

# Your Neural Network for NLP is Probably Wrong

Why You Need to Mask More Than You Think

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Interactions

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## Slides

This QR code has a link to the slides if you want to follow along.



- This has links to the slides, the slides include the speaker notes which has a lot of the extra information I am going to say in this talk.
- The speaker notes also have links to where I got the images and a few extra blog posts that can supplement my explanations on how these layers work.

## Cats



- This QR code as a link to a compilation video of cats getting scared by cucumbers.
- I figured I would give it to you in case you want to watch this instead of paying attention to the talk.
- Side Note: I have heard that this is very upsetting to the cats so I wouldn't recommend doing this to your cat, but, given that these people already filmed it... I don't think it is *that* bad to have a laugh at it.

## Me

- I work on NLP at Interactions.
- My specialization is in Deep Learning.
- I maintain **Mead-Baseline**.

- This is my “Why you should listen to me” slide.
- I do NLP at interactions where we use NLP to “facilitate customer care interactions” aka build better chat bots.
- My specialization is DL and we use DL for most models so I have a lot of experience.
- I help maintain Mead-Baseline, our open-source Deep Learning modeling software so I have a lot of experience writing Neural Networks for NLP.
- I have batched some complex operations for Mead-Baseline.
  - Batched the CRF
  - Batched Beam Search
- I have tracked down a lot of hidden batch instability problems.
- I am trying to get more involved with research so if anyone needs help with some idea, especially if the implementation is tricky, hit me up.

## Binary Logistic Regression

$n :=$  number of features

$c :=$  number of class  $:= 1$

$f \in \mathbb{R}^n$

$w \in \mathbb{R}^n$

$$s = \sum_{i=0}^n f_i * w_i$$

- First I'm going to build up to batching by doing more and more work in a single function call. We start with Binary LR
- In binary LR we have a vector representing features and a vector of the same size representing weights
- (Normally in NLP we have sparse features where we don't have a actual vector of features, we would use something like a hashmap instead, but inside a neural network everything is dense vectors so we will use dense vectors here because it is just a pedagogical example)
- We take the sum of the features weighted by the weight to create a score which we call a logit.
- This operation is called the dot product.
- Normally we then use a function like sigmoid to create a final probability (and then calculate losses and gradients) but we don't need to worry about that for this example.

## Binary Logistic Regression

```
s = np.dot(f, w)
```

- In this example we are representing our feature and weights as numpy vectors
- We can use the dot product function to calculate the sum of the elementwise product of the vectors.

## Binary Logistic Regression

$$f = \begin{bmatrix} 1 & 2 & 3 & 4 \end{bmatrix}$$
$$w = \begin{bmatrix} 5 & 6 & 7 & 8 \end{bmatrix}$$
$$s =$$

- Here we can see our feature vector and our weight vector
- Now lets look at the mechanics of what is happening in our dot product calculation

## Binary Logistic Regression

$$\begin{aligned}f &= \begin{bmatrix} 1 & 2 & 3 & 4 \end{bmatrix} \\w &= \begin{bmatrix} 5 & 6 & 7 & 8 \end{bmatrix} \\s &= 5\end{aligned}$$

- We multiply these terms together and add the result into the s



## Binary Logistic Regression

$$\begin{aligned} f &= \begin{bmatrix} 1 & 2 & 3 & 4 \end{bmatrix} \\ w &= \begin{bmatrix} 5 & 6 & 7 & 8 \end{bmatrix} \\ s &= 17 \end{aligned}$$

- And it continues on like this.

## Binary Logistic Regression

$$\begin{aligned} f &= [ \boxed{1 \ 2 \ 3 \ 4} ] \\ w &= [ \boxed{5 \ 6 \ 7 \ 8} ] \\ s &= 70 \end{aligned}$$

- Here we highlight the elements in these arrays that interact with each other in our dot product.

## Multi-class Logistic Regression

$n$  := number of features

$c$  := number of class

$$f \in \mathbb{R}^n$$

$$W \in \mathbb{R}^{n \times c}$$

$$s \in \mathbb{R}^c$$

$$\forall j \in c \quad s_j = \sum_{i=0}^n f_i * W_{ij}$$

- We now extend to multi-class LR
  - In binary LR we are deciding between two things, for example Positive or Negative sentiment
  - In multi-class LR we are deciding between multi classes, for example Science, Technology, Politics, cooking, and video games for document classification
- We still have a feature vector for some document.
- We now have multiple weight vectors, one for each class
- We pack our weight vectors into a single matrix of size n (the number of features we have) by c (which is the number of classes we have).
- Now our logits are a vector, we have a score for each class
- for each class we do this dot product between the feature vector and one of the weight vectors
- Once we have the score we normally turn it into probabilities with the softmax function but we can skip this for this example.

## Multi-class Logistic Regression

```
s = []  
for w in W.T:  
    s.append(np.dot(f, w))  
s = np.array(s)
```

- Here we see code that calculates the logits how I described it before.
- where we explicitly do the dot between the feature and each weight vector in turn
- We turn the results into a vector at the end

## Multi-class Logistic Regression

```
s = np.dot(f, W)
```

- But it turns out we can actually do this in a single dot product!  
Lets see why

## Multi-class Logistic Regression

$$f = \begin{bmatrix} 1 & 2 & 3 & 4 \end{bmatrix}$$
$$W = \begin{bmatrix} 5 & 9 \\ 6 & 10 \\ 7 & 11 \\ 8 & 12 \end{bmatrix}$$
$$s = \begin{bmatrix} \end{bmatrix}$$

- Here we can see our feature vector again and see our new weight matrix
- Note how the weight matrix is two weight vectors next to each other
- The first column is the same weight vector we saw earlier
- Lets explore how the elements of the feature vector and weight matrix are combined

## Multi-class Logistic Regression

$$f = \begin{bmatrix} 1 & 2 & 3 & 4 \end{bmatrix}$$
$$W = \begin{bmatrix} 5 & 9 \\ 6 & 10 \\ 7 & 11 \\ 8 & 12 \end{bmatrix}$$
$$s = \begin{bmatrix} 5 \end{bmatrix}$$

- Here we can see that we multiply an element from the feature vector and the first column of the weight matrix
- We save this value into the score vector

## Multi-class Logistic Regression

$$f = \begin{bmatrix} 1 & 2 & 3 & 4 \end{bmatrix}$$
$$W = \begin{bmatrix} 5 & 9 \\ 6 & 10 \\ 7 & 11 \\ 8 & 12 \end{bmatrix}$$
$$s = \begin{bmatrix} 17 \end{bmatrix}$$

- And it goes on like this...



## Multi-class Logistic Regression

$$f = [1 \ 2 \ 3 \ 4]$$
$$W = \begin{bmatrix} 5 & 9 \\ 6 & 10 \\ 7 & 11 \\ 8 & 12 \end{bmatrix}$$
$$s = [70]$$

- We see here that the value in this first spot of the logits was calculated between the feature vector and the first column in the weight matrix

## Multi-class Logistic Regression

- Here we see the elements of the feature vector and the second column weight matrix are combined and saved into the second element of the logit vector

$$f = [ \textcircled{1} \ 2 \ 3 \ 4 ]$$
$$W = \begin{bmatrix} 5 & \textcircled{9} \\ 6 & 10 \\ 7 & 11 \\ 8 & 12 \end{bmatrix}$$
$$s = [ \ 70 \ \textcircled{9} ]$$

## Multi-class Logistic Regression

$$f = \begin{bmatrix} 1 & 2 & 3 & 4 \end{bmatrix}$$
$$W = \begin{bmatrix} 5 & 9 \\ 6 & 10 \\ 7 & 11 \\ 8 & 12 \end{bmatrix}$$
$$s = \begin{bmatrix} 70 & 29 \end{bmatrix}$$

- And it goes on like this...

## Multi-class Logistic Regression

$$f = \begin{bmatrix} 1 & 2 & 3 & 4 \end{bmatrix}$$
$$W = \begin{bmatrix} 5 & 9 \\ 6 & 10 \\ 7 & 11 \\ 8 & 12 \end{bmatrix}$$
$$s = \begin{bmatrix} 70 & 110 \end{bmatrix}$$

- We see here that the value in the second spot of the logits was calculated between the feature vector and the second column in the weight matrix
- So we can see here that we can calculate the logits for multiple classes at once between there is no interaction between the columns of the matrix in vector-matrix multiplication.

