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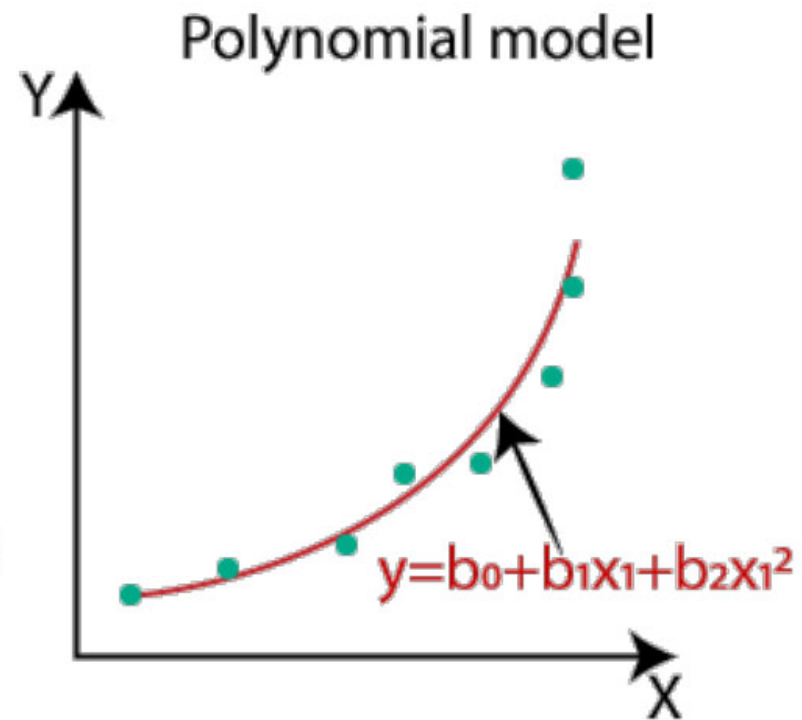
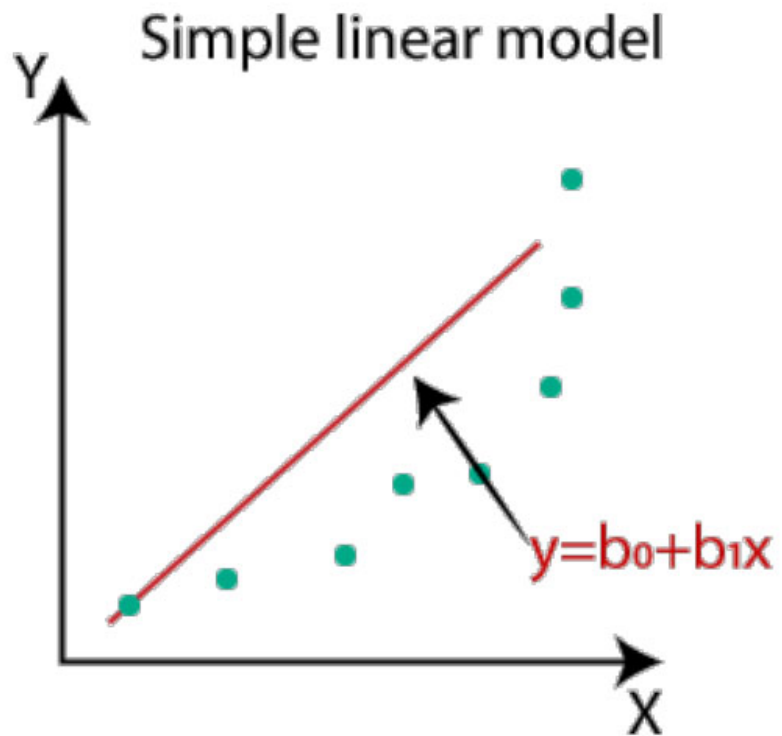
# **Deep Learning**

Winter 2026

## Regularization

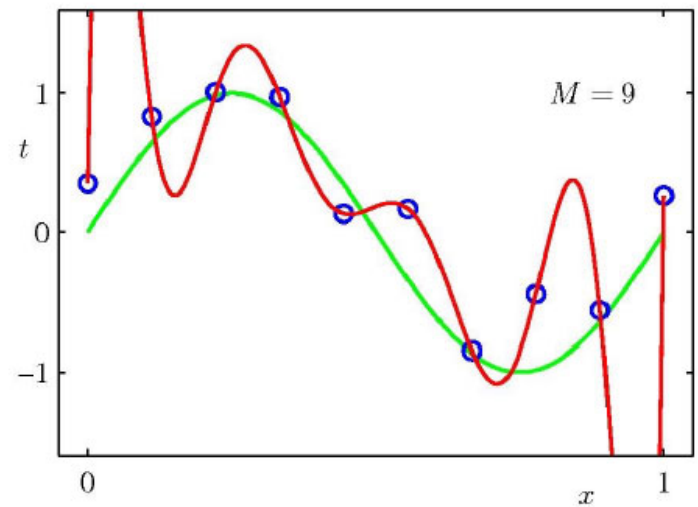
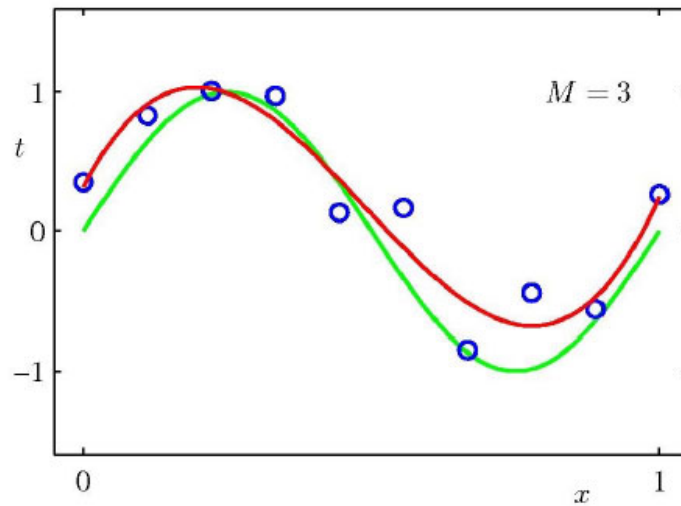
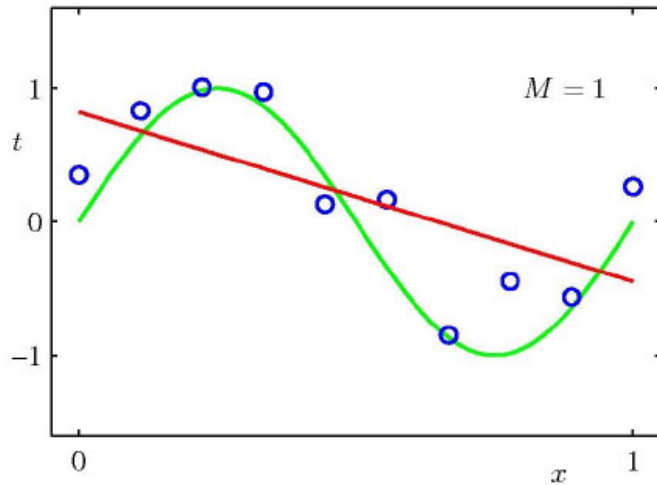
# Polynomial Regression

