1 Class-Conditional Densities for Binary Data

Question A:

$$p(x|y=c) = p(x_1|y=c)p(x_2|x_1, y=c) \cdots p(x_D|x_1, \cdots, x_{D-1}, y=c) = \prod_{i=1}^{D} \theta_{x_i i}$$

Since all the D features are binary, $x_j \in \{0, 1\}$,

for each class of C we need 2^{j-1} parameters for $p(x_j|x_1,\cdots,x_{j-1},y=c)$. In total $\sum_{j=1}^D 2^{j-1} = O(2^D)$. Therefore, we need $O(C\times 2^D)$ parameters for C classes.

Question B: Without factorization, since all the D features are binary, $x_j \in \{0, 1\}$, and C classes for y, in total we need $O(C \times 2^D)$, which is the same as that with factorization.

Question C: For a small N, Naive Bayes is likely to give lower test set error. This is because full models are more likely to overfit with a small N.

Question D: For a large N, full models are likely to give lower test set error. This is because Naive Bayes is simple and is likely to underfit with a large N, while full models have more parameters and will do better.

Question E: For Naive Bayes,

$$p(y|x) = \frac{p(x,y)}{p(x)} = \frac{p(y)}{p(x)} \prod_{d} p(x^{d}|y) = \frac{p(y)}{\sum_{i=1}^{C} p(x|y=c_i)} \prod_{d} p(x^{d}|y)$$

Since we assumed a uniform class prior p(y), O(p(y)) = O(1), and computation complexity of p(y|x) = O(CD).

For full model,

$$p(y|x) = \frac{p(x,y)}{p(x)} = \frac{p(y)}{p(x)}p(x|y)$$

, and computation complexity of D-dimensional vector is O(D). Different from Naive Bayes, full model doesn't take into account every y = c, but their overall probability. Therefore, computation complexity of p(y|x) = O(D).

2 Sequence Prediction

Question A:

Runnin	######################################		

File #0:			
Emission Sequence	Max Probability State Sequence		

25421	31033		
01232367534	22222100310		
5452674261527433	1031003103222222		
7226213164512267255	1310331000033100310		
0247120602352051010255241	22222222222222222222103		
File #1:			
Emission Sequence	Max Probability State Sequence		
,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,			
77550	22222		
7224523677	2222221000		
505767442426747	222100003310031		
72134131645536112267	10310310000310333100		
4733667771450051060253041	2221000003222223103222223		
File #2:		File #4:	
File #2. Emission Sequence	Max Probability State Sequence	Emission Sequence	Max Probability State Sequence
######################################			
60622	11111	23664	01124
4687981156	2100202111	3630535602	0111201112
815833657775062	02101111111111	350201162150142	011244012441112
21310222515963505015	0202011111111111021	00214005402015146362	11201112412444011112
6503199452571274006320025	1110202111111102021110211	2111266524665143562534450	2012012424124011112411124
		File #5:	
File #3:		Emission Sequence	Max Probability State Sequence
Emission Sequence Max Probability State Sequence		#######################################	
***************************************		68535	10111
13661	00021	4546566636	1111111111
2102213421	3131310213	638436858181213	110111010000011
166066262165133 53164662112162634156	133333133133100	13240338308444514688	00010000000111111100
03104002112102034150	20000021313131002133 1310021333133133133133133	0111664434441382533632626	21111111111111001111110101

${\bf Question}~{\bf B:}$ The results using Forward algorithm are as follows:

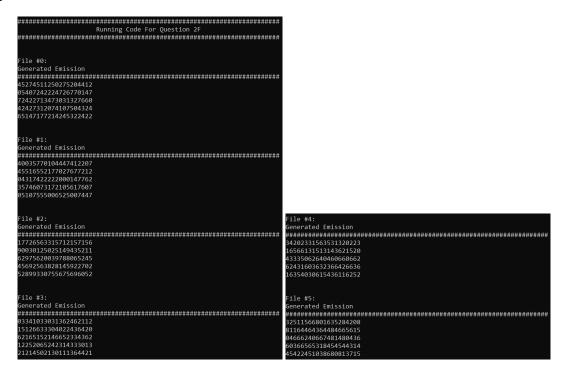
The results using Backward algorithm are as follows:

Question C:

Question D:

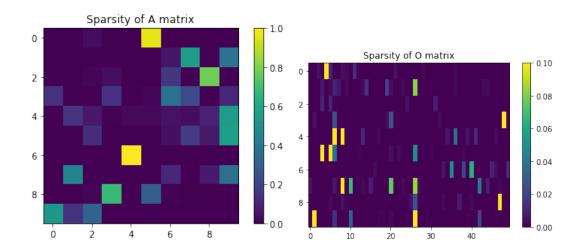
Question E: The transition and emission matrices from 2C and 2D are different, supervised matrices are more uniform, while unsupervised are sparse and nonuniform. 2C (supervised HMM) provides a more accurate representation. This is because for unsupervised 2D, we simply random the initial matrices, which cannot truly reflect Ron's moods. To improve the unsupervised learning data, we might need to initialize the transition and emission matrices A & O more precisely, such as using probability distributions from supervised matrices.

Question F:



Ouestion G:

The trained A and O matrices are both sparse, with most elements near 0. The sparsity of transition matrix A means that for each state, there might be very few states to transit to, and the sparsity of observation matrix O means that for each state, there might be very few observations to belong to.



Question H: As the number of hidden states is increased, sample emission sentences from the HMM becomes more coherent and smooth. When there is only one hidden state, then there is no more transitions, meaning that observation words are just randomly picked from the dataset. Allowing more hidden states will increase the training data likelihood, but when we have too many hidden states for the fixed observation set, this will lead to overfitting and may increase the test error.

Question I: As shown in the figure, I think state 7 is semantically meaningful. This state is filled with



law-related words, such as 'law', 'court', 'representative', 'justice', 'rules'. I think this state represent all law-related words in the dataset, which is distinct from other states.