## Batched Multi-class Logistic Regression

$n$  := number of features

$c$  := number of class

$b$  := number of examples

$$F \in \mathbb{R}^{b \times n}$$

$$W \in \mathbb{R}^{n \times c}$$

$$S \in \mathbb{R}^{b \times c}$$

$$\forall_{k \in b} \forall_{j \in c} S_{kj} = \sum_{i=0}^n F_{ki} * W_{ij}$$

- We just saw how you can calculate the scores for multiple classes for a single example in a single vector-matrix multiply. Now we will see how you can actually do this for multiple examples at once.
- Here we see Batched Multi-class Logistic Regression.
- This means that we are doing multi-class LR but will process multiple examples at once
- Like before we still have a weight matrix of  $n$  by  $c$
- In this case we have  $b$  examples
- Now we have stacked multiple feature vectors into a feature matrix with  $b$  rows where each row represents a different feature vector.
- Now our scores are a matrix where each row is the scores for an example and each column is the score for some class
- We could calculate this by doing the dot product of each feature vector and the weight matrix and stacking the results into our score matrix

## Batched Multi-class Logistic Regression

```
S = []  
for f in F:  
    S.append(np.dot(f, W))  
S = np.stack(S)
```

- Here we see a straight forward way to do this, where we loop over the features and calculate the scores for each example separately.
- The dot product of the feature vector and weight matrix returns a vector
- We collect these vectors and stack them into a single array

## Batched Multi-class Logistic Regression

```
S = np.dot(F, W)
```

- But we can actually do one better, we can calculate the whole thing, all classes for all examples in a single shot!
- Lets take a closer look

## Batched Multi-class Logistic Regression

$$F = \begin{bmatrix} 1 & 2 & 3 & 4 \\ 13 & 14 & 15 & 16 \end{bmatrix}$$
$$W = \begin{bmatrix} 5 & 9 \\ 6 & 10 \\ 7 & 11 \\ 8 & 12 \end{bmatrix}$$
$$s = \begin{bmatrix} \\ \\ \\ \end{bmatrix}$$

- Here we see our new feature matrix is two feature vectors stacked on top of each other
- Our weight matrix is still a n by c matrix of weight vectors standing next to each other
- We are now doing matrix matrix multiplication



## Batched Multi-class Logistic Regression

$$F = \begin{bmatrix} 1 & 2 & 3 & 4 \\ 13 & 14 & 15 & 16 \end{bmatrix}$$
$$W = \begin{bmatrix} 5 & 9 \\ 6 & 10 \\ 7 & 11 \\ 8 & 12 \end{bmatrix}$$
$$s = \begin{bmatrix} 5 \\ \end{bmatrix}$$

- Here we see the values for the first class of the first example are computed using the values in the first row of the feature vector and the first column of the weight matrix

## Batched Multi-class Logistic Regression

$$F = \begin{bmatrix} 1 & 2 & 3 & 4 \\ 13 & 14 & 15 & 16 \end{bmatrix}$$
$$W = \begin{bmatrix} 5 & 9 \\ 6 & 10 \\ 7 & 11 \\ 8 & 12 \end{bmatrix}$$
$$s = \begin{bmatrix} 17 \\ \end{bmatrix}$$

- Here we see the values for the first class of the first example are computed using the values in the first row of the feature vector and the first column of the weight matrix

## Batched Multi-class Logistic Regression

$$F = \begin{bmatrix} 1 & 2 & 3 & 4 \\ 13 & 14 & 15 & 16 \end{bmatrix}$$
$$W = \begin{bmatrix} 5 & 9 \\ 6 & 10 \\ 7 & 11 \\ 8 & 12 \end{bmatrix}$$
$$s = \begin{bmatrix} 70 \end{bmatrix}$$

- Here we see the values for the first class of the first example are computed using the values in the first row of the feature vector and the first column of the weight matrix

## Batched Multi-class Logistic Regression

$$F = \begin{bmatrix} 1 & 2 & 3 & 4 \\ 13 & 14 & 15 & 16 \end{bmatrix}$$
$$W = \begin{bmatrix} 5 & 9 \\ 6 & 10 \\ 7 & 11 \\ 8 & 12 \end{bmatrix}$$
$$s = \begin{bmatrix} 70 & 9 \end{bmatrix}$$

- Here we see the values for the second class of the first example are computed using the values in the first row of the feature vector and the second column of the weight matrix

## Batched Multi-class Logistic Regression

$$F = \begin{bmatrix} 1 & 2 & 3 & 4 \\ 13 & 14 & 15 & 16 \end{bmatrix}$$
$$W = \begin{bmatrix} 5 & 9 \\ 6 & 10 \\ 7 & 11 \\ 8 & 12 \end{bmatrix}$$
$$s = \begin{bmatrix} 70 & 29 \end{bmatrix}$$

- Here we see the values for the second class of the first example are computed using the values in the first row of the feature vector and the second column of the weight matrix

## Batched Multi-class Logistic Regression

$$F = \begin{bmatrix} 1 & 2 & 3 & 4 \\ 13 & 14 & 15 & 16 \end{bmatrix}$$
$$W = \begin{bmatrix} 5 & 9 \\ 6 & 10 \\ 7 & 11 \\ 8 & 12 \end{bmatrix}$$
$$s = \begin{bmatrix} 70 & 110 \end{bmatrix}$$

- Here we see the values for the second class of the first example are computed using the values in the first row of the feature vector and the second column of the weight matrix

## Batched Multi-class Logistic Regression

$$F = \begin{bmatrix} 1 & 2 & 3 & 4 \\ 13 & 14 & 15 & 16 \end{bmatrix}$$
$$W = \begin{bmatrix} 5 & 9 \\ 6 & 10 \\ 7 & 11 \\ 8 & 12 \end{bmatrix}$$
$$s = \begin{bmatrix} 70 & 110 \\ 65 & \end{bmatrix}$$

- Here we see the values for the first class of the second example are computed using the values in the second row of the feature vector and the first column of the weight matrix

## Batched Multi-class Logistic Regression

$$F = \begin{bmatrix} 1 & 2 & 3 & 4 \\ 13 & 14 & 15 & 16 \end{bmatrix}$$
$$W = \begin{bmatrix} 5 & 9 \\ 6 & 10 \\ 7 & 11 \\ 8 & 12 \end{bmatrix}$$
$$s = \begin{bmatrix} 70 & 110 \\ 149 & \end{bmatrix}$$

- Here we see the values for the first class of the second example are computed using the values in the second row of the feature vector and the first column of the weight matrix



## Batched Multi-class Logistic Regression

$$F = \begin{bmatrix} 1 & 2 & 3 & 4 \\ 13 & 14 & 15 & 16 \end{bmatrix}$$
$$W = \begin{bmatrix} 5 & 9 \\ 6 & 10 \\ 7 & 11 \\ 8 & 12 \end{bmatrix}$$
$$s = \begin{bmatrix} 70 & 110 \\ 38 & 2 \end{bmatrix}$$

- Here we see the values for the first class of the second example are computed using the values in the second row of the feature vector and the first column of the weight matrix

## Batched Multi-class Logistic Regression

$$F = \begin{bmatrix} 1 & 2 & 3 & 4 \\ 13 & 14 & 15 & 16 \end{bmatrix}$$
$$W = \begin{bmatrix} 5 & 9 \\ 6 & 10 \\ 7 & 11 \\ 8 & 12 \end{bmatrix}$$
$$s = \begin{bmatrix} 70 & 110 \\ 382 & 117 \end{bmatrix}$$

- Here we see the values for the second class of the second example are computed using the values in the second row of the feature vector and the second column of the weight matrix

## Batched Multi-class Logistic Regression

- Here we see the values for the second class of the second example are computed using the values in the second row of the feature vector and the second column of the weight matrix

$$F = \begin{bmatrix} 1 & 2 & 3 & 4 \\ 13 & 14 & 15 & 16 \end{bmatrix}$$

$$W = \begin{bmatrix} 5 & 9 \\ 6 & 10 \\ 7 & 11 \\ 8 & 12 \end{bmatrix}$$

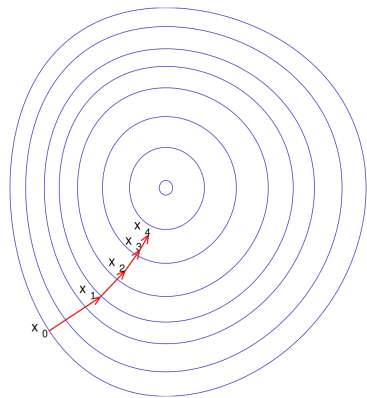
$$s = \begin{bmatrix} 70 & 110 \\ 382 & 117 \end{bmatrix}$$

## Batched Multi-class Logistic Regression

$$F = \begin{bmatrix} 1 & 2 & 3 & 4 \\ 13 & 14 & 15 & 16 \end{bmatrix}$$
$$W = \begin{bmatrix} 5 & 9 \\ 6 & 10 \\ 7 & 11 \\ 8 & 12 \end{bmatrix}$$
$$s = \begin{bmatrix} 70 & 110 \\ 382 & 614 \end{bmatrix}$$

- And finally the last elements uses the two unused rows and columns.
- So here we have seen that we can stack independent examples together just like we could stack together independent weight vectors to calculate the scores for multiple examples at once.
- This works because the rows and columns in matrix-matrix multiplication all act independently of each other.
- This means we can put totally unrelated data an be next to each other and they won't interact as it flows through your network
- Running multiple examples like this is what we mean by batching.

