

Lecture 7 Recap

I2DL: Prof. Niessner, Prof. Leal-Taixé

Naïve Losses: L2 vs L1

• L2 Loss:

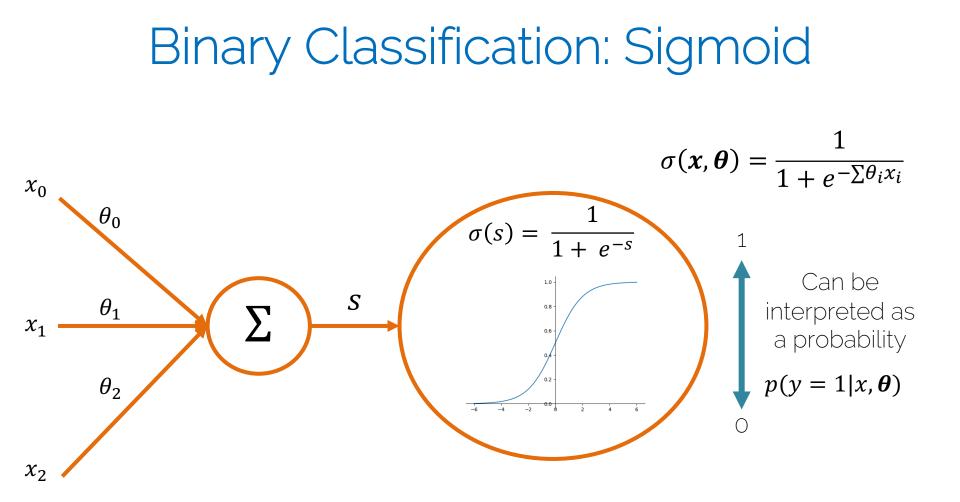
$$-L^{2} = \sum_{i=1}^{n} (y_{i} - f(x_{i}))^{2}$$

- Sum of squared
 differences (SSD)
- Prone to outliers
- Compute-efficient (optimization)
- Optimum is the mean

• L1 Loss:

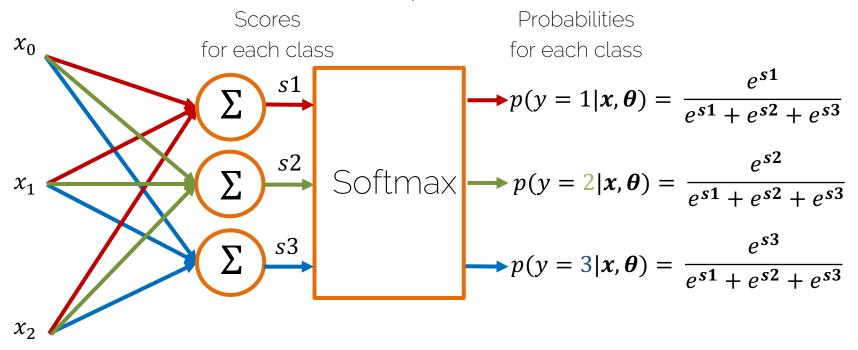
$$-L^{1} = \sum_{i=1}^{n} |y_{i} - f(x_{i})|$$

- Sum of absolute differences
- Robust
- Costly to compute
- Optimum is the median



Softmax Formulation

• What if we have multiple classes?



