#### Topics:

Convolution

## **CS 4644-DL / 7643-A ZSOLT KIRA**

#### Assignment 2

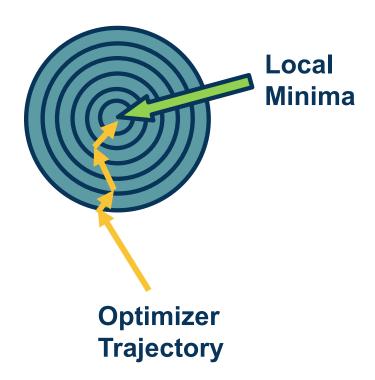
- Implement convolutional neural networks
- Resources (in addition to lectures):
  - DL book: Convolutional Networks
  - CNN notes <a href="https://www.cc.gatech.edu/classes/AY2022/cs7643">https://www.cc.gatech.edu/classes/AY2022/cs7643</a> spring/assets/L10 cnns notes.pdf
  - Backprop notes
    <a href="https://www.cc.gatech.edu/classes/AY2022/cs7643">https://www.cc.gatech.edu/classes/AY2022/cs7643</a> spring/assets/L10 cnns backprop notes.pdf
  - There will be various OH tutorials
  - Slower OMSCS lectures on dropbox: Module 2 Lessons 5-6 (M2L5/M2L6)
     (<a href="https://www.dropbox.com/sh/iviro188gq0b4vs/AADdHxX\_Uy1TkpF\_yvIzX0nPa?dl=0">https://www.dropbox.com/sh/iviro188gq0b4vs/AADdHxX\_Uy1TkpF\_yvIzX0nPa?dl=0</a>)

#### GPU resources

- For assignments, can use CPU or Google Colab
- Projects:
  - Google Cloud Credits

# Even given a good neural network architecture, we need a good optimization algorithm to find good weights

- What optimizer should we use?
  - Different optimizers make different weight updates depending on the gradients
- How should we initialize the weights?
- What regularizers should we use?
- What loss function is appropriate?





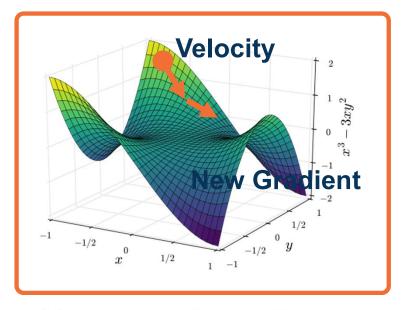
**Key idea:** Rather than combining velocity with current gradient, go along velocity **first** and then calculate gradient at new point

 We know velocity is probably a reasonable direction

$$\widehat{\boldsymbol{w}}_{i-1} = \boldsymbol{w}_{i-1} + \boldsymbol{\beta} \boldsymbol{v}_{i-1}$$

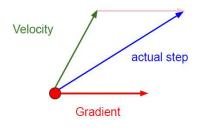
$$v_i = \beta v_{i-1} + \frac{\partial L}{\partial \widehat{w}_{i-1}}$$

$$w_i = w_{i-1} - \alpha v_i$$



#### Momentum update:

#### Nesterov Momentum



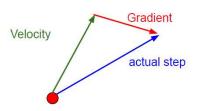


Figure Credit: Fei-Fei Li, Justin Johnson, Serena Yeung, CS 231n

**Nesterov Momentum** 



**Solution:** Time-varying bias correction

Typically  $\beta_1 = 0.9$ ,  $\beta_2 = 0.999$ 

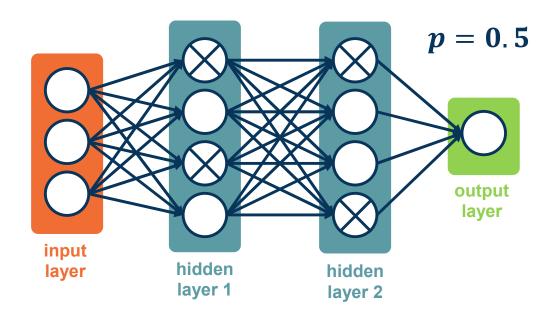
So  $\hat{v_i}$  will be small number divided by (1-0.9=0.1) resulting in more reasonable values (and  $\hat{G}_i$  larger)

$$v_i = \beta_1 v_{i-1} + (1 - \beta_1) \left( \frac{\partial L}{\partial w_{i-1}} \right)$$

$$G_i = \beta_2 G_{i-1} + (1 - \beta_2) \left( \frac{\partial L}{\partial w_{i-1}} \right)^2$$

$$\widehat{v_i} = \frac{v_i}{1 - \beta_1^t} \qquad \widehat{G_i} = \frac{G_i}{1 - \beta_2^t}$$

$$w_i = w_{i-1} - \frac{\alpha \, \widehat{v}_i}{\sqrt{\widehat{G}_i + \epsilon}}$$



