ECE 1508S2: Applied Deep Learning

Chapter 6: Recurrent NNs

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In many applications, we have sequence data, e.g.,

- speech data which is usually a long time series
- text data which is sequence of words and letters
- financial data that is typically a time-dependent sequence of values

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 - → This can easily make the NN size infeasible
- On another hand, we do not have so much features
 - \downarrow Just think of a long text, where we need to predict the next word in it

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We need to develop some techniques to handle such data

First, let's see some examples!



- This is a classification problem
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- What about the data?
 - → We sample the audio signal at rate 44.1 kHz and quantize the samples

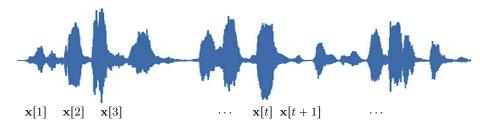


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 - \downarrow We put every N successive samples in a frames \equiv vector of samples



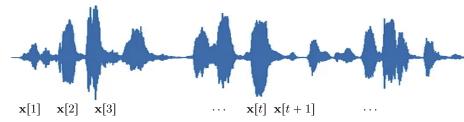
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Assume we listen to a 15-minute talk: we want to find out whether it is about sport or science

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 - □ Processing all samples together seems to be an unnecessary hardness

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But do we need to pass them all together through an NN?

- It does not seem to be!
- It does not generalize well
 - \downarrow We want to classify shorter and longer talks as well



Now let's consider another example: we have a long text and want to learn what is the next word in the sentence

- This is a prediction task
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...therapy. UofT undergraduate students explore the use of AI to treat speech

$$\mathbf{x}[1] \quad \mathbf{x}[2] \quad \mathbf{x}[3]$$

$$\mathbf{x}[t] \mathbf{x}[t+1] \cdots \mathbf{x}[T] \mathbf{y} =?$$

$$\mathbf{x}[T]$$



Another example: we have a sequence of stock prices and are interested in the future price

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Let's start with a simple example: we want to train a neural network that gets a sentence and complete the next word

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- + How does the training dataset look like?
- We are given with several long texts in the same context: at each entry of sequence in each of these texts the whole sequence is the data-point and the next word is label

Applied Deep Learning

Let's make some specification to clarify the problem

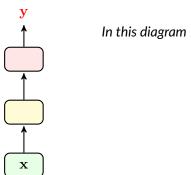
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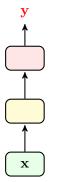
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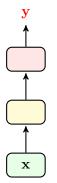
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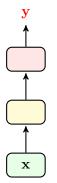
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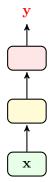
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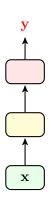


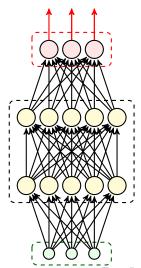
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- Red box is the output layer
- Arrows refer to all links between the layers



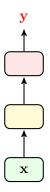
For instance, we could think of following equivalence





Predicting Next Word: MLP

Let's try solving this problem with a simple MLP



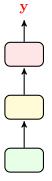
We have a fully-connected FNN that

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We train this MLP

we go over all text

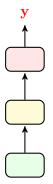
Predicting Next Word: MLP



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. Julia has been nominated to receive Alexander von Humboldt Prize for her

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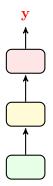


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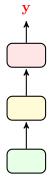
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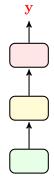
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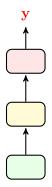
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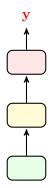
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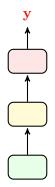
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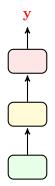
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Does FNN predict "her"?



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Applied Deep Learning Chapter 6: RNNs © A.



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Does FNN predict "her"? No! How can it remember we are talking about Julia?!

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Applied Deep Learning

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 many of these predictions are irrelevant, e.g.,
 - inominated to" is followed by "receive": has nothing to say about "her"

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This indicates that we need to make a memory component for our NN

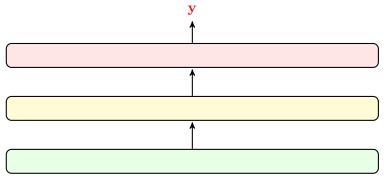
Predicting Next Word: Large MLP

Maybe we could give more inputs to the FNN!



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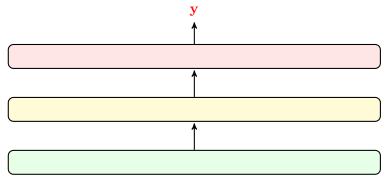
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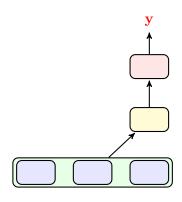
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But, what is Julia has been mentioned 10 pages ago? Forget about large MLPs!

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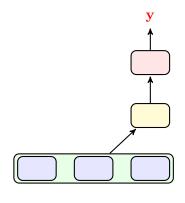
Let's now think about CNNs



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- ullet takes N inputs, i.e., single entry
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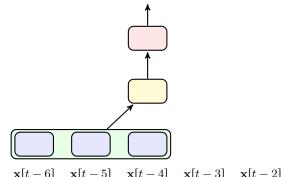


We have a fully-connected FNN that

- takes N inputs, i.e., single entry
- ullet returns N outputs, i.e., predicted next word

We use convolution to

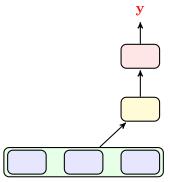
- look into a larger input by a filter of size N
- extract features from a larger part of text and pass it to hidden layers



 $\mathbf{x}[t-6]$ $\mathbf{x}[t-5]$ $\mathbf{x}[t-4]$ $\mathbf{x}[t-3]$ $\mathbf{x}[t-2]$ $\mathbf{x}[t-1]$ $\mathbf{x}[t]$

.. Julia has been nominated to receive Alexander von Humboldt Prize for her

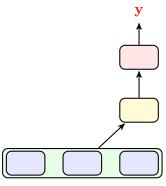
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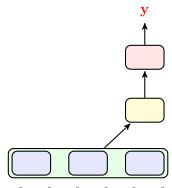
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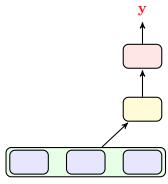
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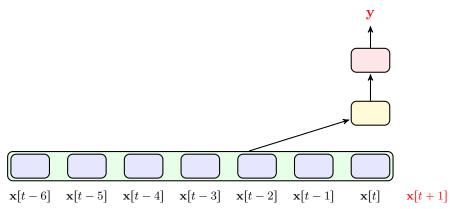
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Yet, it doesn't seem to remember Julia! Unless we slide over the whole text!

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Predicting Next Word: Large CNN



... Julia has been nominated to receive Alexander von Humboldt Prize for her

Though better than MLP, it is still infeasible to track the whole text

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$$\mathbf{x}[t-6]$$
 $\mathbf{x}[t-5]$ $\mathbf{x}[t-4]$ $\mathbf{x}[t-3]$ $\mathbf{x}[t-2]$ $\mathbf{x}[t-1]$ $\mathbf{x}[t]$ $\mathbf{x}[t+1]$... Julia has been nominated to receive Alexander von Humboldt Prize for her

The problem with all architectures we know is that they have finite memory



Applied Deep Learning

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The problem with all architectures we know is that they have finite memory

- They can only remember from their input
 - we always give them independent inputs with similar features
 - they gradually learn to connect any of such inputs toits label
- If we need to remember for long time we have to give them huge inputs
- But the memory component does not seem to be so huge

 - ↓ If text now switches to Theodore we should refresh our memory that we are talking about a "single person" and "male"

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Finite Memory Component with Infinite Response

Component We Miss

We need to extract a memory component from our data that is finite in size but has been influenced (at least theoretically) infinitely



Applied Deep Learning

Finite Memory Component with Infinite Response

Component We Miss

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State-space model can help us building such memory component: it is widely used in control theory to describe evolution of a system over time

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State-space model can help us building such memory component: it is widely used in control theory to describe evolution of a system over time

Assume $\mathbf{x}[t]$ is an input to a system at time t: the system returns an output $\mathbf{y}[t]$ to this input and a state variable $\mathbf{s}[t]$. The output in the next time, i.e., t+1, depends on the new input and current state, i.e.,

$$y[t+1], s[t+1] = f(x[t+1], s[t])$$

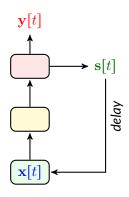
In the above representation: the state is a finite-size variable that carries information for infinitely long time

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Applied Deep Learning Chapter 6: RNNs

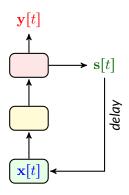
We can look at a NN as a state-dependent system



In this architecture, the NN

- takes $\mathbf{x}[t]$ and previous state $\mathbf{s}[t-1]$ as inputs
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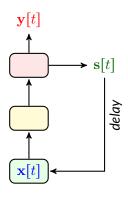
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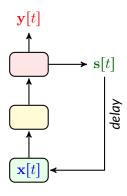
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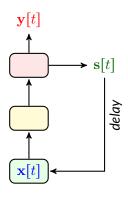
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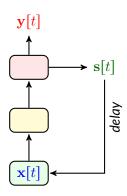
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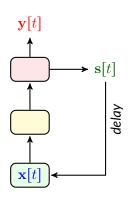
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Theoretically, this NN has infinite time response



Say NN initiates with some s[0]. It takes only sample x[1] and no other time samples is given to it. At time t,

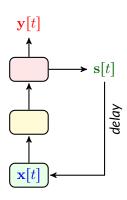
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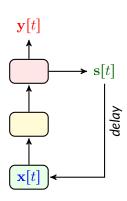
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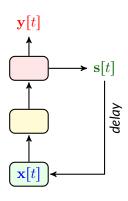
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- $\mathbf{y}[t]$ depends on $\mathbf{s}[t-1]$
- $\mathbf{s}[t-1]$ depends on $\mathbf{s}[t-2]$
- $\mathbf{s}[1]$ depends on $\mathbf{x}[1]$

This means that $\mathbf{y}[t]$ still remembers $\mathbf{x}[1]$

It seems that for our purpose state-space model helps extracting a good memory components. The challenge is to design a good state-dependent NN



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- + Why is it a challenge? We make an NN with input (x, s) and output (y, s') and then train it!
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 - by how should we specify the state as we do not have a clear clue about it

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 - □ empirical risk depends on weights in the NN
 - it also depends previous memory components which can be learnable
 - if we want to train the model efficiently, we need to get back over time and update all those memory components!

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Several attempts have been done: we look briefly into two of them

- Jordan Network
 - ightharpoonup proposed by Michael Jordan¹ in 1986
 - it computes the state variable to be a simple moving average

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¹Professor at CS Department of University of Berkeley; check his page

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Recurrent NNs \equiv RNNs

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RNN

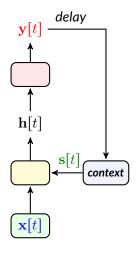
RNN is an NN with state-variable: the term recurrent refers to the connection between former state and new output

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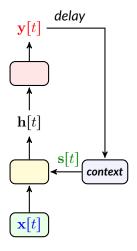
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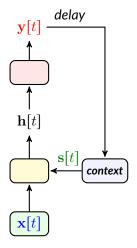
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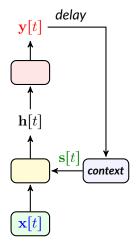
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The output is

