# ASEN 6519 - Assignment 1

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February 26, 2019

## Problem 1

#### State Trace of Short HMM Dataset

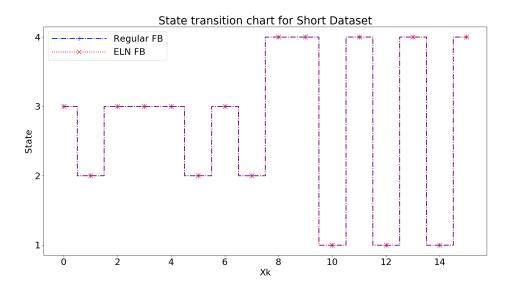


Figure 1: For the short dataset, the state trace is in lockstep for both the normal forward backwards and the logarithmic-based forward backward algorithm.

#### Posterior Probabilities

States	X0	X1	X2	X3	X4	X5	X6	X7	X8	X9	X11	X11	X12	X13	X14	X15
1	0	0	0	0	0	0	0	0	0.183297	0.0100234	0.976834	0.00115028	0.930981	0.00233115	0.950946	0.0548677
2	0	0.951268	0.0122975	0.153837	0.105652	0.89824	0.00223589	0.941818	0	0	0	0	0	0	0	0
3	1	0.0487323	0.987703	0.846163	0.894348	0.10176	0.997764	0.0581818	0	0	0	0	0	0	0	0
4	0	0	0	0	0	0	0	0	0.816703	0.989977	0.0231662	0.99885	0.0690191	0.997669	0.049054	0.945132

## Log Likelihood of Short Data

-28.13852617314715

# Problem 2

Sampling was performed by creating a chain of sampled state values for each of the random variables  $x_k$ , which in the case of a HMM, are already conveniently in topological ordering. Weights were determined for each sample by the following update function:

$$w_k = w_{k-1}P(y = ev|x_k = s) \tag{1}$$

The current weight at timestep k is the prior weight  $w_{k-1}$  times the probability that the sampled state for  $x_k$  emits the observed evidence y = ev. This iterative update is done until the end of the chain to give the weight for the entire sample.

Posterior distribution estimates for each state of  $X_k$  were generated by using the following formula:

$$P(x_k = s | Y_{1:T}) = \frac{\sum_{i=1}^{N} w_i \delta(x_i = s)}{\sum_{i=1}^{N} w_i}$$
 (2)

Where N is number of samples,  $\delta$  is to Kronecker Delta function / indicator that sifts out the desired states across all samples for a given state random variable  $x_k$ . This is a weighted frequency count of the sampled states for each state random variable  $x_k$ .

#### Exact Inference vs. Approximate Inference (LW) with 100 Samples

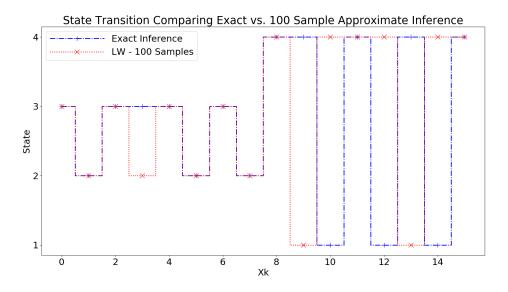


Figure 2: For 100 samples, oftentimes there are mismatched states in a max-likelihood state trace.

#### Exact Inference vs. Approximate Inference (LW) with 1000 Samples

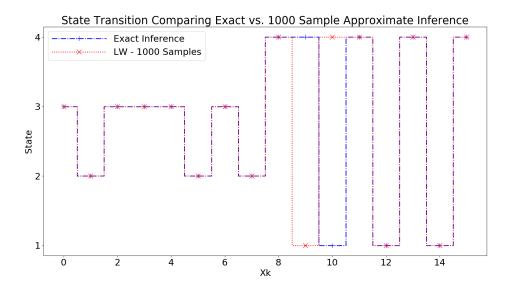


Figure 3: For 1000 samples, oftentimes there are mismatched states in a max-likelihood state trace.

#### Exact Inference vs. Approximate Inference (LW) with 10000 Samples

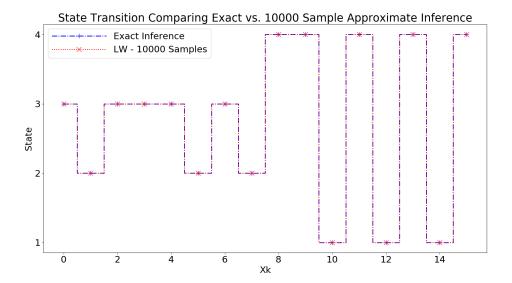


Figure 4: For 10000 samples, most runs result in perfect state trace alignment.