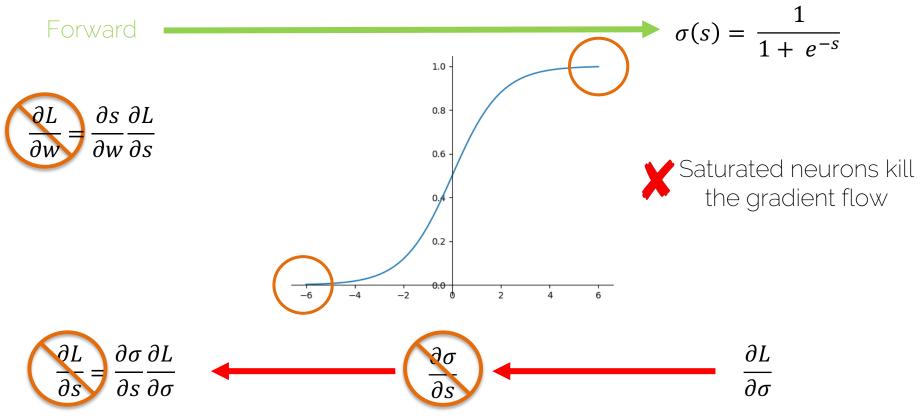
Example: Hinge vs Cross-Entropy

Hinge Loss:
$$L_i = \sum_{k \neq y_i} \max(0, s_k - s_{y_i} + 1)$$

Cross Entropy : $L_i = -\log(\frac{e^{s_{y_i}}}{\sum_k e^{s_k}})$

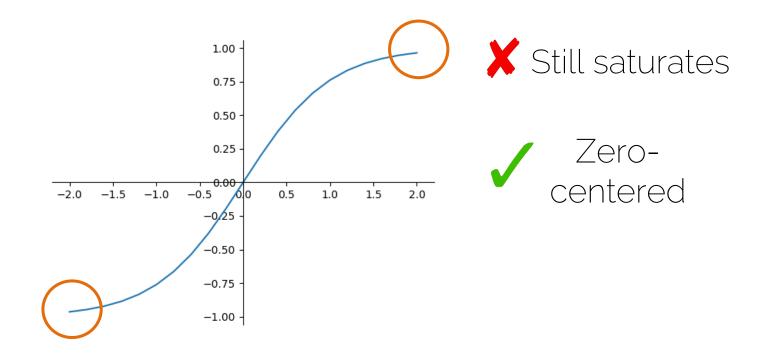
Given the	following scores for $oldsymbol{x}_i$:	Hinge loss:	Cross Entropy loss:
Model 1	s = [5, -3, 2]	max(0, -3 - 5 + 1) + max(0, 2 - 5 + 1) = 0	$-\ln\left(\frac{e^5}{e^5 + e^{-3} + e^2}\right) = 0.05$
Model 2	<i>s</i> = [5, 10, 10]	max(0, 10 - 5 + 1) + max(0, 10 - 5 + 1) = 12	$-\ln\left(\frac{e^{5}}{e^{5}+e^{10}+e^{10}}\right) = 5.70$
Model 3	s = [5, -20, -20] $y_i = 0$	$\max(0, -20 - 5 + 1) + \\\max(0, -20 - 5 + 1) = 0$	$-\ln\left(\frac{e^5}{e^5 + e^{-20} + e^{-20}}\right)$ = 2 * 10 ⁻¹¹
	- Cross	s Entropy *always* wants t	o improve! (loss never 0)
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Sigmoid Activation



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TanH Activation



[LeCun et al. 1991] Improving Generalization Performance in Character Recognition

Rectified Linear Units (ReLU)



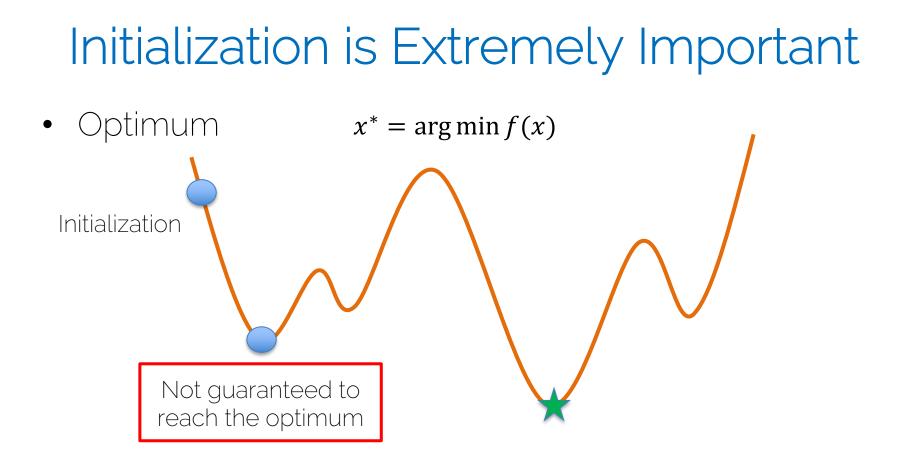
Quick Guide

• Sigmoid is not really used.

• ReLU is the standard choice.

• Second choice are the variants of ReLU or Maxout.

• Recurrent nets will require TanH or similar.



