Building your Recurrent Neural Network - Step by Step

Welcome to Course 5's first assignment! In this assignment, you will implement your first Recurrent Neural Network in numpy.

Recurrent Neural Networks (RNN) are very effective for Natural Language Processing and other sequence tasks because they have "memory". They can read inputs $x^{\langle t \rangle}$ (such as words) one at a time, and remember some information/context through the hidden layer activations that get passed from one time-step to the next. This allows a uni-directional RNN to take information from the past to process later inputs. A bidirection RNN can take context from both the past and the future.

Notation:

- ullet Superscript [l] denotes an object associated with the l^{th} layer.
 - ullet Example: $a^{[4]}$ is the 4^{th} layer activation. $W^{[5]}$ and $b^{[5]}$ are the 5^{th} layer parameters.
- ullet Superscript (i) denotes an object associated with the i^{th} example.
 - ullet Example: $x^{(i)}$ is the i^{th} training example input.
- ullet Superscript $\langle t
 angle$ denotes an object at the t^{th} time-step.
 - Example: $x^{\langle t \rangle}$ is the input x at the t^{th} time-step. $x^{(i)\langle t \rangle}$ is the input at the t^{th} timestep of example i.
- ullet Lowerscript i denotes the i^{th} entry of a vector.
 - lacksquare Example: $a_i^{[l]}$ denotes the i^{th} entry of the activations in layer l.

We assume that you are already familiar with numpy and/or have completed the previous courses of the specialization. Let's get started!

Let's first import all the packages that you will need during this assignment.

```
In [1]: import numpy as np
from rnn_utils import *
```

1 - Forward propagation for the basic Recurrent Neural Network

Later this week, you will generate music using an RNN. The basic RNN that you will implement has the structure below. In this example, $T_x=T_y$.

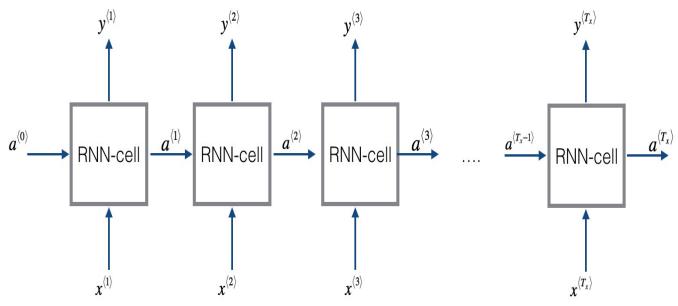


Figure 1: Basic RNN model

Here's how you can implement an RNN:

Steps:

- 1. Implement the calculations needed for one time-step of the RNN.
- 2. Implement a loop over T_x time-steps in order to process all the inputs, one at a time.

Let's go!

1.1 - RNN cell

A Recurrent neural network can be seen as the repetition of a single cell. You are first going to implement the computations for a single time-step. The following figure describes the operations for a single time-step of an RNN cell.

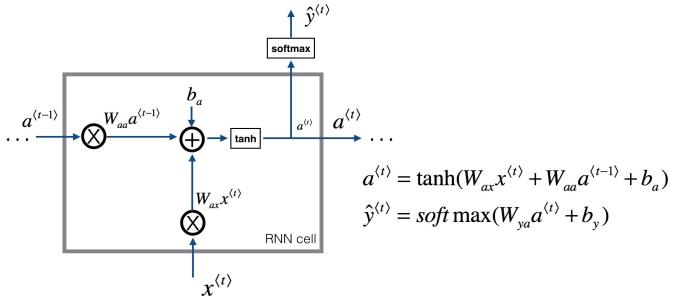


Figure 2: Basic RNN cell. Takes as input $x^{\langle t \rangle}$ (current input) and $a^{\langle t-1 \rangle}$ (previous hidden state containing information from the past), and outputs $a^{\langle t \rangle}$ which is given to the next RNN cell and also used to predict $y^{\langle t \rangle}$

Exercise: Implement the RNN-cell described in Figure (2).

Instructions:

- 1. Compute the hidden state with tanh activation: $a^{\langle t \rangle} = anh(W_{aa}a^{\langle t-1 \rangle} + W_{ax}x^{\langle t \rangle} + b_a)$.
- 2. Using your new hidden state $a^{\langle t \rangle}$, compute the prediction $\hat{y}^{\langle t \rangle} = softmax(W_{ya}a^{\langle t \rangle} + b_y)$. We provided you a function: softmax.
- 3. Store $(a^{\langle t \rangle}, a^{\langle t-1 \rangle}, x^{\langle t \rangle}, parameters)$ in cache
- 4. Return $a^{\langle t
 angle}$, $y^{\langle t
 angle}$ and cache

We will vectorize over m examples. Thus, $x^{\langle t \rangle}$ will have dimension (n_x, m) , and $a^{\langle t \rangle}$ will have dimension (n_a, m) .

```
In [2]: # GRADED FUNCTION: rnn cell forward
        def rnn cell forward(xt, a prev, parameters):
            Implements a single forward step of the RNN-cell as described in Fi
        gure (2)
            Arguments:
            xt -- your input data at timestep "t", numpy array of shape (n x, m
        ).
            a prev -- Hidden state at timestep "t-1", numpy array of shape (n a
        , m)
            parameters -- python dictionary containing:
                                Wax -- Weight matrix multiplying the input, num
        py array of shape (n a, n x)
                                 Waa -- Weight matrix multiplying the hidden sta
        te, numpy array of shape (n a, n a)
                                Wya -- Weight matrix relating the hidden-state
        to the output, numpy array of shape (n y, n a)
                                ba -- Bias, numpy array of shape (n a, 1)
                                by -- Bias relating the hidden-state to the out
        put, numpy array of shape (n y, 1)
            Returns:
            a next -- next hidden state, of shape (n a, m)
            yt pred -- prediction at timestep "t", numpy array of shape (n y, m
            cache -- tuple of values needed for the backward pass, contains (a
        next, a prev, xt, parameters)
            11 11 11
            # Retrieve parameters from "parameters"
            Wax = parameters["Wax"]
            Waa = parameters["Waa"]
            Wya = parameters["Wya"]
            ba = parameters["ba"]
            by = parameters["by"]
            ### START CODE HERE ### (≈2 lines)
            # compute next activation state using the formula given above
            a next = np.tanh(np.dot(Wax, xt) + np.dot(Waa, a prev) + ba)
            # compute output of the current cell using the formula given above
            yt pred = softmax(np.dot(Wya, a next) + by)
            ### END CODE HERE ###
            # store values you need for backward propagation in cache
            cache = (a next, a prev, xt, parameters)
            return a next, yt pred, cache
```

```
In [4]: np.random.seed(1)
       xt = np.random.randn(3,10)
       a prev = np.random.randn(5,10)
       Waa = np.random.randn(5,5)
       Wax = np.random.randn(5,3)
       Wya = np.random.randn(2,5)
       ba = np.random.randn(5,1)
       by = np.random.randn(2,1)
       parameters = {"Waa": Waa, "Wax": Wax, "Wya": Wya, "ba": ba, "by": by}
       a next, yt pred, cache = rnn cell forward(xt, a prev, parameters)
       print("a next[4] = ", a next[4])
       print("a next.shape = ", a next.shape)
       print("yt pred[1] =", yt pred[1])
       print("yt pred.shape = ", yt pred.shape)
       a \text{ next}[4] = [0.59584544 \ 0.18141802 \ 0.61311866 \ 0.99808218 \ 0.8501
       6201 0.99980978
        -0.18887155 0.99815551 0.6531151 0.82872037]
       a next.shape = (5, 10)
       8387 0.88945212
         0.36920224 0.9966312 0.9982559 0.17746526]
       yt pred.shape = (2, 10)
```

a_next[4]:	[0.59584544 0.18141802 0.61311866 0.99808218 0.85016201 0.99980978 -0.18887155 0.99815551 0.6531151 0.82872037]
a_next.shape:	(5, 10)
yt[1]: [0.9888161 0.01682021 0.21140899 0.36817467 0.98988387 0.88945212 0.368	
yt.shape:	(2, 10)

1.2 - RNN forward pass

You can see an RNN as the repetition of the cell you've just built. If your input sequence of data is carried over 10 time steps, then you will copy the RNN cell 10 times. Each cell takes as input the hidden state from the previous cell $(a^{\langle t-1 \rangle})$ and the current time-step's input data $(x^{\langle t \rangle})$. It outputs a hidden state $(a^{\langle t \rangle})$ and a prediction $(y^{\langle t \rangle})$ for this time-step.

