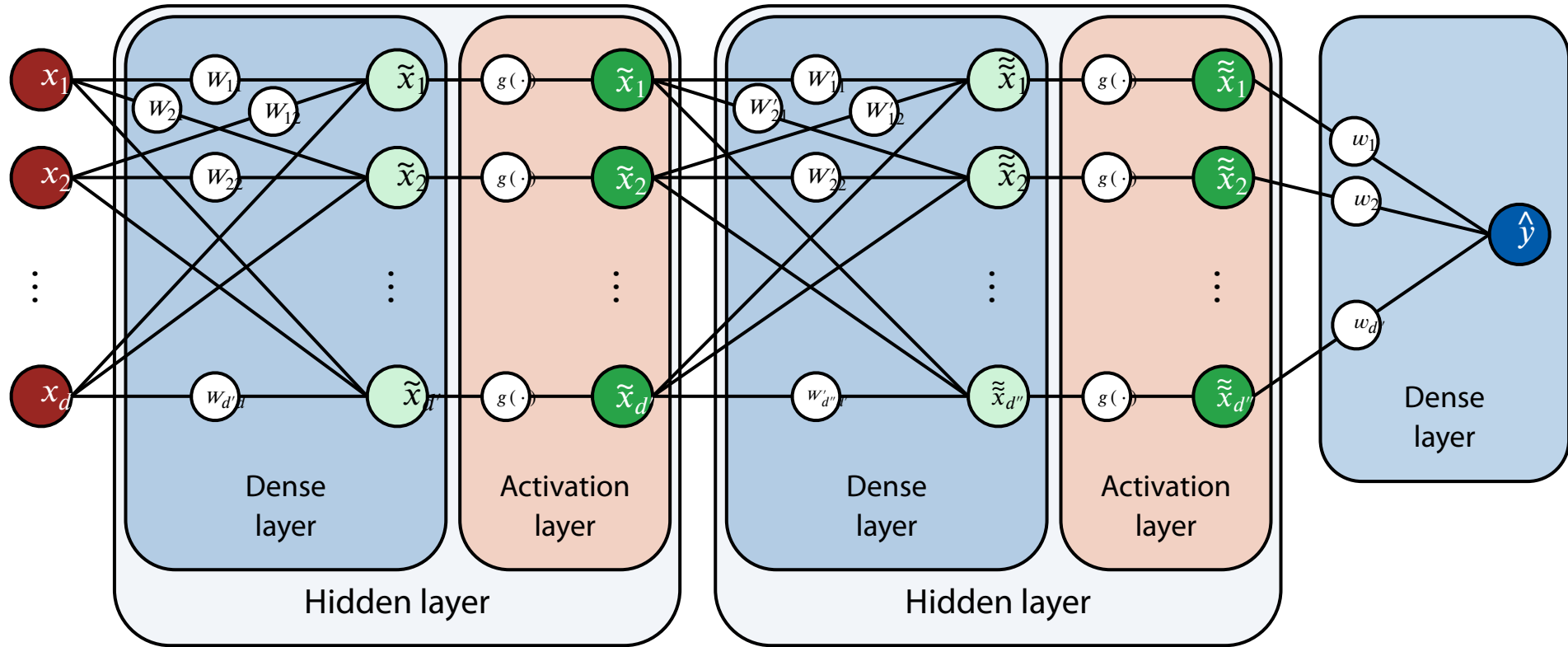
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Why care about weight initialization?

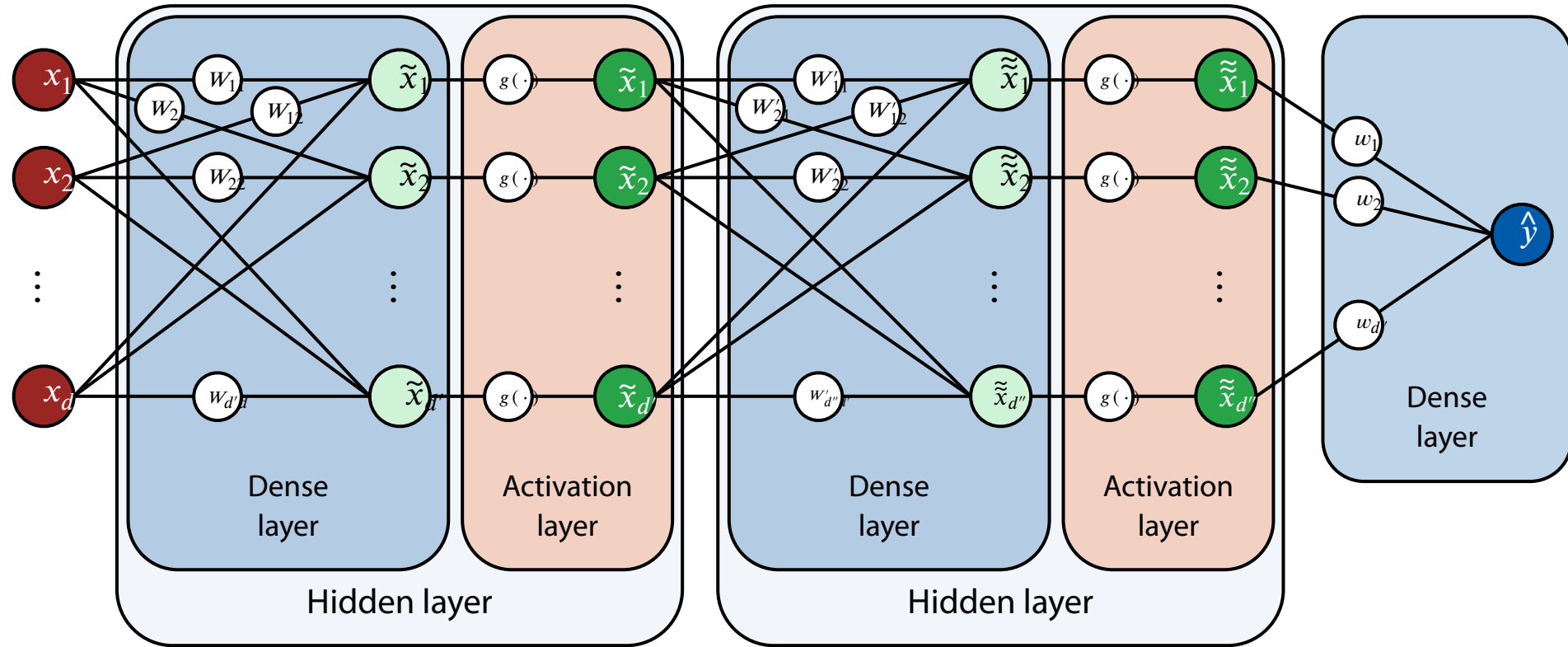


Initialization with a constant (?)