HMM.py

March 5, 2020

```
# CS/CNS/EE 155 2018
    # Problem Set 6
    # Author:
                   Andrew Kang
    # Description: Set 6 skeleton code
    # You can use this (optional) skeleton code to complete the HMM
    # implementation of set 5. Once each part is implemented, you can simply
    # execute the related problem scripts (e.q. run 'python 2G.py') to quickly
    # see the results from your code.
    # Some pointers to get you started:
    #
          - Choose your notation carefully and consistently! Readable
            notation will make all the difference in the time it takes you
            to implement this class, as well as how difficult it is to debug.
    #
          - Read the documentation in this file! Make sure you know what
            is expected from each function and what each variable is.
          - Any reference to "the (i, j) th" element of a matrix T means that
    #
    #
            you should use T[i][j].
    #
          - Note that in our solution code, no NumPy was used. That is, there
            are no fancy tricks here, just basic coding. If you understand HMMs
            to a thorough extent, the rest of this implementation should come
            naturally. However, if you'd like to use NumPy, feel free to.
    #
          - Take one step at a time! Move onto the next algorithm to implement
    #
            only if you're absolutely sure that all previous algorithms are
            correct. We are providing you waypoints for this reason.
    # To get started, just fill in code where indicated. Best of luck!
    import random
```

```
class HiddenMarkovModel:
    Class implementation of Hidden Markov Models.
    def __init__(self, A, 0):
        Initializes an HMM. Assumes the following:
            - States and observations are integers starting from 0.
            - There is a start state (see notes on A_start below). There
              is no integer associated with the start state, only
              probabilities in the vector A_start.
            - There is no end state.
        Arguments:
            A:
                        Transition matrix with dimensions L x L.
                        The (i, j) th element is the probability of
                        transitioning from state i to state j. Note that
                        this does not include the starting probabilities.
            0:
                        Observation matrix with dimensions L x D.
                        The (i, j) th element is the probability of
                        emitting observation j given state i.
        Parameters:
            L:
                        Number of states. word of tags: N, V, adj
            D:
                        Number of observations. fish, sleep
                        The transition matrix. L x L
            A:
                        The observation matrix. L x D
            0:
                        Starting transition probabilities. The i^th element
            A_start:
                        is the probability of transitioning from the start
                        state to state i. For simplicity, we assume that
                        this distribution is uniform.
        111
        self.L = len(A)
        self.D = len(0[0])
        self.A = A
        self.0 = 0
        self.A_start = [1. / self.L for _ in range(self.L)]
```

```
def viterbi(self, x):
       Uses the Viterbi algorithm to find the max probability state
       sequence corresponding to a given input sequence.
       Arguments:
                       Input sequence in the form of a list of length M,
           x:
                       consisting of integers ranging from O to D-1.
       Returns:
           max seq:
                     State sequence corresponding to x with the highest
                       probability.
       111
       M = len(x)
                       # Length of sequence.
       # The (i, j) th elements of probs and segs are the max probability
       # of the prefix of length i ending in state j and the prefix
       # that gives this probability, respectively.
       # For instance, probs[1][0] is the probability of the prefix of
       # length 1 ending in state 0.
       probs = [[0. for _ in range(self.L)] for _ in range(M + 1)] #__
    probability
       seqs = [['' for _ in range(self.L)] for _ in range(M + 1)] # viterbiu
       # L # of states/tags, D: # of observations/ , M: sequence , A: L x L,_{\sqcup}
\rightarrow 0: L \times D
       for i in range(self.L): # initialize 1st state, probs[1][i] A_start⊔
→* 0, probs[0] [0,...,0]
           probs[1][i] = self.A_start[i] * self.O[i][x[0]]
       for seq in range(1, M): # sequence
           for curr in range(self.L): # states/tags
               max value = 0
               \max idx = 0
               for last in range(self.L): # state
                   if (probs[seq][last] * self.A[last][curr] * self.
\rightarrow0[curr][x[seq]] >= max_value):
                       max_value = probs[seq][last] * self.A[last][curr] *__
\rightarrowself.0[curr][x[seq]] #
                       max_idx = last
               probs[seq + 1][curr] = max_value #
                                                         state probs
               seqs[seq + 1][curr] = max_idx
```

```
