#### **CS 179: LECTURE 16**

MODEL COMPLEXITY, REGULARIZATION, AND CONVOLUTIONAL NETS

### LAST WEEK

- Intro to cuDNN
- Deep neural nets using cuBLAS and cuDNN

#### TODAY

- Building a "better" model for image classification
- Overfitting and regularization
- Convolutional neural nets

#### MODEL COMPLEXITY

- Consider a class of models f(x; w)
  - A function f of an input x with parameters w
  - For now, let's just consider  $x \in \mathbb{R}$  (1D input) as a toy example
- Polynomial regression fits a polynomial of degree d to our input, i.e.  $f(x; w) = w_0 + w_1 x + w_2 x^2 + \cdots + w_d x^d$
- Intuitively, a higher degree polynomial is a more complex model function than a lower degree polynomial

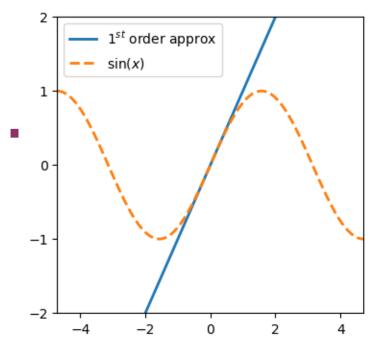
#### INTUITION: TAYLOR SERIES

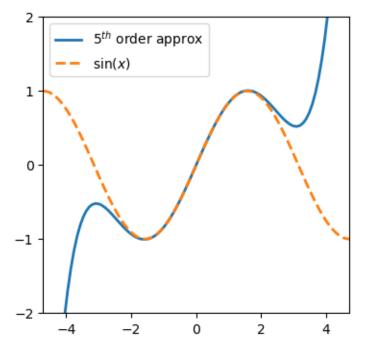
- More formally, one model class is more complex than another if it contains more functions
- If we already know the function g that we want to approximate, we can use Taylor polynomials
  - For many functions g, we have  $g(x) = \sum_{k=0}^{\infty} w_k x^k$
  - One way to approximate is as  $g(x) \approx \sum_{k=0}^{d} w_k x^k$
- Higher degree polynomial gives a better approximation?

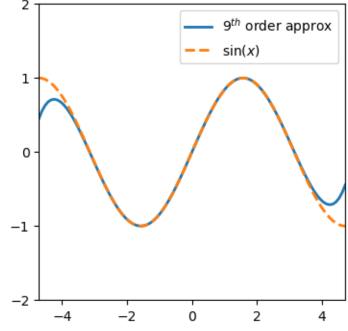
#### INTUITION: TAYLOR SERIES

■ Taylor expansions of sin(x) about 0 for d = 1,5,9

Taylor Approximations of sin(x)







# LEAST SQUARES FITTING

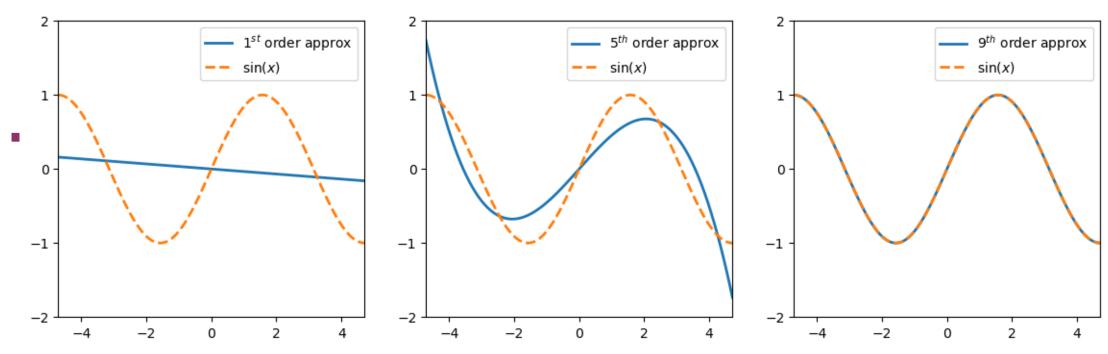
- Generally, we don't know the true function a priori
- Instead, we approximate it with a model function f(x; w)
- Rather than Taylor coefficients, we really want parameters
- $w^*$  that minimize some loss function J(w) on a dataset  $\{(x^{(i)}, y^{(i)})\}_{i=1}^N$ , e.g. mean squared error:

$$w^* = \underset{w}{\operatorname{argmin}} J(w) = \underset{w}{\operatorname{argmin}} \frac{1}{N} \sum_{i=1}^{N} (y^{(i)} - f(x^{(i)}; w))^2$$