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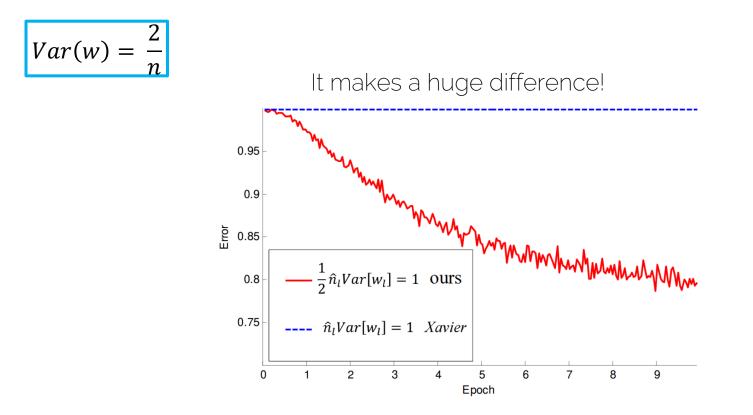
Xavier Initialization

• How to ensure the variance of the output is the same as the input?

$$(nVar(w)Var(x)) = 1$$

$$Var(w) = \frac{1}{n}$$

ReLU Kills Half of the Data



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[He et al., ICCV'15] He Initialization 12



Lecture 8



Data Augmentation

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Data Augmentation

• A classifier has to be invariant to a wide variety of transformations



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And Kittens



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Cute Baby



White Cats And Kittens





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Appearance





Illumination

Data Augmentation

• A classifier has to be invariant to a wide variety of transformations

• Helping the classifier: synthesize data simulating plausible transformations

Data Augmentation

a. No augmentation (= 1 image)



224x224



b. Flip augmentation (= 2 images)



224x224



c. Crop+Flip augmentation (= 10 images)



224x224

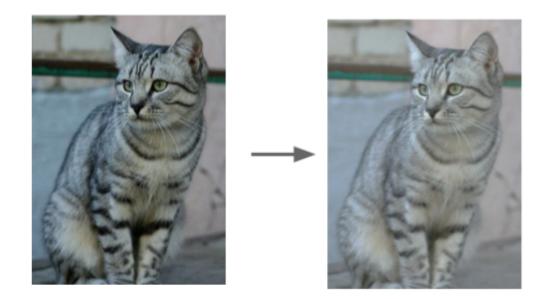


+ flips

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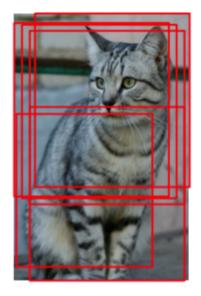
Data Augmentation: Brightness

• Random brightness and contrast changes



Data Augmentation: Random Crops

- Training: random crops
 - Pick a random L in [256,480]
 - Resize training image, short side L
 - Randomly sample crops of 224x224



- Testing: fixed set of crops
 - Resize image at N scales
 - 10 fixed crops of 224x224: (4 corners + 1 center) \times 2 flips

Data Augmentation

• When comparing two networks make sure to use the same data augmentation!

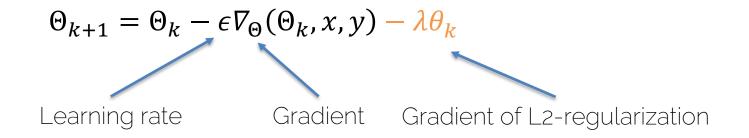
Consider data augmentation a part of your network
 design