$$\mathbf{y}[t] = f\left(\mathbf{W}_2\mathbf{h}[t]\right)$$

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Jordan Network

Jordan Network has a memory component with infinte response: say we set initial state to zero, give $\mathbf{x}[1]$, and keep the input zero for the rest of time; then,

```
\mathbf{y[1]} = f\left(\mathbf{W}_{2}f\left(\mathbf{W}_{1}\mathbf{x[1]}\right)\right)
\mathbf{y[2]} = f\left(\mathbf{W}_{2}f\left(\mathbf{W}_{m}\mathbf{y[1]}\right)\right) = f\left(\mathbf{W}_{2}f\left(\mathbf{W}_{m}f\left(\mathbf{W}_{2}f\left(\mathbf{W}_{1}\mathbf{x[1]}\right)\right)\right)\right)
\vdots
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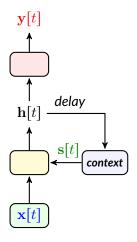
```
\mathbf{y[1]} = f\left(\mathbf{W}_{2}f\left(\mathbf{W}_{1}\mathbf{x[1]}\right)\right)
\mathbf{y[2]} = f\left(\mathbf{W}_{2}f\left(\mathbf{W}_{m}\mathbf{y[1]}\right)\right) = f\left(\mathbf{W}_{2}f\left(\mathbf{W}_{m}f\left(\mathbf{W}_{2}f\left(\mathbf{W}_{1}\mathbf{x[1]}\right)\right)\right)\right)
\vdots
```

But, the network does not learn how to remember

- It takes a fixed moving average as memory component
- It only learns how to use this pre-defined memory for prediction
 - $\,\,\,\,\,\,\,\,$ We could say that it implicitly learn to remember by learning \mathbf{W}_m

Applied Deep Learning

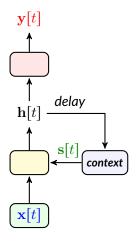
Elman Network uses output of hidden layer as state: also called hidden state



The proposal was a shallow NN

• It starts with some initial hidden state h[0]

Elman Network uses output of hidden layer as state: also called hidden state



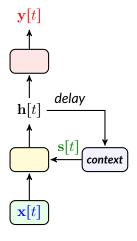
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$$\mathbf{s}[t] = \mathbf{h}[t-1]$$



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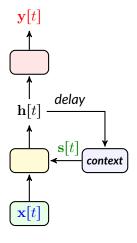
$$\mathbf{s}[t] = \mathbf{h}[t-1]$$

The hidden layer then computes

$$\mathbf{h}[t] = f\left(\mathbf{W}_{1}\mathbf{x}[t] + \mathbf{W}_{m}\mathbf{s}[t]\right)$$

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The output is

$$\mathbf{y}[t] = f\left(\mathbf{W}_2\mathbf{h}[t]\right)$$

←□ → ←□ → ←□ → ←□ → ←□

Similar to Jordan Network, Elman Network has a memory component with infinte response: with zero initial hidden state, we have

$$\mathbf{y}[1] = f\left(\mathbf{W}_{2}f\left(\mathbf{W}_{1}\mathbf{x}[1]\right)\right)$$

$$\mathbf{y}[2] = f\left(\mathbf{W}_{2}f\left(\mathbf{W}_{m}\mathbf{h}[1]\right)\right) = f\left(\mathbf{W}_{2}f\left(\mathbf{W}_{m}f\left(\mathbf{W}_{1}\mathbf{x}[1]\right)\right)\right)$$

$$\vdots$$

Elman network also learns how to remember only implicitly



Though Jordan and Elman Networks had memory, they did not get trained accurately over time, i.e., they simplified the solution to the second challenge

- + What is really this challenge?
- We are going to deal with it in next sections, but let's see it on these simple networks first



Applied Deep Learning

³We will see that this is not always this easy!

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Applied Deep Learning

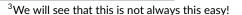
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Let's assume a simple setting: we are to train our NN on single data sequence

- We have the sequence $\mathbf{x}[1], \dots, \mathbf{x}[T]$ as the data-point
- For each entry of this sequence we have the true label
- We are able to compute the loss between outputs and true labels as³

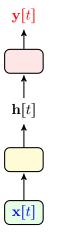
$$\hat{R} = \sum_{t=1}^{T} \mathcal{L}\left(\mathbf{y}[t], \mathbf{v}[t]\right)$$

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³We will see that this is not always this easy!

Recap: Basic FNN

For sake if comparison, let's first train a basic FNN on this data sequence



We have a shallow FNN

The hidden layer computes

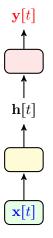
$$\mathbf{h}[t] = f\left(\mathbf{W}_1 \mathbf{x}[t]\right)$$

• The output is

$$\mathbf{y}[t] = f\left(\mathbf{W}_2\mathbf{h}[t]\right)$$

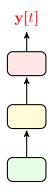
Recap: Basic FNN

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How do we train this FNN?

- We compute the gradients $\nabla_{\mathbf{W}_1} \hat{R}$ and $\nabla_{\mathbf{W}_2} \hat{R}$ \hookrightarrow We do it via backpropagation
- We apply gradient descent



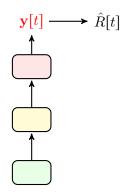
$$\mathbf{x}[1]$$

 $\mathbf{x}[2]$

 \cdots $\mathbf{x}[t-1]$ $\mathbf{x}[t]$

 \cdots $\mathbf{x}[T-1]$ $\mathbf{x}[T]$

$$\hat{R} = \sum_{t=1}^{T} \underbrace{\mathcal{L}\left(\mathbf{y}[t], \mathbf{v}[t]\right)}_{\hat{R}[t]}$$



$$\mathbf{x}[1]$$

$$\mathbf{x}[2]$$

$$\mathbf{x}[t]$$

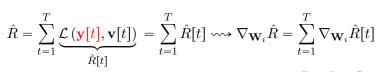
$$\mathbf{x}[2]$$
 \cdots $\mathbf{x}[t-1]$ $\mathbf{x}[t]$ \cdots $\mathbf{x}[T-1]$ $\mathbf{x}[T]$

$$\mathbf{x}[2]$$

$$\mathbf{x}[t-1]$$

$$\mathbf{x}[\iota]$$

$$= \sum_{i=1}^{T} \nabla_{\mathbf{W}_{i}} \hat{R}[i]$$



Let's learn W_1 and to ease computation we use our cheating notation, i.e., use \circ to show any product:



Let's learn \mathbf{W}_1 and to ease computation we use our cheating notation, i.e., use \circ to show any product: to compute the gradient we start with the output

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We know that $\mathbf{y}[t] = f(\mathbf{W}_2\mathbf{h}[t])$, so we can write

$$\nabla_{\mathbf{W}_{1}}\mathbf{y}[t] = \nabla_{\mathbf{h}[t]}\mathbf{y}[t] \circ \nabla_{\mathbf{W}_{1}}\mathbf{h}[t]$$
$$= \left(\dot{f}\left(\mathbf{W}_{2}\mathbf{h}[t]\right) \circ \mathbf{W}_{2}\right) \circ \nabla_{\mathbf{W}_{1}}\mathbf{h}[t]$$

Applied Deep Learning

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What about $\nabla_{\mathbf{W}_1}\mathbf{h}[t]$? We keep on backward!



Applied Deep Learning

Up to now, we have

$$\nabla_{\mathbf{W}_1} \hat{R}[t] = \nabla_{\mathbf{y}[t]} \hat{R}[t] \circ \nabla_{\mathbf{h}[t]} \mathbf{y}[t] \circ \nabla_{\mathbf{W}_1} \mathbf{h}[t]$$



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We use the fact that $\mathbf{h}[t] = f(\mathbf{W}_1 \mathbf{x}[t])$

$$\nabla_{\mathbf{W}_1} \mathbf{h}[t] = \nabla_{\mathbf{W}_1} f(\mathbf{W}_1 \mathbf{x}[t]) + \nabla_{\mathbf{x}[t]} \mathbf{h}[t] \circ \nabla_{\mathbf{W}_1} \mathbf{x}[t]$$



Up to now, we have

$$\nabla_{\mathbf{W}_1} \hat{R}[t] = \nabla_{\mathbf{y}[t]} \hat{R}[t] \circ \nabla_{\mathbf{h}[t]} \mathbf{y}[t] \circ \nabla_{\mathbf{W}_1} \mathbf{h}[t]$$

We use the fact that $\mathbf{h}[t] = f(\mathbf{W}_1 \mathbf{x}[t])$

$$\begin{split} \nabla_{\mathbf{W}_1}\mathbf{h}[t] &= \nabla_{\mathbf{W}_1}f\left(\mathbf{W}_1\mathbf{x}[t]\right) + \nabla_{\mathbf{x}[t]}\mathbf{h}[t] \circ \nabla_{\mathbf{W}_1}\mathbf{x}[t] \\ &= \dot{f}\left(\mathbf{W}_1\mathbf{x}[t]\right) \circ \mathbf{x}[t] + \left(\dot{f}\left(\mathbf{W}_1\mathbf{x}[t]\right) \circ \mathbf{W}_1\right) \circ \underbrace{\mathbf{0}}_{\mathbf{x}[t] \text{ is not a function of } \mathbf{W}_1 \end{split}$$

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$$\nabla_{\mathbf{W}_1} \hat{R}[t] = \nabla_{\mathbf{y}[t]} \hat{R}[t] \circ \nabla_{\mathbf{h}[t]} \mathbf{y}[t] \circ \nabla_{\mathbf{W}_1} \mathbf{h}[t]$$

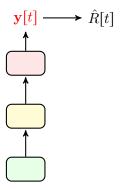
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$$\begin{split} \nabla_{\mathbf{W}_1} \mathbf{h}[t] &= \nabla_{\mathbf{W}_1} f\left(\mathbf{W}_1 \mathbf{x}[t]\right) + \nabla_{\mathbf{x}[t]} \mathbf{h}[t] \circ \nabla_{\mathbf{W}_1} \mathbf{x}[t] \\ &= \dot{f}\left(\mathbf{W}_1 \mathbf{x}[t]\right) \circ \mathbf{x}[t] + \left(\dot{f}\left(\mathbf{W}_1 \mathbf{x}[t]\right) \circ \mathbf{W}_1\right) \circ \underbrace{\mathbf{0}}_{\mathbf{x}[t]} \text{ is not a function of } \mathbf{W}_1 \end{split}$$

Therefore, we end chain rule here

$$\nabla_{\mathbf{W}_1} \hat{R}[t] = \nabla_{\mathbf{y}[t]} \hat{R}[t] \circ \nabla_{\mathbf{h}[t]} \mathbf{y}[t] \circ \nabla_{\mathbf{W}_1} \mathbf{h}[t]$$



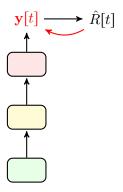


$$\mathbf{x}[1]$$

 $\mathbf{x}[2]$

 \cdots $\mathbf{x}[t-1]$ $\mathbf{x}[t]$ \cdots $\mathbf{x}[T-1]$ $\mathbf{x}[T]$

$$\nabla_{\mathbf{W}_1} \hat{R}[t] =$$

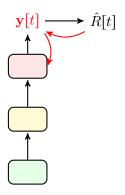


$$\mathbf{x}[1]$$

 $\mathbf{x}[2]$

 \cdots $\mathbf{x}[t-1]$ $\mathbf{x}[t]$ \cdots $\mathbf{x}[T-1]$ $\mathbf{x}[T]$

$$\nabla_{\mathbf{W}_1} \hat{R}[t] = \nabla_{\mathbf{y}[t]} \hat{R}[t]$$

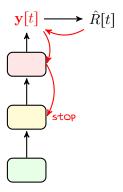


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 \cdots $\mathbf{x}[t-1]$ $\mathbf{x}[t]$ \cdots $\mathbf{x}[T-1]$ $\mathbf{x}[T]$

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$$\mathbf{x}[1]$$
 $\mathbf{x}[2]$

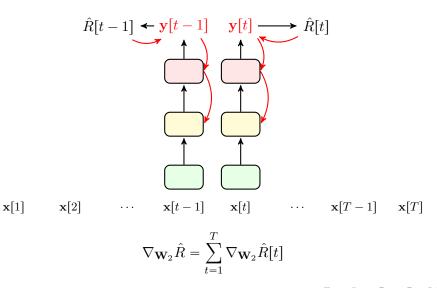
$$\mathbf{x}[t-1]$$

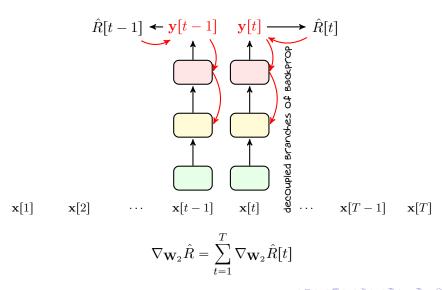
$$\mathbf{x}[t]$$

$$\cdots$$
 $\mathbf{x}[t-1]$ $\mathbf{x}[t]$ \cdots $\mathbf{x}[T-1]$ $\mathbf{x}[T]$

$$\mathbf{x}[T]$$

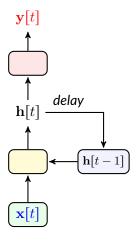
$$\nabla_{\mathbf{W}_1} \hat{R}[t] = \nabla_{\mathbf{y}[t]} \hat{R}[t] \circ \nabla_{\mathbf{h}[t]} \mathbf{y}[t] \circ \nabla_{\mathbf{W}_1} \mathbf{h}[t]$$





Training a Basic RNN

Now, let's train Elman network on this sequence



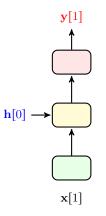
The proposal was a shallow NN

- It starts with some initial hidden state h[0]
- The hidden layer then computes

$$\mathbf{h}[t] = f\left(\mathbf{W}_1 \mathbf{x}[t] + \mathbf{W}_{\mathrm{m}} \mathbf{h}[t-1]\right)$$

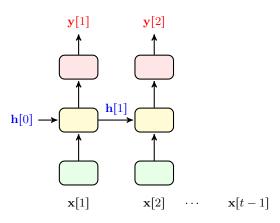
The output is

$$\mathbf{y}[t] = f\left(\mathbf{W}_2\mathbf{h}[t]\right)$$

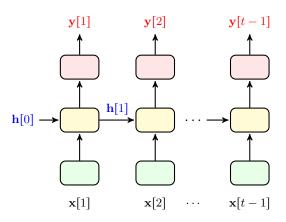


 $\mathbf{x}[2] \quad \cdots \quad \mathbf{x}[t-1]$

 $\mathbf{x}[t] \quad \cdots \quad \mathbf{x}[T]$

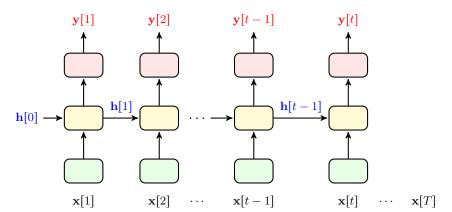


 $\mathbf{x}[t] \quad \cdots \quad \mathbf{x}[T]$

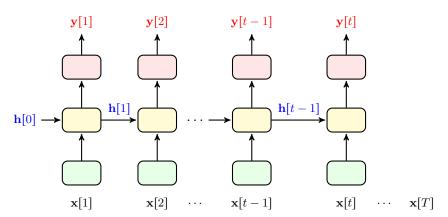


 $\mathbf{x}[t] \quad \cdots \quad \mathbf{x}[T]$

4 D > 4 B > 4 E > 4 E > E *) Q (*

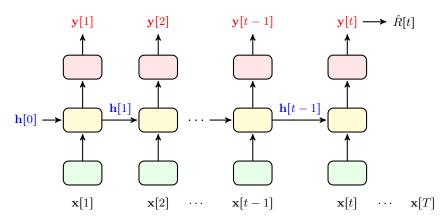


Applied Deep Learning



$$\hat{R} = \sum_{t=1}^{T} \underbrace{\mathcal{L}(\mathbf{y}[t], \mathbf{v}[t])}_{\hat{R}[t]} = \sum_{t=1}^{T} \hat{R}[t]$$

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Let's again try to learn W_1 : we start with the output

$$\nabla_{\mathbf{W}_1} \hat{R}[t] = \nabla_{\mathbf{y}[t]} \hat{R}[t] \circ \nabla_{\mathbf{W}_1} \mathbf{y}[t]$$



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$$\nabla_{\mathbf{W}_1} \hat{R}[t] = \nabla_{\mathbf{y}[t]} \hat{R}[t] \circ \nabla_{\mathbf{W}_1} \mathbf{y}[t]$$

We know that $\mathbf{y}[t] = f(\mathbf{W}_2\mathbf{h}[t])$, so we can write

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$$\nabla_{\mathbf{W}_1}\mathbf{h}[t] =$$

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Next, we note that $\mathbf{h}[t] = f\left(\mathbf{W}_1\mathbf{x}[t] + \mathbf{W}_{\mathrm{m}}\mathbf{h}[t-1]\right)$

$$\nabla_{\mathbf{W}_1} \mathbf{h}[t] = \nabla_{\mathbf{W}_1} f\left(\mathbf{W}_1 \mathbf{x}[t] + \mathbf{W}_{\mathrm{m}} \mathbf{h}[t-1]\right)$$



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Next, we note that $h[t] = f(\mathbf{W}_1 \mathbf{x}[t] + \mathbf{W}_m \mathbf{h}[t-1])$

$$\nabla_{\mathbf{W}_{1}}\mathbf{h}[t] = \nabla_{\mathbf{W}_{1}}f\left(\mathbf{W}_{1}\mathbf{x}[t] + \mathbf{W}_{m}\mathbf{h}[t-1]\right) + \nabla_{\mathbf{x}[t]}\mathbf{h}[t] \circ \underbrace{\nabla_{\mathbf{W}_{1}}\mathbf{x}[t]}_{\mathbf{0}}$$



Let's again try to learn W_1 : we start with the output

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Next, we note that $\mathbf{h}[t] = f\left(\mathbf{W}_1\mathbf{x}[t] + \mathbf{W}_{\mathrm{m}}\mathbf{h}[t-1]\right)$

$$\nabla_{\mathbf{W}_{1}}\mathbf{h}[t] = \nabla_{\mathbf{W}_{1}}f\left(\mathbf{W}_{1}\mathbf{x}[t] + \mathbf{W}_{m}\mathbf{h}[t-1]\right) + \nabla_{\mathbf{x}[t]}\mathbf{h}[t] \circ \underbrace{\nabla_{\mathbf{W}_{1}}\mathbf{x}[t]}_{\mathbf{0}} + \nabla_{\mathbf{h}[t-1]}\mathbf{h}[t] \circ \underbrace{\nabla_{\mathbf{W}_{1}}\mathbf{h}[t-1]}_{\mathbf{2}}$$

Well! We know that
$$\mathbf{h}[t-1]=f\left(\mathbf{W}_1\mathbf{x}[t-1]+\mathbf{W}_{\mathrm{m}}\mathbf{h}[t-2]\right)$$

$$\nabla_{\mathbf{W}_1}\mathbf{h}[t-1]=$$



Well! We know that
$$\mathbf{h}[t-1] = f\left(\mathbf{W}_1\mathbf{x}[t-1] + \mathbf{W}_m\mathbf{h}[t-2]\right)$$

$$\nabla_{\mathbf{W}_1} \mathbf{h}[t-1] = \nabla_{\mathbf{W}_1} f\left(\mathbf{W}_1 \mathbf{x}[t-1] + \mathbf{W}_{\mathrm{m}} \mathbf{h}[t-2]\right)$$



Well! We know that
$$\mathbf{h}[t-1] = f(\mathbf{W}_1 \mathbf{x}[t-1] + \mathbf{W}_m \mathbf{h}[t-2])$$