# LEAST SQUARES FITTING

• Least squares polynomial fits of sin(x) for d = 1,5,9

Least-Squares Polynomial Approximations of sin(x)

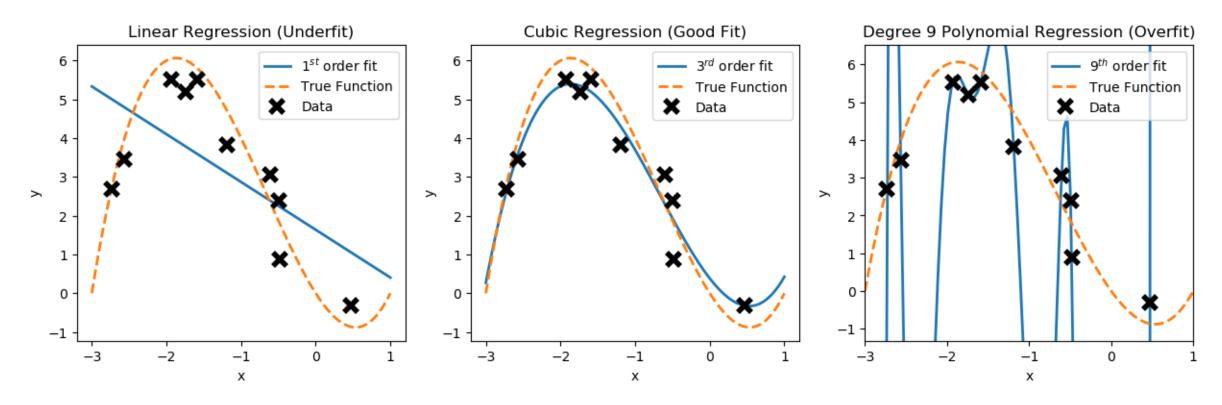


#### WHY SHOULD YOU CARE?

- So far, it seems like you should always prefer the more complex model, right?
- That's because these toy examples assume
  - We have a LOT of data
  - Our data is noiseless
  - Our model function behaves well between our data points
- In the real world, these assumptions are almost always false!

## UNDERFITTING & OVERFITTING

Fitting polynomials to noisy data from the orange function



#### UNDERFITTING & OVERFITTING

- Goal: learn a model that generalizes well to <u>unseen</u> test data
- Underfitting: model is too simple to learn any meaningful patterns in the data – high training error and high test error
- Overfitting: model is so complex that it doesn't generalize well to unseen data because it pays too much attention to the training data low training error but high test error

#### UNDERFITTING & OVERFITTING

- Underfitting is easy to deal with try using a more complex model class because it is more <u>expressive</u>
  - Complexity is roughly the "size" of the function space encoded by a model class (the set of all functions the class can represent)
  - **Expressiveness** is how well that model class can approximate the functions we are interested in
- If a more complex model class overfits, can we reduce its complexity while retaining its expressiveness?

#### REGULARIZATION

- If we make certain structural assumptions about the model we want to learn, we can do just this!
- These assumptions are called <u>regularizers</u>
- Most commonly, we minimize an <u>augmented loss function</u>