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- What happens if we initialize all weights with the same value?
- Within each layer, the gradients for each of the weights will be the same as well \Rightarrow updates will be the same \Rightarrow network degrades!

Initialization with a constant (?)

- ▶ Ok, so constant initialization is a bad idea
- ▶ So, any random initialization should be fine, right?

Some intuition

- ▶ For simplicity, let's omit the activation functions for now
- ▶ Then, the output of a neural network composed of dense layers only is:

$$\hat{y} = W_{out} \cdot \dots \cdot W_{h2} \cdot W_{h1}x$$

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$$g \sim S^{m-1}$$

where m is the **depth** of the network

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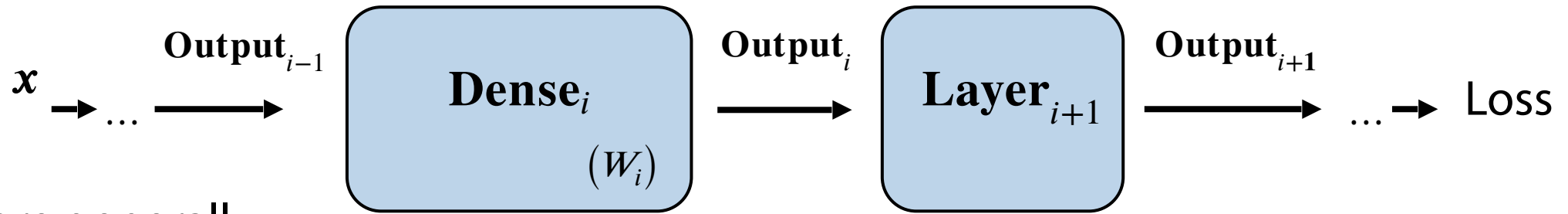
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- ▶ For S too large, the gradients will **explode**; for S too small, they will **vanish**

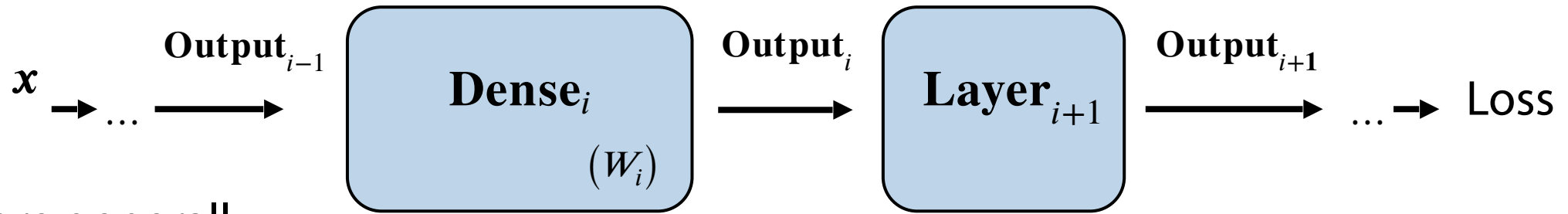
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► More generally:

$$\frac{\partial \text{Loss}}{\partial W_i} = \frac{\partial \text{Loss}}{\partial \text{Output}_i} \cdot \frac{\partial \text{Dense}_i}{\partial W_i} = \frac{\partial \text{Loss}}{\partial \text{Output}_{i+1}} \cdot \frac{\partial \text{Layer}_{i+1}}{\partial \text{Output}_i} \cdot \text{Output}_{i-1}$$

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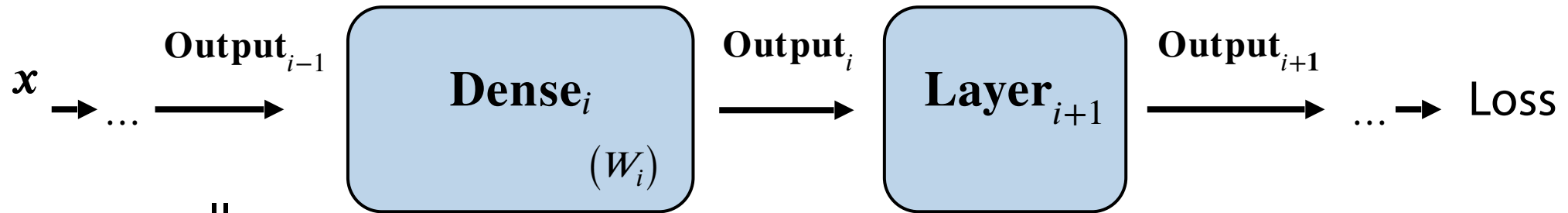


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This will accumulate the product of the gradients
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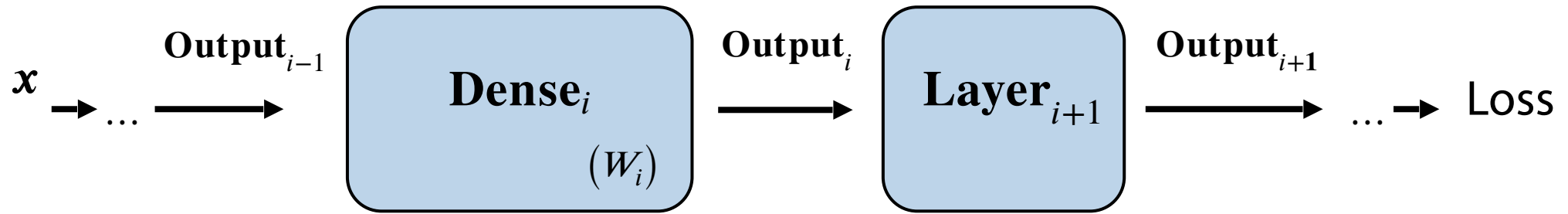
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- Idea: for stable learning we would like to “keep” the scale of the gradients at each step:

$$\text{Var}\left(\frac{\partial \text{Layer}_{i+1}}{\partial \text{Output}_i} \cdot \frac{\partial \text{Layer}_i}{\partial \text{Output}_{i-1}}\right) \approx \text{Var}\left(\frac{\partial \text{Layer}_{i+1}}{\partial \text{Output}_i}\right)$$