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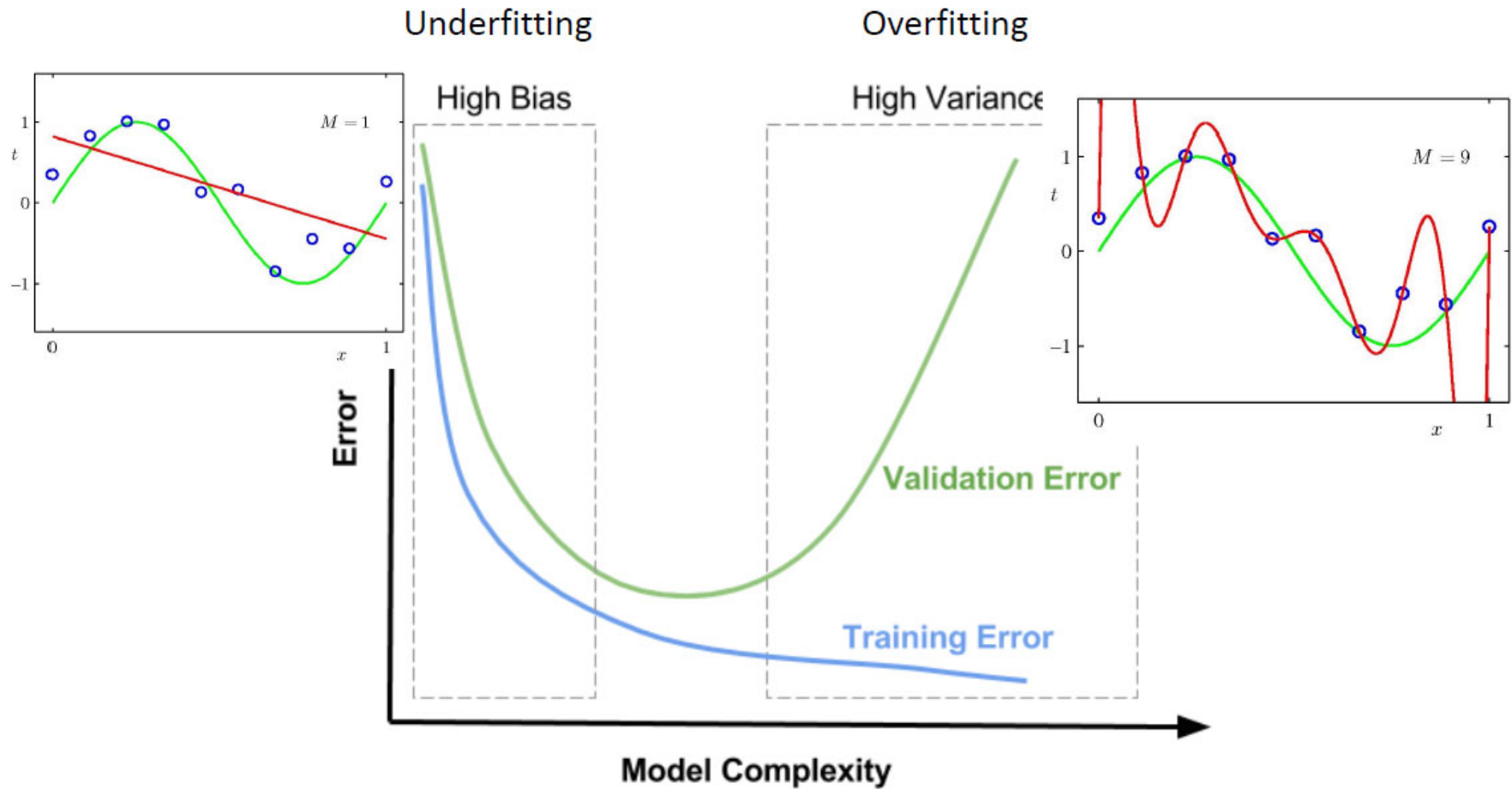


