

CPSC 340: Machine Learning and Data Mining

Convolutional Neural Networks
(and miscellaneous deep learning tricks)

Spring 2022 (2021W2)

Admin

- A6 is out. Get it done early!
 - Due April 8th
- Bonus! 100% grade in class for anyone who attends today's class virtually!
 - Click here to redeem offer within 2 minutes:
 - <https://tinyurl.com/100percentGradeFor340>

Admin

- Plan for April 6th (next wed)
 - An experiment!
- We will watch two videos from my research past
 - With live questions!
- Deep Visualization Toolbox:
<https://www.youtube.com/watch?v=AgkfIQ4IGaM>
- Deep Learning Overview & Visualizing What Deep Neural Networks Learn
<https://www.youtube.com/watch?v=3lp9eN5JE2A>

But first

- A want to briefly revisit two things I flew through

Setting the Step-Size

- Stochastic gradient is **very sensitive to the step size** in deep models.
- **Bias step-size multiplier:** use bigger step-size for the bias variables.
- **Momentum** (stochastic version of “heavy-ball” algorithm):
 - Add term that moves in previous direction:

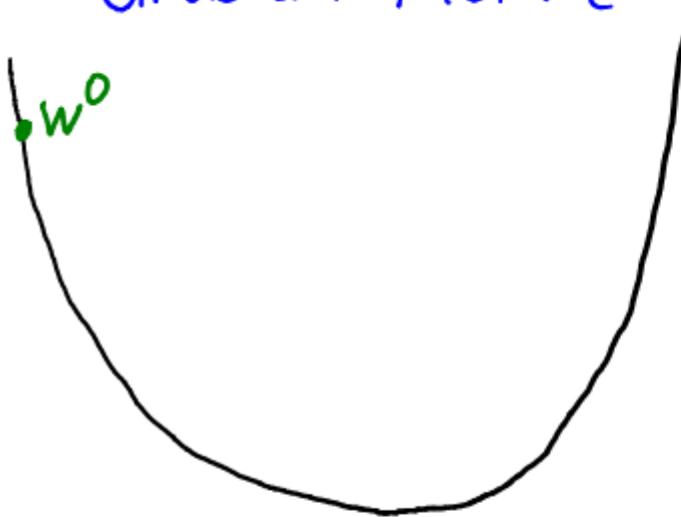
$$w^{t+1} = w^t - \alpha^t \nabla f_i(w^t) + \beta^t (w^t - w^{t-1})$$

Keep going in the old direction

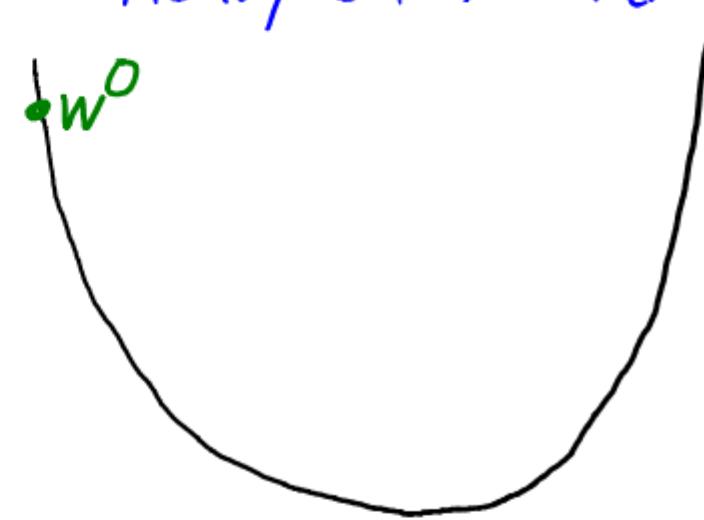
- Usually $\beta^t = 0.9$.

Gradient Descent vs. Heavy-Ball Method

Gradient Method

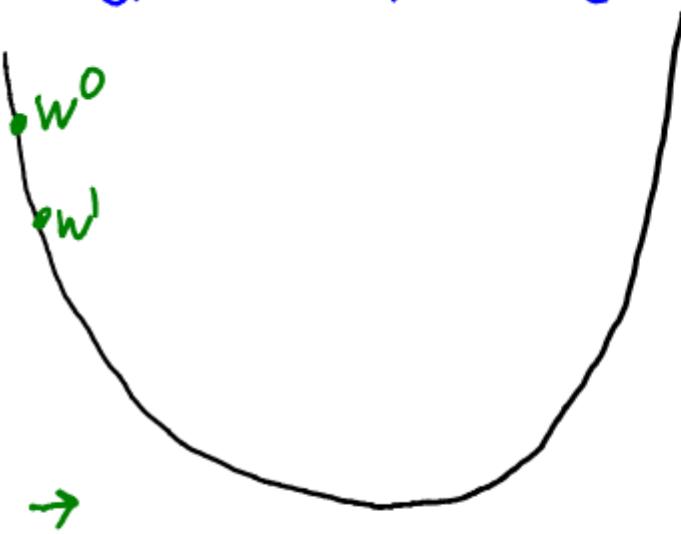


Heavy-ball Method

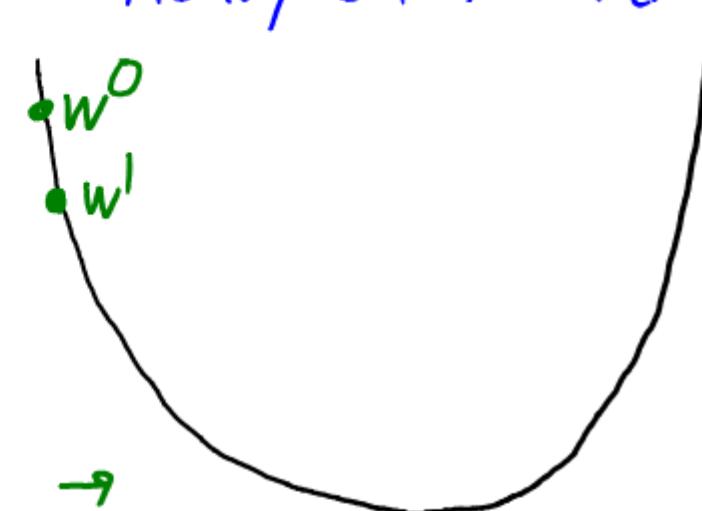


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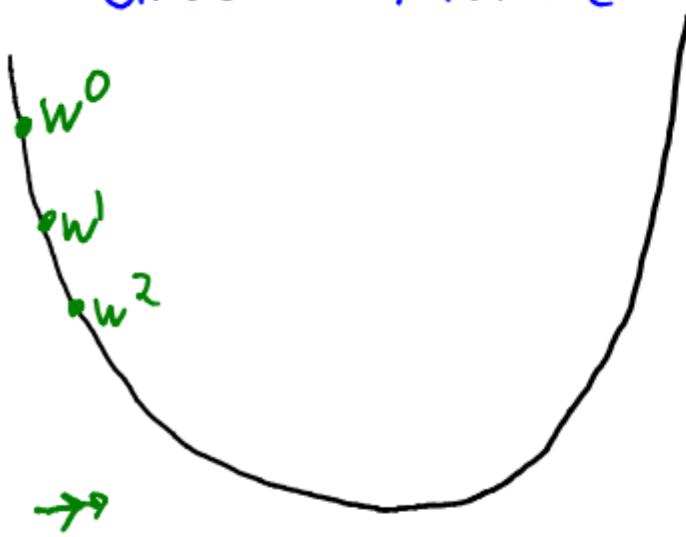


Heavy-ball Method

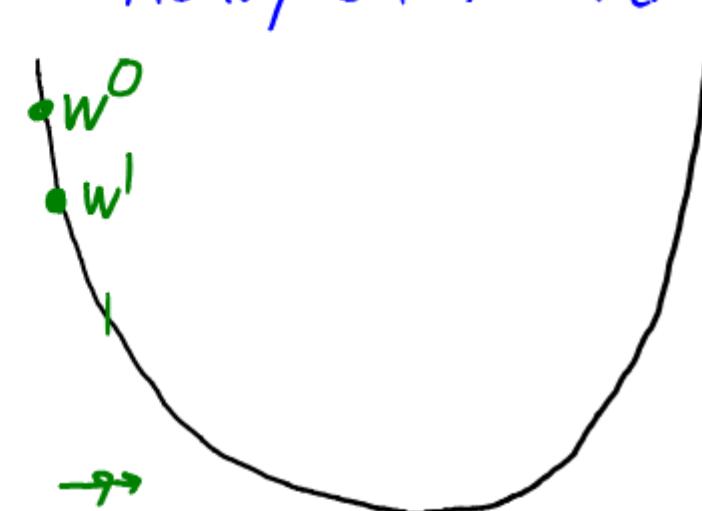


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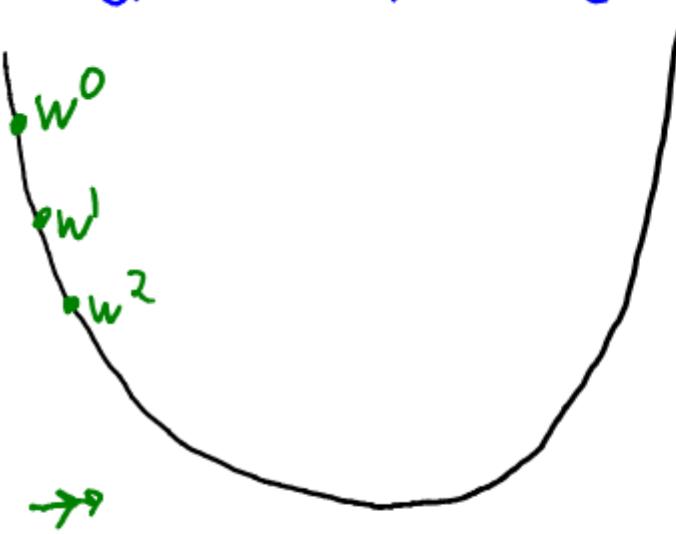