$$\nabla_{\mathbf{W}_1} \mathbf{h}[t-1] = \nabla_{\mathbf{W}_1} f(\mathbf{W}_1 \mathbf{x}[t-1] + \mathbf{W}_m \mathbf{h}[t-2])$$

$$\nabla_{\mathbf{W}_{1}}\mathbf{h}[t-1] = \nabla_{\mathbf{W}_{1}}f\left(\mathbf{W}_{1}\mathbf{x}[t-1] + \mathbf{W}_{m}\mathbf{h}[t-2]\right) + \nabla_{\mathbf{x}[t-1]}\mathbf{h}[t-1] \circ \underbrace{\nabla_{\mathbf{W}_{1}}\mathbf{x}[t-1]}_{\mathbf{0}}$$

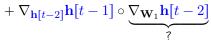


Applied Deep Learning

Chapter 6: RNNs

Well! We know that
$$\mathbf{h}[t-1] = f(\mathbf{W}_1 \mathbf{x}[t-1] + \mathbf{W}_m \mathbf{h}[t-2])$$

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Applied Deep Learning

Chapter 6: RNNs

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$$+ \nabla_{\mathbf{h}[t-2]}\mathbf{h}[t-1] \circ \underbrace{\nabla_{\mathbf{W}_{1}}\mathbf{h}[t-2]}_{?}$$

We are not still done! We know
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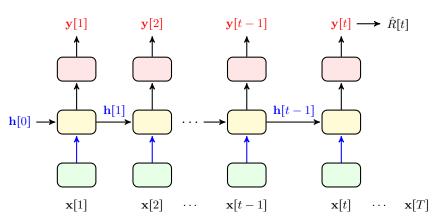
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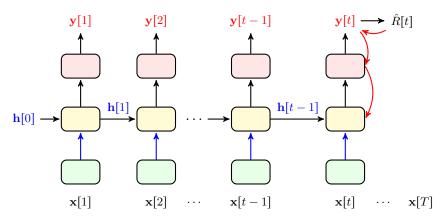
We should in fact pass all the way back to the initial time interval time!



Note that all blue edges are representing \mathbf{W}_1



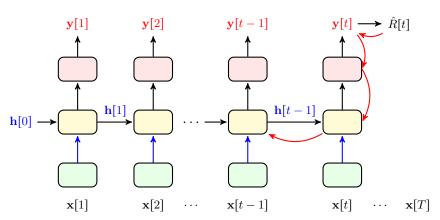
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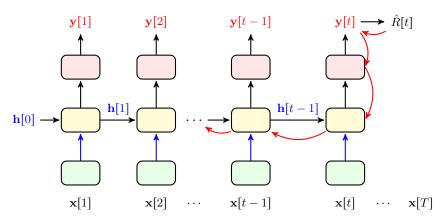
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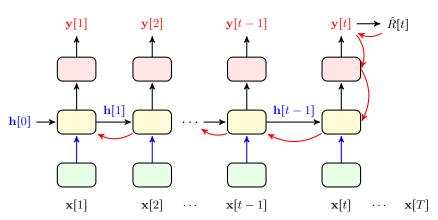


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Learning Through Time: Elman Network

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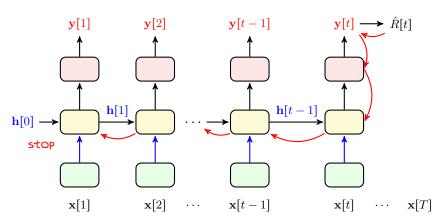


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Applied Deep Learning

Moral of Story

To learn how to remember, we need to train our RNN through time: at each time interval, we should move all the way back to origin to find out how exactly we should change the weights!



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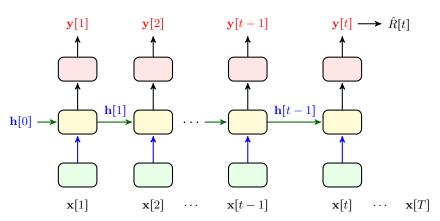
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This means that Elman did not really addressed the second challenge!

Learning Through Time: Elman's Approximation

Elman treated it as a simple FNN with only one extra input!

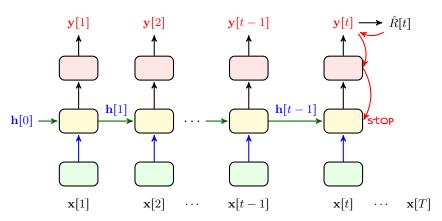


Training \mathbf{W}_1 is exactly as FNN. We just have one extra \mathbf{W}_{m} here!

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RNNs: Need to Learn Memory

Though appreciated, Elman and Jordan Networks did not do the job

- 1 Their memory component is rather simple
 - → We should use deeper models that enable advanced memory components
- 2 They do not really learn how to remember



RNNs: Need to Learn Memory

Though appreciated, Elman and Jordan Networks did not do the job

- Their memory component is rather simple
- 2 They do not really learn how to remember

These led us to development of RNNs!

RNN: Less Generic Definition

An RNN can be designed with any known architecture by letting NN also learn from its past features and outputs. This new enabling is called recurrence

RNN: Dataset and Learning Setting

We are now looking into a supervised learning problem where we are to

learn label from a sequence of data that has generally a temporal correlation



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Let's denote the sequence with $\mathbf{x}[1], \dots, \mathbf{x}[T]$

Temporal Correlation

By temporal correlation we mean that entries at other time instances carry information about one particular entry $\mathbf{x}[t]$

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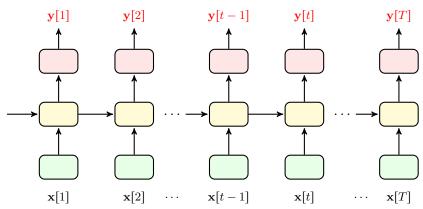
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Temporal Correlation

By temporal correlation we mean that entries at other time instances carry information about one particular entry $\mathbf{x}[t]$

- + But how is a label assigned to this sequence?
- Well, that can be of various forms!

We considered a very simple case: many-to-many type I

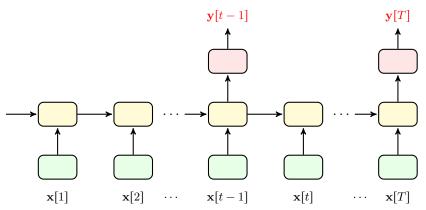


In this case, every entry has a label

4 D P 4 B P 4 E P 4 E P 4 E P 4 C

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We considered a very simple case: many-to-many type II

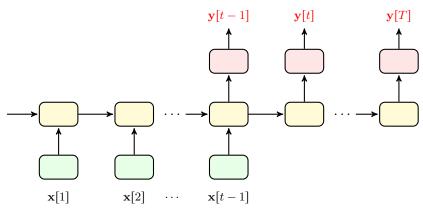


In this case, we get a label once after multiple entries

speech recognition: $\mathbf{x}[t]$ is a small part of speech and $\mathbf{y}[t]$ says what is a every couple of minutes about

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We considered a very simple case: many-to-many type III

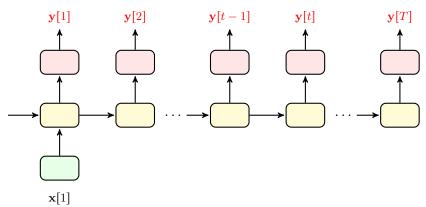


In this case, we start to get labels after some delay

Ly language translation: $\mathbf{x}[1], \dots, \mathbf{x}[t-1]$ is a sentence in German and $\mathbf{y}[t-1], \dots, \mathbf{y}[T]$ is its translation to English

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We considered a very simple case: one-to-many

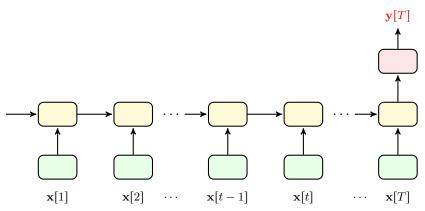


In this case, we get only one input data and have a sequence of labels

 \downarrow image captioning: $\mathbf{x}[1]$ is an image and $\mathbf{y}[1], \dots, \mathbf{y}[T]$ is a caption describing what is inside the image

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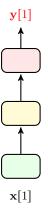
We considered a very simple case: many-to-one



In this case, we get only one label for a whole input sequence

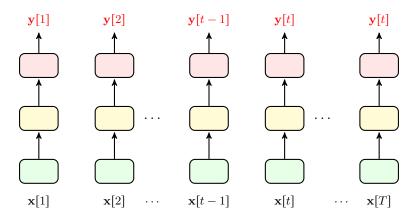
sequence classification: $\mathbf{x}[1], \dots, \mathbf{x}[T]$ is a speech and $\mathbf{y}[T]$ says if this speech is constructive or destructive

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In fact, FNNs are one-to-one RNNs

4 □ ▶ 4 ₫ ▶ 4 ₫ ▶ 4 ₫ ▶ 2 *)Q(*



In fact, FNNs are one-to-one RNNs

- ⇒ we can think of every data-point as a sequence of length one, or

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RNN: General Form

To construct an RNN, we can use any module that we have learned:

- we can use a fully-connected layer
- we can use a convolutional layer
- we can use a residual unit
- we may use an inception unit used in GoogLeNet

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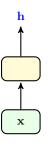
. . .

But, classical implementations use fully-connected layers

We can use one layer to make a shallow RNN or multiple to make a deep RNN

- + Don't we do any change to them?
- Not a serious change
 - \rightarrow We may change the activation to $tanh(\cdot)$: we will see later why
 - → We expand the input dimension: since we need to also give memory as input

Let's break it down a bit: say the yellow box is a fully-connected layer

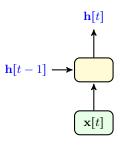


This layer gets input $\mathbf{x} \in \mathbb{R}^N$ and returns activated feature $\mathbf{h} \in \mathbb{R}^M$

- We replace its activation by $tanh(\cdot)$
 - → Not necessary, but usually suggested
- We modify it to get a new input in \mathbb{R}^{N+M}

 - $\rightarrow M$ entries for features **h** in time t-1

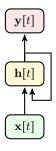
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We show it this way now

- We call $\mathbf{h}[t]$ usually the hidden state
- We pass the hidden state through an output layer
 - ∪ Output is not necessarily corresponding to label

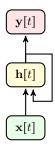
We may show our shallow RNN compactly via the following diagram



In this diagram

- Each edge is a set of weights, e.g., a weight matrix
- The return edge also has a delay in time

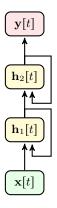
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In this diagram

- Each edge is a set of weights, e.g., a weight matrix
- The return edge also has a delay in time
- + But, isn't that simply Elman Network?
- If we use a fully-connected layer and sigmoid activation; then, Yes! But, Remember that
 - Elman did not train it over time
 - Elman in its model used sigmoid activation

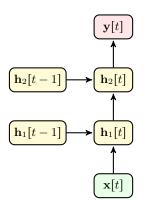
We can add more layers to make a deep RNN



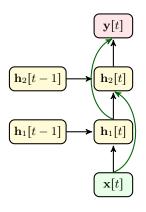
In this diagram

- Each edge is a set of weights, e.g., a weight matrix
- The return edge also has a delay in time
- And, it's no more Elman Network

It might be easier to think of it as the following diagram



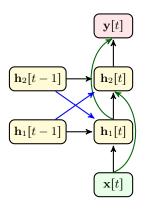
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We can use any module that we like

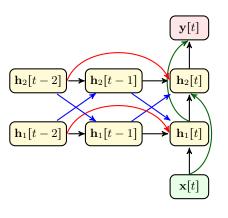
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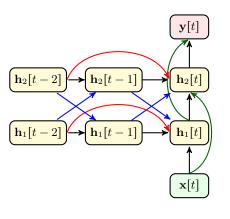
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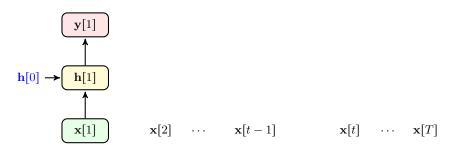


- We may add skip connection
- We could make dense connections
- We may skip over time
- We may replace hidden layers with convolutions or anything else

Shallow RNN: Elman-Like Network

Let's start with simple RNN: a shallow RNN with fully-connected hidden layer

- + You mean Elman Network?
- Yes, but we now try to address the challenges that Elman did not address Let's look ts the flow of information once again



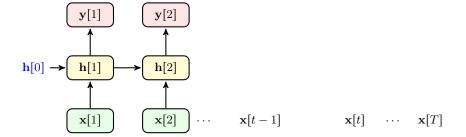
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Chapter 6: RNNs

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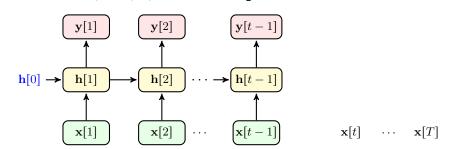


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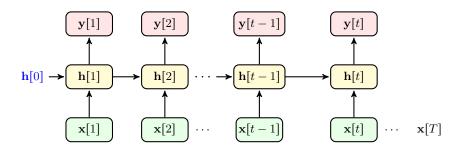
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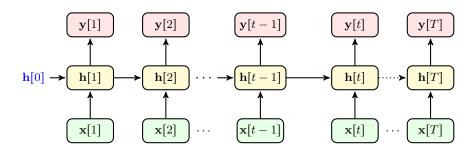
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Applied Deep Learning

Let's specify the learnable parameters with some colors

 $\mathbf{h}[t-1]$

 $\mathbf{h}[t]$

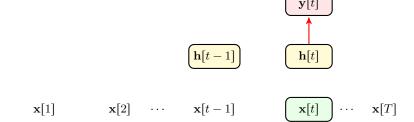
 $\mathbf{x}[1]$

 $\mathbf{x}[2] \quad \cdots \quad \mathbf{x}[t-1]$

 $\mathbf{x}[t]$

Say we set the activation to $f(\cdot)$

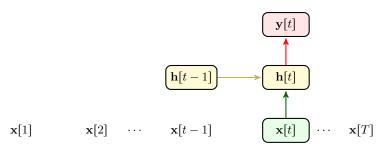
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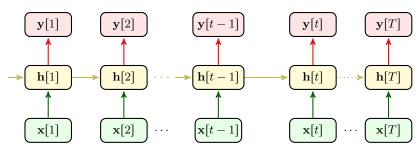


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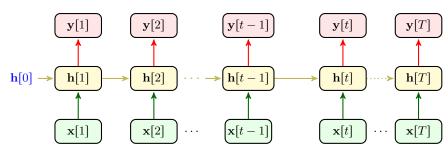
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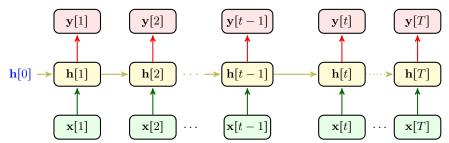


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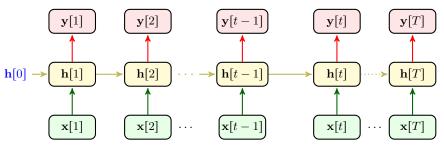
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- ${f 3}$ We can start with any hidden state: this means we can learn ${f h}[0]$ too

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We now want to train this basic RNN



We now want to train this basic RNN



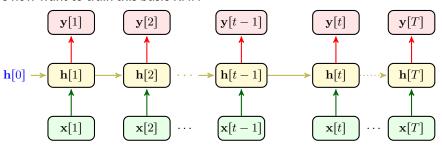
Let's consider training for only one sample

We assume that we have a sequence of labels $\mathbf{v}[1],\dots,\mathbf{v}[T]$

- $oldsymbol{1}$ We know some $\mathbf{v}[t]$ could be empty, e.g., in many-to-one scenario
- 2 So could be some of $\mathbf{x}[t]$'s, e.g., in one-to-many scenario

We however have no problem with that!

We now want to train this basic RNN



Let's consider training for only one sample

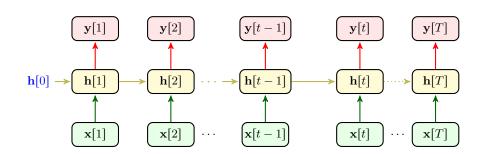
The loss in general can be written as

$$\hat{R} = \mathcal{L}\left(\mathbf{y}[1:T], \mathbf{v}[1:T]\right)$$

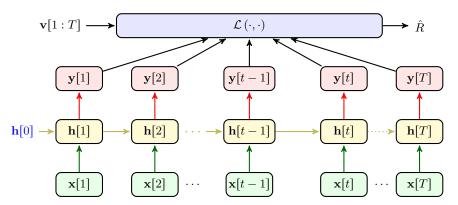
where we use shorten notation $y[1:T] = y[1], \dots, y[T]$

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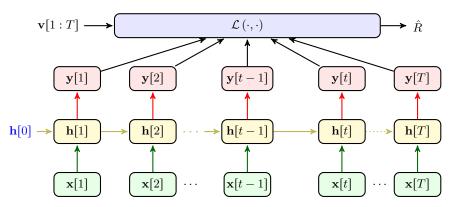
We can think of such a diagram



We can think of such a diagram

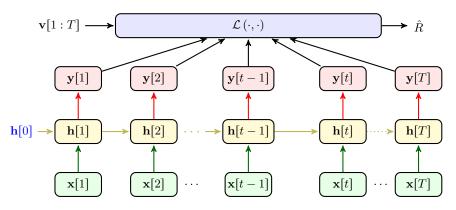


We can think of such a diagram



- + It seems to be hard to get back from \hat{R} to each $\mathbf{y}[t]$
- Well! That's true, but we have some remedy for it

We can think of such a diagram

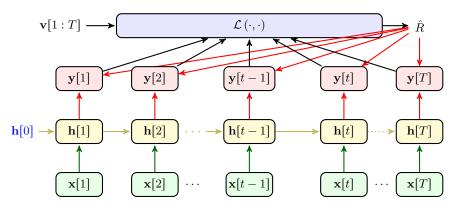


For the moment, we assume that we can compute the $abla_{\mathbf{y}[t]}\hat{R}$ for all t



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We can think of such a diagram



For the moment, we assume that we can

compute the $\nabla_{\mathbf{y}[t]}\hat{R}$ for all $t\equiv$ move backward from \hat{R} to any $\mathbf{y}[t]$



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Starting from \hat{R} , say we want to find $\nabla_{\mathbf{W}_{\mathrm{m}}}\hat{R}$

• Since \hat{R} is a function of $\mathbf{y}[1:T]$, we should write a vectorized chain rule

$$\nabla_{\mathbf{W}_{\mathbf{m}}} \hat{R} =$$

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• Since \hat{R} is a function of $\mathbf{y}[1:T]$, we should write a vectorized chain rule

$$\nabla_{\mathbf{W}_{\mathrm{m}}} \hat{R} = \nabla_{\mathbf{y}[1]} \hat{R} \circ \nabla_{\mathbf{W}_{\mathrm{m}}} \mathbf{y}[1]$$