**An idea:** For each node, keep its output with probability *p* 

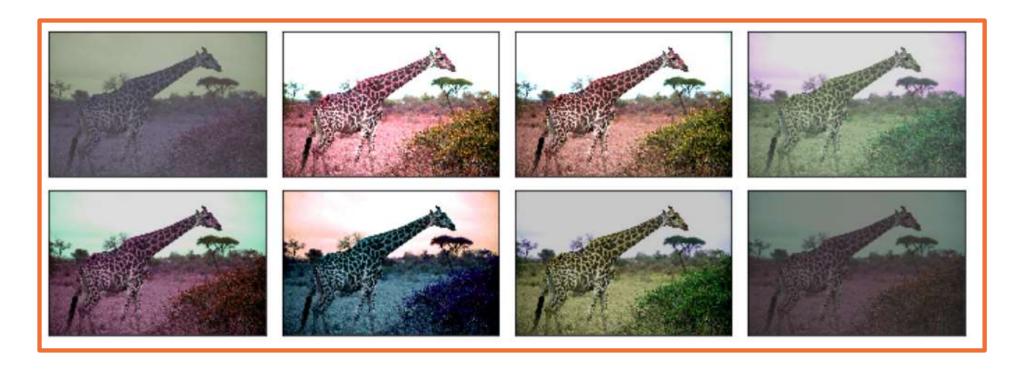
Activations of deactivated nodes are essentially zero

Choose whether to mask out a particular node each iteration

From: Dropout: A Simple Way to Prevent Neural Networks from Overfitting, Srivastava et al.



#### **Color Jitter**



From https://mxnet.apache.org/versions/1.5.0/tutorials/gluon/data\_augmentation.html



- We can give the model flexibility through learnable parameters
   γ (scale) and β (shift)
- Network can learn to not normalize if necessary!
- This layer is called a
   Batch Normalization
   (BN) layer