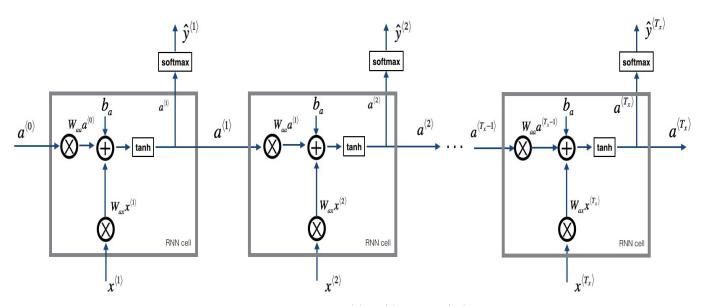


Figure 3: Basic RNN. The input sequence $x=(x^{\langle 1 \rangle},x^{\langle 2 \rangle},\ldots,x^{\langle T_x \rangle})$ is carried over T_x time steps. The network outputs $y=(y^{\langle 1 \rangle},y^{\langle 2 \rangle},\ldots,y^{\langle T_x \rangle})$.

Exercise: Code the forward propagation of the RNN described in Figure (3).

Instructions:

- 1. Create a vector of zeros (a) that will store all the hidden states computed by the RNN.
- 2. Initialize the "next" hidden state as a_0 (initial hidden state).
- 3. Start looping over each time step, your incremental index is t:
 - Update the "next" hidden state and the cache by running rnn cell forward
 - Store the "next" hidden state in a (t^{th} position)
 - Store the prediction in y
 - · Add the cache to the list of caches
- 4. Return a, y and caches

```
In [5]: # GRADED FUNCTION: rnn forward
        def rnn forward(x, a0, parameters):
            Implement the forward propagation of the recurrent neural network d
        escribed in Figure (3).
            Arguments:
            x -- Input data for every time-step, of shape (n x, m, T x).
            a0 -- Initial hidden state, of shape (n a, m)
            parameters -- python dictionary containing:
                                 Waa -- Weight matrix multiplying the hidden sta
        te, numpy array of shape (n a, n a)
                                 Wax -- Weight matrix multiplying the input, num
        py array of shape (n a, n x)
                                 Wya -- Weight matrix relating the hidden-state
        to the output, numpy array of shape (n y, n a)
                                 ba -- Bias numpy array of shape (n a, 1)
                                 by -- Bias relating the hidden-state to the out
        put, numpy array of shape (n y, 1)
            Returns:
            a -- Hidden states for every time-step, numpy array of shape (n a,
        m, T x)
            y pred -- Predictions for every time-step, numpy array of shape (n
        y, m, T \times 1
            caches -- tuple of values needed for the backward pass, contains (1
        ist of caches, x)
             11 11 11
            # Initialize "caches" which will contain the list of all caches
            caches = []
            # Retrieve dimensions from shapes of x and Wy
            n x, m, T x = x.shape
            n y, n a = parameters["Wya"].shape
            ### START CODE HERE ###
            # initialize "a" and "y" with zeros (≈2 lines)
            a = np.zeros((n a, m, T x))
            y \text{ pred} = \text{np.zeros}((n y, m, T x))
            # Initialize a next (≈1 line)
            a next = a0
            # loop over all time-steps
            for t in range(T \times):
                # Update next hidden state, compute the prediction, get the cac
        he (≈1 line)
                a next, yt pred, cache = rnn cell forward(x[:,:,t], a next, par
        ameters)
                # Save the value of the new "next" hidden state in a (≈1 line)
                a[:,:,t] = a next
                # Save the value of the prediction in y (≈1 line)
                y pred[:,:,t] = yt pred
```

```
In [6]: np.random.seed(1)
       x = np.random.randn(3, 10, 4)
       a0 = np.random.randn(5,10)
       Waa = np.random.randn(5,5)
       Wax = np.random.randn(5,3)
       Wya = np.random.randn(2,5)
       ba = np.random.randn(5,1)
       by = np.random.randn(2,1)
        parameters = {"Waa": Waa, "Wax": Wax, "Wya": Wya, "ba": ba, "by": by}
       a, y pred, caches = rnn forward(x, a0, parameters)
       print("a[4][1] = ", a[4][1])
       print("a.shape = ", a.shape)
       print("y pred[1][3] =", y pred[1][3])
       print("y_pred.shape = ", y_pred.shape)
       print("caches[1][1][3] =", caches[1][1][3])
       print("len(caches) = ", len(caches))
       a[4][1] = [-0.99999375 \ 0.77911235 \ -0.99861469 \ -0.99833267]
       a.shape = (5, 10, 4)
       y \text{ pred.shape} = (2, 10, 4)
       caches[1][1][3] = [-1.1425182 -0.34934272 -0.20889423 0.58662319]
       len(caches) = 2
```

a[4][1]:	[-0.99999375 0.77911235 -0.99861469 -0.99833267]
a.shape:	(5, 10, 4)
y[1][3]:	[0.79560373 0.86224861 0.11118257 0.81515947]
y.shape:	(2, 10, 4)
cache[1][1][3]:	[-1.1425182 -0.34934272 -0.20889423 0.58662319]
len(cache):	2

Congratulations! You've successfully built the forward propagation of a recurrent neural network from scratch. This will work well enough for some applications, but it suffers from vanishing gradient problems. So it works best when each output $y^{\langle t \rangle}$ can be estimated using mainly "local" context (meaning information from inputs $x^{\langle t' \rangle}$ where t' is not too far from t).

In the next part, you will build a more complex LSTM model, which is better at addressing vanishing gradients. The LSTM will be better able to remember a piece of information and keep it saved for many timesteps.

2 - Long Short-Term Memory (LSTM) network

This following figure shows the operations of an LSTM-cell.

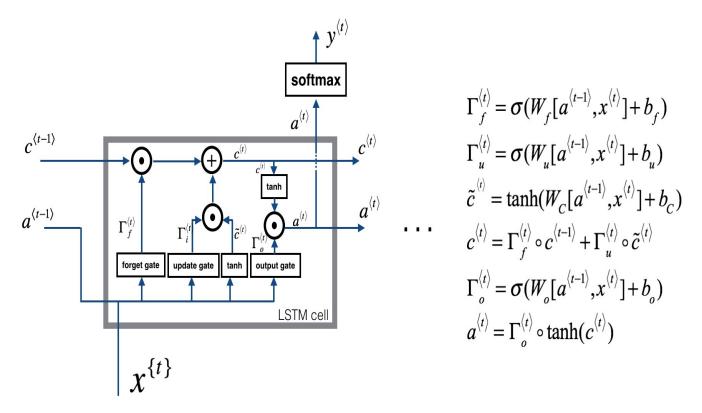


Figure 4: LSTM-cell. This tracks and updates a "cell state" or memory variable $c^{\langle t \rangle}$ at every time-step, which can be different from $a^{\langle t \rangle}$.