# Under/Over Fitting

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# Bias/Variance Trade-off

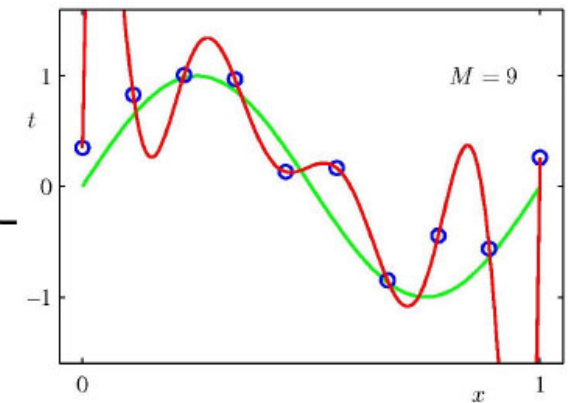
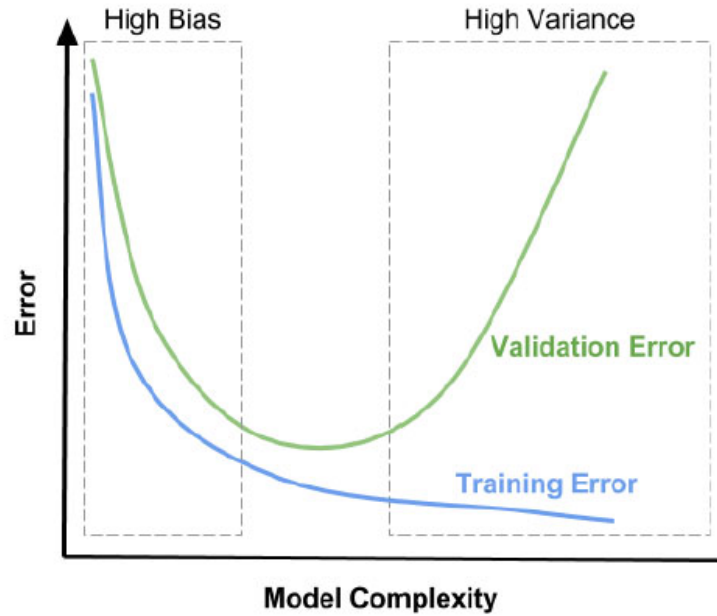
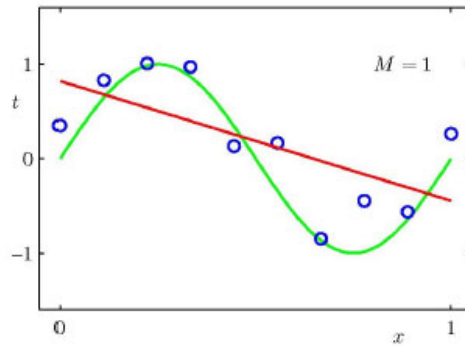


# What Goes Wrong?

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- Underfit: too many assumptions
  - Hypothesis class doesn't contain the optimal hypothesis
    - (or even a "good" hypothesis)
  - Data representation discards essential information
- Overfit: not enough assumptions
  - Hypothesis class is "hard to search"
  - Learning algorithm is inefficient, can't optimize parameters
  - Many hypotheses perfectly fit training data

