50.007 Machine Learning Fall 2020

Design Project

Report Implementation Approach & Results

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Emission parameter estimation using maximum likelihood estimation (MLE)

Looking at the training data provided, each line in the file consists of a token (ie. word), followed by a space, then finally its tag. A single empty line serves to separate sentences in the training data.

To learn the emission parameter from training data for first-order Hidden Markov Model (HMM), first, we wrote a function <code>count_emission</code> to tabularise the number of occurrences of each word emitted by a specific state.

For each line in the file containing a word and its tag, we store the number of occurrences a specific observation (ie. word) in a dictionary as a key-value pair, with key as the word, and value as the number of occurrences. This forms the a dictionary for a specific state (ie. tag).

```
{key = observation: value = count}
```

We then store the dictionary of each state as the value of another dictionary, where the key represents the tag. As a result, count_emission function returns emission_tracker, which is nested dictionary in this format:

This would hence allow us to easily retrieve values required to calculate the emission parameters, using the function <code>emission_para</code>:

$$e(x|y) = \frac{Count(y \to x)}{Count(y)}$$

Based on the dictionary obtained that tabularises the emission counts above, we can easily obtain $Count(y \to x)$ using <code>emission_tracker[y][x]</code>, and also obtaining Count(y) using <code>sum(emission_tracker[y].values())</code>.

Emission parameter estimation with special token

In a modified version of computing emission parameters, the function <code>emission_para_token</code> takes into account words that appear in the test set, but do not appear in the training set. Similarly, this function returns the emission parameters.

$$e(x|y) = \left\{ \begin{array}{ll} \frac{\text{Count}(y \to x)}{\text{Count}(y) + k} & \text{If the word token } x \text{ appears in the training set} \\ \frac{k}{\text{Count}(y) + k} & \text{If word token } x \text{ is the special token } \#\text{UNK} \# \end{array} \right.$$

First, we do a conditional if-else check on the observation - whether the observation exists in the training data. If the observation did not appear in the training data, a special token, #UNK#, is assigned to the observation. The remaining steps remain the same as MLE estimation of emission parameter without special token above. With k=0.5, we can easily obtain emission parameters.

Simple system that produces tag y

A function tag_producer was implemented to take in a sentence as a list object, with each word being a string type element of the list, and return a list of predicted tags y for the sentence.

$$y^* = \operatorname*{arg\,max}_y e(x|y)$$

For each word in a sentence, we loop through all the states that the word can take and calculate the respective emission probabilities. We then return the state that gives the highest emission probability as the predicted tag. The individual predicted tags for each word are then stored in a list containing all predicted tags of a sentence, which is returned by the function.

Results

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Using the evaluation script given, our results are as follow:

EN #Entity in gold data: 13179 #Entity in prediction: 18650 #Correct Entity: 9542 Entity precision: 0.5116 Entity recall: 0.7240 Entity F: 0.5996 #Correct Sentiment: 8456 Sentiment precision: 0.4534 Sentiment recall: 0.6416 Sentiment F: 0.5313 ==== CN #Entity in gold data: 700 #Entity in prediction: 4248 #Correct Entity: 345 Entity precision: 0.0812 Entity recall: 0.4929 Entity F: 0.1395 #Correct Sentiment: 167 Sentiment precision: 0.0393 Sentiment recall: 0.2386 Sentiment F: 0.0675 ----SG #Entity in gold data: 4301 #Entity in prediction: 12237 #Correct Entity: 2386 Entity precision: 0.1950 Entity recall: 0.5548 Entity F: 0.2885 #Correct Sentiment: 1531 Sentiment precision: 0.1251 Sentiment recall: 0.3560 Sentiment F: 0.1851

Transition parameter estimation using maximum likelihood estimation (MLE)

To learn the transition parameter from training data for first-order Hidden Markov Model (HMM), first, we wrote a function $count_transition$ to tabularise the number of occurrences of transiting from a particulate state to another particular state. Here, we define our transition from $state\ u$ to $state\ v$.

For each line in the file containing a word and its tag, we store the number of occurrences of transitioning from one state to another in a nested dictionary format - the key represents $state\ u$, and the value is a dictionary where key represents $state\ v$ and the value is the number of occurrences.