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

Parameters to be learned: \gamma, \beta

Output: \{y_i = \mathrm{BN}_{\gamma,\beta}(x_i)\}

\mu_{\mathcal{B}} \leftarrow \frac{1}{m} \sum_{i=1}^m x_i \qquad // \text{mini-batch mean}
\sigma_{\mathcal{B}}^2 \leftarrow \frac{1}{m} \sum_{i=1}^m (x_i - \mu_{\mathcal{B}})^2 \qquad // \text{mini-batch variance}
\widehat{x}_i \leftarrow \frac{x_i - \mu_{\mathcal{B}}}{\sqrt{\sigma_{\mathcal{B}}^2 + \epsilon}} \qquad // \text{normalize}
y_i \leftarrow \gamma \widehat{x}_i + \beta \equiv \mathrm{BN}_{\gamma,\beta}(x_i) \qquad // \text{scale and shift}
```

From: Batch Normalization: Accelerating Deep Network Training by Reducing Internal Covariate Shift, Sergey Ioffe, Christian Szegedy



Example: Cross entropy loss

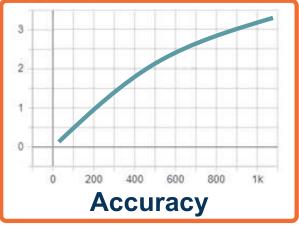
$$L = -\log P(Y = y_i | X = x_i)$$

Accuracy is measured based on:

$$argmax_i(P(Y = y_i | X = x_i))$$

Since the correct class score only has to be slightly higher, we can have flat loss curves but increasing accuracy!







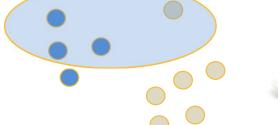
 Precision/Recall curves represent the inherent tradeoff between number of positive predictions and correctness of predictions

#### Definitions

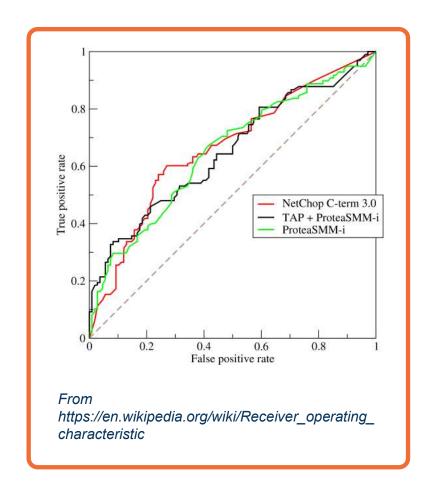
- True Positive Rate:  $TPR = \frac{tp}{tp+fn}$
- False Positive Rate:  $FPR = \frac{fp}{fp+tn}$

$$Accuracy = \frac{tp+tn}{tp+tn+fp+fn}$$



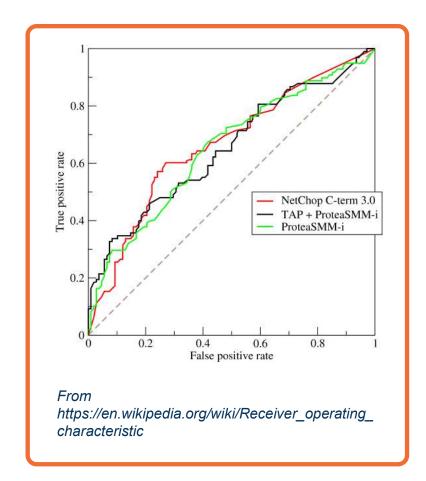








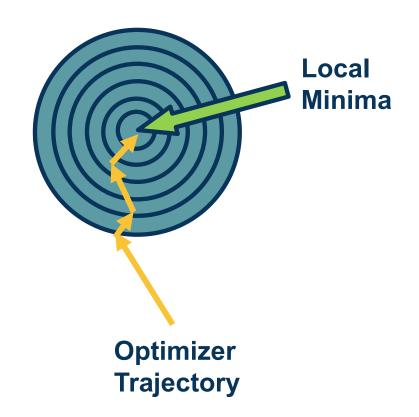
- Precision/Recall curves represent the inherent tradeoff between number of positive predictions and correctness of predictions
- Definitions
  - True Positive Rate:  $TPR = \frac{tp}{tp+fn}$
  - False Positive Rate:  $FPR = \frac{fp}{fp+tn}$
  - $Accuracy = \frac{tp+tn}{tp+tn+fp+fn}$
- We can obtain a curve by varying the (probability) threshold:
  - Area under the curve (AUC) common single-number metric to summarize
- Mapping between this and loss is not simple!





