

Lecture 9: Theoretical basics of machine learning

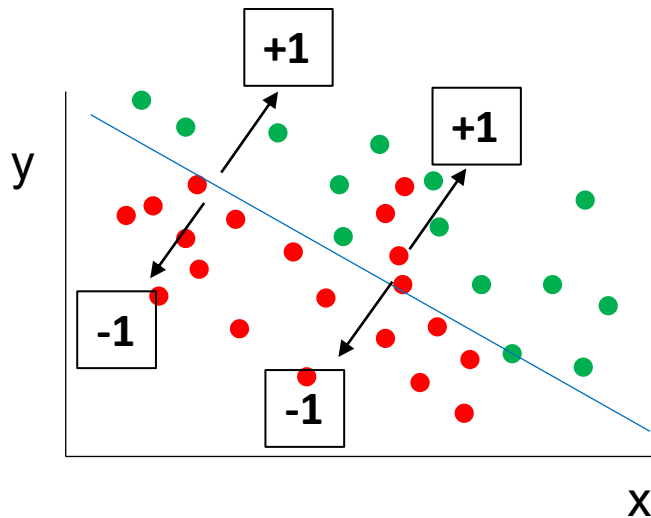
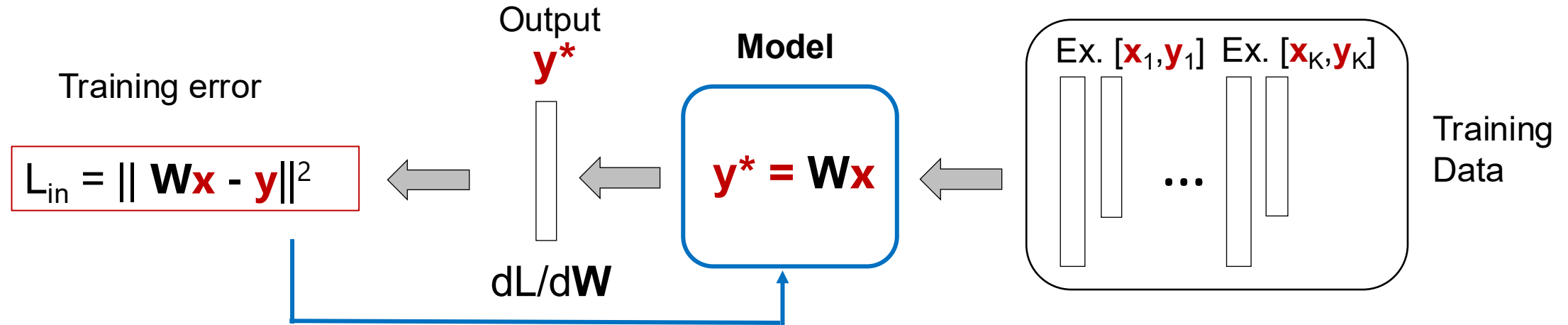
Machine Learning and Imaging

BME 548L
Roarke Horstmeyer

Announcements

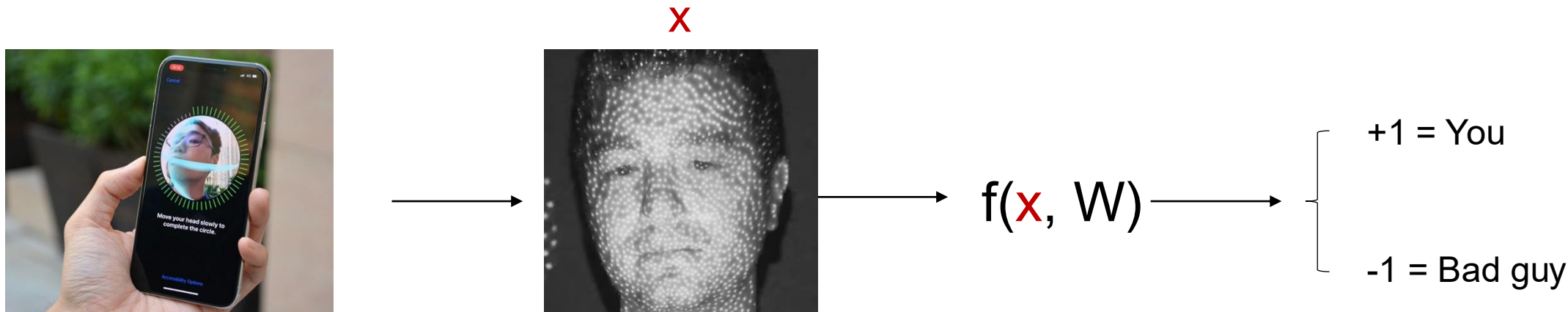
- HW1 due *TODAY*, 2/11 at 11:59pm
 - Submit via Canvas
- Lab workbooks due today
- HW2 will be posted soon, will be due ~**two weeks after**

The linear classification model – what's not to like?



1. Can only separate data with lines (hyper-planes)...
2. We only allowed for binary labels ($y = +/- 1$)
3. Error function L_{in} inherently makes assumptions about statistical distribution of data

Cost functions matter: a simple example



What if you're a CIA agent?

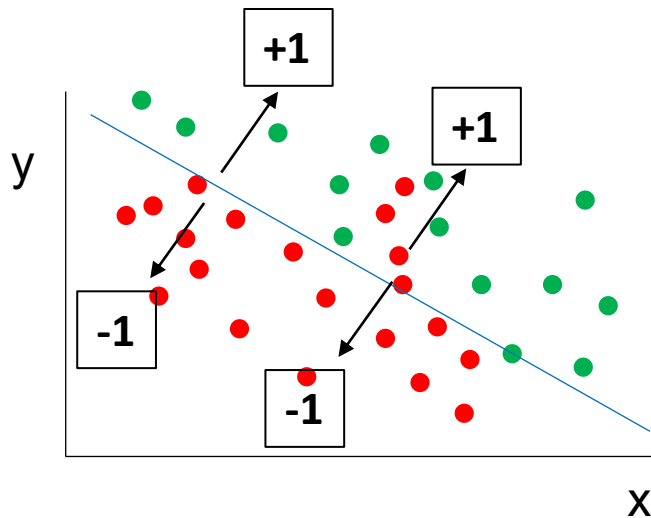
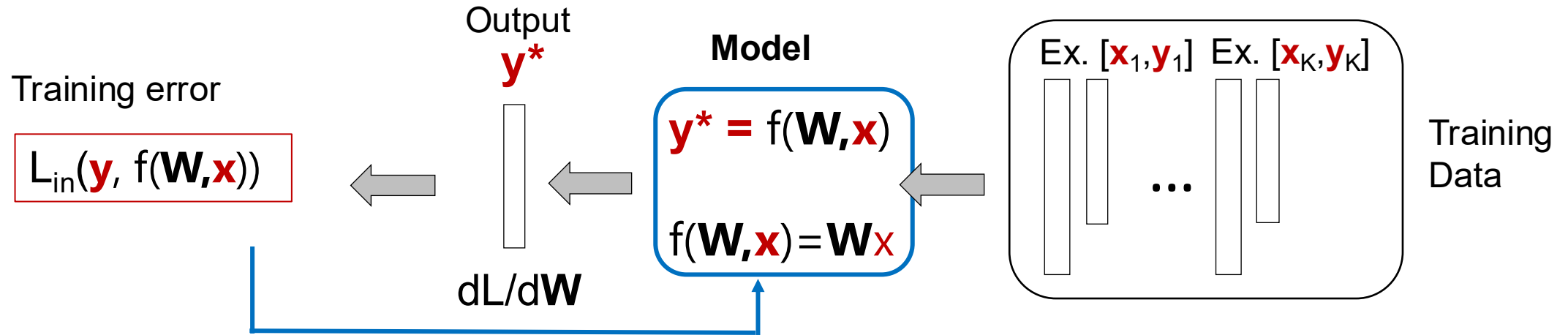
$$L_{in} = \underbrace{100,000 \text{ ReLU}[f(\mathbf{x}, \mathbf{W}) - y]}_{\text{BIG penalty for intruder}} + \underbrace{\text{ReLU}[y - f(\mathbf{x}, \mathbf{W})]}_{\text{Don't mind about annoyance...}}$$

$f(\mathbf{x}, \mathbf{W})$

	+1	-1	
+1	No Error	False reject	It's you, but you can't get in...
-1	False accept	No Error	

Letting an intruder in

The linear classification model – what's not to like?



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Deriving cost function for logistic classification for probabilistic outputs

Similar to the linear classification case, the likelihood of observing N independent outputs is given by,

$$\begin{aligned} P(y_1, y_2 \dots y_N \mid \mathbf{x}_1, \mathbf{x}_2, \dots \mathbf{x}_N) &= \prod_{n=1}^N P(y_n \mid \mathbf{x}_n) \\ &= \prod_{n=1}^N \theta(y_n \mathbf{w}^T \mathbf{x}_n) \end{aligned}$$

The Logistic Function θ

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

Also called
Sigmoid
function

This is the probability of the labels, given the data. We'd like to maximize this probability!

*Like the linear regression case, but now the probability of classes given the data is not Gaussian distributed, but instead follows the sigmoid curve (is bound to $[0, 1]$, which is more realistic)

$$\text{Maximize } P(y_1, y_2 \dots y_N \mid \mathbf{x}_1, \mathbf{x}_2, \dots \mathbf{x}_N) = \prod_{n=1}^N \theta(y_n \mathbf{w}^T \mathbf{x}_n)$$

Deriving cost function for logistic classification for probabilistic outputs

$$\text{Maximize } P(y_1, y_2, \dots, y_N \mid \mathbf{x}_1, \mathbf{x}_2, \dots, \mathbf{x}_N) = \prod_{n=1}^N \theta(y_n \mathbf{w}^T \mathbf{x}_n)$$

$$\text{Minimize } -\frac{1}{N} \ln \left(\prod_{n=1}^N \theta(y_n \mathbf{w}^T \mathbf{x}) \right)$$

$$\text{Minimize } \frac{1}{N} \sum_{n=1}^N \ln \left(\frac{1}{\theta(y_n \mathbf{w}^T \mathbf{x})} \right)$$

$$\text{Use relationship } \theta(a) = \frac{1}{1 + e^{-a}}$$

$$\text{Minimize } L_{in}(\mathbf{w}) = \frac{1}{N} \sum_{n=1}^N \ln(1 + e^{-y_n \mathbf{w}^T \mathbf{x}})$$

Cross entropy error for logistic classification

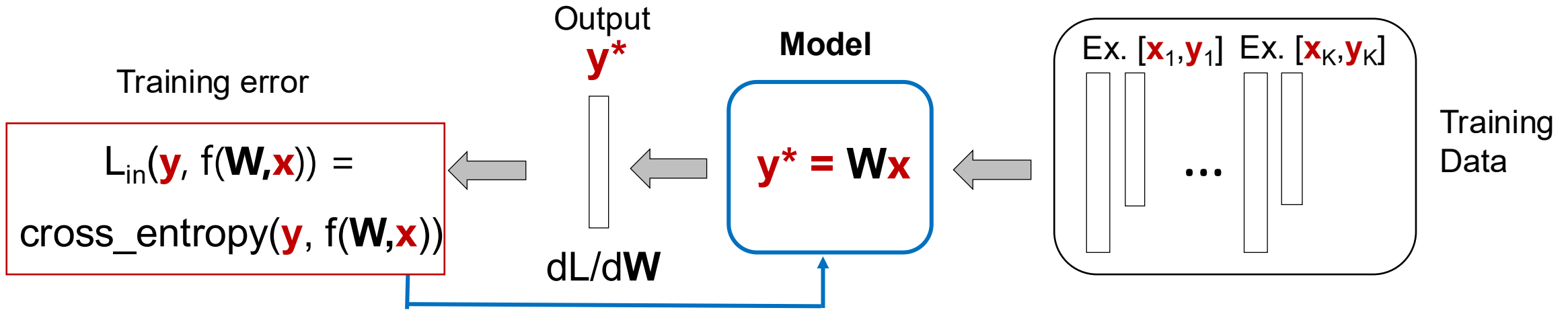
Typically requires iterative solution to minimize

$$L_{in}(\mathbf{w}) = \frac{1}{N} \sum_{n=1}^N (y_n - \mathbf{w}^T \mathbf{x})^2$$

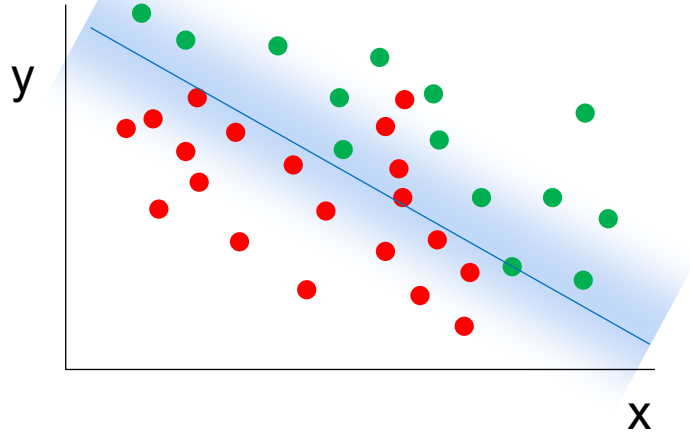
Mean-square error for linear classification

Closed form solution available

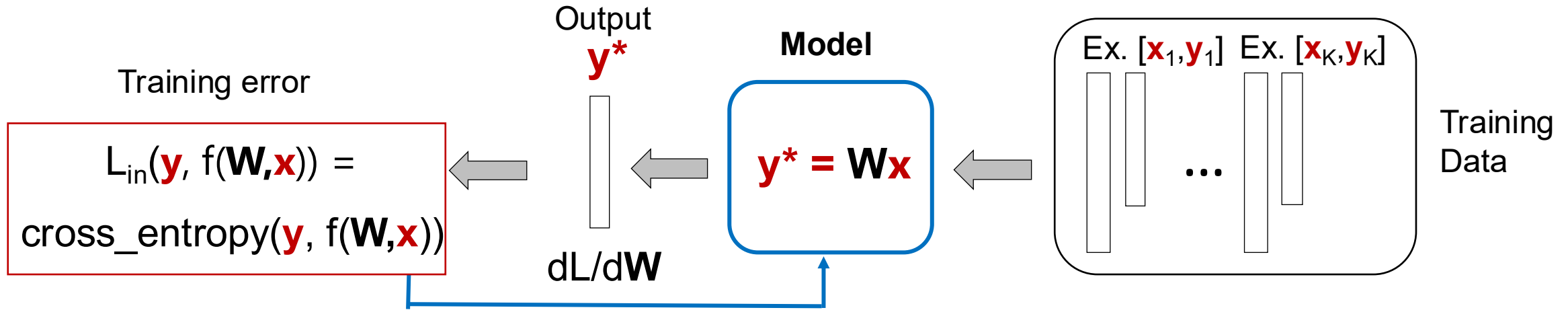
The linear classification model – what's not to like?