Heavy-ball Method

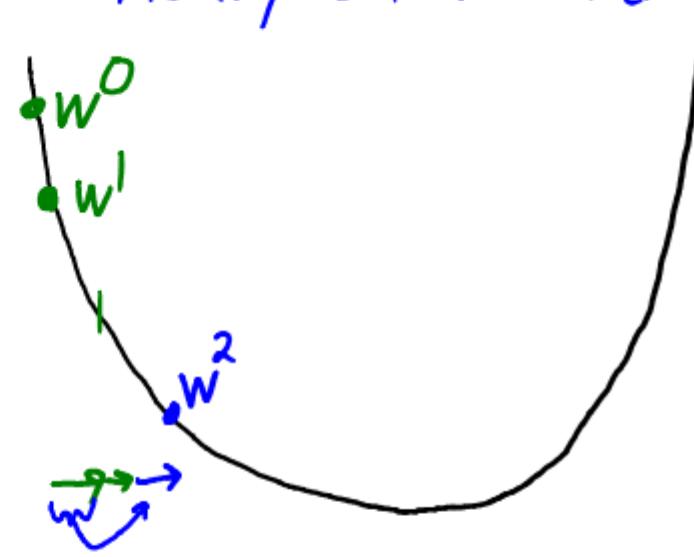


Gradient Descent vs. Heavy-Ball Method

Gradient Method



Heavy-ball Method

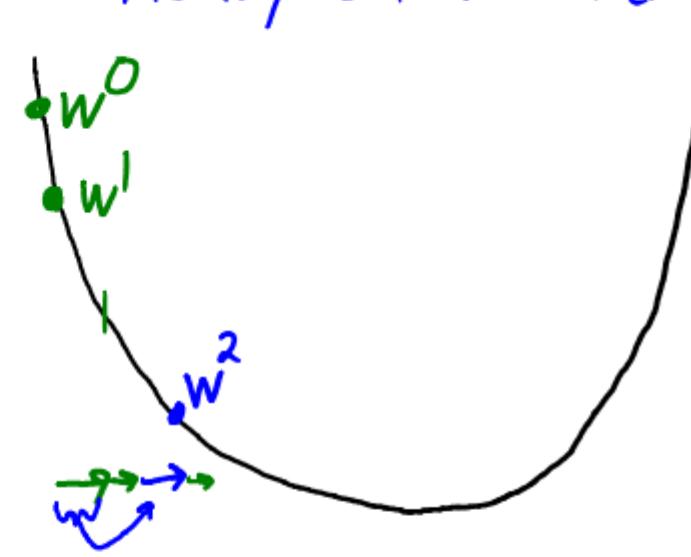


Gradient Descent vs. Heavy-Ball Method

Gradient Method



Heavy-ball Method

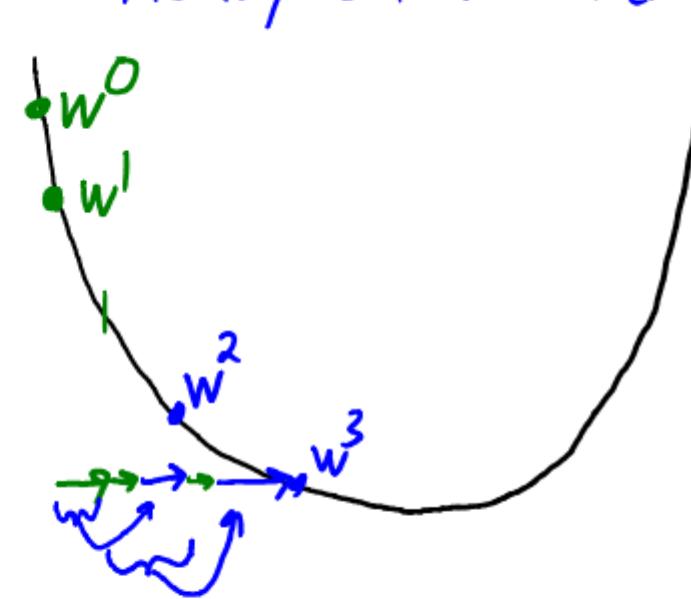


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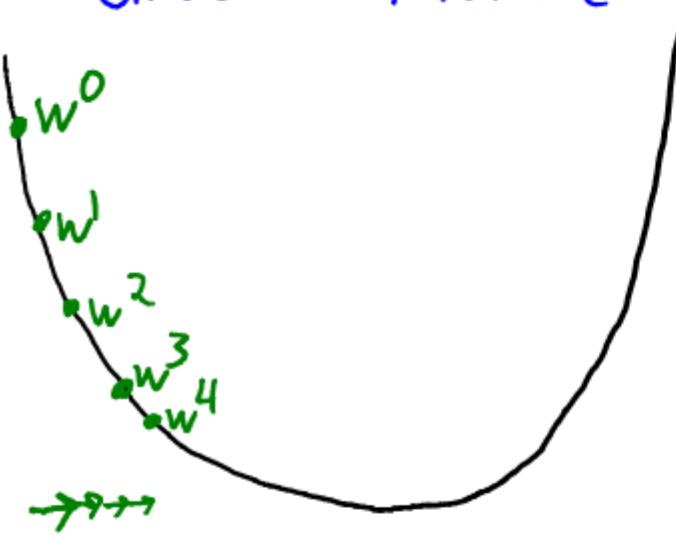


Heavy-ball Method

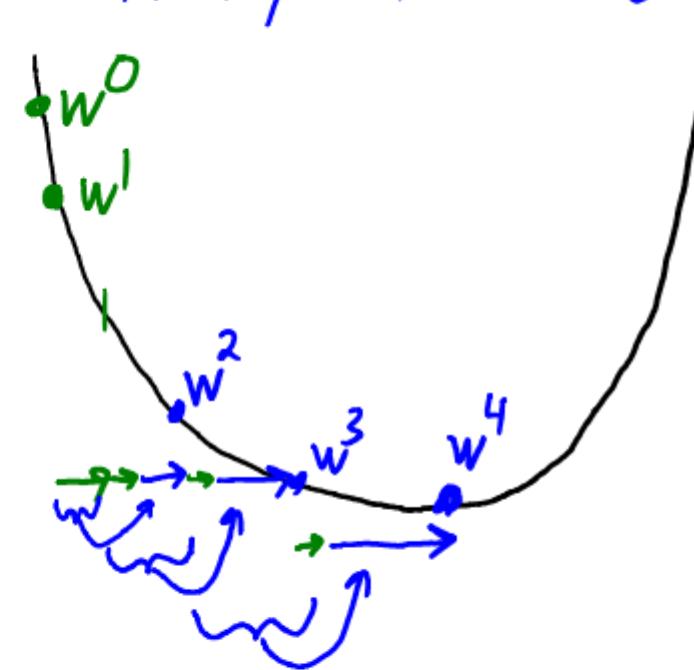


Gradient Descent vs. Heavy-Ball Method

Gradient Method

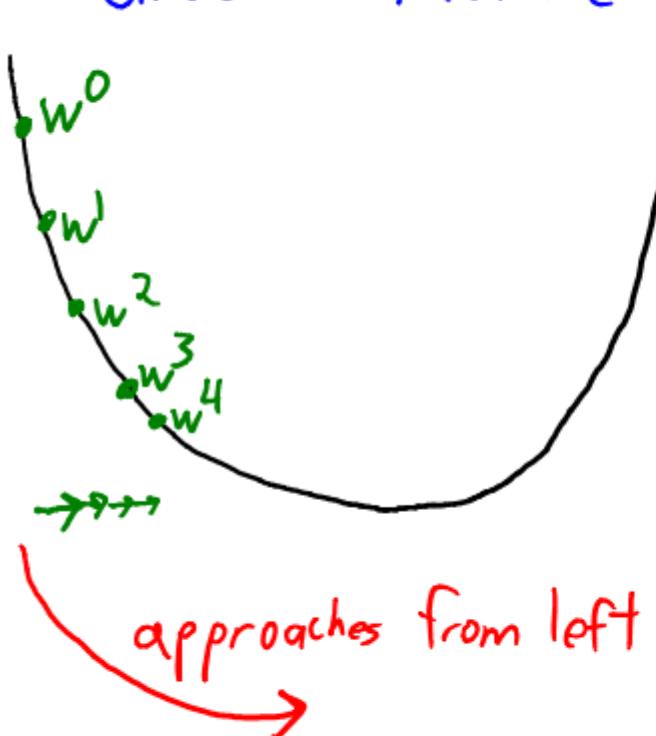


Heavy-ball Method



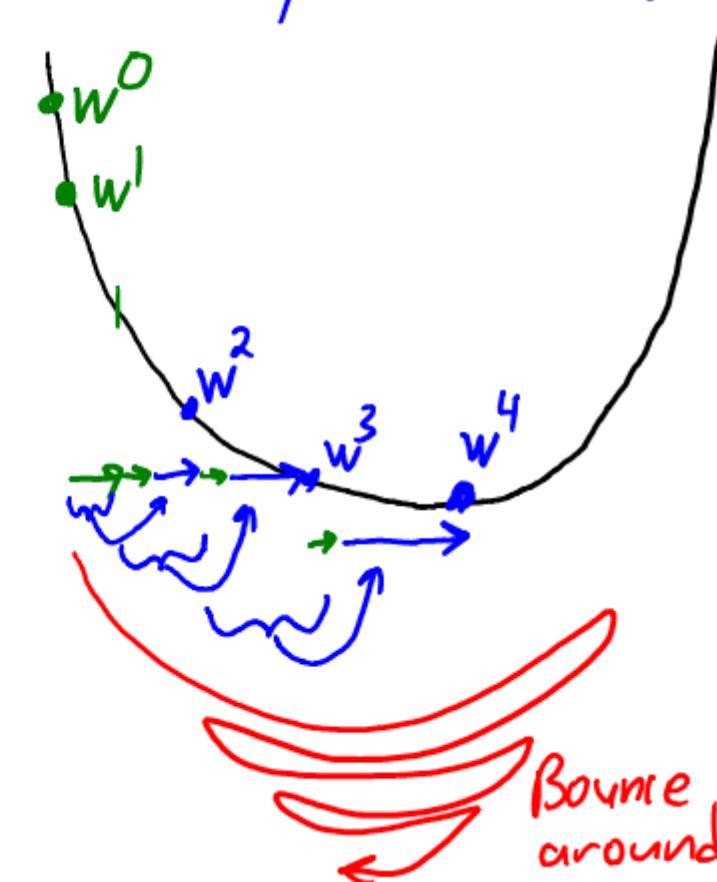
Gradient Descent vs. Heavy-Ball Method

Gradient Method



approaches from left

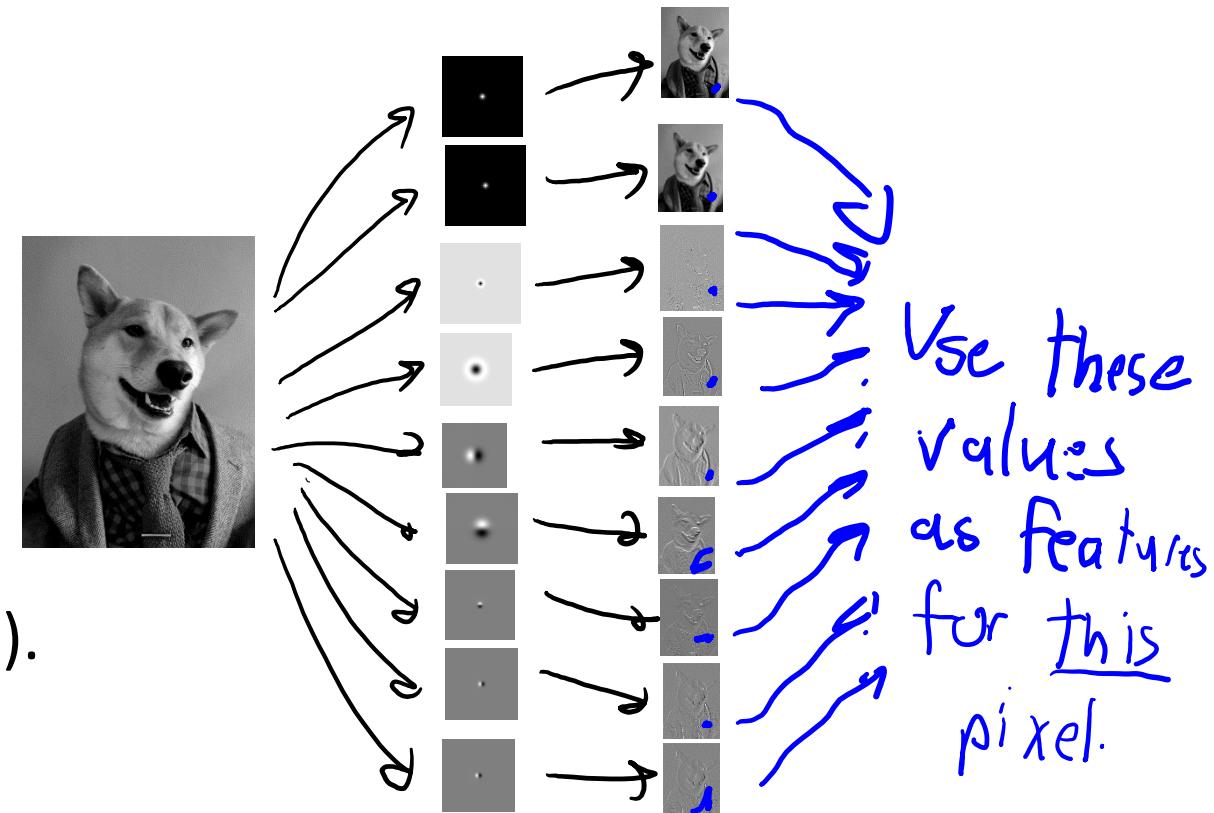
Heavy-ball Method



Good demo to check out: <https://distill.pub/2017/momentum/>

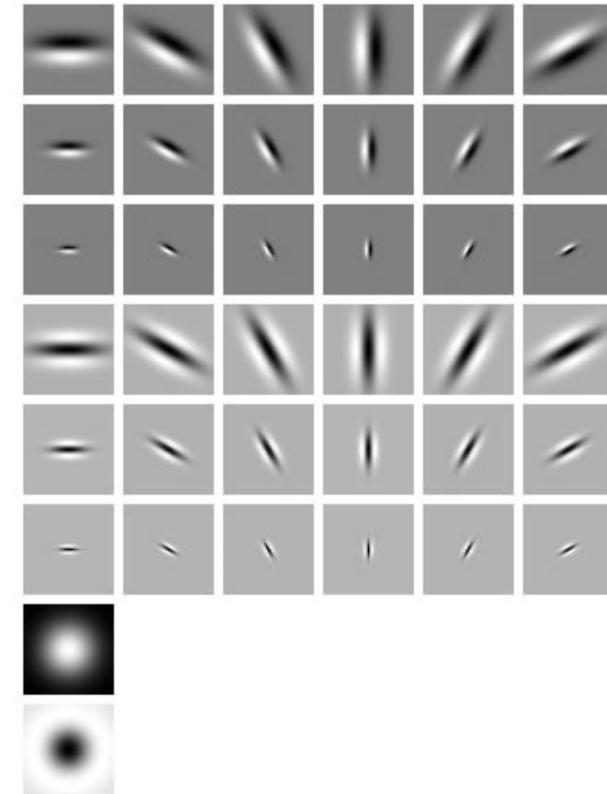
Convolutions as Features

- Classic vision methods use **convolutions as features**:
 - Usually have different types/variances/orientations.
 - Can take maxes across locations/orientations/scales.
- Notable convolutions:
 - Gaussian (blurring/averaging).
 - Laplace of Gaussian (second-derivative).
 - Gabor filters (directional first- or higher-derivative).



Filter Banks

- To characterize context, we used to use **filter banks** like “MR8”:
 - 1 Gaussian filter, 1 Laplacian of Gaussian filter.
 - 6 max(abs(Gabor)) filters:
 - 3 scales of sine/cosine (maxed over 6 orientations).