## Full Gradient Descent



### Pros:

- This is the true gradient
- Your step is guaranteed to yield better performance on all examples

### Cons:

- It is slow
- You need to run each example before you update any parameters

- The way we train these models is with something called gradient descent
- This basically means run an example through, calculate how wrong the predicted answer is (the loss), and use the gradient (the derivative) of the parameters with respect to this loss to calculate how to update the parameters.
- In some cases people do full gradient descent where they calculate the answers for each example and get the gradient for everything
- this is good because this is the true gradient, this is the best step you can take to do better next time
- This is slow because you have to wait to run the network on all your data before you update any parameters
- In this picture we can see the big strides in the right direction
- [Image from here](#)

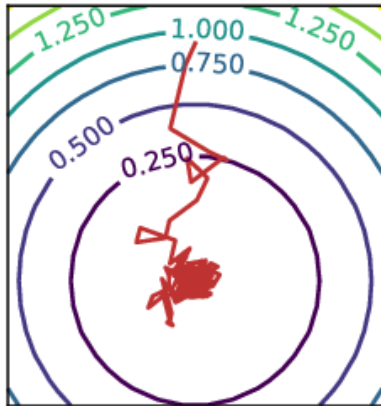
# Stochastic Gradient Descent

## Pros:

- You get to update your parameters each step.
- You can get better generalization because to the randomness in steps.

## Cons:

- Your gradient could be wrong.
- The gradient for one example can cause a change that is detrimental to another example.



- On the opposite end of the spectrum we have stochastic gradient descent
- We calculate a gradient and update parameters after each example
- Talk about pros and cons
- In this image we can see how much more random the gradients are
- [Image from here](#)

## Mini-Batched Gradient Descent

### Pros:

- You get to update your parameters more often.
- You can a better approximation of your gradient.

### Cons:

- You need to run the examples in a batch which can be tricky.
- Minibatch size is now a hyperparameter.

- This is a happy middle ground between the two
- We calculate the gradient for a small batch of examples and update
- Talk about pros and cons
- We can run this mini-batch together via batching

## Gradient Accumulation?

- We saw before that we want to have batches for training dynamics
- Why do you have to actually run things at once as a single batch?
- Can't you just run each example individually and accumulate the gradients in memory and apply them when you get enough?
- Given that I am giving a whole talk on why batching is hard shouldn't you try to avoid it?

- Gradient Accumulation is used when your networks are really large and can't fit on a single GPU
- But we have a good reason to actually batch and that is...



## Speed



- [Image from here](#)
- We showed before that we can express batched operations as matrix matrix multiplications
- Lucky for us humanity has gotten really good at doing these matmuls fast
- We have specialized libraries and even specialized devices, like GPUs, that are designed to do these in a parallel and fast way
- We can stand on the shoulders of gaming to train models really fast

## Batching in NLP

[The dog ran very fast]

- Now that we know batching is a good idea and you are all really jazzed up about it, I'm going to rain on your parade a bit.
- The problem is that batching is pretty hard in NLP
- We saw before that batching involves stacking multiple vectors of the same shape into a matrix
- In NLP we often have examples that have different lengths
  - Because sentences have different lengths
  - Because words have different number of characters in them
- We can't just batch any given pair of sentences.

## Batching in NLP

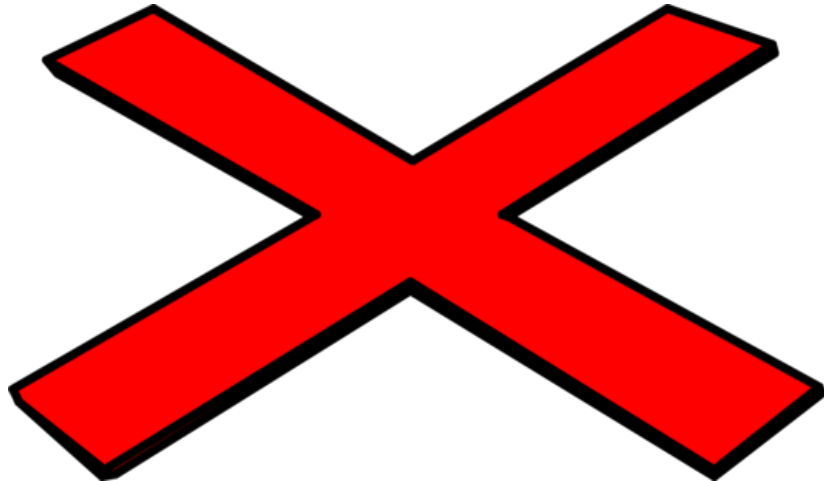
[The dog ran very fast]

[The cat slept]

- Now that we know batching is a good idea and you are all really jazzed up about it, I'm going to rain on your parade a bit.
- The problem is that batching is pretty hard in NLP
- We saw before that batching involves stacking multiple vectors of the same shape into a matrix
- In NLP we often have examples that have different lengths
  - Because sentences have different lengths
  - Because words have different number of characters in them
- We can't just batch any given pair of sentences.

## Batching in NLP

The	dog	ran	very	fast
The	cat	slept		



- Here we can't batch these examples because they are different lengths, this would create a ragged array, we need it to be regular.
- [Image from here](#)

## Introducing <PAD>

[	The	dog	ran	very	fast	]
	The	cat	slept	<PAD>	<PAD>	

- Now let's look at how we can handle batching things of different lengths.
- We are going to use a new symbol called <PAD>.
- We add this to the end of shorter sentence so that all the sentences in a batch have the same length.
- Now that they are all the same length we can turn them into a matrix.

## Embeddings and <PAD>

- In NLP we have this idea of an “embedding”.
- This is a mapping from a word id to a vector that represents the word.
- We call each of the dimensions in the vector a feature.
- The principled approach is force the vector for <PAD> to be all zeros.
- Pytorch `nn.Embedding` has a `padding_idx` parameter that will force the vector with that index to be zero.
- In Tensorflow you need to role your own. **This is an example of doing that in Mead-Baseline.**

- Why is this the principled way?
- In general we don't want the network to know anything about the padding.
- The padding should just be an artifact of batching that doesn't interact with the network or the other values at all.
- Part of hiding the <PAD> from the network is not letting it learn a representation of the vector.
- We do this by forcing it to zero.

## The lengths vector

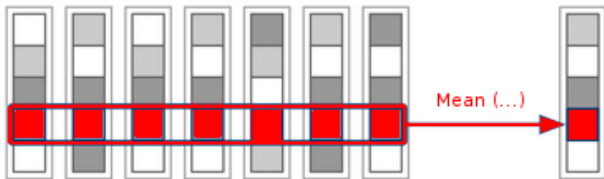
The	dog	ran	very	fast
The	cat	slept	<PAD>	<PAD>

$$L = \begin{bmatrix} 5 & 3 \end{bmatrix}$$

- Now that some of our examples are longer because of this <PAD> we need to track how long the actual data is
- We store this in a lengths vector
- This will be a vector of the same length as the number of items in the batch
- The value will be the number of non padded elements in that example
- You can actually have padding on every example in the batch.
- This means the length of the physical batch is longer than the maximum value in the lengths vector
- This is normally an artifact of being lazy with your data preprocessing, a bit sloppy, and can result in extra computation being done.
- Because doing this is wasteful we won't really consider this situation.