As a result, count_transition function returns transition_tracker, which is a nested dictionary in this format:

This would hence allow us to easily retrieve values required to calculate the transition parameters, using the function transition para:

$$q(y_i|y_{i-1}) = \frac{Count(y_{i-1},y_i)}{Count(y_{i-1})}$$

Based on the dictionary obtained that tabularises the transition counts above, we can easily obtain $Count(y_{i-1}, y_i)$ using transition_tracker[state_u][state_v], and also obtaining $Count(y_{i-1})$ using sum(transition tracker[state u].values()).

Viterbi Algorithm

Using the functions we have implemented in Parts 2 and 3, we implement the viterbi algorithm to return the best path.

$$a_{u,v} = q(y_v|y_u)$$

$$b_v(x_j) = e(x_j|y_j)$$

Keeping in mind that the natural logarithm function is strictly increasing, all values are calculated in natural logarithmic form in our implementation, replacing multiplication by addition. This is based on the logarithmic property ln(ab) = ln(a) + ln(b). The purpose of taking the natural logarithm of probabilities is to reduce the risk of potential numerical underflow. However, one new issue with natural logarithmic calculations is that $ln\ 0 = -\infty$. To deal with probability values that are equal to 0, we simply set $ln\ 0 = C$, where C is a large negative number.

In addition, in view of the fact that the training data is split into sentences as indicated by a single empty line, the base case of $\pi(0, v) = 1$ if v = START, 0 otherwise is already implicitly accounted for. Furthermore, $\ln 1 = 0$. Hence, our implementation's base case is to determine the score at the position of the first word in a sentence, which is not dependent on the START state position.

The score of the best path at each position/time step are stored in a dictionary in such format:

We store the previous state, $state\ u$, as a tuple with the highest score of the path from START to $state\ v$ is for the purpose of backtracking.

At every position/time-step, we calculate the transition and emission probabilities, then take their natural logarithmic values as the score that we are tracking on. Prior to calculating the emission probabilities, we check whether the observation has appeared in the training set or not. If the observation was not seen in the training set, we assign the special token #UNK#.

Moving forward recursively:

We begin by initialising the first word of a sentence, ie. calculating $\pi(1, v)$:

```
\pi(1, v) = ln[a_{START, v} \cdot b_{v}(x_{1})]
= ln(a_{START, v}) + ln([b_{v}(x_{1})])
= ln(1) + ln[b_{v}(x_{1})]
= ln[b_{v}(x_{1})]
```

From the second word to the last word of a sentence, we calculate all the possible paths and keep track of the state that gives the maximum score:

$$\pi(k, v) = \max_{v} \left\{ \pi(k-1, u) + \ln(a_{u,v}) + \ln[b_{v}(x_{k})] \right\}$$

In the final step, we calculate the score of the overall path from START state to STOP state. ie. calculate $\pi(k+1,STOP) = max_v \{\pi(k,u) + ln(a_{uSTOP})\}$.

After calculating the best path from START state to STOP state, ie. the best path for the entire sentence, we then do backtracking to retrieve the states (ie. tags) at each position/time step. During the forward recursive process, we stored the state of position k-1 that gave the maximum score as a tuple along with the score at position k, hence we now retrieve all the states from each tuple to produce the full tag sequence for a sentence in the backtracking process.

Here are our results from our viterbi algorithm:

EN

#Entity in gold data: 13179 #Entity in prediction: 13375

#Correct Entity: 10791
Entity precision: 0.8068
Entity recall: 0.8188
Entity F: 0.8128

#Correct Sentiment : 10268
Sentiment precision: 0.7677
Sentiment recall: 0.7791
Sentiment F: 0.7734
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CN

#Entity in gold data: 700 #Entity in prediction: 850

#Correct Entity: 208
Entity precision: 0.2447
Entity recall: 0.2971
Entity F: 0.2684

#Correct Sentiment : 122
Sentiment precision: 0.1435
Sentiment recall: 0.1743
Sentiment F: 0.1574
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SG

#Entity in gold data: 4301 #Entity in prediction: 4274

#Correct Entity : 2212
Entity precision: 0.5175
Entity recall: 0.5143
Entity F: 0.5159