$$\tilde{J}(w) = J(w) + \lambda R(w)$$

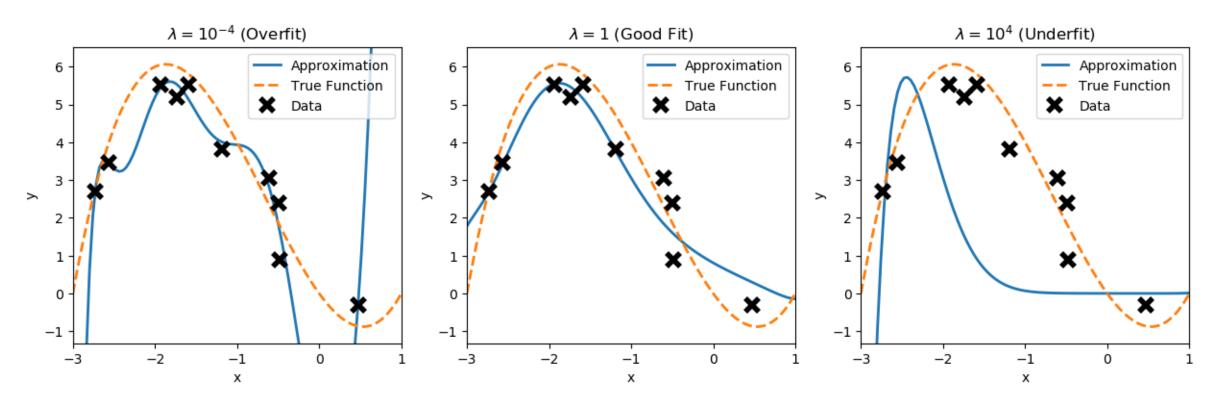
• J(w) is the original loss function,  $\lambda$  is the regularization strength, and R(w) is a regularization term

# ℓ<sub>2</sub> WEIGHT DECAY

- In  $\ell_2$  weight decay regularization,  $R(w) = w^T w = \sum_{k=1}^d w_k^2$
- Minimizing  $\tilde{J}(w) = J(w) + \lambda w^T w$ 
  - Balances the goals of minimizing the loss J(w) and finding a set of weights w that are small in magnitude
  - High  $\lambda$  means we care more about small weights, while low  $\lambda$  means we care more about a low (un-augmented) loss
- Intuitively, small weights  $w \rightarrow$  smoother function (no huge oscillations like the 9<sup>th</sup> degree polynomial we overfit)

# ℓ<sub>2</sub> WEIGHT DECAY

Regularizing a degree 9 polynomial fit with  $\ell_2$  weight decay



#### RETURNING TO NEURAL NETS

- All of the intuition we've built for polynomials is also valid for neural nets!
- The complexity of a deep neural net is related (roughly) to the number of learned parameters and the number of layers
- More complex neural nets, i.e. <u>deeper</u> (more layers) and/or <u>wider</u> (more hidden units) are much more likely to overfit to the training data.

#### RETURNING TO NEURAL NETS

- $\ell_2$  weight decay helps us learn smoother neural nets by encouraging learned weights to be smaller.
- To incorporate  $\ell_2$  weight decay, just do stochastic gradient descent on the augmented loss function

$$\tilde{J}(\mathbf{W}^{(1)}, ..., \mathbf{W}^{(L)}) = J(\mathbf{W}^{(1)}, ..., \mathbf{W}^{(L)}) + \lambda \sum_{i,j,\ell} \mathbf{W}_{ij}^{(\ell)^2}$$

$$\nabla_{\mathbf{W}^{(\ell)}}[\tilde{J}] = \nabla_{\mathbf{W}^{(\ell)}}[J] + 2\lambda \mathbf{W}^{(\ell)}$$

#### NEURAL NETS AND IMAGE DATA

- Let's now consider the special case of doing machine learning on image data with neural nets
- As we've studied them so far, neural nets model relationships between every single pair of pixels
- However, in any image, the color and intensity of neighboring pixels are much more strongly correlated than those of faraway pixels, i.e. images have <u>local structure</u>

#### NEURAL NETS AND IMAGE DATA

- Images are also <u>translation invariant</u>
  - A face is still a face, regardless of whether it's in the top left of an image or the bottom right
- Can we encode these assumptions of local structure into a neural network as a regularizer?
- If we could, we would get models that learned something about our data set <u>as a collection of images</u>.