## Mean Pooling

### Mean pool over filters



- As a warm up we are going to look at Mean Pooling.
- The main use-case of mean pooling is in the Neural Bag of Words model.
- We embed each word resulting in a matrix that is sentence length by embedding size
- We do a mean over time for each feature, this means feature is mean-ed separately by looking at it's value for each individual word.
- This results in a sentence vector that we can then use for classification



## Mean Pooling

$$\begin{bmatrix} 1 & 10 & 8 & 17 & 13 & 17 \end{bmatrix} = \frac{66}{6} = 11.0$$

- Mean pooling gives us a clear example of how padding can get us in trouble
- Side Note: In most of these examples there is actually another dimension of features that these operations are actually acting on but we are going to ignore it and only really look at a single feature in this dimension for simplicity
- Here we see the mean pooling for a vector of length 6
- Here we see the mean pooling for a length of 4 (shorter)

## Mean Pooling

$$\begin{bmatrix} 1 & 10 & 8 & 17 & 13 & 17 \end{bmatrix} = \frac{66}{6} = 11.0$$

$$\begin{bmatrix} 22 & 24 & 9 & 13 \end{bmatrix} = \frac{68}{4} = 17.0$$

- Mean pooling gives us a clear example of how padding can get us in trouble
- Side Note: In most of these examples there is actually another dimension of features that these operations are actually acting on but we are going to ignore it and only really look at a single feature in this dimension for simplicity
- Here we see the mean pooling for a vector of length 6
- Here we see the mean pooling for a length of 4 (shorter)
- How we see that in order to batch we need to pad the shorter vector
- Now our denominator is much larger so our result of mean-ing is smaller!
- Before I said that batching allows us to combine unrelated examples and calculate them in one shot and get the same results as if you had calculated them independently
- When your padding goes wrong the results for an example will be dependent on the other things in the batch. Padding can cause bleeding of information examples and our independence is broken thus our calculations will be wrong.

## Mean Pooling

$$\begin{bmatrix} 1 & 10 & 8 & 17 & 13 & 17 \end{bmatrix} = \frac{66}{6} = 11.0$$

$$\begin{bmatrix} 22 & 24 & 9 & 13 \end{bmatrix} = \frac{68}{4} = 17.0$$

$$\begin{bmatrix} 5 & 4 & 8 & 9 & 10 & 34 \\ 6 & 3 & 1 & 4 & 0 & 0 \end{bmatrix} = \begin{bmatrix} \frac{66}{6} \\ \frac{68}{6} \end{bmatrix} = \begin{bmatrix} 11.0 \\ 11.\bar{3} \end{bmatrix}$$

- Mean pooling gives us a clear example of how padding can get us in trouble
- Side Note: In most of these examples there is actually another dimension of features that these operations are actually acting on but we are going to ignore it and only really look at a single feature in this dimension for simplicity
- Here we see the mean pooling for a vector of length 6
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- When your padding goes wrong the results for an example will be dependent on the other things in the batch. Padding can cause bleeding of information examples and our independence is broken thus our calculations will be wrong.

## Mean Pooling

```
>>> x
array([[ 1, 10,  8, 17, 13, 17],
       [22, 24,  9, 13,  0,  0]])
>>> np.mean(x, axis=1)
array([11.          , 11.33333333])
>>> lengths
array([6, 4])
>>> np.sum(x, axis=1) / lengths
array([11., 17.])
```

- We can solve this by explicitly using the lengths vector.
- We can manually do our sum and then divide by the actual length instead of the length of the padded vector
- This assumes that the padded value is zero. Soon we will see how to insure that.
- This sort of thing is important because if you don't do it then values will dependent on what else is in the batch.
- This problem is especially important in training vs production environments
  - when training it an example is almost always in a batch
  - When you are in a batch is is relatively rare to be the longest so being correct in the presence of padding is important
  - In prod you are often a single example and you are by definition the longest thing in the batch
  - This means that you need to have the same results with and without padding

## Cross Entropy

$$L_{\text{cross-entropy}}(\hat{y}, y) = - \sum_i y_i \log(\hat{y}_i)$$

- Our next example of where padding can get up to no good is in our loss functions
- Our normal loss is the cross entropy loss between two vectors
- $y$  is the vector that has our real labels and  $\hat{y}$  is our predicted labels
- cross entropy is a measure of the difference between two distributions
- The cross entropy is the sum of the true value times the log of the predicted values
- In classification where we output a probability distribution over classes we can use this loss function to train the model

## Sparse Cross Entropy

$$L_{\text{cross-entropy}}(\hat{y}, y) = - \sum_i y_i \log(\hat{y}_i)$$

$$L_{\text{sparse-cross-entropy}}(\hat{y}, y) = - \log(\hat{y}_i[y])$$

- Because our labels are always a one hot vector (it only has a 1 at the location that represents the label index) most of these elements in this sum will be zero
- In fact the only value will be the probability for the predicted class
- We can short cut this calculation by just indexing the predicted vector with the true class index
- This is what we call sparse cross entropy

## Token-Level (Sparse) Cross Entropy Loss

$T :=$  length of example

$V :=$  number of possible labels

$$s \in \mathbb{R}^{T \times V}$$

$$l \in \mathbb{N}^T$$

$$J = - \sum_i^T s_i[l_i]$$

- This idea of sparse cross entropy can be extended to the idea of a token level loss.
- This means that a decision is made for every token and we calculate the loss for the probability vector that is output for each token in a sentence.
- We will see soon how this loss at the padding positions can mess us up.
- Here we see the cross entropy loss where we have a matrix of scores. Each row represents the scores of a token and each column is the score for some class.
- Our loss is then the sum (or mean) of the score for each token.
- This kind of loss is often used for language modeling where the label is the index of the next word or else sequence tagging where the label is the tokens tag.

## (Sparse) Cross Entropy Loss

$$s = \begin{bmatrix} -0.34 & 0.1 & -0.11 & \cdots & 0.001 \\ 0.93 & -8.88 & -0.39 & \cdots & 0.12 \\ & & \vdots & & \\ -0.45 & 0.23 & 1.1 & \cdots & -0.3 \end{bmatrix}$$

$$l = \begin{bmatrix} 2 & 1 & 10 & 5 & 1 \end{bmatrix}$$

$$J = \begin{bmatrix} -0.11 & -8.88 & 0.28 & 0.43 & 0.23 \end{bmatrix}$$

- Here we see some examples of score, labels, and the selected scores.



## (Sparse) Cross Entropy Loss

$$s = \begin{bmatrix} -0.34 & 0.1 & -0.11 & \cdots & 0.001 \\ 0.93 & -8.88 & -0.39 & \cdots & 0.12 \\ & & \vdots & & \\ -0.45 & 0.23 & 1.1 & \cdots & -0.3 \end{bmatrix}$$
$$l = \begin{bmatrix} 2 & 1 & 10 & 5 & 1 \\ 5 & 6 & 3 & & \end{bmatrix}$$
$$J = \begin{bmatrix} -0.11 & -8.88 & 0.28 & 0.43 & 0.23 \\ -3.4 & 0.6 & 0.37 & -956.0 & 10000.0 \end{bmatrix}$$

- Here we see what can happen when you are batching and have a shorter sentence.
- When you are selecting the scores for the padding tokens they really could be anything
- We don't really want the network to have to learn to output a padding symbol after the normal length, that isn't a situation you would really have a use for.
- We also don't want random lucky values to make our loss look really good.
- We need a way to zero out the effects of these padded locations

## Making a Mask

```
>>> lengths = np.array([5, 3, 4])
>>> steps = np.arange(np.max(lengths))
>>> steps
array([0, 1, 2, 3, 4])
>>> mask = (np.reshape(lengths, (-1, 1)) > steps).astype(np.int)
>>> mask
array([[1, 1, 1, 1, 1],
       [1, 1, 1, 0, 0],
       [1, 1, 1, 1, 0]], dtype=int32)
```

- We will zero out these values by using a mask
- This mask has ones where the valid tokens are and a zero where the padding tokens are.
- Here we see how to create a mask.
- We create this with broadcasting.
- We create a vector of incrementing steps along the dimension that includes padding and compare them to the lengths vector.
- When the length is longer than the current steps we are still valid so the value is one.
- Once the step is longer than the length we are in the pad and will output zeros.