#Correct Sentiment: 1821
Sentiment precision: 0.4261
Sentiment recall: 0.4234
Sentiment F: 0.4247

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Modified viterbi algorithm to obtain the third best output sequence

To obtain the third best output sequence, we find the top three best paths and store these values at each position/time-step. For node i at position k, we obtain all the possible paths from all nodes in position k-1 and calculate the scores of these paths. We then keep track of the scores of the best three paths, along with the path from START up till position k, passing through the node that gives one of the best three scores at position k-1. This process is repeated for all nodes at position k to obtain the best three scores at each node. One numerical example would be - given position k has j nodes, we will keep track of $k \cdot j$ number of best paths and scores.

Similar to the Viterbi algorithm to get the best output sequence that was implemented in part 3, the base case of $\pi(0, v) = 1$ if v = START, 0 otherwise is already implicitly accounted for, as $\ln 1 = 0$. Hence, our implementation's base case is to determine the score at the position of the first word in a sentence, which is not dependent on the START state position.

At each position/time step, the scores of the best three paths for each node are stored. All the scores across all positions are stored in the scores dictionary in such format:

We store the previous state, $state\ u$, as a tuple with the highest score of the path from START to $state\ v$ is for the purpose of backtracking.

At every position/time-step, we calculate the transition and emission probabilities, then take their natural logarithmic values as the score that we are tracking on. Prior to calculating the emission probabilities, we check whether the observation has appeared in the training set or not. If the observation was not seen in the training set, we assign the special token #UNK#.

The forward process to calculate and determine the best three paths described above is split into three portions:

First, we initialise the first word of a sentence, ie. calculating $\pi(1, \nu)$. There is no dependent score from the previous position, which is the START state, because of ln(1) = 0.

```
\pi(1, v) = ln[a_{START, v} \bullet b_{v}(x_{1})]
= ln(a_{START, v}) + ln([b_{v}(x_{1})])
= ln(1) + ln[b_{v}(x_{1})]
= ln[b_{v}(x_{1})]
```

From the second word to the last word of a sentence, we calculate all possible paths and keep track of the three states that give the top three best paths.

ie. for each node v at position k, we calculate all possible values of $\pi(k, v)$.

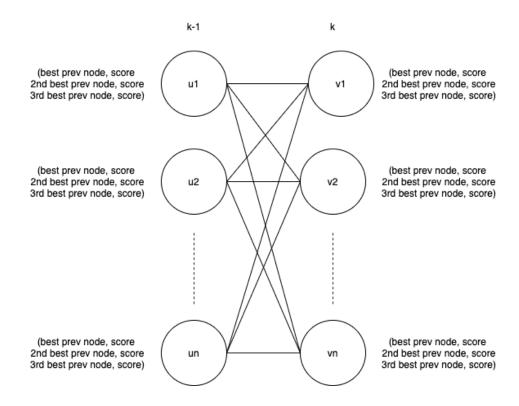
The calculation of $\pi(k, v)$ is based on $\pi(k-1, u) + ln(a_{u,v}) + ln[b_v(x_k)]$, where node u is all the possible states in position k-1. We then obtain the top three best scores for each node v. We store these three scores and its corresponding preceding node u as a tuple in the scores dictionary as described above.

In the final step, we calculate the score of the best 3 paths for each node at position n, where n is the length of a sentence, to the STOP state.

ie. calculate $\pi(n+1, STOP)$ based on $\pi(n, u) + ln(a_{u,STOP})$ for the best three paths across all node u at position n.

After obtaining all the scores, we then filtered out the best three paths. The scores and its corresponding preceding node are then stored in a tuple.

In the backtracking process, we hold N best path sequences in the list N_bestP aths in decreasing order: N_bestP aths = [[1st best path sequence], [2nd best path sequence], ..., [N best path sequence]]. During the forward algorithm, in each node at position k, we kept track of the 3 best previous nodes (at position k-1) with their respective scores as illustrated in the figure below.



For each path, we need to keep track of the corresponding previous node's 3 best paths. Therefore at each current node at position k, we consult the scores dictionary at the current position k with the key of the best previous node (at position k-1) for each N best paths. Now, we have the best previous node to the current node and we add it to the corresponding path in the list N_bestP aths . We iterate through this process for all positions until position=1.