Similar to the RNN example above, you will start by implementing the LSTM cell for a single time-step. Then you can iteratively call it from inside a for-loop to have it process an input with T_x time-steps.

About the gates

- Forget gate

For the sake of this illustration, lets assume we are reading words in a piece of text, and want use an LSTM to keep track of grammatical structures, such as whether the subject is singular or plural. If the subject changes from a singular word to a plural word, we need to find a way to get rid of our previously stored memory value of the singular/plural state. In an LSTM, the forget gate lets us do this:

$$\Gamma_f^{\langle t \rangle} = \sigma(W_f[a^{\langle t-1 \rangle}, x^{\langle t \rangle}] + b_f)$$
 (1)

Here, W_f are weights that govern the forget gate's behavior. We concatenate $[a^{\langle t-1\rangle},x^{\langle t\rangle}]$ and multiply by W_f . The equation above results in a vector $\Gamma_f^{\langle t\rangle}$ with values between 0 and 1. This forget gate vector will be multiplied element-wise by the previous cell state $c^{\langle t-1\rangle}$. So if one of the values of $\Gamma_f^{\langle t\rangle}$ is 0 (or close to 0) then it means that the LSTM should remove that piece of information (e.g. the singular subject) in the corresponding component of $c^{\langle t-1\rangle}$. If one of the values is 1, then it will keep the information.

-- - . .

2.1 - LSTM cell

Exercise: Implement the LSTM cell described in the Figure (3).

Instructions:

- 1. Concatenate $a^{\langle t-1
 angle}$ and $x^{\langle t
 angle}$ in a single matrix: $concat=\left[egin{array}{c}a^{\langle t-1
 angle}\\x^{\langle t
 angle}\end{array}
 ight]$
- 2. Compute all the formulas 1-6. You can use sigmoid() (provided) and np.tanh().
- 3. Compute the prediction $y^{\langle t \rangle}$. You can use $\mathtt{softmax}$ () (provided).

```
In [7]: # GRADED FUNCTION: lstm cell forward
        def lstm cell forward(xt, a prev, c prev, parameters):
            Implement a single forward step of the LSTM-cell as described in Fi
        gure (4)
            Arguments:
            xt -- your input data at timestep "t", numpy array of shape (n x, m
        ).
            a prev -- Hidden state at timestep "t-1", numpy array of shape (n a
        , m)
            c prev -- Memory state at timestep "t-1", numpy array of shape (n a
        , m)
            parameters -- python dictionary containing:
                                 Wf -- Weight matrix of the forget gate, numpy a
        rray of shape (n \ a, n \ a + n \ x)
                                 bf -- Bias of the forget gate, numpy array of s
        hape (n a, 1)
                                 Wi -- Weight matrix of the update gate, numpy a
        rray of shape (n \ a, n \ a + n \ x)
                                 bi -- Bias of the update gate, numpy array of s
        hape (n a, 1)
                                 Wc -- Weight matrix of the first "tanh", numpy
        array of shape (n \ a, n \ a + n \ x)
                                 bc -- Bias of the first "tanh", numpy array of
        shape (n a, 1)
                                 Wo -- Weight matrix of the output gate, numpy a
        rray of shape (n \ a, n \ a + n \ x)
                                 bo -- Bias of the output gate, numpy array of
        shape (n a, 1)
                                 Wy -- Weight matrix relating the hidden-state t
        o the output, numpy array of shape (n y, n a)
                                 by -- Bias relating the hidden-state to the out
        put, numpy array of shape (n y, 1)
            Returns:
            a next -- next hidden state, of shape (n a, m)
            c next -- next memory state, of shape (n a, m)
            yt pred -- prediction at timestep "t", numpy array of shape (n y, m
            cache -- tuple of values needed for the backward pass, contains (a
        next, c next, a prev, c prev, xt, parameters)
            Note: ft/it/ot stand for the forget/update/output gates, cct stands
        for the candidate value (c tilde),
                  c stands for the memory value
            ,, ,, ,,
            # Retrieve parameters from "parameters"
            Wf = parameters["Wf"]
            bf = parameters["bf"]
            Wi = parameters["Wi"]
            bi = parameters["bi"]
            Wc = parameters["Wc"]
            bc = parameters["bc"]
```

```
In [8]: | np.random.seed(1)
        xt = np.random.randn(3,10)
        a prev = np.random.randn(5,10)
        c prev = np.random.randn(5,10)
        Wf = np.random.randn(5, 5+3)
        bf = np.random.randn(5,1)
        Wi = np.random.randn(5, 5+3)
        bi = np.random.randn(5,1)
        Wo = np.random.randn(5, 5+3)
        bo = np.random.randn(5,1)
        Wc = np.random.randn(5, 5+3)
        bc = np.random.randn(5,1)
        Wy = np.random.randn(2,5)
        by = np.random.randn(2,1)
        parameters = {"Wf": Wf, "Wi": Wi, "Wo": Wo, "Wc": Wc, "Wy": Wy, "bf": b
        f, "bi": bi, "bo": bo, "bc": bc, "by": by}
        a next, c next, yt, cache = lstm cell forward(xt, a prev, c prev, param
        eters)
        print("a_next[4] = ", a next[4])
        print("a next.shape = ", c next.shape)
        print("c next[2] = ", c next[2])
        print("c next.shape = ", c_next.shape)
        print("yt[1] =", yt[1])
        print("yt.shape = ", yt.shape)
        print("cache[1][3] =", cache[1][3])
        print("len(cache) = ", len(cache))
        a \text{ next}[4] = [-0.66408471 \ 0.0036921 \ 0.02088357 \ 0.22834167 \ -0.8557
        5339 0.00138482
          0.76566531 0.34631421 -0.00215674 0.43827275]
        a next.shape = (5, 10)
        c next[2] = [0.63267805 1.00570849 0.35504474 0.20690913 -1.6456]
        6718 0.11832942
          0.76449811 - 0.0981561 - 0.74348425 - 0.26810932
        c next.shape = (5, 10)
        yt[1] = [0.79913913 0.15986619 0.22412122 0.15606108 0.97057211
        0.31146381
          0.00943007 0.12666353 0.39380172 0.07828381]
        yt.shape = (2, 10)
        cache[1][3] = [-0.16263996   1.03729328   0.72938082   -0.54101719   0.027
        52074 -0.30821874
          0.07651101 -1.03752894 1.41219977 -0.376474221
        len(cache) = 10
```

a_next[4]:	[-0.66408471 0.0036921 0.02088357 0.22834167 -0.85575339 0.00138482 0.76566531 0.34631421 -0.00215674 0.43827275]
a_next.shape:	(5, 10)
c_next[2]:	[0.63267805 1.00570849 0.35504474 0.20690913 -1.64566718 0.11832942 0.76449811 -0.0981561 -0.74348425 -0.26810932]
c_next.shape:	(5, 10)
yt[1]:	[0.79913913 0.15986619 0.22412122 0.15606108 0.97057211 0.31146381 0.00943007 0.12666353 0.39380172 0.07828381]
yt.shape:	(2, 10)
cache[1][3]:	[-0.16263996 1.03729328 0.72938082 -0.54101719 0.02752074 -0.30821874 0.07651101 -1.03752894 1.41219977 -0.37647422]
len(cache):	10

2.2 - Forward pass for LSTM

Now that you have implemented one step of an LSTM, you can now iterate this over this using a for-loop to process a sequence of T_x inputs.