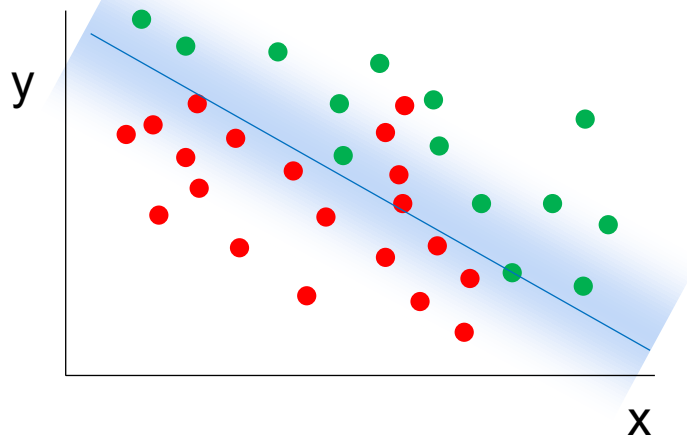
Probabilistic mapping to y



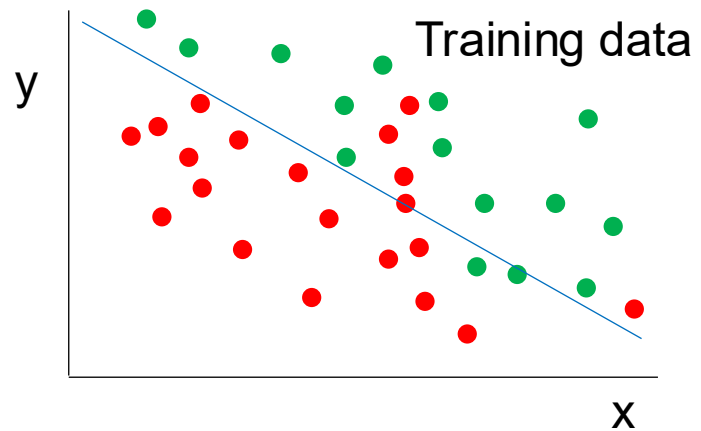
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Probabilistic mapping to y

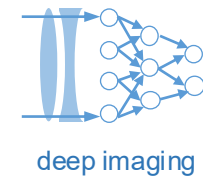
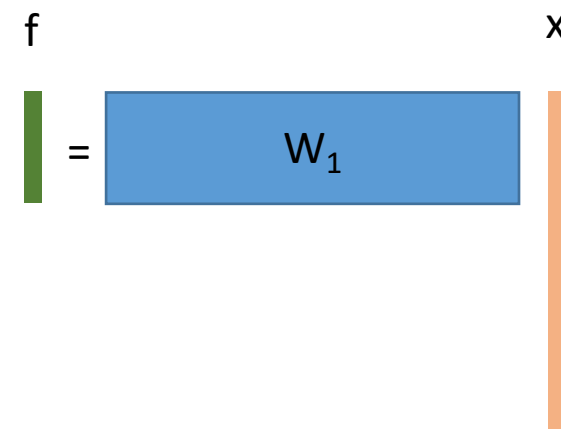


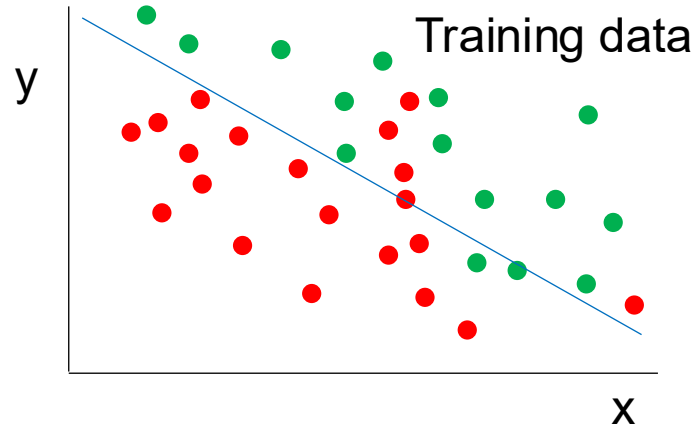
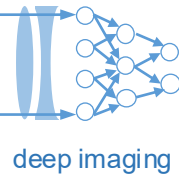
1. Can only separate data with lines (hyper-planes)...
2. We only allowed for binary labels ($y = \pm 1$)
3. Error function L_{in} inherently makes assumptions about statistical distribution of data



$$f = W_1 x$$

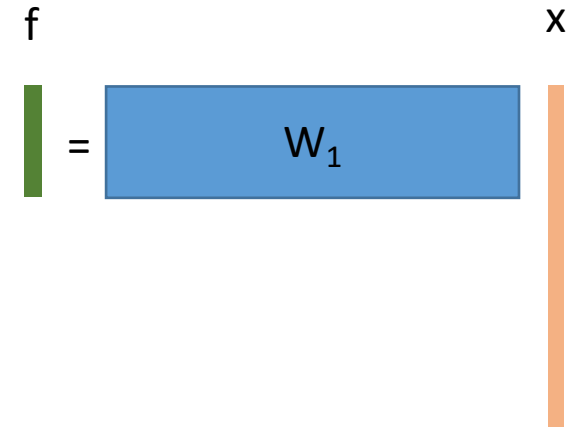
Learned f : not flexible





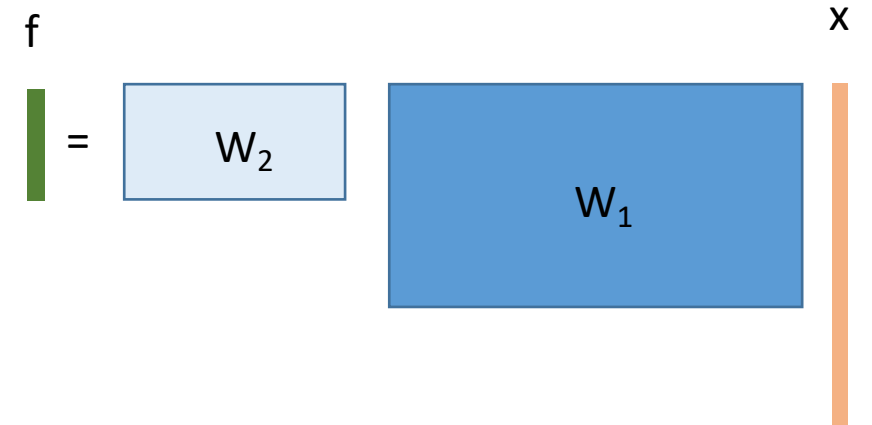
$$f = W_1 x$$

Learned f : not flexible



Can we add flexibility by multiplying with another weight matrix?

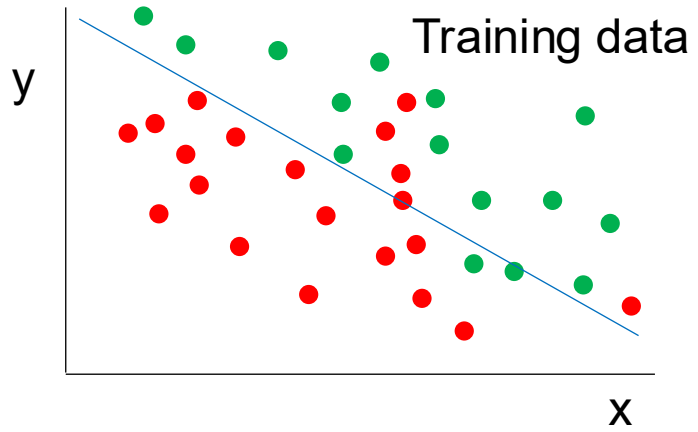
$$\begin{cases} f_1 = W_1 x + b_1 \\ f_2 = W_2 f_1 + b_2 \end{cases}$$



$$f_2 = W_2(W_1 x + b_1) + b_2$$

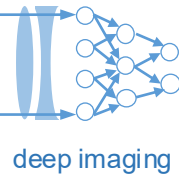
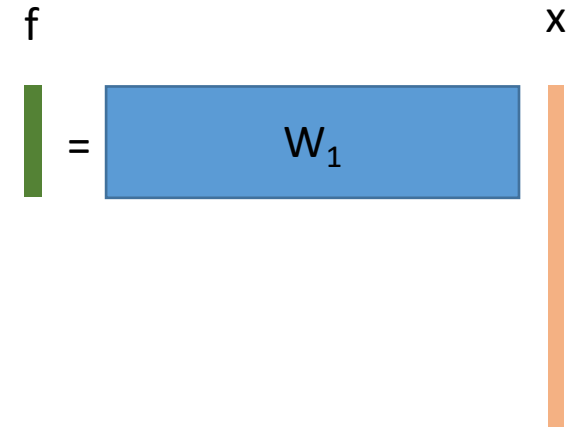
$$f_2 = W' x + b'$$

Unfortunately not...



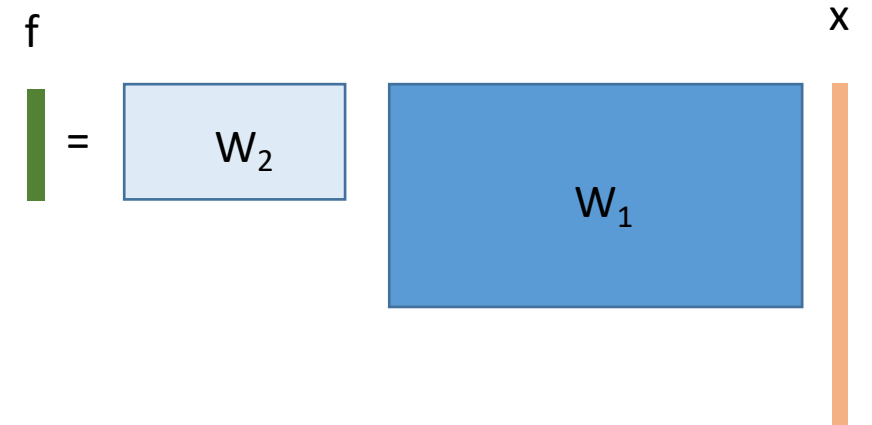
$$f = W_1 x$$

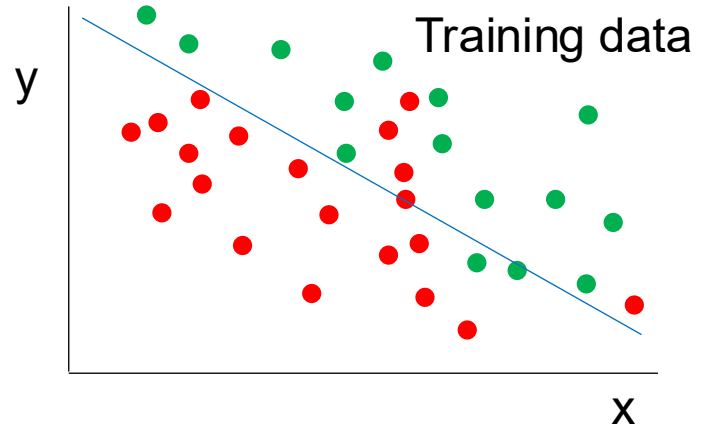
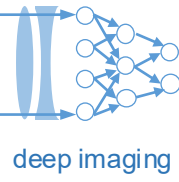
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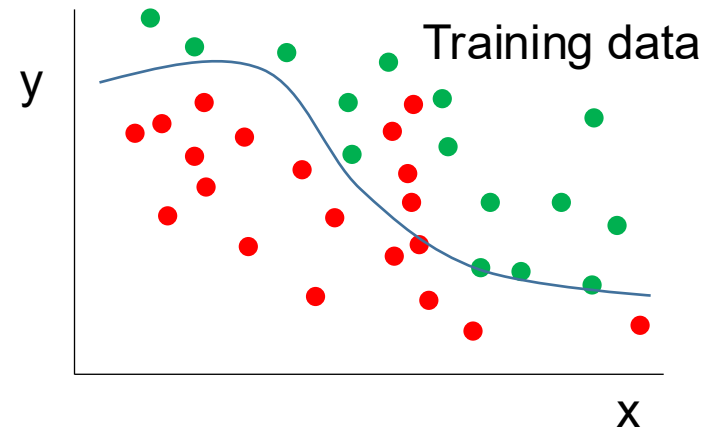
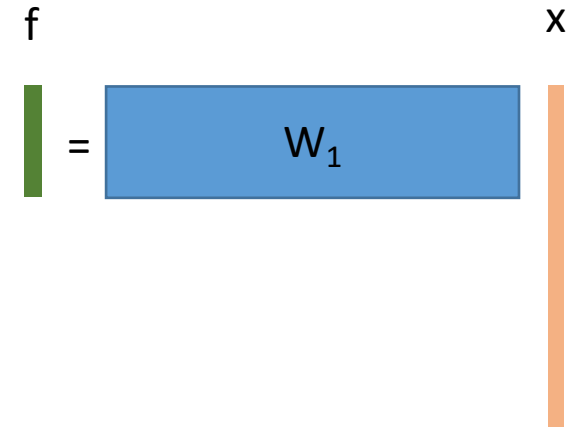
$$\begin{cases} f_1 = W_1 x + b_1 \\ f_2 = W_2 f_1 + b_2 \end{cases}$$





$$f = W_1 x$$

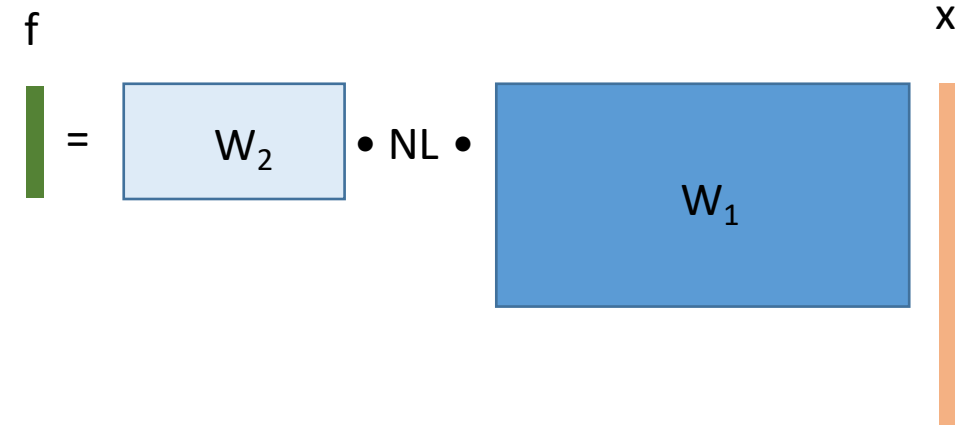
Learned f : not flexible

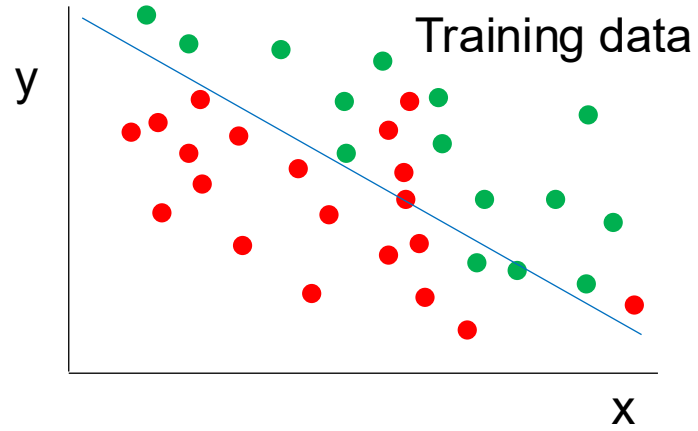
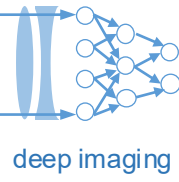