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• Since \hat{R} is a function of $\mathbf{y}[1:T]$, we should write a vectorized chain rule

$$\nabla_{\mathbf{W}_{\mathrm{m}}} \hat{R} = \nabla_{\mathbf{y}[1]} \hat{R} \circ \nabla_{\mathbf{W}_{\mathrm{m}}} \mathbf{y}[1] + \ldots + \nabla_{\mathbf{y}[T]} \hat{R} \circ \nabla_{\mathbf{W}_{\mathrm{m}}} \mathbf{y}[T]$$

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• Since we assumed that we have $\nabla_{\mathbf{y}[t]}\hat{R}$, our main task is to find

$$abla_{\mathbf{W}_{\mathrm{m}}}\mathbf{y}[t]$$

for all t: we could hence compute it for a general t

 $\,\,\,\,\,\,\,\,\,$ This is a tensor-like gradient, i.e., $\left[\nabla_{\mathbf{W}_{\mathrm{m}}}y_{1}[t],\ldots,\nabla_{\mathbf{W}_{\mathrm{m}}}y_{M}[t]\right]$



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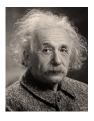
- $\,\,\,\,\,\,\,\,\,$ This is a tensor-like gradient, i.e., $\left[\nabla_{\mathbf{W}_{\mathrm{m}}}y_{1}[t],\ldots,\nabla_{\mathbf{W}_{\mathrm{m}}}y_{M}[t]\right]$
- Apparently, we should apply chain rule for several times!

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+ But, this is going to be exhausting?



- + But, this is going to be exhausting?
- Well, we could again follow Albert Einstein advice!



"Everything should be made as simple as possible, but not simpler!"

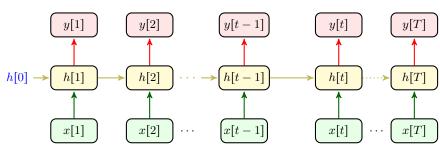
Let's consider a dummy RNN with all variables being scalar

- **1** We have $y[t] = f(w_2h[t])$
- ② We have $h[t] = f(w_1x[t] + w_mh[t-1])$
- **3** We start with hidden state h[0]

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So the diagram gets simplified as below



Starting from \hat{R} , say we want to find $\nabla_{w_{\mathrm{m}}}\hat{R}$: our main task is to find

$$rac{\partial y[t]}{\partial w_{
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To go backward, we note that $y[t] = f(w_2h[t])$

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• we now have y[t], h[t] and x[t] for all t

To go backward, we note that $y[t] = f(w_2h[t])$

• y[t] is a function of h[t]: so we write the chain rule as

$$\frac{\partial y[t]}{\partial w_{\rm m}} = \frac{\partial y[t]}{\partial h[t]} \frac{\partial h[t]}{\partial w_{\rm m}}$$

$$\frac{\partial \mathbf{y[t]}}{\partial w_{\mathrm{m}}} = w_{2} \dot{f}(w_{2} h[t]) \frac{\partial h[t]}{\partial w_{\mathrm{m}}}$$

We keep going backward, by noting that $h[t] = f(w_1x[t] + w_mh[t-1])$

• h[t] is a function of $w_{\rm m}$ and h[t-1]:⁴

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 $^{^4}$ We ignore w_1 and x[t], as they are obviously not functions of $w_{\mathbf{m}}$ (\mathbb{Z}) \mathbb{Z} \mathbb{Z} \mathbb{Z} \mathbb{Z} \mathbb{Z} \mathbb{Z}

$$\frac{\partial \mathbf{y[t]}}{\partial w_{\mathrm{m}}} = w_{2} \dot{f}(w_{2} h[t]) \frac{\partial h[t]}{\partial w_{\mathrm{m}}}$$

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$$\frac{\partial \mathbf{y[t]}}{\partial w_{\mathrm{m}}} = w_{2} \dot{f}(w_{2} h[t]) \frac{\partial h[t]}{\partial w_{\mathrm{m}}}$$

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- So we write the chain rule as

$$\frac{\partial h[t]}{\partial w_{\rm m}} =$$

 4 We ignore w_1 and x[t], as they are obviously not functions of $w_{
m m}$ (3) + (2) + (3) + (3) + (3) + (4) + (5) + (5) + (6) + (7) + (7) + (8) + (8) + (9) + (9) + (10) +

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$$\frac{\partial \mathbf{y[t]}}{\partial w_{\mathrm{m}}} = w_{2} \dot{f}(w_{2} h[t]) \frac{\partial h[t]}{\partial w_{\mathrm{m}}}$$

We keep going backward, by noting that $h[t] = f(w_1x[t] + w_mh[t-1])$

- h[t] is a function of w_{m} and h[t-1]:⁴ \downarrow let's write $h[t] = g\left(w_{\mathrm{m}}, h[t-1]\right)$
- So we write the chain rule as

$$\frac{\partial h[t]}{\partial w_{\mathbf{m}}} = \frac{\partial g}{\partial w_{\mathbf{m}}} \underbrace{\frac{\partial w_{\mathbf{m}}}{\partial w_{\mathbf{m}}}}_{1}$$

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$$\frac{\partial \mathbf{y[t]}}{\partial w_{\mathrm{m}}} = w_{2} \dot{f}(w_{2} h[t]) \frac{\partial h[t]}{\partial w_{\mathrm{m}}}$$

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- h[t] is a function of $w_{\mathbf{m}}$ and h[t-1]:⁴ \downarrow let's write $h[t] = g(w_{\mathbf{m}}, h[t-1])$
- So we write the chain rule as

$$\frac{\partial h[t]}{\partial w_{\mathbf{m}}} = \frac{\partial g}{\partial w_{\mathbf{m}}} \underbrace{\frac{\partial w_{\mathbf{m}}}{\partial w_{\mathbf{m}}}}_{1} + \frac{\partial g}{\partial h[t-1]} \frac{\partial h[t-1]}{\partial w_{\mathbf{m}}}$$

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 $^{^4}$ We ignore w_1 and x[t], as they are obviously not functions of $w_{\mathbf{m}} \cdot \mathbb{Z} \times \mathbb{$

$$\frac{\partial \mathbf{y[t]}}{\partial w_{\mathrm{m}}} = w_{2} \dot{f}(w_{2} h[t]) \frac{\partial h[t]}{\partial w_{\mathrm{m}}}$$

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- So we write the chain rule as

$$\begin{split} \frac{\partial h[t]}{\partial w_{\rm m}} &= \frac{\partial g}{\partial w_{\rm m}} \underbrace{\frac{\partial w_{\rm m}}{\partial w_{\rm m}}}_{1} + \frac{\partial g}{\partial h[t-1]} \frac{\partial h[t-1]}{\partial w_{\rm m}} \\ &= \frac{\partial g}{\partial w_{\rm m}} + \frac{\partial g}{\partial h[t-1]} \frac{\partial h[t-1]}{\partial w_{\rm m}} \end{split}$$

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 $^{^4}$ We ignore w_1 and x[t], as they are obviously not functions of $w_{\mathbf{m}}$ $_{\bullet}$ $_{\bullet}$ $_{\bullet}$ $_{\bullet}$ $_{\bullet}$ $_{\bullet}$ $_{\bullet}$ $_{\bullet}$ $_{\bullet}$ $_{\bullet}$

$$\frac{\partial \mathbf{y[t]}}{\partial w_{\mathrm{m}}} = w_{2} \dot{f}(w_{2} h[t]) \frac{\partial h[t]}{\partial w_{\mathrm{m}}}$$

We keep going backward, by noting that h[t] = f(z[t])

• Let's define $z[t] = w_1x[t] + w_mh[t-1]$



Applied Deep Learning

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- Let's define $z[t] = w_1x[t] + w_mh[t-1]$
- $h[t] = g\left(\mathbf{w_m}, h[t-1]\right)$
 - \rightarrow we can compute $\partial g/\partial w_{\mathrm{m}}=h[t-1]\dot{f}(z[t])$
 - \rightarrow we can compute $\partial g/\partial h[t-1] = w_{\rm m}\dot{f}(z[t])$

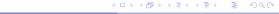
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 - \rightarrow we can compute $\partial g/\partial h[t-1] = w_{\rm m} \dot{f}(z[t])$
- So we can simplify the chain rule as

$$\frac{\partial h[t]}{\partial w_{\rm m}} = h[t-1]\dot{f}(z[t]) + w_{\rm m}\dot{f}(z[t]) \frac{\partial h[t-1]}{\partial w_{\rm m}}$$

$$= \dot{f}(z[t]) \left(h[t-1] + w_{\rm m}\frac{\partial h[t-1]}{\partial w_{\rm m}}\right)$$



$$\frac{\partial y[t]}{\partial w_{\rm m}} = w_2 \dot{f}(w_2 h[t]) \dot{f}(z[t]) \left(h[t-1] + w_{\rm m} \frac{\partial h[t-1]}{\partial w_{\rm m}} \right)$$

We keep going backward:
$$h[t-1] = f(z[t-1])$$

$$\rightarrow$$
 Recall that $z[t-1] = w_1x[t-1] + w_mh[t-2]$



$$\frac{\partial y[t]}{\partial w_{\rm m}} = w_2 \dot{f}(w_2 h[t]) \dot{f}(z[t]) \left(h[t-1] + w_{\rm m} \frac{\partial h[t-1]}{\partial w_{\rm m}} \right)$$

We keep going backward: h[t-1] = f(z[t-1])

- ightharpoonup Recall that $z[t-1]=w_1x[t-1]+w_{\mathrm{m}}h[t-2]$
- We just need to replace t with t-1 in the last derivation

$$\frac{\partial h[t-1]}{\partial w_{\rm m}} = \dot{f}\left(z[t-1]\right) \left(h[t-2] + w_{\rm m} \frac{\partial h[t-2]}{\partial w_{\rm m}}\right)$$

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Backpropagation at time t is hence described by a pair of recursive equations

• At time t, we compute

$$\frac{\partial y[t]}{\partial w_{\rm m}} = w_2 \dot{f}(w_2 h[t]) \frac{\partial h[t]}{\partial w_{\rm m}}$$



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• For each i = t, t - 1, ..., 1, we use

$$\frac{\partial h[i]}{\partial w_{\rm m}} = \dot{f}(z[i]) \left(h[i-1] + w_{\rm m} \frac{\partial h[i-1]}{\partial w_{\rm m}} \right)$$

Backpropagation at time t is hence described by a pair of recursive equations

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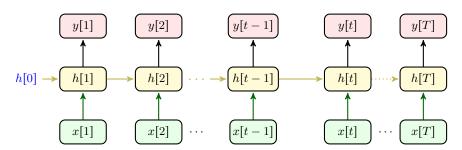
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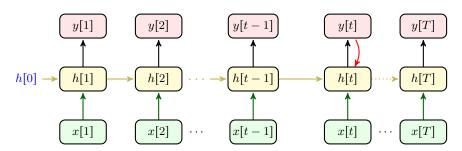
$$\frac{\partial h[i]}{\partial w_{\rm m}} = \dot{f}(z[i]) \left(h[i-1] + w_{\rm m} \frac{\partial h[i-1]}{\partial w_{\rm m}} \right)$$

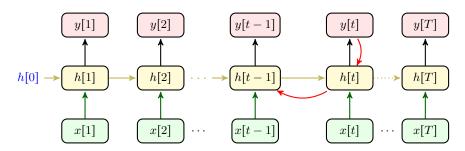
• We stop at i = 1, where we get

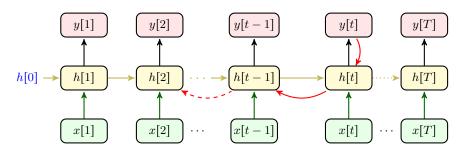
$$\frac{\partial h[1]}{\partial w_{\mathbf{m}}} = \dot{f}(z[1]) \left(h[0] + w_{\mathbf{m}} \underbrace{\frac{\partial h[0]}{\partial w_{\mathbf{m}}}}_{\mathbf{0}} \right) = \dot{f}(z[1]) h[0]$$

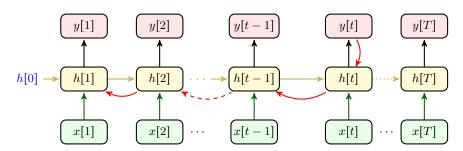
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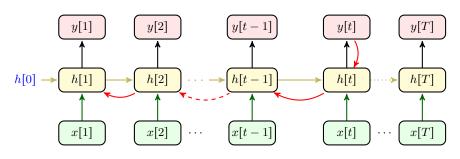








Applied Deep Learning Chapter 6: RNNs



Key Point

Propagating back in time is described via a recursive equation

Recursive equation has an interesting feature

ullet Say we have backpropagated from time t

$$\frac{\partial y[t]}{\partial w_{\rm m}} = w_2 \dot{f}(w_2 h[t]) \frac{\partial h[t]}{\partial w_{\rm m}}$$

 $\,\,\,\,\,\,\,\,\,\,$ We hence have $\partial h[t]/\partial w_{\mathrm{m}}, \partial h[t-1]/\partial w_{\mathrm{m}}, \ldots, \partial h[1]/\partial w_{\mathrm{m}}$



Recursive equation has an interesting feature

Say we have backpropagated from time t

$$\frac{\partial \mathbf{y[t]}}{\partial w_{\mathbf{m}}} = w_2 \dot{f}(w_2 h[t]) \frac{\partial h[t]}{\partial w_{\mathbf{m}}}$$

- $\,\,\,\,\,\,\,\,\,\,$ We hence have $\partial h[t]/\partial w_{\mathrm{m}}, \partial h[t-1]/\partial w_{\mathrm{m}}, \ldots, \partial h[1]/\partial w_{\mathrm{m}}$
- Now, if we want to backpropagate from t-1
 - $\,\,\,\,\,\,\,\,\,$ We already have $\partial h[t-1]/\partial w_{\mathrm{m}}, \partial h[t-2]/\partial w_{\mathrm{m}}, \ldots, \partial h[1]/\partial w_{\mathrm{m}}$

Recursive equation has an interesting feature

ullet Say we have backpropagated from time t

$$\frac{\partial y[t]}{\partial w_{\rm m}} = w_2 \dot{f}(w_2 h[t]) \frac{\partial h[t]}{\partial w_{\rm m}}$$

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- Now, if we want to backpropagate from t-1
 - $\,\,\,\,\,\,\,\,\,$ We already have $\partial h[t-1]/\partial w_{\mathrm{m}}, \partial h[t-2]/\partial w_{\mathrm{m}}, \ldots, \partial h[1]/\partial w_{\mathrm{m}}$

Moral of Story

Pass forward till end of sequence and backpropagate to the beginning just once



Backpropagation Through Time (BPTT)

BPTT():

- 1 Start at t with $\nabla_{\mathbf{W}_{\mathbf{m}}}\mathbf{y}[t] = \mathbf{W}_2 \circ \dot{f}(\mathbf{W}_2\mathbf{h}[t]) \circ \nabla_{\mathbf{W}_{\mathbf{m}}}\mathbf{h}[t]$
- 2 Go back in time as $\nabla_{\mathbf{W}_{\mathbf{m}}}\mathbf{h}[i] = \dot{f}\left(\mathbf{z}[i]\right) \circ \left(\mathbf{h}[i-1] + \mathbf{W}_{\mathbf{m}} \circ \nabla_{\mathbf{W}_{\mathbf{m}}}\mathbf{h}[i-1]\right)$
- 3 Stop at i=1 with $\nabla_{\mathbf{W}_{\mathbf{m}}}\mathbf{h}[1]=\dot{f}\left(\mathbf{z}[1]\right)\circ\mathbf{h}[0]$

Backpropagation Through Time (BPTT)

BPTT():

- 1 Start at t with $\nabla_{\mathbf{W}_{\mathbf{m}}}\mathbf{y}[t] = \mathbf{W}_2 \circ \dot{f}(\mathbf{W}_2\mathbf{h}[t]) \circ \nabla_{\mathbf{W}_{\mathbf{m}}}\mathbf{h}[t]$
- 2 Go back in time as $\nabla_{\mathbf{W}_{\mathbf{m}}}\mathbf{h}[i] = \dot{f}\left(\mathbf{z}[i]\right) \circ \left(\mathbf{h}[i-1] + \mathbf{W}_{\mathbf{m}} \circ \nabla_{\mathbf{W}_{\mathbf{m}}}\mathbf{h}[i-1]\right)$
- 3 Stop at i=1 with $\nabla_{\mathbf{W_m}} \mathbf{h}[1] = \dot{f}(\mathbf{z}[1]) \circ \mathbf{h}[0]$
- + It looks very similar to backpropagation in deep NNs!
- Exactly! Even simple RNN is very deep through time
- Don't we experience vanishing or exploding gradient then!
- Yes! Let's check it out

Let's expand the gradient in our example

$$\frac{\partial y[t]}{\partial w_{\rm m}} = w_2 \dot{f}(w_2 h[t]) \dot{f}(z[t]) h[t-1]$$

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Let's expand the gradient in our example

$$\frac{\partial y[t]}{\partial w_{\rm m}} = w_2 \dot{f}(w_2 h[t]) \dot{f}(z[t]) h[t-1]
+ w_2 \dot{f}(w_2 h[t]) w_{\rm m} \dot{f}(z[t]) \dot{f}(z[t-1]) h[t-2]$$

Let's expand the gradient in our example

$$\frac{\partial y[t]}{\partial w_{\rm m}} = w_2 \dot{f}(w_2 h[t]) \dot{f}(z[t]) h[t-1]
+ w_2 \dot{f}(w_2 h[t]) w_{\rm m} \dot{f}(z[t]) \dot{f}(z[t-1]) h[t-2]
+ w_2 \dot{f}(w_2 h[t]) w_{\rm m}^2 \dot{f}(z[t]) \dot{f}(z[t-1]) \dot{f}(z[t-2]) h[t-3]
+ \cdots$$

Let's expand the gradient in our example