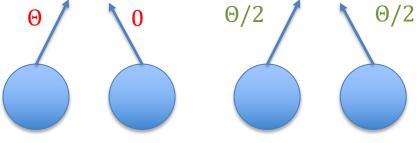
Advanced Regularization

Weight Decay

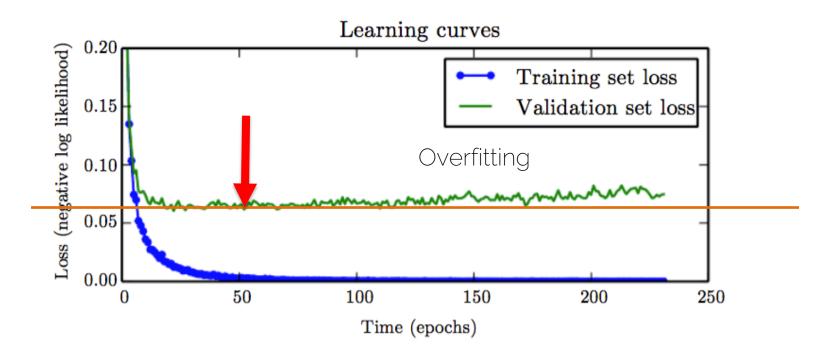
• L2 regularization



- Penalizes large weights
- Improves generalization

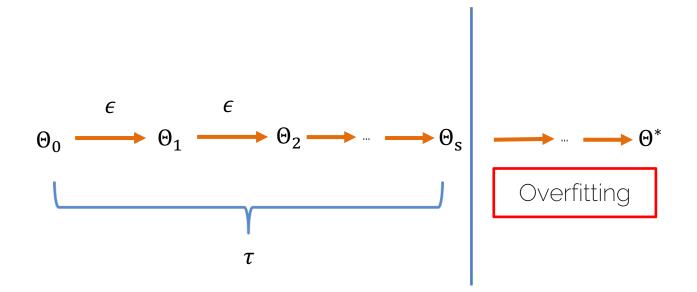


Early Stopping



Early Stopping

• Easy form of regularization



Bagging and Ensemble Methods

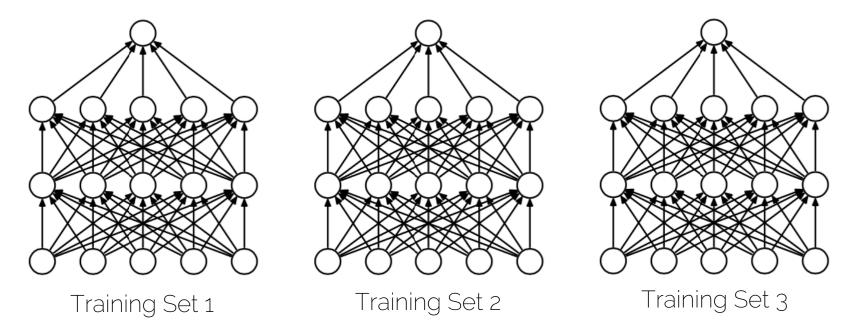
• Train multiple models and average their results

• E.g., use a different algorithm for optimization or change the objective function / loss function.

• If errors are uncorrelated, the expected combined error will decrease linearly with the ensemble size

Bagging and Ensemble Methods

• Bagging: uses k different datasets



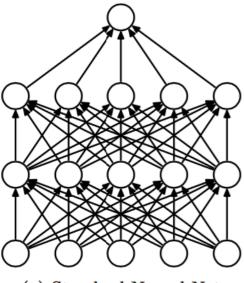


Dropout

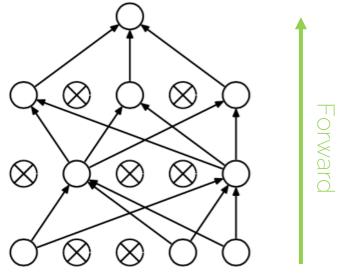
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Dropout

• Disable a random set of neurons (typically 50%)

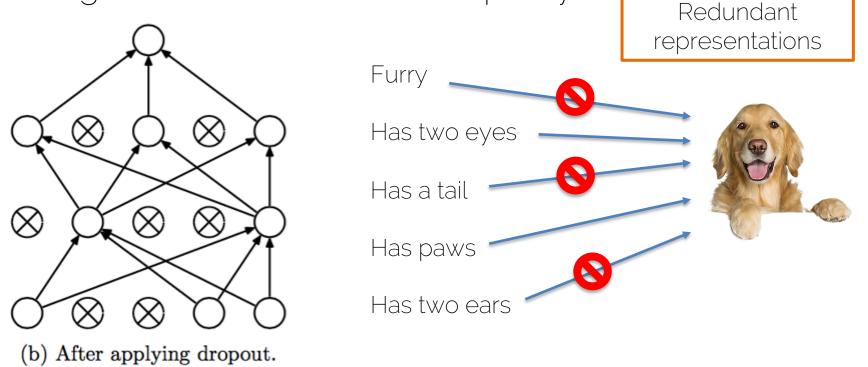


(a) Standard Neural Net



(b) After applying dropout.