max_seq_rev = []
   max_value = max(probs[M])
   max_idx = probs[M].index(max_value) #__
 state max_value( max_prob) max_idx
   max_seq_rev.append( str(max_idx) )
   for i in range(M, 1, -1):
        max_seq_rev.append( str(seqs[i][max_idx]) )
       max_idx = seqs[i][max_idx]
   max_seq = max_seq_rev[::-1]
   return "".join(max_seq)
def forward(self, x, normalize=False):
    Uses the forward algorithm to calculate the alpha probability
    vectors corresponding to a given input sequence.
    Arguments:
                    Input sequence in the form of a list of length M,
        x:
                    consisting of integers ranging from 0 to D - 1.
        normalize: Whether to normalize each set of alpha_j(i) vectors
                    at each i. This is useful to avoid underflow in
                    unsupervised learning.
    Returns:
        alphas:
                    Vector of alphas.
                    The (i, j) th element of alphas is alpha_j(i),
                    i.e. the probability of observing prefix x^1:i
                    and state y^i = j.
                    e.g. alphas[1][0] corresponds to the probability
                    of observing x^1:1, i.e. the first observation,
                    given that y^1 = 0, i.e. the first state is 0.
    111
   M = len(x)
                    # Length of sequence.
    alphas = [[0. for _ in range(self.L)] for _ in range(M + 1)]
   for i in range(self.L):
```

```
alphas[1][i] = self.A_start[i] * self.O[i][x[0]]
       for seq in range(1, M): # sequence
           for curr in range(self.L): # states/tags
               sum_value = 0
               for last in range(self.L): # state
                   sum_value += alphas[seq][last] * self.A[last][curr] * self.
\hookrightarrow 0[curr][x[seq]] # Viterbi replaces sum with max
               alphas[seq + 1][curr] = sum_value
           if normalize:
               sum_alpha = sum(alphas[seq + 1])
               for curr in range(self.L):
                   alphas[seq + 1][curr] /= sum_alpha
       return alphas
   def backward(self, x, normalize=False):
       Uses the backward algorithm to calculate the beta probability
       vectors corresponding to a given input sequence.
       Arguments:
                       Input sequence in the form of a list of length M,
           x:
                       consisting of integers ranging from 0 to D - 1.
           normalize: Whether to normalize each set of beta j(i) vectors
                        at each i. This is useful to avoid underflow in
                        unsupervised learning.
       Returns:
           betas:
                       Vector of betas.
                        The (i, j) th element of betas is beta_j(i), i.e.
                        the probability of observing prefix x^{(i+1)}M and
                        state y \hat{i} = j.
                        e.g. betas[M][O] corresponds to the probability
                        of observing x^M+1:M, i.e. no observations,
                        given that y \mathcal{M} = 0, i.e. the last state is 0.
       111
       M = len(x)
                        # Length of sequence.
       betas = [[0. for _ in range(self.L)] for _ in range(M + 1)]
```

```
for i in range(self.L):
           betas[-1][i] = 1 # PPT 74, beta(M) = 1
       for seq in range(-1, -M-1, -1): # sequence
           for curr in range(self.L): # states/tags
               sum_value = 0
               for nxt in range(self.L): # state
                   if seq != -M:
                       sum_value += betas[seq][nxt] * self.A[curr][nxt] * self.
\rightarrow0[nxt][x[seq]] # PPT 74, A_{z,j} A_{j,z}
                   else:
                       sum_value += betas[seq][nxt] * self.A_start[nxt] * self.
\rightarrow0[nxt][x[seq]]
               betas[seq - 1][curr] = sum_value
           if normalize:
               sum_beta = sum(betas[seq - 1])
               for curr in range(self.L):
                   betas[seq - 1][curr] /= sum_beta
       return betas
   def supervised_learning(self, X, Y):
       Trains the HMM using the Maximum Likelihood closed form solutions
       for the transition and observation matrices on a labeled
       datset (X, Y). Note that this method does not return anything, but
       instead updates the attributes of the HMM object.
       Arguments:
           X:
                       A dataset consisting of input sequences in the form
                       of lists of variable length, consisting of integers
                       ranging from 0 to D-1. In other words, a list of
                       lists.
           Y:
                       A dataset consisting of state sequences in the form
                       of lists of variable length, consisting of integers
                       ranging from 0 to L-1. In other words, a list of
                       lists.
                       Note that the elements in X line up with those in Y.
       111
       # Calculate each element of A using the M-step formulas.
```

```
A_count = [[0. for i in range(self.L)] for j in range(self.L)]
       A_sum = [0. for i in range(self.L)]
       # For each input sequence:
       for y in Y: # Y is a list of lists of length L, so y should be a list, u
\rightarrow each y[i] is a mood tag
