Building your Recurrent Neural Network - Step by Step

Welcome to Course 5's first assignment! In this assignment, you will implement key components of a 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 unidirectional RNN to take information from the past to process later inputs. A bidirectional RNN can take context from both the past and the future.

Notation:

- Superscript $[l]$ denotes an object associated with the l^{th} layer.
- Superscript (i) denotes an object associated with the i^{th} example.
- Superscript $\langle t \rangle$ denotes an object at the t^{th} time-step.
- **Sub**script i denotes the i^{th} entry of a vector.

Example:

- $a_5^{(2)[3]<4>}$ denotes the activation of the 2nd training example (2), 3rd layer [3], 4th time step , and 5th entry in the vector.

Pre-requisites

- We assume that you are already familiar with numpy.
- To refresh your knowledge of numpy, you can review course 1 of this specialization "Neural Networks and Deep Learning".
  - Specifically, review the week 2 assignment "Python Basics with numpy (optional)" (https://www.coursera.org/learn/neural-networks-deep-learning/item/Zh0CU).

Be careful when modifying the starter code

- When working on graded functions, please remember to only modify the code that is between the

  ```
  #### START CODE HERE
  ```

  and

  ```
  #### END CODE HERE
  ```

- In particular, Be careful to not modify the first line of graded routines. These start with:

  ```
  # GRADED FUNCTION: routine_name
  ```

- The automatic grader (autograder) needs these to locate the function.

- Even a change in spacing will cause issues with the autograder.
- It will return 'failed' if these are modified or missing."

Updates for 3b

If you were working on the notebook before this update...

- The current notebook is version "3b".
- You can find your original work saved in the notebook with the previous version name ("v3a")
- To view the file directory, go to the menu "File->Open", and this will open a new tab that shows the file directory.

List of updates

- `rnn_cell_backward`
  - fixed error in equations
  - harmonize rnn backward diagram with rnn_forward diagram and fixed Wax multiple (changed from at to xt).
  - clarified dba batch as summing 'm' examples
  - aligned equations
- `lstm_cell_backward`
  - aligned equations
- `lstm_forward`
  - fixed typo, Wb to bf
- `lstm_cell_forward`
  - changed c_next_tmp.shape to a_next_tmp.shape in test case
  - clarified dbxx batch as summing 'm' examples

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

In [3]:
```
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, $T_x = T_y$.

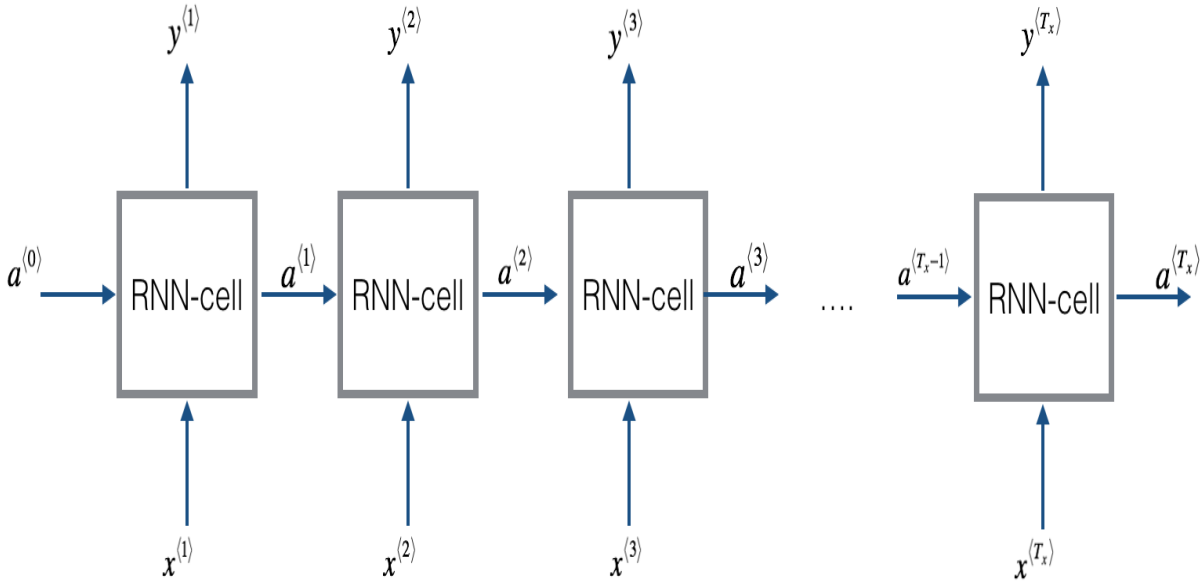

Figure 1: Basic RNN model

Dimensions of input x

Input with n_x number of units

- For a single timestep of a single input example, $x^{(i)\langle t \rangle}$ is a one-dimensional input vector.
- Using language as an example, a language with a 5000 word vocabulary could be one-hot encoded into a vector that has 5000 units. So $x^{(i)\langle t \rangle}$ would have the shape (5000,).
- We'll use the notation n_x to denote the number of units in a single timestep of a single training example.

Time steps of size T_x

- A recurrent neural network has multiple time steps, which we'll index with t.
- In the lessons, we saw a single training example $x^{(i)}$ consist of multiple time steps T_x. For example, if there are 10 time steps, $T_x = 10$

Batches of size m

- Let's say we have mini-batches, each with 20 training examples.
- To benefit from vectorization, we'll stack 20 columns of $x^{(i)}$ examples.
- For example, this tensor has the shape (5000,20,10).
- We'll use m to denote the number of training examples.
- So the shape of a mini-batch is (n_x, m, T_x)

3D Tensor of shape (n_x, m, T_x)

- The 3-dimensional tensor x of shape (n_x, m, T_x) represents the input x that is fed into the RNN.

Taking a 2D slice for each time step: $x^{\langle t \rangle}$

- At each time step, we'll use a mini-batches of training examples (not just a single example).
- So, for each time step t, we'll use a 2D slice of shape (n_x, m).
- We're referring to this 2D slice as $x^{\langle t \rangle}$. The variable name in the code is `xt`.

Definition of hidden state a

- The activation $a^{\langle t \rangle}$ that is passed to the RNN from one time step to another is called a "hidden state."

Dimensions of hidden state a

- Similar to the input tensor x, the hidden state for a single training example is a vector of length n_a.
- If we include a mini-batch of m training examples, the shape of a mini-batch is (n_a, m).
- When we include the time step dimension, the shape of the hidden state is (n_a, m, T_x)
- We will loop through the time steps with index t, and work with a 2D slice of the 3D tensor.
- We'll refer to this 2D slice as $a^{\langle t \rangle}$.
- In the code, the variable names we use are either `a_prev` or `a_next`, depending on the function that's being implemented.
- The shape of this 2D slice is (n_a, m)

Dimensions of prediction $\hat{y}$

- Similar to the inputs and hidden states, $\hat{y}$ is a 3D tensor of shape (n_y, m, T_y).
    - n_y: number of units in the vector representing the prediction.
    - m: number of examples in a mini-batch.
    - T_y: number of time steps in the prediction.
- For a single time step t, a 2D slice $\hat{y}^{\langle t \rangle}$ has shape (n_y, m).
- In the code, the variable names are:
    - `y_pred`: $\hat{y}$
    - `yt_pred`: $\hat{y}^{\langle t \rangle}$

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.

1.1 - RNN cell

A recurrent neural network can be seen as the repeated use 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.

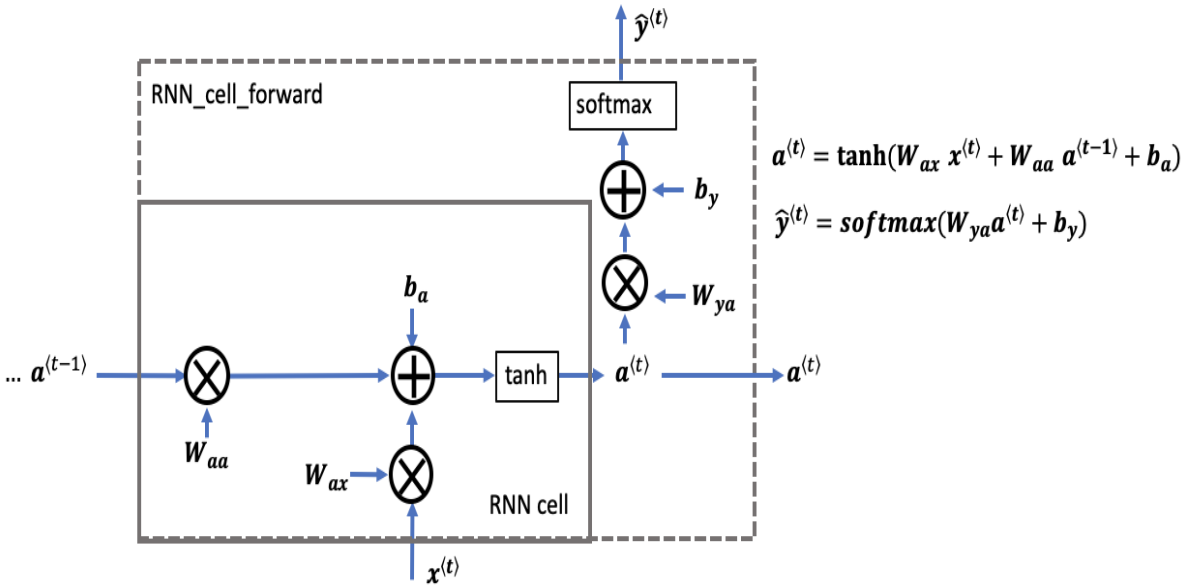

Figure 2: Basic RNN cell. Takes as input $x^{\langle t \rangle}$ (current input) and $a^{\langle t-1 \rangle}$ (previous hidden state containing information from the past), and outputs $a^{\langle t \rangle}$ which is given to the next RNN cell and also used to predict $\hat{y}^{\langle t \rangle}$

rnn cell versus rnn_cell_forward

- Note that an RNN cell outputs the hidden state $a^{\langle t \rangle}$.
    - The rnn cell is shown in the figure as the inner box which has solid lines.
- The function that we will implement, `rnn_cell_forward`, also calculates the prediction $\hat{y}^{\langle t \rangle}$
    - The rnn_cell_forward is shown in the figure as the outer box that has dashed lines.

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

Instructions:

1. Compute the hidden state with tanh activation: $a^{\langle t \rangle} = \tanh(W_{aa} a^{\langle t-1 \rangle} + W_{ax} x^{\langle t \rangle} + b_a)$.
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 the function `softmax`.
3. Store $(a^{\langle t \rangle}, a^{\langle t-1 \rangle}, x^{\langle t \rangle}, parameters)$ in a `cache`.
4. Return $a^{\langle t \rangle}$, $\hat{y}^{\langle t \rangle}$ and `cache`

Additional Hints

- numpy.tanh (https://www.google.com/search?q=numpy+tanh&rlz=1C5CHFA_enUS854US855&oq=numpy+tanh&aqs=chrome..69i57j0l5.134
8)

- We've created a `softmax` function that you can use. It is located in the file 'rnn_utils.py' and has been imported.
- For matrix multiplication, use numpy.dot (https://docs.scipy.org/doc/numpy/reference/generated/numpy.dot.html)

In [4]:
```
# 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 Figure (