## (Sparse) Cross Entropy Loss

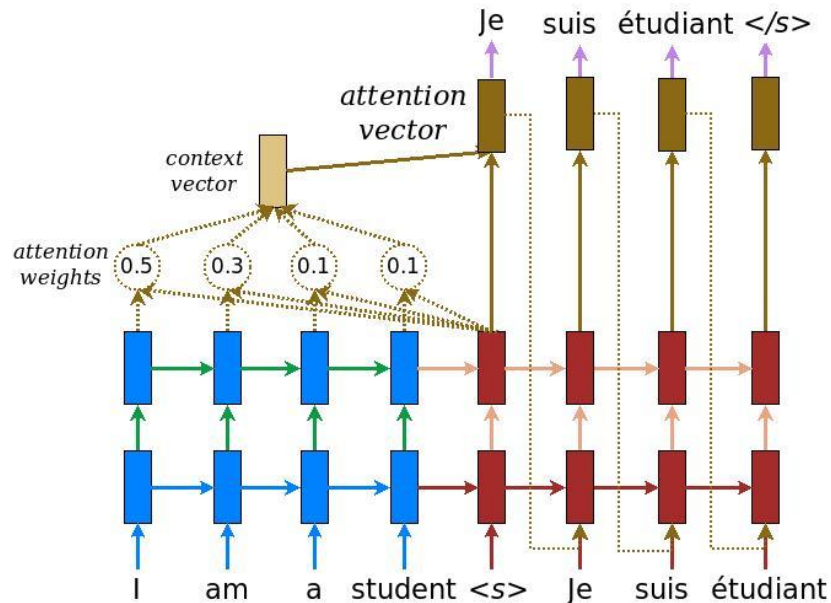
$$J = \begin{bmatrix} -0.11 & -8.88 & 0.28 & 0.43 & 0.23 \\ -3.4 & 0.6 & 0.37 & -956.0 & 10000.0 \end{bmatrix}$$

$$\text{mask} = \begin{bmatrix} 1 & 1 & 1 & 1 & 1 \\ 1 & 1 & 1 & 0 & 0 \end{bmatrix}$$

$$J * \text{mask} = \begin{bmatrix} -0.11 & -8.88 & 0.28 & 0.43 & 0.23 \\ -3.4 & 0.6 & 0.37 & 0 & 0 \end{bmatrix}$$

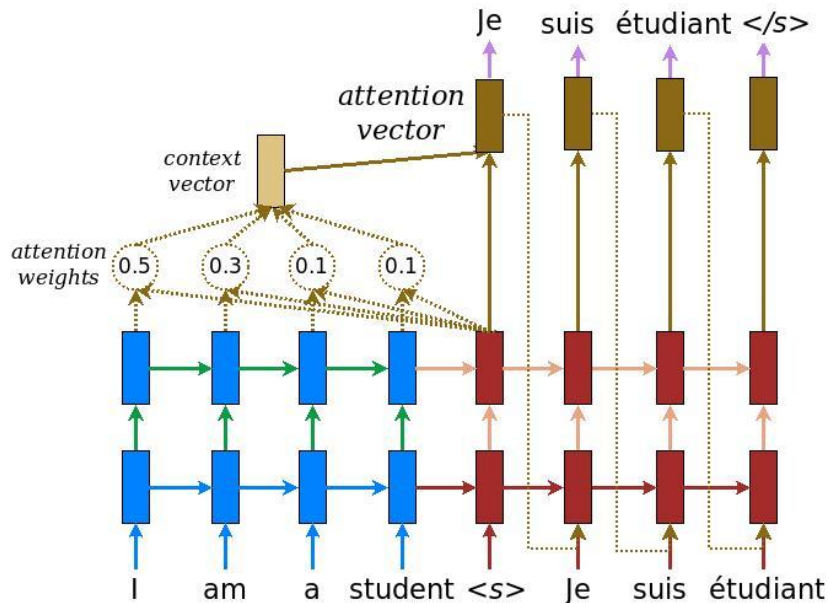
- Here we can use our mask to force the padding locations to zero.
- We can then sum the values (or mean them with the lengths if you want).
- Some toolkits like Pytorch can handle this for you, you give a value to the `padding_idx` parameter and it will ignore values where the label is that padding index
- Toolkits like tensorflow will probably make you do this yourself.

## Attention



- Attention is used to contextualize a vector relative to others.
- We have some vector here in the decoder and want to create a new vector that summarizes the information in these vectors created by the encoder that is relevant to decoder vector.
- We do this by creating a similarity score between the vector we care about and each vector in the memory bank of vectors
- We then normalize these values and create new vector by taking a weight average of the memory bank
- This is often used in seq2seq models to short cut information from the encoder into the decoder
- [Image From Here](#)

## Attention



- In a more general format we have 3 sets of vectors
  - Query Vectors: the ones we are processing right now
  - Key Vectors: The ones we are doing the similarity score with
  - Value Vectors: The ones that are actually mixed together
- In some cases these vectors can be the same (for example RNN Seq2seq normally use  $K = Q$ ) in other case they are either different or the same thing projected into different spaces
- The point of this is that the representation of a vector that is useful down stream might not be one that is similar to the query vectors so the keys can be the similar ones.
- [Image From Here](#)

## Attention

$$q = \begin{bmatrix} \text{---} & q & \text{---} \end{bmatrix}$$
$$K = \begin{bmatrix} | & | & & | \\ k_1 & k_2 & \cdots & k_n \\ | & | & & | \end{bmatrix}$$
$$s = \text{sim}(q, K) = \begin{bmatrix} s_1 & s_2 & \cdots & s_n \end{bmatrix}$$

- First we calculate a score between the vectors
- There are multiple ways to calculate these similarities, the original paper used a small neural network, later architectures (notably the transformer) use the dot product to calculate the score.
- The transformer actually uses the scaled dot product where it divides the score by  $\sqrt{d_k}$  because the dot product will get larger as the vectors get larger
- In this case we are just using a single vector for  $q$  but we can actually use a matrix  $Q$  to calculate the attention between multiple vectors at once.
- This vector version is common for RNN seq2seq models (where you have single vector at each timestep)
- The Matrix version is used in the transformer where we want to do attention for the whole sentence at once

## Attention

$$s = \text{sim}(q, K) = [0.21 \quad -0.34 \quad \dots \quad 0.55]$$

$$\text{softmax}(x) = \frac{e^{x_i}}{\sum_j e^{x_j}}$$

$$\alpha = \text{softmax}(s) = [0.1684 \quad 0.0977 \quad \dots \quad 0.2379]$$

- Then we use the softmax so the scores are all positive and they sum to one.
- The softmax is done by exponentiating each element and then normalizing the elements.
- I have a blog post about how you can do that in a numerically stable way if you decide to roll your own implementation.
- We can then use these scores to do a weight average.

Numerically stable softmax blog post.

## Attention

$$V = \begin{bmatrix} | & | & & | \\ v_1 & v_2 & \cdots & v_n \\ | & | & & | \end{bmatrix}$$

$$s = \text{sim}(q, K) = [s_1 \quad s_2 \quad \cdots \quad s_n]$$

$$\alpha = \text{softmax}(s)$$

$$c = V * \alpha^T$$

- Then we do a weighted average of the value vectors based on these calculated (and normalized) scores.
- Now  $c$  is a combination of the vectors in  $V$  weighted by their attention scores which were calculated based on the similarity of the query vectors to  $q$



## Attention

$$K = \begin{bmatrix} | & | & & | & | \\ k_1 & k_2 & \cdots & \langle \text{PAD} \rangle & \langle \text{PAD} \rangle \\ | & | & & | & | \end{bmatrix}$$
$$s = [s_1 \quad s_2 \quad \cdots \quad 0 \quad 0]$$

- But what happens when we calculate the similarity between vectors that have padding on them?
- With most similarity metrics (dot/scaled dot product) we do get zeros for the similarity between a padding vector and a valid vector (assuming the pad is zeros)
- Is this enough so that our calculations will be correct?
- Would I have brought this up if it was?

