Artem Maevskiy



Network Regularization

Weight initialization, dropout, batch normalization

2021









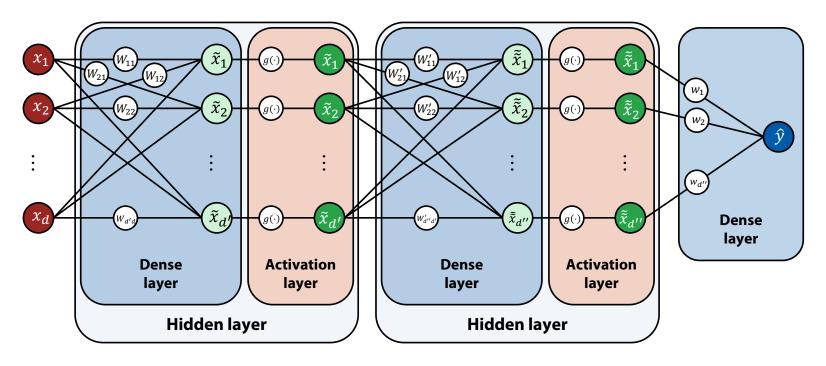






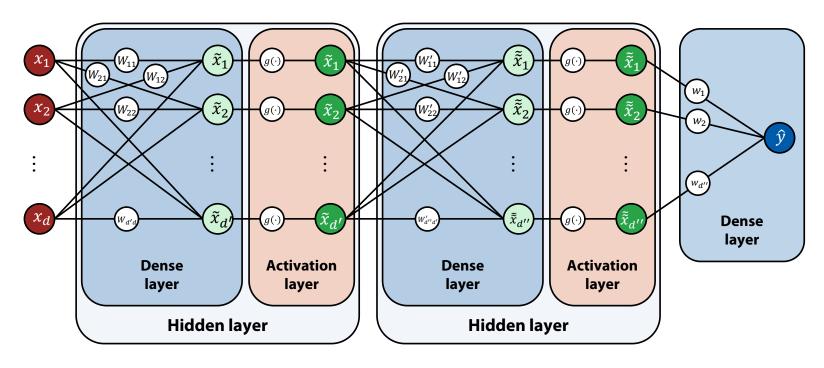
Why care about weight initialization?

Initialization with a constant (?)



What happens if we initialize all weights with the same value?

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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 ⇒ updates will be the same ⇒ network degrades!

Initialization with a constant (?)

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

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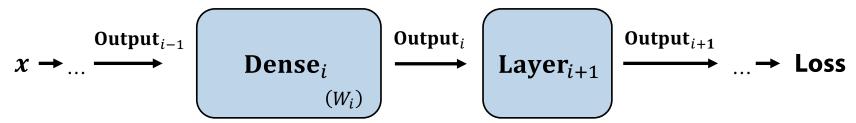
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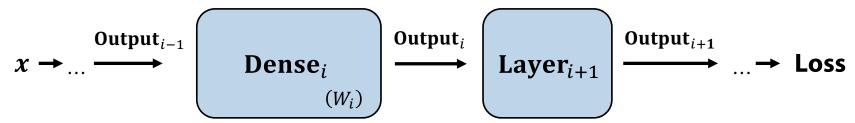
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► For *S* too large, the gradients will **explode**; for *S* too small, they will **vanish**



More generally:

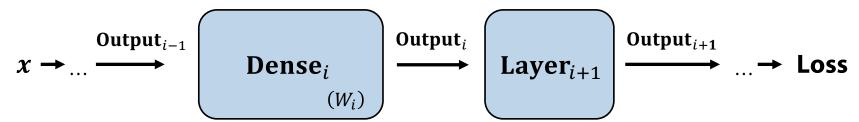
$$\frac{\partial \mathbf{Loss}}{\partial W_i} = \frac{\partial \mathbf{Loss}}{\partial \mathbf{Output}_i} \cdot \frac{\partial \mathbf{Dense}_i}{\partial W_i} = \frac{\partial \mathbf{Loss}}{\partial \mathbf{Output}_{i+1}} \cdot \frac{\partial \mathbf{Layer}_{i+1}}{\partial \mathbf{Output}_i} \cdot \mathbf{Output}_{i-1}$$



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This will accumulate the product of the gradients for the subsequent layers



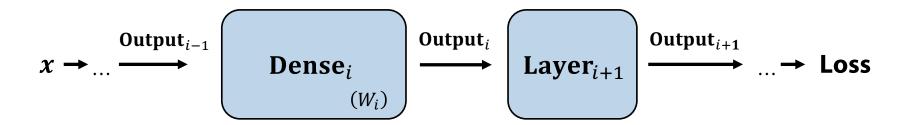
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Idea: for stable learning we would like to "keep" the scale of the gradients at each step:

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



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

$$Var\left(Layer_{i+1}\left(Layer_{i}\left(Output_{i-1}\right)\right)\right) \approx Var\left(Layer_{i}\left(Output_{i-1}\right)\right)$$