Add a non-linearity!

$$f = W_2 \max(W_1 x, 0)$$

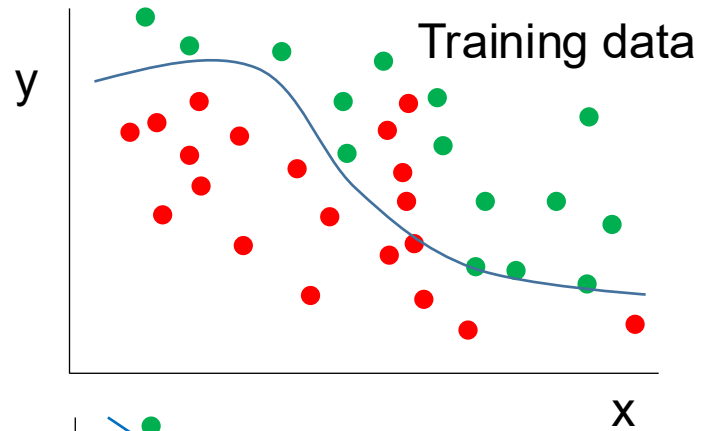
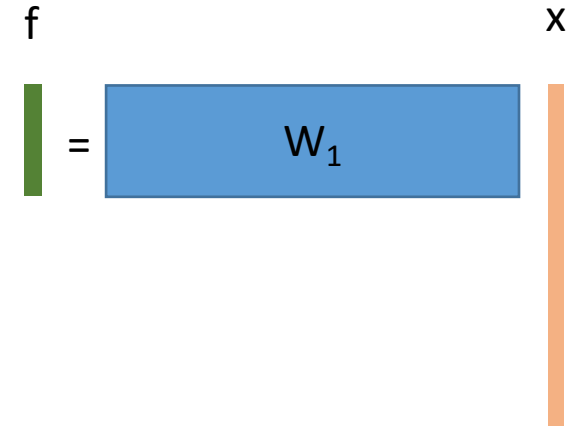
Learned f : a bit flexible





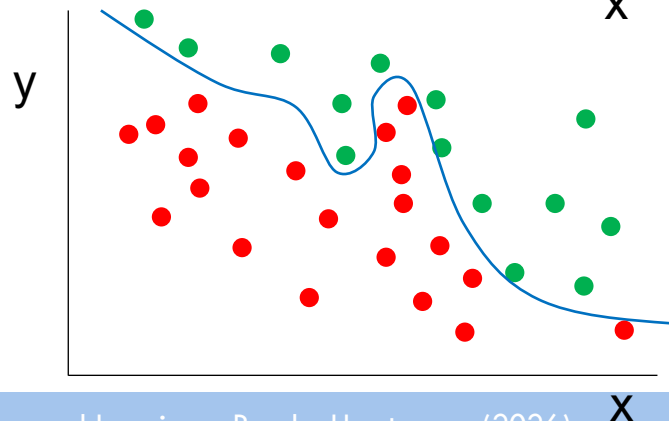
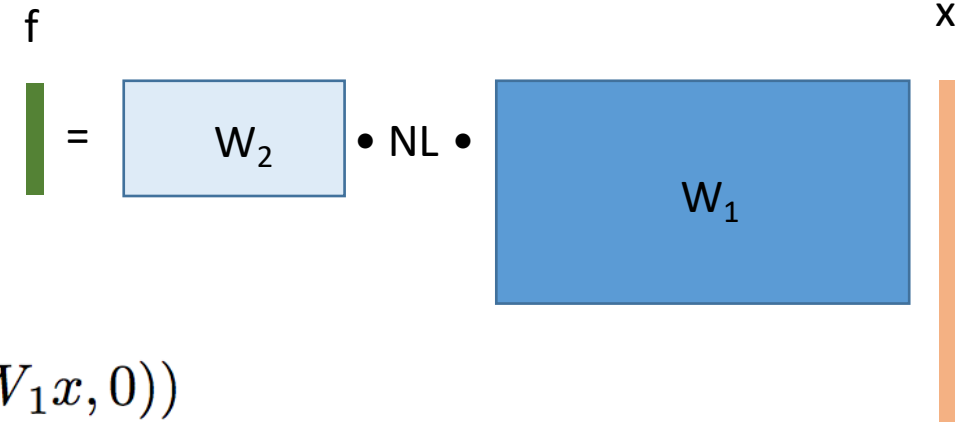
$$f = W_1 x$$

Learned f : not flexible



$$f = W_2 \max(W_1 x, 0)$$

Learned f : a bit flexible



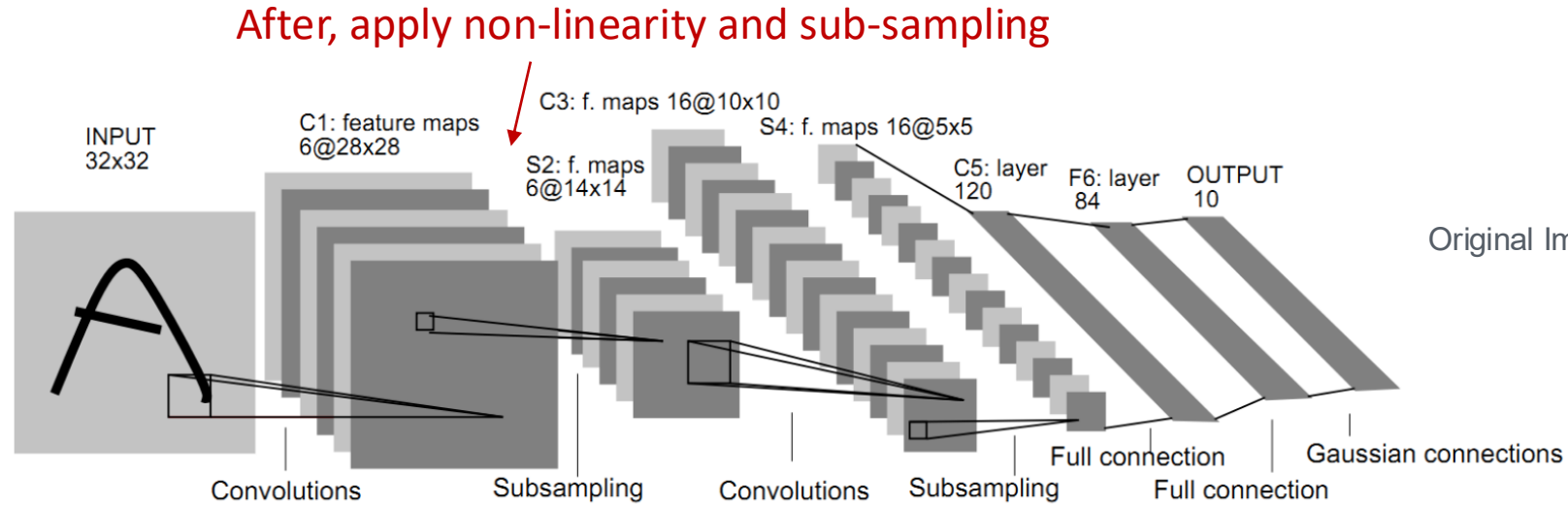
$$f = W_3 \max(0, W_2 \max(W_1 x, 0))$$

Learned f : more flexible

Does it generalize???

↓
We can keep adding these “layers”...

Getting us to Convolutional Neural Networks



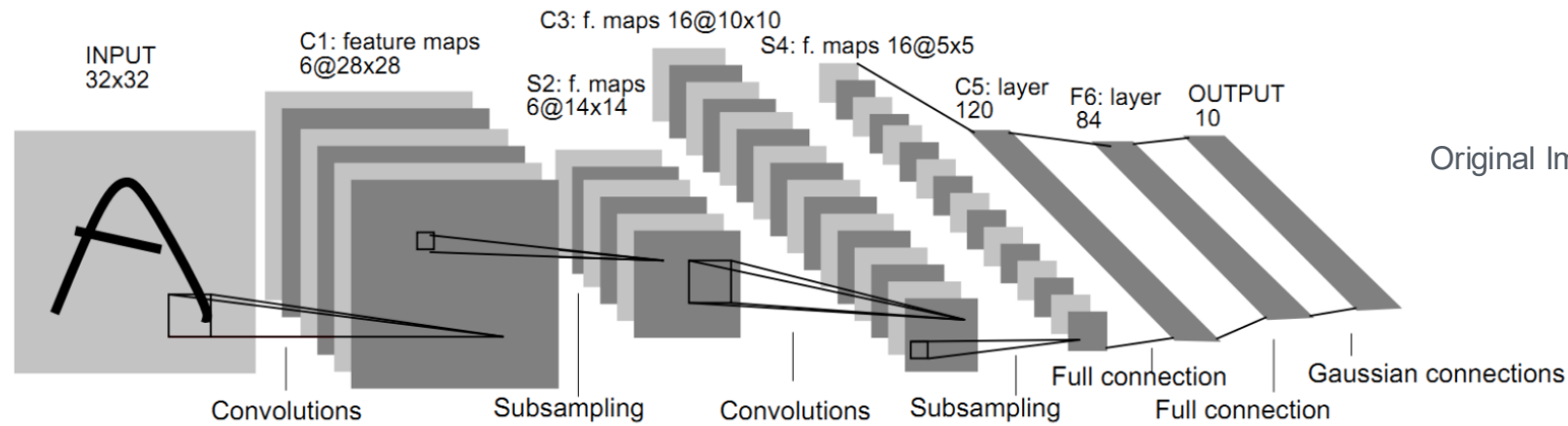
Original Image published in [LeCun et al., 1998]

Each matrix W is a convolution matrix

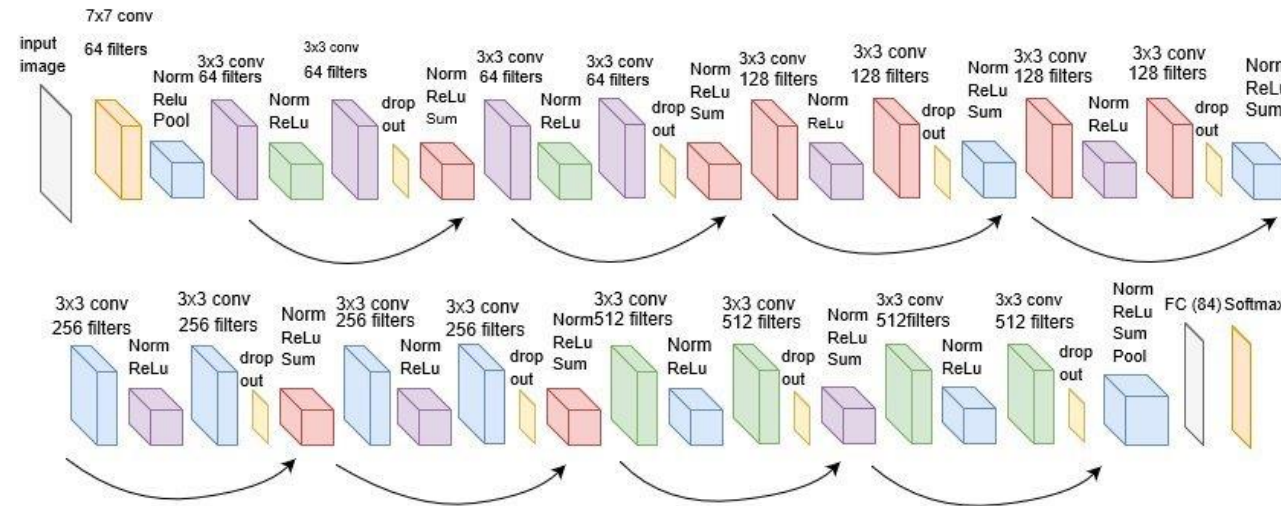
Repeat a few times

At the end, use a full W for a final matrix multiplication

Getting us to Convolutional Neural Networks

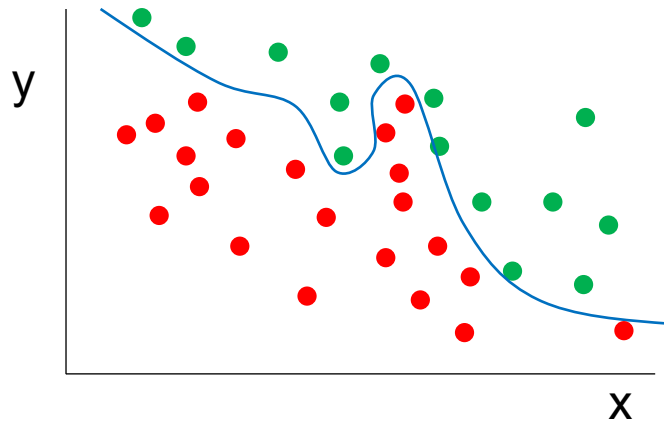


In practice, this process is repeated many times:



Aside #1 before convolutional neural network details

Q: Can we try to avoid making these learning models too complicated?

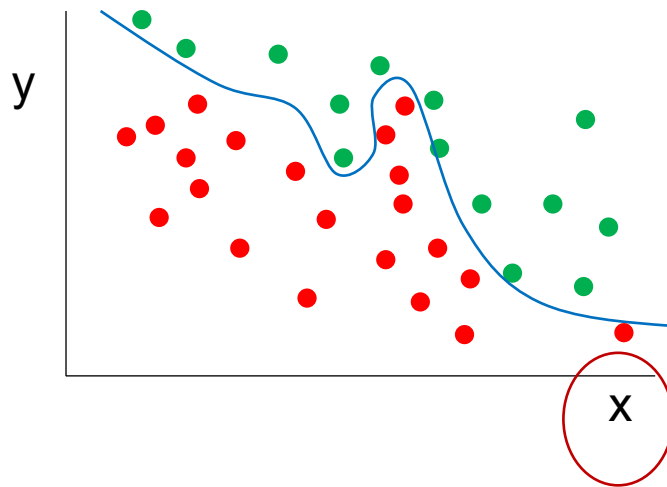


Learned f : more flexible

Does it generalize???