## Attention

```
>>> x = np.array([
    [0.32, 0.11, 0.83, 0.9, 0.4],
    [0.5, 0.2, -1.0, 0, 0]
])
>>> x
array([[ 0.32,  0.11,  0.83,  0.9 ,  0.4 ],
       [ 0.5 ,  0.2 , -1.   ,  0.   ,  0.   ]])
>>> softmax(x)
array([[0.15759942, 0.12774761, 0.26244893, 0.28147862, 0.17072542],
       [0.31476139, 0.23318098, 0.07023276, 0.19091244, 0.19091244]])
```

- Here we can see that when the scores going into the softmax are zero we will get non-zero scores out!
- This means that when we do the mixing of the value vectors were are going to include the padding vectors in our weighted average.
- What we need to have there be zeros **after** the softmax which will zero out the contribution of the padding.

## Attention

```
>>> mask = make_mask(np.array([5, 3]))
>>> mask
array([[1, 1, 1, 1, 1],
       [1, 1, 1, 0, 0]], dtype=uint8)
>>> attn = softmax(x) * mask
>>> attn
array([[0.15759942, 0.12774761, 0.26244893, 0.28147862, 0.17072542],
       [0.31476139, 0.23318098, 0.07023276, 0.          , 0.          ]])
>>> np.sum(attn, axis=1)
array([1.          , 0.61817513])
>>> softmax(x[1:np.newaxis, 0:3])
>>> array([[0.50917835, 0.3772086 , 0.11361305]])
```

- Ok, our first cut is just masking out those padding weights, now their values are zero so they won't contribute to the weighted average anymore.
- Is this ok?
- We can see that the values for this example don't sum to one anymore, this solution isn't enough
- We we can also see that the values calculated for the shorter example are very different than when that example is calculated by itself.
- Remember the whole point of batching is that we should be able to run multiple examples at once and get the same values as if we ran them through one at a time

## Attention

```
>>> inv_mask = (1 - mask) * -1e9
>>> masked_x = (x * mask) + inv_mask
>>> masked_x
array([[ 3.2e-01,  1.1e-01,  8.3e-01,  9.0e-01,  4.0e-01],
       [ 5.0e-01,  2.0e-01, -1.0e+00, -1.0e+09, -1.0e+09]])
>>> softmax(masked_x)
array([[0.15759942, 0.12774761, 0.26244893, 0.28147862, 0.17072542],
       [0.50917835, 0.3772086 , 0.11361305, 0.          , 0.          ]])
>>> softmax(x[1:np.newaxis, 0:3])
>>> array([[0.50917835, 0.3772086 , 0.11361305]])
>>> softmax(masked_x).sum(axis=1)
array([1., 1.]
```

- We need to hack softmax to force it to output zeros for the padded values while the other values sum to one.
- We know that the softmax will exponentiate each element and we know that exponentiation by a negative number of will approach zero as the exponent grows.
- With a sufficiently negative number we can force an element to zero
- On this first line  $(1 - \text{mask})$  will flip a mask of ones and zeros.
  - $1 - 1 = 0$  and  $1 - 0 = 1$
  - This will mean that the ones are now where the pads and zero where the valid elements are
- Then we multiply in a very negative number so out mask is all zeros and this neg
- We multiply by the mask to make sure the padding locations are zero (not strictly necessary) then we add the inverse mask in.
- the inverse mask has all zeros for the valid elements so they don't change but it has  $-1e^9$  for the pad so zero + that negative is that negative
- Now when we do the softmax we see that out values match and they sum to one in each axis

## Attention



Encoder-Decoder Attention



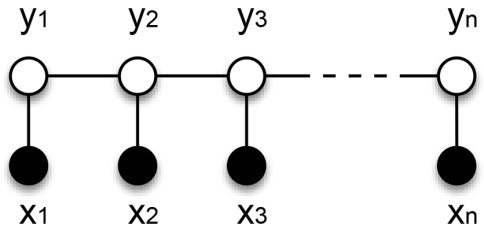
Encoder Self-Attention



Masked Decoder Self-Attention

- We have been looking at encoder-decoder attention where we compare a vector in the decoder to the values in the encoder.
- There is also self attention where we compare to all the tokens around us.
- There is also a special case of self attention where we can use a mask to enforce causality
- This mask would stop the attention from looking forward in time at words that haven't been generated yet.
- This mask works the same way as the length mask.
- Some times you need to combine these masks.
- Attention is an example where just masking out the padding value isn't enough, we actually have to manipulate the padding values in a specific way.
- [Image is from here](#)

## Conditional Random Field

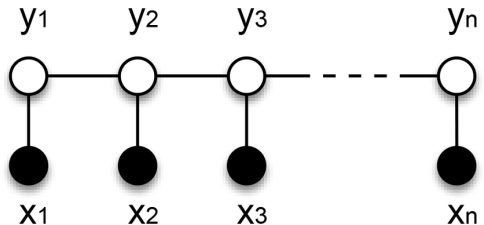


$$P(y|x; \theta) = \frac{\exp(\sum_i \sum_j w_j f_j(y_{i-1}, y_i, x, i))}{\sum_{y' \in Y} \exp(\sum_i \sum_j w_j f_j(y'_{i-1}, y'_i, x, i))}$$

Great Tutorial on CRFs

- This is our most complex example so it is ok if you don't understand everything
- In our token level loss from before we made a decision about the label of a token without considering the label we chose for the last token.
- Here is this factor graph diagram of a CRF we see that the label decisions are conditioned on the previous label. This makes the CRF a global model.
- In the token level loss our scores and therefore the label decision were contextual in that they were based on the whole input (from an RNN or something) but they don't consider the past labels so it doesn't have any idea of global coherence.
- These scores are called the local representation of the token.
- A trend in research is to buff these local representations while not caring about the global part too much
- Problems that deal with making a sequence of decisions like are called structured prediction problems
- [Image from Here](#)

## Conditional Random Field



$$P(y|x; \theta) = \frac{\exp(\sum_i \sum_j w_j f_j(y_{i-1}, y_i, x, i))}{\sum_{y' \in Y} \exp(\sum_i \sum_j w_j f_j(y'_{i-1}, y'_i, x, i))}$$

Great Tutorial on CRFs

- Our loss function here can help illustrate the differences
- Thinking back to our token level loss function our loss was the sum of the losses at each step
- This is explicitly optimizing this local decision making process.
- In this loss function we see that we are actually acting on the sequence level
- Our loss is now the score you model gives the gold sequence  $y$
- Normalized but the score of the possible sequences.
- We want to optimize it so more probability is placed on the gold sentence
- [Image from Here](#)

## Conditional Random Field - Best Path

$t :=$  number of tokens

$c :=$  number of classes

$$e \in \mathbb{R}^{t \times c}$$

$$t \in \mathbb{R}^{c \times c}$$

$$y \in \mathbb{N}^t$$

$$\text{gold score} = \sum_i e_i[y_i] + t[y_{i-1}, y_i]$$

- $e$  is our emission scores, this tells us the class scores for each token. This is the local representation of the token
  - One thing to note is that these scores are unnormalized (not softmaxed). This lets us avoid some problems locally normalized models have but that is out of scope
- $t$  is our transition scores, this tells us the score of moving from one class to another at the next step.
- $e$  is a score tensor that is output by the network whereas  $t$  is a parameter that is learned
- Our score is sum over tokens of emission score for that label plus the score of transitioning to that label from the previous label.
- It is easy to see how if you have padding you are going to get junk in your score, The emission score and the transistors involving padding will get in there
- This is an easy fix we have seen before, before you sum across the tokens use your mask to set the padding values to zero and you'll get the right score.



## Conditional Random Field - All Paths

$t$  := number of tokens

$c$  := number of classes

$\alpha \in \mathbb{R}^c$

$e \in \mathbb{R}^{t \times c}$

$t \in \mathbb{R}^{c \times c}$

$$\forall_{i \in t} \alpha_i = \text{logsumexp}(\alpha_{i-1} + e_i^T + t)$$

- When calculating all paths you can't actually enumerate all the possible paths. There are  $t^c$  possible paths.
- We can use a dynamic programming to do this, it is out of scope though.
- We have  $\alpha$  which represents the total score for any path that ends up in this label
- We often initialize this  $\alpha$  with all the score on a fake  $\langle \text{START} \rangle$  that helps us represent which labels often start a sequence.
- We calculate a new version of these scores at each time step and overwrite the previous  $\alpha$ s
- We need to make sure that we only update the  $\alpha$  if we are transitioning to a valid token
- If we are looking at a  $\langle \text{PAD} \rangle$  we don't want to update the alpha.