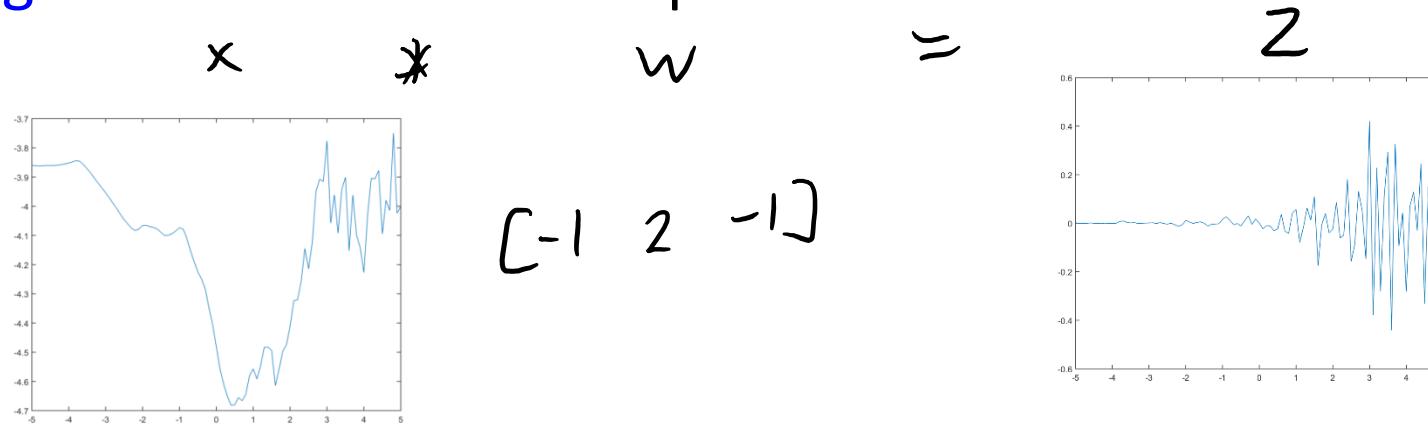
- **Convolutional neural networks** (next time!) are replacing filter banks.

Now back to our regularly scheduled program

1D Convolution as Matrix Multiplication

- **1D convolution:**

- Takes **signal 'x'** and **filter 'w'** to produces vector 'z':



- Can be written as a **matrix multiplication**:

$$W_x = \begin{bmatrix} 1 & -2 & 1 & 0 & 0 & 0 & \dots & 0 & 0 & 0 & 0 \\ 0 & 1 & -2 & 1 & 0 & 0 & \dots & 0 & 0 & 0 & 0 \\ 0 & 0 & 1 & -2 & 1 & 0 & \dots & 0 & 0 & 0 & 0 \\ \vdots & & & & & & & & & & \\ 0 & 0 & 0 & 0 & 0 & 0 & \dots & 0 & 1 & -2 & 1 \end{bmatrix} x = z$$

Typo: the signs of -1 and 2 are wrong in the matrix

1D Convolution as Matrix Multiplication

- Each element of a convolution is an **inner product**:

$$\begin{aligned} z_i &= \sum_{j=-m}^m w_j x_{i+j} \\ &= w^T x_{(i-m:i+m)} \\ &= \tilde{w}^T x \quad \text{where } \tilde{w} = [0 \ 0 \ 0 \ \underbrace{\dots}_w \ 0 \ 0] \end{aligned}$$

positions $i-m$ through $i+m$

- So **convolution is a matrix multiplication** (I'm ignoring boundaries):

$$z = \tilde{W}x \quad \text{where } \tilde{W} = \left[\begin{array}{cccccc} \overbrace{w} & & & & 0 & 0 & 0 \\ 0 & \overbrace{w} & & & 0 & 0 & \\ 0 & 0 & \overbrace{w} & & 0 & & \\ 0 & 0 & 0 & \overbrace{w} & & & \end{array} \right]$$

} matrix can be
very sparse and
only has $2m+1$ variables.

- The shorter 'w' is, the more sparse the matrix is.

2D Convolution as Matrix Multiplication

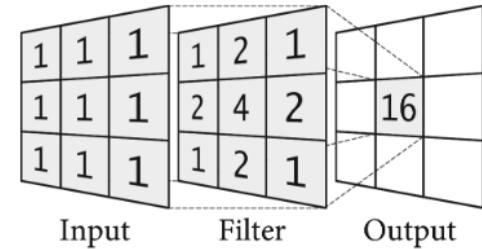
- 2D convolution:

- Signal 'x', filter 'w', and output 'z' are now all **images/matrices**:

$$x * w = z$$



$$\begin{bmatrix} -2 & -1 & 0 \\ -1 & 0 & 1 \\ 0 & 1 & 2 \end{bmatrix}$$



- Vectorized 'z' can be written as a **matrix multiplication** with vectorized 'x':

$$W = \begin{bmatrix} -2 & -1 & 0 & 0 & 0 & \dots & 0 & -1 & 0 & 1 & 0 & 0 & \dots & 0 & 0 & 1 & 2 & 0 & 0 & \dots & 0 & 0 \\ 0 & -2 & -1 & 0 & 0 & \dots & 0 & 0 & -1 & 0 & 1 & 0 & 0 & \dots & 0 & 0 & 0 & 1 & 2 & 0 & \dots & 0 & 0 \\ \vdots & \vdots & \vdots & \vdots & \vdots & \ddots & \vdots & \ddots & \vdots & \vdots & \vdots & \vdots & \vdots & \ddots & \vdots & \vdots & \ddots & \vdots & \vdots \\ 0 & 0 & 0 & 0 & 0 & \dots & 0 & 2 & -1 & 0 & -6 & 0 & 0 & \dots & 0 & -1 & 0 & 1 & 0 & \dots & 0 & 0 & \dots & 0 & 1 & 2 & 0 \\ 0 & 0 & 0 & 0 & 0 & \dots & 0 & 0 & 2 & -1 & 0 & 0 & 0 & \dots & 0 & -1 & 0 & 1 & 0 & \dots & 0 & 0 & \dots & 0 & 1 & 2 & 0 \\ 0 & 0 & 0 & 0 & 0 & \dots & 0 & 0 & 0 & 2 & -1 & 0 & 0 & \dots & 0 & 0 & -1 & 0 & 1 & \dots & 0 & 0 & \dots & 0 & 1 & 2 & 0 \\ \vdots & \vdots & \vdots & \vdots & \vdots & \ddots & \vdots & \ddots & \vdots & \vdots & \vdots & \vdots & \vdots & \ddots & \vdots & \vdots & \ddots & \vdots & \vdots \end{bmatrix}$$

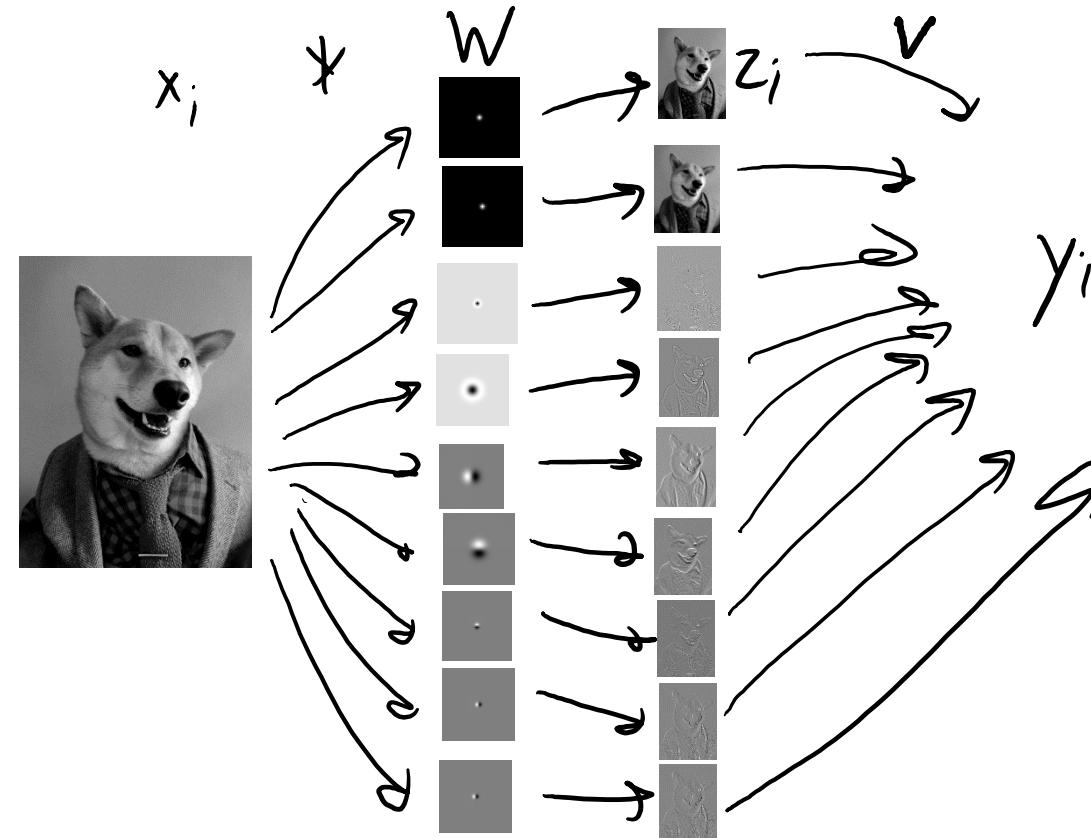
Motivation for Convolutional Neural Networks

- Consider training neural networks on 256 by 256 images.
 - This is 256 by 256 by 3 \approx 200,000 inputs.
- If first layer has $k=10,000$, then it has **about 2 billion parameters**.
 - We want to avoid this huge number (due to storage and overfitting).
- Key idea: make Wx_i act like several convolutions (to make it sparse):
 1. Each row of W only applies to part of x_i .
 2. Use the same parameters between rows.
- Forces most weights to be zero, reduces number of parameters.

$$w_1 = [0 \quad 0 \quad 0 \quad \dots \quad w \quad \dots \quad 0 \quad 0 \quad 0]$$
$$w_2 = [0 \quad \dots \quad w \quad \dots \quad 0 \quad 0 \quad 0 \quad 0]$$

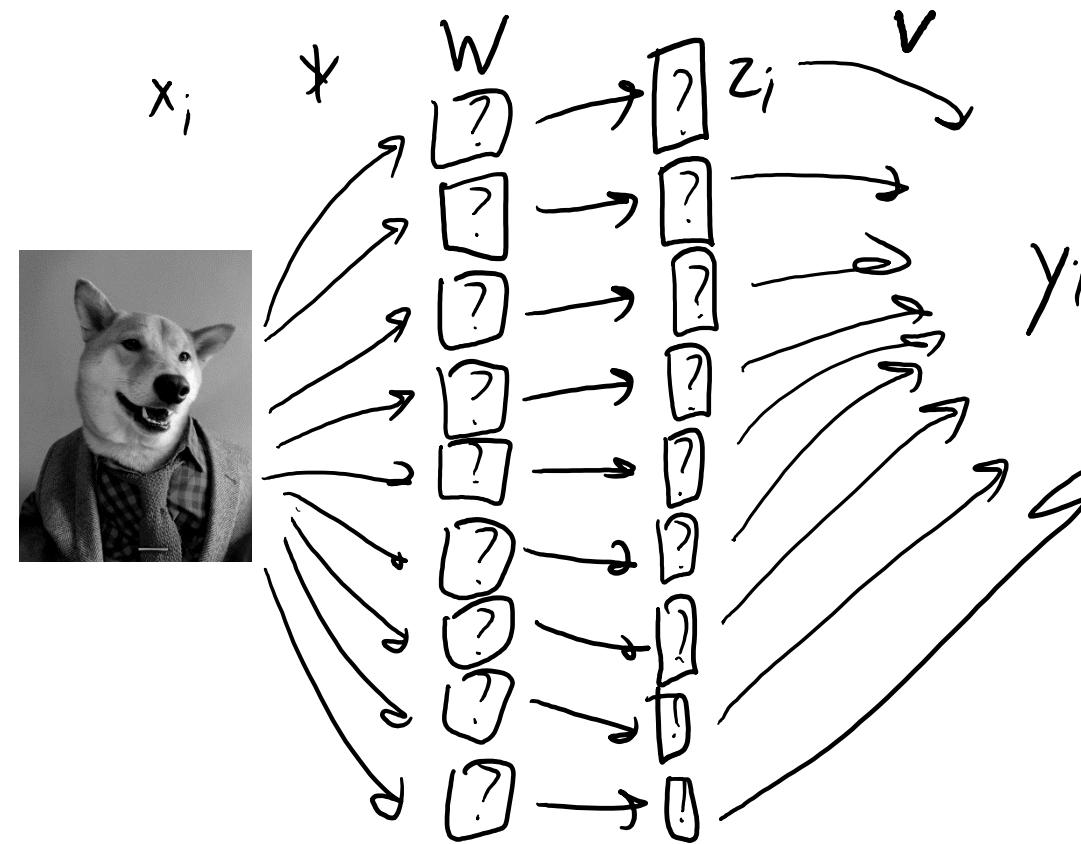
Motivation for Convolutional Neural Networks

- Classic vision methods uses **fixed convolutions** as features:
 - Usually have **different types/variances/orientations**.
 - Can do subsampling or take **maxes** across locations/orientations/scales.