$$\frac{\partial y[t]}{\partial w_{\rm m}} = w_2 \dot{f}(w_2 h[t]) \dot{f}(z[t]) h[t-1]
+ w_2 \dot{f}(w_2 h[t]) w_{\rm m} \dot{f}(z[t]) \dot{f}(z[t-1]) h[t-2]
+ w_2 \dot{f}(w_2 h[t]) w_{\rm m}^2 \dot{f}(z[t]) \dot{f}(z[t-1]) \dot{f}(z[t-2]) h[t-3]
+ \cdots
+ w_2 \dot{f}(w_2 h[t]) w_{\rm m}^{i-1} \left(\prod_{j=0}^{i-1} \dot{f}(z[t-j]) \right) h[t-i]
+ \cdots$$

Let's expand the gradient in our example

$$\frac{\partial y[t]}{\partial w_{\rm m}} = w_2 \dot{f}(w_2 h[t]) \dot{f}(z[t]) h[t-1]
+ w_2 \dot{f}(w_2 h[t]) w_{\rm m} \dot{f}(z[t]) \dot{f}(z[t-1]) h[t-2]
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+ \cdots
+ w_2 \dot{f}(w_2 h[t]) w_{\rm m}^{t-1} \left(\prod_{j=0}^{i-1} \dot{f}(z[t-j]) \right) h[0]$$

This says that gradient of hidden state in i steps back in time is multiplied by

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- If $\dot{f}(\cdot) > 1$ most of the time

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Recall deep NNs: we could see two cases

- If $\dot{f}(\cdot) > 1$ most of the time
 - very old hidden states can explode the gradient
 - → this usually does not happen, because most activations are not like that
- If $f(\cdot) < 1$ most of the time
 - very old hidden states have pretty much no impact on the gradient
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Solutions to vanishing gradient through time are mainly two approaches

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- Use truncated BPTT
 - □ Repeat short-term BPTTs every couple of time intervals

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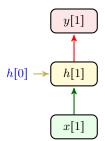
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Solutions to vanishing gradient through time are mainly two approaches

- Use $tanh(\cdot)$ activation
 - $\,\,\,\downarrow\,\,\, \tanh\left(\cdot\right)$ is able to keep memory for a longer period
- Use truncated BPTT
 - □ Repeat short-term BPTTs every couple of time intervals
- Invoke gating approach
 - → This is the most sophisticated approach
 - It was known for a long time, but received attention much later!

To understand the idea of Gating, let's get back to our basic RNN

- **1** We start with hidden state h[0]
- **2** We have $h[t] = f(w_1x[t] + w_mh[t-1])$
- **3** We have $y[t] = f(w_2h[t])$

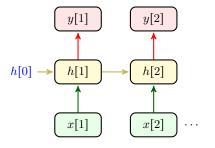


 $x[2] \cdots x[t-1]$

 $x[t] \quad \cdots \quad x[T]$

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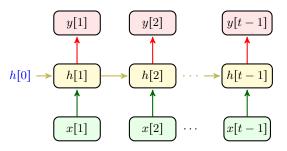


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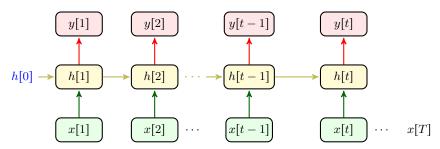
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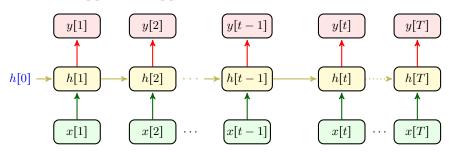
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Looking at h[t] as memory, we can say we are always updating the memory

Recall our motivating example: we wanted to predict the next word

$$\mathbf{x}[t-6]$$
 $\mathbf{x}[t-5]$ $\mathbf{x}[t-4]$ $\mathbf{x}[t-3]$ $\mathbf{x}[t-2]$ $\mathbf{x}[t-1]$ $\mathbf{x}[t]$

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Should we updating the memory all the way from Julia?

- Obviously No!
- How can we do it? Let's try a thought experiment



Say we access to a sequence $u[t] \in [0,1]$: assume the following happens

- ightharpoonup At time t_0 , we have $u[t_0] = 1$



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Let's see what happens to the memory



Gating

Principle of Gating

At time t_0 , we can say

- We have $x[t_0] \propto$ "Julia" and compute $h[t_0] = f(w_1 x[t_0] + w_m h[t_0 1])$
 - $\rightarrow \tilde{h}[t_0]$ has fresh memory about "Julia"
 - \hookrightarrow Since $u[t_0]=1$, we update as $h[t_0]=1\times \tilde{h}[t_0]+0\times h[t_0-1]=\tilde{h}[t_0]$
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This repeats from $t_0 + 1$ till t, so at time t

- No matter what h[t], we have u[t] = 0
 - \downarrow We update as $h[t] = 0 \times \tilde{h}[t] + 1 \times h[t-1] = h[t-1] = \dots = \tilde{h}[t_0]$

u[t] gates the memory: it decides how much memory we should pass and forget

• We update $h[t] = u[t]\tilde{h}[t] + (1 - u[t])h[t - 1]$



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- + How does it help with vanishing gradient?
- Well! It is implicitly making skip connections through time

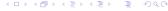


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Recall that with standard BPTT, we have for $i=t,t-1,\ldots,t_0$

$$\frac{\partial h[i]}{\partial w_{\rm m}} = \dot{f}(z[i]) \left(h[i-1] + w_{\rm m} \frac{\partial h[i-1]}{\partial w_{\rm m}} \right)$$



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But, now we skip multiple time slots, as we have for $i=t,t-1,\ldots,t_0$

$$\frac{\partial h[i]}{\partial w_{\rm m}} = \frac{\partial h[i-1]}{\partial w_{\rm m}}$$



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+ Sounds inspiring! But, how could you get u[t]?



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81/157

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Gate

Let $\mathbf{x}[t]$ be input and $\mathbf{h}[t-1]$ be last hidden state: a gate $\mathbf{\Gamma}[t]$ is computed as

$$\mathbf{\Gamma}[t] = \sigma \left(\mathbf{W}_{\Gamma, \text{in}} \mathbf{x}[t] + \mathbf{W}_{\Gamma, \text{m}} \mathbf{h}[t-1] + \mathbf{b}_{\Gamma} \right)$$

where $\sigma\left(\cdot\right)$ is sigmoid function, and $\mathbf{W}_{\Gamma,\mathrm{in}}$, $\mathbf{W}_{\Gamma,\mathrm{m}}$ and \mathbf{b}_{Γ} are learnable⁵

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- + Why do we use sigmoid function?
- Simply because it is between 0 and 1
- + What should we set the dimension of $\Gamma[t]$?
- Same as the variable (memory component) that we want to gate

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There are various gated architectures: we look into two of them

- Gated Recurrent Unit (GRU)
- 2 Long Short-Term Memory (LSTM)

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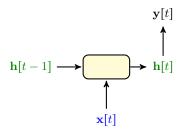
Before we start, let's recall their basic RNN counterpart

Basic RNN Counterpart

Say we set activation to $f\left(\cdot\right) \leadsto$ we usually set it to $\tanh\left(\cdot\right)$

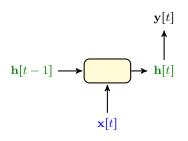
- **1** Start with an initial hidden state
 - \rightarrow we can learn $\mathbf{h}[0]$
- 2 Compute memory as $\mathbf{h}[t] = f(\mathbf{W}_1 \mathbf{x}[t] + \mathbf{W}_m \mathbf{h}[t-1])$
 - $\,\,\,\,\,\,\,\,\,\,$ we can learn ${f W}_1$ and ${f W}_m$
- f 3 Compute output ${f y}[t]=f_{
 m out}({f W}_2{f h}[t]) \leadsto f_{
 m out}$ and f could be different
 - $\,\,\,\downarrow\,\,\,$ we can learn ${f W}_2$

When we study a gated architecture: it is common to look at the hidden layer as a unit which takes some inputs and returns some outputs



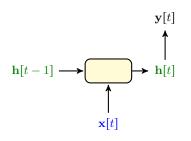
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We are mainly interested on this block: we want to know that given last state and new input

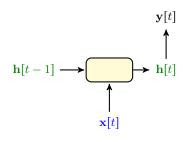
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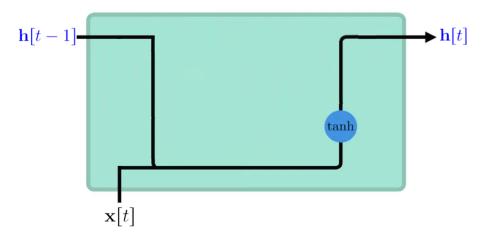
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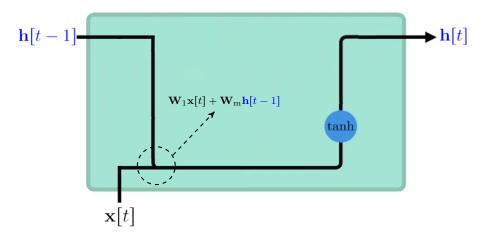
We are mainly interested on this block: we want to know that given last state and new input

- 1 How does this unit update hidden state?
- What components are passed to the next time interval?
 - \rightarrow Here, we have only $\mathbf{h}[t]$
 - But, we may have other components

Classical Diagram: Basic RNN



Classical Diagram: Basic RNN



Gated Recurrent Unit (GRU)

Say we set activation to $f(\cdot) \leadsto$ we usually set it to $\tanh(\cdot)$

- 1 Start with an initial hidden state
- 2 Compute update gate $\Gamma_{\rm u}[t] = \sigma \left(\mathbf{W}_{\rm u,in} \mathbf{x}[t] + \mathbf{W}_{\rm u,m} \mathbf{h}[t-1] \right)$
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- **4** Compute actual memory $\tilde{\mathbf{h}}[t] = f(\mathbf{W}_1 \mathbf{x}[t] + \mathbf{W}_m \Gamma_r[t] \odot \mathbf{h}[t-1])$



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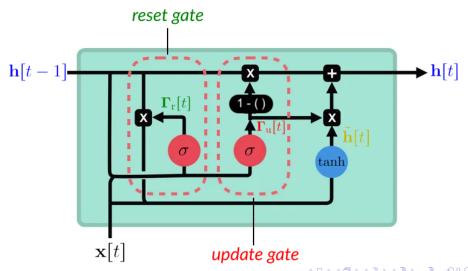
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- **6** Compute output $\mathbf{y}[t] = f_{\mathrm{out}}(\mathbf{W}_2\mathbf{h}[t]) \leadsto f_{\mathrm{out}}$ and f could be different

Or we could give $\mathbf{h}[t]$ to a new layer: for instance a new GRU whose input is $\mathbf{h}[t]$ and has its own state

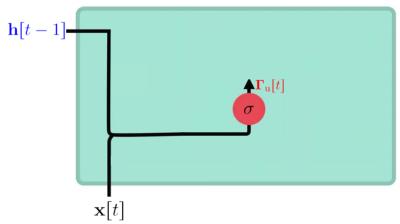
GRU

Practical Gated Architectures: GRU

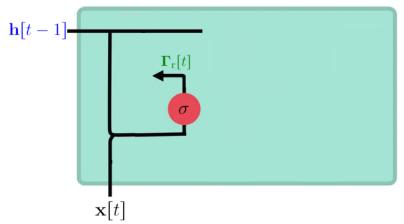
This is what's going on in a GRU cell



Compute update gate $\Gamma_{\rm u}[t] = \sigma \left(\mathbf{W}_{\rm u,in} \mathbf{x}[t] + \mathbf{W}_{\rm u,m} \mathbf{h}[t-1] \right)$

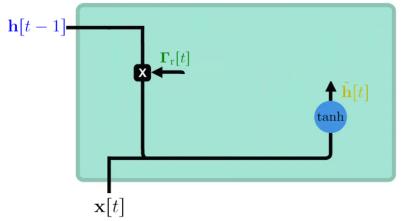


Compute reset gate $\Gamma_{\rm r}[t] = \sigma \left(\mathbf{W}_{\rm r,in} \mathbf{x}[t] + \mathbf{W}_{\rm r,m} \mathbf{h}[t-1] \right)$



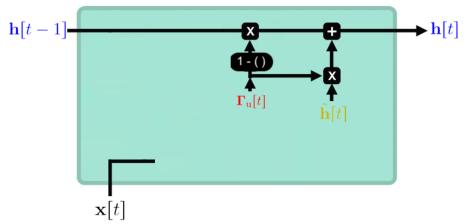


Compute actual memory $\tilde{\mathbf{h}}[t] = f(\mathbf{W}_1 \mathbf{x}[t] + \mathbf{W}_m \Gamma_r[t] \odot \mathbf{h}[t-1])$



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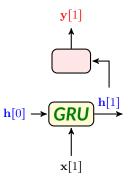
Update hidden state as $\mathbf{h}[t] = (1 - \mathbf{\Gamma}_{\mathbf{u}}[t]) \odot \mathbf{h}[t-1] + \mathbf{\Gamma}_{\mathbf{u}}[t] \odot \tilde{\mathbf{h}}[t]$



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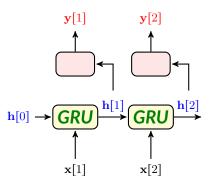
GRU

GRU: Forward Pass

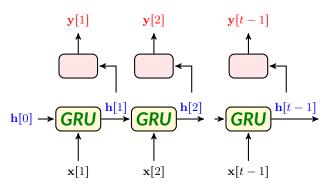


GRU

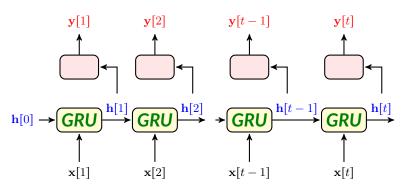
GRU: Forward Pass



GRU: Forward Pass



GRU: Forward Pass



GRU

GRU: Backward Pass

Say we finished forward pass at time t. We now want to find $\nabla_{\mathbf{W}} \hat{R}$ for some \mathbf{W} that is inside GRU, e.g., $\mathbf{W}_{\mathrm{u,m}}$:



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Say we finished forward pass at time t. We now want to find $\nabla_{\mathbf{W}} \hat{R}$ for some \mathbf{W} that is inside GRU, e.g., $\mathbf{W}_{\mathrm{u,m}}$: we start backpropagating from $\nabla_{\mathbf{y}[t]} \hat{R}$

$$\nabla_{\mathbf{W}} \hat{R} = \nabla_{\mathbf{y}[t]} \hat{R} \circ \nabla_{\mathbf{W}} \mathbf{y}[t]$$



Applied Deep Learning

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$$\nabla_{\mathbf{W}}\mathbf{h}[t] = \nabla_{\mathbf{\Gamma}_{\mathbf{u}}[t]}\mathbf{h}[t] \circ \nabla_{\mathbf{W}}\mathbf{\Gamma}_{\mathbf{u}}[t] + \nabla_{\mathbf{h}[t-1]}\mathbf{h}[t] \circ \nabla_{\mathbf{W}}\mathbf{h}[t-1]$$



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3 ...



Practical Gated Architectures: Long Short-Term Memory

Long Short-Term Memory (LSTM)

Say we set activation to $f\left(\cdot\right) \leadsto$ we usually set it to $\tanh\left(\cdot\right)$

- 1 Start with initial hidden state and cell state
- 2 Compute forget gate $\Gamma_f[t] = \sigma \left(\mathbf{W}_{f,in} \mathbf{x}[t] + \mathbf{W}_{f,m} \mathbf{h}[t-1] \right)$
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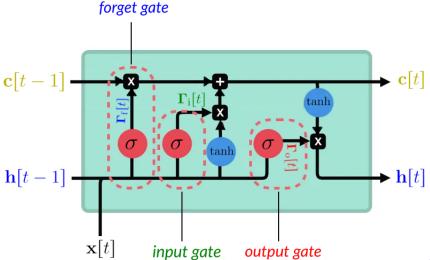
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- **7** Update hidden state as $\mathbf{h}[t] = \Gamma_{\mathbf{o}}[t] \odot f(\mathbf{c}[t])$
- 8 Compute output $\mathbf{y}[t] = f_{\text{out}}(\mathbf{W}_2\mathbf{h}[t]) \rightsquigarrow f_{\text{out}}$ and f could be different

Or we could give $\mathbf{h}[t]$ to a new layer: for instance a new LSTM whose input is h[t] and has its own hidden and cell states

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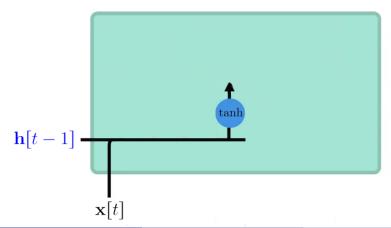
Practical Gated Architectures: LSTM

This is how inside an LSTM unit looks like



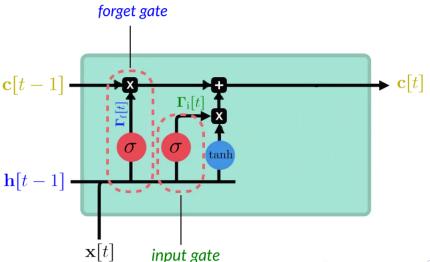
Practical Gated Architectures: LSTM

Actual cell state
$$\tilde{\mathbf{c}}[t] = f(\mathbf{W}_1 \mathbf{x}[t] + \mathbf{W}_m \mathbf{h}[t-1])$$



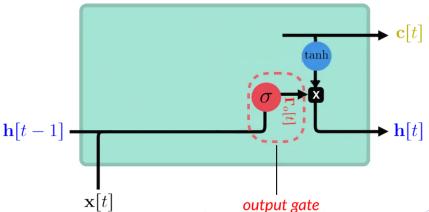
Practical Gated Architectures: LSTM

We use forget gate and update gate to update cell state



Practical Gated Architectures: LSTM

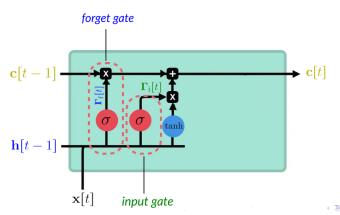
We use output gate to control fellow of memory to the hidden state



Practical Gated Architectures: LSTM

Intuitively, the gates in LSTM impact the flow of information as follows

- Forget gate controls how much we forget from last state
 - Arr Assume Arr Arr Arr Arr Arr 1 Arr 4 Arr Arr 4 Arr 4 Arr Arr 4 Arr 6 Arr 7 Arr 7 Arr 8 Arr 7 Arr 7 Arr 8 Arr 8 Arr 9 9 Arr 9 9 Arr 9 9 Arr 9 9 Arr 9 9

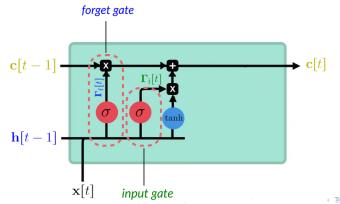