Some intuition



- ▶ Similarly, we would also like to not scale the outputs at each step of the forward pass:

$$\text{Var}\left(\text{Layer}_{i+1}\left(\text{Layer}_i\left(\text{Output}_{i-1}\right)\right)\right) \approx \text{Var}\left(\text{Layer}_i\left(\text{Output}_{i-1}\right)\right)$$

Random initialization

$$\begin{aligned} \text{Var}\left(\frac{\partial \mathbf{Layer}_{i+1}}{\partial \mathbf{Output}_i} \cdot \frac{\partial \mathbf{Layer}_i}{\partial \mathbf{Output}_{i-1}}\right) &\approx \text{Var}\left(\frac{\partial \mathbf{Layer}_{i+1}}{\partial \mathbf{Output}_i}\right) \\ \text{Var}\left(\mathbf{Layer}_{i+1}\left(\mathbf{Layer}_i\left(\mathbf{Output}_{i-1}\right)\right)\right) &\approx \text{Var}\left(\mathbf{Layer}_i\left(\mathbf{Output}_{i-1}\right)\right) \end{aligned}$$

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- ▶ Generally, these two requirements may contradict each other
- ▶ E.g. for ReLU activation they result in initialization requirements, respectively:

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- ▶ Typically you can just choose one of them, or alternatively average them out:

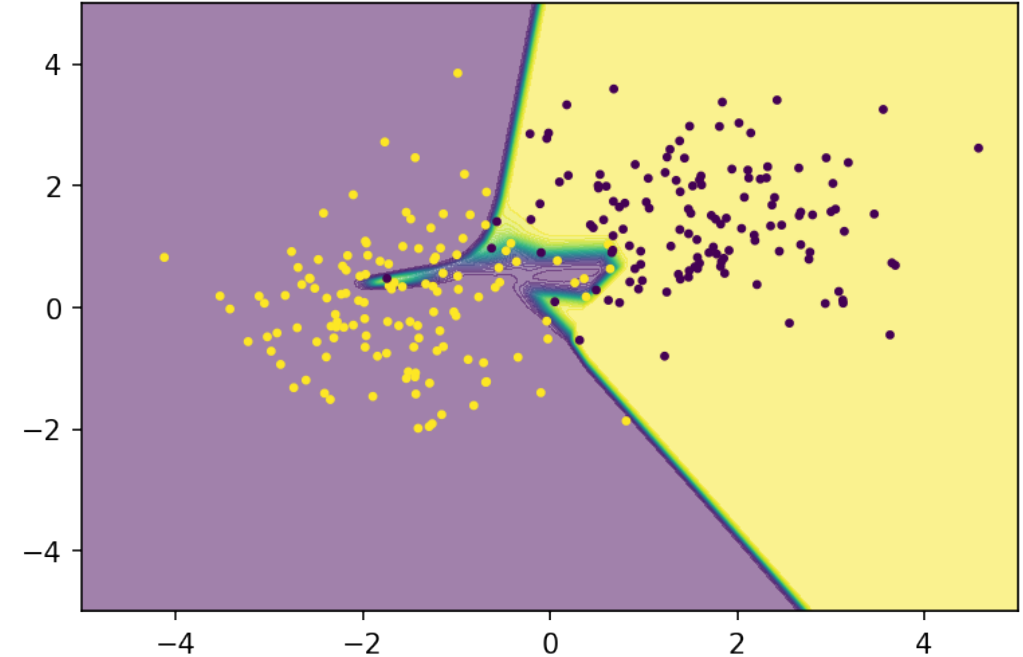
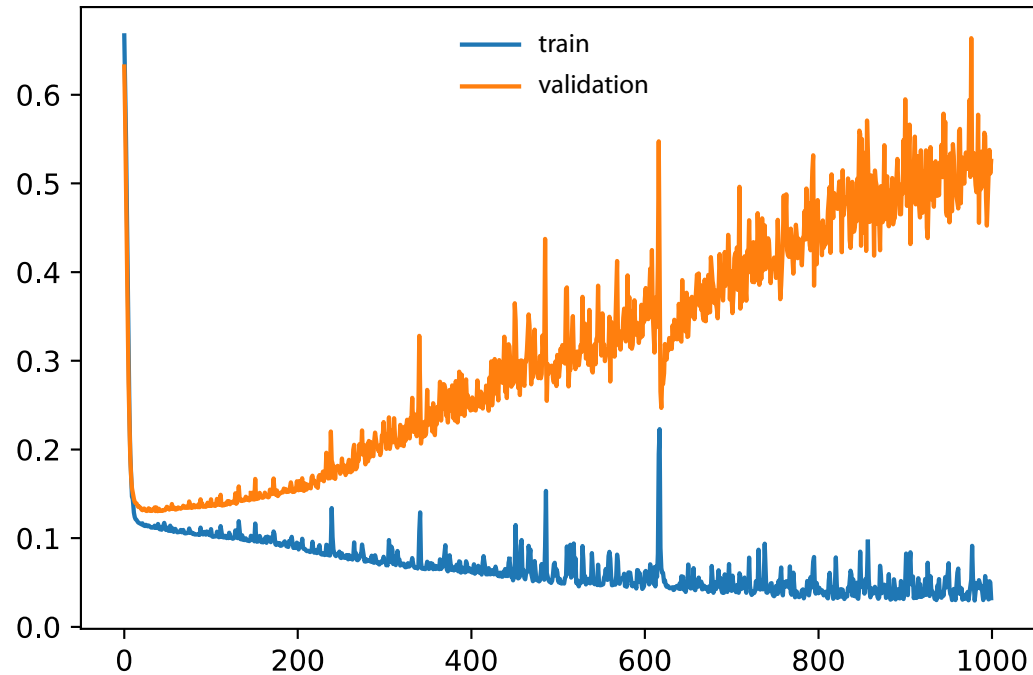
$$\text{Var}\left(W_{ij}\right) = \frac{4}{(\# \text{ outgoing connections}) + (\# \text{ incoming connections})}$$