We have to modify the backtracking algorithm when picking the best path from position k=1 as nodes at this position only have the a singular score (ie. there is only 1 way to traverse from START state to nodes at this position).

With N_bestP aths , we can return the Nth best path.

This is our result for the third best output sequence, based on our modified viterbi algorithm:

EN

#Entity in gold data: 13179 #Entity in prediction: 13556

#Correct Entity: 10788
Entity precision: 0.7958
Entity recall: 0.8186
Entity F: 0.8070

#Correct Sentiment: 10100
Sentiment precision: 0.7451
Sentiment recall: 0.7664
Sentiment F: 0.7556

We ran the Nviterbi algorithm with N=1 and we got the same result as in Part 3. This proves that the algorithm is consistent regardless of what N is provided, and will work for any specified N.

Second-Order Hidden Markov Model

The second-order hidden Markov model works in a similar fashion compared to the first order HMM. The only difference now is that we consider second-order dependencies, in a trigram format - take into account second-order dependencies where the prediction of a tag is dependent on not just one, but two preceding tags. Here, we define the order of states that we calculate at each position/time step to be $state\ u \rightarrow state\ v \rightarrow state\ w$.

Mathematical Formulation

$$a_{u,v,w} = q(w|u,v) = \frac{Count(u,v,w)}{Count(u,v)}$$

$$b_w(x_j) = e(x|y) = \frac{Count(w \to x_j)}{Count(w)}$$

Base case:

$$\pi(0, v, w) = 1$$
 if $v, w = START, 0$ otherwise

Moving forward recursively:

For any
$$k \in \{1, ..., n\}$$
, for any $v, w \in \tau$,
$$\pi(k, v, w) = \max_{v \in \tau} \{\pi(k-1), u, v) \cdot a_{u,v,w} \cdot b_w(x_k)\}$$

Transition from y_n to STOP:

$$\max_{y_1, ..., y_n} p(x_1, ..., x_n, y_0 = START, y_1, ...y_n, y_{n+1} = STOP)$$

= $\max_{v,w \in \tau} \{\pi(n, v, w) \cdot a_{v,w,STOP}\}$

Backtracking

$$y_{n-1}^*, y_n^* = argmax_{v,w} \{\pi(n, v, w) \cdot a_{v,w,STOP}\}$$

For
$$k = n - 2$$
 to $k = 1$:

$$y_k^* = argmax_u \{ \pi(k+1, u, y_{k+1}^*) \cdot a_{u, y_{k+1}^*, y_{k+2}^*} \cdot b_{y_{k+2}^*}(x_{k+2}) \}$$

= $argmax_u \{ \pi(k+1, u, y_{k+1}^*) \cdot a_{u, y_{k+1}^*, y_{k+2}^*} \}$

Code Implementation

Similarly, we also assign the special token #UNK#, to words that did not appear in the training set, but yet appearing in the test set. This smoothing parameter helps to optimise our predictions further, especially for unseen words.

We implemented two new functions, HMM2_count_transition and HMM2_transition_para to tabularise transition counts and to calculate the transition probabilities respectively.

HMM2_count_transition function returns transition_tracker, which is a nested dictionary in this format:

This would hence allow us to easily retrieve numerator and denominator values required to calculate the transition parameters, using the function HMM2 transition para.

Viterbi algorithm for second-order HMM

Keeping in mind that the natural logarithm function is strictly increasing, all values are calculated in natural logarithmic form in our implementation, replacing multiplication by addition. This is based on the logarithmic property $\ln{(ab)} = \ln(a) + \ln(b)$. The purpose of taking the natural logarithm of probabilities is to reduce the risk of potential numerical underflow. However, one new issue with natural logarithmic calculations is that $\ln{0} = -\infty$. To deal with probability values that are equal to 0, we simply set $\ln{0} = C$, where C is a large negative number.

Similar to the viterbi-related algorithms implemented in Parts 3 and 4 above, the base case of $\pi(0, START, START) = 1$ if v, w = START, 0 otherwise is already accounted for, due to the structure of the datasets and that $\ln 1 = 0$. Hence, our implementation's base case is to determine the score at $\pi(1, START, w)$, ie. the score at the position of the first word in a sentence.