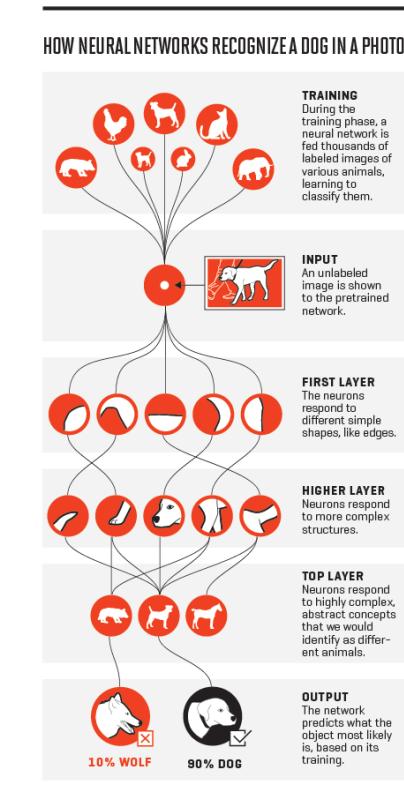
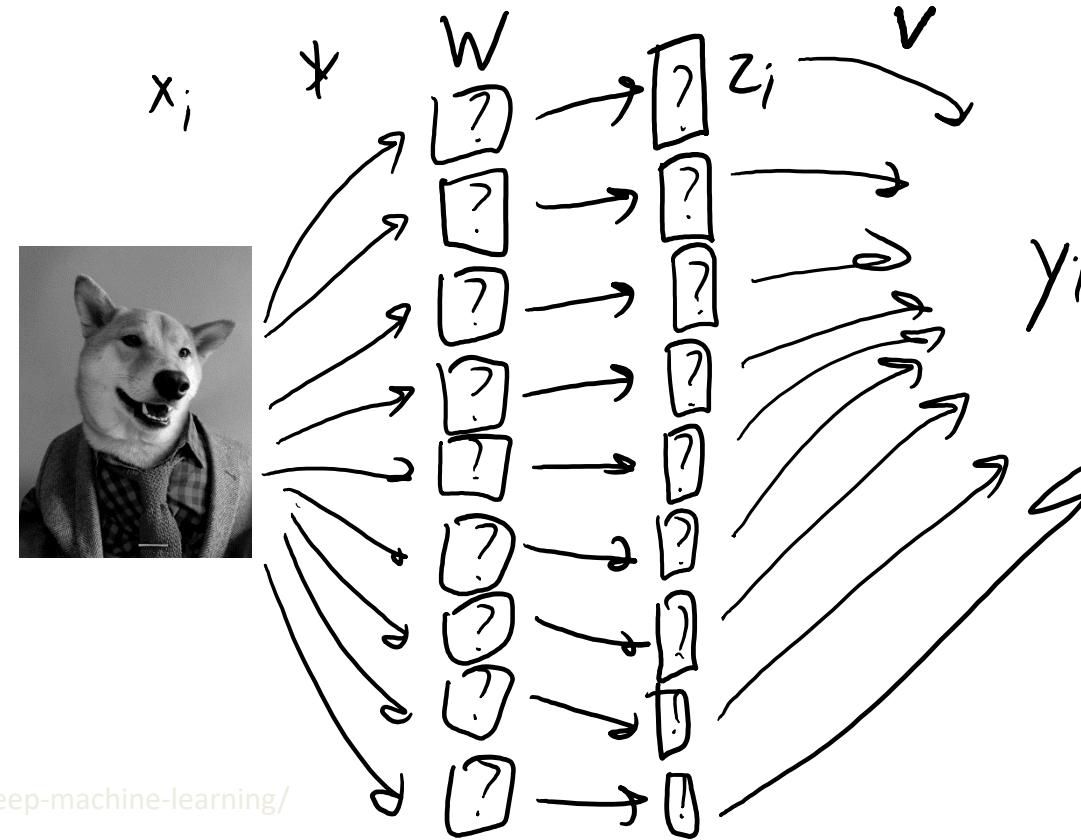
Motivation for Convolutional Neural Networks

- Convolutional neural networks learn the convolutions:
 - Learning ‘W’ and ‘v’ automatically chooses types/variances/orientations.
 - Don’t pick from fixed convolutions, but learn the elements of the filters.



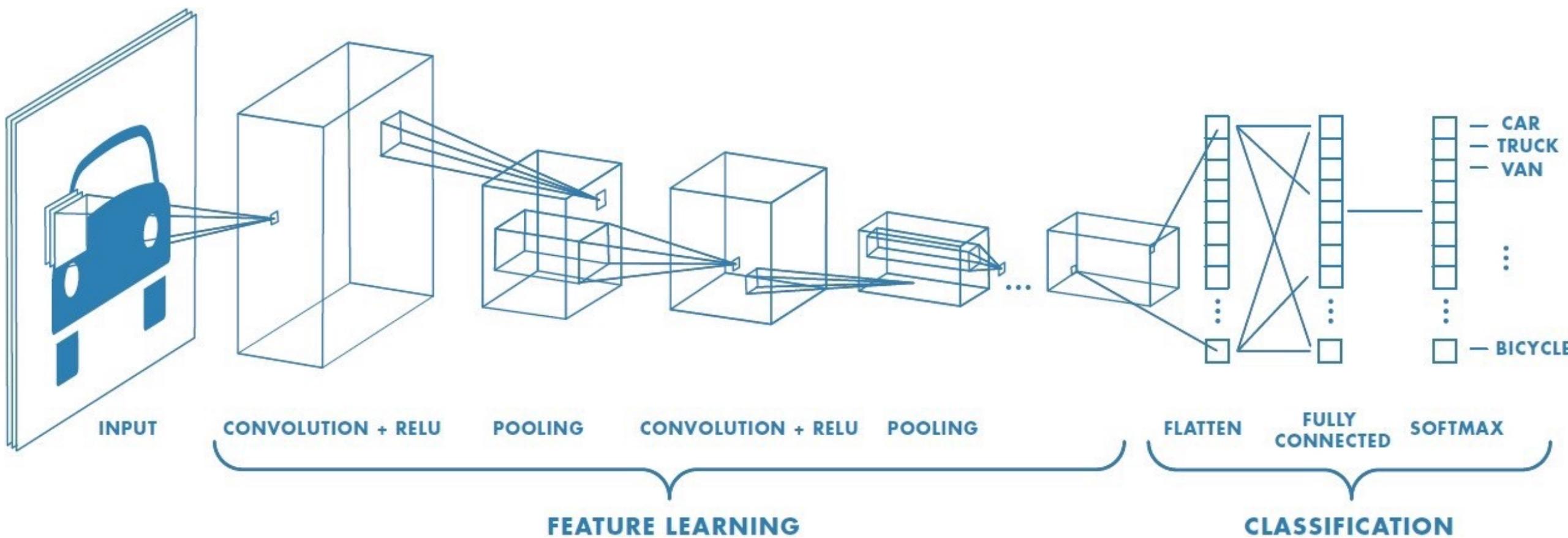
Motivation for Convolutional Neural Networks

- Convolutional neural networks learn the convolutions:
 - Learning ‘W’ and ‘v’ automatically chooses types/variances/orientations.
 - Can do multiple layers of convolution to get deep hierarchical features.

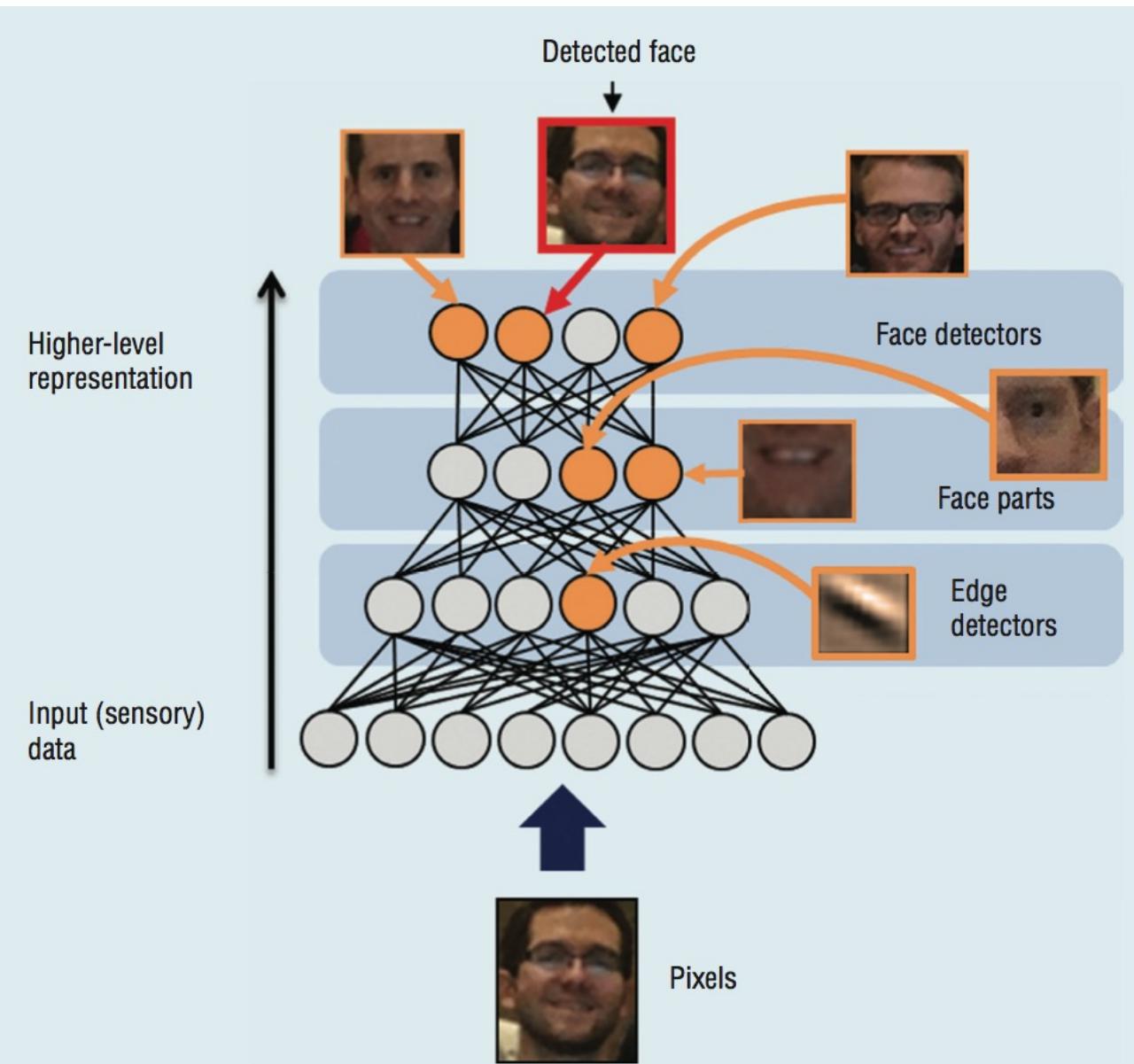


Two Main Motivations

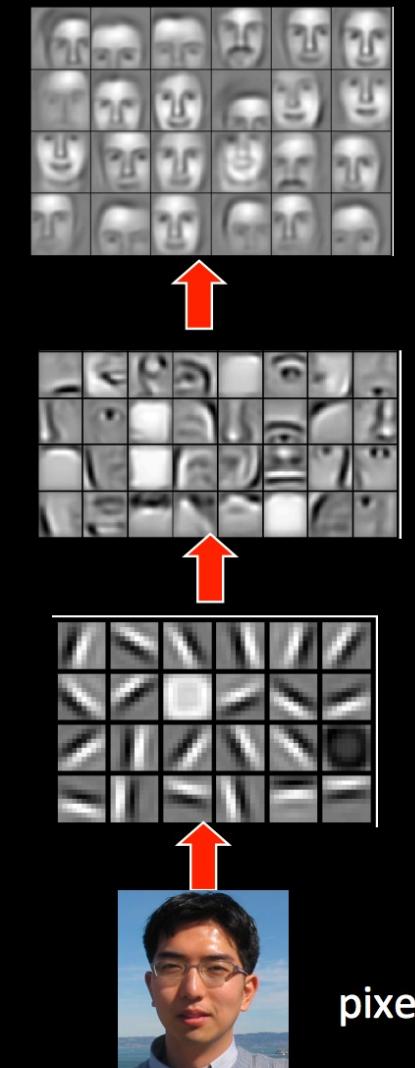
- Translation invariance (data-efficient to learn, less likely to overfit)
- Hierarchy