#### Resource:

A disciplined approach to neural network hyper-parameters: Part 1 -- learning rate, batch size, momentum, and weight decay, Leslie N. Smith

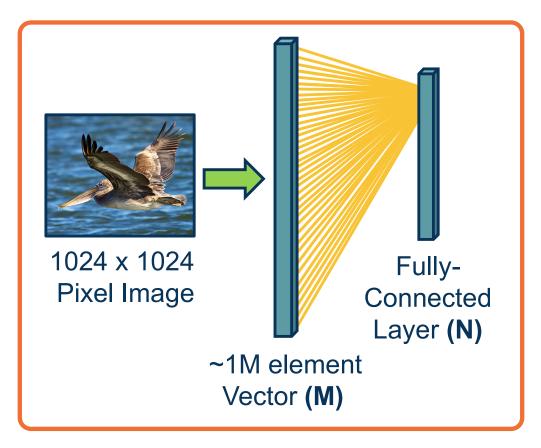




# Convolution & Pooling



#### The connectivity in linear layers doesn't always make sense



**How many parameters?** 

M\*N (weights) + N (bias)

Hundreds of millions of parameters for just one layer

More parameters => More data needed

Is this necessary?



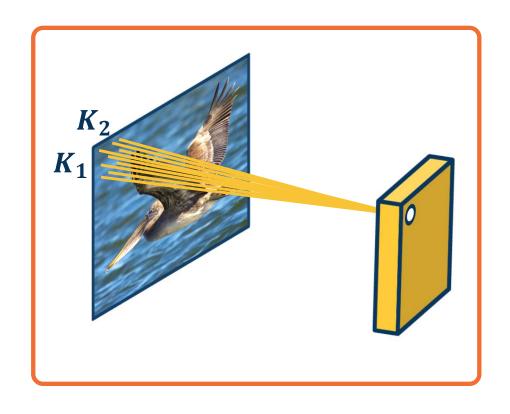
## Image features are spatially localized!

- Smaller features repeated across the image
  - Edges
  - Color
  - Motifs (corners, etc.)
- No reason to believe one feature tends to appear in one location vs. another (stationarity)



Can we induce a *bias* in the design of a neural network layer to reflect this?





Each node only receives input from  $K_1 \times K_2$  window (image patch)

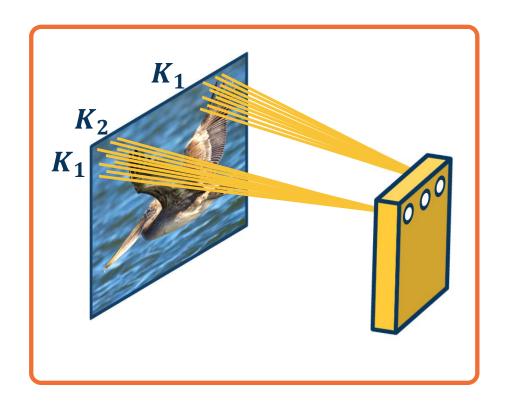
Region from which a node receives input from is called its receptive field

#### **Advantages:**

- Reduce parameters to (K<sub>1</sub>× K<sub>2</sub> + 1)
   \* N where N is number of output nodes
- Explicitly maintain spatial information

Do we need to learn location-specific features?





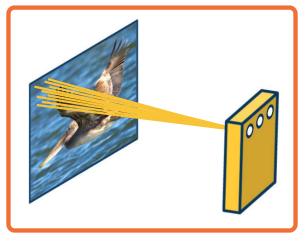
### Nodes in different locations can **share** features

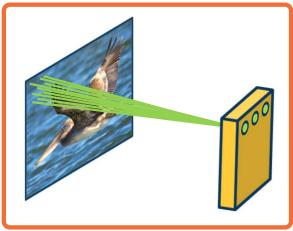
- No reason to think same feature (e.g. edge pattern) can't appear elsewhere
- Use same weights/parameters in computation graph (shared weights)

#### **Advantages:**

- Reduce parameters to  $(K_1 \times K_2 + 1)$
- Explicitly maintain spatial information





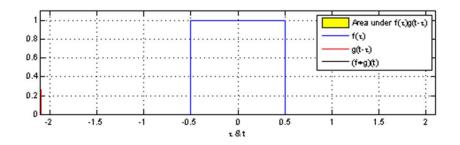


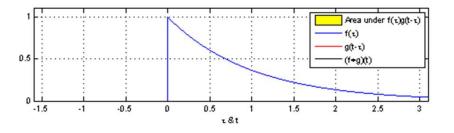
## We can learn **many** such features for this one layer

- Weights are **not** shared across different feature extractors
- Parameters:  $(K_1 \times K_2 + 1)$ \* *M* where *M* is number of features we want to learn