Random initialization

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

$$Var(W_{ij}) = \frac{2}{\text{(# outgoing connections)}}$$

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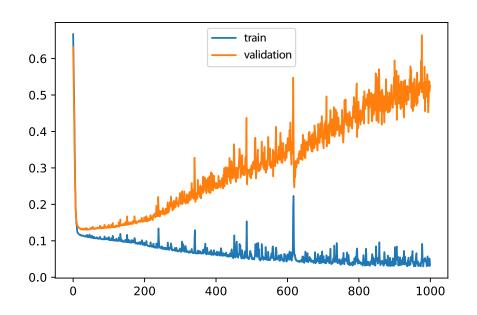
Typically you can just choose one of them, or alternatively average them out:

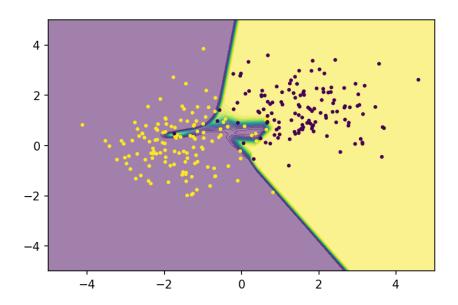
$$Var(W_{ij}) = \frac{4}{(\text{# outgoing connections}) + (\text{# incoming connections})}$$

Overfitting with neural networks

The problem of overfitting

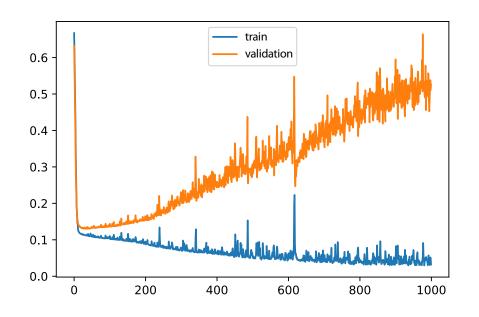
Being highly complex models, neural networks are prone to overfitting

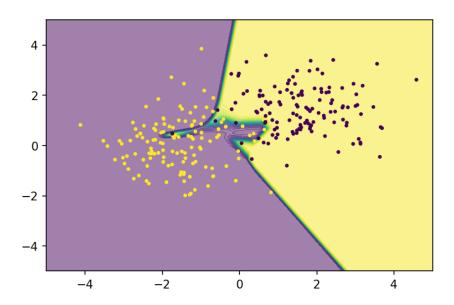




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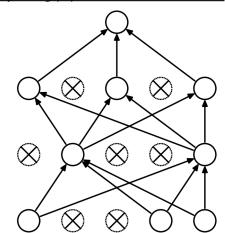
- ▶ Regularization techniques like L1/L2 regularization are also available for neural networks
- We also discussed early stopping (i.e. stop the training before validation error grows)

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

(a) Standard Neural Net

(b) After applying dropout.

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



- At train time sets neuron activations to 0 with a given probability p
- At test time multiplies the activation
 by p
 - i.e. sets it to the expected value

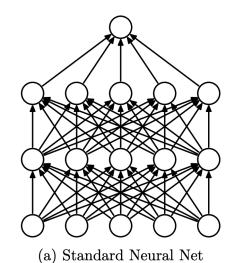
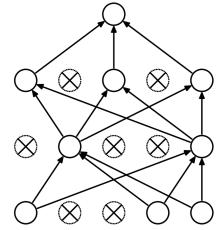


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(b) After applying dropout.

- At train time sets neuron activations to 0 with a given probability p
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- Makes neuron learn to work with a randomly chosen sample of other neurons

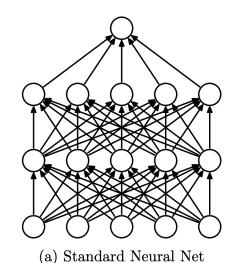
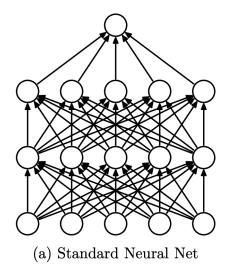


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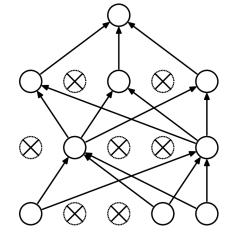
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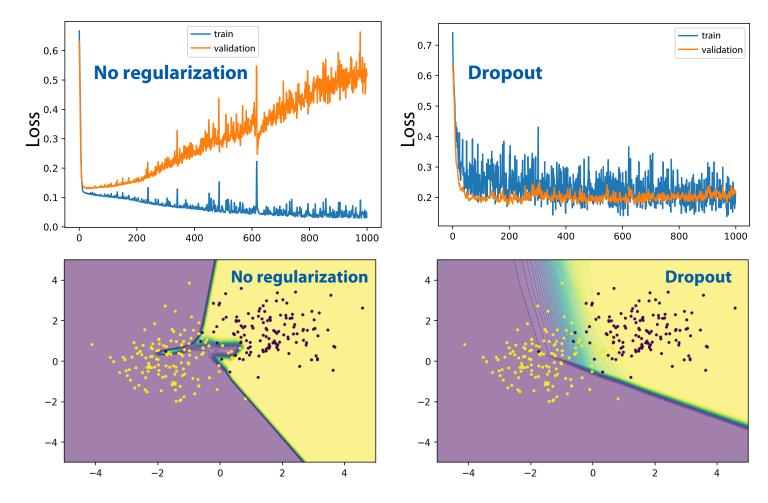
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(b) After applying dropout.