Aside #1 before convolutional neural network details

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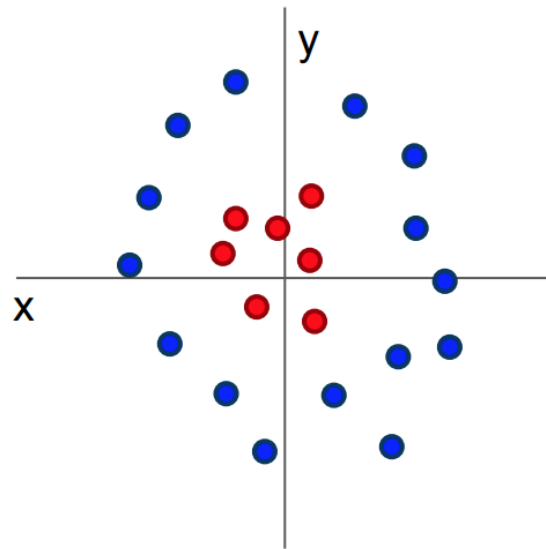


Learned f : more flexible

Does it generalize???

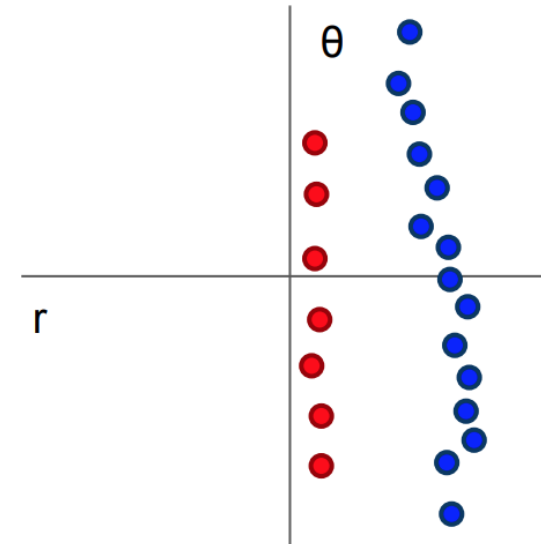
A: Yes, by transforming the data coordinates *before* classification

Image Features: Motivation



Cannot separate red and blue points with linear classifier

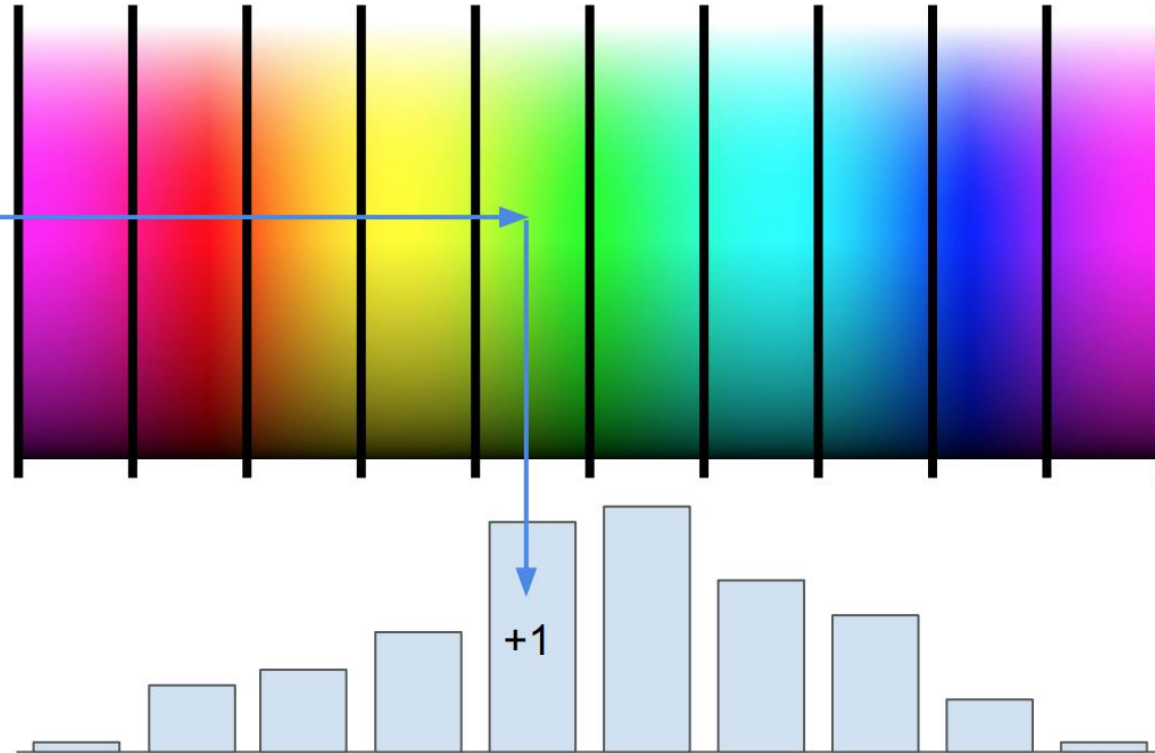
$$f(x, y) = (r(x, y), \theta(x, y))$$



After applying feature transform, points can be separated by linear classifier

From Stanford CS231: <http://cs231n.stanford.edu/>

Example: Color Histogram

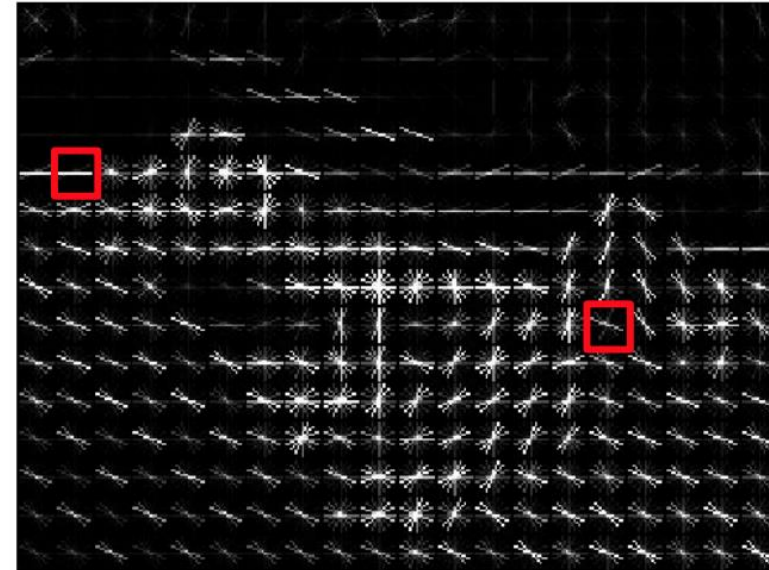


From Stanford CS231: <http://cs231n.stanford.edu/>

Example: Histogram of Oriented Gradients (HoG)



Divide image into 8x8 pixel regions
Within each region quantize edge
direction into 9 bins



Example: 320x240 image gets divided
into 40x30 bins; in each bin there are
9 numbers so feature vector has
 $30 \times 40 \times 9 = 10,800$ numbers

Lowe, "Object recognition from local scale-invariant features", ICCV 1999
Dalal and Triggs, "Histograms of oriented gradients for human detection," CVPR 2005

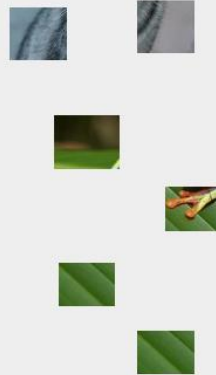
From Stanford CS231: <http://cs231n.stanford.edu/>

Example: Bag of Words

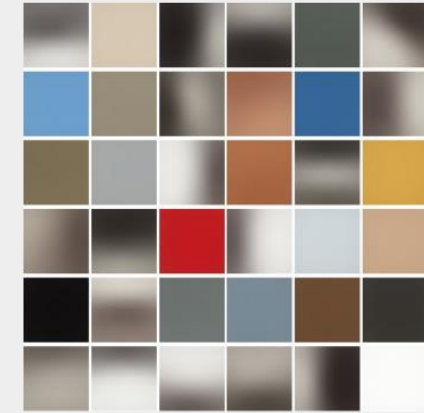
Step 1: Build codebook



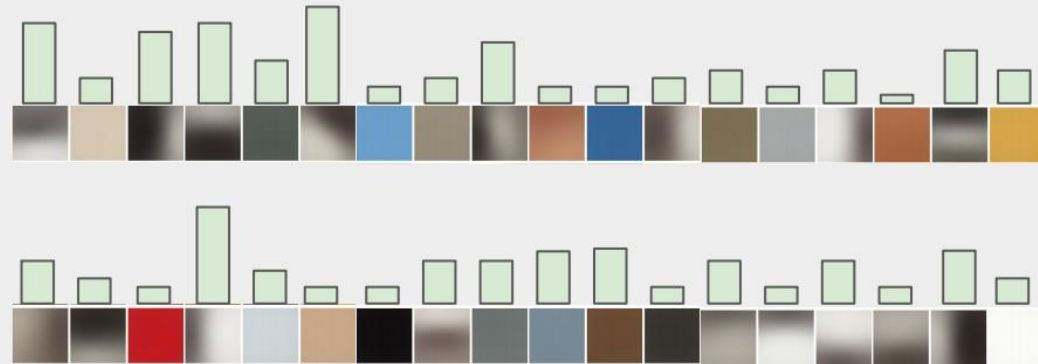
Extract random patches



Cluster patches to form "codebook" of "visual words"



Step 2: Encode images

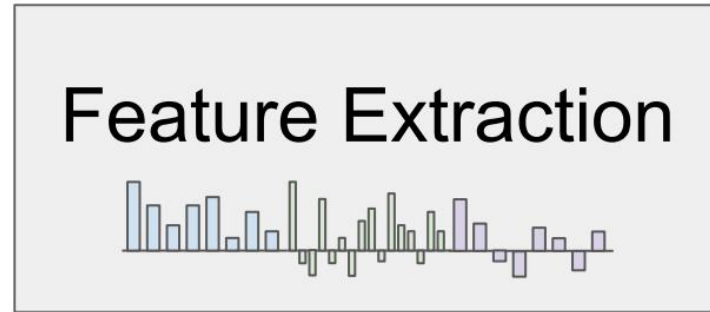


Fei-Fei and Perona, "A bayesian hierarchical model for learning natural scene categories", CVPR 2005

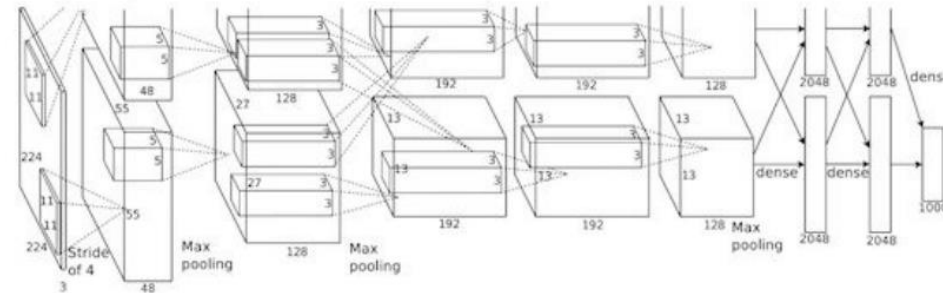
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Image features vs ConvNets

“Hand crafted”



10 numbers giving scores for classes



Krizhevsky, Sutskever, and Hinton, “Imagenet classification with deep convolutional neural networks”, NIPS 2012.
Figure copyright Krizhevsky, Sutskever, and Hinton, 2012.
Reproduced with permission.

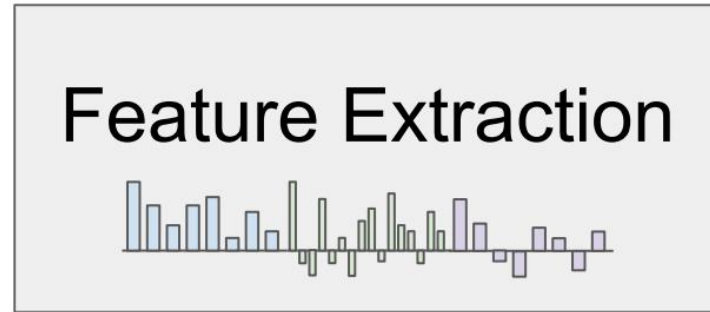
10 numbers giving scores for classes



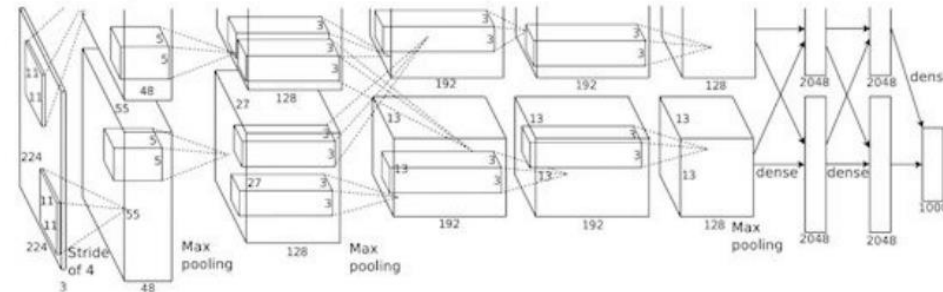
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Krizhevsky, Sutskever, and Hinton, “Imagenet classification with deep convolutional neural networks”, NIPS 2012.
Figure copyright Krizhevsky, Sutskever, and Hinton, 2012.
Reproduced with permission.

10 numbers giving scores for classes



training

History has now proven – bottom approach works better!