#### This operation is **extremely common** in electrical/computer engineering!

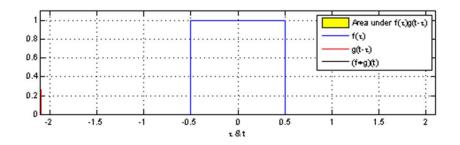


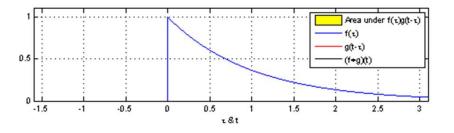


From https://en.wikipedia.org/wiki/Convolution



#### This operation is **extremely common** in electrical/computer engineering!





From https://en.wikipedia.org/wiki/Convolution

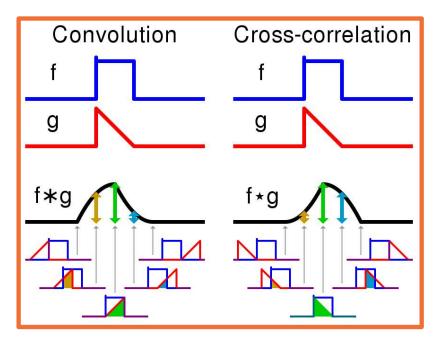


#### This operation is extremely common in electrical/computer engineering!

In mathematics and, in particular, functional analysis, **convolution** is a mathematical operation on two functions f and g producing a third function that is typically viewed as a modified version of one of the original functions, giving the area overlap between the two functions as a function of the amount that one of the original functions is translated.

Convolution is similar to cross-correlation.

It has **applications** that include probability, statistics, computer vision, image and signal processing, electrical engineering, and differential equations.



Visual comparison of **convolution** and **cross-correlation**.

From https://en.wikipedia.org/wiki/Convolution



**Notation:** 

$$F \otimes (G \otimes I) = (F \otimes G) \otimes I$$

1D Convolution 
$$y_k = \sum_{n=0}^{N-1} h_n \cdot x_{k-n}$$

$$y_{0} = h_{0} \cdot x_{0}$$

$$y_{1} = h_{1} \cdot x_{0} + h_{0} \cdot x_{1}$$

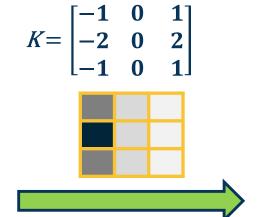
$$y_{2} = h_{2} \cdot x_{0} + h_{1} \cdot x_{1} + h_{0} \cdot x_{2}$$

$$y_{3} = h_{3} \cdot x_{0} + h_{2} \cdot x_{1} + h_{1} \cdot x_{2} + h_{0} \cdot x_{3}$$

$$\vdots$$

**2D** Convolution









**I**mage

Kernel (or filter)

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



Output / filter / feature map



2D Convolution





We will make this convolution operation a layer in the neural network

- Initialize kernel values randomly and optimize them!
- These are our parameters (plus a bias term per filter)

**Image** 

Kernel (or filter)