At each position/time step, the score of the best path are stored in a dictionary, scores, in such format:

We store the highest score of previous 2 states, $state\ u\ and\ state\ v$, as the value of the nested dictionary from START to $state\ w$ is for the purpose of backtracking.

At every position/time-step, we calculate the transition and emission probabilities, then take their natural logarithmic values as the score that we are tracking on. Prior to calculating the emission probabilities, we check whether the observation has appeared in the training set or not. If the observation was not seen in the training set, we assign the special token #UNK#.

Moving forward recursively:

1. We begin by initialising the first word of a sentence, ie. calculating $\pi(1, v, w)$:

$$\pi(1, v, w) = ln[a_{START, START, w} \bullet b_w(x_1)]$$

$$= ln(a_{START, START, w}) + ln([b_w(x_1)])$$

$$= ln(1) + ln[b_w(x_1)]$$

$$= ln[b_w(x_1)]$$

- 2. Case: 1 Word Sentence and 2 Word Sentences
 - a. In both cases, we calculate the possible paths and keep track of the state that yields the maximum score. It is important to look up the scores dictionary for the correct previous states $\pi(k-1,u,v)$.

$$\pi(k, v, w) = \max_{v} \left\{ \pi(k-1, u, v) + \ln(a_{u,v,w}) + \ln[b_w(x_k)] \right\}$$

An example for 1 word sentences is as follows:

$$\pi(0, v, w) = \max_{v} \{ \pi(-1, START, START) + \ln(a_{START, START, START}) \}$$

$$\pi(1, v, w) = \max_{v} \{ \pi(1, START, START) + \ln(a_{START, START, w}) + \ln[b_{w}(x_{1})] \}$$

$$\pi(2, v, w) = \max_{v} \{ \pi(1, START, v) + \ln(a_{START, v, w}) \}$$

- 3. Case: >2 Word Sentences
 - a. From the second word to the last word of a sentence, we calculate all the possible paths and keep track of the state that gives the maximum score:

$$\pi(k, v, w) = \max_{v} \left\{ \pi(k-1, u, v) + \ln(a_{u,v,w}) + \ln[b_w(x_k)] \right\}$$

4. In the final step, we calculate the score of the overall path from START state to STOP state.

ie. calculate
$$\pi(k+1, w, STOP) = max_v \{\pi(k, v, w) + ln(a_{v,w,STOP})\}$$
.

After calculating the best path from START state to STOP state, ie. the best path for the entire sentence, we then do backtracking to retrieve the states (ie. tags) at each position/time step. During the forward recursive process, we stored the state of position k-1 that gave the maximum score as a tuple along with the score at position k, hence we now retrieve all the states from each tuple to produce the full tag sequence for a sentence in the backtracking process.

This is our result for the second order hidden markov model, based on our modified viterbi algorithm:

EN

#Entity in gold data: 13179 #Entity in prediction: 13374

#Correct Entity : 10835
Entity precision: 0.8102
Entity recall: 0.8221
Entity F: 0.8161

#Correct Sentiment: 10416 Sentiment precision: 0.7788 Sentiment recall: 0.7903

Sentiment F: 0.7845

Tuning k value (Optional)

We tune the value of k to select the ideal balance between replacing unseen words with the #UNK# token and reducing the amount of vocabulary that our HMM model can capture in its parameters. In this case, we simply select the value of k that produces the best results on the development set.

We tested a range of k values from 2 to 9 and found that the higher the k, the better the **Part 3 Viterbi** Algorithm model performed in terms of F scores.