           for i in range(len(y) - 1): # A is calculated by
               A_{count}[y[i]][y[i+1]] += 1
               A_sum[y[i]] += 1
       for curr in range(self.L): # to normalize the A matrix
           for nxt in range(self.L):
               self.A[curr][nxt] = A_count[curr][nxt] / A_sum[curr]
       # Calculate each element of O using the M-step formulas.
       O_{count} = [[0. for i in range(self.D)] for j in range(self.L)] # <math>O = L_{L}
\rightarrow x D, (i, j) is the probability of emitting observation j given state i.
       0_sum = [0. for i in range(self.L)]
       for x, y in zip(X,Y):
           for i in range(len(y)):
               O_{count[y[i]][x[i]] += 1}
               0_sum[y[i]] += 1
       for curr in range(self.L):
           for nxt in range(self.D):
               self.O[curr][nxt] = O_count[curr][nxt] / O_sum[curr]
       pass
   def unsupervised_learning(self, X, N_iters):
       Trains the HMM using the Baum-Welch algorithm on an unlabeled
       datset X. Note that this method does not return anything, but
       instead updates the attributes of the HMM object.
       Arguments:
                       A dataset consisting of input sequences in the form
           X:
                        of lists of length M, consisting of integers ranging
                       from 0 to D - 1. In other words, a list of lists.
           N_iters:
                       The number of iterations to train on.
       ,,,
```

```
for iteration in range(N_iters):
           A_count = [[0. for i in range(self.L)] for j in range(self.L)]
           A_sum = [0. for i in range(self.L)]
           O_count = [[0. for i in range(self.D)] for j in range(self.L)]
           0_sum = [0. for i in range(self.L)]
           # for each list
           for x in X:
               M = len(x)
                              # Length of sequence.
                # expectation step: given A \& O matrix, predict probs of y's
\rightarrow for each training x
                # use forward-backward algorithm, alpha & beta [1,...,M]
                alphas = self.forward(x, normalize=True) # (M+1)xL, (i, j)is_{\perp}
\rightarrow alpha j(i), probability of observing prefix x^1:i and state y^i = j.
                betas = self.backward(x, normalize=True)
                Marginals = [0. for _ in range(self.L)]
                for seq in range(1, M + 1): # for each prefix x^1:i
                    for curr in range(self.L): # states/tags
                        Marginals[curr] = alphas[seq][curr] * betas[seq][curr]
                    Marginals_sum = sum(Marginals)
                    for curr in range(self.L):
                        Marginals[curr] /= Marginals_sum # normalized P(y^i_{\perp})
\hookrightarrow (curr) | x),
                        # Maximization Step: Use y's to estimate new (A,O)
                        O_count[curr][ x[seq - 1] ] += Marginals[curr] #__
\hookrightarrow seq 1 , x[] 0
                        O_sum[curr] += Marginals[curr]
                        if seq != M: # A is calculated by
                            A_sum[curr] += Marginals[curr] # for unsupervised_
                               A\_count
\rightarrow learning, there is no tag,
                   P(y^i, y^i+1 \mid x) A
                for seq in range(1, M):
                    A_update = [[0. for _ in range(self.L)] for _ in range(self.
for curr in range(self.L):
                        for nxt in range(self.L):
                            A_update[curr][nxt] = alphas[seq][curr] * self.
\rightarrowA[curr][nxt] * self.O[nxt][x[seq]] * betas[seq + 1][nxt]
                             \# seq 1 , x[] 0
```

```
A_update_sum = sum( [sum(A_update[i]) for i in_u
→range(len(A_update)) ] )
                   for curr in range(self.L):
                       for nxt in range(self.L):
                           A_update[curr][nxt] /= A_update_sum # normalized_
\rightarrow P(y^i, y^i+1 / x)
                   for curr in range(self.L):
                       for nxt in range(self.L):
                           A_count[curr][nxt] += A_update[curr][nxt]
           for curr in range(self.L): # to normalize the A & O matrix
               for nxt in range(self.L):
                   self.A[curr][nxt] = A_count[curr][nxt] / A_sum[curr]
           for curr in range(self.L):
               for nxt in range(self.D):
                   self.O[curr][nxt] = O_count[curr][nxt] / O_sum[curr]
       pass
   def get_data_with_distribute(self, dist): #
       r = random.random()
       for i, p in enumerate(dist):
           if r < p:
               return i
           r -= p
   def generate_emission(self, M):
       111
       Generates an emission of length M, assuming that the starting state
       is chosen uniformly at random.
       Arguments:
           M:
                       Length of the emission to generate.
       Returns:
           emission:
                       The randomly generated emission as a list.
                       The randomly generated states as a list.
           states:
       111
       emission = []
       states = []
```

```
y_start = random.randint(0, self.L -1) # starting state is chosen_
→uniformly at random
      states.append(y_start)
      for i in range(M):
           # Generate observation/emission x