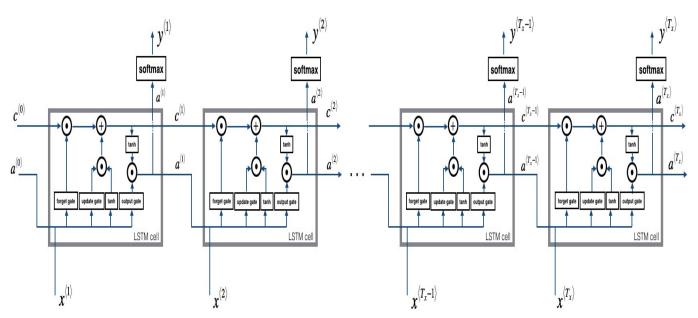


Figure 4: LSTM over multiple time-steps.

Exercise: Implement $lstm_forward()$ to run an LSTM over T_x time-steps.

Note: $c^{\langle 0 \rangle}$ is initialized with zeros.

```
In [9]: # GRADED FUNCTION: lstm forward
        def lstm forward(x, a0, parameters):
            Implement the forward propagation of the recurrent neural network u
        sing an LSTM-cell described in Figure (3).
            Arguments:
            x -- Input data for every time-step, of shape (n x, m, T x).
            a0 -- Initial hidden state, of shape (n a, m)
            parameters -- python dictionary containing:
                                 Wf -- Weight matrix of the forget gate, numpy a
        rray of shape (n \ a, n \ a + n \ x)
                                 bf -- Bias of the forget gate, numpy array of s
        hape (n a, 1)
                                 Wi -- Weight matrix of the update gate, numpy a
        rray of shape (n \ a, n \ a + n \ x)
                                 bi -- Bias of the update gate, numpy array of s
        hape (n a, 1)
                                 Wc -- Weight matrix of the first "tanh", numpy
        array of shape (n_a, n_a + n_x)
                                 bc -- Bias of the first "tanh", numpy array of
        shape (n \ a, 1)
                                 Wo -- Weight matrix of the output gate, numpy a
        rray of shape (n \ a, n \ a + n \ x)
                                 bo -- Bias of the output gate, numpy array of s
        hape (n a, 1)
                                 Wy -- Weight matrix relating the hidden-state t
        o the output, numpy array of shape (n y, n a)
                                 by -- Bias relating the hidden-state to the out
        put, numpy array of shape (n y, 1)
            Returns:
            a -- Hidden states for every time-step, numpy array of shape (n a,
        m, T x)
            y -- Predictions for every time-step, numpy array of shape (n y, m,
            caches -- tuple of values needed for the backward pass, contains (1
        ist of all the caches, x)
             # Initialize "caches", which will track the list of all the caches
            caches = []
            ### START CODE HERE ###
            # Retrieve dimensions from shapes of x and Wy (\approx2 lines)
            n x, m, T x = x.shape
            n y, n a = parameters["Wy"].shape
            # initialize "a", "c" and "y" with zeros (\approx 3 lines)
            a = np.zeros((n a, m, T x))
            c = np.zeros((n a, m, T x))
            y = np.zeros((n y, m, T x))
            # Initialize a next and c next (≈2 lines)
            a next = a0
```

```
In [10]: np.random.seed(1)
        x = np.random.randn(3, 10, 7)
        a0 = np.random.randn(5,10)
        Wf = np.random.randn(5, 5+3)
        bf = np.random.randn(5,1)
        Wi = np.random.randn(5, 5+3)
        bi = np.random.randn(5,1)
        Wo = np.random.randn(5, 5+3)
        bo = np.random.randn(5,1)
        Wc = np.random.randn(5, 5+3)
        bc = np.random.randn(5,1)
        Wy = np.random.randn(2,5)
        by = np.random.randn(2,1)
        parameters = {"Wf": Wf, "Wi": Wi, "Wo": Wo, "Wc": Wc, "Wy": Wy, "bf": b
        f, "bi": bi, "bo": bo, "bc": bc, "by": by}
        a, y, c, caches = lstm forward(x, a0, parameters)
        print("a[4][3][6] = ", a[4][3][6])
        print("a.shape = ", a.shape)
        print("y[1][4][3] =", y[1][4][3])
        print("y.shape = ", y.shape)
        print("caches[1][1[1]] =", caches[1][1][1])
        print("c[1][2][1]", c[1][2][1])
        print("len(caches) = ", len(caches))
        a[4][3][6] = 0.172117767533
        a.shape = (5, 10, 7)
        y[1][4][3] = 0.95087346185
        y.shape = (2, 10, 7)
        .20075807 0.18656139
          0.410051651
        c[1][2][1] -0.855544916718
        len(caches) = 2
```

a[4][3][6] =	0.172117767533
a.shape =	(5, 10, 7)
y[1][4][3] =	0.95087346185
y.shape =	(2, 10, 7)
caches[1][1][1] =	[0.82797464 0.23009474 0.76201118 -0.22232814 -0.20075807 0.18656139 0.41005165]
c[1][2][1] =	-0.855544916718
len(caches) =	2

Congratulations! You have now implemented the forward passes for the basic RNN and the LSTM. When using a deep learning framework, implementing the forward pass is sufficient to build systems that achieve great performance.

The rest of this notebook is optional, and will not be graded.

3 - Backpropagation in recurrent neural networks (OPTIONAL / UNGRADED)

In modern deep learning frameworks, you only have to implement the forward pass, and the framework takes care of the backward pass, so most deep learning engineers do not need to bother with the details of the backward pass. If however you are an expert in calculus and want to see the details of backprop in RNNs, you can work through this optional portion of the notebook.

When in an earlier course you implemented a simple (fully connected) neural network, you used backpropagation to compute the derivatives with respect to the cost to update the parameters. Similarly, in recurrent neural networks you can to calculate the derivatives with respect to the cost in order to update the parameters. The backprop equations are quite complicated and we did not derive them in lecture. However, we will briefly present them below.

3.1 - Basic RNN backward pass

We will start by computing the backward pass for the basic RNN-cell.

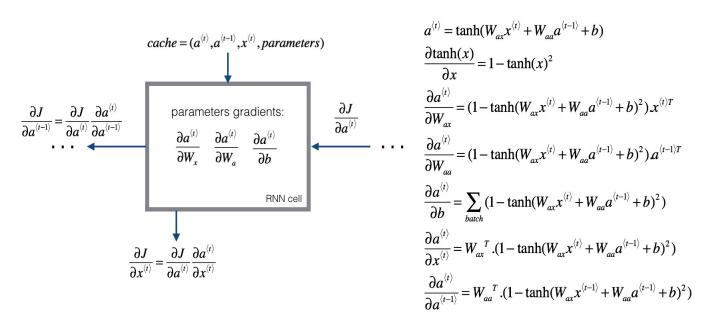


Figure 5: RNN-cell's backward pass. Just like in a fully-connected neural network, the derivative of the cost function J backpropagates through the RNN by following the chain-rule from calculas. The chain-rule is also used to calculate $(\frac{\partial J}{\partial W_{ax}}, \frac{\partial J}{\partial W_{aa}}, \frac{\partial J}{\partial b})$ to update the parameters (W_{ax}, W_{aa}, b_a) .