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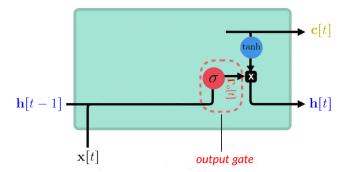
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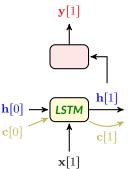
Practical Gated Architectures: LSTM

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LSTM: Forward Pass

Starting from initial hidden and cell state: LSTM passes forward as

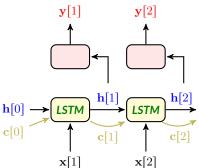


Pay Attention

Note that unlike other architectures, LSTM does not keep all memory inside hidden state but it carries it also in cell state. This state is only for memory and is not directly used by higher layers, e.g., output layer of the NN

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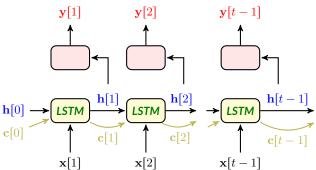
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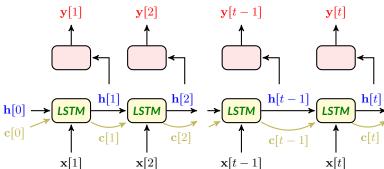
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Applied Deep Learning

Chapter 6: RNNs

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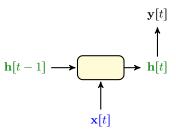
Suggestion

Try writing it once to see the impact of gates!



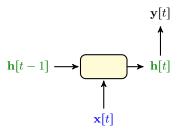
We have up to now considered unidirectional RNNs

we start from beginning of the sequence and move in one direction



We have up to now considered unidirectional RNNs

we start from beginning of the sequence and move in one direction



But, can't we learn from future input as well?

Future entries can have information about past: say our RNN wants to fill the empty field

... the color that many people assume is the color of sun ...

Obviously, future input in the sequence is helping in this example!

Future entries can have information about past: say our RNN wants to fill the empty field

... the color that many people assume is the color of sun ...

Obviously, future input in the sequence is helping in this example!

- + But, how can we get information from future?
- Well, we have the whole sequence: we could move once from left to right and once from right to left

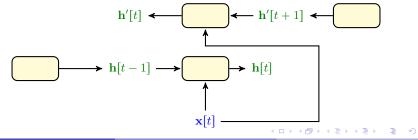
Bidirectional RNNs

Bidirectional RNNs

A bidirectional RNN (BRNN) consists of two RNNs

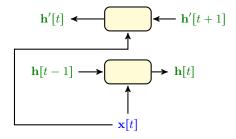
- one that starts with an initial hidden state at t=0 and computes $\mathbf{h}[t]$ from $\mathbf{h}[t-1]$ and $\mathbf{x}[t]$
- another that starts with an initial hidden state at t=T+1 and computes $\mathbf{h'}[t]$ from $\mathbf{h'}[t+1]$ and $\mathbf{x}[t]$

Output at time t is determined from merged version of the two hidden states



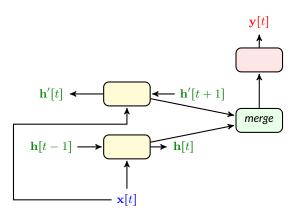
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Bidirectional RNNs



Applied Deep Learning

Bidirectional RNNs



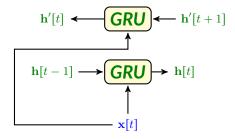
+ What exactly is this merge block?

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It gets the two states and returns a vector that matches output layer

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Bidirectional GRU

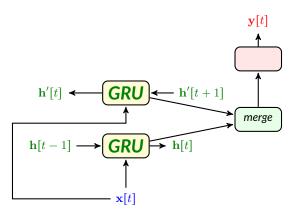


- + Should we use any RNN here?
- Sure! We may use GRU



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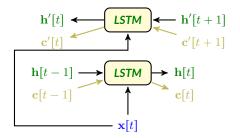
Bidirectional GRU



- + Should we use any RNN here?
- Sure! We may use GRU



Bidirectional LSTM

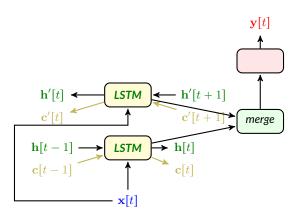


- + Should we use any RNN here?
- Sure! We may use LSTM



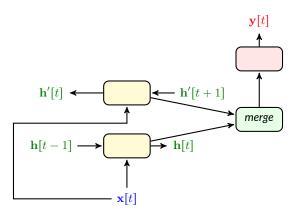
Applied Deep Learning

Bidirectional LSTM



- + Should we use any RNN here?
- Sure! We may use LSTM





- + Any suggestion for merging the hidden states?
- Sure! Let's see some code



Applied Deep Learning

In PyTorch, we can access a basic RNN in torch.nn module

torch.nn.RNN()

We can make it deep by simply choosing num_layers more than one and bidirectional by setting bidirectional to True.

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torch.nn.LSTM()
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```

In bidirectional case, we get access to both states. To merge them, we could

- add the two states
- average them
- concatenate them, i.e., $\mathbf{h}_{\mathrm{c}}[t] = (\mathbf{h}[t], \mathbf{h}'[t])$

or do any other operation that we find useful



Computing Loss: Challenge

We mentioned several times in this chapter that we assume

we can compute the loss between RNN's output sequence and label sequence

However, it is in general a challenge!

Computing Loss: Challenge

We mentioned several times in this chapter that we assume

we can compute the loss between RNN's output sequence and label sequence

However, it is in general a challenge!

- + Why is it a challenge? We did it easily in FNN and CNN chapters!
- Because the problem there was already properly segmented!
- + What do you mean by segmented?
- Let's break it down!

Let's consider a simple example: we have an image that includes a sequence of handwritten digits, e.g.,

- The sequence includes five digits
- Each digit is either 1, 2, 3, or 4



Let's consider a simple example: we have an image that includes a sequence of handwritten digits, e.g.,

- The sequence includes five digits
- Each digit is either 1, 2, 3, or 4

Our task is to recognize this sequence, i.e., return the five digits in correct order

- This is a classification task
- How can we do it? We use NNs
 - We train an NN over lots of images: we have lots of images of sequence of digits

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Let's say we are going to use a CNN



- We segment an input image into a sequence of five images
 - → These images are all as large as CNN's input size



To use a CNN, we need to specify our input size

- We segment an input image into a sequence of five images
 - → These images are all as large as CNN's input size

23 241 ~~ 2, 3, 2, 4, 1

• We label each image with its label, e.g., $\mathcal Q$ is labeled as 2

- We segment an input image into a sequence of five images
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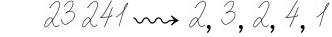
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- We label each image with its label, e.g., 2 is labeled as 2
- We give these five images to our CNN and get five outputs
 - → Assume we use softmax at the output layer
 - → For each image, we get a vector of size 4 as output
 - ightharpoonup Each entry represents probability of image being one of digits 1, 2, 3, and 4
- To compute loss, we compare each output with its corresponding label

$$\mathbf{v}[1] = \begin{bmatrix} 0\\1\\0\\0\\0 \end{bmatrix} \quad \mathbf{v}[2] = \begin{bmatrix} 0\\0\\1\\0 \end{bmatrix} \quad \mathbf{v}[3] = \begin{bmatrix} 0\\1\\0\\0 \end{bmatrix} \quad \mathbf{v}[4] = \begin{bmatrix} 0\\0\\0\\0\\1 \end{bmatrix} \quad \mathbf{v}[5] = \begin{bmatrix} 1\\0\\0\\0\\0 \end{bmatrix}$$

$$\begin{bmatrix} y_1[1]\\y_2[1]\\y_3[1]\\y_4[1] \end{bmatrix} = \mathbf{y}[1] \quad \mathbf{y}[2] \quad \mathbf{y}[3] \quad \mathbf{y}[4] \quad \mathbf{y}[5]$$

 $\downarrow y_j[t]$ is the probability of digit in time t being j

Here, we already have the data segmented into

a sequence that for each time step has a label

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a sequence that for each time step has a label

So, computing loss is easy as pie!

$$\hat{R} = \mathcal{L}(\mathbf{y}[1], \dots, \mathbf{y}[5], \mathbf{v}[1], \dots, \mathbf{v}[5])$$

$$= \sum_{t=1}^{5} \mathcal{L}(\mathbf{y}[t], \mathbf{v}[t]) = \sum_{t=1}^{5} \hat{R}[t]$$

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When we compute gradients, we note that only $\hat{R}[t]$ depends on $\mathbf{y}[t]$: so, for a given output at time t=i we can simply write

$$\nabla_{\mathbf{y}[i]} \hat{R} = \sum_{t=1}^{5} \nabla_{\mathbf{y}[i]} \hat{R}[t] = \nabla_{\mathbf{y}[i]} \hat{R}[i] = \nabla_{\mathbf{y}[i]} \mathcal{L}(\mathbf{y}[i], \mathbf{v}[i])$$

- + But is it practical to do segmentation by hand?
- No! This is why we built RNNs!



- + But is it practical to do segmentation by hand?
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With RNNs, we address this learning task as bellow

- We look at the complete image as a sequence of data
- We go over each frame separately

If we are extremely lucky; then, our segmentation looks like this

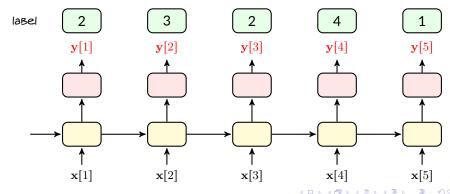
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and we have a label for each time step

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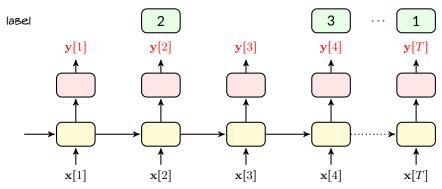


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But, that's too good to happen! Usually we have

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$$\longrightarrow \mathbf{x}[1], \mathbf{x}[2], \dots, \mathbf{x}[T]$$

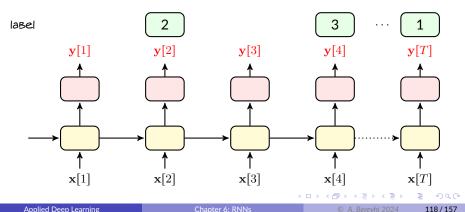
and we have a label in some time steps



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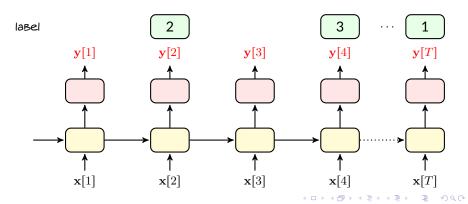
In this typical case, two questions seem non-trivial

1 Where should we put each label? \equiv Where should we read each label?



In this typical case, two questions seem non-trivial

- 1 Where should we put each label? \equiv Where should we read each label?
- 2 What should we do with non-labeled outputs, e.g., y[1]?



The key challenge in computing the loss is that we do not have necessarily one-to-one correspondence with sequence data



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Correspondence Problem

With sequence data, we could have a data-sequence of time T that is labeled by a sequence of size K < T where

no time index is specified for any label in the K-long label sequence

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Correspondence Problem

With sequence data, we could have a data-sequence of time T that is labeled by a sequence of size K < T where

no time index is specified for any label in the K-long label sequence

Correspondence problem exists pretty much in all practical sequence data

- In speech recognition, multiple time frames correspond to a single word
- In text recognition, multiple image frames correspond to a single letter
- ...

Correspondence Problem: Formulation

Let's formulate the problem clearly: Say we have

A sequence of data

$$\mathbf{x}[1:T] = \mathbf{x}[1], \dots, \mathbf{x}[T]$$

that is labeled with the sequence of K true labels

$$\mathbf{v}[1:K] = \mathbf{v}[1], \dots, \mathbf{v}[K]$$

where K and T can be different

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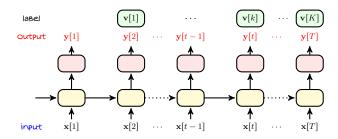
where K and T can be different

For this setting, we want to train an RNN with this data sequence: starting with an initial state, this RNN returns an output sequence

$$\mathbf{y}[1:T] = \mathbf{y}[1], \dots, \mathbf{y}[T]$$



Correspondence Problem: Formulation



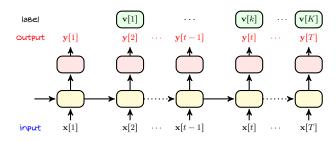
To be able to train this RNN, we need to

- **1** define a loss function that computes $\hat{R} = \mathcal{L}(\mathbf{y}[1:T], \mathbf{v}[1:K])$

$$\nabla_{\mathbf{y}[1]}\hat{R}, \dots, \nabla_{\mathbf{y}[T]}\hat{R}$$



Correspondence Problem: Formulation



To use this RNN after training, i.e., for inferring, we need to

- 2 know how to map outputs to predicted labels
 - $\,\,\,\,\,\,\,\,\,$ We need to extract K labels from $\mathbf{y}[1:T]$, i.e.,

$$\mathbf{y}[1], \dots, \mathbf{y}[T] \mapsto \hat{\mathbf{v}}[1], \dots, \hat{\mathbf{v}}[K]$$

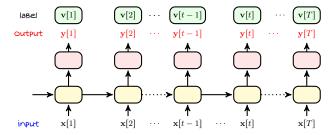
Let's look into different settings



Setting I: Perfectly Segmented

In some problems, we have our data perfectly segmented

• There is a separate label for each time step, i.e., K=T \rightarrow many-to-many type I and one-to-many



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 - lasel v[1] v[2] v[t-1] v[t] v[T] Output v[T] v[T]

Attention

We can always treat a non-existing input entry as an empty

igspace We are good as long as we have a label at each time t

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Setting I: Defining Loss

In such settings, we define the loss to be aggregated loss over time

$$\hat{R} = \sum_{t=1}^{T} \mathcal{L}\left(\mathbf{y}[t], \mathbf{v}[t]\right)$$

for some loss function $\mathcal{L}\left(\cdot,\cdot\right)$

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The gradients are then trivially computed

Gradient with respect to particular output $\mathbf{y}[t]$ is

$$\nabla_{\mathbf{y}[t]} \hat{R} = \nabla_{\mathbf{y}[t]} \hat{R}[t] = \nabla_{\mathbf{y}[t]} \mathcal{L}\left(\mathbf{y}[t], \mathbf{v}[t]\right)$$

Setting I: Inference