Overfitting with neural networks



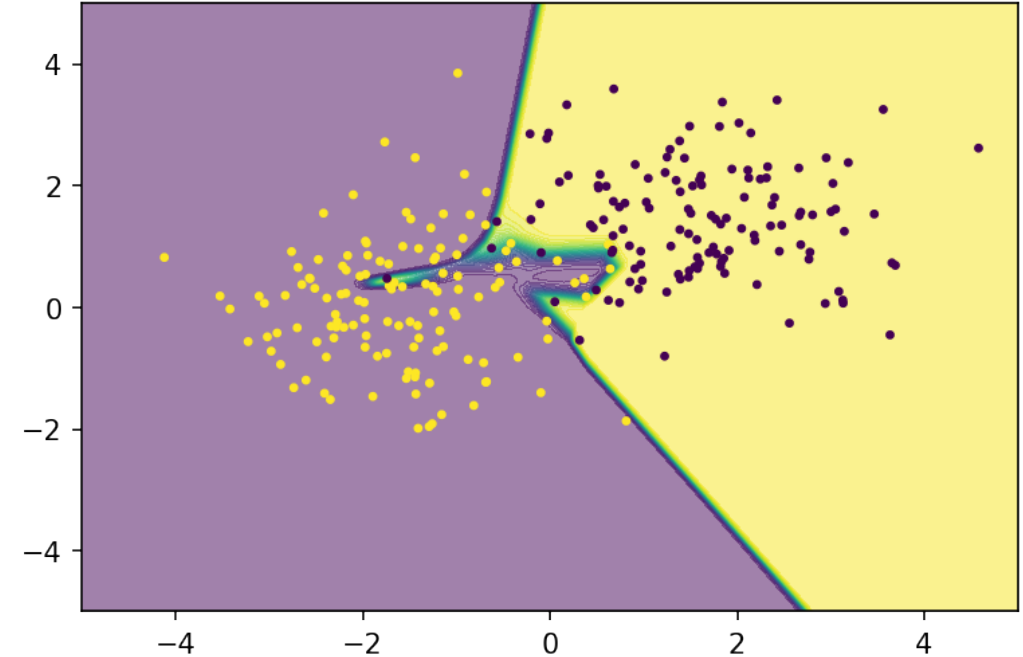
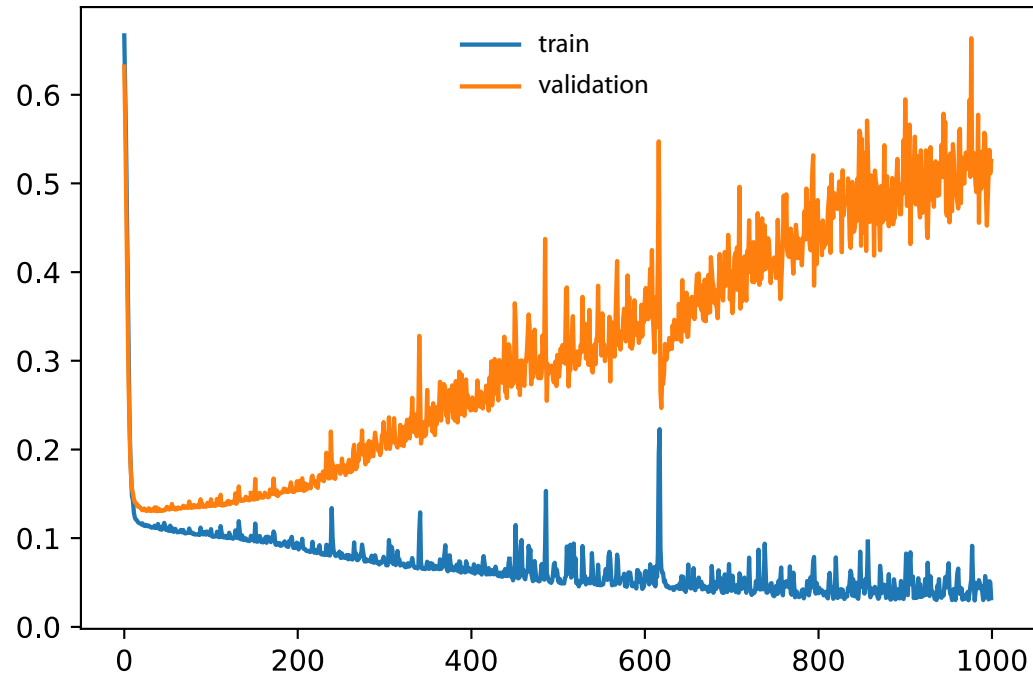
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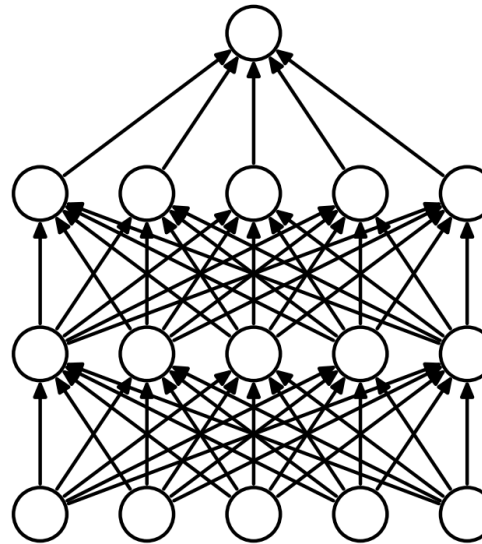


- ▶ Regularization techniques like L1/L2 regularization are also available for neural networks
- ▶ We also discussed **early stopping** (i.e. stop the training before validation error grows)