    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, numpy arr
                        Waa -- Weight matrix multiplying the hidden state, nu
                        Wya -- Weight matrix relating the hidden-state to the
                        ba --  Bias, numpy array of shape (n_a, 1)
                        by -- Bias relating the hidden-state to the output, n
    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,
    """

    # 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(Waa, a_prev) + np.dot(Wax, xt) + 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 [5]:
```
np.random.seed(1)
xt_tmp = np.random.randn(3,10)
a_prev_tmp = np.random.randn(5,10)
parameters_tmp = {}
parameters_tmp['Waa'] = np.random.randn(5,5)
parameters_tmp['Wax'] = np.random.randn(5,3)
parameters_tmp['Wya'] = np.random.randn(2,5)
parameters_tmp['ba'] = np.random.randn(5,1)
parameters_tmp['by'] = np.random.randn(2,1)

a_next_tmp, yt_pred_tmp, cache_tmp = rnn_cell_forward(xt_tmp, a_prev_tmp, par
print("a_next[4] = \n", a_next_tmp[4])
print("a_next.shape = \n", a_next_tmp.shape)
print("yt_pred[1] =\n", yt_pred_tmp[1])
print("yt_pred.shape = \n", yt_pred_tmp.shape)
```

```
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_pred[1] =
 [ 0.9888161   0.01682021  0.21140899  0.36817467  0.98988387  0.88945212
  0.36920224  0.9966312   0.9982559   0.17746526]
yt_pred.shape =
 (2, 10)
```

Expected Output:

```
a_next[4] =
 [ 0.59584544  0.18141802  0.61311866  0.99808218  0.85016201  0.9998
0978
 -0.18887155  0.99815551  0.6531151   0.82872037]
a_next.shape =
 (5, 10)
yt_pred[1] =
 [ 0.9888161   0.01682021  0.21140899  0.36817467  0.98988387  0.8894
5212
  0.36920224  0.9966312   0.9982559   0.17746526]
yt_pred.shape =
 (2, 10)
```

1.2 - RNN forward pass

- A recurrent neural network (RNN) is a repetition of the RNN cell that you've just built.
  - If your input sequence of data is 10 time steps long, then you will re-use the RNN cell 10 times.
- Each cell takes two inputs at each time step:

- $a^{\langle t-1 \rangle}$: The hidden state from the previous cell.
- $x^{\langle t \rangle}$: The current time-step's input data.

- It has two outputs at each time step:
  - A hidden state ($a^{\langle t \rangle}$)
  - A prediction ($y^{\langle t \rangle}$)
- The weights and biases (W_{aa}, b_a, W_{ax}, b_x) are re-used each time step.
  - They are maintained between calls to rnn_cell_forward in the 'parameters' dictionary.

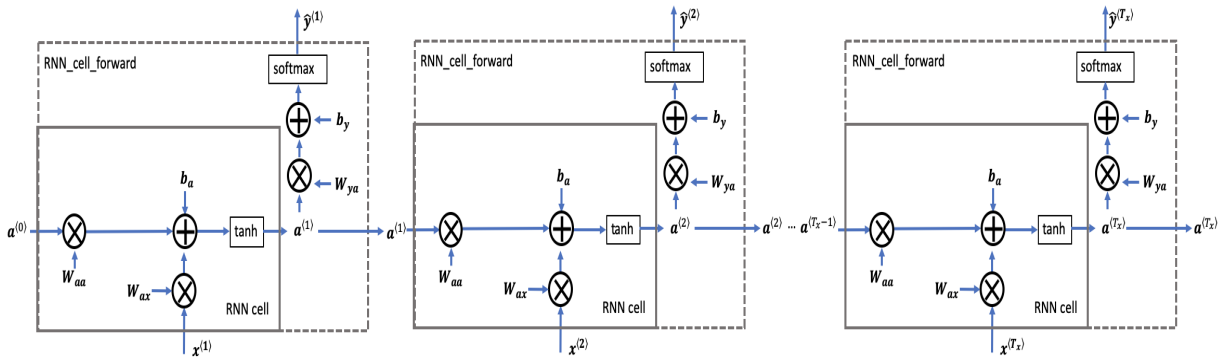

Figure 3: Basic RNN. The input sequence $x = (x^{\langle 1 \rangle}, x^{\langle 2 \rangle}, \ldots, x^{\langle T_x \rangle})$ is carried over T_x time steps. The network outputs $y = (y^{\langle 1 \rangle}, y^{\langle 2 \rangle}, \ldots, y^{\langle T_x \rangle})$.

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

Instructions:

- Create a 3D array of zeros, a of shape (n_a, m, T_x) that will store all the hidden states computed by the RNN.
- Create a 3D array of zeros, $\hat{y}$, of shape (n_y, m, T_x) that will store the predictions.
  - Note that in this case, $T_y = T_x$ (the prediction and input have the same number of time steps).
- Initialize the 2D hidden state `a_next` by setting it equal to the initial hidden state, a_0.
- At each time step t:
  - Get $x^{\langle t \rangle}$, which is a 2D slice of x for a single time step t.
    - $x^{\langle t \rangle}$ has shape (n_x, m)
    - x has shape (n_x, m, T_x)
  - Update the 2D hidden state $a^{\langle t \rangle}$ (variable name `a_next`), the prediction $\hat{y}^{\langle t \rangle}$ and the cache by running `rnn_cell_forward`.
    - $a^{\langle t \rangle}$ has shape (n_a, m)
  - Store the 2D hidden state in the 3D tensor a, at the t^{th} position.
    - a has shape (n_a, m, T_x)
  - Store the 2D $\hat{y}^{\langle t \rangle}$ prediction (variable name `yt_pred`) in the 3D tensor $\hat{y}_{pred}$ at the t^{th} position.
    - $\hat{y}^{\langle t \rangle}$ has shape (n_y, m)
    - $\hat{y}$ has shape (n_y, m, T_x)

    - Append the cache to the list of caches.
- Return the 3D tensor a and $\hat{y}$, as well as the list of caches.

Additional Hints

- np.zeros (https://docs.scipy.org/doc/numpy/reference/generated/numpy.zeros.html)
- If you have a 3 dimensional numpy array and are indexing by its third dimension, you can use array slicing like this: `var_name[:,:,i]`.

In [6]:
```
# GRADED FUNCTION: rnn_forward

def rnn_forward(x, a0, parameters):
    """
    Implement the forward propagation of the recurrent neural network describ

    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 state, nu
                        Wax -- Weight matrix multiplying the input, numpy arr
                        Wya -- Weight matrix relating the hidden-state to the
                        ba --  Bias numpy array of shape (n_a, 1)
                        by -- Bias relating the hidden-state to the output, n

    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,
    caches -- tuple of values needed for the backward pass, contains (list of
    """

    # Initialize "caches" which will contain the list of all caches
    caches = []

    # Retrieve dimensions from shapes of x and parameters["Wya"]
    n_x, m, T_x = x.shape
    n_y, n_a = parameters["Wya"].shape

    ### START CODE HERE ###

    # initialize "a" and "y_pred" with zeros (≈2 lines)
    a = np.zeros((n_a, m, T_x))
    y_pred = np.zeros((n_y, m, T_x))

    # Initialize a_next (≈1 line)
    a_next = a0

    # loop over all time-steps of the input 'x' (1 line)
    for t in range(T_x):
        # Update next hidden state, compute the prediction, get the cache (≈2
        xt = x[:, :, t]
        a_next, yt_pred, cache = rnn_cell_forward(xt, a_next, parameters)
        # 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
        # Append "cache" to "caches" (≈1 line)
        caches.append(cache)

    ### END CODE HERE ###

    # store values needed for backward propagation in cache
    caches = (caches, x)

    return a, y_pred, caches
```

In [7]:
```
np.random.seed(1)
x_tmp = np.random.randn(3,10,4)
a0_tmp = np.random.randn(5,10)
parameters_tmp = {}
parameters_tmp['Waa'] = np.random.randn(5,5)
parameters_tmp['Wax'] = np.random.randn(5,3)
parameters_tmp['Wya'] = np.random.randn(2,5)
parameters_tmp['ba'] = np.random.randn(5,1)
parameters_tmp['by'] = np.random.randn(2,1)

a_tmp, y_pred_tmp, caches_tmp = rnn_forward(x_tmp, a0_tmp, parameters_tmp)
print("a[4][1] = \n", a_tmp[4][1])
print("a.shape = \n", a_tmp.shape)
print("y_pred[1][3] =\n", y_pred_tmp[1][3])
print("y_pred.shape = \n", y_pred_tmp.shape)
print("caches[1][1][3] =\n", caches_tmp[1][1][3])
print("len(caches) = \n", len(caches_tmp))
```

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

Expected Output:

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

Congratulations! You've successfully built the forward propagation of a recurrent neural network from scratch.