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

Example from before



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

Normalization layers

This technique was originally proposed to mitigate the "internal covariate shift"

internal covariate shift

the updates in one layer change the input distributions of the subsequent layers

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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

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

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

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

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

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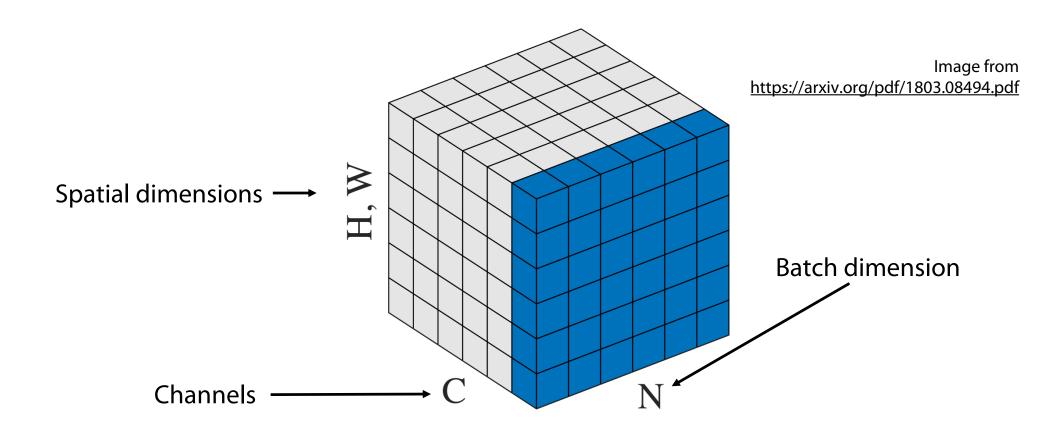
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Effectively removes the 'shift' and 'scale' degrees of freedom from the previous layer

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- Which dimension to normalize over? Typically like this:
 - Batch of 1D vectors [Batch_dim x Features_dim]
 - separately for each component in Features_dim, i.e. over Batch_dim

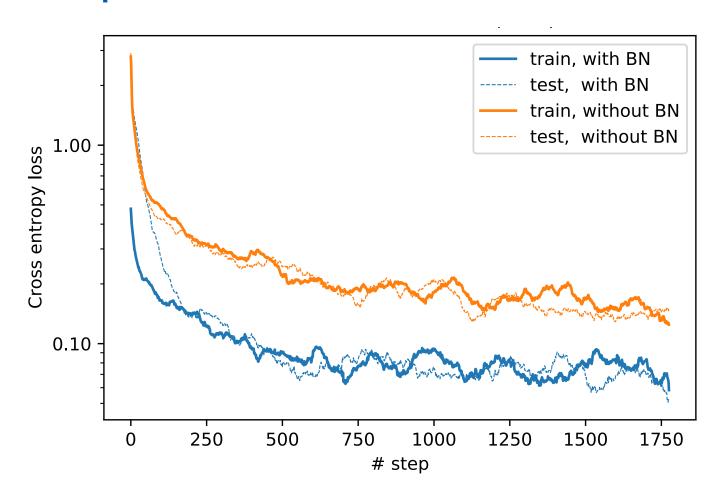
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 - Batch of ND objects [Batch_dim x Spacial_dim1 x ... x Channel_dim]
 - separately for each component in Channel_dim, i.e. over Batch_dim x Spacial_dim1 x ...



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



(shown: moving average loss)

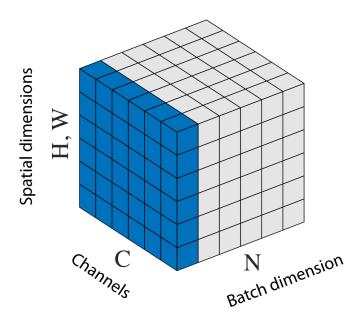
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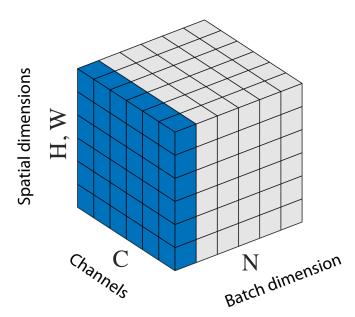




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 - the effect is quite different though
 - e.g. Layer Normalization "entangles" different neurons within a layer

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



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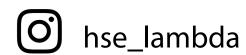
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Food for thought: how exactly would you implement an early stopping rule?

Thank you!





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