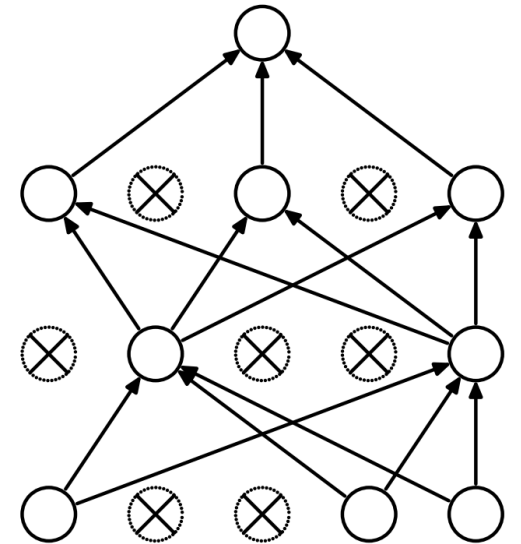
Dropout

- ▶ At train time – sets neuron activations to 0 with a given probability p

Image from:
<http://jmlr.org/papers/v15/srivastava14a.html>



(a) Standard Neural Net

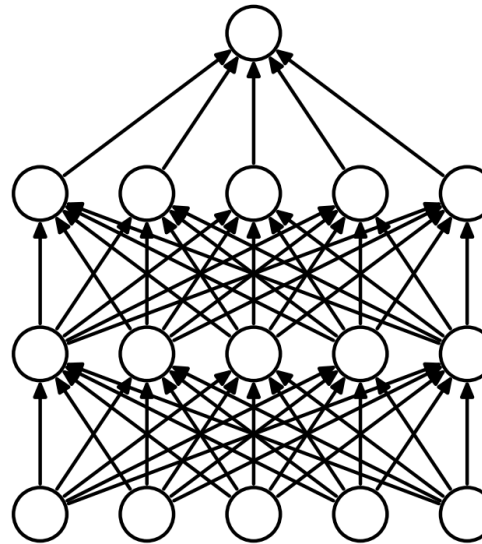


(b) After applying dropout.

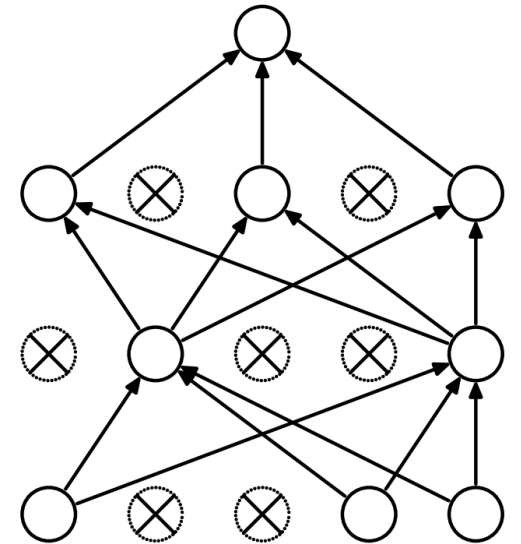
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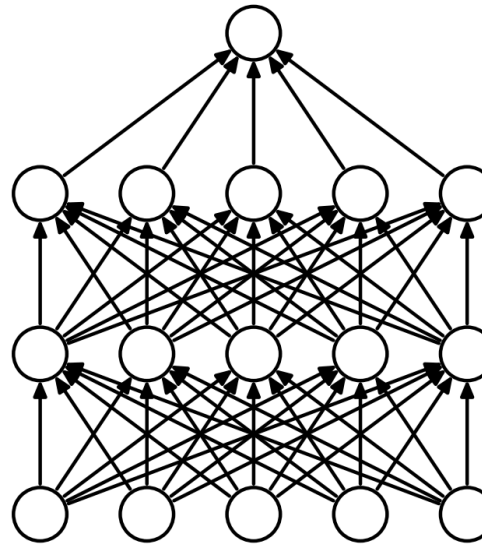


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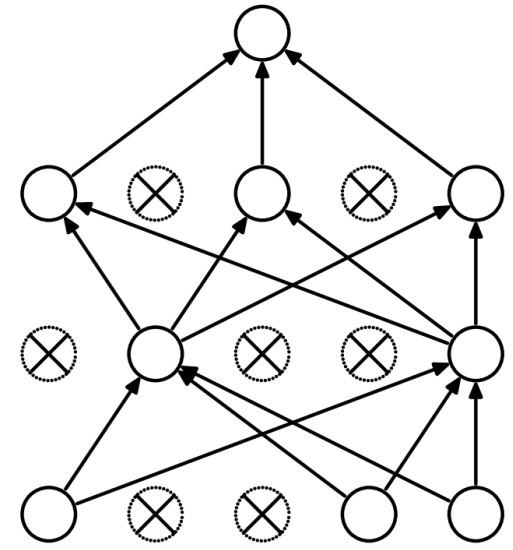
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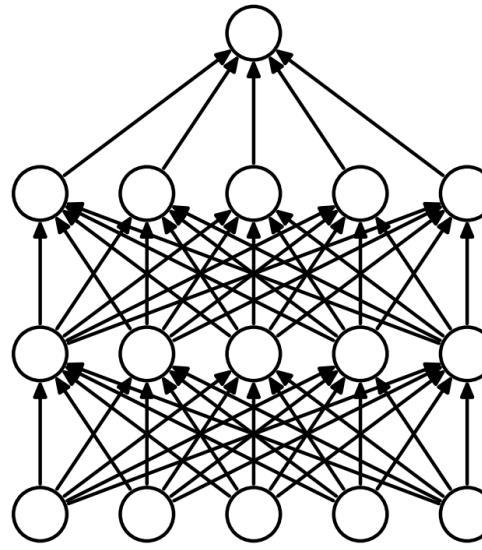


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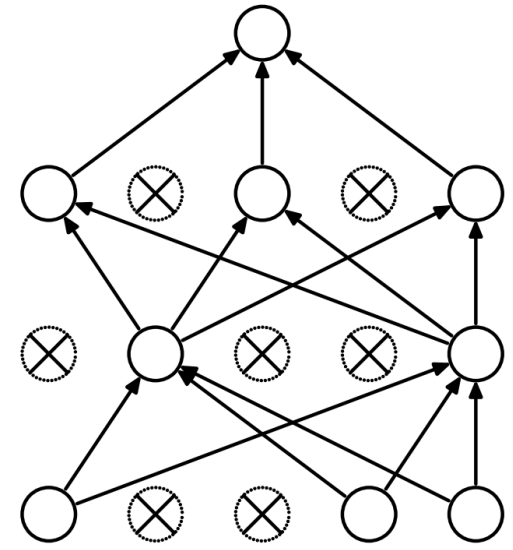
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- ▶ Drives it towards **creating useful features** rather than relying on other neurons to correct its mistakes