From Stanford CS231: <http://cs231n.stanford.edu/>

Statistical Machine Learning in ~30 minutes

Two competing goals in machine learning:

1. Can we make sure the in-sample error $L_{\text{in}}(y, f(x, W))$ is small enough during network training?
 - Appropriate cost function
 - “complex enough” model

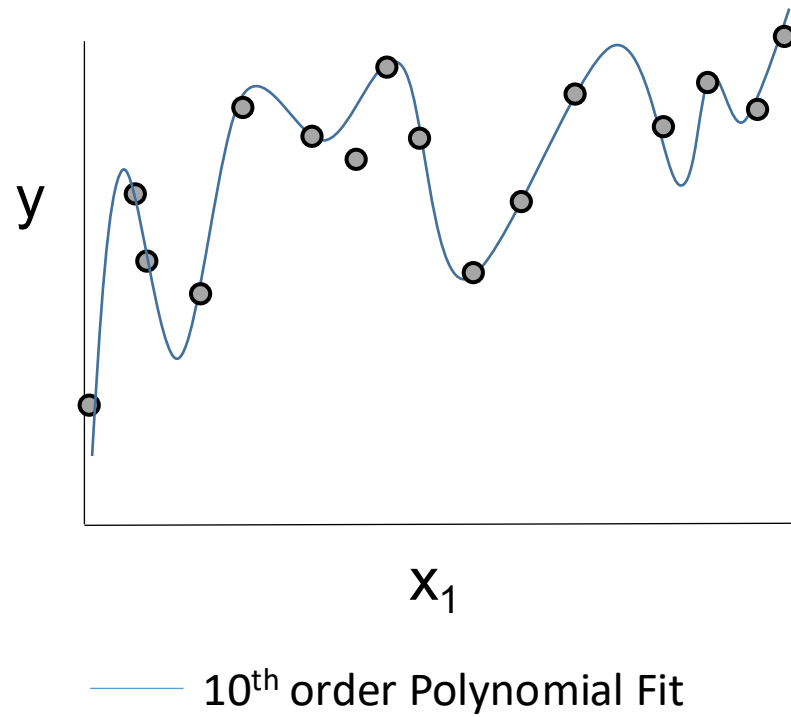
Statistical Machine Learning in ~30 minutes

Two competing goals in machine learning:

1. Can we make sure the in-sample error $L_{\text{in}}(y, f(x, W))$ is small enough during network training?
 - Appropriate cost function
 - “complex enough” model
2. Can we make sure that $L_{\text{out}}(y, f(x, W))$ is close enough to $L_{\text{in}}(y, f(x, W))$ during network testing?
 - Probabilistic analysis says yes!
 - $|L_{\text{in}} - L_{\text{out}}|$ bounded from above
 - Bound grows with model capacity (i.e., complexity - bad)
 - Bound shrinks with # of training examples (good)

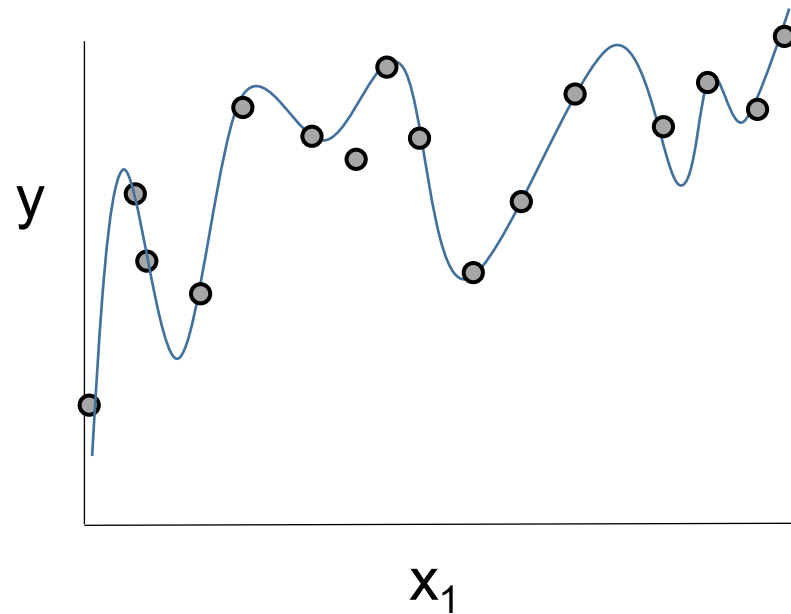
Model overfitting versus underfitting – a thought exercise

Let's fit these “training” data points:



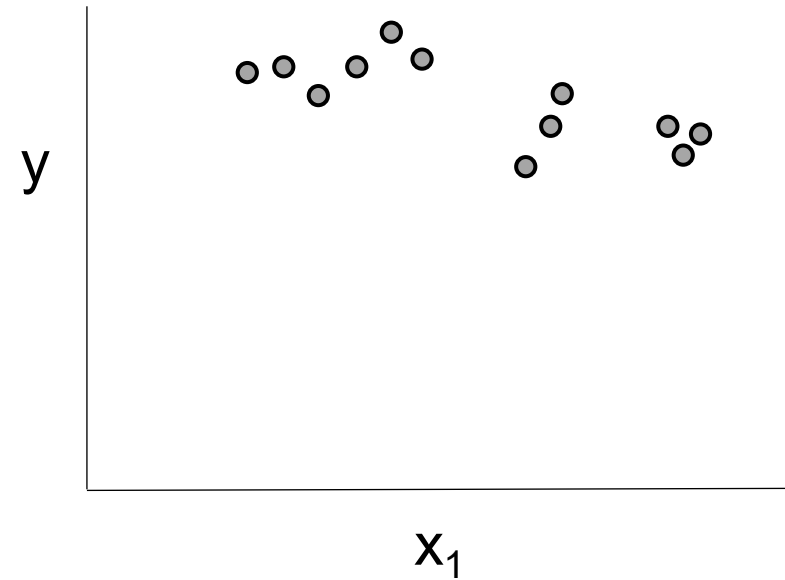
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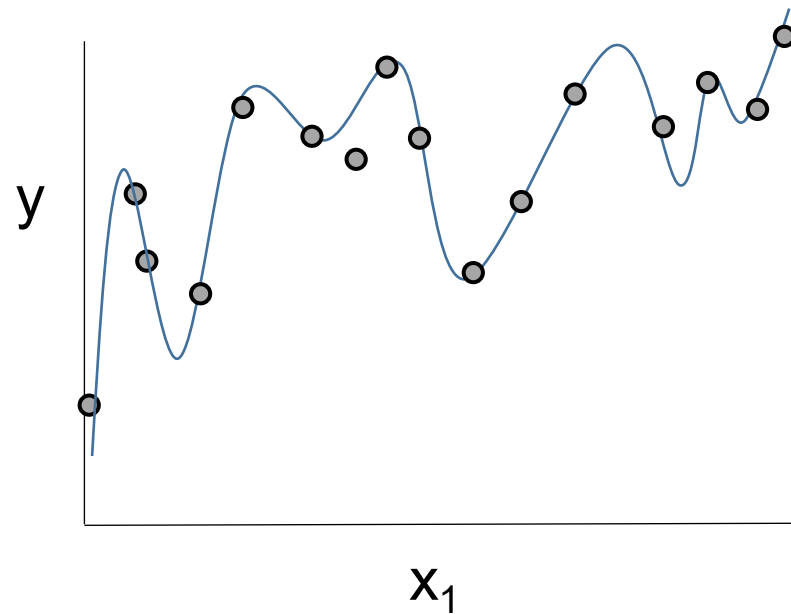
— 10th order Polynomial Fit

And then here's our testing dataset – good?



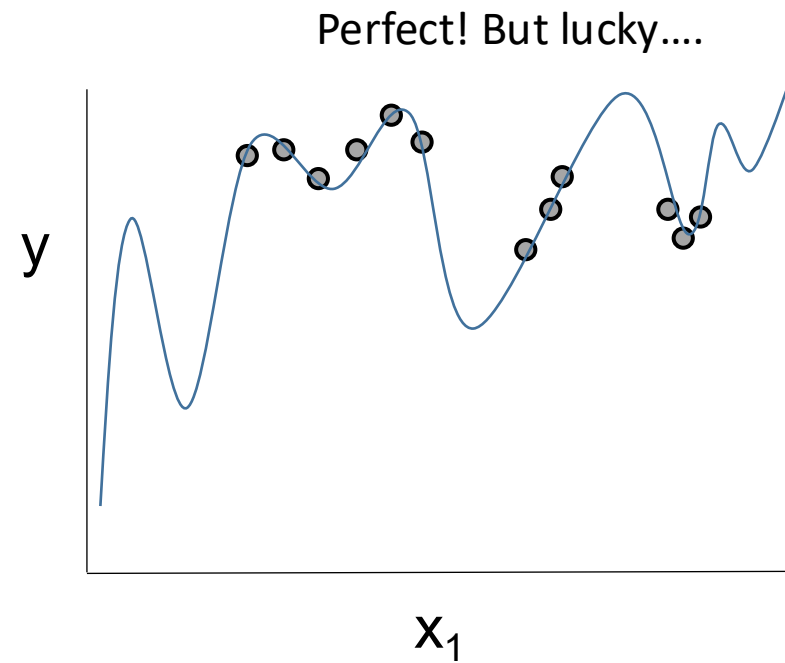
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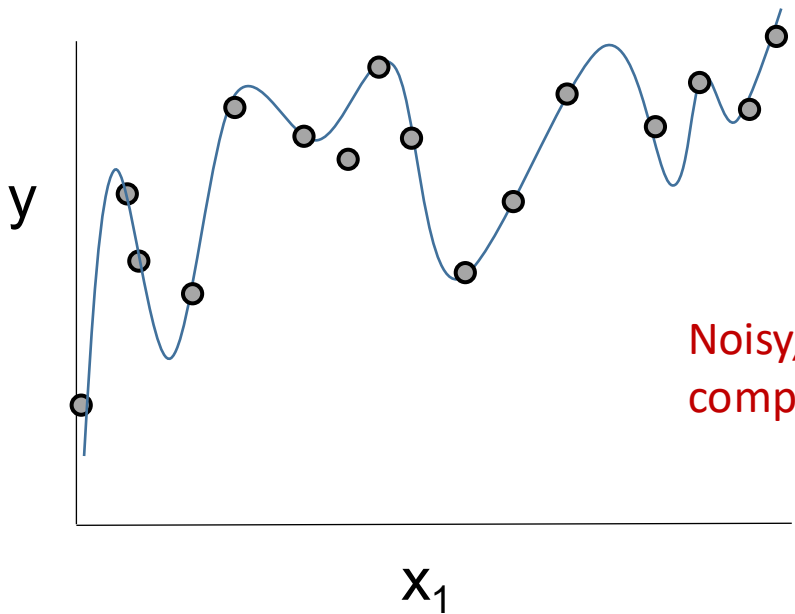
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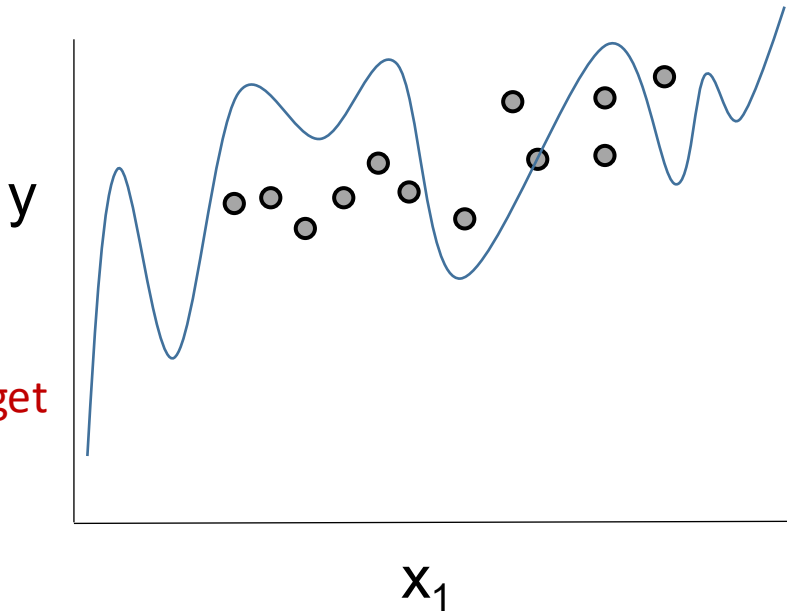
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— 10th order Polynomial Fit

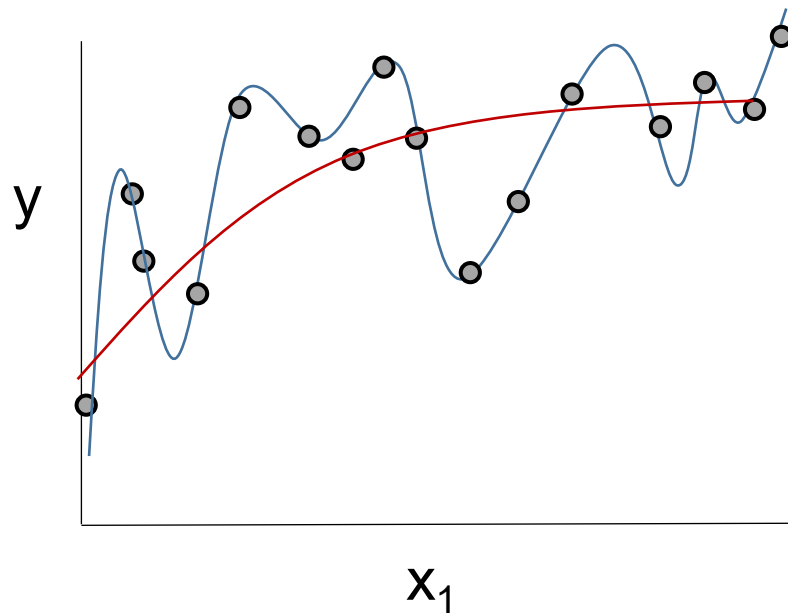
What if our test dataset was this :



↔
Noisy, low
complexity target

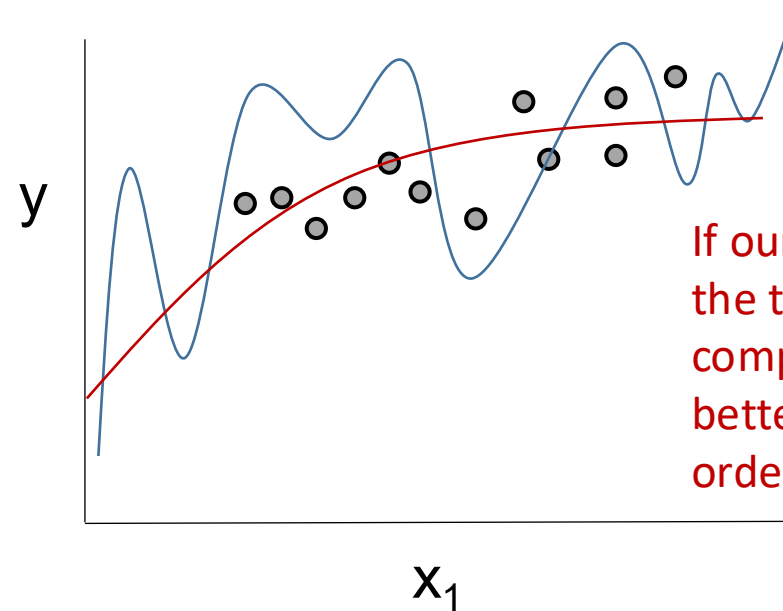
Model overfitting versus underfitting – a thought exercise

Let's fit these "training" data points:



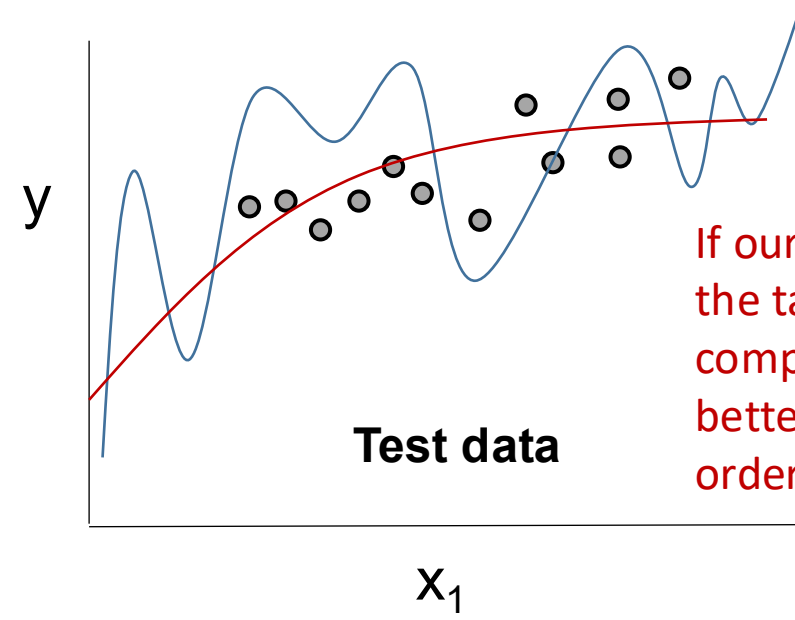
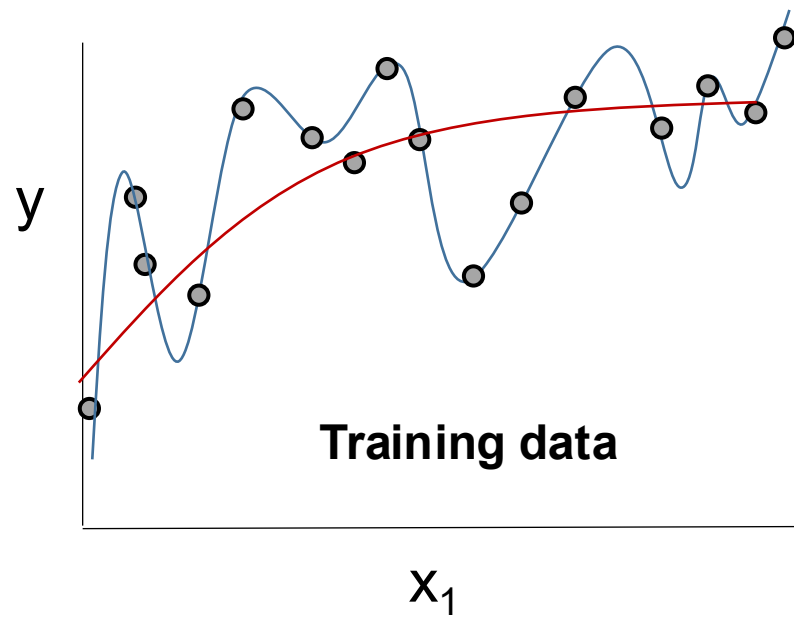
- 10th order Polynomial Fit
- 2nd order Polynomial Fit

What if our test dataset was this :



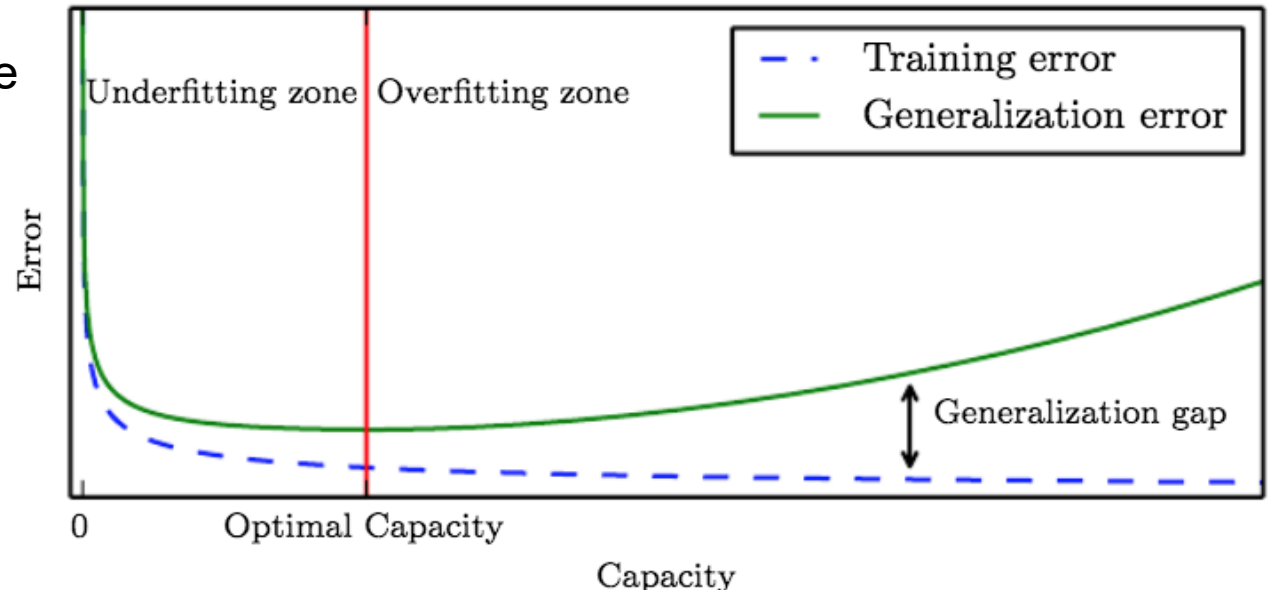
If our data was noisy and the target followed a low-complexity model, we'd be better off with a second order fit!

Model overfitting versus underfitting – a thought exercise



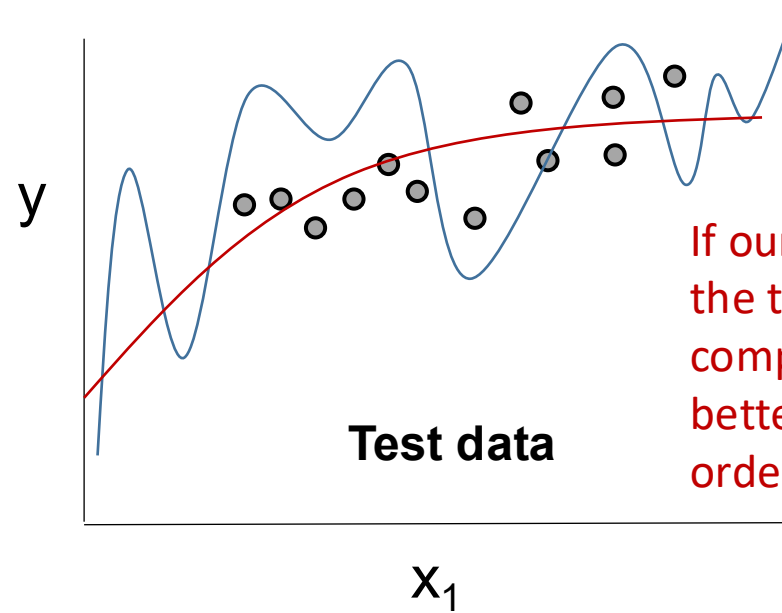
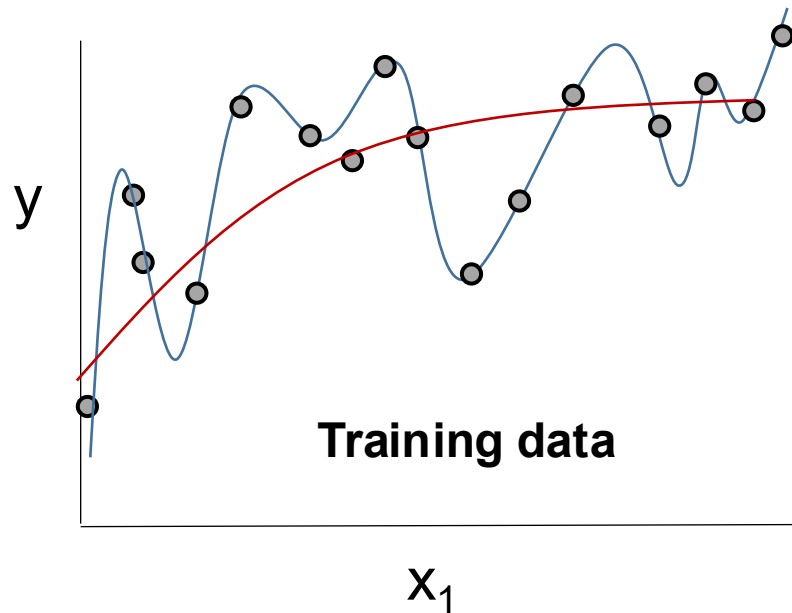
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Model capacity: ability to fit a wide range of functions



Deep Learning, I. Goodfellow et al., Fig. 5.3

Model overfitting versus underfitting – a thought exercise

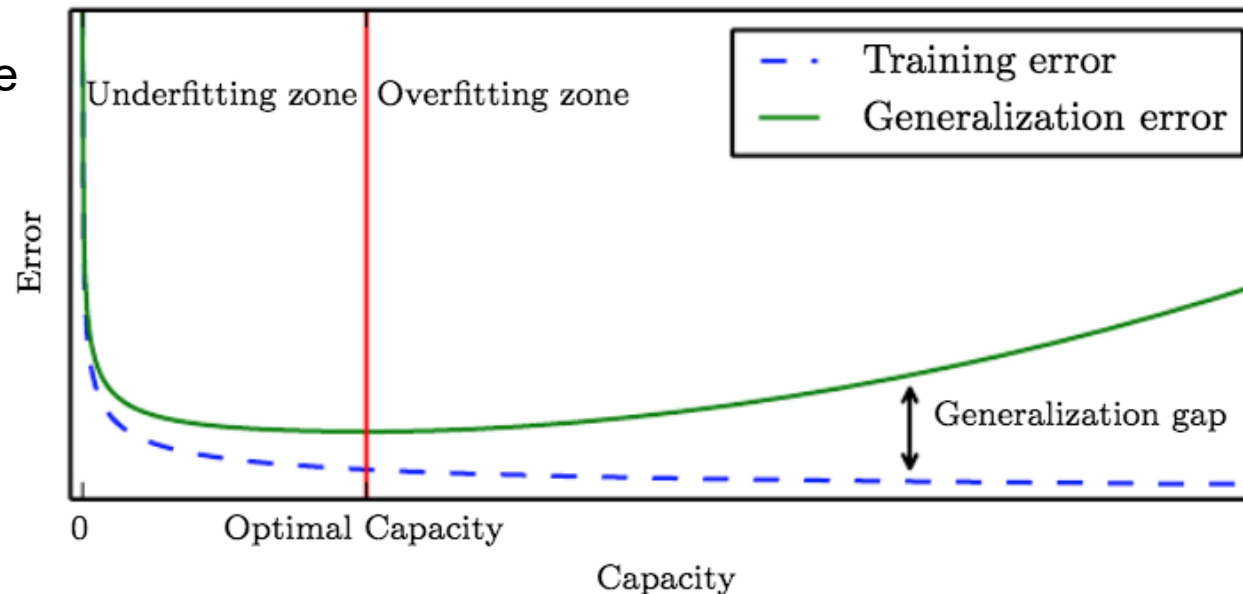


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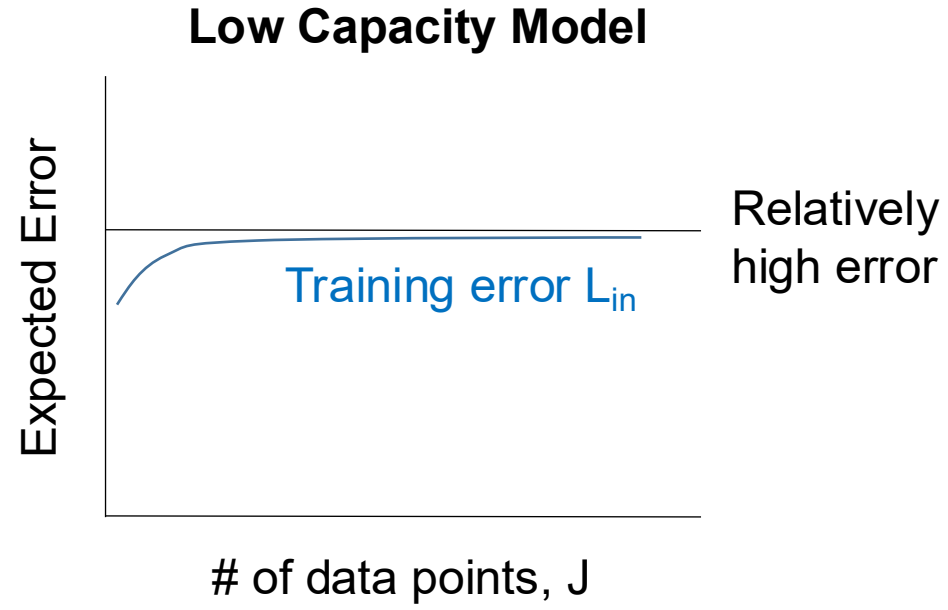
Model capacity: ability to fit a wide range of functions

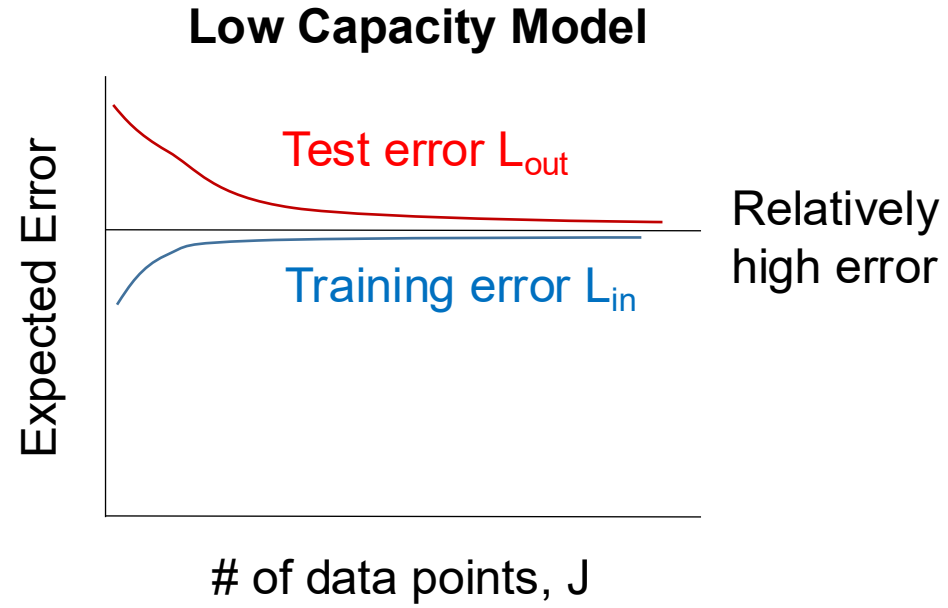
Control capacity through model's hypothesis space (set of functions model can take)

Hard to know ahead of time!