           #idx_random = random.randint(0, self.D -1)
           #max_value = max(self.O[ states[i] ])
           #x_index = self.O[ states[i] ].index(max_value)
           x_index = self.get_data_with_distribute(self.O[ states[i] ])
           emission.append(x_index)
           # Generate next state y.
           \#idx\_random = random.randint(0, self.L -1)
           y_index = self.get_data_with_distribute(self.A[ states[i] ])
           states.append(y_index)
      return emission, states[:-1]
  def probability_alphas(self, x):
      Finds the maximum probability of a given input sequence using
       the forward algorithm.
       Arguments:
                       Input sequence in the form of a list of length M,
           x:
                       consisting of integers ranging from 0 to D - 1.
       Returns:
          prob:
                       Total probability that x can occur.
       # Calculate alpha vectors.
      alphas = self.forward(x)
       # alpha_j(M) gives the probability that the state sequence ends
       # in j. Summing this value over all possible states j gives the
       # total probability of x paired with any state sequence, i.e.
       # the probability of x.
      prob = sum(alphas[-1])
      return prob
  def probability_betas(self, x):
```

```
Finds the maximum probability of a given input sequence using
        the backward algorithm.
        Arguments:
                        Input sequence in the form of a list of length M,
            x:
                        consisting of integers ranging from 0 to D - 1.
        Returns:
                        Total probability that x can occur.
            prob:
       betas = self.backward(x)
        # beta j(1) gives the probability that the state sequence starts
        # with j. Summing this, multiplied by the starting transition
        # probability and the observation probability, over all states
        # gives the total probability of x paired with any state
        # sequence, i.e. the probability of x.
       prob = sum([betas[1][j] * self.A_start[j] * self.O[j][x[0]] \
                    for j in range(self.L)])
       return prob
def supervised_HMM(X, Y):
   Helper function to train a supervised HMM. The function determines the
    number of unique states and observations in the given data, initializes
    the transition and observation matrices, creates the HMM, and then runs
    the training function for supervised learning.
   Arguments:
        X:
                    A dataset consisting of input sequences in the form
                    of lists of variable length, consisting of integers
                    ranging from 0 to D - 1. In other words, a list of lists.
        Y:
                    A dataset consisting of state sequences in the form
                    of lists of variable length, consisting of integers
                    ranging from 0 to L - 1. In other words, a list of lists.
                    Note that the elements in X line up with those in Y.
    # Make a set of observations.
   observations = set()
   for x in X:
        observations |= set(x)
    # Make a set of states.
```

```
states = set()
   for y in Y:
        states |= set(y)
   # Compute L and D.
   L = len(states)
   D = len(observations)
   # Randomly initialize and normalize matrix A.
   A = [[random.random() for i in range(L)] for j in range(L)]
   for i in range(len(A)):
       norm = sum(A[i])
        for j in range(len(A[i])):
            A[i][j] /= norm
    # Randomly initialize and normalize matrix O.
   0 = [[random.random() for i in range(D)] for j in range(L)]
   for i in range(len(0)):
       norm = sum(0[i])
       for j in range(len(O[i])):
            0[i][j] /= norm
    # Train an HMM with labeled data.
   HMM = HiddenMarkovModel(A, 0)
   HMM.supervised_learning(X, Y)
   return HMM
def unsupervised_HMM(X, n_states, N_iters):
   Helper function to train an unsupervised HMM. The function determines the
    number of unique observations in the given data, initializes
    the transition and observation matrices, creates the HMM, and then runs
    the training function for unsupervised learing.
   Arguments:
        X:
                    A dataset consisting of input sequences in the form
                    of lists of variable length, consisting of integers
                    ranging from 0 to D - 1. In other words, a list of lists.
        n_states: Number of hidden states to use in training.
       N_iters:
                   The number of iterations to train on.
    111
```

```
# Make a set of observations.
observations = set()
for x in X:
    observations |= set(x)
# Compute L and D.
L = n_states
D = len(observations)
# Randomly initialize and normalize matrix A.
random.seed(2020)
A = [[random.random() for i in range(L)] for j in range(L)]
for i in range(len(A)):
    norm = sum(A[i])
    for j in range(len(A[i])):
        A[i][j] /= norm
# Randomly initialize and normalize matrix O.
random.seed(155)
0 = [[random.random() for i in range(D)] for j in range(L)]
for i in range(len(0)):
   norm = sum(0[i])
    for j in range(len(O[i])):
        0[i][j] /= norm
# Train an HMM with unlabeled data.
HMM = HiddenMarkovModel(A, 0)
HMM.unsupervised_learning(X, N_iters)
return HMM
```