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



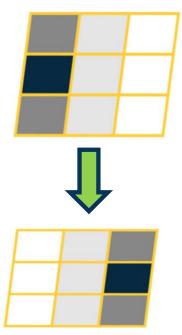
Output / filter / feature map



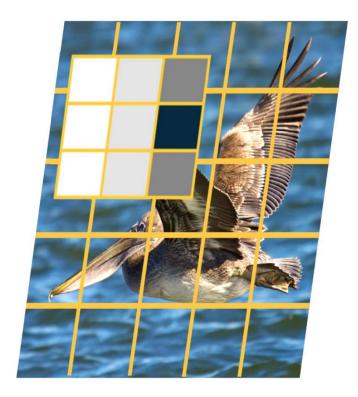
**2D** Convolution

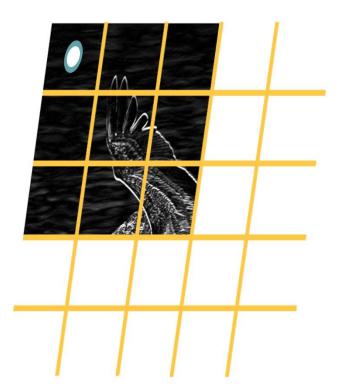


#### 1. Flip kernel (rotate 180 degrees)



## 2. Stride along image







$$y(r,c) = (x*k)(r,c) = \sum_{a=-\frac{H-1}{2}}^{\frac{H-1}{2}} \sum_{b=-\frac{W-1}{2}}^{\frac{W-1}{2}} x(a,b) k(r-a,c-b)$$

$$(0,0)$$

$$k_1 = 3$$

$$w = 5 \qquad (\frac{H-1}{2}, \frac{W-1}{2})$$

$$y(0,0) = x(-2,-2)k(2,2) + x(-2,-1)k(2,1) + x(-2,0)k(2,0) + x(-2,1)k(2,-1) + x(-2,2)k(2,-2) + \dots$$



$$y(r,c) = (x*k)(r,c) = \sum_{a=-\frac{k_1-1}{2}}^{\frac{k_1-1}{2}} \sum_{b=-\frac{k_2-1}{2}}^{\frac{k_2-1}{2}} x(r-a,c-b) k(a,b)$$

$$(0,0)$$

$$(-\frac{k_1-1}{2}, -\frac{k_2-1}{2})$$

$$k_1 = 3$$

$$k_2 = 3 \quad \underbrace{k_1-1}_{2}, \underbrace{k_2-1}_{2}$$

$$W = 5 \quad (H-1,W-1)$$

#### As we have seen:

- Convolution: Start at end of kernel and move back
- Cross-correlation: Start in the beginning of kernel and move forward (same as for image)

#### An **intuitive interpretation** of the relationship:

- Take the kernel, and rotate 180 degrees along center (sometimes referred to as "flip")
- Perform cross-correlation
- (Just dot-product filter with image!)

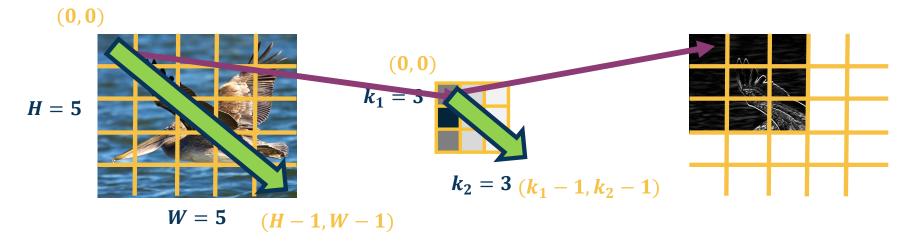
$$K = \begin{bmatrix} 1 & 2 & 3 \\ 4 & 5 & 6 \\ 7 & 8 & 9 \end{bmatrix}$$



$$K' = \begin{bmatrix} 9 & 8 & 7 \\ 6 & 5 & 4 \\ 3 & 2 & 1 \end{bmatrix}$$



$$y(r,c) = (x * k)(r,c) = \sum_{a=0}^{k_1-1} \sum_{b=0}^{k_2-1} x(r+a,c+b) k(a,b)$$



Since we will be learning these kernels, this change does not matter!



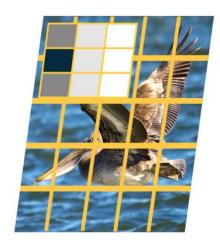
$$X(0:2,0:2) = \begin{bmatrix} 200 & 150 & 150 \\ 100 & 50 & 100 \\ 25 & 25 & 10 \end{bmatrix} \qquad K' = \begin{bmatrix} 1 & 0 & -1 \\ 2 & 0 & -2 \\ 1 & 0 & -1 \end{bmatrix}$$

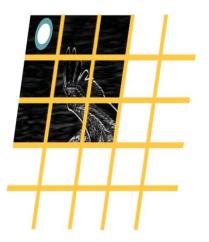
$$\mathsf{K}' = \begin{bmatrix} 1 & 0 & -1 \\ 2 & 0 & -2 \\ 1 & 0 & -1 \end{bmatrix}$$