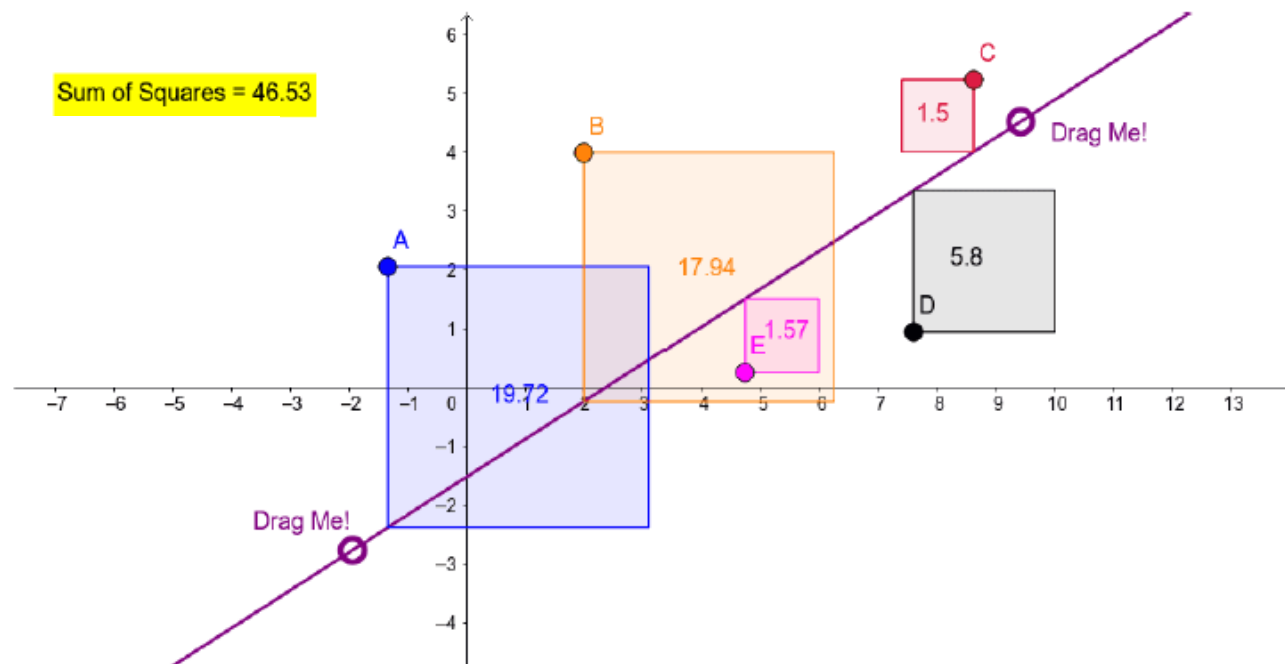
# How Do We Stop Overfitting



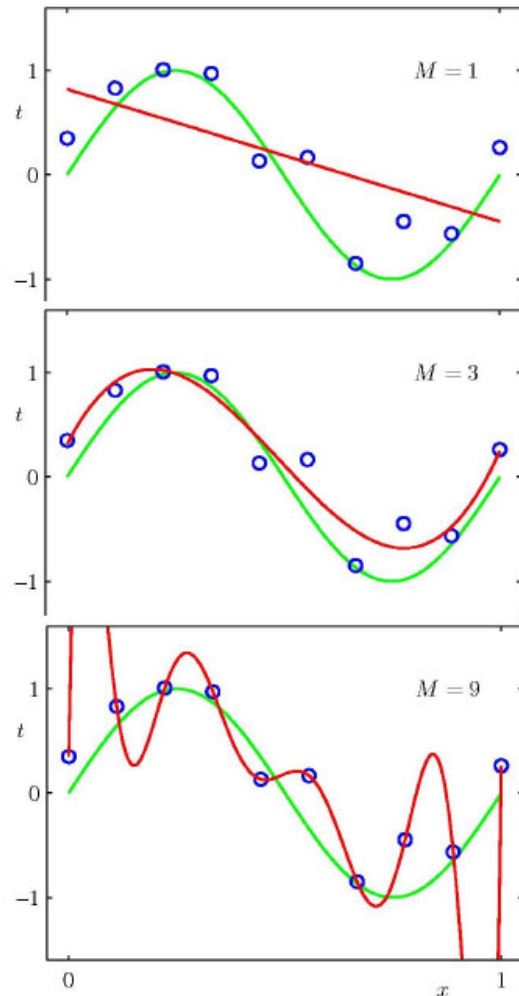
# Motivating Regularization

- Our predictor  $f$  with parameters  $w$  has loss  $L$ :

$$L(y, f(x; w)) = (y - w \cdot x)^2$$



# Motivating Regularization



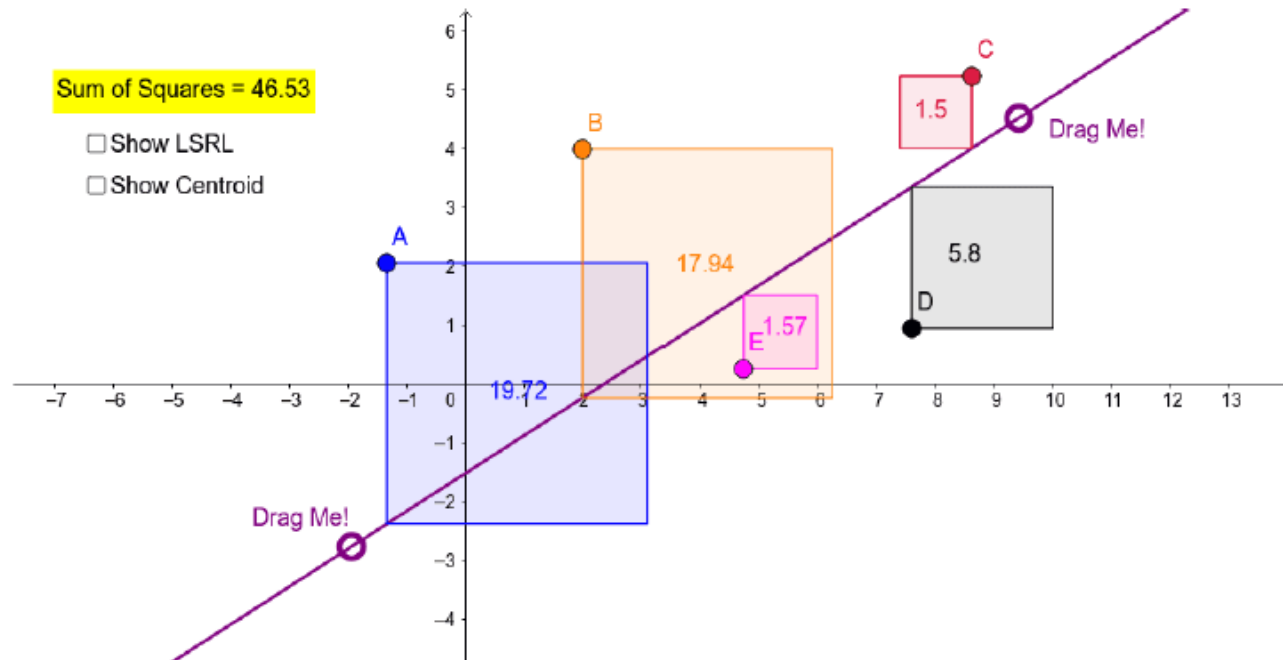
	$M = 1$	$M = 3$	$M = 9$
$w_0^*$	0.82	0.31	0.35
$w_1^*$	-1.27	7.99	232.37
$w_2^*$		-25.43	-5321.83
$w_3^*$		17.37	48568.31
$w_4^*$			-231639.30
$w_5^*$			640042.26
$w_6^*$			-1061800.52
$w_7^*$			1042400.18
$w_8^*$			-557682.99
$w_9^*$			125201.43



# Motivating Regularization

- What if we add a term to our loss?