#### RECAP: CONVOLUTIONS

- Consider a c-by-h-by-w convolutional  $\frac{\text{kernel}}{\text{kernel}}$  or  $\frac{\text{filter}}{\text{filter}}$  array K and a C-by-H-by-W array representing an  $\frac{\text{image}}{\text{kernel}}$
- The convolution (technically cross-correlation)  $\mathbf{Z} = \mathbf{K} \otimes \mathbf{X}$  is

$$\mathbf{Z}[i,j,k] = \sum_{\ell=0}^{c-1} \sum_{m=0}^{h-1} \sum_{n=0}^{w-1} \mathbf{K}[\ell,m,n] \, \mathbf{X}[i+\ell,j+m,k+n]$$

There are multiple ways to deal with boundary conditions;
 for now, ignore any indices that are out of bounds

# RECAP: CONVOLUTIONS (c = 1)

-	100					
0	0	0	0	0	0	
0	105	102	100	97	96	
0	103	99	103	101	102	7
0	101	98	104	102	100	
0	99	101	106	104	99	
0	104	104	104	100	98	

Kernel	Matrix
IXCI IICI	IVIGUIA

0	-1	0
-1	5	-1
0	-1	0

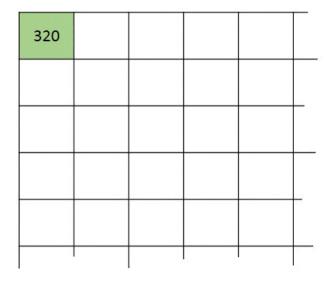


Image Matrix

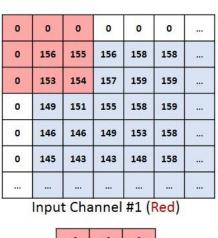
$$0*0+0*-1+0*0$$

$$+0*-1+105*5+102*-1$$

$$+0*0+103*-1+99*0=320$$

**Output Matrix** 

# RECAP: CONVOLUTIONS (c = 3)



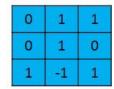
0	0	0	0	0	0	•
0	167	166	167	169	169	
0	164	165	168	170	170	
0	160	162	166	169	170	
0	156	156	159	163	168	
0	155	153	153	158	168	

0	0	0	0	0	0	
0	163	162	163	165	165	
0	160	161	164	166	166	
0	156	158	162	165	166	
0	155	155	158	162	167	
0	154	152	152	157	167	
	744					

Input Channel #2 (Green)

Input Channel #3 (Blue)

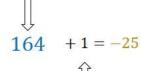
-1	-1	1
0	1	-1
0	1	1



Kernel Channel #1

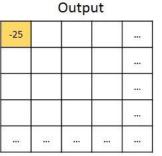
Kernel Channel #2

Kernel Channel #3





Bias = 1



Same source as last figure



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

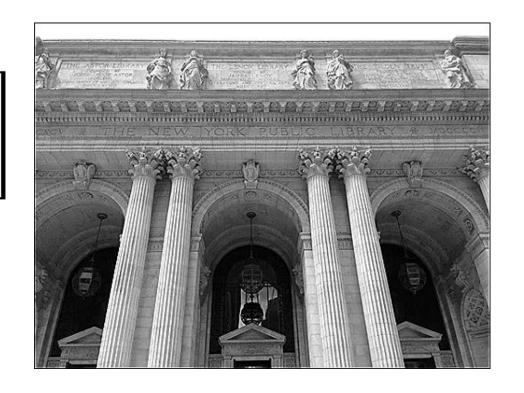






$$\begin{bmatrix} 0 & -1 & 0 \\ -1 & 5 & -1 \\ 0 & -1 & 0 \end{bmatrix}$$

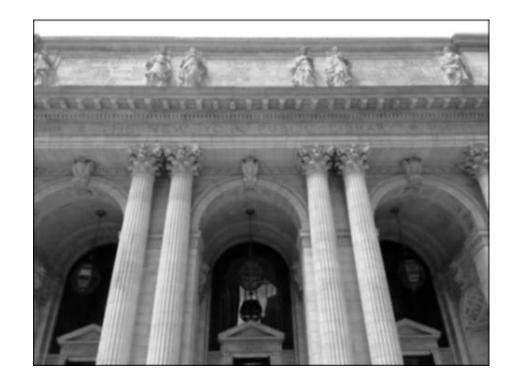






$$\frac{1}{16} \begin{bmatrix} 1 & 2 & 1 \\ 2 & 4 & 2 \\ 1 & 2 & 1 \end{bmatrix}$$