## Conditional Random Field - All Paths

---

### Algorithm 1: The Forward Algorithm

---

$T$  := number of tokens in batch;

$c$  := number of classes;

$t$  := number of valid tokens in example;

$\alpha \in \mathbb{R}^c$ ;

$e \in \mathbb{R}^{t \times c}$ ;

$t \in \mathbb{R}^{c \times c}$ ;

**for**  $i \leftarrow 0$  **to**  $T$  **do**

**if**  $i < t$  **then**

$\alpha = \text{logsumexp}(\alpha + e_i^T + t)$ ;

**else**

$\alpha = \alpha$

**end**

**end**

---

- I have removed the  $\langle \text{START} \rangle$  and  $\langle \text{END} \rangle$  from this algorithm but you should use them, they are a higher-fidelity representation and let you model which starts the likely to start of end a sequence
- This boils down to a simple conditional.
- If we are still inside the valid tokens we assign alpha based on the results of the dynamic programming algorithm
- If we are in the padding step we just reuse the old  $\alpha$ s
- How do we implement this conditional with masking?

## Conditional Random Field

```
>>> lengths = np.array([3, 1, 5])
>>> lengths
array([3, 1, 5])
>>> for i in range(np.max(lengths)):
...     new_alphas = calculate_step(alpha, emission, transition, i)
...     mask = (i < lengths).astype(np.uint8)
...     print(mask)
...     alphas = (alpha * mask) + (new_alphas * (1 - mask))
[1 1 1]
[1 0 1]
[1 0 1]
[0 0 1]
[0 0 1]
```

- In this example we can see how we can use create masks that represent conditionals.
- First we track how far we are into the loop with the  $i$  variable.
- First we calculate the updated  $\alpha$ s regardless of whether we are padding or not
- Then we make a mask where we have ones if we are inside the valid lengths
- Then we update the  $\alpha$ s based on the mask so that we only update the ones we should.

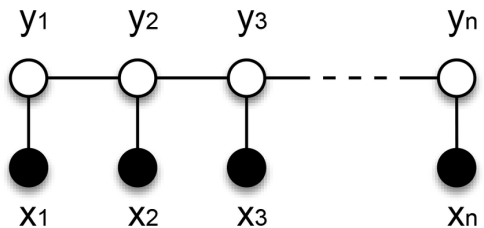
## Conditional Random Field

```
>>> a
array([1, 2, 3, 4, 5])
>>> b
array([-1, -2, -3, -4, -5])
>>> mask = np.where(a >= 4, 1, 0)
>>> mask
array([0, 0, 0, 1, 1])
>>> (a * mask) + (b * (1 - mask))
array([-1, -2, -3, 4, 5])
```

- This is a more illustrative example.
- We see positive and negative values.
- We make a mask that represents “if  $a \geq 4$  then return  $a$  else return  $b$ ”.
- We multiply  $a$  by the mask so now all the values are zero when the conditional is false.
- We multiply  $b$  by the inverse mask, this makes the elements where  $a$  is non zero 0.
- No the non-zero values are disjoint between  $a$  and  $b$ .
- We add the vectors to combine the results.

## Conditional Random Field

A complex layer with complex padding



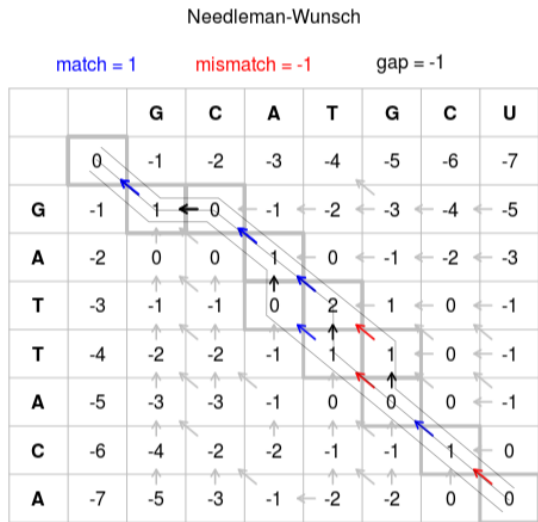
- The CRF is a rather complex layer but the main reason I brought it up here is that it introduces a new way that we need to handle padding.
- We are using the mask to simulate a conditional.
- This is common in parallel processing where we compute both branches of a conditional and use a mask to combine the result according to that conditional.

## Viterbi

- When we predict actual tags from a CRF we use the Viterbi algorithm to find the maximum scoring path.
- The Viterbi algorithm is the exact same as the forward algorithm except that it uses a  $\max$  instead of a  $\log \text{sum} \exp$  when computing the new  $\alpha$ s.
- $\alpha$  now represents the maximum score of any path that ends at each class.
- We can use the exact same padding trick when doing Viterbi decoding!

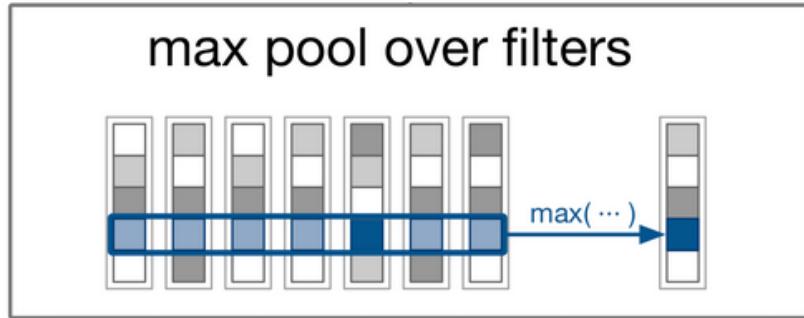
- A way to look at these two algorithms is that they are the same algorithm using different semirings. A semiring is basically a re-definition of math in some context that uses different operations for sums and additions
  - The CRF forward is on the log semiring where the addition is a  $\log \text{sum} \exp$  addition
  - The Viterbi algorithm is on the Viterbi semiring where the addition is a maximum
- in the CRF  $\alpha$  was the sum of scores for all paths that end in some class
- in Viterbi  $\alpha$  is the maximum score of any path that ends in some class

## Viterbi - Getting a Path



- Like most dynamic programming algorithms you need to store your decisions so you can back track through to get your actual answer
- In this picture we can see these arrows that tell us where we came from and can be used to recover the edit path.
- When doing Viterbi decoding we calculate these backtracks with the  $\text{argmax}$
- As we calculate the back tracking we need to not overwrite the final decision once we hit the end of the valid lengths
- We use the same kind of padding we used above for this.
- When we are following the backtrack we need to do something similar but flip the lengths we care about because we are reading from the back of the tensor.
- [Image from here](#)
- Now that we have talked about some complex cases lets look at some subtle ones.

## Max Pooling



- Max pooling is often applied across a feature dimension in the same way that mean pooling is
- Here we see a set of vectors, each one is a featured representation of a word in a sentence
- We do max pooling across a single feature dimension.
- We end up with a single vector where each feature is the largest value that feature took in the course of the sentence
- We can also do this at a word level where each vector represents a character in a word
- A common use case is to encode words with a Convolutional Network (which we will get to soon) and then aggregate the results with max pooling.
- [Image from here](#)



## Max Pooling

$$x = [1 \quad 0 \quad 4 \quad 3 \quad 2 \quad 0 \quad 0]$$
$$\text{lengths} = [5]$$
$$\max(x) = 4$$

- Here we are zoomed in a single feature over time, these are the numbers in the blue highlighted area on the last slide
- Lets say these vectors we are zoomed in one just went through a ReLU  $\max(0, x)$  so it will never have a negative number in it.
- This example has a length of 5 which means the last two zero features are padding
- Pop Quiz: Is it safe to do a max without thinking about padding?
- **YES**

## Max Pooling

$$x = [-3 \quad -1 \quad -4 \quad -3 \quad -2 \quad 0 \quad 0]$$

$$\text{lengths} = [5]$$

$$\max(x) = 0$$

$$\max(x[:5]) = -1$$

- What about a vector that looks like this?
- Lets say here we are looking at a feature of a word right after embedding.
- Your max pooling it going to output zero for this feature!
- Your network will never see a negative value for most examples while training! (most examples are padded)
- Your network will see these negative values at test time and have no idea what to do!
- How do we fix this?