$$L(y, f(x; w)) = (y - w \cdot x)^2 + \lambda w^2$$



# How Do We Stop Overfitting?

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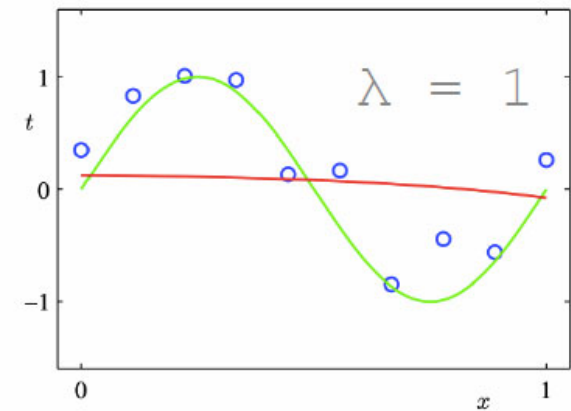
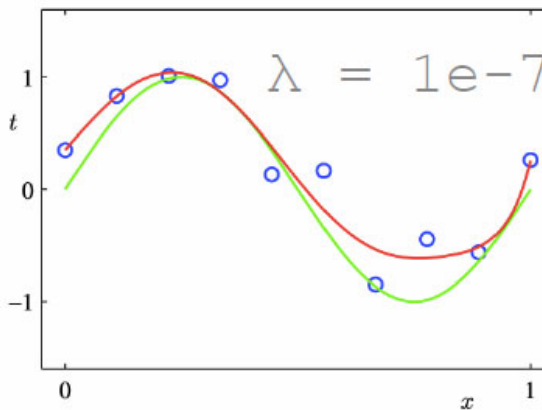
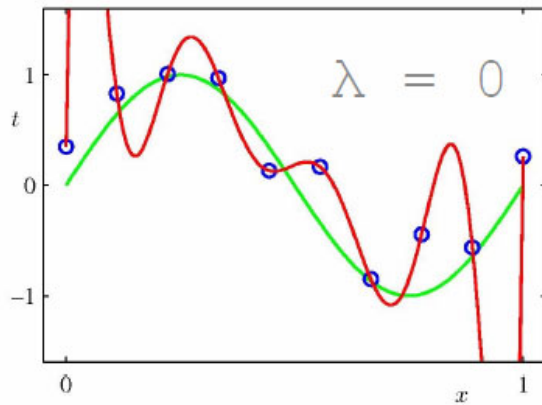
$$L(y, f(x; w)) = (y - w \cdot x)^2 + \lambda w^2$$

	$\lambda = 0$	$1e-7$	1
$w_0^*$	0.35	0.35	0.13
$w_1^*$	232.37	4.74	-0.05
$w_2^*$	-5321.83	-0.77	-0.06
$w_3^*$	48568.31	-31.97	-0.05
$w_4^*$	-231639.30	-3.89	-0.03
$w_5^*$	640042.26	55.28	-0.02
$w_6^*$	-1061800.52	41.32	-0.01
$w_7^*$	1042400.18	-45.95	-0.00
$w_8^*$	-557682.99	-91.53	0.00
$w_9^*$	125201.43	72.68	0.01

# How Do We Stop Overfitting?

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$$L(y, \hat{f}(x; w)) = (y - w \cdot x)^2 + \lambda w^2$$



# Regularized Least Squares

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$$\mathcal{L}(\mathbf{w}) = \frac{1}{2} \sum_{i=1}^N \{y_i - \mathbf{w}^T \cdot \mathbf{x}_i\}^2 \quad \mathcal{L}(\mathbf{w}) = \frac{1}{2} \sum_{i=1}^N \{y_i - \mathbf{w}^T \cdot \mathbf{x}_i\}^2 + \lambda \frac{1}{2} \mathbf{w}^T \mathbf{w}$$

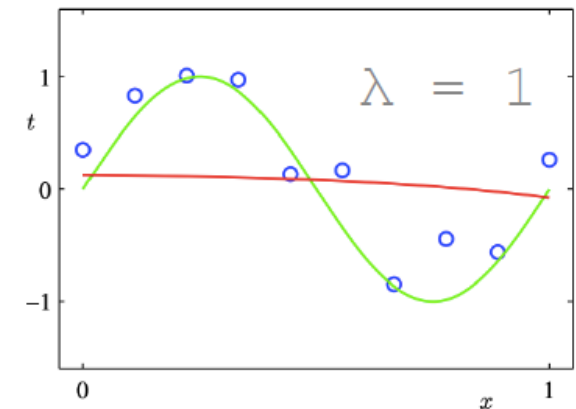
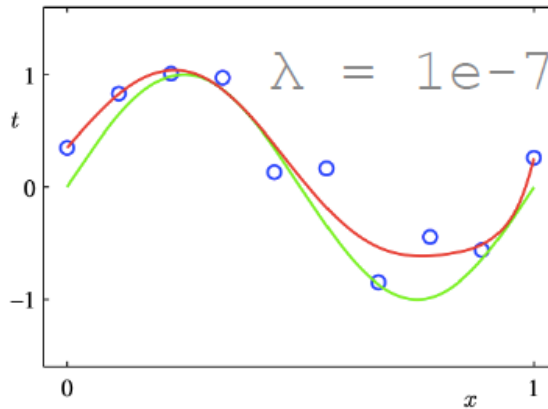
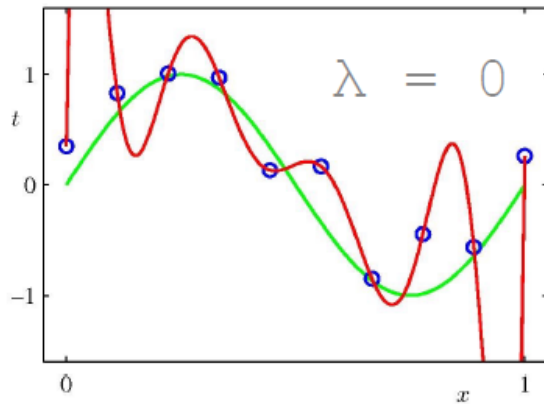
$$\mathbf{w} = (\mathbf{X}^T \mathbf{X})^{-1} \mathbf{X}^T \mathbf{Y}$$

$$\mathbf{w} = (\lambda \mathbf{I} + \mathbf{X}^T \mathbf{X})^{-1} \mathbf{X}^T \mathbf{Y}$$