Inference in such setting is performed by one-to-one mapping: at time t, we predict based on $\mathbf{y}[t]$

$$\mathbf{y}[1] \mapsto \hat{\mathbf{v}}[1], \dots, \mathbf{y}[T] \mapsto \hat{\mathbf{v}}[T]$$

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For instance, assume $\mathbf{y}[t]$ is output of a softmax activation; then, we set

$$\hat{\mathbf{v}}[t] = \operatorname{argmax} \mathbf{y}[t]$$

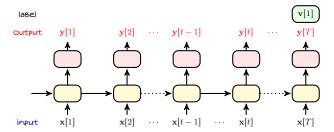
where argmax returns the index of the largest entry, e.g.,

$$\operatorname{argmax} \begin{bmatrix} 0.1\\0.7\\0.2\\0 \end{bmatrix} = 2$$

Setting II: Known Segments

In some problems, we have only one label for the whole sequence, i.e., K=1

- ↓ It corresponds to many-to-one type of problems
 - ☐ This can be that we have really only one label, e.g., content classification



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Setting II: Defining Loss for Dumb NN

A naive approach to define loss is to set it be the loss between last output and label, i.e.,

$$\hat{R} = \mathcal{L}(\mathbf{y}[T], \mathbf{v}[1])$$

Applied Deep Learning

Chapter 6: RNNs

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With this loss, gradient with respect to particular output $\mathbf{y}[t]$ is

$$\nabla_{\mathbf{y}[t]} \hat{R} = \begin{cases} \nabla_{\mathbf{y}[T]} \mathcal{L} \left(\mathbf{y}[T], \mathbf{v}[1] \right) & t = T \\ 0 & t \neq T \end{cases}$$

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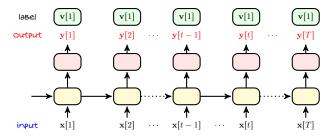
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- + But does it make sense to ignore all other outputs?
- Not at all! We are training a dumb NN that can respond only when it's over with the whole sequence!

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Setting II: Loss for Smarter Training

An extremely smart NN is the one who knows the label before the input speaks!

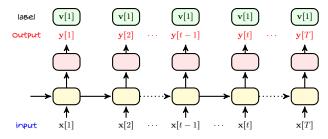


For this NN, the loss is

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For this NN, the loss is

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But, we should be careful! We should not expect NN to know everything from potentially irrelevant input!

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Setting II: Defining Proper Loss

A realistic approach is to define the loss via a weighted sum, i.e.,

$$\hat{R} = \sum_{t=1}^{T} w_t \mathcal{L}(\mathbf{y}[t], \mathbf{v}[1])$$

where w_t is the weight at time t

- initially w_t is small
- ullet it gradually increases up to its maximum w_T
 - $\,\,\,\,\,\,\,\,\,\,$ by time T the NN should know the label

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Setting II: Inference

Inference in such setting is performed by many-to-one mapping: we only predict based on $\mathbf{y}[1:T]$

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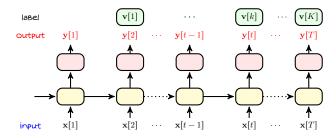
$$\tilde{\mathbf{v}}[t] = \operatorname{argmax} \mathbf{y}[t]$$

and then take a (potentially weighted) majority vote: $\hat{\mathbf{v}}[1]$ is the class that most often estimated with occurrence at each time being weighted by some weight

Setting III: Unknown Segments

Most common case is that we have a label sequence shorter than our data

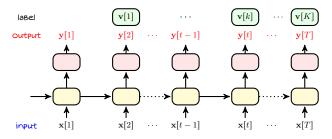
- □ Each label in this sequence is corresponding to a segment of input



Setting III: Unknown Segments

Most common case is that we have a label sequence shorter than our data

- □ Each label in this sequence is corresponding to a segment of input



Note that we are dealing with a sequence to sequence model: we want to learn

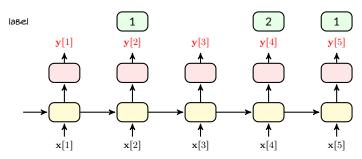
relation between sequence x[1:T] and sequence v[1:K]!

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Assume we have image 121 that is divided into a sequence of five pixel vectors

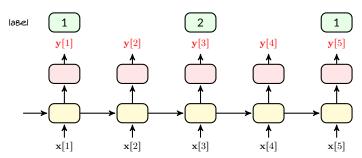
- Since it is a training data, it is labeled as 121
 - We do not know after which output we should expect RNN to know first, second or third digit!



Applied Deep Learning

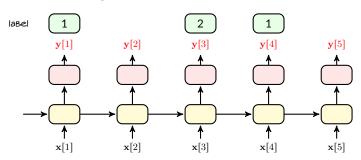
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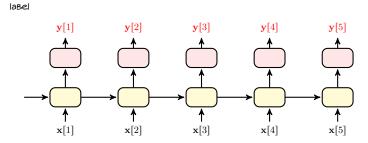
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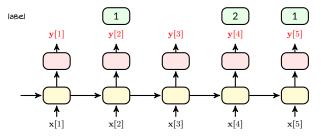


- + Sounds impossible!
- Only impossible is impossible! Let's carry on and see what we can do!

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Setting II: Genie-Defined Loss

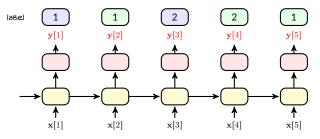
Assume a genie has told us end of each segment



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Setting II: Genie-Defined Loss

Assume a genie has told us end of each segment



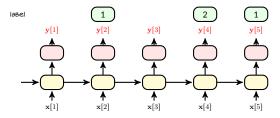
We can fill the empty labels with repetition, and then define the loss as

$$\hat{R} = \sum_{k=1}^{K} \sum_{t=i_{k-1}+1}^{i_k} w_t \mathcal{L}(\mathbf{y[t]}, \mathbf{v[k]})$$

where i_k is where label \mathbf{v}_k ends, e.g., in above diagram $i_1=2$

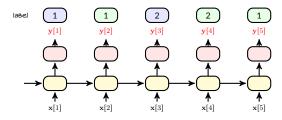
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Setting II: Defining Loss



We don't have the genie:

Setting II: Defining Loss



We don't have the genie: we could assume that i_k is something to learn!

$$\hat{R}(\mathbf{i}) = \sum_{k=1}^{K} \sum_{t=i_{k-1}+1}^{i_k} w_t \mathcal{L}(\mathbf{y}[t], \mathbf{v}[k])$$

where $\mathbf{i} = [0, i_1, \dots, i_K]$ is something we need to learn

Setting II: Optimal Segmentation

- + How could we learn i? Should we compute also $\nabla_i \hat{R}$?
- Well! You may try! But, obviously i_k is an <code>integer!</code>



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Optimal Segmentation

Optimal approach for finding i is to train the NN for all possible choice for i and then find the final training loss $\hat{R}(i)$. The optimal segmentation is then given by

$$\mathbf{i}^{\star} = \operatorname*{argmin}_{\mathbf{i}} \hat{R}\left(\mathbf{i}\right)$$

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$$\mathbf{i}^{\star} = \underset{\mathbf{i}}{\operatorname{argmin}} \hat{R}(\mathbf{i})$$

- + Is it computationally feasible?
- No! The number of possible choice for i grows exponentially with T! We need to go for sub-optimal approaches

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- + How is it exponentially large?
- Let's look at our example

In our last example, we should assign label sequence 121 to a sequence of length 5: each entry of output sequence in this case can be labeled by 1 (the first one), 2 or 1 (the last one). This means that we have 3 choices of label for each time interval; thus, the total number of possible segmentations is around 3^5 .

In general number of segmentations grows exponentially with T

- + But wait a moment! We have also counted the case of labeling all outputs with 1! This cannot be the case!
- This is right! It is in general much less than 3^5 but it's still exponential Let's see the exact possible segmentations!

We intend to compare each of $y[1], \dots, y[5]$ with a label

• We know that the label sequence is 121

 $\mathbf{y}[1]$ $\mathbf{y}[2]$ $\mathbf{y}[3]$ $\mathbf{y}[4]$ $\mathbf{y}[5]$

label

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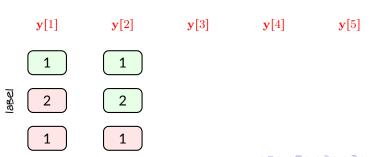
We intend to compare each of $y[1], \dots, y[5]$ with a label

- We know that the label sequence is 121
 - → First output is definitely in the first segment: it's label is definitely 1



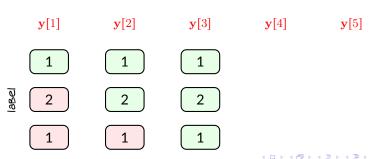
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- We know that the label sequence is 121
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 - Second output could be still in the first segment or in the second segment



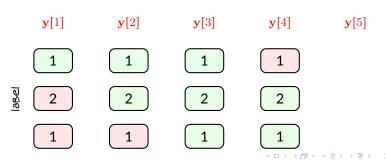
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 - ⇒ Second output could be still in the first segment or in the second segment
 - → Third output could be in the first, second, or third segment



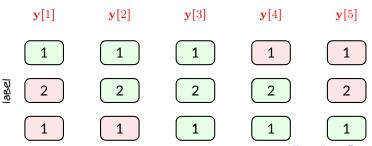
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 - Our labels should finish by the end of output sequence: fourth output cannot be in first segment



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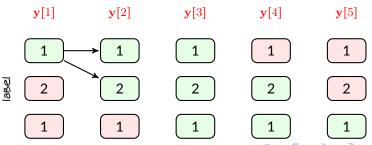
- We know that the label sequence is 121
 - → First output is definitely in the first segment: it's label is definitely 1
 - Second output could be still in the first segment or in the second segment
 - → Third output could be in the first, second, or third segment
 - Our labels should finish by the end of output sequence: fourth output cannot be in first segment
 - Last output could be only in the third segment



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We intend to compare each of $y[1], \dots, y[5]$ with a label

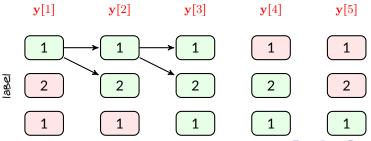
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 - ☐ Third output could be in the first, second, or third segment
 - Our labels should finish by the end of output sequence: fourth output cannot be in first segment
 - Last output could be only in the third segment



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We intend to compare each of $y[1], \dots, y[5]$ with a label

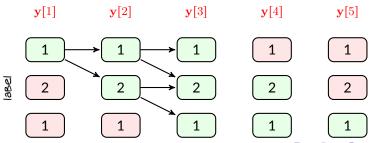
- We know that the label sequence is 121
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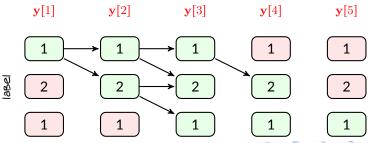
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Applied Deep Losephs Chapter (c DNN)s

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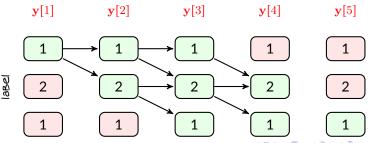
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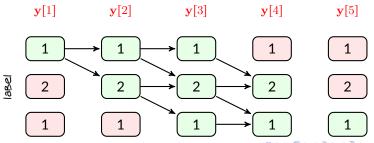
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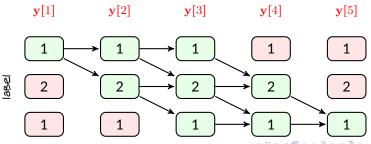
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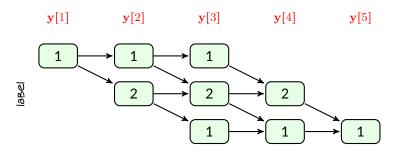
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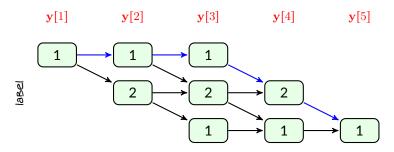
Setting II: Showing Segmentations on Graph



Though it's exponentially large: we see that each segmentation corresponds to one path on this graph

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Setting II: Showing Segmentations on Graph

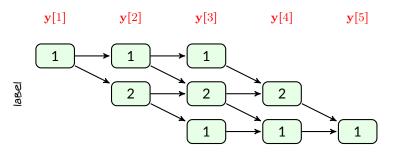


Though it's exponentially large: we see that each segmentation corresponds to one path on this graph

Blue path corresponds to $i_1 = 3$, $i_2 = 4$, and $i_3 = 5$, i.e., $\mathbf{i} = [0, 3, 4, 5]$

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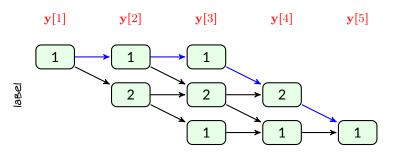
Applied Deep Learning



We can compute the loss for each segmentation directly on this graph: let's say that we have L different paths on the graph. For each path, we can write an expanded label sequence, e.g.,

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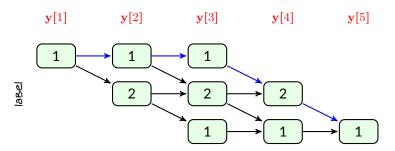
Applied Deep Learning



We can compute the loss for each segmentation directly on this graph: let's say that we have L different paths on the graph. For each path, we can write an expanded label sequence, e.g.,

Expanded label sequence of blue path is { 1, 1, 1, 2, 1}

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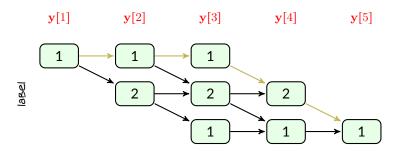
We can compute the loss for each segmentation directly on this graph: let's say that we have L different paths on the graph. For each path, we can write an expanded label sequence, e.g.,

Expanded label sequence of blue path is { 1, 1, 1, 2, 1}

This sequence is of length T and we show it for path ℓ with $\tilde{\mathbf{v}}_{\ell}[t]$

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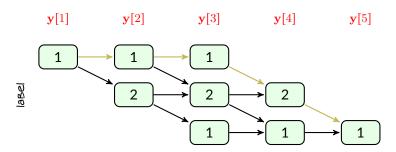
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For each path $\ell=1,\ldots,L$, the loss is computed by aggregating the losses between outputs and extended labels

$$\hat{R}_{\ell} = \sum_{t=1}^{T} w_t \mathcal{L}\left(\mathbf{y[t]}, \tilde{\mathbf{v}}_{\ell}[t]\right) =$$

Applied Deep Learning



For each path $\ell=1,\ldots,L$, the loss is computed by aggregating the losses between outputs and extended labels

$$\hat{R}_{\ell} = \sum_{t=1}^{T} w_{t} \mathcal{L}\left(\mathbf{y}[t], \tilde{\mathbf{v}}_{\ell}[t]\right) = \sum_{t=1}^{T} \hat{R}_{\ell}[t]$$

It again decomposes into sum of T terms with only one being function of $\mathbf{y}[t]$

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Setting II: Optimal Segmentation on Graph

Applied Deep Learning

We can represent the optimal segmentation on the graph as below