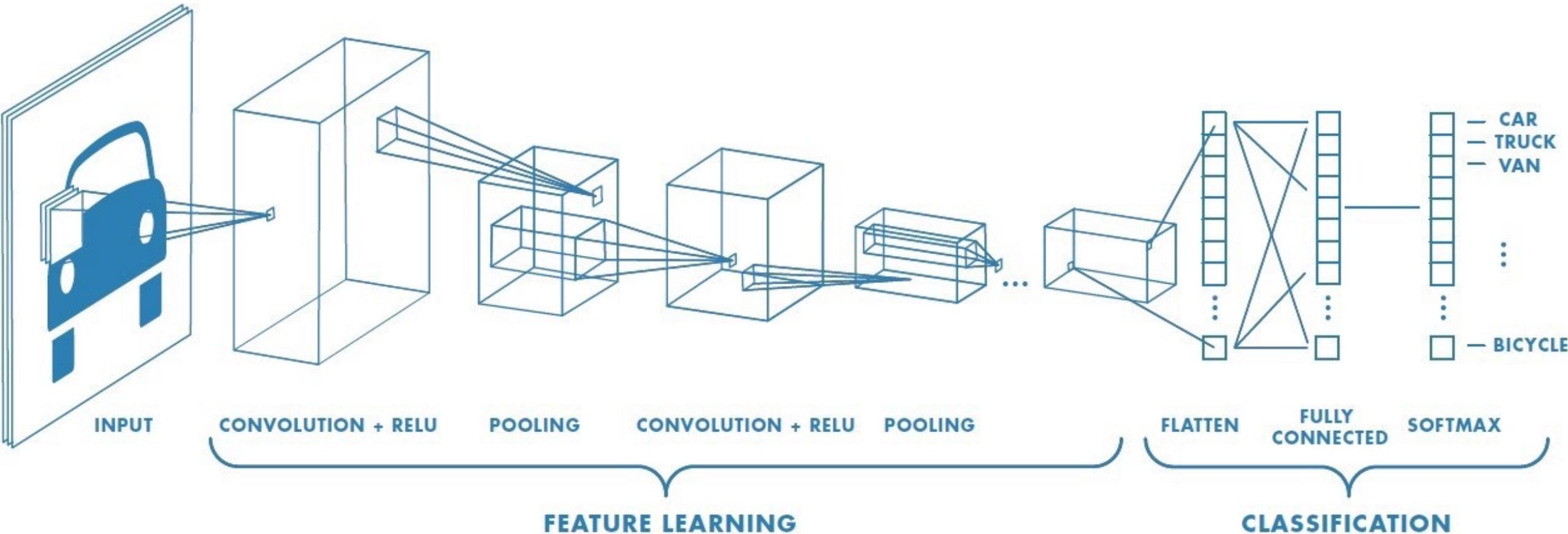


Hierarchically composed feature representations



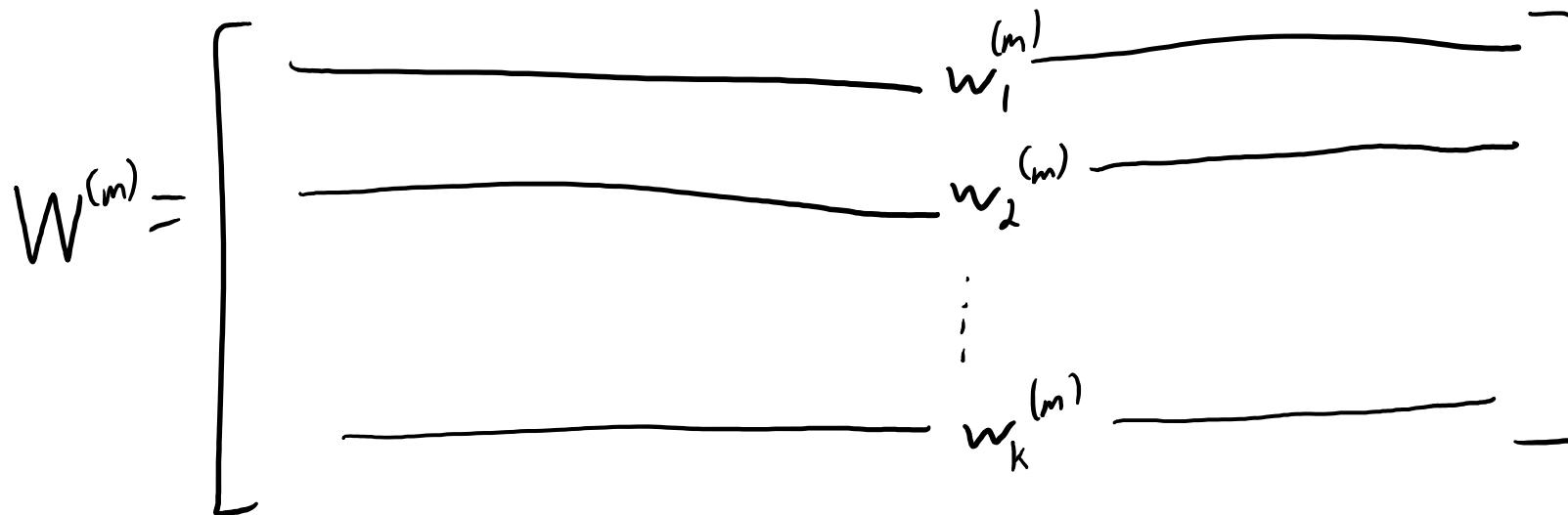
Hierarchy of feature representations





Convolutional Neural Networks

- Convolutional Neural Networks classically have 3 layer “types”:
 - Fully connected layer: usual neural network layer with unrestricted W.



Convolutional Neural Networks

- Convolutional Neural Networks classically have 3 layer “types”:
 - Fully connected layer: usual neural network layer with unrestricted W.
 - Convolutional layer: restrict W to act like several convolutions.

1D example

$$W^{(m)} = \left[\begin{array}{ccccccccc} & w_1^{(m)} & & & & & & & \\ 0 & 0 & 0 & & w_1^{(m)} & & & & \\ & & & & & & & & \\ 0 & 0 & 0 & 0 & 0 & 0 & & & \\ & & & & & & w_1^{(m)} & & \\ \hline & w_2^{(m)} & & & & & & & \\ 0 & 0 & 0 & & w_2^{(m)} & & & & \\ & & & & & & & & \\ 0 & 0 & 0 & 0 & 0 & 0 & & & \\ & & & & & & w_2^{(m)} & & \\ \hline 0 & 0 & 0 & & & & & & \\ 0 & 0 & 0 & & & & & & \\ 0 & 0 & 0 & & & & & & \end{array} \right]$$

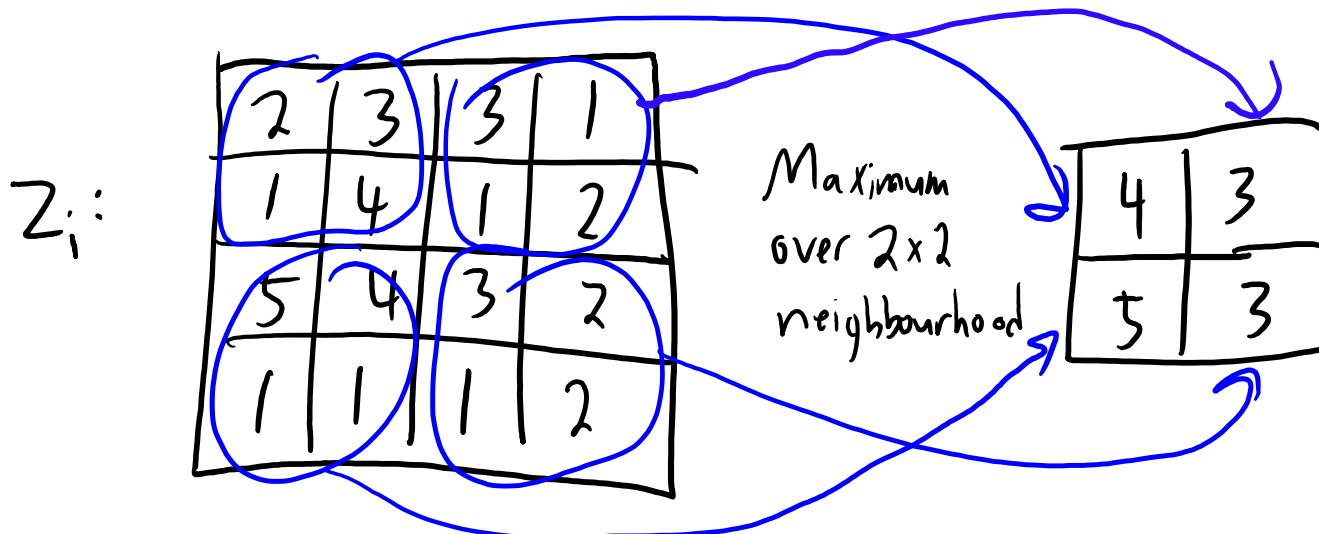
distance between centers of convolution is called "stride"

Same $w_i^{(m)}$ used across multiple rows.

Sparse and small number of parameters

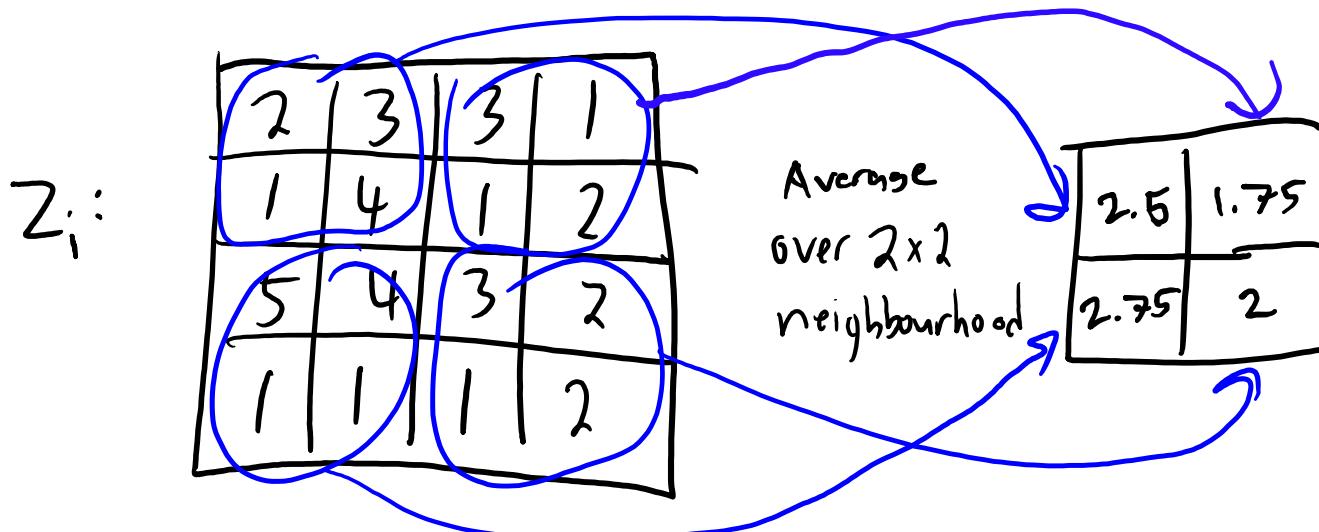
Convolutional Neural Networks

- Convolutional Neural Networks classically have 3 layer “types”:
 - Fully connected layer: usual neural network layer with unrestricted W.
 - Convolutional layer: restrict W to act like several convolutions.
 - Pooling layer: combine results of convolutions.
 - Can add some invariance or just make the number of parameters smaller.
 - Often ‘max pooling’:



Convolutional Neural Networks

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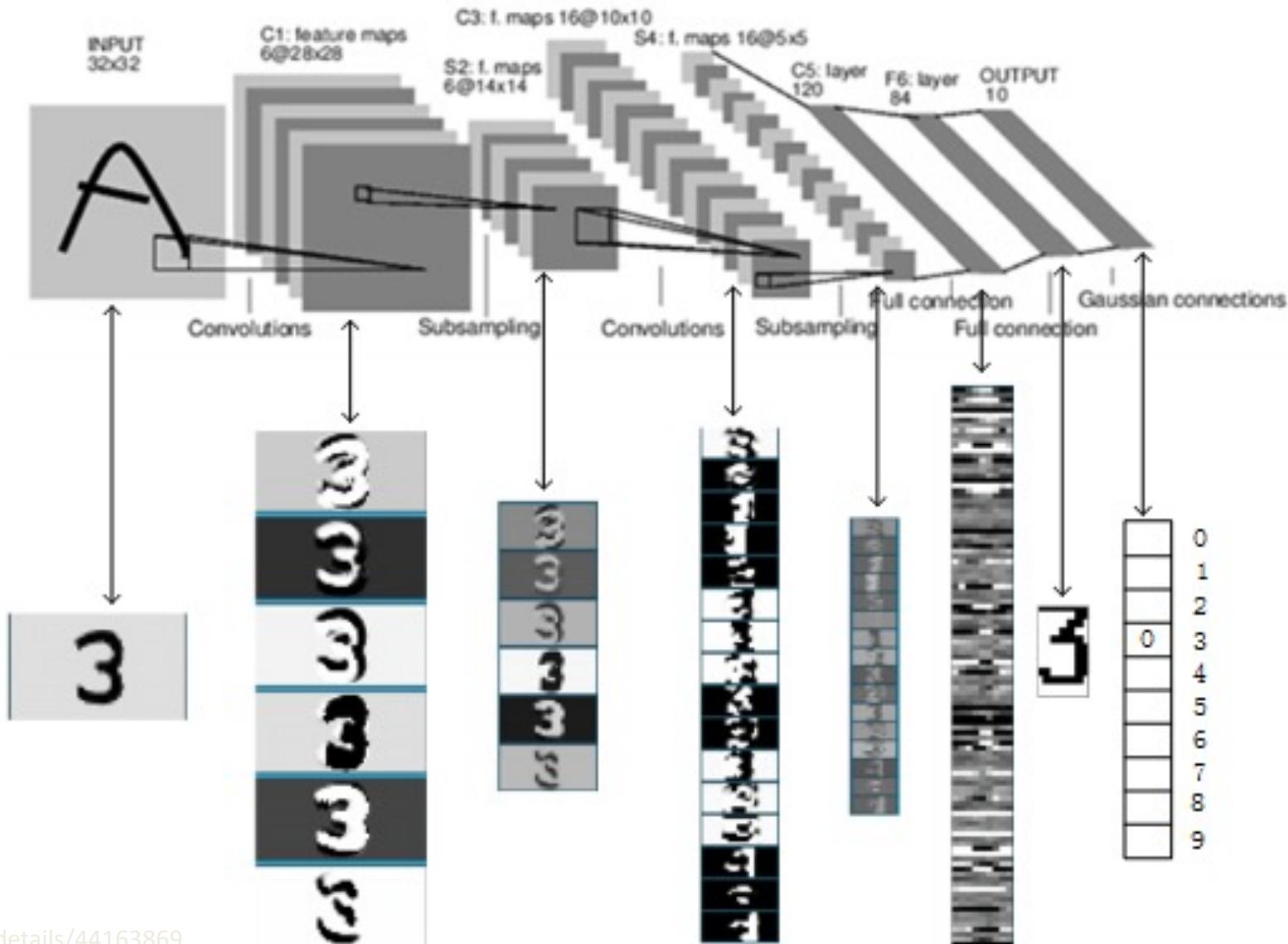


bonus!

Max Pooling vs Average Pooling

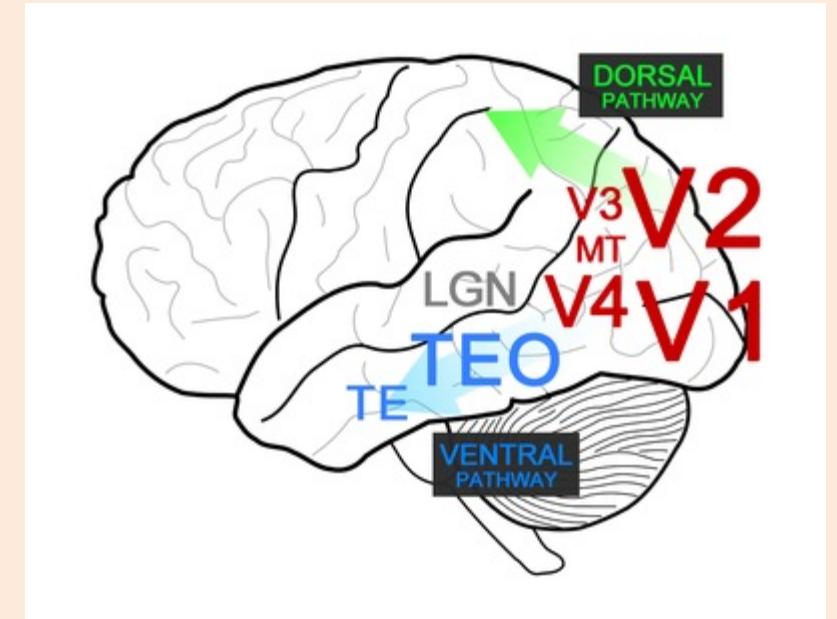
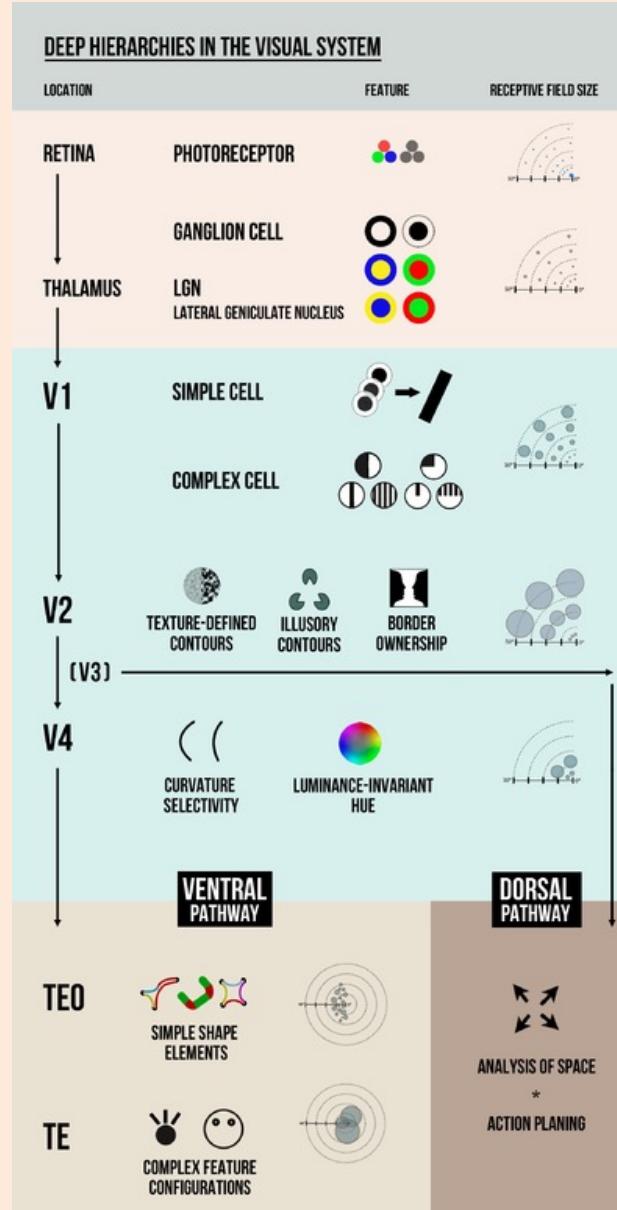
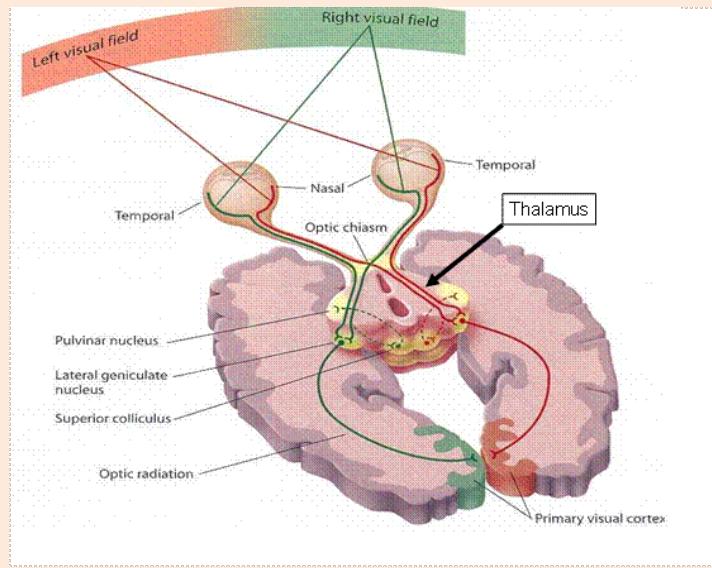
- Both downsample the image
- Max pooling: “any of these options is present”
 - Much more common, especially in early layers
 - “There’s an edge here, but I don’t really care how thick it is”
- Average pooling: “all/most of these options are present”
 - If used, more often at the end of the network
 - “Most of the big patches look like a picture of a train”

LeNet for Optical Character Recognition



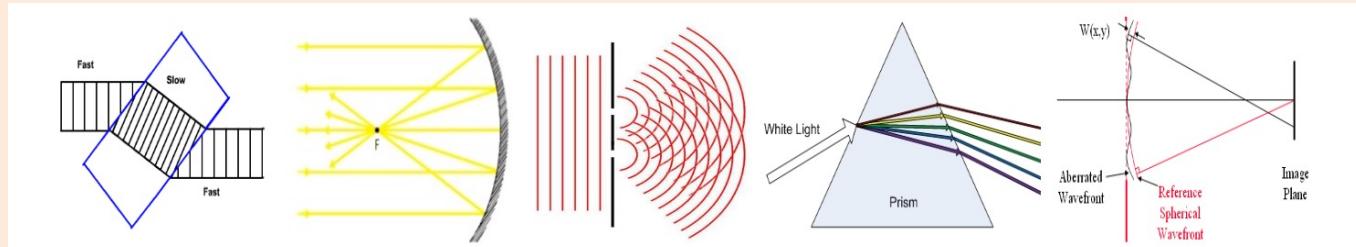
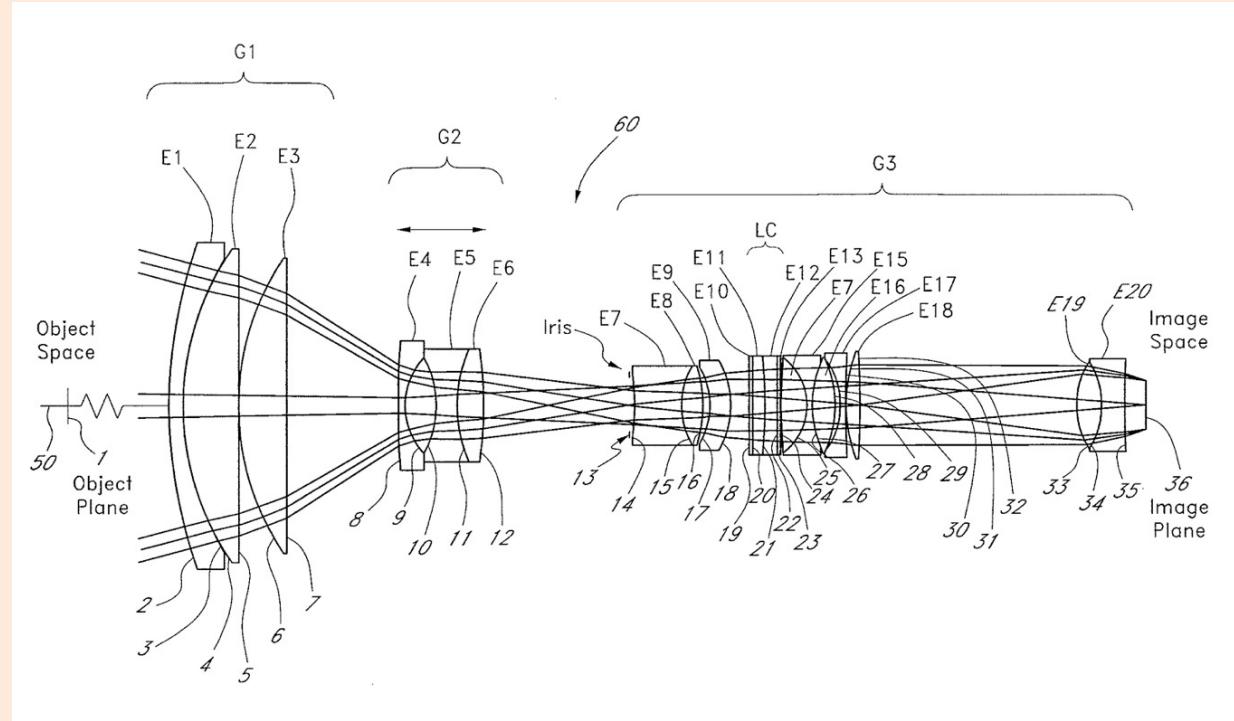
bonus!