2 notebook

March 5, 2020

1 Problem 2

In this Jupyter notebook, we visualize how HMMs work. This visualization corresponds to problem 2 in set 6.

Assuming your HMM module is complete and saved at the correct location, you can simply run all cells in the notebook without modification.

```
[1]: import os
  import numpy as np
  from IPython.display import HTML

from HMM import unsupervised_HMM
  from HMM_helper import (
        text_to_wordcloud,
        states_to_wordclouds,
        parse_observations,
        sample_sentence,
        visualize_sparsities,
        animate_emission
)
```

1.1 Visualization of the dataset

We will be using the Constitution as our dataset. First, we visualize the entirety of the Constitution as a wordcloud:

```
[2]: text = open(os.path.join(os.getcwd(), 'data/constitution.txt')).read()
wordcloud = text_to_wordcloud(text, title='Constitution')
```

Constitution



1.2 Training an HMM

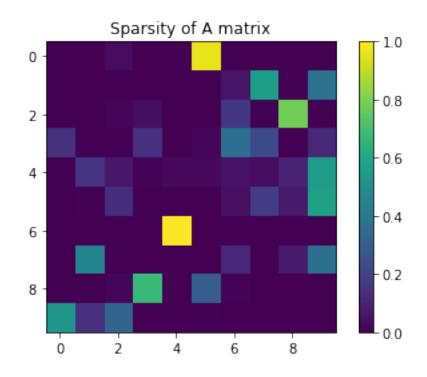
Now we train an HMM on our dataset. We use 10 hidden states and train over 100 iterations:

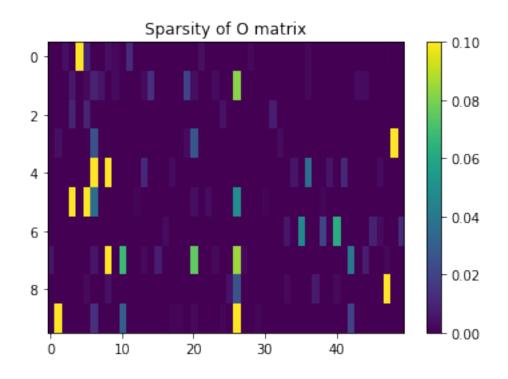
```
[3]: obs, obs_map = parse_observations(text)
hmm8 = unsupervised_HMM(obs, 10, 100)
```

1.3 Part G: Visualization of the sparsities of A and O

We can visualize the sparsities of the A and O matrices by treating the matrix entries as intensity values and showing them as images. What patterns do you notice?

[4]: visualize_sparsities(hmm8, O_max_cols=50)





1.4 Generating a sample sentence

As you have already seen, an HMM can be used to generate sample sequences based on the given dataset. Run the cell below to show a sample sentence based on the Constitution.

```
[5]: print('Sample Sentence:\n=========')
print(sample_sentence(hmm8, obs_map, n_words=25))
```

Sample Sentence:

War office the concurrence necessary and unless in state obliged to exports thirds be public thereby the elections states and expiration of an and counterfeiting...