Situations when this RNN will perform better:

- This will work well enough for some applications, but it suffers from the vanishing gradient problems.
- The RNN works best when each output $\hat{y}^{\langle t \rangle}$ can be estimated using "local" context.
- "Local" context refers to information that is close to the prediction's time step t.
- More formally, local context refers to inputs $x^{\langle t' \rangle}$ and predictions $\hat{y}^{\langle t \rangle}$ where t' is close to 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

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

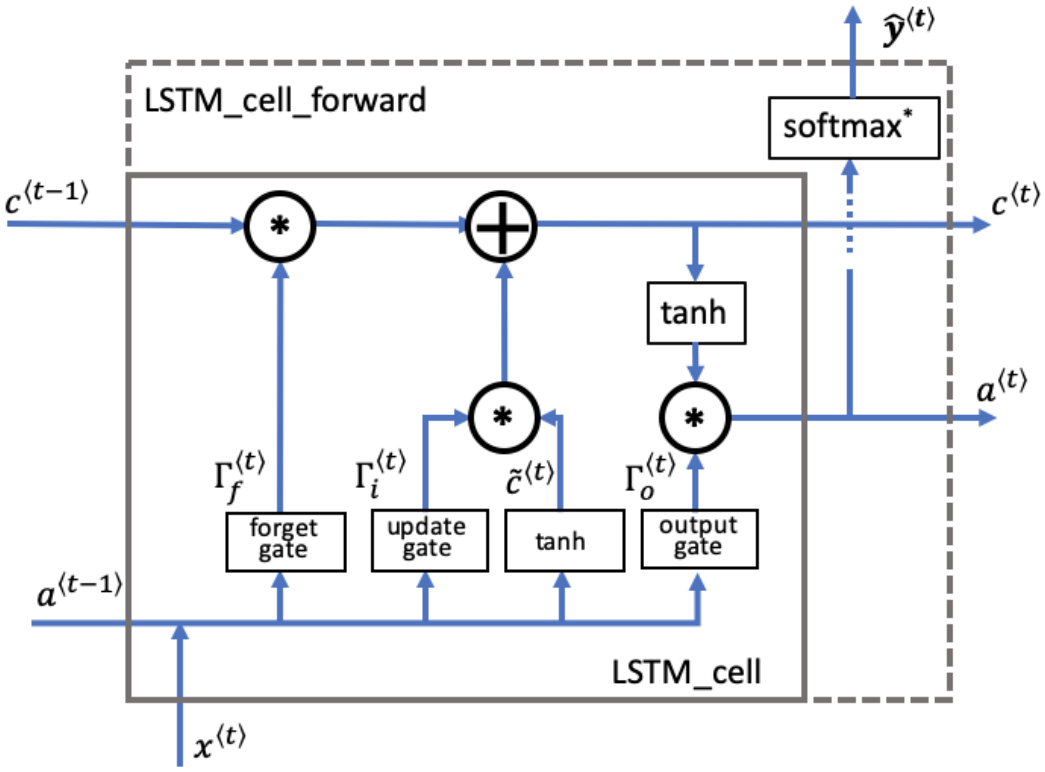

Figure 4: LSTM-cell. This tracks and updates a "cell state" or memory variable $c^{\langle t \rangle}$ at every time-step, which can be different from $a^{\langle t \rangle}$.
Note, the $softmax^*$ includes a dense layer and softmax

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.

Overview of gates and states

- Forget gate Γ_f

- Let's assume we are reading words in a piece of text, and plan to use an LSTM to keep track of grammatical structures, such as whether the subject is singular ("puppy") or plural ("puppies").
- If the subject changes its state (from a singular word to a plural word), the memory of the previous state becomes outdated, so we "forget" that outdated state.
- The "forget gate" is a tensor containing values that are between 0 and 1.
  - If a unit in the forget gate has a value close to 0, the LSTM will "forget" the stored state in the corresponding unit of the previous cell state.
  - If a unit in the forget gate has a value close to 1, the LSTM will mostly remember the corresponding value in the stored state.

Equation

$$\mathbf{\Gamma}_f^{\langle t \rangle} = \sigma(\mathbf{W}_f[\mathbf{a}^{\langle t-1 \rangle}, \mathbf{x}^{\langle t \rangle}] + \mathbf{b}_f) \tag{1}$$

Explanation of the equation:

- $\mathbf{W_f}$ contains weights that govern the forget gate's behavior.
- The previous time step's hidden state [$a^{\langle t-1 \rangle}$ and current time step's input $x^{\langle t \rangle}$] are concatenated together and multiplied by $\mathbf{W_f}$.
- A sigmoid function is used to make each of the gate tensor's values $\mathbf{\Gamma}_f^{\langle t \rangle}$ range from 0 to 1.
- The forget gate $\mathbf{\Gamma}_f^{\langle t \rangle}$ has the same dimensions as the previous cell state $c^{\langle t-1 \rangle}$.
- This means that the two can be multiplied together, element-wise.
- Multiplying the tensors $\mathbf{\Gamma}_f^{\langle t \rangle} * \mathbf{c}^{\langle t-1 \rangle}$ is like applying a mask over the previous cell state.
- If a single value in $\mathbf{\Gamma}_f^{\langle t \rangle}$ is 0 or close to 0, then the product is close to 0.
  - This keeps the information stored in the corresponding unit in $\mathbf{c}^{\langle t-1 \rangle}$ from being remembered for the next time step.
- Similarly, if one value is close to 1, the product is close to the original value in the previous cell state.
  - The LSTM will keep the information from the corresponding unit of $\mathbf{c}^{\langle t-1 \rangle}$, to be used in the next time step.

Variable names in the code

The variable names in the code are similar to the equations, with slight differences.

- `Wf`: forget gate weight $\mathbf{W}_f$
- `bf`: forget gate bias $\mathbf{b}_f$
- `ft`: forget gate $\Gamma_f^{\langle t \rangle}$

Candidate value $\tilde{\mathbf{c}}^{\langle t \rangle}$

- The candidate value is a tensor containing information from the current time step that **may** be stored in the current cell state $\mathbf{c}^{\langle t \rangle}$.
- Which parts of the candidate value get passed on depends on the update gate.

- The candidate value is a tensor containing values that range from -1 to 1.
- The tilde "~" is used to differentiate the candidate $\tilde{\mathbf{c}}^{\langle t \rangle}$ from the cell state $\mathbf{c}^{\langle t \rangle}$.

Equation

$$\tilde{\mathbf{c}}^{\langle t \rangle} = \tanh\left(\mathbf{W}_c [\mathbf{a}^{\langle t-1 \rangle}, \mathbf{x}^{\langle t \rangle}] + \mathbf{b}_c\right) \tag{3}$$

Explanation of the equation

- The 'tanh' function produces values between -1 and +1.

Variable names in the code

- `cct`: candidate value $\tilde{\mathbf{c}}^{\langle t \rangle}$

- Update gate $\mathbf{\Gamma}_i$

- We use the update gate to decide what aspects of the candidate $\tilde{\mathbf{c}}^{\langle t \rangle}$ to add to the cell state $c^{\langle t \rangle}$.
- The update gate decides what parts of a "candidate" tensor $\tilde{\mathbf{c}}^{\langle t \rangle}$ are passed onto the cell state $\mathbf{c}^{\langle t \rangle}$.
- The update gate is a tensor containing values between 0 and 1.
  - When a unit in the update gate is close to 1, it allows the value of the candidate $\tilde{\mathbf{c}}^{\langle t \rangle}$ to be passed onto the hidden state $\mathbf{c}^{\langle t \rangle}$
  - When a unit in the update gate is close to 0, it prevents the corresponding value in the candidate from being passed onto the hidden state.
- Notice that we use the subscript "i" and not "u", to follow the convention used in the literature.

Equation

$$\mathbf{\Gamma}_i^{\langle t \rangle} = \sigma(\mathbf{W}_i [a^{\langle t-1 \rangle}, \mathbf{x}^{\langle t \rangle}] + \mathbf{b}_i) \tag{2}$$

Explanation of the equation

- Similar to the forget gate, here $\mathbf{\Gamma}_i^{\langle t \rangle}$, the sigmoid produces values between 0 and 1.
- The update gate is multiplied element-wise with the candidate, and this product ($\mathbf{\Gamma}_i^{\langle t \rangle} * \tilde{c}^{\langle t \rangle}$) is used in determining the cell state $\mathbf{c}^{\langle t \rangle}$.

Variable names in code (Please note that they're different than the equations)

In the code, we'll use the variable names found in the academic literature. These variables don't use "u" to denote "update".

- `Wi` is the update gate weight $\mathbf{W}_i$ (not "Wu")
- `bi` is the update gate bias $\mathbf{b}_i$ (not "bu")
- `it` is the forget gate $\mathbf{\Gamma}_i^{\langle t \rangle}$ (not "ut")

- Cell state $\mathbf{c}^{\langle t \rangle}$

- The cell state is the "memory" that gets passed onto future time steps.
- The new cell state $\mathbf{c}^{\langle t \rangle}$ is a combination of the previous cell state and the candidate value.

Equation

$$\mathbf{c}^{\langle t \rangle} = \mathbf{\Gamma}_f^{\langle t \rangle} * \mathbf{c}^{\langle t-1 \rangle} + \mathbf{\Gamma}_i^{\langle t \rangle} * \tilde{\mathbf{c}}^{\langle t \rangle} \tag{4}$$

Explanation of equation

- The previous cell state $\mathbf{c}^{\langle t-1 \rangle}$ is adjusted (weighted) by the forget gate $\mathbf{\Gamma}_f^{\langle t \rangle}$
- and the candidate value $\tilde{\mathbf{c}}^{\langle t \rangle}$, adjusted (weighted) by the update gate $\mathbf{\Gamma}_i^{\langle t \rangle}$

Variable names and shapes in the code

- `c`: cell state, including all time steps, $\mathbf{c}$ shape (n_a, m, T)
- `c_next`: new (next) cell state, $\mathbf{c}^{\langle t \rangle}$ shape (n_a, m)
- `c_prev`: previous cell state, $\mathbf{c}^{\langle t-1 \rangle}$, shape (n_a, m)

- Output gate $\mathbf{\Gamma}_o$

- The output gate decides what gets sent as the prediction (output) of the time step.
- The output gate is like the other gates. It contains values that range from 0 to 1.

Equation

$$\mathbf{\Gamma}_o^{\langle t \rangle} = \sigma(\mathbf{W}_o[\mathbf{a}^{\langle t-1 \rangle}, \mathbf{x}^{\langle t \rangle}] + \mathbf{b}_o) \tag{5}$$

Explanation of the equation

- The output gate is determined by the previous hidden state $\mathbf{a}^{\langle t-1 \rangle}$ and the current input $\mathbf{x}^{\langle t \rangle}$
- The sigmoid makes the gate range from 0 to 1.

Variable names in the code

- `Wo`: output gate weight, $\mathbf{W_o}$
- `bo`: output gate bias, $\mathbf{b_o}$
- `ot`: output gate, $\mathbf{\Gamma}_o^{\langle t \rangle}$

- Hidden state $\mathbf{a}^{\langle t \rangle}$

- The hidden state gets passed to the LSTM cell's next time step.
- It is used to determine the three gates ($\mathbf{\Gamma}_f, \mathbf{\Gamma}_u, \mathbf{\Gamma}_o$) of the next time step.
- The hidden state is also used for the prediction $y^{\langle t \rangle}$.