## Max Pooling

$$x' = [-3 \quad -1 \quad -4 \quad -3 \quad -2 \quad -10000 \quad -10000]$$

$$\text{lengths} = [5]$$

$$\max(x') = -1$$

- We use the same trick we used in attention, by setting the pads so a very negative value we can ensure they will never be picked

## Min Pooling

$$x = \begin{bmatrix} 3 & 1 & 4 & 3 & 2 & 0 & 0 \end{bmatrix}$$

$$x' = \begin{bmatrix} 3 & 1 & 4 & 3 & 2 & 10000 & 10000 \end{bmatrix}$$

$$\text{lengths} = \begin{bmatrix} 5 \end{bmatrix}$$

$$\min(x) = 0$$

$$\min(x[:5]) = 1$$

$$\min(x') = 1$$

- This same thing applies in Min pooling, we where the padding can be selected as the minimum.
- We can solve this similarly but set the padding to a very high value.

## Max Pooling Results

Model	mean	std	min	max
Masking	83.60	0.22	83.31	83.86
Naïve	82.90	0.43	82.15	83.20

- Here we can see a small selection of results comparing these differences.
- There were 5 runs of a simple Max Over Time Neural Bag of Words on the SST2 dataset
- As we can see the mean result of the version where we masked correctly is higher than not doing it.
- We also see that the variability is lower. I would imagine that this is because we no longer have randomness in the results depending on what things are batched together.

## Convolution 1D

I  
like  
this  
movie  
very  
much  
!

0.6	0.5	0.2	-0.1	0.4
0.8	0.9	0.1	0.5	0.1
0.4	0.6	0.1	-0.1	0.7
...	...	...	...	...
...	...	...	...	...
...	...	...	...	...
...	...	...	...	...

0.2	0.1	0.2	0.1	0.1
0.1	0.1	0.4	0.1	0.1

0.51

I  
like  
this  
movie  
very  
much  
!

0.6	0.5	0.2	-0.1	0.4
0.8	0.9	0.1	0.5	0.1
0.4	0.6	0.1	-0.1	0.7
...	...	...	...	...
...	...	...	...	...
...	...	...	...	...
...	...	...	...	...

0.2	0.1	0.2	0.1	0.1
0.1	0.1	0.4	0.1	0.1

0.51
0.53

- Convolutions are basically used to process short n-grams.
- This is commonly used in text classification
- The conv will look at multiple words at a time, and create a new representation of the word contextualized by the words around it.
- Once we have contextualized the representation we will aggregate into a single vector with some sort of pooling.
- It is also common in taggers to use a convolution to create a representation of a word based on the characters.
- [Image from here](#)

## Convolution 1D

$$w = [1 \quad 2 \quad 3]$$
$$\begin{bmatrix} 4 & 5 & 6 & 7 & 8 \end{bmatrix}$$
$$\begin{bmatrix} 32 & 38 & 44 \end{bmatrix}$$

- So this shows us how a convolution works, we look at a window, scale the values by a learned weight matrix and sum them to create a new value for this position.
- Unfortunately here we can see that a convolution will actually create a new, shorter sentence
- We need to introduce a new form of padding that I will call conv-padding to distinguish from the padding we have been talking about before

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## Padded Convolution 1D

$$w = \begin{bmatrix} 1 & 2 & 3 \end{bmatrix}$$
$$\begin{bmatrix} 0 & 4 & 5 & 6 & 7 & 8 & 0 \end{bmatrix}$$
$$\begin{bmatrix} 23 & 32 & 38 & 44 & 23 \end{bmatrix}$$

- Here we can see how conv-padding can be used to stop the sequence from getting shorter
- Here we can see we are adding a padding of zeros on each side
- There are filtersz // 2 units on each side
- Another advantage of this conv-padding is that without padding there is only a single calculation that involves the first word. conv-padding makes it so that there is more information from the tokens on each end.
- We see here how we now get a vector that is the same length as the input (before padding)

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## Batched Convolution 1D

$$w = \begin{bmatrix} 1 & 2 & 3 \end{bmatrix}$$
$$\begin{bmatrix} 6 & 14 & 7 & 0 & 0 & \end{bmatrix}$$
$$\begin{bmatrix} 0 & 6 & 14 & 7 & 0 & 0 & 0 \end{bmatrix}$$
$$\begin{bmatrix} 54 & 55 & 28 & 7 & 0 \end{bmatrix}$$

- In this example we are zoomed in on an example that has already been padded, the two zeros at the end
- Now we are going to pad it again because the conv-padding is applied to the whole batch at once
- As we slide the conv we can see that at the end we have spots where the only part of the window is on the valid tokens but it is writing into the padding section of the output vector.
- This is introducing a new value what could be selected by the max pool during training time that would never happen at test time.

## Batched Convolution 1D

$$w = [1 \quad 2 \quad 3]$$

When we were already padded our convolution will bleed into the padding

$$\begin{bmatrix} 0 & 6 & 14 & 7 & 0 & 0 \\ 54 & 55 & 28 & 7 & 0 \end{bmatrix}$$

When we do this convolution by itself

$$\begin{bmatrix} 0 & 6 & 14 & 7 & 0 \\ 54 & 55 & 28 \end{bmatrix}$$

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## Batched Convolution 1D

These Convolutions are commonly used for:

- Encoding word n-grams which are then max pooled to create sentence representations for text classification
- Encoding character n-grams and max pooling to create word representations. These “char compositional” features are common for sequence tagging
- Creating a windowed context for tagging

- Does this matter? This seems rare enough that you shouldn't care
- Yes it does
  - I first noticed this error in a tagger where the output tags were different between running in batched mode and running in single mode.
  - The very first tagger I trained to try to recreate this had the exact same error so it is rather prevalent
  - These taggers only use these convs in the character compositional word representations. The fact that this errors can propagate all the way to a label flip is worrying for the robustness of Neural Networks

## Batched Convolution 1D

These Convolutions are commonly used for:

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- Creating a windowed context for tagging

- Elmo has this batch instability bug too, but given their cavalier attitude to the RNN hidden states I think the engineering in Elmo is pretty from the hip so they probably don't care. This is a rant for another time so I will move on.
- An Interesting observation is that for the last use case I don't think we really need to care, padding in a later stage (CRF or the token level loss) should save us.
- "But how do I fix this one Brian?"
- This is one is pretty easy you just need to zero out the padded values after they go through the conv. If you are following it with maxpool you might want to just jump right to masking the pad values with a very negative number

## Unit Tests

```
def test_batch_stability():  
    ex1 = generate_example(length=random.randint(5, 10))  
    ex2 = generate_example(length=random.randint(10, 15))  
    res1 = network(ex1, lengths=[len(res1)])  
    res2 = network(ex2, lengths=[len(res2)])  
    batch, lengths = batch_examples(ex1, ex2)  
    res = network(batch, lengths=lengths)  
    batch1, batch2 = extract_results(res, lengths)  
    np.testing.assert_allclose(batch1, ex1)  
    np.testing.assert_allclose(batch2, ex2)
```

An example of these tests in our CRF

- The easiest way to avoid these errors is a unittest, There is a nice recipe for these batch instability tests
  1. Generate two examples with different lengths
  2. Run each example through the thing you want to test and record results
  3. Batch the two examples together
  4. Run the batch through the thing you want to test
  5. Extract the values that represent each example and compare them to the values from running solo

## Conclusion

- Almost nothing is a safe when you are padding.
- You should be passing lengths around for most things.
- Some frameworks help manage padding and you should use them when you can.
- These bugs are hard to find so you should be writing unittests designed to check for them.

- As mentioned pytorch losses can handle this for you with the `padding_idx` parameter.
- Most RNN implementations take a length parameter and handle this for you.
- DyNet has an autobatching feature where they basically have a computation graph compiler.
  - You can write forward passes that are not batched and the compiler will batch them on the fly.
  - Unfortunately DyNet isn't very active these days.

## Contact

- [Slides](#)
- [Twitter: @blester125](#)
- [Github](#)
- [Mead-Baseline](#)

- Here is a link to these slides which are in a git repo with the latex to make them.
- This is my twitter, feel free to talk to me about batching or any NLP stuff.
- This is my github which has a lot of (i think) cool NLP stuff on it
- And this is a link to Mead-Baseline, we currently have a lot of layers that don't have these batch stability tests. So if this talked interested you and you want to try your hand at it make a PR maybe?
- Thank you, any questions?