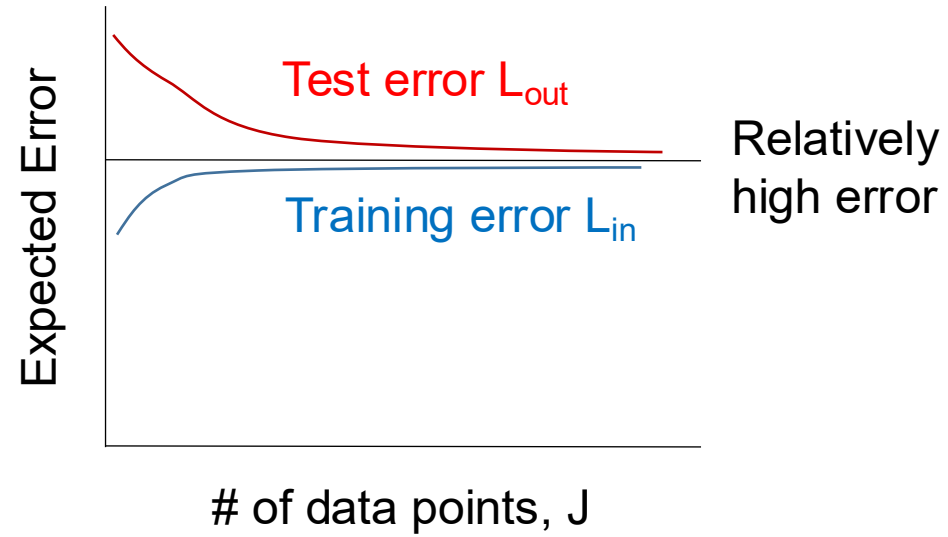


Deep Learning, I. Goodfellow et al., Fig. 5.3

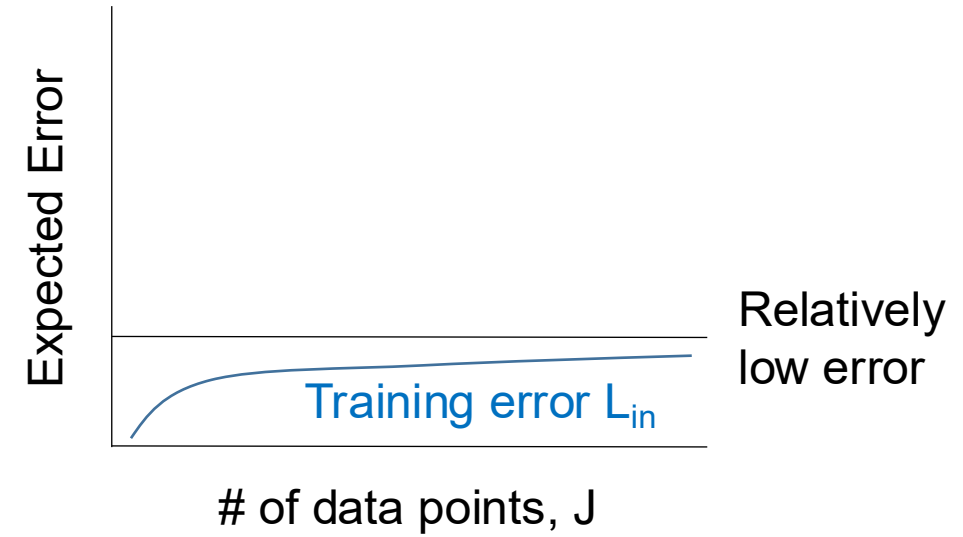




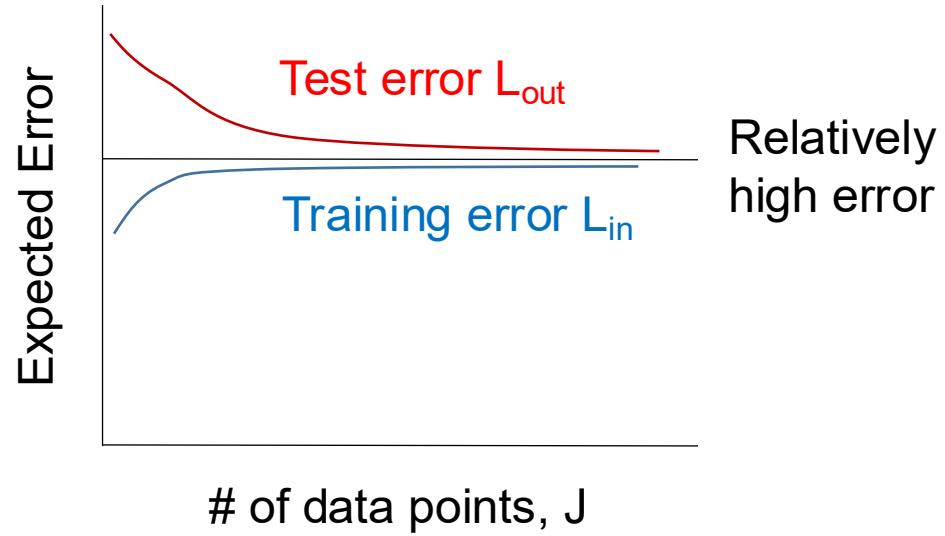
Low Capacity (complexity) Model



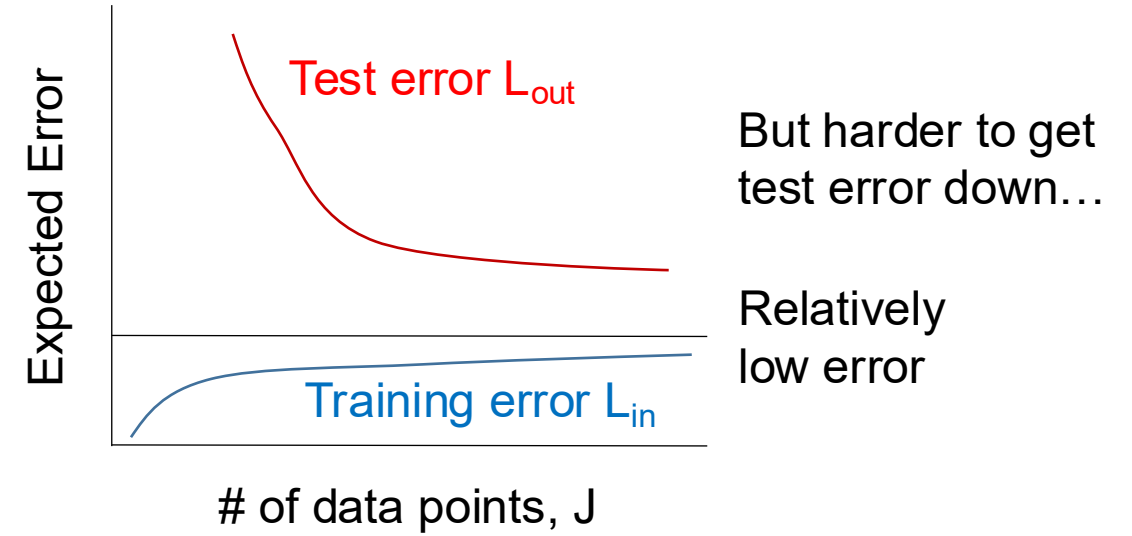
High Capacity (complexity) Model



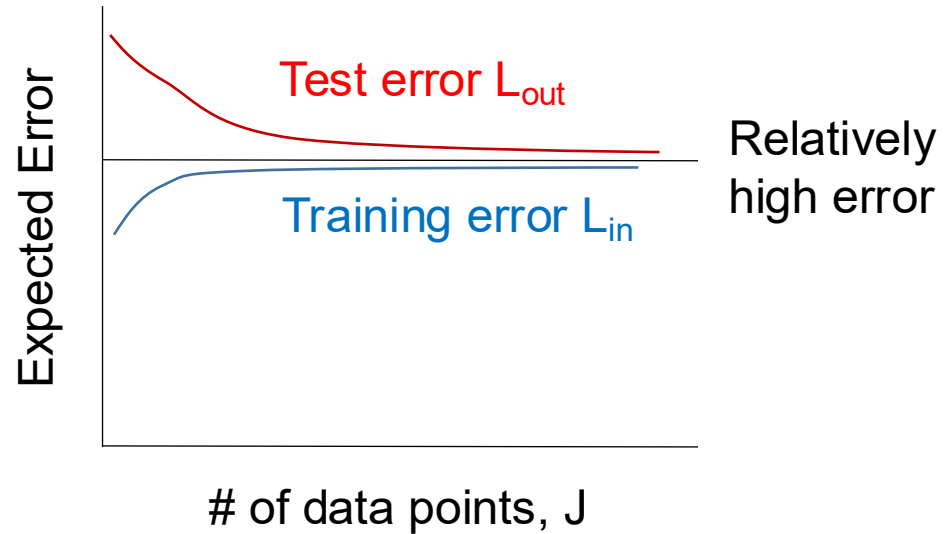
Low Capacity (complexity) Model



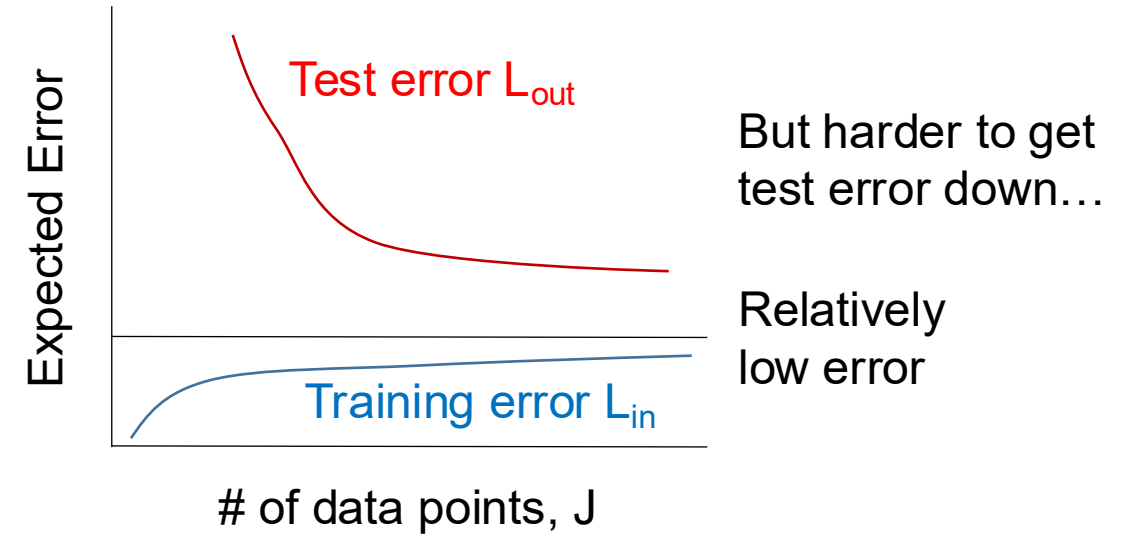
High Capacity (complexity) Model



Low Capacity (complexity) Model



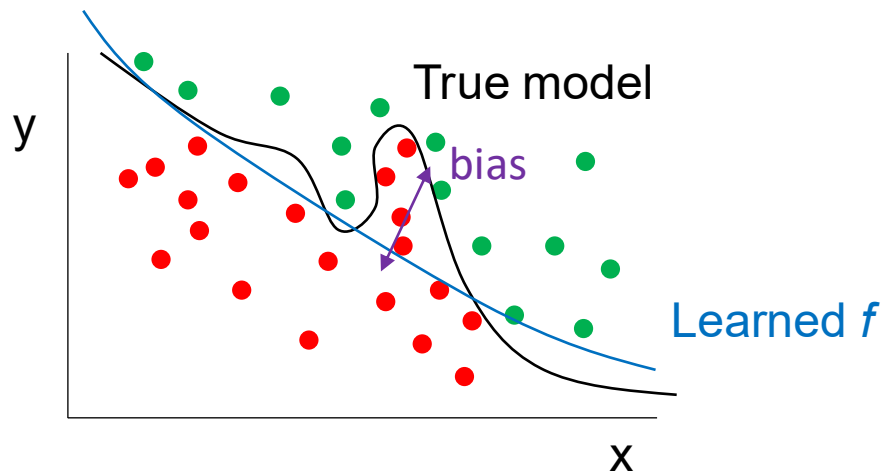
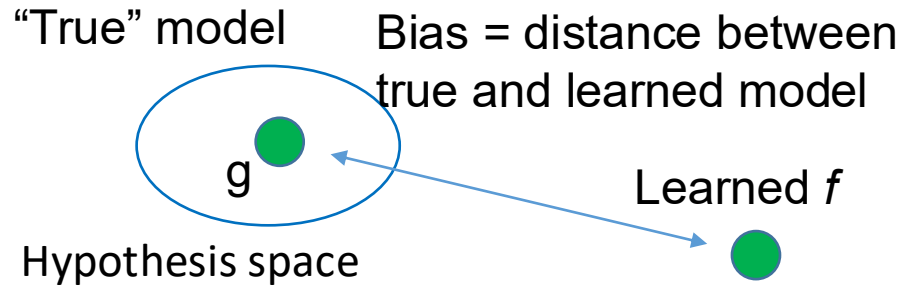
High Capacity (complexity) Model



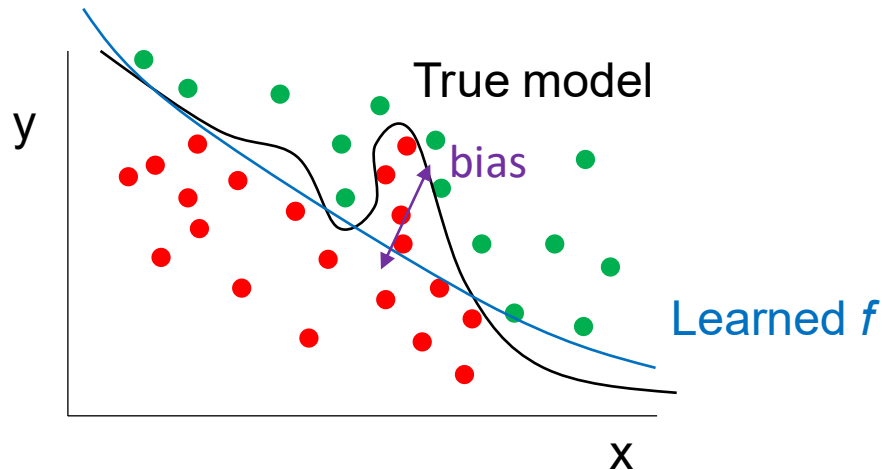
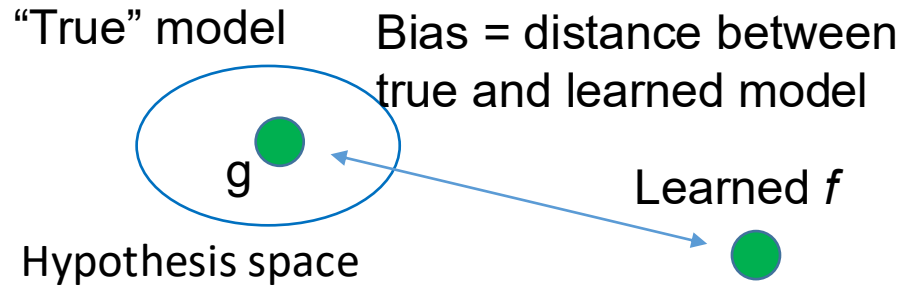
Take away concepts:

- Can't ever really expect test error to be less than training error
- Complicated models tend to appear to “do better” during training, before trying test data
- When the model gets complicated and you don't have enough data, challenging to get test error down

Model bias versus variance



Model bias versus variance

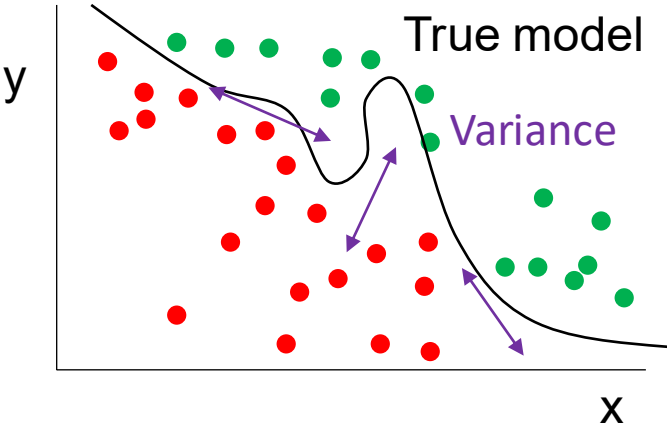
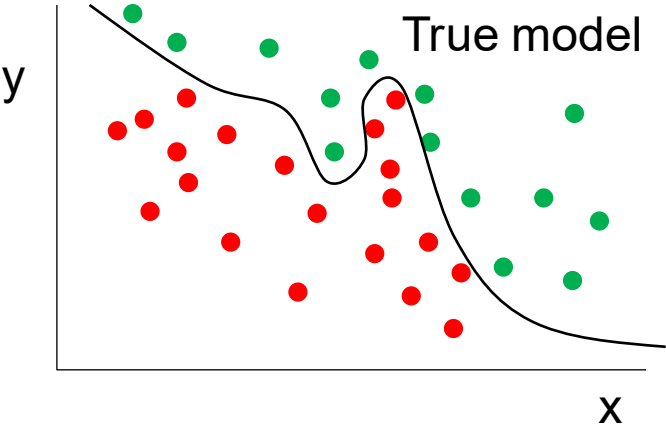
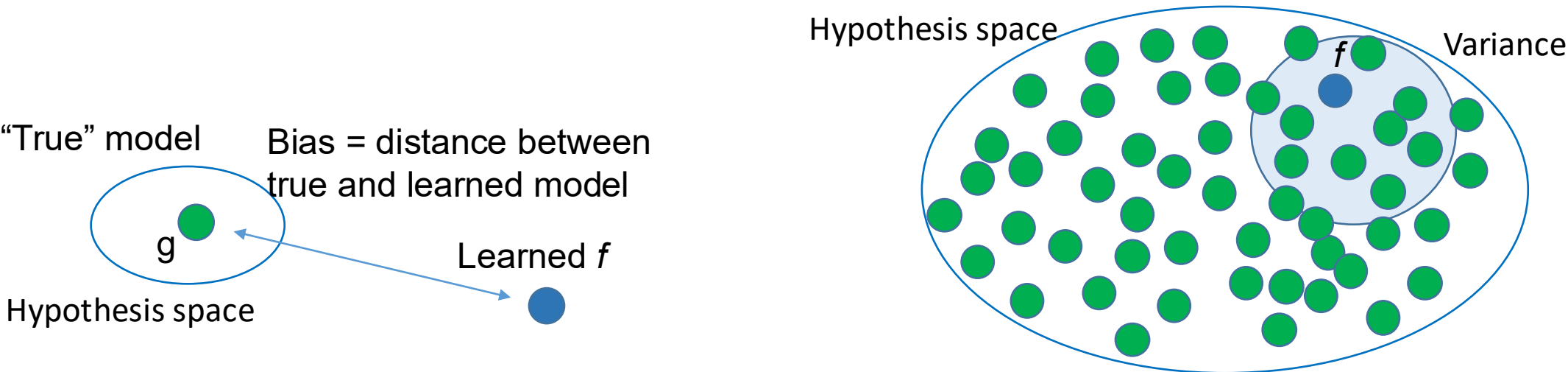


Models that tend to be “a bit too simple” are biased away from “true” model

$$\text{Bias} = (g(\mathbf{x}) - f(\mathbf{x}))^2$$

Measures how far our learning model f is biased away from target function g (for perfect training data classification)

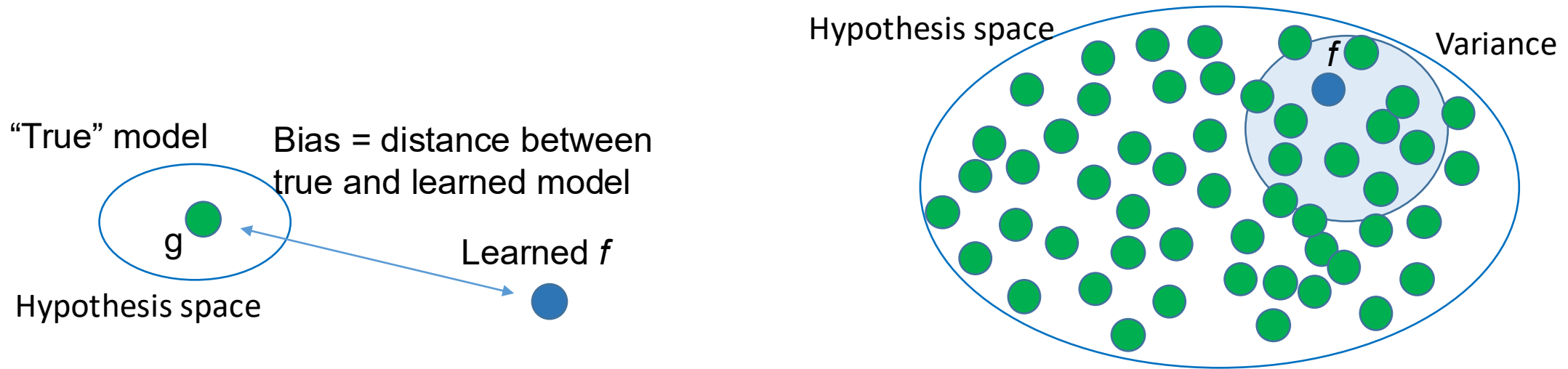
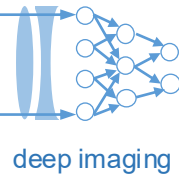
Model bias versus variance



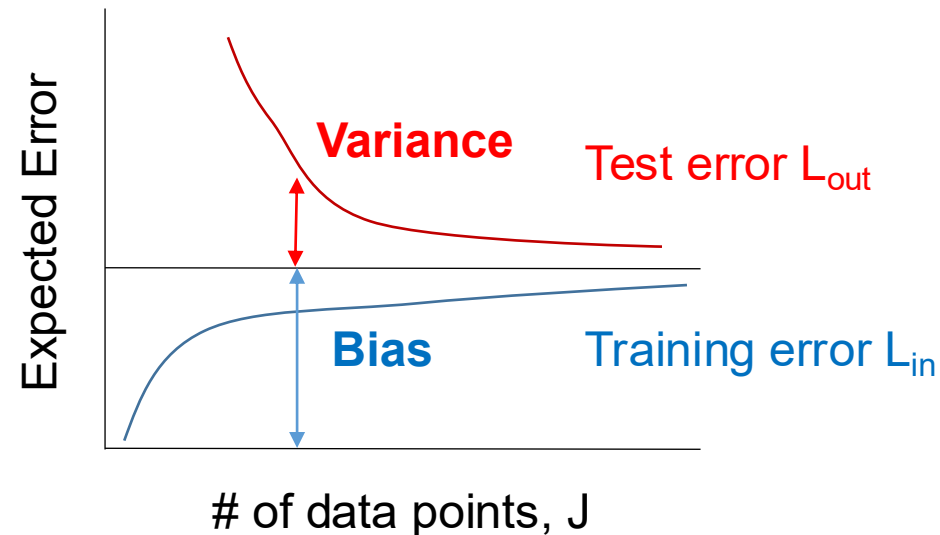
$$\text{Variance} = \text{Var}[g(\mathbf{x})]$$

More complicated datasets exhibit lots of variance between ideal boundary for training and testing

Model bias versus variance



Error vs. # data points



Test Error is sum of model bias and variance!

Goal is to find a model f that balances between these two quantities for a given dataset

How to formally define capacity and complexity?

- Short answer: it's complicated...
- Related to something called the *VC Dimension*
 - Can provide theoretical bounds on performance
 - Dimensional bounds rather than scalar bounds...
- I decided not to go into it, but please do take a look at the following lecture material to learn more!

Learning From Data (Caltech, Prof. Y Abu-Mostafa)

<https://www.youtube.com/watch?v=Dc0sr0kdBVI#t=3m24s>

Conclusions from statistical machine learning

- Conclusion: you want a model that is complex enough to capture variations within high-dimensional space, but not too complex such that it overfits the data
- Want a model with a high capacity, but can still *generalize* to data outside training set
 - More data -> less overfitting, complex target -> more overfitting
- For simple models, we can measure complexity via degrees of freedom, the VC bound and so-on to help us nail down ideal models that can generalize well

Conclusions from statistical machine learning

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- For simple models, we can measure complexity via degrees of freedom, the VC bound and so-on to help us nail down ideal models that can generalize well
- **For DL models:** this will get too hard...here's a few counter-intuitive properties:
 1. A fixed DL *architecture* exhibits data-dependent complexities
 - e.g., “good” DL networks achieve 0 training error on images with random labels, so cannot generalize at all in this case, and are too complex
 2. DL networks with more hidden units leads to *better* generalization (the main finding of the last few years). So deeper models tend to be less complex, actually...
 3. Complexity depends upon loss function and optimization method...

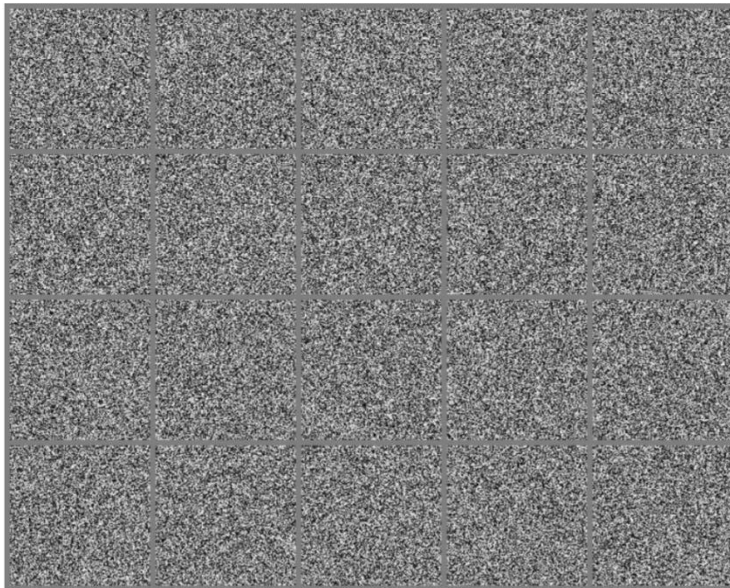
Important to remember: “No Free Lunch Theorem”

- *“Averaged over all possible data-generating distributions, every classification algorithm has the same error rate when classifying previously unobserved points.”*
- The most sophisticated DL algorithm has same average performance (averaged over all possible tasks) as the simplest.

Important to remember: “No Free Lunch Theorem”

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- The most sophisticated DL algorithm has same average performance (averaged over all possible tasks) as the simplest.
- Must make assumptions about probability distributions of inputs we’ll encounter in real-world

Set of 20 “images”, random Gaussian distribution



Face at different orientations =
manifold n-D space

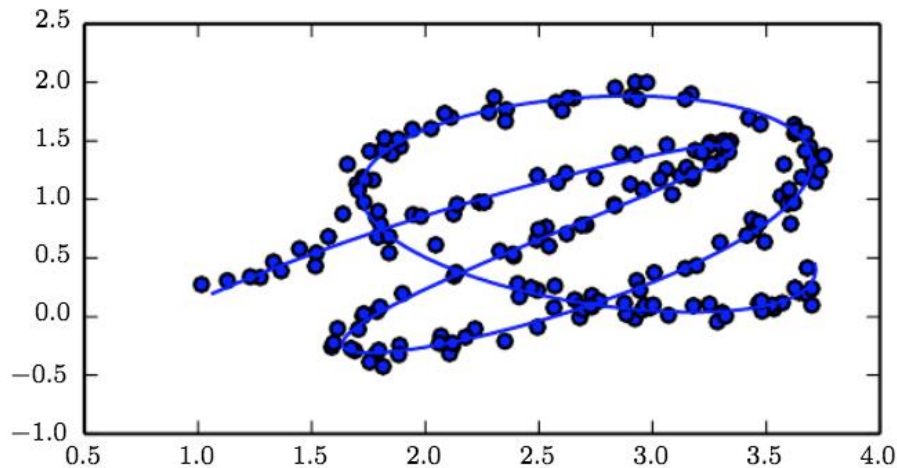


Deep Learning, I. Goodfellow et al., Fig. 5.12-13

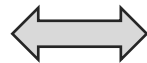
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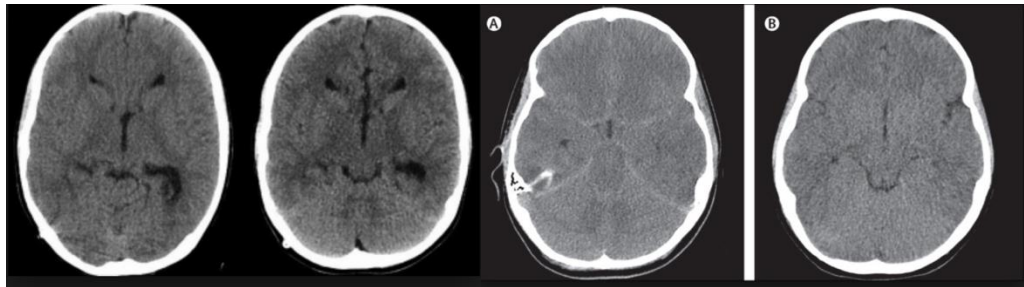
1D Manifold in 2D space



Manifold
Hypothesis

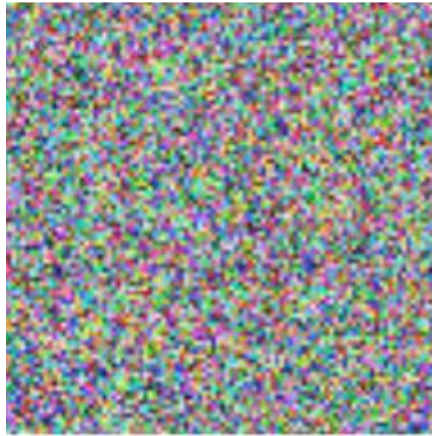


CT reconstructions of every brain in the world = kD manifold in nD space?



Deep Learning, I. Goodfellow et al., Fig. 5.11

Noise $\sim N(0,1)$



Generative
Model