Equation

$$\mathbf{a}^{\langle t \rangle} = \mathbf{\Gamma}_o^{\langle t \rangle} * \tanh(\mathbf{c}^{\langle t \rangle}) \tag{6}$$

Explanation of equation

- The hidden state $\mathbf{a}^{\langle t \rangle}$ is determined by the cell state $\mathbf{c}^{\langle t \rangle}$ in combination with the output gate $\mathbf{\Gamma}_o$.
- The cell state is passed through the "tanh" function to rescale values between -1 and +1.
- The output gate acts like a "mask" that either preserves the values of $\tanh(\mathbf{c}^{\langle t \rangle})$ or keeps those values from being included in the hidden state $\mathbf{a}^{\langle t \rangle}$

Variable names and shapes in the code

- a: hidden state, including time steps. $\mathbf{a}$ has shape (n_a, m, T_x)
- 'a_prev`: hidden state from previous time step. $\mathbf{a}^{\langle t-1 \rangle}$ has shape (n_a, m)
- a_next: hidden state for next time step. $\mathbf{a}^{\langle t \rangle}$ has shape (n_a, m)

- Prediction $\mathbf{y}_{pred}^{\langle t \rangle}$

- The prediction in this use case is a classification, so we'll use a softmax.

The equation is:

$$\mathbf{y}_{pred}^{\langle t \rangle} = \text{softmax}(\mathbf{W}_y \mathbf{a}^{\langle t \rangle} + \mathbf{b}_y)$$

Variable names and shapes in the code

- y_pred: prediction, including all time steps. $\mathbf{y}_{pred}$ has shape (n_y, m, T_x). Note that $(T_y = T_x)$ for this example.
- yt_pred: prediction for the current time step t. $\mathbf{y}_{pred}^{\langle t \rangle}$ has shape (n_y, m)

2.1 - LSTM cell

Exercise: Implement the LSTM cell described in the Figure (4).

Instructions:

1. Concatenate the hidden state $a^{\langle t-1 \rangle}$ and input $x^{\langle t \rangle}$ into a single matrix:

$$concat = \begin{bmatrix} a^{\langle t-1 \rangle} \\ x^{\langle t \rangle} \end{bmatrix}$$

1. Compute all the formulas 1 through 6 for the gates, hidden state, and cell state.
2. Compute the prediction $y^{\langle t \rangle}$.

Additional Hints

- You can use numpy.concatenate (https://docs.scipy.org/doc/numpy/reference/generated/numpy.concatenate.html). Check which value to use for the `axis` parameter.
- The functions `sigmoid()` and `softmax` are imported from `rnn_utils.py`.
- numpy.tanh (https://docs.scipy.org/doc/numpy/reference/generated/numpy.tanh.html)
- Use np.dot (https://docs.scipy.org/doc/numpy/reference/generated/numpy.dot.html) for matrix multiplication.
- Notice that the variable names `Wi`, `bi` refer to the weights and biases of the **update** gate. There are no variables named "Wu" or "bu" in this function.

In [8]:
```
# 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 Figure (

    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 array o
                        bf -- Bias of the forget gate, numpy array of shape (
                        Wi -- Weight matrix of the update gate, numpy array o
                        bi -- Bias of the update gate, numpy array of shape (
                        Wc -- Weight matrix of the first "tanh", numpy array
                        bc --  Bias of the first "tanh", numpy array of shape
                        Wo -- Weight matrix of the output gate, numpy array o
                        bo --  Bias of the output gate, numpy array of shape
                        Wy -- Weight matrix relating the hidden-state to the
                        by -- Bias relating the hidden-state to the output, n

    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,

    Note: ft/it/ot stand for the forget/update/output gates, cct stands for t
          c stands for the cell state (memory)
    """

    # Retrieve parameters from "parameters"
    Wf = parameters["Wf"] # forget gate weight
    bf = parameters["bf"]
    Wi = parameters["Wi"] # update gate weight (notice the variable name)
    bi = parameters["bi"] # (notice the variable name)
    Wc = parameters["Wc"] # candidate value weight
    bc = parameters["bc"]
    Wo = parameters["Wo"] # output gate weight
    bo = parameters["bo"]
    Wy = parameters["Wy"] # prediction weight
    by = parameters["by"]

    # Retrieve dimensions from shapes of xt and Wy
    n_x, m = xt.shape
    n_y, n_a = Wy.shape

    ### START CODE HERE ###
    # Concatenate a_prev and xt (≈1 line)
    concat = np.concatenate((a_prev, xt), axis=0)

    # Compute values for ft (forget gate), it (update gate),
    # cct (candidate value), c_next (cell state),
    # ot (output gate), a_next (hidden state) (≈6 lines)
    ft = sigmoid(np.dot(Wf, concat) + bf)        # forget gate
```

```
    it = sigmoid(np.dot(Wi, concat) + bi)        # update gate
    cct = np.tanh(np.dot(Wc, concat) + bc)       # candidate value
    c_next = ft * c_prev + it * cct              # cell state
    ot = sigmoid(np.dot(Wo, concat) + bo)        # output gate
    a_next = ot * np.tanh(c_next)                # hidden state

    # Compute prediction of the LSTM cell (≈1 line)
    yt_pred = softmax(np.dot(Wy, a_next) + by)
    ### END CODE HERE ###

    # store values needed for backward propagation in cache
    cache = (a_next, c_next, a_prev, c_prev, ft, it, cct, ot, xt, parameters)

    return a_next, c_next, yt_pred, cache
```

In [9]:
```
np.random.seed(1)
xt_tmp = np.random.randn(3,10)
a_prev_tmp = np.random.randn(5,10)
c_prev_tmp = np.random.randn(5,10)
parameters_tmp = {}
parameters_tmp['Wf'] = np.random.randn(5, 5+3)
parameters_tmp['bf'] = np.random.randn(5,1)
parameters_tmp['Wi'] = np.random.randn(5, 5+3)
parameters_tmp['bi'] = np.random.randn(5,1)
parameters_tmp['Wo'] = np.random.randn(5, 5+3)
parameters_tmp['bo'] = np.random.randn(5,1)
parameters_tmp['Wc'] = np.random.randn(5, 5+3)
parameters_tmp['bc'] = np.random.randn(5,1)
parameters_tmp['Wy'] = np.random.randn(2,5)
parameters_tmp['by'] = np.random.randn(2,1)

a_next_tmp, c_next_tmp, yt_tmp, cache_tmp = lstm_cell_forward(xt_tmp, a_prev_
print("a_next[4] = \n", a_next_tmp[4])
print("a_next.shape = ", a_next_tmp.shape)
print("c_next[2] = \n", c_next_tmp[2])
print("c_next.shape = ", c_next_tmp.shape)
print("yt[1] =", yt_tmp[1])
print("yt.shape = ", yt_tmp.shape)
print("cache[1][3] =\n", cache_tmp[1][3])
print("len(cache) = ", len(cache_tmp))
```

```
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  0.3114
6381
  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
```

Expected Output:

```
a_next[4] =
 [-0.66408471   0.0036921    0.02088357   0.22834167 -0.85575339   0.0013
8482
  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   0.1183
2942
  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 -0.3082
1874
  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.

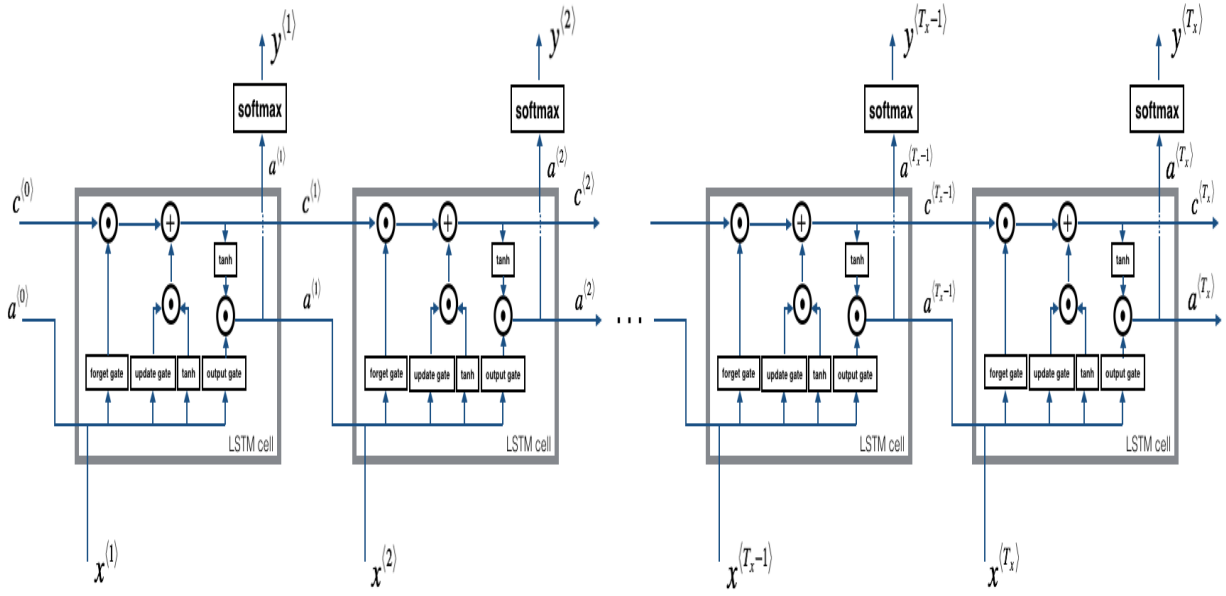

Figure 5: LSTM over multiple time-steps.

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

Instructions

- Get the dimensions n_x, n_a, n_y, m, T_x from the shape of the variables: `x` and `parameters`.

- Initialize the 3D tensors a, c and y.
    - a: hidden state, shape (n_a, m, T_x)
    - c: cell state, shape (n_a, m, T_x)
    - y: prediction, shape (n_y, m, T_x) (Note that $T_y = T_x$ in this example).
    - **Note** Setting one variable equal to the other is a "copy by reference". In other words, don't do `c = a', otherwise both these variables point to the same underlying variable.
- Initialize the 2D tensor $a^{\langle t \rangle}$
    - $a^{\langle t \rangle}$ stores the hidden state for time step t. The variable name is a_next.
    - $a^{\langle 0 \rangle}$, the initial hidden state at time step 0, is passed in when calling the function. The variable name is a0.
    - $a^{\langle t \rangle}$ and $a^{\langle 0 \rangle}$ represent a single time step, so they both have the shape (n_a, m)
    - Initialize $a^{\langle t \rangle}$ by setting it to the initial hidden state ($a^{\langle 0 \rangle}$) that is passed into the function.
- Initialize $c^{\langle t \rangle}$ with zeros.
    - The variable name is `c_next`.
    - $c^{\langle t \rangle}$ represents a single time step, so its shape is (n_a, m)
    - **Note**: create `c_next` as its own variable with its own location in memory. Do not initialize it as a slice of the 3D tensor c. In other words, **don't** do `c_next = c[:,:,0]`.
- For each time step, do the following:
    - From the 3D tensor x, get a 2D slice $x^{\langle t \rangle}$ at time step t.
    - Call the `lstm_cell_forward` function that you defined previously, to get the hidden state, cell state, prediction, and cache.
    - Store the hidden state, cell state and prediction (the 2D tensors) inside the 3D tensors.
    - Also append the cache to the list of caches.

In [10]:
```
# GRADED FUNCTION: lstm_forward

def lstm_forward(x, a0, parameters):
    """
    Implement the forward propagation of the recurrent neural network using a