Deriving the one step backward functions:

To compute the $rnn_cell_backward$ you need to compute the following equations. It is a good exercise to derive them by hand.

The derivative of \tanh is $1-\tanh(x)^2$. You can find the complete proof <u>here (https://www.wyzant.com/resources/lessons/math/calculus/derivative_proofs/tanx)</u>. Note that: $\mathrm{sech}(x)^2=1-\tanh(x)^2$

Similarly for
$$\frac{\partial a^{\langle t \rangle}}{\partial W_{ax}}, \frac{\partial a^{\langle t \rangle}}{\partial W_{aa}}, \frac{\partial a^{\langle t \rangle}}{\partial b}$$
, the derivative of $\tanh(u)$ is $(1 - \tanh(u)^2)du$.

The final two equations also follow same rule and are derived using the tanh derivative. Note that the arrangement is done in a way to get the same dimensions to match.

```
In [11]: def rnn cell backward(da next, cache):
             Implements the backward pass for the RNN-cell (single time-step).
             Arguments:
             da next -- Gradient of loss with respect to next hidden state
             cache -- python dictionary containing useful values (output of rnn
         cell forward())
             Returns:
             gradients -- python dictionary containing:
                                  dx -- Gradients of input data, of shape (n x, m
                                  da prev -- Gradients of previous hidden state,
         of shape (n a, m)
                                  dWax -- Gradients of input-to-hidden weights, o
         f shape (n \ a, n \ x)
                                  dWaa -- Gradients of hidden-to-hidden weights,
         of shape (n a, n a)
                                 dba -- Gradients of bias vector, of shape (n a,
         1)
             11 11 11
             # Retrieve values from cache
             (a next, a prev, xt, parameters) = cache
             # Retrieve values from parameters
             Wax = parameters["Wax"]
             Waa = parameters["Waa"]
             Wya = parameters["Wya"]
             ba = parameters["ba"]
             by = parameters["by"]
             ### START CODE HERE ###
             # compute the gradient of tanh with respect to a next (≈1 line)
             dtanh = (1 - np.square(a next)) * da next
             # compute the gradient of the loss with respect to Wax (pprox 2 lines)
             dxt = np.dot(Wax.T, dtanh)
             dWax = np.dot(dtanh, xt.T)
             # compute the gradient with respect to Waa (≈2 lines)
             da prev = np.dot(Waa.T, dtanh)
             dWaa = np.dot(dtanh, a prev.T)
             # compute the gradient with respect to b (≈1 line)
             dba = np.sum(dtanh, axis = 1, keepdims=True)
             ### END CODE HERE ###
             # Store the gradients in a python dictionary
             gradients = {"dxt": dxt, "da prev": da prev, "dWax": dWax, "dWaa":
         dWaa, "dba": dba}
             return gradients
```

```
In [12]: | np.random.seed(1)
         xt = np.random.randn(3,10)
         a prev = np.random.randn(5,10)
         Wax = np.random.randn(5,3)
         Waa = np.random.randn(5,5)
         Wya = np.random.randn(2,5)
         b = np.random.randn(5,1)
         by = np.random.randn(2,1)
         parameters = {"Wax": Wax, "Waa": Waa, "Wya": Wya, "ba": ba, "by": by}
         a next, yt, cache = rnn cell forward(xt, a prev, parameters)
         da next = np.random.randn(5,10)
         gradients = rnn cell backward(da next, cache)
         print("gradients[\"dxt\"][1][2] =", gradients["dxt"][1][2])
         print("gradients[\"dxt\"].shape =", gradients["dxt"].shape)
         print("gradients[\"da prev\"][2][3] =", gradients["da prev"][2][3])
         print("gradients[\"da prev\"].shape =", gradients["da prev"].shape)
         print("gradients[\"dWax\"][3][1] = ", gradients["dWax"][3][1])
         print("gradients[\"dWax\"].shape =", gradients["dWax"].shape)
         print("gradients[\"dWaa\"][1][2] =", gradients["dWaa"][1][2])
         print("gradients[\"dWaa\"].shape =", gradients["dWaa"].shape)
         print("gradients[\"dba\"][4] =", gradients["dba"][4])
         print("gradients[\"dba\"].shape =", gradients["dba"].shape)
```

```
gradients["dxt"][1][2] = -0.460564103059
gradients["dxt"].shape = (3, 10)
gradients["da_prev"][2][3] = 0.0842968653807
gradients["da_prev"].shape = (5, 10)
gradients["dWax"][3][1] = 0.393081873922
gradients["dWax"].shape = (5, 3)
gradients["dWaa"][1][2] = -0.28483955787
gradients["dWaa"].shape = (5, 5)
gradients["dba"][4] = [ 0.80517166]
gradients["dba"].shape = (5, 1)
```

gradients["dxt"][1][2] =	-0.460564103059
gradients["dxt"].shape =	(3, 10)
gradients["da_prev"][2][3] =	0.0842968653807
gradients["da_prev"].shape =	(5, 10)
gradients["dWax"][3][1] =	0.393081873922
gradients["dWax"].shape =	(5, 3)
gradients["dWaa"][1][2] =	-0.28483955787
gradients["dWaa"].shape =	(5, 5)
gradients["dba"][4] =	[0.80517166]
gradients["dba"].shape =	(5, 1)

Backward pass through the RNN

Computing the gradients of the cost with respect to $a^{\langle t \rangle}$ at every time-step t is useful because it is what helps the gradient backpropagate to the previous RNN-cell. To do so, you need to iterate through all the time steps starting at the end, and at each step, you increment the overall db_a , dW_{aa} , dW_{ax} and you store dx.

Instructions:

Implement the $rnn_backward$ function. Initialize the return variables with zeros first and then loop through all the time steps while calling the $rnn_cell_backward$ at each time timestep, update the other variables accordingly.

```
In [13]: def rnn backward(da, caches):
              Implement the backward pass for a RNN over an entire sequence of in
         put data.
             Arguments:
             da -- Upstream gradients of all hidden states, of shape (n a, m, T
         _{X})
             caches -- tuple containing information from the forward pass (rnn f
         orward)
             Returns:
             gradients -- python dictionary containing:
                                  dx -- Gradient w.r.t. the input data, numpy-arr
         ay of shape (n \times, m, T \times)
                                  da0 -- Gradient w.r.t the initial hidden state,
         numpy-array of shape (n a, m)
                                  dWax -- Gradient w.r.t the input's weight matri
         x, numpy-array of shape (n a, n x)
                                  dWaa -- Gradient w.r.t the hidden state's weigh
         t matrix, numpy-arrayof shape (n a, n a)
                                  dba -- Gradient w.r.t the bias, of shape (n a,
         1)
              11 11 11
             ### START CODE HERE ###
             # Retrieve values from the first cache (t=1) of caches (≈2 lines)
             (caches, x) = caches
             (a1, a0, x1, parameters) = caches[1]
             # Retrieve dimensions from da's and x1's shapes (≈2 lines)
             n a, m, T x = da.shape
             n \times m = x1.shape
             # initialize the gradients with the right sizes (≈6 lines)
             dx = np.zeros((n x, m, T x))
             dWax = np.zeros((n a , n x))
             dWaa = np.zeros((n a, n a))
             dba = np.zeros((n a, 1))
             da0 = np.zeros((n a, m))
             da prevt = np.zeros((n a, m))
              # Loop through all the time steps
             for t in reversed(range(T x)):
                  # Compute gradients at time step t. Choose wisely the "da next"
         and the "cache" to use in the backward propagation step. (\approx 1 line)
                 gradients = rnn cell backward(da[:,:,t] + da prevt, caches[t])
                  # Retrieve derivatives from gradients (≈ 1 line)
                 dxt, da prevt, dWaxt, dWaat, dbat = gradients["dxt"], gradients
         ["da prev"], gradients["dWax"], gradients["dWaa"], gradients["dba"]
                  # Increment global derivatives w.r.t parameters by adding their
         derivative at time-step t (≈4 lines)
                 dx[:, :, t] = dxt
                 dWax += dWaxt
                 dWaa += dWaat
                 dba += dbat
```

```
In [14]: np.random.seed(1)
         x = np.random.randn(3, 10, 4)
         a0 = np.random.randn(5,10)
         Wax = np.random.randn(5,3)
         Waa = np.random.randn(5,5)
         Wya = np.random.randn(2,5)
         ba = np.random.randn(5,1)
         by = np.random.randn(2,1)
         parameters = {"Wax": Wax, "Waa": Waa, "Wya": Wya, "ba": ba, "by": by}
         a, y, caches = rnn forward(x, a0, parameters)
         da = np.random.randn(5, 10, 4)
         gradients = rnn backward(da, caches)
         print("gradients[\"dx\"][1][2] =", gradients["dx"][1][2])
         print("gradients[\"dx\"].shape =", gradients["dx"].shape)
         print("gradients[\"da0\"][2][3] =", gradients["da0"][2][3])
         print("gradients[\"da0\"].shape =", gradients["da0"].shape)
         print("gradients[\"dWax\"][3][1] = ", gradients["dWax"][3][1])
         print("gradients[\"dWax\"].shape =", gradients["dWax"].shape)
         print("gradients[\"dWaa\"][1][2] =", gradients["dWaa"][1][2])
         \verb|print("gradients[\"dWaa\"].shape = ", gradients["dWaa"].shape)|\\
         print("gradients[\"dba\"][4] =", gradients["dba"][4])
         print("gradients[\"dba\"].shape =", gradients["dba"].shape)
         gradients["dx"][1][2] = [-2.07101689 -0.59255627 0.02466855 0.01483
         317]
         gradients["dx"].shape = (3, 10, 4)
         gradients["da0"][2][3] = -0.314942375127
```

```
gradients["dx"][1][2] = [-2.07101689 -0.59255627 0.02466855 0.01483
317]
gradients["dx"].shape = (3, 10, 4)
gradients["da0"][2][3] = -0.314942375127
gradients["da0"].shape = (5, 10)
gradients["dWax"][3][1] = 11.2641044965
gradients["dWax"].shape = (5, 3)
gradients["dWaa"][1][2] = 2.30333312658
gradients["dWaa"].shape = (5, 5)
gradients["dba"][4] = [-0.74747722]
gradients["dba"].shape = (5, 1)
```

gradients["dx"][1][2] =	[-2.07101689 -0.59255627 0.02466855 0.01483317]
gradients["dx"].shape =	(3, 10, 4)
gradients["da0"][2][3] =	-0.314942375127
gradients["da0"].shape =	(5, 10)
gradients["dWax"][3][1] =	11.2641044965
gradients["dWax"].shape =	(5, 3)
gradients["dWaa"][1][2] =	2.30333312658
gradients["dWaa"].shape =	(5, 5)
gradients["dba"][4] =	[-0.74747722]
gradients["dba"].shape =	(5, 1)

3.2 - LSTM backward pass

3.2.1 One Step backward

The LSTM backward pass is slightly more complicated than the forward one. We have provided you with all the equations for the LSTM backward pass below. (If you enjoy calculus exercises feel free to try deriving these from scratch yourself.)

3.2.2 gate derivatives

$$d\Gamma_o^{\langle t \rangle} = da_{next} * \tanh(c_{next}) * \Gamma_o^{\langle t \rangle} * (1 - \Gamma_o^{\langle t \rangle})$$
 (7)

$$d ilde{c}^{\langle t
angle} = (dc_{next} * \Gamma_u^{\langle t
angle} + \Gamma_o^{\langle t
angle} * (1 - anh(c_{next})^2) * \Gamma_u^{\langle t
angle} * da_{next}) * (1 - (ilde{c}^{\langle t
angle})^2)$$

$$d\Gamma_{u}^{\langle t \rangle} = (dc_{next} * \tilde{c}^{\langle t \rangle} + \Gamma_{o}^{\langle t \rangle} * (1 - \tanh(c_{next})^{2}) * \tilde{c}^{\langle t \rangle} * da_{next}) * \Gamma_{u}^{\langle t \rangle} * (1 - \Gamma_{u}^{\langle t \rangle})$$
(9)

$$d\Gamma_f^{\langle t \rangle} = (dc_{next} * c_{prev} + \Gamma_o^{\langle t \rangle} * (1 - \tanh(c_{next})^2) * c_{prev} * da_{next}) * \Gamma_f^{\langle t \rangle} * (1 - \Gamma_f^{\langle t \rangle})$$
 (10)

3.2.3 parameter derivatives

$$dW_f = d\Gamma_f^{\langle t \rangle} \cdot \left(rac{a_{prev}}{x_t}
ight)^T$$
 (11)

$$dW_u = d\Gamma_u^{\langle t \rangle}. \left(rac{a_{prev}}{x_t}
ight)^T$$
 (12)

$$dW_o = d\Gamma_o^{\langle t \rangle} \cdot \left(rac{a_{prev}}{x_t}
ight)^T$$
 (14)

To calculate db_f, db_u, db_c, db_o you just need to sum across the horizontal (axis= 1) axis on $d\Gamma_f^{\langle t \rangle}, d\Gamma_u^{\langle t \rangle}, d\tilde{c}^{\langle t \rangle}, d\Gamma_o^{\langle t \rangle}$ respectively. Note that you should have the <code>keep_dims=True</code> option.

Finally, you will compute the derivative with respect to the previous hidden state, previous memory state, and input.

$$da_{prev} = W_f^T. d\Gamma_f^{\langle t \rangle} + W_u^T. d\Gamma_u^{\langle t \rangle} + W_c^T. d\tilde{c}^{\langle t \rangle} + W_o^T. d\Gamma_o^{\langle t \rangle}$$
 (15)

Here, the weights for equations 15 are the first n_a, (i.e. $W_f = W_f[:,:na]$ etc...)

$$dc_{prev} = dc_{next} * \Gamma_f^{\langle t \rangle} + \Gamma_o^{\langle t \rangle} * (1 - \tanh(c_{next})^2) * \Gamma_f^{\langle t \rangle} * da_{next}$$
(16)

$$dx^{\langle t \rangle} = W_f^T. d\Gamma_f^{\langle t \rangle} + W_u^T. d\Gamma_u^{\langle t \rangle} + W_c^T. d\tilde{c}_t + W_o^T. d\Gamma_o^{\langle t \rangle}$$
(17)

where the weights for equation 17 are from n_a to the end, (i.e. $W_f = W_f[:,na:]$ etc...)