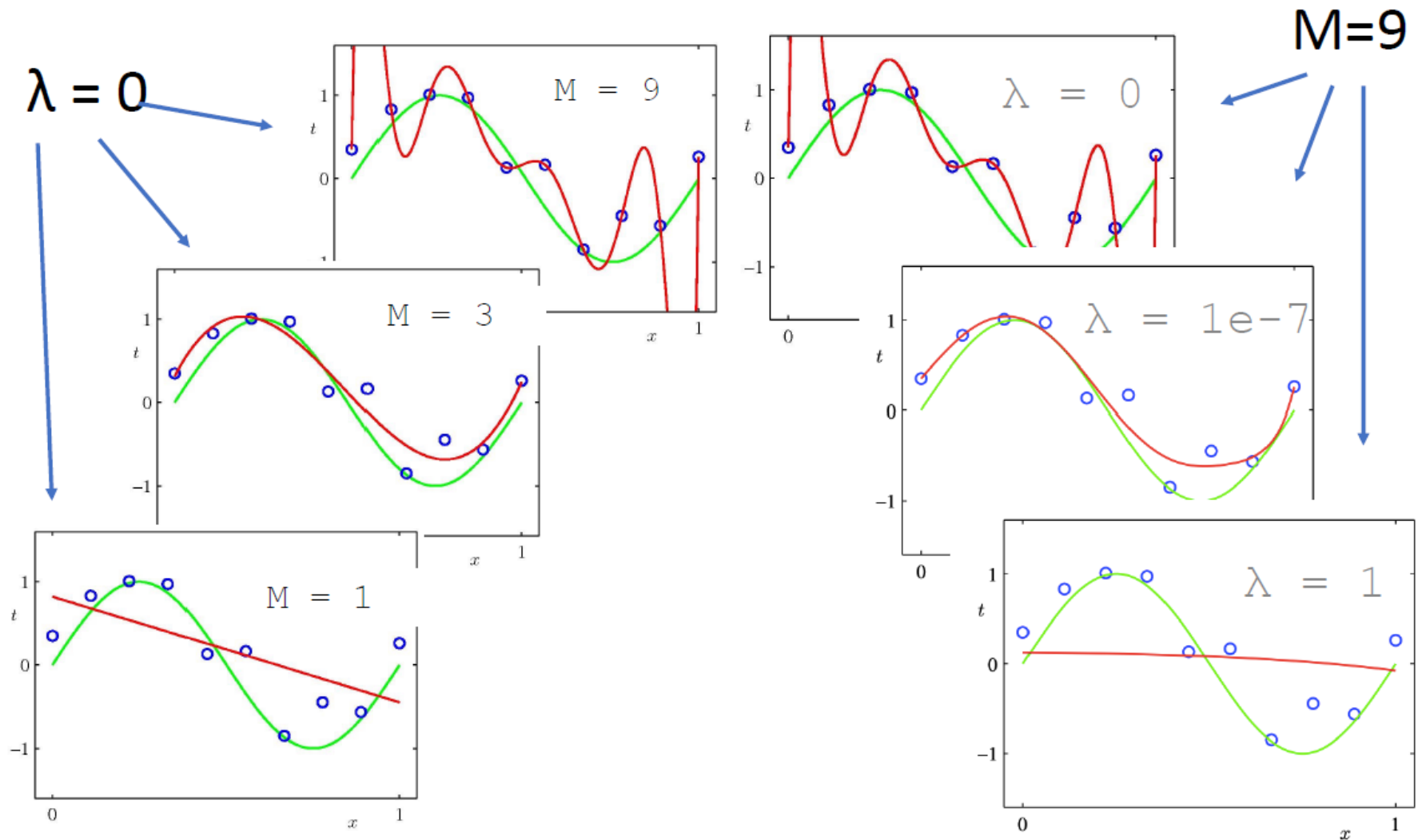
# How Do We Stop Overfitting?