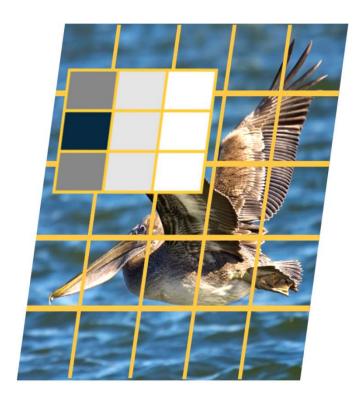
 $X(0:2,0:2) \cdot K' = 65 + bias$ 

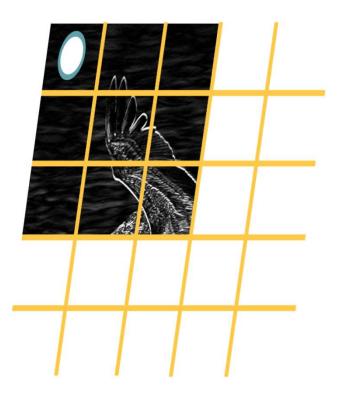
Dot product (element-wise multiply and sum)



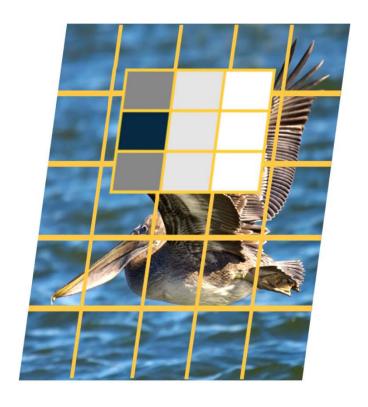


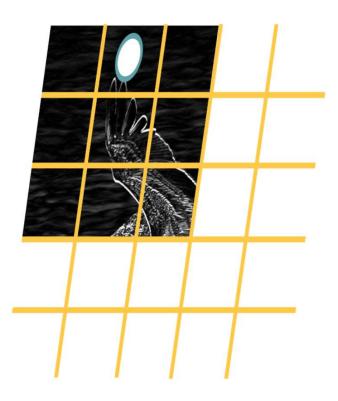






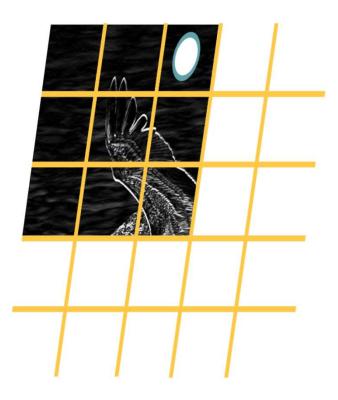




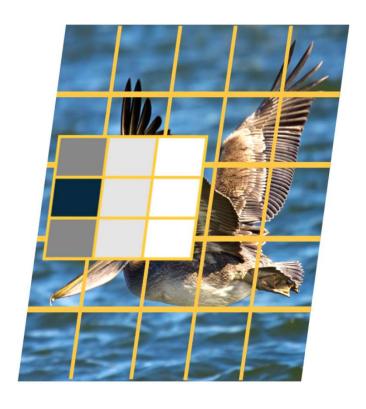


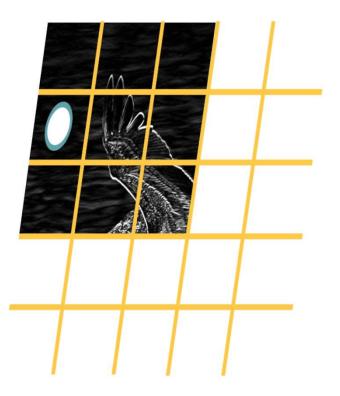




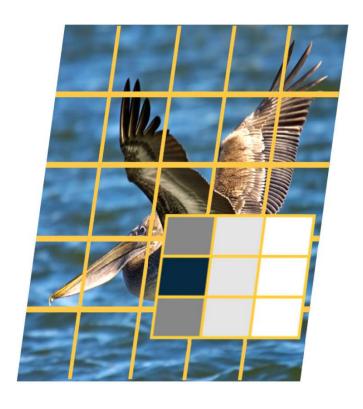


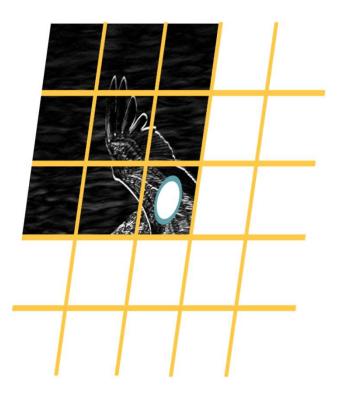












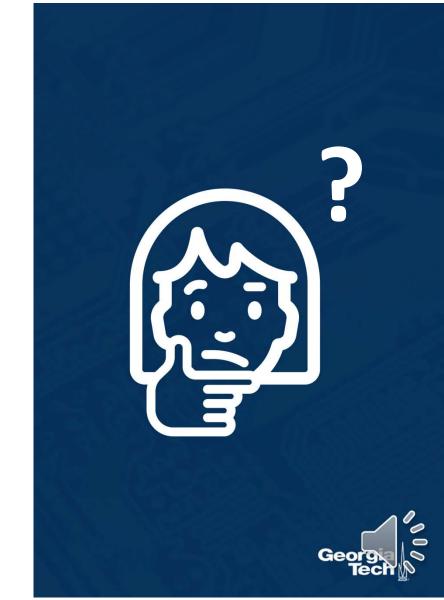


#### **Why Bother with Convolutions?**

Convolutions are just simple linear operations

Why bother with this and not just say it's a linear layer with small receptive field?

- There is a duality between them during backpropagation
- Convolutions have various mathematical properties people care about
- This is historically how it was inspired



# Input & Output Sizes

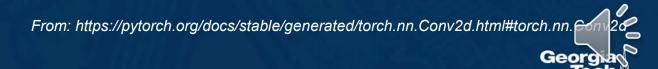


## **Convolution Layer Hyper-Parameters**

#### **Parameters**

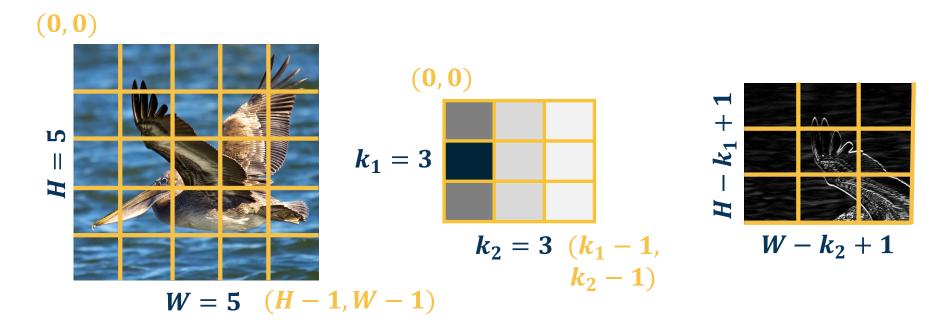
- in\_channels (int) Number of channels in the input image
- out\_channels (int) Number of channels produced by the convolution
- kernel\_size (int or tuple) Size of the convolving kernel
- stride (int or tuple, optional) Stride of the convolution. Default: 1
- padding (int or tuple, optional) Zero-padding added to both sides of the input. Default: 0
- padding\_mode (string, optional) 'zeros', 'reflect', 'replicate' or 'circular'. Default: 'zeros'

Convolution operations have several hyper-parameters



**Output size** of vanilla convolution operation is  $(H - k_1 + 1) \times (W - k_2 + 1)$ 

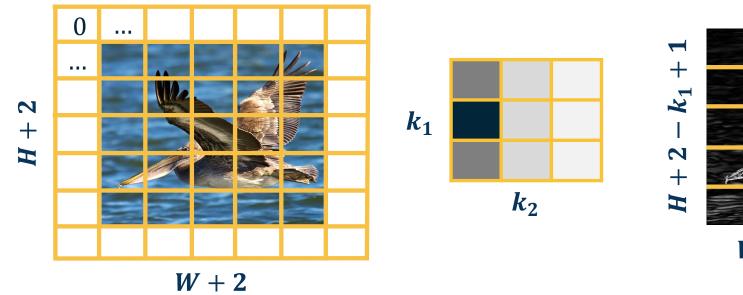
This is called a "valid" convolution and only applies kernel within image

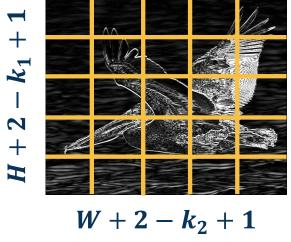




We can pad the images to make the output the same size:

- Zeros, mirrored image, etc.