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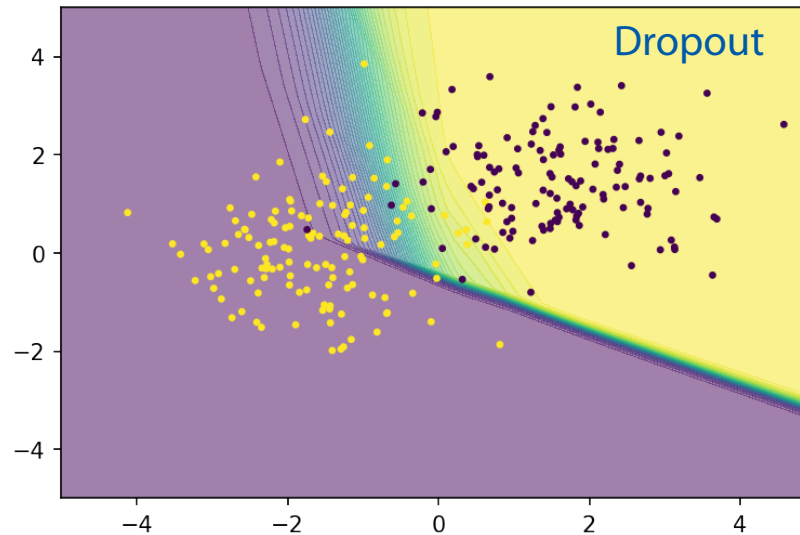
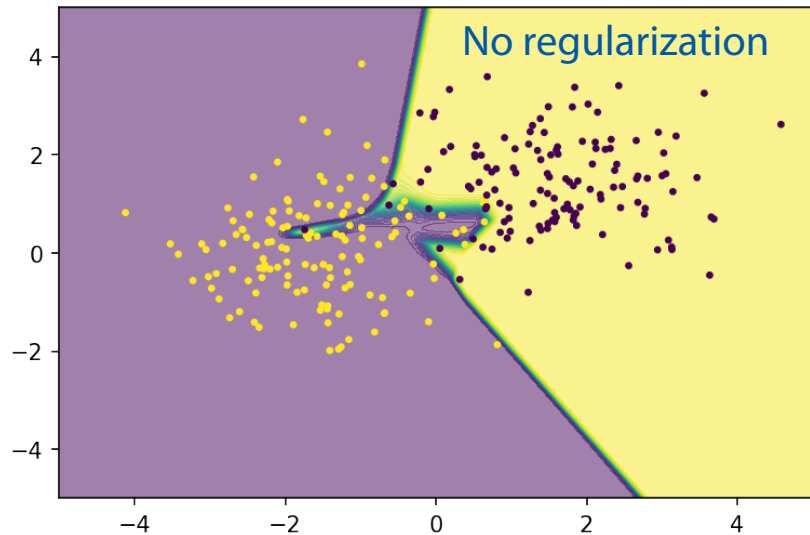
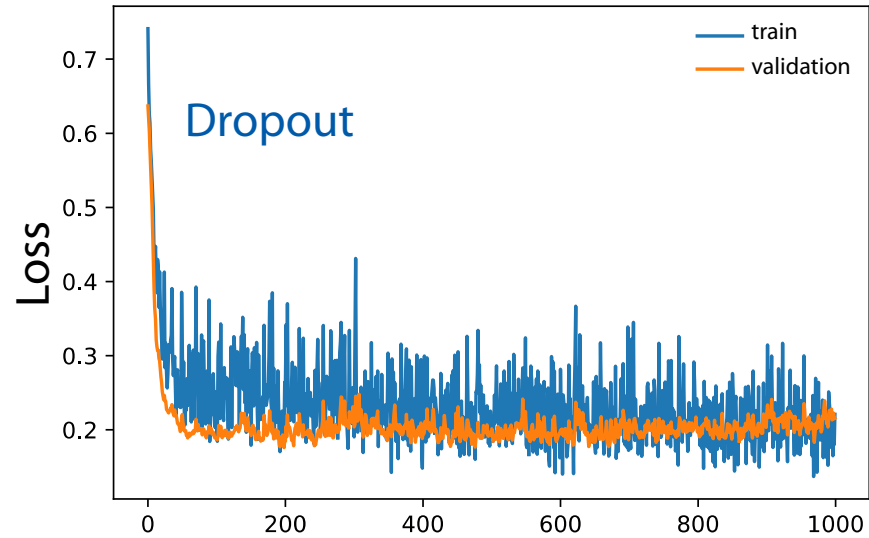
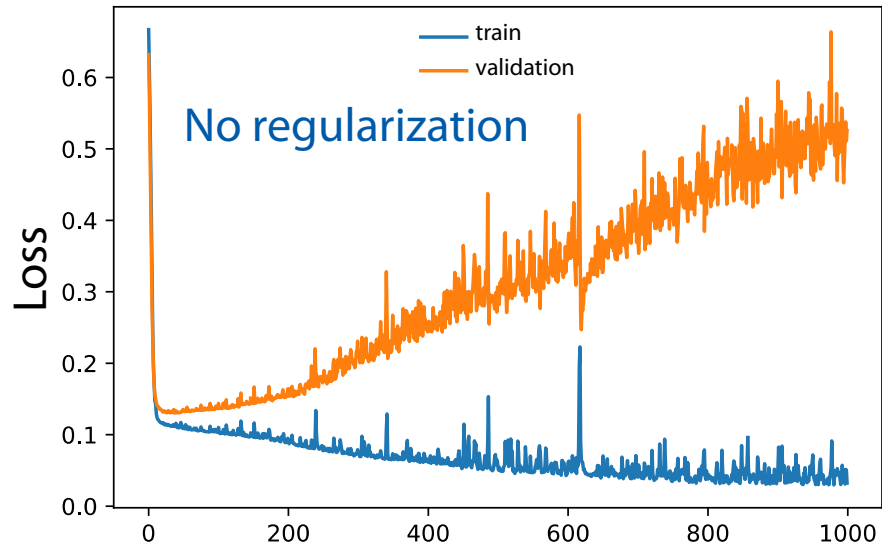


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Example from before



In this example, dropout results in a much better (though still not perfect) fit with lower test error

Normalization layers



Batch normalization

- ▶ This technique was originally proposed to mitigate the “internal covariate shift”

internal covariate shift

the updates in one layer change
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- ▶ Works as follows (layer inputs x_i , outputs y_i):
 - calculate sample **mean** and **variance** of the input on a single batch B