```
OptimalSegmentTraining():

1: Initiate with \hat{R} = +\infty and some random \ell^* = \emptyset

2: for \ell = 1, \dots, L do

3: Let the loss be \hat{R}_{\ell}

4: Train for sufficient epochs

5: if After training \hat{R}_{\ell} < \hat{R} then

6: \hat{R} \leftarrow \hat{R}_{\ell} and \ell^* \leftarrow \ell

7: end if

8: end for

9: Return learnable parameters and \ell^*
```

- + Say we could be over with this infeasible training! How do we use the trained RNN for inference?
- In this case, we have ℓ^* which gives us optimal segmentation: we infer label of each segment based on its corresponding outputs

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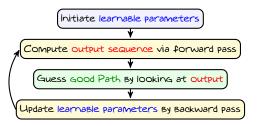
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Setting II: Maximum-Likelihood Segmentation

Since optimal segmentation is infeasible, people uses maximum-likelihood approach that is well-known in detection and coding theory

Maximum-Likelihood Segmentation

Start with an initial guess for optimal path on segmentation graph and do one step of training; then, improve the guess based on the outputs of next forward pass and go for next step of training



Let's look at its pseudo-code

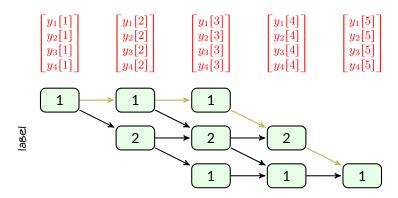
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Setting II: Maximum-Likelihood Segmentation

```
\begin{array}{lll} \operatorname{MaxLikelihoodTraining}(): \\ 1: & \text{ for } \operatorname{Iteration } i=1,\ldots,I \text{ do} \\ 2: & \operatorname{Pass } \operatorname{forward } \operatorname{through } \operatorname{time: Compute } \operatorname{output } \operatorname{sequence } \mathbf{y}[1:T] \\ 3: & \operatorname{Compute } p\left(\tilde{\mathbf{v}}_{\ell}[1:T]|\ell\right) \operatorname{for } \operatorname{each } \operatorname{path } \ell \operatorname{on } \operatorname{segmentation } \operatorname{graph} \\ 4: & \operatorname{Update } \ell^* = \operatorname{argmax}_{\ell} p\left(\tilde{\mathbf{v}}_{\ell}[1:T]|\ell\right) \\ 5: & \operatorname{Set } \operatorname{loss } \operatorname{to } \hat{R}_{\ell^*} \operatorname{and } \operatorname{backpropagate } \operatorname{over } \operatorname{RNN} \\ 6: & \operatorname{Update } \operatorname{learnable } \operatorname{parameters} \\ 7: & \operatorname{end } \operatorname{for} \\ 8: & \operatorname{Return } \operatorname{learnable } \operatorname{parameters } \operatorname{and } \ell^* \end{array}
```

- + Why we call it maximum likelihood?
- Because we guess path by maximizing the likelihood $p\left(ilde{\mathbf{v}}_{\ell}[1:T] | \ell
 ight)$
- + But how can find likelihood of a path?
- We can use output sequence y[1:T]

Setting II: Finding Likelihood on Segmentation Graph



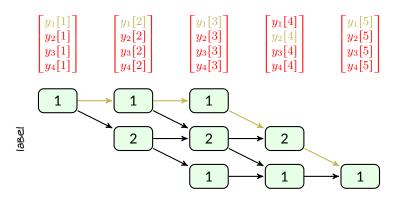
Assume that each label could be 1, 2, 3, or 4: at each time t the RNN returns a 4-dimensional vector whose entries are probability of each class

we can multiply the probabilities of classes on the path

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Setting II: Finding Likelihood on Segmentation Graph

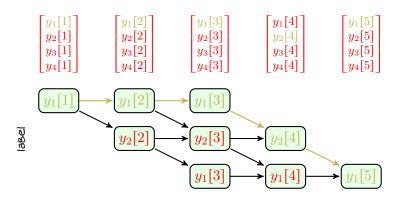


For instance, the yellow path has a likelihood

$$p\left(\tilde{\mathbf{v}}_{\ell}[1:T]|\ell\right) = \prod_{t=1}^{T} p\left(\tilde{\mathbf{v}}_{\ell}[t]|\ell\right) = y_{1}[1]y_{1}[2]y_{1}[3]y_{2}[4]y_{1}[5]$$

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Setting II: Finding Likelihood on Segmentation Graph



Or better to say: we just put output entries in graph and move on the path

$$p\left(\tilde{\mathbf{v}}_{\ell}[1:T]|\ell\right) = \prod_{t=1}^{T} y_{\tilde{v}_{\ell}[t]}[t] = y_{1}[1]y_{1}[2]y_{1}[3]y_{2}[4]y_{1}[5]$$

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Setting II: Maximum-Likelihood Segmentation

+ OK! We can find the likelihood, but how can we maximize it? It's again an exponentially large search!

$$\ell^* = \operatorname*{argmax}_{\ell} p\left(\tilde{\mathbf{v}}_{\ell}[1:T]|\ell\right)$$

Well! If we only need the maximum, it turns not to be exponential

Setting II: Maximum-Likelihood Segmentation

+ OK! We can find the likelihood, but how can we maximize it? It's again an exponentially large search!

$$\ell^* = \operatorname*{argmax} p\left(\tilde{\mathbf{v}}_{\ell}[1:T]|\ell\right)$$

- Well! If we only need the maximum, it turns not to be exponential

We can readily show that finding maximum likelihood on the graph is a dynamic programming problem and can be solved by the Viterbi algorithm

Maximum likelihood training can be implemented efficiently

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Applied Deep Learning

Setting II: Maximum-Likelihood Inference

```
MaxLikelihoodTraining():

1: for Iteration i=1,\ldots,I do

2: Pass forward through time: Compute output sequence \mathbf{y}[1:T]

3: Compute p\left(\tilde{\mathbf{v}}_{\ell}[1:T]|\ell\right) for each path \ell on segmentation graph

4: Update \ell^* = \operatorname{argmax}_{\ell} p\left(\tilde{\mathbf{v}}_{\ell}[1:T]|\ell\right)

5: Set loss to \hat{R}_{\ell^*} and backpropagate over RNN

6: Update learnable parameters

7: end for

8: Return learnable parameters and \ell^*
```

- + How can we use our RNN for inference after training via maximum likelihood segmentation?
- We have access to ℓ^* : we predict the label of each segment based on its corresponding outputs

Setting II: Connectionist Temporal Classification

It turns out that maximum-likelihood could stick to a bad local minimum, i.e.,

it quickly converges to a path ℓ^* that is much different from ℓ^*

- + Is there any solution to this?
- Yes! We can use connectionist temporal classification (CTC) loss

CTC Loss

Instead of searching for a best segmentation and then minimizing its loss, we learn directly from unsegmented data by minimizing the average loss over all possible segmentations, i.e., we define loss to be

$$\hat{R} = \mathbb{E}_{\ell} \left\{ \hat{R}_{\ell} \right\} = \sum_{\ell=1}^{L} p\left(\ell | \tilde{\mathbf{v}}_{\ell}[1:T] \right) \hat{R}_{\ell}$$

and train the RNN by finding learnable parameters that minimize this loss

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Setting II: CTC Loss

- + But, why should it be a better choice of loss?
- Because we are sure that optimal segmentation is contributing to our loss

$$\hat{R} = \sum_{\ell=1}^{L} p\left(\ell \middle| \tilde{\mathbf{v}}_{\ell}[1:T]\right) \hat{R}_{\ell} = p\left(\ell^{\star} \middle| \tilde{\mathbf{v}}_{\ell}[1:T]\right) \hat{R}_{\ell^{\star}} + \sum_{\ell \neq \ell^{\star}} p\left(\ell \middle| \tilde{\mathbf{v}}_{\ell}[1:T]\right) \hat{R}_{\ell}$$

- + Agreed! Now, how should we determine $p\left(\ell^{\star}|\tilde{\mathbf{v}}_{\ell}[1:T]\right)$?
- Just use the Bayes rule!
- + What about the expectation? It is at the end sum of exponentially large number of terms!
- We can again go on the graph and determine it via dynamic programming

Setting II: CTC Loss

The CTC loss can be written as

$$\hat{R} = \sum_{\ell=1}^{L} p\left(\ell \middle| \tilde{\mathbf{v}}_{\ell}[1:T]\right) \hat{R}_{\ell} = \sum_{\ell=1}^{L} p\left(\ell \middle| \tilde{\mathbf{v}}_{\ell}[1:T]\right) \sum_{t=1}^{T} w_{t} \mathcal{L}\left(\mathbf{y}[t], \tilde{\mathbf{v}}_{\ell}[t]\right)$$

$$= \sum_{t=1}^{T} w_{t} \sum_{\ell=1}^{L} p\left(\ell \middle| \tilde{\mathbf{v}}_{\ell}[1:T]\right) \mathcal{L}\left(\mathbf{y}[t], \tilde{\mathbf{v}}_{\ell}[t]\right) = \sum_{t=1}^{T} w_{t} \breve{R}[t]$$

$$\tilde{R}[t]$$

This has been shown that $\check{R}[t]$ can be recursively computed⁶:

by some approximation we are able to readily compute $\nabla_{\mathbf{y}[t]} \breve{R}[t]$

and we set
$$\nabla_{\mathbf{v}[t']} \breve{R}[t] \approx \mathbf{0}$$
 for $t' \neq t$



⁶Check out the original paper

Setting II: Training with CTC Loss

```
CTC_Training():

1: for iteration i = 1, ..., I do

2: Pass forward through time: Compute output sequence y[1:T]

3: Compute CTC loss \hat{R} and \nabla_{y[t]}\hat{R} by recursion

4: Backpropagate through time and update learnable parameters

5: end for

6: Return learnable parameters
```

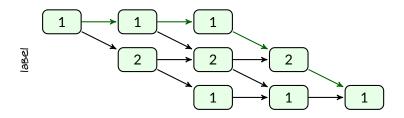
This looks like standard training loop now

the loss is only replaced with CTC loss

- + What about inference?
- Well! We should figure it out, since the training loop does not compute any segmentation path!

Let's get back to our simple example: assume that after training with CTC loss we give an image of handwritten 121 to the RNN

- RNN divides it into 5 frames and is able to track optimal segmentation

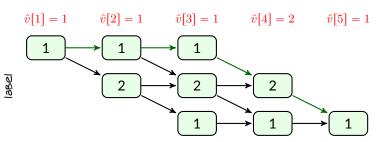


Applied Deep Learning

Let's get back to our simple example: assume that after training with CTC loss we give an image of handwritten 121 to the RNN

- RNN divides it into 5 frames and is able to track optimal segmentation

 - \downarrow The remaining frames belong to the second and third segments
- RNN infers from output sequence $\hat{v}[1:5]$ but does not return optimal path



Applied Deep Learning

We can conclude from $\hat{v}[1:5]$ that the sequence is $\{1,2,1\}$ if we are sure the sequence has no repetition

Label Encoding and Decoding

CTC uses this fact and constructs following encoding and decoding method: it introduces a new label called "blank:-" which does not belong to set of classes

We can conclude from $\hat{v}[1:5]$ that the sequence is $\{1,2,1\}$ if we are sure the sequence has no repetition

Label Encoding and Decoding

CTC uses this fact and constructs following encoding and decoding method: it introduces a new label called "blank:-" which does not belong to set of classes

- While training, it adds blank between any two repetitions
 - \rightarrow For instance, we encode 112 \mapsto 1-12, or 111 \mapsto 1-1-1

We can conclude from $\hat{v}[1:5]$ that the sequence is $\{1,2,1\}$ if we are sure the sequence has no repetition

Label Encoding and Decoding

CTC uses this fact and constructs following encoding and decoding method: it introduces a new label called "blank:-" which does not belong to set of classes

- While training, it adds blank between any two repetitions
 - \rightarrow For instance, we encode 112 \mapsto 1-12, or 111 \mapsto 1-1-1
- For inference, it removes any repetition in inferred sequence $\hat{v}[1:T]$ and then drops blanks

Applied Deep Learning

Setting II: Training and Inference with CTC

```
CTC_Training():

1: for iteration i = 1, ..., I do

2: Add blanks to the label sequences with repetition

3: Pass forward through time: Compute output sequence \mathbf{y}[1:T]

4: Compute CTC loss \hat{R} and \nabla_{\mathbf{y}[t]}\hat{R} by recursion

5: Backpropagate through time and update learnable parameters

6: end for

7: Return learnable parameters
```

Setting II: Training and Inference with CTC

```
CTC_Training():
1: for iteration i = 1, \ldots, I do
      Add blanks to the label sequences with repetition
      Pass forward through time: Compute output sequence y[1:T]
      Compute CTC loss \hat{R} and \nabla_{\mathbf{v}[t]}\hat{R} by recursion
4:
5:
      Backpropagate through time and update learnable parameters
6. end for
7: Return learnable parameters
```

```
CTC Inference():
1: Pass forward through time the input and compute output y[1:T]
2: Infere encoded sequence \hat{v}[1:T] from y[1:T]
3: Remove repetitions from \hat{v}[1:T] \mapsto \hat{v}[1:T']
4: Remove blanks from \hat{v}[1:T'] \mapsto \hat{v}[1:K]
5: Return \hat{v}[1:K]
```

In PyTorch: CTC Loss

We can access CTC loss in torch.nn module as

```
torcn.nn.CTCLoss()
```

Few notes about CTC loss implementation

- We need to specify the index of blank label
 - It should be out of our set of classes
 - \rightarrow By default, it is set to blank = 0
- When we define our model, we should always take blank label into account
- PyTorch considers cross-entropy loss function, i.e., $\mathcal{L}\left(\mathbf{y}, \tilde{\mathbf{v}}\right) = \mathrm{CE}\left(\mathbf{y}, \tilde{\mathbf{v}}\right)$
- As input to CTC loss: y should be logarithm of probabilities