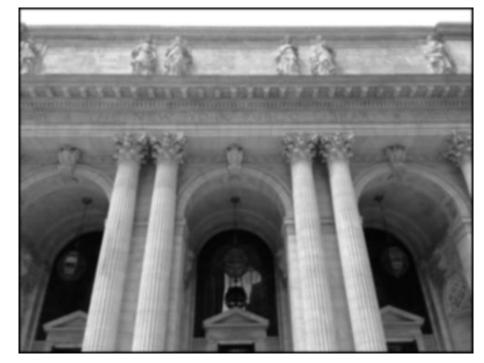






$$\frac{1}{256} \begin{bmatrix}
1 & 4 & 6 & 4 & 1 \\
4 & 16 & 24 & 16 & 4 \\
6 & 24 & 36 & 24 & 6 \\
4 & 16 & 24 & 16 & 4 \\
1 & 4 & 6 & 4 & 1
\end{bmatrix}$$







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







$$\begin{bmatrix} -1 & -1 & -1 \\ -1 & 8 & -1 \\ -1 & -1 & -1 \end{bmatrix}$$





#### ADVANTAGES OF CONVOLUTION

- By sliding the kernel along the image, we can extract the image's <u>local structure</u>!
  - Large objects (by blurring)
  - Sharp edges and outlines
- Since each output pixel of the convolution is highly local, the whole process is also <u>translation invariant</u>!
- Convolution is a <u>linear operation</u>, like matrix multiplication

- So far, the main downside of convolutions is that the coefficients of the kernels seem like magic numbers
- But if we fit a 1D quadratic regression and get the model  $f(x) = 0.382x^2 15.4x + 7$ , then aren't the coefficients 0.382, -15.4, and 7 just magic numbers too?
- Idea: <u>learn convolutional kernels instead of matrices</u> to extract something meaningful from our image data, and then feed that into a dense neural network (with matrices)

- We can do this by creating a new kind of layer, and adding it to the front (closer to the input) of our neural network
- In the forward pass, we convolve our input  $\mathbf{X}^{(\ell-1)}$  with a learned kernel  $\mathbf{K}^{(\ell)}$ , add a scalar bias  $b^{(\ell)}$  to every element of  $\mathbf{Z}^{(\ell)}$ , and apply a nonlinearity  $\theta$  to obtain our output  $\mathbf{X}^{(\ell)}$

$$\mathbf{Z}^{(\ell)} = \mathbf{K}^{(\ell)} \otimes \mathbf{X}^{(\ell-1)} + b^{(\ell)}$$
$$\mathbf{X}^{(\ell)} = \theta(\mathbf{Z}^{(\ell)})$$

- Note that we will actually be attempting to learn multiple (specifically  $c_{\ell}$ ) kernels of shape  $c_{\ell-1} \times h_{\ell} \times w_{\ell}$  per layer  $\ell$ !
  - $c_{\ell-1}$  is the number of channels in input  $\mathbf{X}^{(\ell-1)}$ , so convolving any individual kernel with  $\mathbf{X}^{(\ell-1)}$  will yield 1 output channel
- The output  $\mathbf{X}^{(\ell)}$  is the result of all  $c_{\ell}$  of these convolutions stacked on top of each other (1 output channel per kernel)
- If input  $\mathbf{X}^{(\ell-1)}$  has shape  $c_{\ell-1} \times H_{\ell} \times W_{\ell}$ , then output  $\mathbf{X}^{(\ell)}$  will have shape  $c_{\ell} \times (H_{\ell} h_{\ell} + 1) \times (W_{\ell} w_{\ell} + 1)$

- We then feed the output  $\mathbf{X}^{(\ell)}$  into the next layer as its input
  - If the next layer is a dense layer, we will re-shape  $\mathbf{X}^{(\ell)}$  into a vector (instead of a multi-dimensional array)
- If the next layer is also convolutional, we can pass  $X^{(\ell)}$  as is
- To actually learn good kernels that stage well with the layers we feed them into, we can just use the backpropagation algorithm to do stochastic gradient descent!