$$L(y, \hat{f}(x; w)) = (y - w \cdot x)^2 + \lambda w^2$$

$$w = (\lambda I + X^T X)^{-1} X^T Y$$



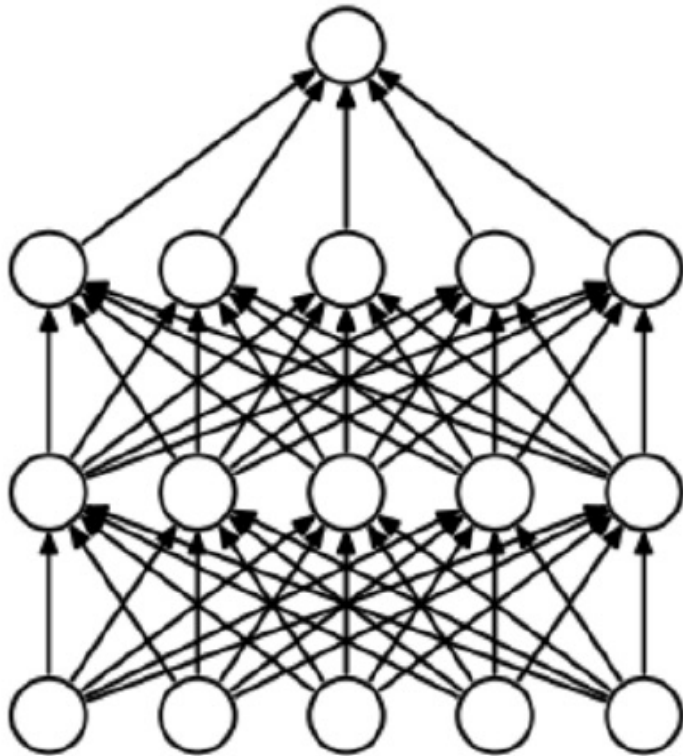
# Summary of Regularization



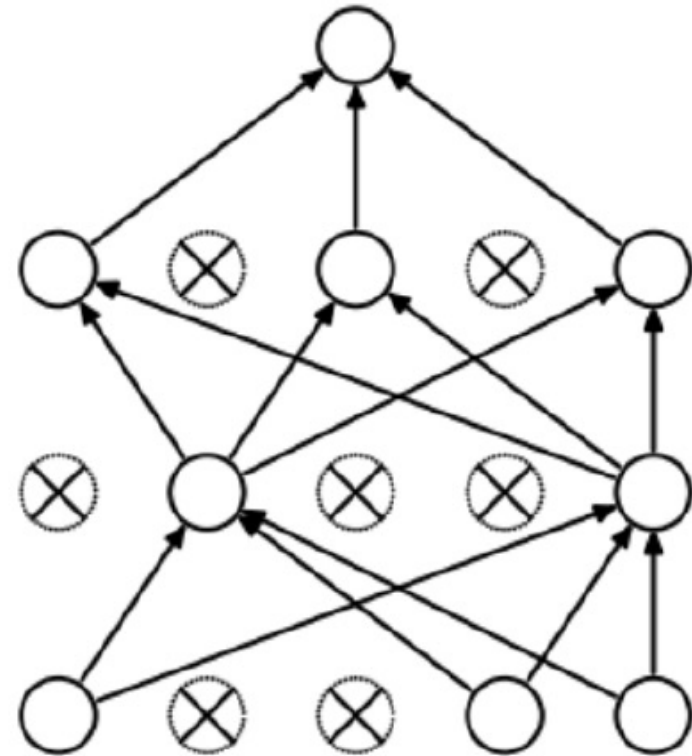
# Dropout

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- Instead of modifying objective function (regularization), modify the **network structure**



(a) Standard Neural Net

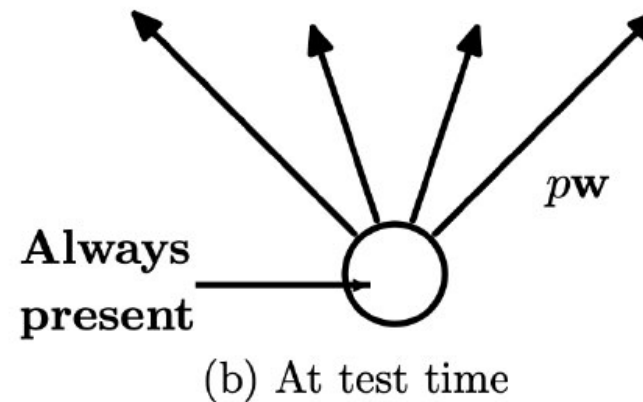
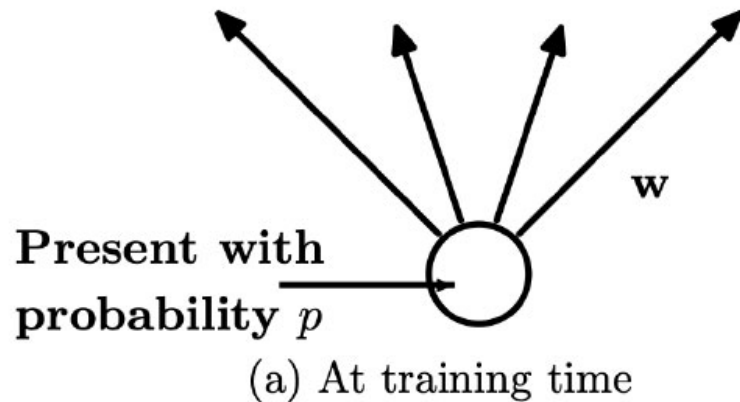


(b) After applying dropout.

# Dropout

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- Randomly “dropout” some neurons with probability  $p$
- The chosen neurons change on each mini-batch
- All other learning remains unchanged



Srivastava, Nitish, et al. "Dropout: a simple way to prevent neural networks from overfitting." *Journal of machine learning research* 15.1 (2014)



# Why Does Dropout Work?

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- **Noise robustness**

- Dropout is similar to making random perturbations in the input
- These small changes shouldn't change the prediction for an input
- Think “reduces variance”

# Why Does Dropout Work?

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- Prevents neuron co-adaptation
  - Neurons, especially near the top of the network, suffer from co-adaptation
  - The neurons take on values that rely on other neurons for correct predictions
  - Multiple points of failure: any single neuron failing can invalidate the others
  - Dropout prevents co-adaptation by randomly removing some of these neurons
  - Network is forced to learn more robust features that are useful with many different random subsets of other neurons

# Why Does Dropout Work?

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- Ensemble training
  - Similar to random forests or boosted models: we can combine many classifiers to produce better output
  - During training, each random permutation is one of  $2^n$  possible neural networks
  - All networks share weights so number of parameters is still  $O(n^2)$
  - Dropout training is like training a collection of  $2^n$  "thinned" networks with extensive weight sharing
  - Test time: approximate the average prediction of all networks by using a single network without dropout

# Dropout Takeaways

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- Works really well for a wide variety of network types and domains
- Very easy to implement with no modifications in training
  - Some minor details (omitted) are important
- Drawbacks
  - Need to select dropout rate
  - This can depend on types of layers and network structure