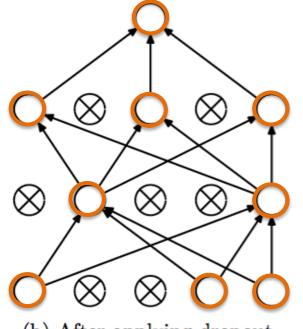
• Using half the network = half capacity

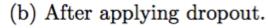


- Using half the network = half capacity
 - Redundant representations
 - Base your scores on more features

• Consider it as a model ensemble

• Two models in one

















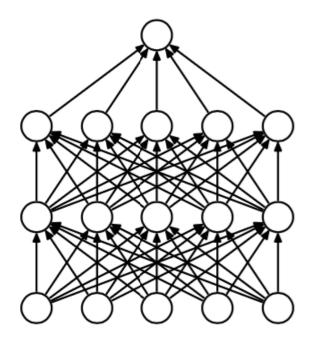
- Using half the network = half capacity
 - Redundant representations
 - Base your scores on more features

- Consider it as two models in one
 - Training a large ensemble of models, each on different set of data (mini-batch) and with SHARED parameters

Reducing co-adaptation between neurons

Dropout: Test Time

• All neurons are "turned on" – no dropout



Conditions at train and test time are not the same

Dropout: Test Time Dropout probability p = 0.5 $z = (\theta_1 x_1 + \theta_2 x_2) \cdot p$ • Test: $E[z] = \frac{1}{4}(\theta_1 0 + \theta_2 0 + \theta_1 x_1 + \theta_2 0 + \theta_1 x_1 + \theta_2 0 + \theta_1 0 + \theta_2 x_2$ • Train: Z θ_1 θ_2 $+ \theta_1 x_1 + \theta_2 x_2)$ x_2 x_1 $\theta_1 x_1 + \theta_2 x_2)$ Weight scaling inference rule

Dropout: Verdict

• Efficient bagging method with parameter sharing

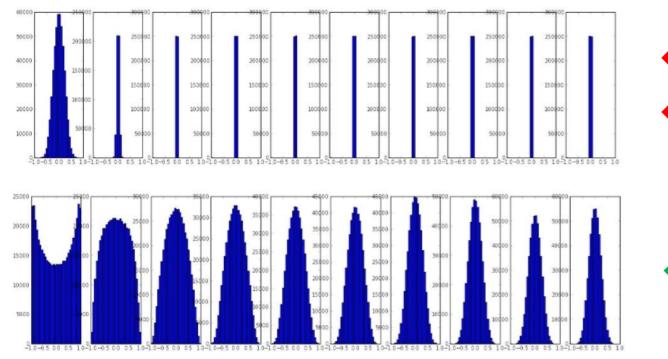
• Try it!

 Dropout reduces the effective capacity of a model → larger models, more training time

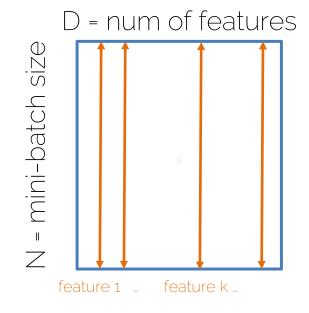


Our Goal

• All we want is that our activations do not die out



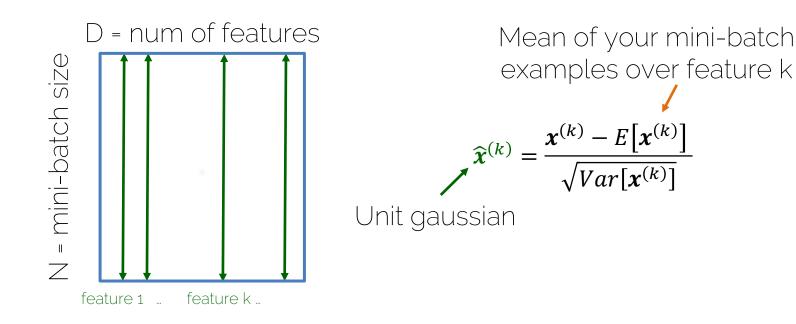
- Wish: Unit Gaussian activations (in our example)
- Solution: let's do it



Mean of your mini-batch examples over feature k $\hat{x}^{(k)} = \frac{x^{(k)} - E[x^{(k)}]}{\sqrt{Var[x^{(k)}]}}$

[loffe and Szegedy, PMLR'15] Batch Normalization $_{40}$

• In each dimension of the features, you have a unit gaussian (in our example)