    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 array o
                        bf -- Bias of the forget gate, numpy array of shape (
                        Wi -- Weight matrix of the update gate, numpy array o
                        bi -- Bias of the update gate, numpy array of shape (
                        Wc -- Weight matrix of the first "tanh", numpy array
                        bc -- Bias of the first "tanh", numpy array of shape
                        Wo -- Weight matrix of the output gate, numpy array o
                        bo -- Bias of the output gate, numpy array of shape (
                        Wy -- Weight matrix relating the hidden-state to the
                        by -- Bias relating the hidden-state to the output, n

    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, T_x)
    c -- The value of the cell state, numpy array of shape (n_a, m, T_x)
    caches -- tuple of values needed for the backward pass, contains (list of
    """

    # Initialize "caches", which will track the list of all the caches
    caches = []

    ### START CODE HERE ###
    Wy = parameters['Wy'] # saving parameters['Wy'] in a local variable in ca
    # Retrieve dimensions from shapes of x and parameters['Wy'] (≈2 lines)
    n_x, m, T_x = x.shape
    n_y, n_a = Wy.shape

    # initialize "a", "c" and "y" with zeros (≈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
    c_next = np.zeros((n_a, m))

    # loop over all time-steps
    for t in range(T_x):
        # Get the 2D slice 'xt' from the 3D input 'x' at time step 't'
        xt = x[:, :, t]
        # Update next hidden state, next memory state, compute the prediction
        a_next, c_next, yt, cache = lstm_cell_forward(xt, a_next, c_next, par
        # Save the value of the new "next" hidden state in a (≈1 line)
        a[:,:,t] = a_next
        # Save the value of the next cell state (≈1 line)
        c[:,:,t]  = c_next
```

```
        # Save the value of the prediction in y (≈1 line)
        y[:,:,t] = yt
        # Append the cache into caches (≈1 line)
        caches.append(cache)

    ### END CODE HERE ###

    # store values needed for backward propagation in cache
    caches = (caches, x)

    return a, y, c, caches
```

In [11]:
```
np.random.seed(1)
x_tmp = np.random.randn(3,10,7)
a0_tmp = np.random.randn(5,10)
parameters_tmp = {}
parameters_tmp['Wf'] = np.random.randn(5, 5+3)
parameters_tmp['bf'] = np.random.randn(5,1)
parameters_tmp['Wi'] = np.random.randn(5, 5+3)
parameters_tmp['bi']= np.random.randn(5,1)
parameters_tmp['Wo'] = np.random.randn(5, 5+3)
parameters_tmp['bo'] = np.random.randn(5,1)
parameters_tmp['Wc'] = np.random.randn(5, 5+3)
parameters_tmp['bc'] = np.random.randn(5,1)
parameters_tmp['Wy'] = np.random.randn(2,5)
parameters_tmp['by'] = np.random.randn(2,1)

a_tmp, y_tmp, c_tmp, caches_tmp = lstm_forward(x_tmp, a0_tmp, parameters_tmp)
print("a[4][3][6] = ", a_tmp[4][3][6])
print("a.shape = ", a_tmp.shape)
print("y[1][4][3] =", y_tmp[1][4][3])
print("y.shape = ", y_tmp.shape)
print("caches[1][1][1] =\n", caches_tmp[1][1][1])
print("c[1][2][1]", c_tmp[1][2][1])
print("len(caches) = ", len(caches_tmp))
```

```
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
```

Expected Output:

```
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  0.1865
6139
  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 (https://www.coursera.org/learn/neural-networks-deep-learning/lecture/0VSHe/derivatives-with-a-computation-graph) 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 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.

Note that this notebook does not implement the backward path from the Loss 'J' backwards to 'a'. This would have included the dense layer and softmax which are a part of the forward path. This is assumed to be calculated elsewhere and the result passed to rnn_backward in 'da'. It is further assumed that loss has been adjusted for batch size (m) and division by the number of examples is not required here.

This section is optional and ungraded. It is more difficult and has fewer details regarding its implementation. This section only implements key elements of the full path.

3.1 - Basic RNN backward pass

We will start by computing the backward pass for the basic RNN-cell and then in the following sections, iterate through the cells.

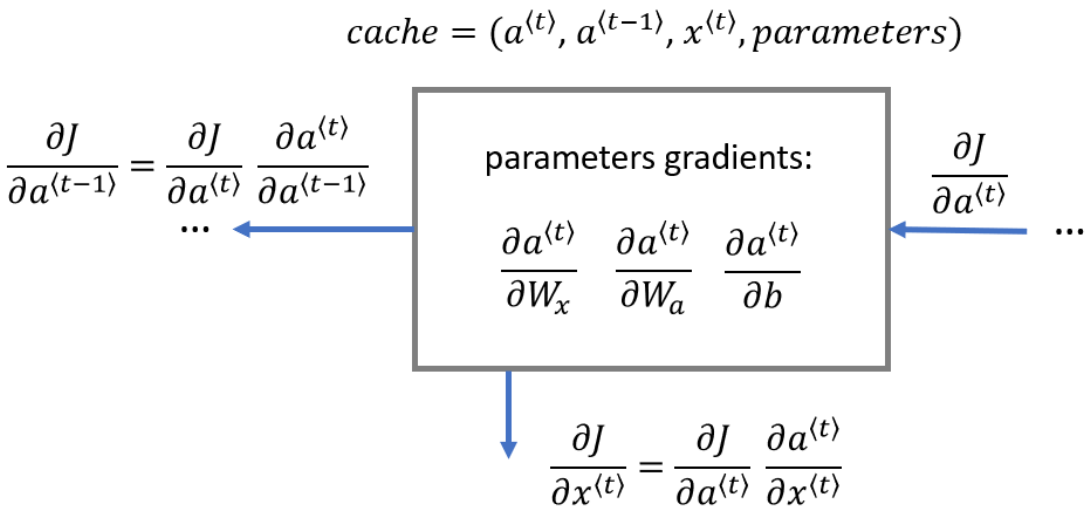

Figure 6: RNN-cell's backward pass. Just like in a fully-connected neural network, the derivative of the cost function J backpropagates through the time steps of the RNN by following the chain-rule from calculus. Internal to the cell, the chain-rule is also used to calculate $(\frac{\partial J}{\partial W_{ax}}, \frac{\partial J}{\partial W_{aa}}, \frac{\partial J}{\partial b})$ to update the parameters (W_{ax}, W_{aa}, b_a). The operation can utilize the cached results from the forward path.

Recall from lecture, the shorthand for the partial derivative of cost relative to a variable is dVariable. For example, $\frac{\partial J}{\partial W_{ax}}$ is dW_{ax}. This will be used throughout the remaining sections.

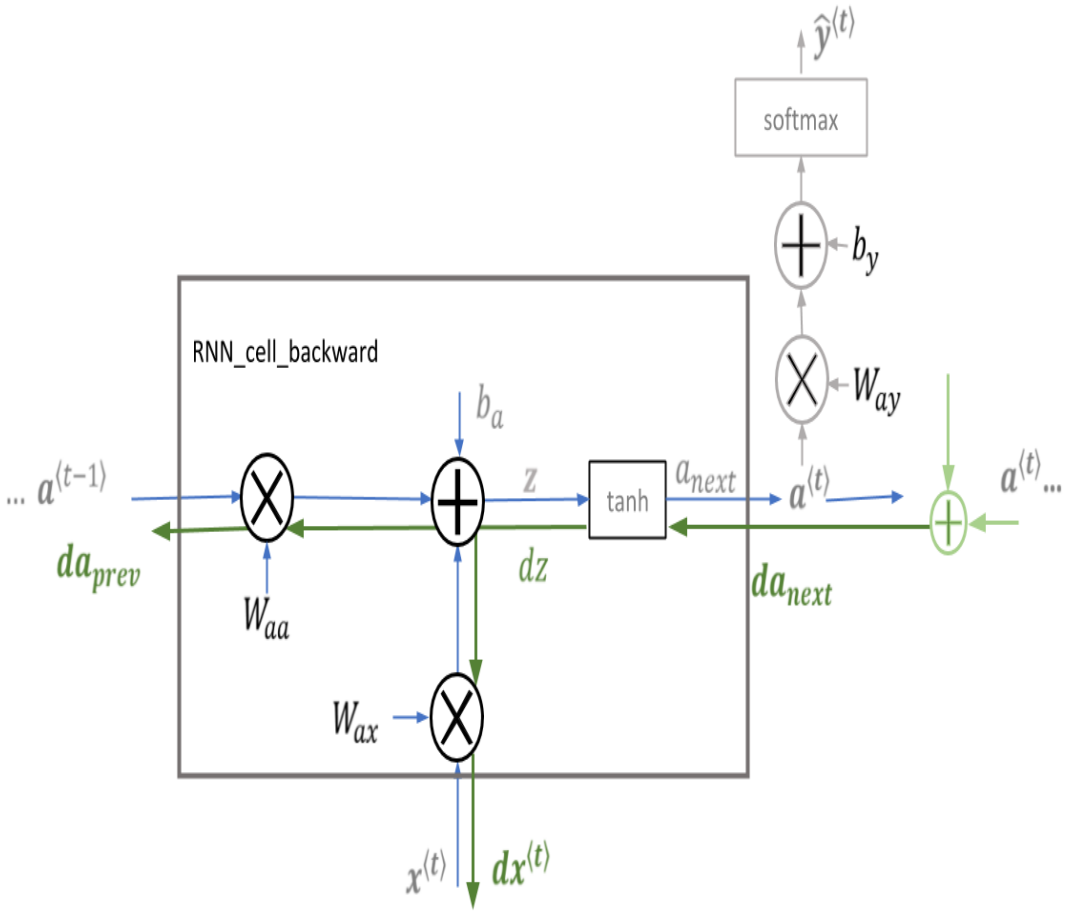

Figure 7: This implementation of rnn_cell_backward does **not** include the output dense layer and softmax which are included in rnn_cell_forward.

da_{next} is $\frac{\partial J}{\partial a^{\langle t \rangle}}$ and includes loss from previous stages and current stage output logic. The addition shown in green will be part of your implementation of rnn_backward.