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$$\mu_B = \frac{1}{|B|} \sum_{i \in B} x_i \quad \sigma_B^2 = \frac{1}{|B|} \sum_{i \in B} (x_i - \mu_B)^2$$

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- **normalize** the input, then **scale and shift** (with the trainable parameters γ, β):

$$y_i = \gamma \cdot \frac{x_i - \mu_B}{\sqrt{\sigma_B^2 + \epsilon}} + \beta$$

Batch normalization

- ▶ Turned out to be **extremely powerful** in many cases
 - Faster and more stable convergence

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Batch normalization

- ▶ Turned out to be **extremely powerful** in many cases
 - Faster and more stable convergence
- ▶ Later was proved to not reduce the internal covariate shift
- ▶ Effectively **removes** the 'shift' and 'scale' degrees of freedom from the previous layer

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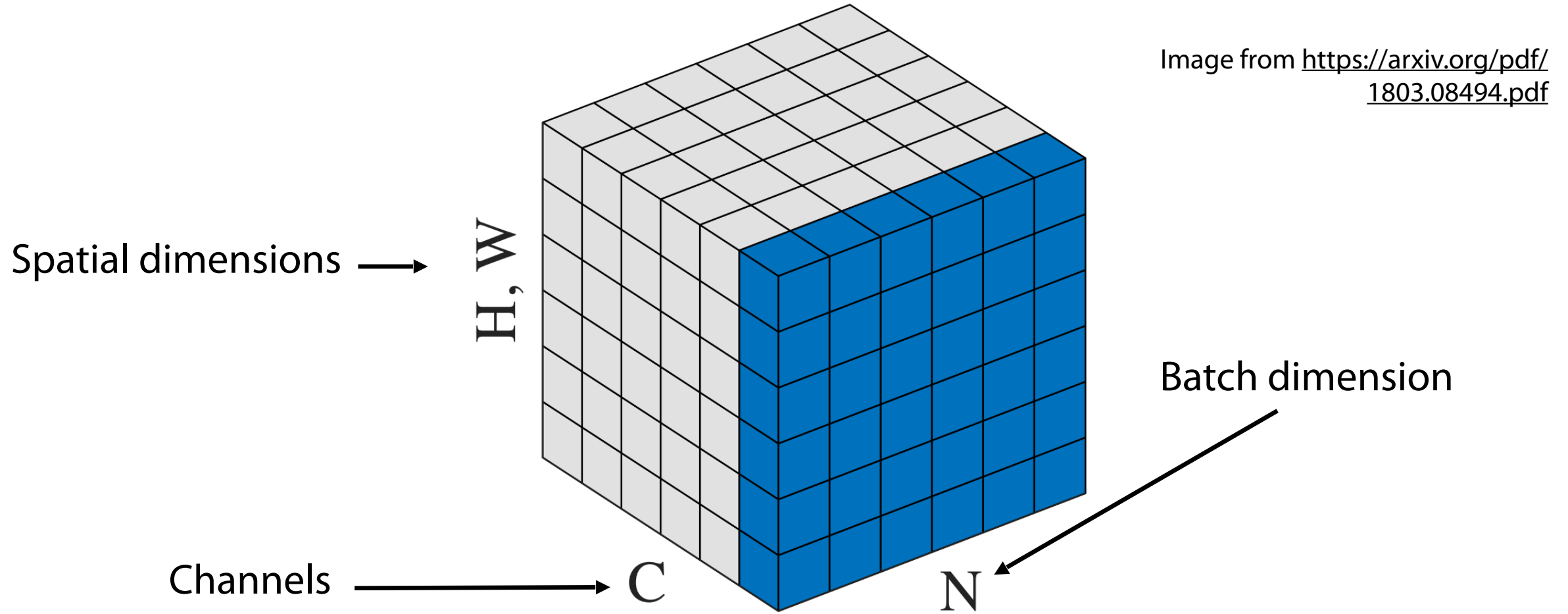
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 - separately for each component in Channel_dim , i.e., over Batch_dim x Spacial_dim1 x ...