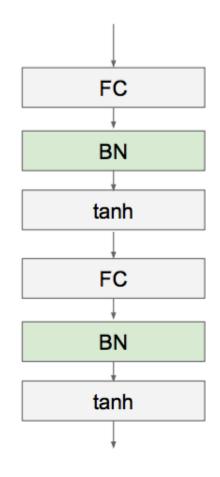


• In each dimension of the features, you have a unit gaussian (in our example)

• For NN in general \rightarrow BN normalizes the mean and variance of the inputs to your activation functions

BN Layer

• A layer to be applied after Fully Connected (or Convolutional) layers and before non-linear activation functions



• 1. Normalize

$$\widehat{\boldsymbol{x}}^{(k)} = \frac{\boldsymbol{x}^{(k)} - E[\boldsymbol{x}^{(k)}]}{\sqrt{Var[\boldsymbol{x}^{(k)}]}} \qquad \qquad \text{Differentiable function so we} \\ \text{can backprop through it}...$$

• 2. Allow the network to change the range

$$\mathbf{y}^{(k)} = \boldsymbol{\gamma}^{(k)} \hat{\mathbf{x}}^{(k)} + \boldsymbol{\beta}^{(k)} \quad \boldsymbol{\bullet}$$

These parameters will be optimized during backprop

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[loffe and Szegedy, PMLR'15] Batch Normalization 44

• 1. Normalize

$$\widehat{\boldsymbol{x}}^{(k)} = \frac{\boldsymbol{x}^{(k)} - E[\boldsymbol{x}^{(k)}]}{\sqrt{Var[\boldsymbol{x}^{(k)}]}}$$

• 2. Allow the network to change the range

$$y^{(k)} = \gamma^{(k)} \hat{x}^{(k)} + \beta^{(k)}$$

backprop

The network *can* learn to undo the normalization

$$\gamma^{(k)} = \sqrt{Var[\mathbf{x}^{(k)}]}$$
$$\beta^{(k)} = E[\mathbf{x}^{(k)}]$$

• Ok to treat dimensions separately? Shown empirically that even if features are not correlated, convergence is still faster with this method

• You can set all biases of the layers before BN to zero, because they will be cancelled out by BN anyway

BN: Train vs Test

 Train time: mean and variance is taken over the minibatch

$$\widehat{\boldsymbol{x}}^{(k)} = \frac{\boldsymbol{x}^{(k)} - \boldsymbol{E}[\boldsymbol{x}^{(k)}]}{\sqrt{Var[\boldsymbol{x}^{(k)}]}}$$

- Test-time: what happens if we can just process one image at a time?
 - No chance to compute a meaningful mean and variance

BN: Train vs Test

Training: Compute mean and variance from mini-batch 1,2,3 ...

Testing: Compute mean and variance by running an exponentially weighted averaged across training minibatches. Use them as σ_{test}^2 and μ_{test} .

 $\begin{aligned} Var_{running} &= \beta_m * Var_{running} + (1 - \beta_m) * Var_{minibatch} \\ \mu_{running} &= \beta_m * \mu_{running} + (1 - \beta_m) * \mu_{minibatch} \\ \beta_m : \text{momentum (hyperparameter)} \end{aligned}$

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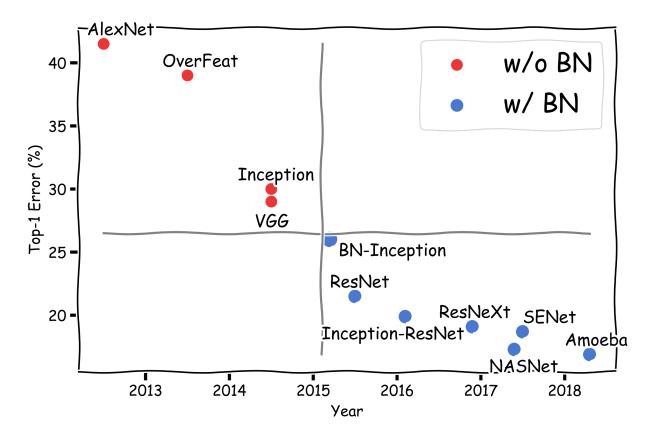
[loffe and Szegedy, PMLR'15] Batch Normalization 48