Exercise: Implement $lstm_cell_backward$ by implementing equations 7-17 below. Good luck!:)

```
In [21]: def lstm cell backward(da next, dc next, cache):
             Implement the backward pass for the LSTM-cell (single time-step).
             Arguments:
             da next -- Gradients of next hidden state, of shape (n a, m)
             dc next -- Gradients of next cell state, of shape (n a, m)
             cache -- cache storing information from the forward pass
             Returns:
             gradients -- python dictionary containing:
                                  dxt -- Gradient of input data at time-step t, o
         f shape (n \times, m)
                                  da prev -- Gradient w.r.t. the previous hidden
         state, numpy array of shape (n a, m)
                                  dc prev -- Gradient w.r.t. the previous memory
         state, of shape (n \ a, \ m, \ T \ x)
                                  dWf -- Gradient w.r.t. the weight matrix of the
         forget gate, numpy array of shape (n a, n a + n x)
                                 dWi -- Gradient w.r.t. the weight matrix of the
         update gate, numpy array of shape (n a, n a + n x)
                                  dWc -- Gradient w.r.t. the weight matrix of the
         memory gate, numpy array of shape (n \ a, n \ a + n \ x)
                                  dWo -- Gradient w.r.t. the weight matrix of the
         output gate, numpy array of shape (n a, n a + n x)
                                  dbf -- Gradient w.r.t. biases of the forget gat
         e, of shape (n a, 1)
                                  dbi -- Gradient w.r.t. biases of the update gat
         e, of shape (n a, 1)
                                  dbc -- Gradient w.r.t. biases of the memory gat
         e, of shape (n_a, 1)
                                 dbo -- Gradient w.r.t. biases of the output gat
         e, of shape (n a, 1)
              # Retrieve information from "cache"
              (a next, c next, a prev, c prev, ft, it, cct, ot, xt, parameters) =
         cache
             ### START CODE HERE ###
             # Retrieve dimensions from xt's and a next's shape (*2 lines)
             n \times m = xt.shape
             n a, m = a next.shape
             concat = np.vstack((a prev, xt))
              # Compute gates related derivatives, you can find their values can
         be found by looking carefully at equations (7) to (10) (\approx4 lines)
             dot = None
             dcct = None
             dit = None
             dft = None
             # Code equations (7) to (10) (\approx 4 lines)
             dit = ((dc next * cct) + (ot * (1.0 - np.square(np.tanh(c next))) *
         cct * da next)) * it * (1.0 - it)
```

```
In [22]: | np.random.seed(1)
         xt = np.random.randn(3,10)
         a prev = np.random.randn(5,10)
         c prev = np.random.randn(5,10)
         Wf = np.random.randn(5, 5+3)
         bf = np.random.randn(5,1)
         Wi = np.random.randn(5, 5+3)
         bi = np.random.randn(5,1)
         Wo = np.random.randn(5, 5+3)
         bo = np.random.randn(5,1)
         Wc = np.random.randn(5, 5+3)
         bc = np.random.randn(5,1)
         Wy = np.random.randn(2,5)
         by = np.random.randn(2,1)
         parameters = {"Wf": Wf, "Wi": Wi, "Wo": Wo, "Wc": Wc, "Wy": Wy, "bf": b
         f, "bi": bi, "bo": bo, "bc": bc, "by": by}
         a next, c next, yt, cache = lstm cell forward(xt, a prev, c prev, param
         eters)
         da next = np.random.randn(5,10)
         dc next = np.random.randn(5,10)
         gradients = lstm cell backward(da next, dc next, cache)
         print("gradients[\"dxt\"][1][2] =", gradients["dxt"][1][2])
         print("gradients[\"dxt\"].shape =", gradients["dxt"].shape)
         print("gradients[\"da prev\"][2][3] =", gradients["da prev"][2][3])
         print("gradients[\"da prev\"].shape =", gradients["da prev"].shape)
         print("gradients[\"dc prev\"][2][3] =", gradients["dc prev"][2][3])
         print("gradients[\"dc prev\"].shape =", gradients["dc prev"].shape)
         print("gradients[\"dWf\"][3][1] = ", gradients["dWf"][3][1])
         print("gradients[\"dWf\"].shape =", gradients["dWf"].shape)
         print("gradients[\"dWi\"][1][2] =", gradients["dWi"][1][2])
         print("gradients[\"dWi\"].shape =", gradients["dWi"].shape)
         print("gradients[\"dWc\"][3][1] =", gradients["dWc"][3][1])
         print("gradients[\"dWc\"].shape =", gradients["dWc"].shape)
         print("gradients[\"dWo\"][1][2] =", gradients["dWo"][1][2])
         print("gradients[\"dWo\"].shape =", gradients["dWo"].shape)
         print("gradients[\"dbf\"][4] =", gradients["dbf"][4])
         print("gradients[\"dbf\"].shape =", gradients["dbf"].shape)
         print("gradients[\"dbi\"][4] =", gradients["dbi"][4])
         print("gradients[\"dbi\"].shape =", gradients["dbi"].shape)
         print("gradients[\"dbc\"][4] =", gradients["dbc"][4])
         print("gradients[\"dbc\"].shape =", gradients["dbc"].shape)
         print("gradients[\"dbo\"][4] =", gradients["dbo"][4])
         print("gradients[\"dbo\"].shape =", gradients["dbo"].shape)
```

```
gradients["dxt"][1][2] = 3.23055911511
gradients["dxt"].shape = (3, 10)
gradients["da prev"][2][3] = -0.0639621419711
gradients["da prev"].shape = (5, 10)
gradients["dc prev"][2][3] = 0.797522038797
gradients["dc prev"].shape = (5, 10)
gradients["dWf"][3][1] = -0.147954838164
gradients["dWf"].shape = (5, 8)
gradients["dWi"][1][2] = 1.05749805523
gradients["dWi"].shape = (5, 8)
gradients["dWc"][3][1] = 2.30456216369
gradients["dWc"].shape = (5, 8)
gradients["dWo"][1][2] = 0.331311595289
gradients["dWo"].shape = (5, 8)
gradients["dbf"][4] = [ 0.18864637]
gradients["dbf"].shape = (5, 1)
gradients["dbi"][4] = [-0.40142491]
gradients["dbi"].shape = (5, 1)
gradients["dbc"][4] = [ 0.25587763]
gradients["dbc"].shape = (5, 1)
gradients["dbo"][4] = [ 0.13893342]
gradients["dbo"].shape = (5, 1)
```

gradients["dxt"][1][2] =	3.23055911511
gradients["dxt"].shape =	(3, 10)
gradients["da_prev"][2][3] =	-0.0639621419711
gradients["da_prev"].shape =	(5, 10)
gradients["dc_prev"][2][3] =	0.797522038797
gradients["dc_prev"].shape =	(5, 10)
gradients["dWf"][3][1] =	-0.147954838164
gradients["dWf"].shape =	(5, 8)
gradients["dWi"][1][2] =	1.05749805523
gradients["dWi"].shape =	(5, 8)
gradients["dWc"][3][1] =	2.30456216369
gradients["dWc"].shape =	(5, 8)
gradients["dWo"][1][2] =	0.331311595289
gradients["dWo"].shape =	(5, 8)
gradients["dbf"][4] =	[0.18864637]
gradients["dbf"].shape =	(5, 1)
gradients["dbi"][4] =	[-0.40142491]
gradients["dbi"].shape =	(5, 1)
gradients["dbc"][4] =	[0.25587763]
gradients["dbc"].shape =	(5, 1)
gradients["dbo"][4] =	[0.13893342]
gradients["dbo"].shape =	(5, 1)
	

3.3 Backward pass through the LSTM RNN

This part is very similar to the rnn_backward function you implemented above. You will first create variables of the same dimension as your return variables. You will then iterate over all the time steps starting from the end and call the one step function you implemented for LSTM at each iteration. You will then update the parameters by summing them individually. Finally return a dictionary with the new gradients.