Equations

To compute the rnn_cell_backward you can utilize the following equations. It is a good exercise to derive them by hand. Here, $*$ denotes element-wise multiplication while the absence of a symbol indicates matrix multiplication.

$$a^{\langle t \rangle} = \tanh(W_{ax} x^{\langle t \rangle} + W_{aa} a^{\langle t-1 \rangle} + b_a) \tag{-}$$

$$\frac{\partial \tanh(x)}{\partial x} = 1 - \tanh^2(x) \tag{-}$$

$$dW_{ax} = (da_{next} * (1 - \tanh^2(W_{ax} x^{\langle t \rangle} + W_{aa} a^{\langle t-1 \rangle} + b_a)))x^{\langle t \rangle T} \tag{1}$$

$$dW_{aa} = (da_{next} * (1 - \tanh^2(W_{ax} x^{\langle t \rangle} + W_{aa} a^{\langle t-1 \rangle} + b_a)))a^{\langle t-1 \rangle T} \tag{2}$$

$$db_a = \sum_{batch} (da_{next} * (1 - \tanh^2(W_{ax} x^{\langle t \rangle} + W_{aa} a^{\langle t-1 \rangle} + b_a))) \tag{3}$$

$$dx^{\langle t \rangle} = {W_{ax}}^T (da_{next} * (1 - \tanh^2(W_{ax} x^{\langle t \rangle} + W_{aa} a^{\langle t-1 \rangle} + b_a))) \tag{4}$$

$$da_{prev} = {W_{aa}}^T (da_{next} * (1 - \tanh^2(W_{ax} x^{\langle t \rangle} + W_{aa} a^{\langle t-1 \rangle} + b_a))) \tag{5}$$

Implementing rnn_cell_backward

The results can be computed directly by implementing the equations above. However, the above can optionally be simplified by computing 'dz' and utlilizing the chain rule.
This can be further simplified by noting that $\tanh(W_{ax} x^{\langle t \rangle} + W_{aa} a^{\langle t-1 \rangle} + b_a)$ was computed and saved in the forward pass.

To calculate dba, the 'batch' above is a sum across all 'm' examples (axis= 1). Note that you should use the keepdims = True option.

It may be worthwhile to review Course 1 Derivatives with a computational graph (https://www.coursera.org/learn/neural-networks-deep-learning/lecture/0VSHe/derivatives-with-a-computation-graph) through Backpropagation Intuition (https://www.coursera.org/learn/neural-networks-deep-learning/lecture/6dDj7/backpropagation-intuition-optional), which decompose the calculation into steps using the chain rule.
Matrix vector derivatives are described here (http://cs231n.stanford.edu/vecDerivs.pdf), though the equations above incorporate the required transformations.

Note rnn_cell_backward does **not** include the calculation of loss from $y\langle t \rangle$, this is incorporated into the incoming da_next. This is a slight mismatch with rnn_cell_forward which includes a dense layer and softmax.

Note: in the code:
$dx^{\langle t \rangle}$ is represented by dxt,
dW_{ax} is represented by dWax,
da_{prev} is represented by da_prev,
dW_{aa} is represented by dWaa,
db_a is represented by dba,
dz is not derived above but can optionally be derived by students to simplify the repeated calculations.

In [12]:
```
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_f

    Returns:
    gradients -- python dictionary containing:
                         dx -- Gradients of input data, of shape (n_x, m)
                         da_prev -- Gradients of previous hidden state, of sha
                         dWax -- Gradients of input-to-hidden weights, of shap
                         dWaa -- Gradients of hidden-to-hidden weights, of sha
                         dba -- Gradients of bias vector, of shape (n_a, 1)
    """

    # 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 the loss with respect to z (optional) (≈1 line)
    dz = None
    temp = da_next *(1 - np.power(np.tanh(np.dot(Wax, xt) + np.dot(Waa, a_pre
    # compute the gradient of the loss with respect to Wax (≈2 lines)
    dxt = np.dot(Wax.T, temp)
    dWax = np.dot(temp, xt.T)

    # compute the gradient with respect to Waa (≈2 lines)
    da_prev = np.dot(Waa.T, temp)
    dWaa = np.dot(temp, a_prev.T)

    # compute the gradient with respect to b (≈1 line)
    dba = np.sum(temp, 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,

    return gradients
```

In [13]:
```
np.random.seed(1)
xt_tmp = np.random.randn(3,10)
a_prev_tmp = np.random.randn(5,10)
parameters_tmp = {}
parameters_tmp['Wax'] = np.random.randn(5,3)
parameters_tmp['Waa'] = np.random.randn(5,5)
parameters_tmp['Wya'] = np.random.randn(2,5)
parameters_tmp['ba'] = np.random.randn(5,1)
parameters_tmp['by'] = np.random.randn(2,1)

a_next_tmp, yt_tmp, cache_tmp = rnn_cell_forward(xt_tmp, a_prev_tmp, paramete

da_next_tmp = np.random.randn(5,10)
gradients_tmp = rnn_cell_backward(da_next_tmp, cache_tmp)
print("gradients[\"dxt\"][1][2] =", gradients_tmp["dxt"][1][2])
print("gradients[\"dxt\"].shape =", gradients_tmp["dxt"].shape)
print("gradients[\"da_prev\"][2][3] =", gradients_tmp["da_prev"][2][3])
print("gradients[\"da_prev\"].shape =", gradients_tmp["da_prev"].shape)
print("gradients[\"dWax\"][3][1] =", gradients_tmp["dWax"][3][1])
print("gradients[\"dWax\"].shape =", gradients_tmp["dWax"].shape)
print("gradients[\"dWaa\"][1][2] =", gradients_tmp["dWaa"][1][2])
print("gradients[\"dWaa\"].shape =", gradients_tmp["dWaa"].shape)
print("gradients[\"dba\"][4] =", gradients_tmp["dba"][4])
print("gradients[\"dba\"].shape =", gradients_tmp["dba"].shape)
```

```
gradients["dxt"][1][2] = -1.3872130506
gradients["dxt"].shape = (3, 10)
gradients["da_prev"][2][3] = -0.152399493774
gradients["da_prev"].shape = (5, 10)
gradients["dWax"][3][1] = 0.410772824935
gradients["dWax"].shape = (5, 3)
gradients["dWaa"][1][2] = 1.15034506685
gradients["dWaa"].shape = (5, 5)
gradients["dba"][4] = [ 0.20023491]
gradients["dba"].shape = (5, 1)
```

Expected Output:

gradients["dxt"][1][2] =	-1.3872130506
gradients["dxt"].shape =	(3, 10)
gradients["da_prev"][2][3] =	-0.152399493774
gradients["da_prev"].shape =	(5, 10)
gradients["dWax"][3][1] =	0.410772824935
gradients["dWax"].shape =	(5, 3)
gradients["dWaa"][1][2] =	1.15034506685
gradients["dWaa"].shape =	(5, 5)
gradients["dba"][4] =	[0.20023491]
gradients["dba"].shape =	(5, 1)

Backward pass through the RNN

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.

- Note that this notebook does not implement the backward path from the Loss 'J' backwards to 'a'.
  - This would have included the dense layer and softmax which are a part of the forward path.
  - This is assumed to be calculated elsewhere and the result passed to rnn_backward in 'da'.
  - You must combine this with the loss from the previous stages when calling rnn_cell_backward (see figure 7 above).
- It is further assumed that loss has been adjusted for batch size (m).
  - Therefore, division by the number of examples is not required here.

In [22]:
```
def rnn_backward(da, caches):
    """
    Implement the backward pass for a RNN over an entire sequence of input da

    Arguments:
    da -- Upstream gradients of all hidden states, of shape (n_a, m, T_x)
    caches -- tuple containing information from the forward pass (rnn_forward

    Returns:
    gradients -- python dictionary containing:
                         dx -- Gradient w.r.t. the input data, numpy-array of
                         da0 -- Gradient w.r.t the initial hidden state, numpy
                         dWax -- Gradient w.r.t the input's weight matrix, num
                         dWaa -- Gradient w.r.t the hidden state's weight matr
                         dba -- Gradient w.r.t the bias, of shape (n_a, 1)
    """

    ### START CODE HERE ###

    # Retrieve values from the first cache (t=1) of caches (≈2 lines)
    #cache = (a_next, a_prev, xt, parameters)
    (caches, x) = caches
    (a1, a0, x1, parameters) = caches[0]

    # Retrieve dimensions from da's and x1's shapes (≈2 lines)
    n_a, m, T_x = da.shape
    n_x, 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(a0.shape)

    # Loop through all the time steps
    for t in reversed(range(T_x)):
        # Compute gradients at time step t.
        # Remember to sum gradients from the output path (da) and the previou
        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_p
        # Increment global derivatives w.r.t parameters by adding their deriv
        dx[:, :, t] = dxt
        dWax += dWaxt
        dWaa += dWaat
        dba += dbat

    # Set da0 to the gradient of a which has been backpropagated through all
    da0 = da_prevt
    ### END CODE HERE ###

    # Store the gradients in a python dictionary
    gradients = {"dx": dx, "da0": da0, "dWax": dWax, "dWaa": dWaa,"dba": dba}
```

```
    return gradients
```

In [23]:
```
np.random.seed(1)
x_tmp = np.random.randn(3,10,4)
a0_tmp = np.random.randn(5,10)
parameters_tmp = {}
parameters_tmp['Wax'] = np.random.randn(5,3)
parameters_tmp['Waa'] = np.random.randn(5,5)
parameters_tmp['Wya'] = np.random.randn(2,5)
parameters_tmp['ba'] = np.random.randn(5,1)
parameters_tmp['by'] = np.random.randn(2,1)

a_tmp, y_tmp, caches_tmp = rnn_forward(x_tmp, a0_tmp, parameters_tmp)
da_tmp = np.random.randn(5, 10, 4)
gradients_tmp = rnn_backward(da_tmp, caches_tmp)

print("gradients[\"dx\"][1][2] =", gradients_tmp["dx"][1][2])
print("gradients[\"dx\"].shape =", gradients_tmp["dx"].shape)
print("gradients[\"da0\"][2][3] =", gradients_tmp["da0"][2][3])
print("gradients[\"da0\"].shape =", gradients_tmp["da0"].shape)
print("gradients[\"dWax\"][3][1] =", gradients_tmp["dWax"][3][1])
print("gradients[\"dWax\"].shape =", gradients_tmp["dWax"].shape)
print("gradients[\"dWaa\"][1][2] =", gradients_tmp["dWaa"][1][2])
print("gradients[\"dWaa\"].shape =", gradients_tmp["dWaa"].shape)
print("gradients[\"dba\"][4] =", gradients_tmp["dba"][4])
print("gradients[\"dba\"].shape =", gradients_tmp["dba"].shape)
```

```
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)
```

Expected Output:

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 pass.