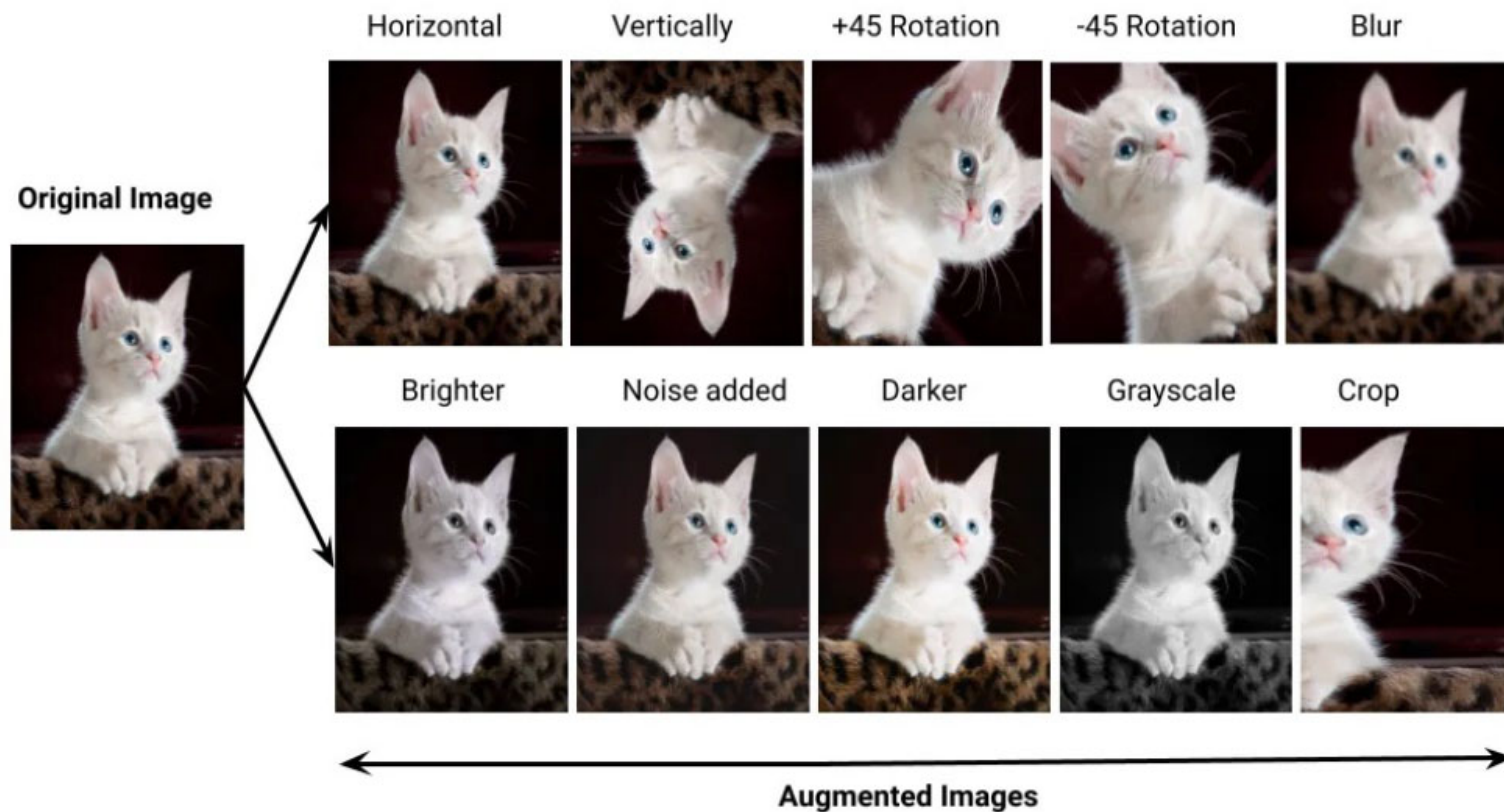
# Data Augmentation

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- Make perturbed copies of your data that vary in ways that should not change the value nature of the output function.
- This can help prevent spurious correlations between data and output.
- Example: Distinguishing clarinet sounds from flute sounds
  - Vary the pitch of each note by + or – 1%, 2%, 3%, 4%....
  - Add background noise of different kinds and at different dB
  - Time-stretch each note a bit
  - Delay or advance the onset of the note
- This can turn 1000 data points into 100,000.

# Image Augmentation

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# More Image Augmentation

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# Image Augmentation -- What if....

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Your training data  
looks like this?

High dynamic range  
Very bright



Your testing data  
looks like this?

Low dynamic range  
Very dark



# Image Augmentation -- What if...

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- Standardize your data:
  - Make sure that you have unit variance in your batch/dataset.
- Give your data the same range overall (e.g. center your values around the same center point)
- Decorrelate your variables (can be harder for images, if every pixel is a variable)

# Batch Normalization

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- Normalize the data scale input to each node
- Subtract the mean value of the data
- Do this on a dimension-by-dimension basis
- Do this at every training step in gradient descent

# Batch Normalization

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**Input:** Values of  $x$  over a mini-batch:  $\mathcal{B} = \{x_{1\dots m}\};$

Parameters to be learned:  $\gamma, \beta$

**Output:**  $\{y_i = \text{BN}_{\gamma, \beta}(x_i)\}$

$$\mu_{\mathcal{B}} \leftarrow \frac{1}{m} \sum_{i=1}^m x_i \quad // \text{ mini-batch mean}$$

$$\sigma_{\mathcal{B}}^2 \leftarrow \frac{1}{m} \sum_{i=1}^m (x_i - \mu_{\mathcal{B}})^2 \quad // \text{ mini-batch variance}$$

$$\hat{x}_i \leftarrow \frac{x_i - \mu_{\mathcal{B}}}{\sqrt{\sigma_{\mathcal{B}}^2 + \epsilon}} \quad // \text{ normalize}$$

$$y_i \leftarrow \gamma \hat{x}_i + \beta \equiv \text{BN}_{\gamma, \beta}(x_i) \quad // \text{ scale and shift}$$

# Batch Normalization

## Network with BatchNorm



Faster Convergence



Robust to Hyperparameters