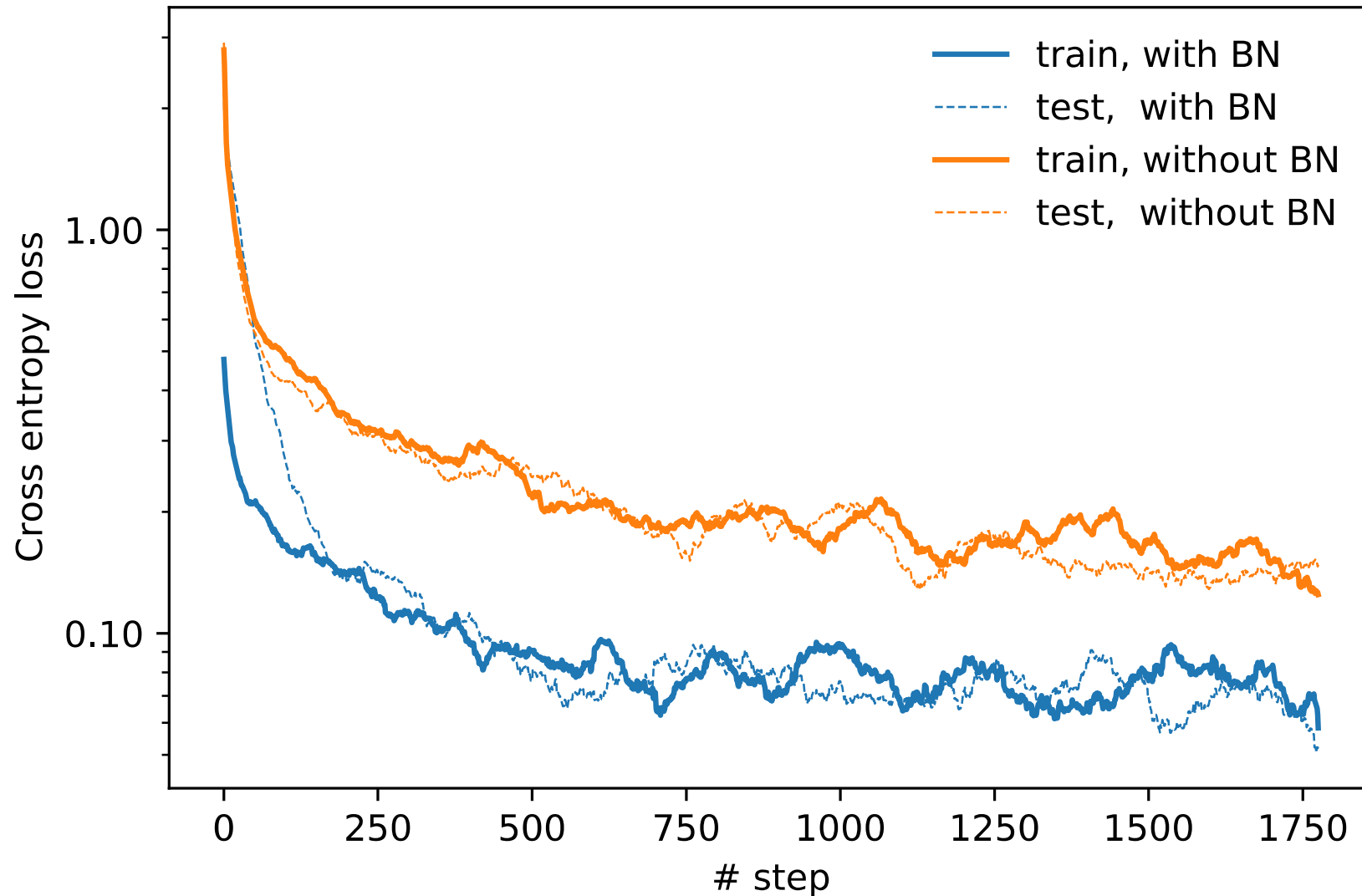
Batch normalization



Batch normalization at inference time

- ▶ Calculating batch statistics at test time may be problematic
 - e.g. when there's a single object to predict
- ▶ Instead: calculate running mean and variance during training, apply at test time

Example: CNN on MNIST



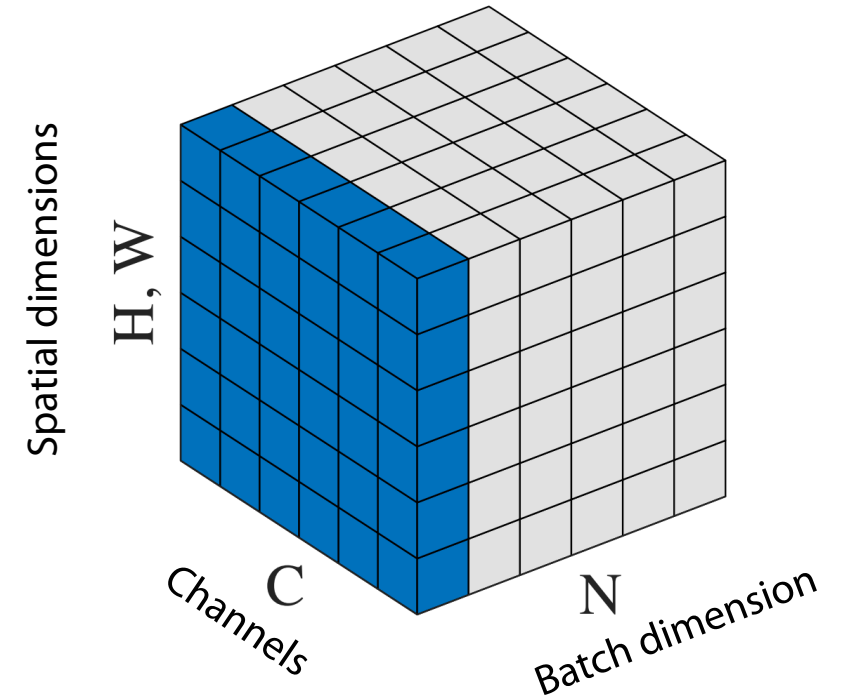
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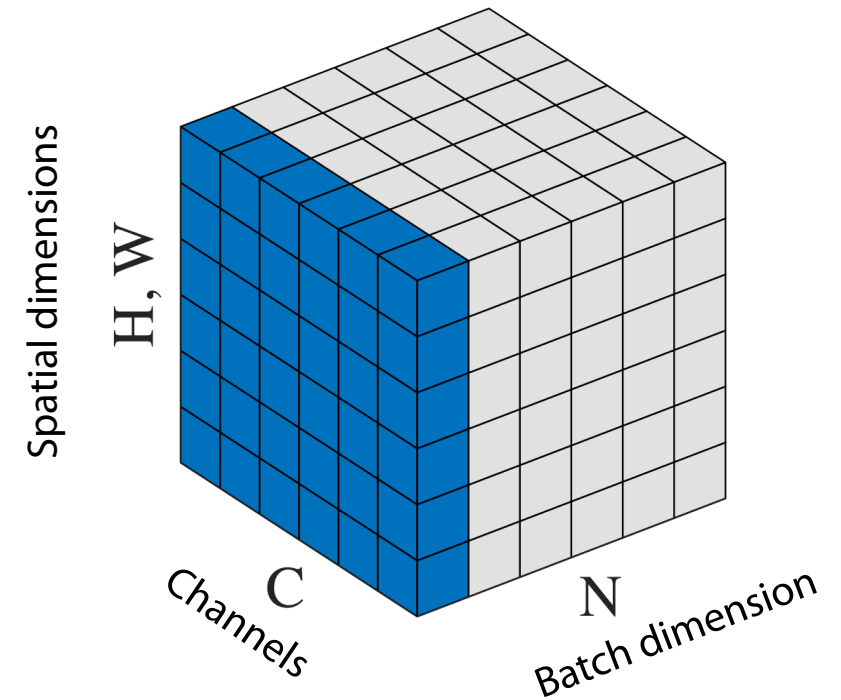
Image from <https://arxiv.org/pdf/1803.08494.pdf>



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- ▶ Alternative: **Layer Normalization**
 - the math is same, except statistics is calculated over channels rather than batch elements
 - the effect is quite different though
 - e.g. Layer Normalization “entangles” different neurons within a layer

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- ▶ Batch normalization is an **extremely powerful** regularization technique, though the reason for that is not entirely clear

- ▶ Food for thought: how exactly would you implement an early stopping rule?

Thank you!



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