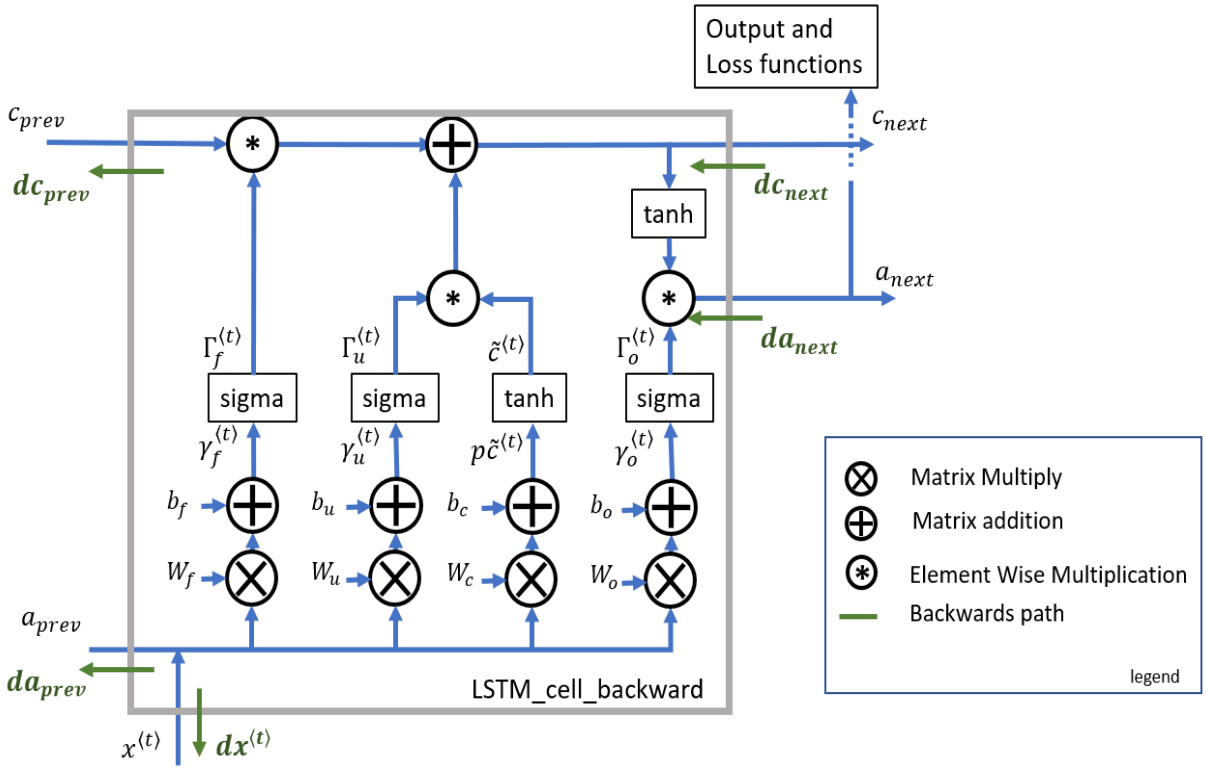

Figure 8: lstm_cell_backward. Note the output functions, while part of the lstm_cell_forward, are not included in lstm_cell_backward

The equations for the LSTM backward pass are provided below. (If you enjoy calculus exercises feel free to try deriving these from scratch yourself.)

3.2.2 gate derivatives

Note the location of the gate derivatives (γ..) between the dense layer and the activation function (see graphic above). This is convenient for computing parameter derivatives in the next step.

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

$$dp\widetilde{c}^{\langle t \rangle} = \left(dc_{next} * \Gamma_u^{\langle t \rangle} + \Gamma_o^{\langle t \rangle} * (1 - \tanh^2(c_{next})) * \Gamma_u^{\langle t \rangle} * da_{next}\right) * \left(1 - \left(\widetilde{c}^{\langle t \rangle}\right)^2\right)$$

$$d\gamma_u^{\langle t \rangle} = \left(dc_{next} * \widetilde{c}^{\langle t \rangle} + \Gamma_o^{\langle t \rangle} * (1 - \tanh^2(c_{next})) * \widetilde{c}^{\langle t \rangle} * da_{next}\right) * \Gamma_u^{\langle t \rangle} * \left(1 - \Gamma_u^{\langle t \rangle}\right)$$

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

3.2.3 parameter derivatives

$$dW_f = d\gamma_f^{\langle t \rangle} \begin{bmatrix} a_{prev} \\ x_t \end{bmatrix}^T \tag{11}$$

$$dW_u = d\gamma_u^{\langle t \rangle} \begin{bmatrix} a_{prev} \\ x_t \end{bmatrix}^T \tag{12}$$

$$dW_c = dp\widetilde{c}^{\langle t \rangle} \begin{bmatrix} a_{prev} \\ x_t \end{bmatrix}^T \tag{13}$$

$$dW_o = d\gamma_o^{\langle t \rangle} \begin{bmatrix} a_{prev} \\ x_t \end{bmatrix}^T \tag{14}$$

To calculate db_f, db_u, db_c, db_o you just need to sum across all 'm' examples (axis= 1) on $d\gamma_f^{\langle t \rangle}, d\gamma_u^{\langle t \rangle}, dp\widetilde{c}^{\langle t \rangle}, d\gamma_o^{\langle t \rangle}$ respectively. Note that you should have the `keepdims = True` option.

$$db_f = \sum_{batch} d\gamma_f^{\langle t \rangle} \tag{15}$$

$$db_u = \sum_{batch} d\gamma_u^{\langle t \rangle} \tag{16}$$

$$db_c = \sum_{batch} d\gamma_c^{\langle t \rangle} \tag{17}$$

$$db_o = \sum_{batch} d\gamma_o^{\langle t \rangle} \tag{18}$$

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

$$da_{prev} = W_f^T d\gamma_f^{\langle t \rangle} + W_u^T d\gamma_u^{\langle t \rangle} + W_c^T dp\widetilde{c}^{\langle t \rangle} + W_o^T d\gamma_o^{\langle t \rangle} \tag{19}$$

Here, to account for concatenation, the weights for equations 19 are the first n_a, (i.e. $W_f = W_f[:, :n_a]$ etc...)

$$dc_{prev} = dc_{next} * \Gamma_f^{\langle t \rangle} + \Gamma_o^{\langle t \rangle} * (1 - \tanh^2(c_{next})) * \Gamma_f^{\langle t \rangle} * da_{next} \tag{20}$$

$$dx^{\langle t \rangle} = W_f^T d\gamma_f^{\langle t \rangle} + W_u^T d\gamma_u^{\langle t \rangle} + W_c^T dp\widetilde{c}^{\langle t \rangle} + W_o^T d\gamma_o^{\langle t \rangle} \tag{21}$$

where the weights for equation 21 are from n_a to the end, (i.e. $W_f = W_f[:, n_a :]$ etc...)

Exercise: Implement `lstm_cell_backward` by implementing equations $7 - 21$ below.

Note: In the code:

$d\gamma_o^{\langle t \rangle}$ is represented by `dot`,
$dp\widetilde{c}^{\langle t \rangle}$ is represented by `dcct`,
$d\gamma_u^{\langle t \rangle}$ is represented by `dit`,
$d\gamma_f^{\langle t \rangle}$ is represented by `dft`

In [44]:
```
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, of shap
                         da_prev -- Gradient w.r.t. the previous hidden state,
                         dc_prev -- Gradient w.r.t. the previous memory state,
                         dWf -- Gradient w.r.t. the weight matrix of the forge
                         dWi -- Gradient w.r.t. the weight matrix of the updat
                         dWc -- Gradient w.r.t. the weight matrix of the memor
                         dWo -- Gradient w.r.t. the weight matrix of the outpu
                         dbf -- Gradient w.r.t. biases of the forget gate, of
                         dbi -- Gradient w.r.t. biases of the update gate, of
                         dbc -- Gradient w.r.t. biases of the memory gate, of
                         dbo -- Gradient w.r.t. biases of the output gate, of
    """

    # 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_x, m = None
#    n_a, m = None
    n_x, m = xt.shape
    n_a, m = a_next.shape

    # Compute gates related derivatives, you can find their values can be fou
#    dot = None
#    dcct = None
#    dit = None
#    dft = None
    dot = da_next * np.tanh(c_next) * ot * (1 - ot)
    dcct = (da_next * ot * (1 - np.tanh(c_next) ** 2) + dc_next) * it * (1 -
    dit = (da_next * ot * (1 - np.tanh(c_next) ** 2) + dc_next) * cct * (1 -
    dft = (da_next * ot * (1 - np.tanh(c_next) ** 2) + dc_next) * c_prev * ft

    # Compute parameters related derivatives. Use equations (11)-(18) (≈8 lin
    """    dWf = None
    dWi = None
    dWc = None
    dWo = None
    dbf = None
    dbi = None
    dbc = None
    dbo = None
    """
    dWf = np.dot(dft, np.hstack([a_prev.T, xt.T]))
    dWi = np.dot(dit, np.hstack([a_prev.T, xt.T]))
```

```
    dWc = np.dot(dcct, np.hstack([a_prev.T, xt.T]))
    dWo = np.dot(dot, np.hstack([a_prev.T, xt.T]))
    dbf = np.sum(dft, axis=1, keepdims=True)
    dbi = np.sum(dit, axis=1, keepdims=True)
    dbc = np.sum(dcct, axis=1, keepdims=True)
    dbo = np.sum(dot, axis=1, keepdims=True)

    # Compute derivatives w.r.t previous hidden state, previous memory state
    """    da_prev = None
    dc_prev = None
    dxt = None
    """
    da_prev = np.dot(parameters_tmp['Wf'][:, :n_a].T, dft) + np.dot(parameter
    dc_prev = (da_next * ot * (1 - np.tanh(c_next) ** 2) + dc_next) * ft
    dxt = np.dot(parameters_tmp['Wf'][:, n_a:].T, dft) + np.dot(parameters_tm
    ### END CODE HERE ###

    # Save gradients in dictionary
    gradients = {"dxt": dxt, "da_prev": da_prev, "dc_prev": dc_prev, "dWf": d
                "dWc": dWc,"dbc": dbc, "dWo": dWo,"dbo": dbo}

    return gradients
```