- Assume that we have  $\Delta^{(\ell)} = \nabla_{\mathbf{X}^{(\ell)}}[J]$  (the gradient with respect to the input of the next layer, which is also the output of this layer)
- **Proof.** By the chain rule, for each kernel  $\mathbf{K}^{(\ell)}$  at this layer  $\ell$ ,

$$\frac{\partial J}{\partial \mathbf{K}_{ijk}^{(\ell)}} = \sum_{a=1}^{c_{\ell}} \sum_{b=1}^{w_{\ell}} \sum_{c=1}^{h_{\ell}} \frac{\partial J}{\partial \mathbf{Z}_{abc}^{(\ell)}} \frac{\partial \mathbf{Z}_{abc}^{(\ell)}}{\partial \mathbf{K}_{ijk}^{(\ell)}}$$

By the chain rule (again)

$$\frac{\partial J}{\partial \mathbf{Z}_{abc}^{(\ell)}} = \frac{\partial J}{\partial \mathbf{X}_{abc}^{(\ell)}} \frac{\partial \mathbf{X}_{abc}^{(\ell)}}{\partial \mathbf{Z}_{abc}^{(\ell)}} = \Delta_{abc}^{(\ell)} \theta' \left( \mathbf{Z}_{abc}^{(\ell)} \right)$$

- This gives us  $\nabla_{\mathbf{Z}^{(\ell)}}[J]$ , the gradient with respect to the output of the convolution
- We can find this with cudnnActivationBackward() (see Lecture 15) ©

- If you give cuDNN the
  - Gradient with respect to the convolved output  $\nabla_{\mathbf{Z}^{(\ell)}}[J]$
  - Input to the convolution  $\mathbf{X}^{(\ell-1)}$
- cuDNN can compute each  $\nabla_{\mathbf{K}^{(\ell)}}[J]$ , the gradient of the loss with respect to each kernel  $\mathbf{K}^{(\ell)}$  (Lecture 17)  $\odot$
- With the  $\nabla_{\mathbf{K}^{(\ell)}}[J]$ 's computed, we can do gradient descent!

- All that remains is for us to find the gradient with respect to the input to this layer  $\Delta^{(\ell-1)} = \nabla_{\mathbf{X}^{(\ell-1)}}[J]$ 
  - This is also the gradient with respect to the output of the next layer, and will be used to continue doing backpropagation.
- Again, cuDNN has a function for it (Lecture 17)
  - You need to provide it the kernels  $\mathbf{K}^{(\ell)}$  and the gradient with respect to the output  $\Delta^{(\ell)} = \nabla_{\mathbf{X}^{(\ell)}}[J]$  (like a dense neural net)