Deep Hierarchies in the Visual System



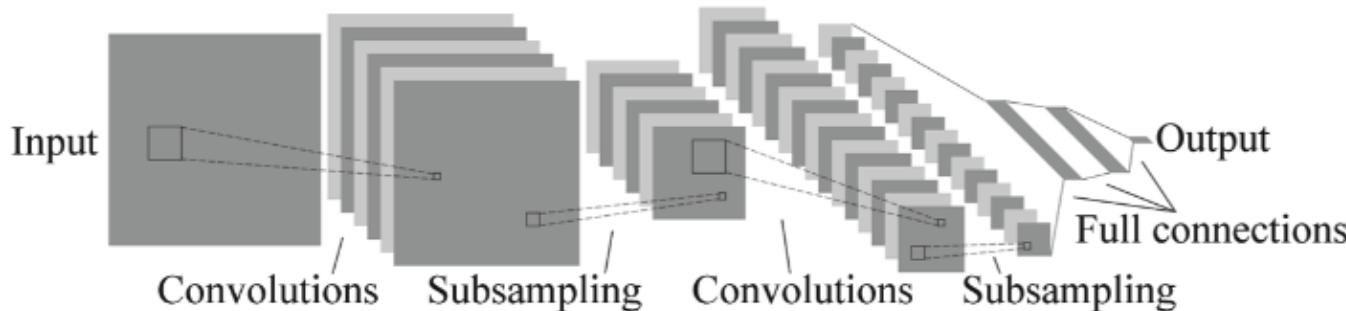
bonus!

Deep Hierarchies in Optics



Convolutional Neural Networks

- Classic convolutional neural network (LeNet):



- Visualizing the “activations” of the layers:

- <http://scs.ryerson.ca/~aharley/vis/conv>
- <https://youtu.be/AgkfIQ4IGaM>



Next

- A very small selection of key advances, things you should know, and tricks of the trade

Recent Lectures: Deep Learning

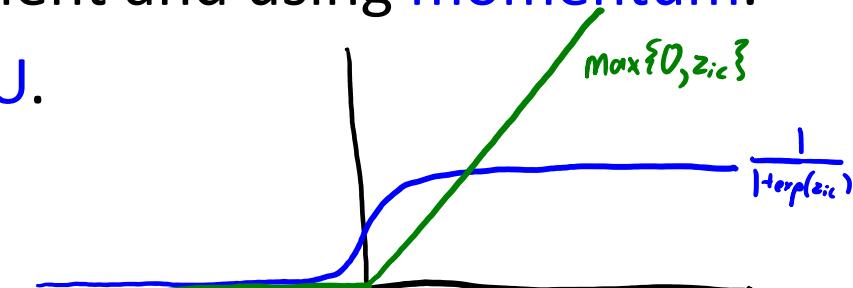
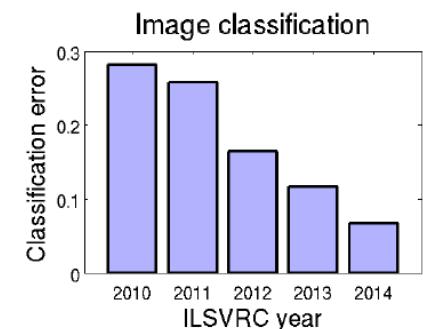
- We've been discussing **neural network / deep learning** models:

$$\hat{y}_i = v^T h(W^{(m)} h(W^{(m-1)} h(\dots \dots W^{(2)} h(W^{(1)} x_i)) \dots))$$

- We discussed unprecedented vision/speech performance.

- We discussed methods to make SGD work better:

- Parameter initialization and data transformations.
- Setting the step size(s) in stochastic gradient and using momentum.
- Alternative non-linear functions like ReLU.



“Residual” Networks (ResNets)

- Impactful recent idea is residual networks (**ResNets**):

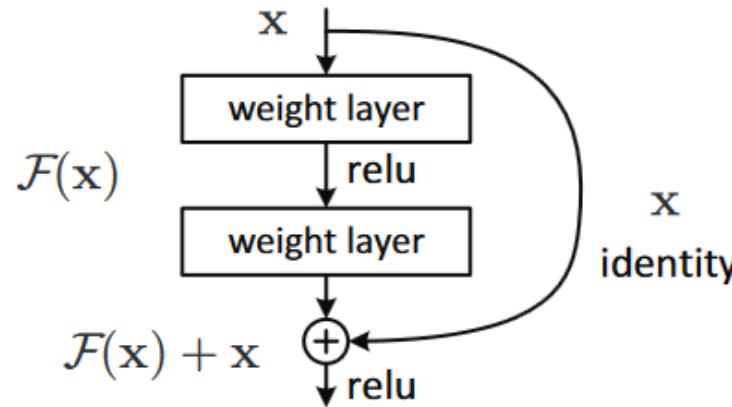


Figure 2. Residual learning: a building block.

- You can **take previous (non-transformed) layer as input** to current layer.
 - Also called “skip connections” or “highway networks”.
- **Non-linear part of the network only needs to model residuals.**
 - Non-linear parts are just “pushing up or down” a linear model in various places.
- This was a key idea behind first methods that used 100+ layers.
 - Evidence that biological networks have skip connections like this.

bonus!

DenseNet

- More recent variation is “DenseNets”:
 - Each layer can see all the values from many previous layers.
 - Gets rid of vanishing gradients.
 - May get same performance with fewer parameters/layers.

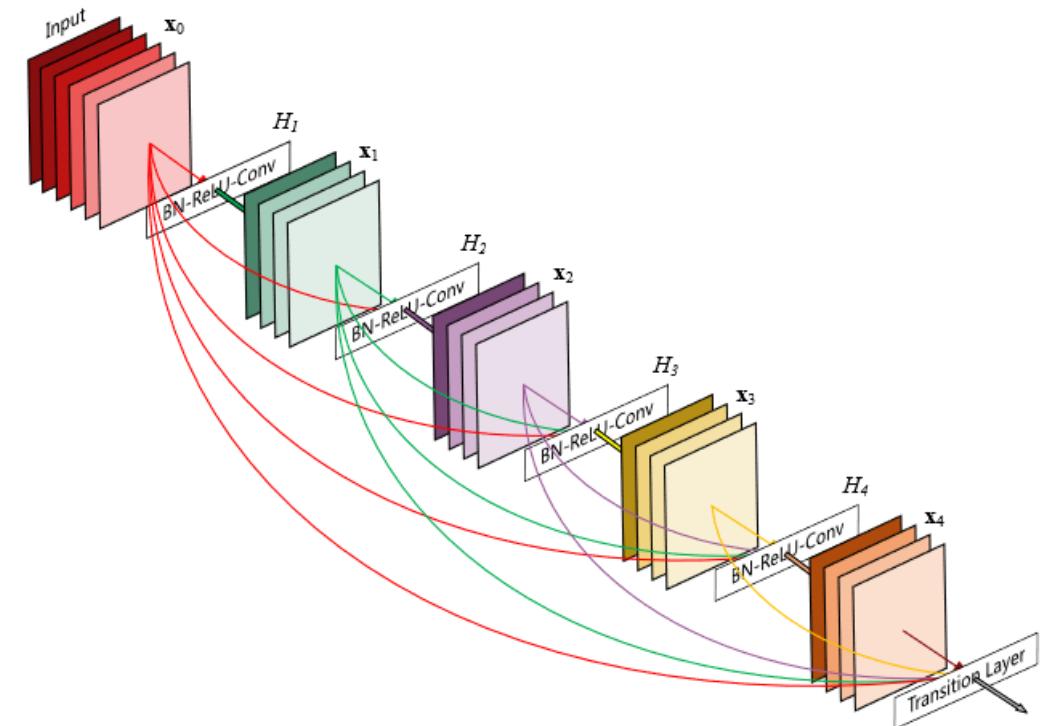


Figure 1: A 5-layer dense block with a growth rate of $k = 4$. Each layer takes all preceding feature-maps as input.

Deep Learning and the Fundamental Trade-Off

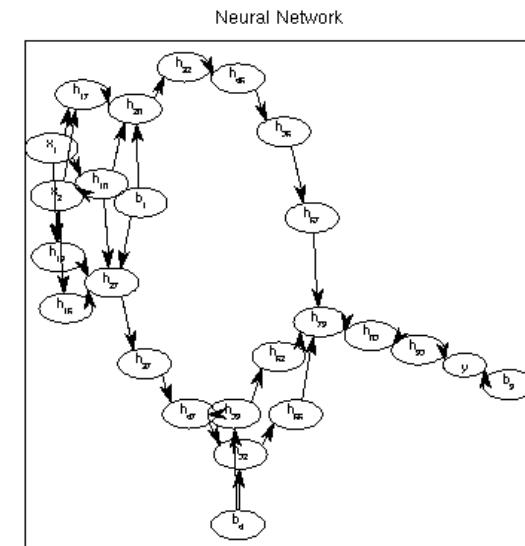
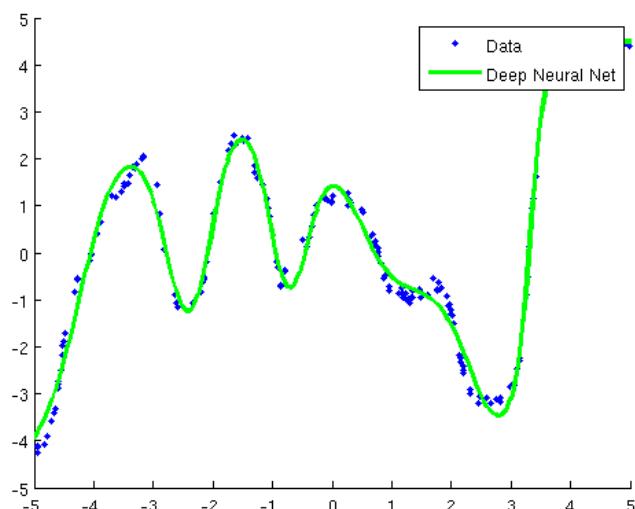
- Neural networks are subject to the fundamental trade-off:
 - With increasing depth, training error of global optima decreases.
 - With increasing depth, training error may poorly approximate test error.
- We want deep networks to model highly non-linear data.
 - But increasing the depth can lead to **overfitting**.
- How could GoogLeNet use 22 layers?
 - Many forms of **regularization** and keeping model complexity under control.
 - Unlike linear models, typically use **multiple types of regularization**.

Standard Regularization

- Traditionally, we've added our usual L2-regularizers:

$$f(v, W^{(3)}, W^{(2)}, W^{(1)}) = \frac{1}{2} \sum_{i=1}^n (v^\top h(W^{(3)} h(W^{(2)} h(W^{(1)} x_i))) - y_i)^2 + \frac{\lambda_4}{2} \|v\|^2 + \frac{\lambda_3}{2} \|W^{(3)}\|_F^2 + \frac{\lambda_2}{2} \|W^{(2)}\|_F^2 + \frac{\lambda_1}{2} \|W^{(1)}\|_F^2$$

- L2-regularization often called “weight decay” in this context.
 - Could also use L1-regularization: gives sparse network.



Standard Regularization

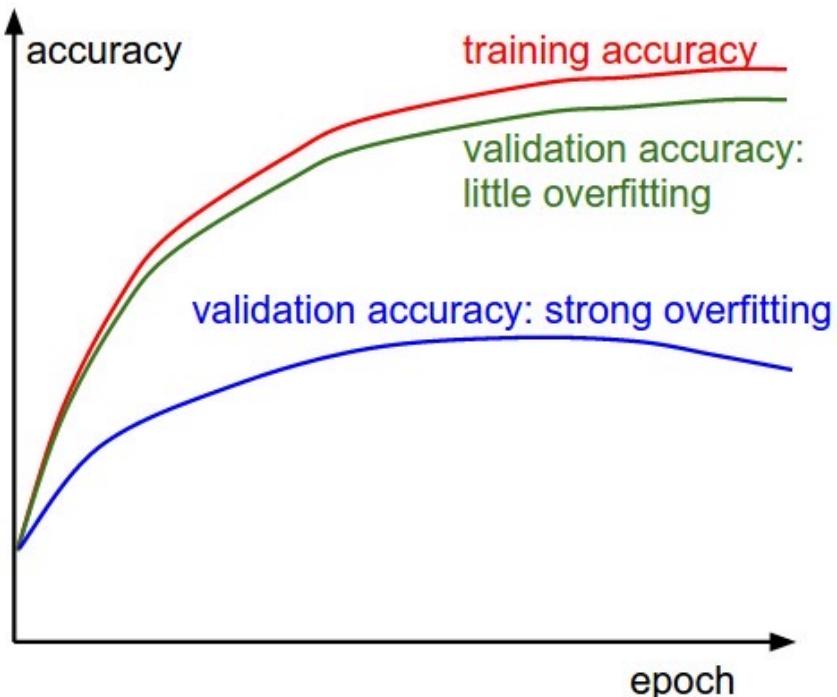
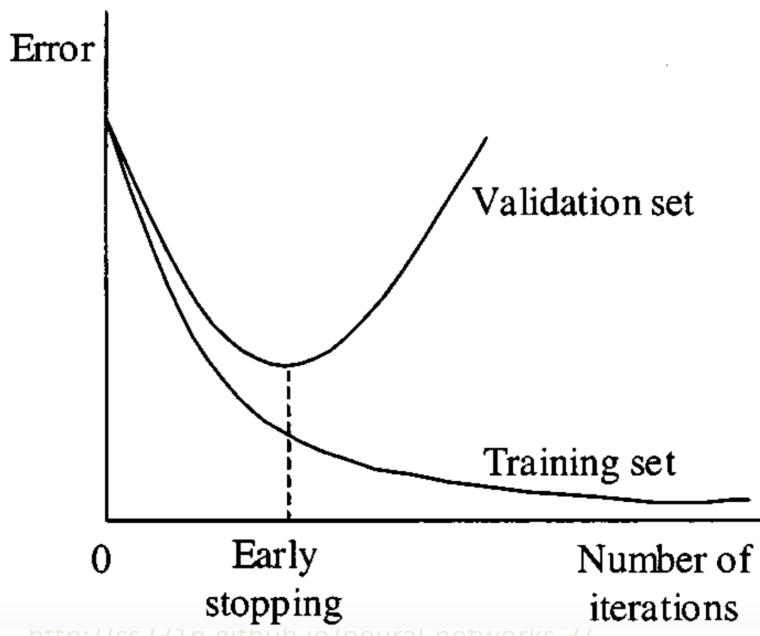
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- L2-regularization often called “weight decay” in this context.
 - Adds λW to gradient, so (S)GD “decays” the weights ‘W’ at each step
 - Could also use L1-regularization: gives sparse network.
- Hyper-parameter optimization gets expensive:
 - Try to optimize validation error in terms of $\lambda_1, \lambda_2, \lambda_3, \lambda_4$.
 - In addition to step-size, number of layers, size of layers, initialization.
- Recent result:
 - Adding a regularizer in this way can create bad local optima.