In [45]:
```
np.random.seed(1)
xt_tmp = np.random.randn(3,10)
a_prev_tmp = np.random.randn(5,10)
c_prev_tmp = np.random.randn(5,10)
parameters_tmp = {}
parameters_tmp['Wf'] = np.random.randn(5, 5+3)
parameters_tmp['bf'] = np.random.randn(5,1)
parameters_tmp['Wi'] = np.random.randn(5, 5+3)
parameters_tmp['bi'] = np.random.randn(5,1)
parameters_tmp['Wo'] = np.random.randn(5, 5+3)
parameters_tmp['bo'] = np.random.randn(5,1)
parameters_tmp['Wc'] = np.random.randn(5, 5+3)
parameters_tmp['bc'] = np.random.randn(5,1)
parameters_tmp['Wy'] = np.random.randn(2,5)
parameters_tmp['by'] = np.random.randn(2,1)

a_next_tmp, c_next_tmp, yt_tmp, cache_tmp = lstm_cell_forward(xt_tmp, a_prev_

da_next_tmp = np.random.randn(5,10)
dc_next_tmp = np.random.randn(5,10)
gradients_tmp = lstm_cell_backward(da_next_tmp, dc_next_tmp, cache_tmp)
print("gradients[\"dxt\"][1][2] =", gradients_tmp["dxt"][1][2])
print("gradients[\"dxt\"].shape =", gradients_tmp["dxt"].shape)
print("gradients[\"da_prev\"][2][3] =", gradients_tmp["da_prev"][2][3])
print("gradients[\"da_prev\"].shape =", gradients_tmp["da_prev"].shape)
print("gradients[\"dc_prev\"][2][3] =", gradients_tmp["dc_prev"][2][3])
print("gradients[\"dc_prev\"].shape =", gradients_tmp["dc_prev"].shape)
print("gradients[\"dWf\"][3][1] =", gradients_tmp["dWf"][3][1])
print("gradients[\"dWf\"].shape =", gradients_tmp["dWf"].shape)
print("gradients[\"dWi\"][1][2] =", gradients_tmp["dWi"][1][2])
print("gradients[\"dWi\"].shape =", gradients_tmp["dWi"].shape)
print("gradients[\"dWc\"][3][1] =", gradients_tmp["dWc"][3][1])
print("gradients[\"dWc\"].shape =", gradients_tmp["dWc"].shape)
print("gradients[\"dWo\"][1][2] =", gradients_tmp["dWo"][1][2])
print("gradients[\"dWo\"].shape =", gradients_tmp["dWo"].shape)
print("gradients[\"dbf\"][4] =", gradients_tmp["dbf"][4])
print("gradients[\"dbf\"].shape =", gradients_tmp["dbf"].shape)
print("gradients[\"dbi\"][4] =", gradients_tmp["dbi"][4])
print("gradients[\"dbi\"].shape =", gradients_tmp["dbi"].shape)
print("gradients[\"dbc\"][4] =", gradients_tmp["dbc"][4])
print("gradients[\"dbc\"].shape =", gradients_tmp["dbc"].shape)
print("gradients[\"dbo\"][4] =", gradients_tmp["dbo"][4])
print("gradients[\"dbo\"].shape =", gradients_tmp["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)
```

Expected Output:

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 [50]:
```
def lstm_backward(da, caches):

    """
    Implement the backward pass for the RNN with LSTM-cell (over a whole sequ

    Arguments:
    da -- Gradients w.r.t the hidden states, numpy-array of shape (n_a, m, T_
    caches -- cache storing information from the forward pass (lstm_forward)

    Returns:
    gradients -- python dictionary containing:
                         dx -- Gradient of inputs, of shape (n_x, m, T_x)
                         da0 -- Gradient w.r.t. the previous hidden state, num
                         dWf -- Gradient w.r.t. the weight matrix of the forge
                         dWi -- Gradient w.r.t. the weight matrix of the updat
                         dWc -- Gradient w.r.t. the weight matrix of the memor
                         dWo -- Gradient w.r.t. the weight matrix of the save
                         dbf -- Gradient w.r.t. biases of the forget gate, of
                         dbi -- Gradient w.r.t. biases of the update gate, of
                         dbc -- Gradient w.r.t. biases of the memory gate, of
                         dbo -- Gradient w.r.t. biases of the save gate, of sh
    """

    # 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 (≈2 lines)
#    n_a, m, T_x = None
#    n_x, m = None
    n_a, m, T_x = da.shape
    n_x, m = x1.shape

    # initialize the gradients with the right sizes (≈12 lines)
    """dx = None
    da0 = None
    da_prevt = None
    dc_prevt = None
    dWf = None
    dWi = None
    dWc = None
    dWo = None
    dbf = None
    dbi = None
    dbc = None
    dbo = None """
    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))
    dbf = np.zeros((n_a, 1))
```

```
    dbi = np.zeros((n_a, 1))
    dbc = np.zeros((n_a, 1))
    dbo = np.zeros((n_a, 1))

    # loop back over the whole sequence
#    for t in reversed(range(None)):
    for t in reversed(range(T_x)):
        # Compute all gradients using lstm_cell_backward
        #gradients = None
        gradients = lstm_cell_backward(da[:,:,t] + da_prevt, dc_prevt, caches
        # Store or add the gradient to the parameters' previous step's gradie
        """da_prevt = None
        dc_prevt = None
        dx[:,:,t] = None
        dWf += None
        dWi += None
        dWc += None
        dWo += None
        dbf += None
        dbi += None
        dbc += None
        dbo += None """
        da_prevt = gradients["da_prev"]
        dc_prevt = gradients["dc_prev"]
        dx[:,:,t] = gradients["dxt"]
        dWf += gradients["dWf"]
        dWi += gradients["dWi"]
        dWc += gradients["dWc"]
        dWo += gradients["dWo"]
        dbf += gradients["dbf"]
        dbi += gradients["dbi"]
        dbc += gradients["dbc"]
        dbo += gradients["dbo"]

    # Set the first activation's gradient to the backpropagated gradient da_p
#    da0 = None
    da0 = gradients["da_prev"]

    ### END CODE HERE ###

    # Store the gradients in a python dictionary
    gradients = {"dx": dx, "da0": da0, "dWf": dWf,"dbf": dbf, "dWi": dWi,"dbi
                "dWc": dWc,"dbc": dbc, "dWo": dWo,"dbo": dbo}

    return gradients
```

In [51]:
```
np.random.seed(1)
x_tmp = np.random.randn(3,10,7)
a0_tmp = np.random.randn(5,10)

parameters_tmp = {}
parameters_tmp['Wf'] = np.random.randn(5, 5+3)
parameters_tmp['bf'] = np.random.randn(5,1)
parameters_tmp['Wi'] = np.random.randn(5, 5+3)
parameters_tmp['bi'] = np.random.randn(5,1)
parameters_tmp['Wo'] = np.random.randn(5, 5+3)
parameters_tmp['bo'] = np.random.randn(5,1)
parameters_tmp['Wc'] = np.random.randn(5, 5+3)
parameters_tmp['bc'] = np.random.randn(5,1)
parameters_tmp['Wy'] = np.zeros((2,5))       # unused, but needed for lstm_fo
parameters_tmp['by'] = np.zeros((2,1))       # unused, but needed for lstm_fo

a_tmp, y_tmp, c_tmp, caches_tmp = lstm_forward(x_tmp, a0_tmp, parameters_tmp)

da_tmp = np.random.randn(5, 10, 4)
gradients_tmp = lstm_backward(da_tmp, caches_tmp)

print("gradients[\"dx\"][1][2] =", gradients_tmp["dx"][1][2])
print("gradients[\"dx\"].shape =", gradients_tmp["dx"].shape)
print("gradients[\"da0\"][2][3] =", gradients_tmp["da0"][2][3])
print("gradients[\"da0\"].shape =", gradients_tmp["da0"].shape)
print("gradients[\"dWf\"][3][1] =", gradients_tmp["dWf"][3][1])
print("gradients[\"dWf\"].shape =", gradients_tmp["dWf"].shape)
print("gradients[\"dWi\"][1][2] =", gradients_tmp["dWi"][1][2])
print("gradients[\"dWi\"].shape =", gradients_tmp["dWi"].shape)
print("gradients[\"dWc\"][3][1] =", gradients_tmp["dWc"][3][1])
print("gradients[\"dWc\"].shape =", gradients_tmp["dWc"].shape)
print("gradients[\"dWo\"][1][2] =", gradients_tmp["dWo"][1][2])
print("gradients[\"dWo\"].shape =", gradients_tmp["dWo"].shape)
print("gradients[\"dbf\"][4] =", gradients_tmp["dbf"][4])
print("gradients[\"dbf\"].shape =", gradients_tmp["dbf"].shape)
print("gradients[\"dbi\"][4] =", gradients_tmp["dbi"][4])
print("gradients[\"dbi\"].shape =", gradients_tmp["dbi"].shape)
print("gradients[\"dbc\"][4] =", gradients_tmp["dbc"][4])
print("gradients[\"dbc\"].shape =", gradients_tmp["dbc"].shape)
print("gradients[\"dbo\"][4] =", gradients_tmp["dbo"][4])
print("gradients[\"dbo\"].shape =", gradients_tmp["dbo"].shape)
```

```
gradients["dx"][1][2] = [ 0.00218254  0.28205375 -0.48292508 -0.43281115]
gradients["dx"].shape = (3, 10, 4)
gradients["da0"][2][3] = 0.312770310257
gradients["da0"].shape = (5, 10)
gradients["dWf"][3][1] = -0.0809802310938
gradients["dWf"].shape = (5, 8)
gradients["dWi"][1][2] = 0.40512433093
gradients["dWi"].shape = (5, 8)
gradients["dWc"][3][1] = -0.0793746735512
gradients["dWc"].shape = (5, 8)
gradients["dWo"][1][2] = 0.038948775763
gradients["dWo"].shape = (5, 8)
gradients["dbf"][4] = [-0.15745657]
gradients["dbf"].shape = (5, 1)
```

```
gradients["dbi"][4] = [-0.50848333]
gradients["dbi"].shape = (5, 1)
gradients["dbc"][4] = [-0.42510818]
gradients["dbc"].shape = (5, 1)
gradients["dbo"][4] = [-0.17958196]
gradients["dbo"].shape = (5, 1)
```

Expected Output:

gradients["dx"][1][2] =	[0.00218254 0.28205375 -0.48292508 -0.43281115]
gradients["dx"].shape =	(3, 10, 4)
gradients["da0"][2][3] =	0.312770310257
gradients["da0"].shape =	(5, 10)
gradients["dWf"][3][1] =	-0.0809802310938
gradients["dWf"].shape =	(5, 8)
gradients["dWi"][1][2] =	0.40512433093
gradients["dWi"].shape =	(5, 8)
gradients["dWc"][3][1] =	-0.0793746735512
gradients["dWc"].shape =	(5, 8)
gradients["dWo"][1][2] =	0.038948775763
gradients["dWo"].shape =	(5, 8)
gradients["dbf"][4] =	[-0.15745657]
gradients["dbf"].shape =	(5, 1)
gradients["dbi"][4] =	[-0.50848333]
gradients["dbi"].shape =	(5, 1)
gradients["dbc"][4] =	[-0.42510818]
gradients["dbc"].shape =	(5, 1)
gradients["dbo"][4] =	[-0.17958196]
gradients["dbo"].shape =	(5, 1)

Congratulations !

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

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