EN EN #Entity in gold data: 13179 #Entity in gold data: 13179 #Entity in gold data: 13179 #Entity in prediction: 13341 #Entity in prediction: 13344 #Entity in prediction: 13341 #Correct Entity: 10816 #Correct Entity : 10813 #Correct Entity : 10814 Entity precision: 0.8107 Entity recall: 0.8207 Entity precision: 0.8103 Entity precision: 0.8106 Entity recall: 0.8205 Entity recall: 0.8205 Entity F: 0.8157 Entity F: 0.8155 Entity F: 0.8154 #Correct Sentiment : 10290 #Correct Sentiment : 10292 #Correct Sentiment : 10294 Sentiment precision: 0.7715 Sentiment recall: 0.7809 Sentiment precision: 0.7711 Sentiment precision: 0.7716 Sentiment recall: 0.7808 Sentiment recall: 0.7811 Sentiment F: 0.7759 Sentiment F: 0.7762 Sentiment F: 0.7763 k = 2k = 3k = 4FN #Entity in gold data: 13179 #Entity in gold data: 13179 #Entity in gold data: 13179 #Entity in prediction: 13338 #Entity in prediction: 13329 #Entity in prediction: 13330 #Correct Entity: 10827 #Correct Entity: 10817 #Correct Entity: 10830 Entity precision: 0.8123 Entity recall: 0.8215 Entity precision: 0.8110 Entity precision: 0.8125 Entity recall: 0.8218 Entity recall: 0.8208 Entity F: 0.8169 Entity F: 0.8159 Entity F: 0.8171 #Correct Sentiment: 10304 #Correct Sentiment: 10294 #Correct Sentiment : 10307 Sentiment precision: 0.7731 Sentiment precision: 0.7718 Sentiment precision: 0.7732 Sentiment recall: 0.7818 Sentiment recall: 0.7811 Sentiment recall: 0.7821 Sentiment F: 0.7774 Sentiment F: 0.7764 Sentiment F: 0.7776 k = 5k = 6k = 7#Entity in gold data: 13179 #Entity in gold data: 13179 #Entity in prediction: 13326 #Entity in prediction: 13327

k = 8

#Correct Entity: 10833

Entity recall: 0.8220

Entity F: 0.8174

Sentiment F: 0.7780

Entity precision: 0.8129

#Correct Sentiment : 10311

Sentiment precision: 0.7737 Sentiment recall: 0.7824

k = 9

#Correct Entity: 10834

Entity recall: 0.8221

Entity F: 0.8175

Entity precision: 0.8130

#Correct Sentiment : 10312

Sentiment recall: 0.7825

Sentiment F: 0.7781

Sentiment precision: 0.7738

We then tested a few k values from range of 2 to 9 and found that the higher the k, the better the **Part 5 Second Order HMM** model seems to perform in terms of F scores.

#Entity in gold data: 13179 #Entity in gold data: 13179 #Entity in gold data: 13179 #Entity in prediction: 13376 #Entity in prediction: 13377 #Entity in prediction: 13377 #Correct Entity : 10836 #Correct Entity: 10835 #Correct Entity: 10836 Entity precision: 0.8100 Entity recall: 0.8221 Entity F: 0.8160 Entity precision: 0.8100 Entity recall: 0.8222 Entity precision: 0.8100 Entity recall: 0.8222 Entity F: 0.8161 Entity F: 0.8161 #Correct Sentiment : 10414 #Correct Sentiment : 10420 #Correct Sentiment : 10420 Sentiment precision: 0.7789
Sentiment recall: 0.7907
Sentiment F: 0.7848 Sentiment precision: 0.7786 Sentiment recall: 0.7902 Sentiment precision: 0.7789 Sentiment recall: 0.7907 Sentiment F: 0.7843 Sentiment F: 0.7848 k = 2k = 6k = 7ΕN #Entity in gold data: 13179 #Entity in gold data: 13179 #Entity in prediction: 13381 #Entity in prediction: 13381 #Correct Entity : 10840 #Correct Entity : 10840 Entity precision: 0.8101 Entity precision: 0.8101 Entity recall: 0.8225 Entity recall: 0.8225 Entity F: 0.8163 Entity F: 0.8163 #Correct Sentiment : 10422 Sentiment precision: 0.7789 #Correct Sentiment : 10422 Sentiment precision: 0.7789 Sentiment recall: 0.7908 Sentiment recall: 0.7908 Sentiment F: 0.7848 Sentiment F: 0.7848

To prevent overfitting of the model, k= 8 will be preferred instead of k=9. In addition, this model is for reference and its output is not considered for the evaluation on the test sets provided in the zip file.

k = 9

k = 8