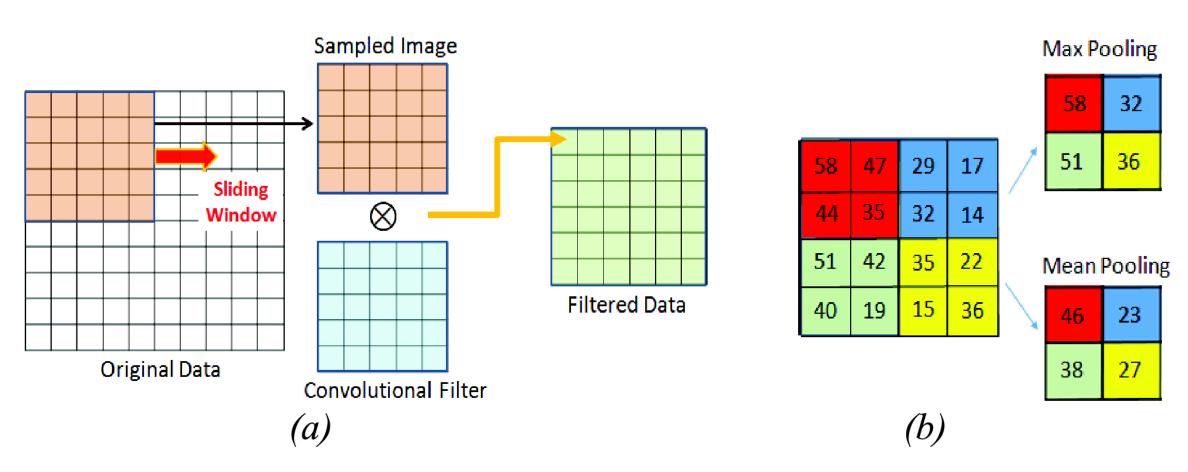
#### POOLING LAYERS

- After each convolutional layer, it is common to add a <u>pooling</u> layer to down-sample the input
- Most commonly, one would take every non-overlapping  $n \times n$  window of a convolved output, and replace each window with
- a single pixel whose intensity is either
  - The maximum intensity found in that  $n \times n$  window
  - The mean intensity of the pixels in that  $n \times n$  window

#### QUICK LINKS FOR CONVOLUTIONAL NNS AND POOLING LAYER

- https://en.wikipedia.org/wiki/Convolutional\_neural\_network
- https://towardsdatascience.com/a-comprehensive-guide-to-convolutional-neural-networks-the-eli5-way-3bd2b1164a53
- https://machinelearningmastery.com/pooling-layers-for-convolutional-neural-networks/
- http://ieeexplore.ieee.org/document/7590035
  - Hand Gesture Recognition Using Micro-Doppler Signatures With Convolutional Neural Network

#### EXAMPLE OF $2 \times 2$ POOLING



http://ieeexplore.ieee.org/document/7590035/all-figures

#### POOLING LAYERS

- Motivation: convolution compresses the amount of information in the image spatially
  - Blur → nearby pixels are more similar
  - Edge  $\rightarrow$  "important" pixels are brighter than their surroundings
- Why not use that compression to reduce dimensionality?
- Forward and backwards propagation for pooling layers are fairly straightforward, and cuDNN can do both (Lecture 17)

#### WHY BOTHER?

- Consider the MNIST dataset of handwritten digits
  - Each image is  $28 \times 28$  pixels  $\rightarrow 784$  input dimensions, and it can be one of 10 output classes
- A neural net with 1 dense hidden layer with 200 units and an dense output layer with 10 units will have  $200 \times (784 + 1) + 10 \times (200 + 1) = 159,010$  parameters
  - We're modeling relationships between every pair of pixels; most of the relationships we learn probably aren't meaningful

- Let's instead consider the following convolutional net:
  - Layer I: Twenty  $(1 \times 5 \times 5)$  kernels
  - Layer 2: 2 × 2 pooling
- Layer 3: Five  $(20 \times 3 \times 3)$  kernels
  - Layer 4: 2 × 2 pooling
  - Layer 5: Dense layer with 10 output units

- Input shape  $(1 \times 28 \times 28)$  (MNIST image)
- Twenty  $(1 \times 5 \times 5)$  kernels
  - $20 \times ((1 \times 5 \times 5) + 1) = 520 \text{ parameters}$
- Output shape  $(20 \times 24 \times 24)$
- 2 × 2 pooling
  - Output shape  $(20 \times 12 \times 12)$

- Input shape  $(20 \times 12 \times 12)$  (conv 1 + pool 1)
- Five  $(20 \times 3 \times 3)$  kernels
  - $5 \times ((20 \times 3 \times 3) + 1) = 905$  parameters
- Output shape  $(5 \times 10 \times 10)$
- $2 \times 2$  pooling
  - Output shape  $(5 \times 5 \times 5)$

- Input shape  $(5 \times 5 \times 5)$  (conv 2 + pool 2)
- Flatten into a 125-dimensional vector
- Dense layer with 50 hidden units
- $10 \times (125 + 1) = 1260$  parameters
  - Output is a 10-dimensional vector

- This gives us a total of 520 + 905 + 1260 = 2685 parameters, far fewer than the dense net's 159,010
- However, with far fewer parameters, this model
- Learns something more meaningful about image structure
  - Achieves a significantly better accuracy on unseen data
- We've effectively regularized the neural net to perform well on image data! HW6: implement it and see for yourself.