- ullet Note padding often refers to pixels added to **one size** (P = 1 here)



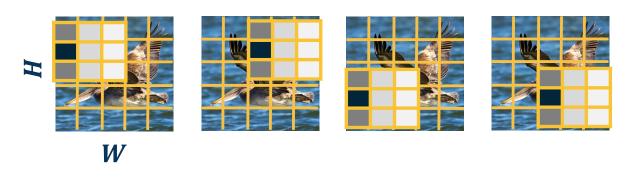


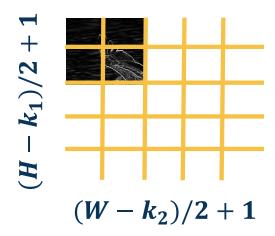


We can move the filter along the image using larger steps (stride)

- This can potentially result in loss of information
- Can be used for dimensionality reduction (not recommended)

### Stride = 2 (every other pixel)







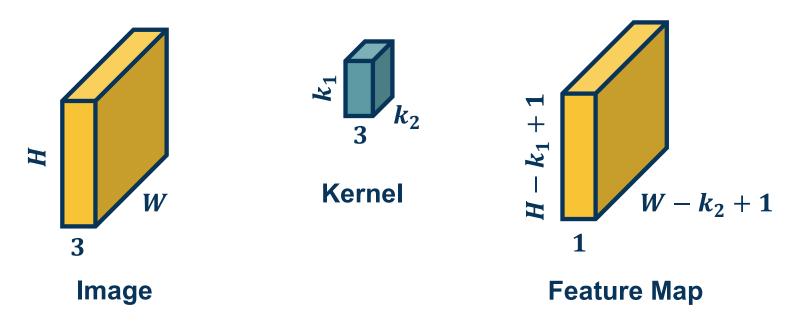
# Stride can result in **skipped pixels**, e.g. stride of 3 for 5x5 input





We have shown inputs as a **one-channel image** but in reality they have three channels (red, green, blue)

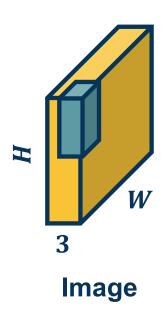
In such cases, we have 3-channel kernels!





We have shown inputs as a **one-channel image** but in reality they have three channels (red, green, blue)

In such cases, we have 3-channel kernels!



Similar to before, we perform **element-wise multiplication** between kernel and image patch, summing them up **(dot product)** 

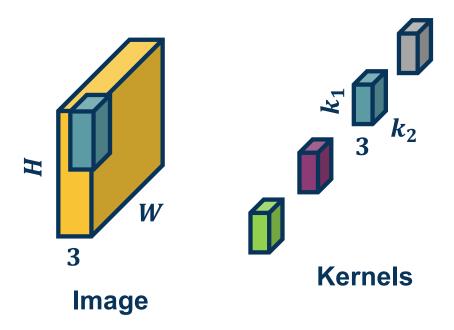
Except with  $k_1 * k_2 * 3$  values

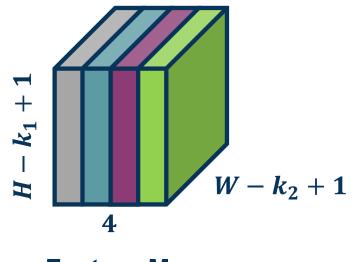


#### We can have multiple kernels per layer

We stack the feature maps together at the output

Number of channels in output is equal to *number* of kernels





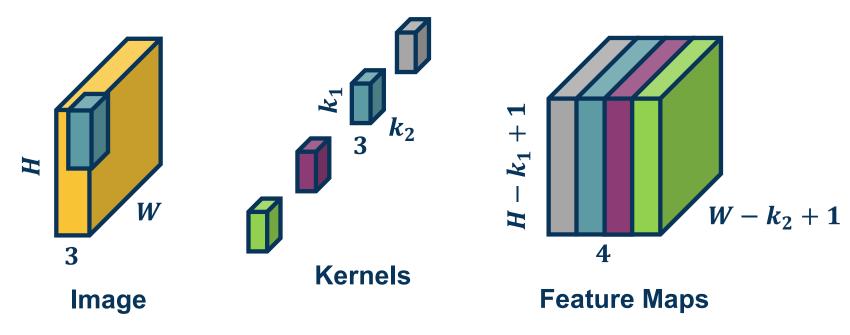
**Feature Maps** 



Number of parameters with N filters is:  $N * (k_1 * k_2 * 3 + 1)$ 

#### Example:

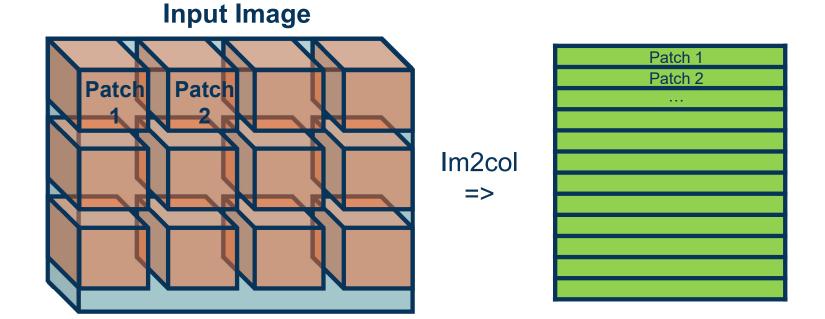
$$k_1 = 3, k_2 = 3, N = 4 input channels = 3, then  $(3 * 3 * 3 + 1) * 4 = 112$$$





Just as before, in practice we can vectorize this operation

Step 1: Lay out image patches in vector form (note can overlap!)

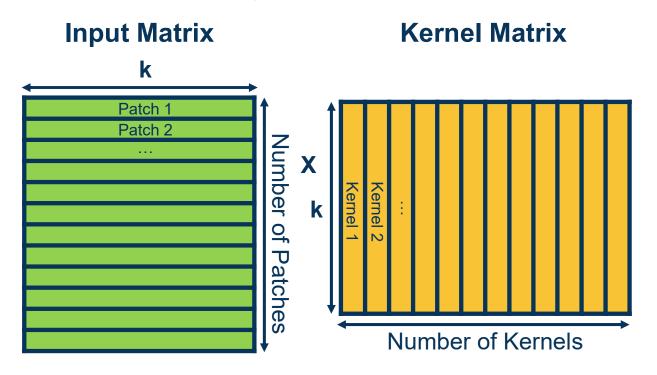


Adapted from: https://petewarden.com/2015/04/20/why-gemm-is-at-the-heart-of-deep-learning/



Just as before, in practice we can vectorize this operation

Step 2: Multiple patches by kernels



Adapted from: https://petewarden.com/2015/04/20/why-gemm-is-at-the-heart-of-deep-learning/