1.5 Part H: Using varying numbers of hidden states

Using different numbers of hidden states can lead to different behaviours in the HMMs. Below, we train several HMMs with 1, 2, 4, and 16 hidden states, respectively. What do you notice about their emissions? How do these emissions compare to the emission above?

```
[6]: hmm1 = unsupervised_HMM(obs, 1, 100)
print('\nSample Sentence:\n==========')
print(sample_sentence(hmm1, obs_map, n_words=25))
```

Sample Sentence:

Or justice the shall state the shall emit court effect together any pay shall the and bound of states originated such the member capitation congress...

```
[7]: hmm2 = unsupervised_HMM(obs, 2, 100)
print('\nSample Sentence:\n==========')
print(sample_sentence(hmm2, obs_map, n_words=25))
```

Sample Sentence:

Legislature tonnage and and president and electors representative deprived effect or section and number between to that and shall the senate before by the whom...

```
[8]: hmm4 = unsupervised_HMM(obs, 4, 100)
print('\nSample Sentence:\n==========')
print(sample_sentence(hmm4, obs_map, n_words=25))
```

Sample Sentence:

Not cases have of judicial blood impeachment law their the fourths or be shall of states of the place adjournment concurrence of representatives or on...

```
[9]: hmm16 = unsupervised_HMM(obs, 16, 100)
print('\nSample Sentence:\n===========')
print(sample_sentence(hmm16, obs_map, n_words=25))
```

Sample Sentence:

To inferior of a monday maritime the judges of behaviour and arsenals and taxes been yeas concur departments and diminished to work different may a...

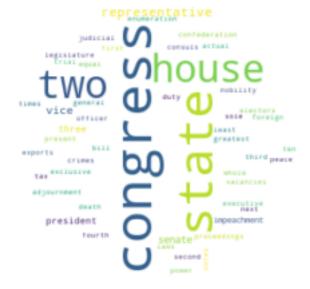
1.6 Part I: Visualizing the wordcloud of each state

Below, we visualize each state as a wordcloud by sampling a large emission from the state:

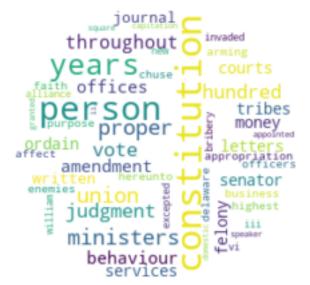
[10]: wordclouds = states_to_wordclouds(hmm8, obs_map)

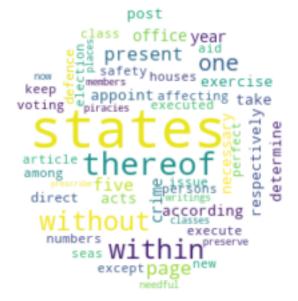








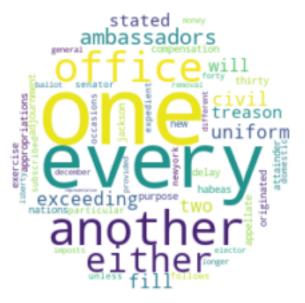












1.7 Visualizing the process of an HMM generating an emission

The visualization below shows how an HMM generates an emission. Each state is shown as a wordcloud on the plot, and transition probabilities between the states are shown as arrows. The darker an arrow, the higher the transition probability.

At every frame, a transition is taken and an observation is emitted from the new state. A red arrow indicates that the transition was just taken. If a transition stays at the same state, it is represented as an arrowhead on top of that state.

Use fullscreen for a better view of the process.

```
[11]: anim = animate_emission(hmm8, obs_map, M=8)
HTML(anim.to_html5_video())
```

Animating...

[11]: <IPython.core.display.HTML object>

Nor bound in the form determine the consent