Early Stopping

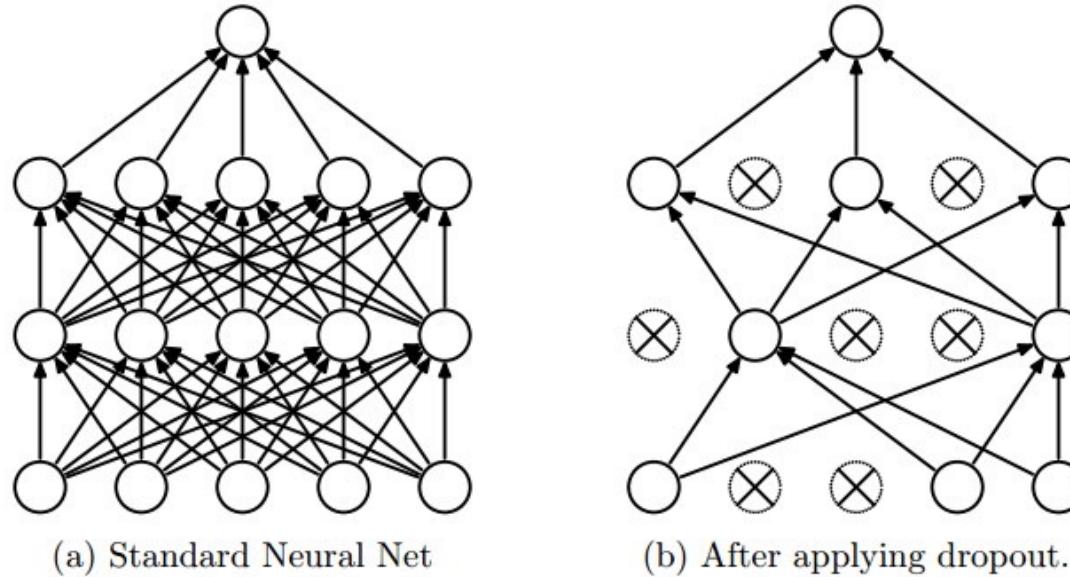
- Another common type of regularization is “early stopping”:
 - Monitor the validation error as we run stochastic gradient.
 - Stop the algorithm if validation error starts increasing.



bonus!

Dropout

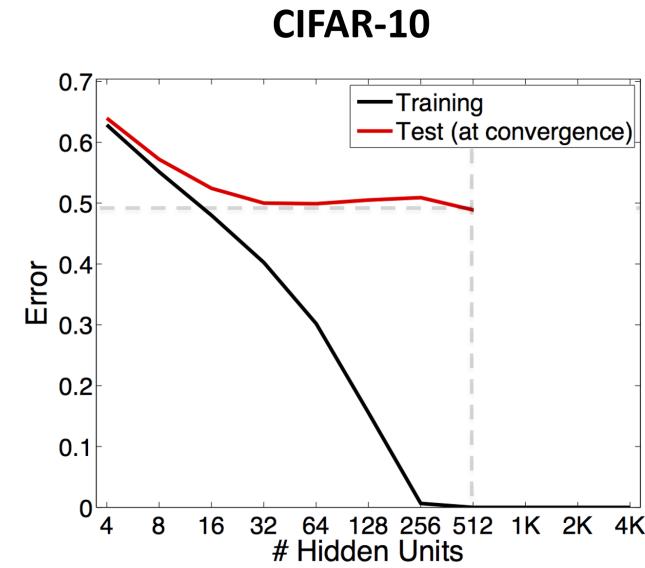
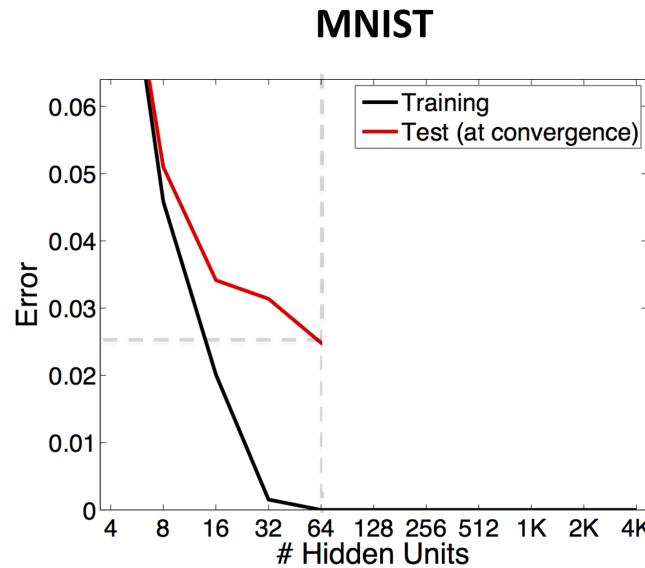
- **Dropout** is a more recent form of explicit regularization:
 - On each iteration, randomly set some x_i and z_i to zero (often use 50%).



- Adds invariance to missing inputs or latent factors
 - Encourages distributed representation rather than relying on specific z_i .
- Can be interpreted as an ensemble over networks with different parts missing.
- After a lot of early success, dropout is already kind of going out of fashion.

“Hidden” Regularization in Neural Networks

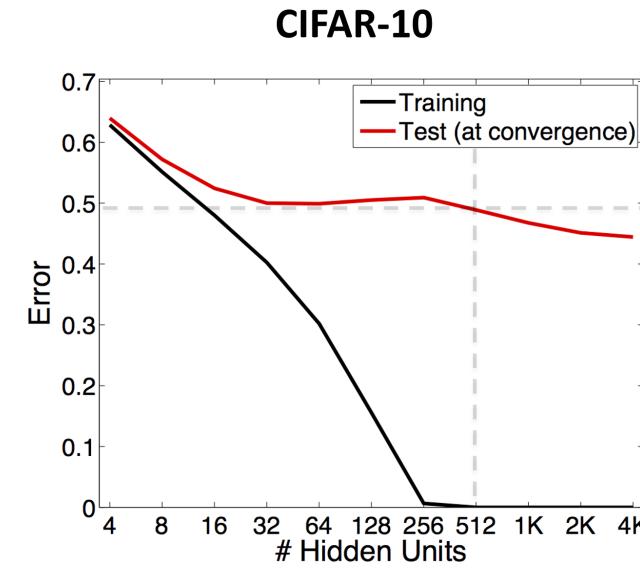
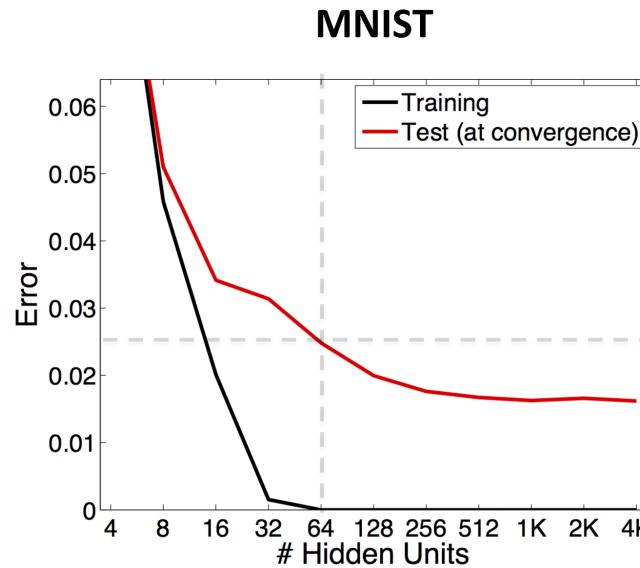
- Fitting single-layer neural network with SGD and no regularization:



- Training goes to 0 with enough units: we’re finding a global min.
- What should happen to training and test error for larger #hidden?

“Hidden” Regularization in Neural Networks

- Fitting single-layer neural network with SGD and no regularization:



- Test error continues to go down!?! Where is fundamental trade-off??
- There exist global mins with large #hidden units have test error = 1.
 - But among the global minima, SGD is somehow converging to “good” ones.

bonus!

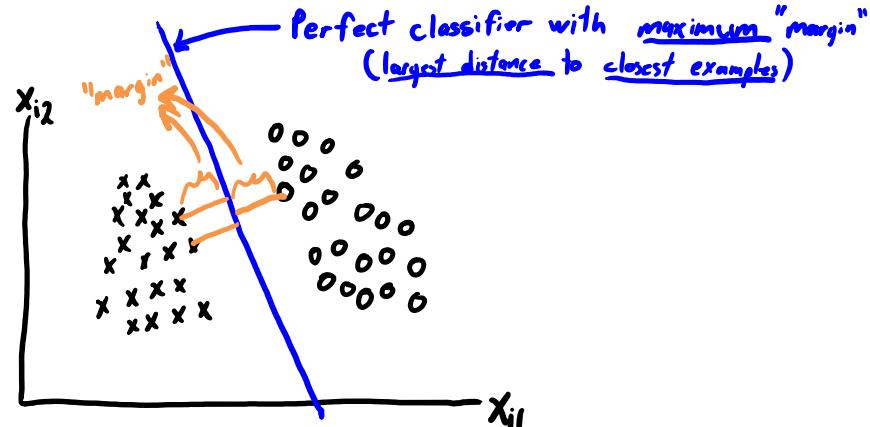
Implicit Regularization of SGD

- There is growing evidence that using SGD regularizes parameters.
 - We call this the “implicit regularization” of the optimization algorithm.
- Beyond empirical evidence, we know this happens in simpler cases.
- Example of implicit regularization:
 - Consider a least squares problem where there exists a ‘w’ where $X w = y$.
 - Residuals are all zero, we fit the data exactly.
 - You run [stochastic] gradient descent starting from $w=0$.
 - Converges to solution $X w = y$ that has the minimum L2-norm.
 - So using SGD is equivalent to L2-regularization here, but regularization is “implicit”.

bonus!

Implicit Regularization of SGD

- Example of implicit regularization:
 - Consider a **logistic regression** problem where data is linearly separable.
 - We can fit the data exactly.
 - You run gradient descent from any starting point.
 - Converges to **max-margin solution** of the problem.
 - So using gradient descent is equivalent to encouraging large margin.



- Similar result known for **boosting**.

(pause)

Deep Learning “Tricks of the Trade”

- We've discussed **heuristics to make deep learning work:**
 - Parameter initialization and data transformations.
 - Setting the **step size(s)** in stochastic gradient and using **momentum**.
 - **ResNets** and alternative non-linear functions like **ReLU**.
 - Different forms of regularization:
 - L2-regularization, early stopping, dropout, implicit regularization from SGD.
- These are often **still not enough** to get deep models working.
- Deep computer vision models are all **convolutional neural networks**:
 - The $W^{(m)}$ are **very sparse** and have **repeated parameters** (“tied weights”).
 - Drastically reduces number of parameters (speeds training, reduces overfitting).

Summary

- **ResNets** include untransformed previous layers.
 - Network focuses non-linearity on residual, allows huge number of layers.
- **Regularization** is crucial to neural net performance:
 - L2-regularization, early stopping, dropout, implicit regularization of SGD.
- **Convolutional neural networks:**
 - Restrict $W^{(m)}$ matrices to represent sets of convolutions.
 - Often combined with max (pooling).
- Next time: modern convolutional neural networks and applications.
 - Image segmentation, depth estimation, image colorization, artistic style.