Instructions: Implement the $lstm_backward$ function. Create a for loop starting from T_x and going backward. For each step call $lstm_cell_backward$ and update the your old gradients by adding the new gradients to them. Note that dxt is not updated but is stored.

```
In [37]: def lstm backward(da, caches):
              Implement the backward pass for the RNN with LSTM-cell (over a whol
         e sequence).
             Arguments:
             da -- Gradients w.r.t the hidden states, numpy-array of shape (n a,
         m, T x)
             dc -- Gradients w.r.t the memory states, numpy-array of shape (n a,
         m, T x)
             caches -- cache storing information from the forward pass (1stm for
         ward)
             Returns:
             gradients -- python dictionary containing:
                                  dx -- Gradient of inputs, of shape (n x, m, T x
                                  da0 -- Gradient w.r.t. the previous hidden stat
         e, numpy array of shape (n a, m)
                                  dWf -- Gradient w.r.t. the weight matrix of the
         forget gate, numpy array of shape (n a, n a + n x)
                                  dWi -- Gradient w.r.t. the weight matrix of the
         update gate, numpy array of shape (n \ a, \ n \ a + n \ x)
                                  dWc -- Gradient w.r.t. the weight matrix of the
         memory gate, numpy array of shape (n_a, n_a + n_x)
                                  dWo -- Gradient w.r.t. the weight matrix of the
         save gate, numpy array of shape (n \ a, \ n \ a + n \ x)
                                  dbf -- Gradient w.r.t. biases of the forget gat
         e, of shape (n a, 1)
                                  dbi -- Gradient w.r.t. biases of the update gat
         e, of shape (n a, 1)
                                  dbc -- Gradient w.r.t. biases of the memory gat
         e, of shape (n a, 1)
                                  dbo -- Gradient w.r.t. biases of the save gate,
         of shape (n a, 1)
             11 11 11
              # Retrieve values from the first cache (t=1) of caches.
              (caches, x) = caches
             (a1, c1, a0, c0, f1, i1, cc1, o1, x1, parameters) = caches[0]
              ### START CODE HERE ###
             # Retrieve dimensions from da's and x1's shapes (\approx 2 lines)
             n_a, m, T_x = da.shape
             n \times m = x1.shape
             # initialize the gradients with the right sizes (≈12 lines)
             dx = np.zeros((n x, m, T x))
             da0 = np.zeros((n a, m))
             da prevt = np.zeros((n a, m))
             dc prevt = np.zeros((n a, m))
             dWf = np.zeros((n a, n a + n x))
             dWi = np.zeros((n a, n a + n x))
             dWc = np.zeros((n a, n a + n x))
             dWo = np.zeros((n a, n a + n x))
```

```
In [38]: np.random.seed(1)
         x = np.random.randn(3, 10, 7)
         a0 = np.random.randn(5,10)
         Wf = np.random.randn(5, 5+3)
         bf = np.random.randn(5,1)
         Wi = np.random.randn(5, 5+3)
         bi = np.random.randn(5,1)
         Wo = np.random.randn(5, 5+3)
         bo = np.random.randn(5,1)
         Wc = np.random.randn(5, 5+3)
         bc = np.random.randn(5,1)
         parameters = {"Wf": Wf, "Wi": Wi, "Wo": Wo, "Wc": Wc, "Wy": Wy, "bf": b
         f, "bi": bi, "bo": bo, "bc": bc, "by": by}
         a, y, c, caches = lstm forward(x, a0, parameters)
         da = np.random.randn(5, 10, 4)
         gradients = lstm backward(da, caches)
         print("gradients[\"dx\"][1][2] =", gradients["dx"][1][2])
         print("gradients[\"dx\"].shape =", gradients["dx"].shape)
         print("gradients[\"da0\"][2][3] =", gradients["da0"][2][3])
         print("gradients[\"da0\"].shape =", gradients["da0"].shape)
         print("gradients[\"dWf\"][3][1] = ", gradients["dWf"][3][1])
         print("gradients[\"dWf\"].shape =", gradients["dWf"].shape)
         print("gradients[\"dWi\"][1][2] = ", gradients["dWi"][1][2])
         print("gradients[\"dWi\"].shape =", gradients["dWi"].shape)
         print("gradients[\"dWc\"][3][1] =", gradients["dWc"][3][1])
         print("gradients[\"dWc\"].shape =", gradients["dWc"].shape)
         print("gradients[\"dWo\"][1][2] =", gradients["dWo"][1][2])
         print("gradients[\"dWo\"].shape =", gradients["dWo"].shape)
         print("gradients[\"dbf\"][4] =", gradients["dbf"][4])
         print("gradients[\"dbf\"].shape =", gradients["dbf"].shape)
         print("gradients[\"dbi\"][4] =", gradients["dbi"][4])
         print("gradients[\"dbi\"].shape =", gradients["dbi"].shape)
         print("gradients[\"dbc\"][4] =", gradients["dbc"][4])
         print("gradients[\"dbc\"].shape =", gradients["dbc"].shape)
         print("gradients[\"dbo\"][4] =", gradients["dbo"][4])
         print("gradients[\"dbo\"].shape =", gradients["dbo"].shape)
```

```
gradients["dx"][1][2] = [-0.00173313  0.08287442  -0.30545663  -0.43281
115]
gradients["dx"].shape = (3, 10, 4)
gradients["da0"][2][3] = -0.095911501954
gradients["da0"].shape = (5, 10)
gradients["dWf"][3][1] = -0.0698198561274
gradients["dWf"].shape = (5, 8)
gradients["dWi"][1][2] = 0.102371820249
gradients["dWi"].shape = (5, 8)
gradients["dWc"][3][1] = -0.0624983794927
gradients["dWc"].shape = (5, 8)
gradients["dWo"][1][2] = 0.0484389131444
gradients["dWo"].shape = (5, 8)
gradients["dbf"][4] = [-0.0565788]
gradients["dbf"].shape = (5, 1)
gradients["dbi"][4] = [-0.15399065]
gradients["dbi"].shape = (5, 1)
gradients["dbc"][4] = [-0.29691142]
gradients["dbc"].shape = (5, 1)
gradients["dbo"][4] = [-0.29798344]
gradients["dbo"].shape = (5, 1)
```

gradients["dx"][1][2] =	[-0.00173313 0.08287442 -0.30545663 -0.43281115]
gradients["dx"].shape =	(3, 10, 4)
gradients["da0"][2][3] =	-0.095911501954
gradients["da0"].shape =	(5, 10)
gradients["dWf"][3][1] =	-0.0698198561274
gradients["dWf"].shape =	(5, 8)
gradients["dWi"][1][2] =	0.102371820249
gradients["dWi"].shape =	(5, 8)
gradients["dWc"][3][1] =	-0.0624983794927
gradients["dWc"].shape =	(5, 8)
gradients["dWo"][1][2] =	0.0484389131444
gradients["dWo"].shape =	(5, 8)
gradients["dbf"][4] =	[-0.0565788]
gradients["dbf"].shape =	(5, 1)
gradients["dbi"][4] =	[-0.06997391]
gradients["dbi"].shape =	(5, 1)
gradients["dbc"][4] =	[-0.27441821]
gradients["dbc"].shape =	(5, 1)
gradients["dbo"][4] =	[0.16532821]
gradients["dbo"].shape =	(5, 1)
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Congratulations!

Congratulations on completing this assignment. You now understand how recurrent neural networks work!

Lets go on to the next exercise, where you'll use an RNN to build a character-level language model.