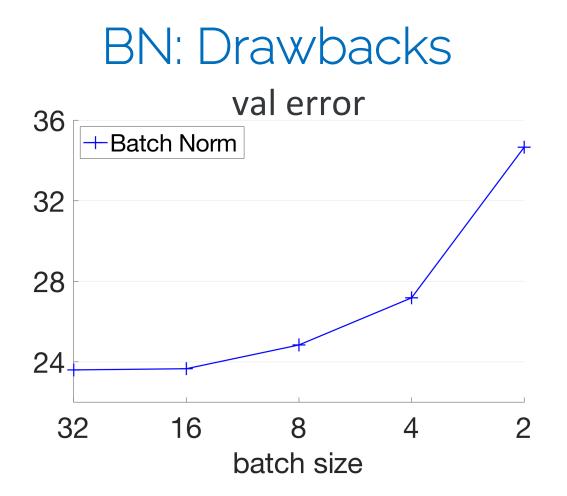
BN: What do you get?

Very deep nets are much easier to train → more stable gradients

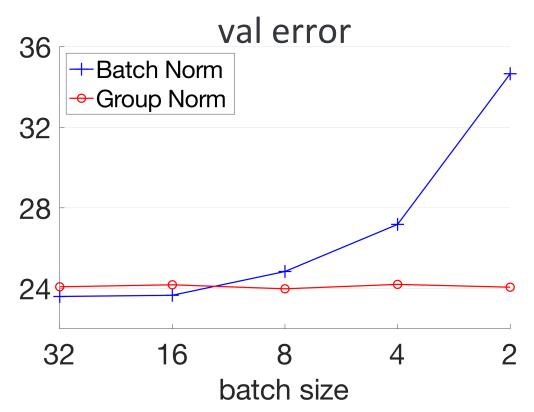
• A much larger range of hyperparameters works similarly when using BN

BN: A Milestone

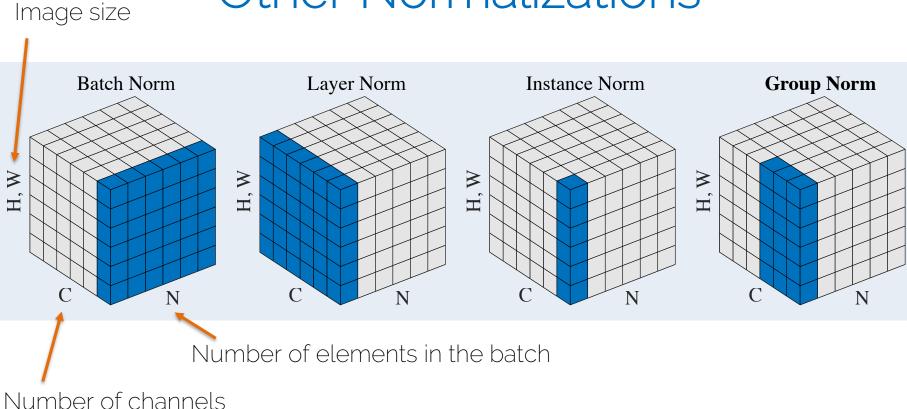




Other Normalizations

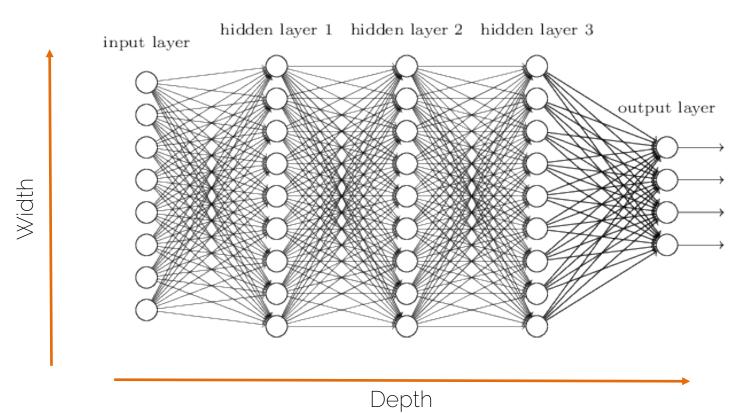


Other Normalizations

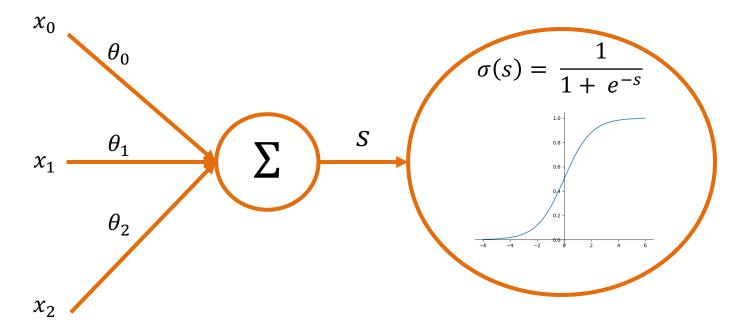




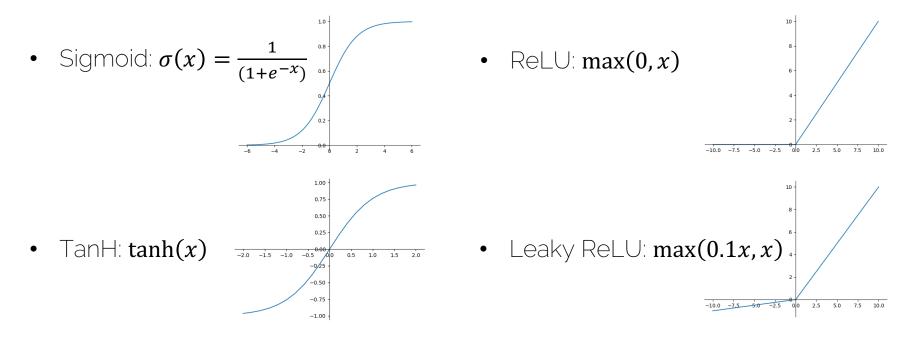
What We Know



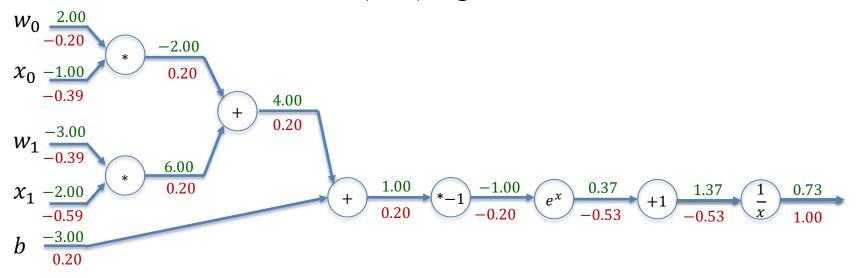
Concept of a 'Neuron'



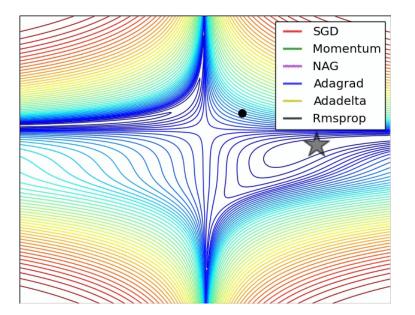
Activation Functions (non-linearities)



Backpropagation



SGD Variations (Momentum, etc.)



Data Augmentation

a. No augmentation (= 1 image)





b. Flip augmentation (= 2 images)



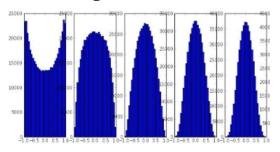




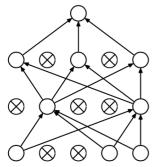
Weight Regularization e.g., L^2 -reg: $R^2(W) = \sum_{i=1}^N w_i^2$

Batch-Norm

$$\widehat{\boldsymbol{x}}^{(k)} = \frac{\boldsymbol{x}^{(k)} - E[\boldsymbol{x}^{(k)}]}{\sqrt{Var[\boldsymbol{x}^{(k)}]}}$$



Dropout



⁽b) After applying dropout.

Why not simply more Layers?

• We cannot make networks arbitrarily complex

- Why not just go deeper and get better?
 - No structure!!
 - It is just brute force!
 - Optimization becomes hard
 - Performance plateaus / drops!



See you next week!

References

- Goodfellow et al. "Deep Learning" (2016),
 Chapter 6: Deep Feedforward Networks
- Bishop "Pattern Recognition and Machine Learning" (2006),
 Chapter 5.5: Regularization in Network Nets
- <u>http://cs231n.github.io/neural-networks-1/</u>
- <u>http://cs231n.github.io/neural-networks-2/</u>
- <u>http://cs231n.github.io/neural-networks-3/</u>