As seen in the above state traces, the likelihood weighting algorithm does provide fairly accurate state distribution inference. As the number of samples increases, the accuracy generally increases. However, if the long HMM dataset is substituted in place of the short dataset, the likelihood weighting (at least with the chosen q function) makes no inference. This is likely because emission probabilities have a zero probability for at least one of the states which makes it very, very unlikely to sample a full state sequence that does not 'zero out' the final weight.

## Problem 3

For the long dataset, the state trace is in lockstep early in the sequence, but starts to diverge later in the sequence. This is likely the the standard forward-backward algorithm (without Mann's logarithmic approach) is suffering from floating point underflow which creates incorrect values during the calculation of alpha and beta.

## State Trace for Long HMM Dataset

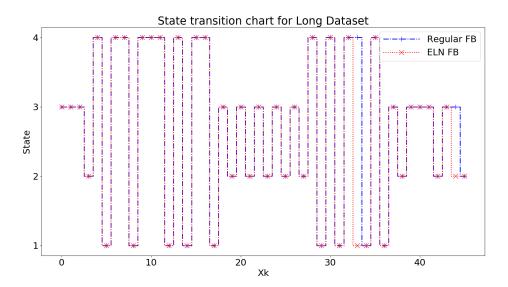


Figure 5: For the long dataset, the state trace is in lockstep for both the normal forward backwards and the logarithmic-based forward backward algorithm earlier in the sequence, but diverges later in the sequence.

## Posterior Probabilities for Long Dataset

State	X1	X2	Х3	X4	X5	X42	X43	X44	X45	X46
1	0	0	0	0	0.00618432	0	0	0	0	0
2	0	0.172346	0.00578533	0.995108	0	0.0132082	0.943541	0.0133714	0.157476	0.831321
3	1	0.827654	0.994215	0.00489207	0	0.986792	0.0564594	0.986629	0.842524	0.168679
4	0	0	0	0	0.993816	0	0	0	0	0

# Log Likelihood of Long Data

-84.74339048594653

## Extra Credit: Viterbi

#### State Trace for Short HMM Dataset

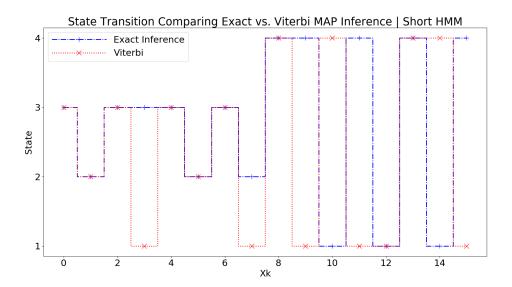


Figure 6: For the short dataset, the Viterbi algorithm is not as accurate as exact inference.

## State Trace for Long HMM Dataset

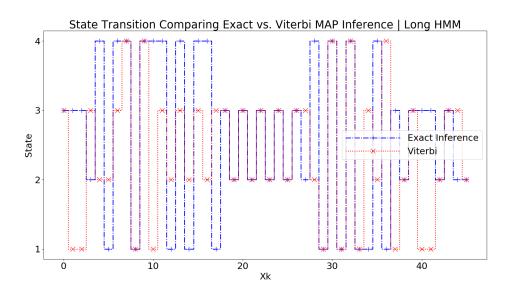


Figure 7: For the long dataset, the Viterbi algorithm is not as accurate as exact inference.

#### Code:

```
# forward_backward.py
import numpy as np
from exponential_scaling import eexp, eln, elnsum, elnproduct
import colored_traceback
colored_traceback.add_hook()

class ForwardBackwardHMM():
```

```
def __init__(self , transition_probs , evidence_probs , initial_probs , obs):
    self.n_states = len(initial_probs)
    self.init = initial_probs
    self.emis = evidence_probs
    self.trans = transition_probs
    self.n_obs = len(obs)
    self.obs = obs
def max_likelihood_state_estimate(self, probs):
    probs = probs.transpose()
    states = [np.argmax(prob_vec) + 1 for prob_vec in probs]
    return states
def forward_backward(self):
    Implements the forward backward algorithm for a simple HMM example. This
        function calls
    private methods foward and backwards to obtain the state probabilities for the
         forward
    and backward passes. They are then multiplied and normalized to get the full
        probabilities
    for each state at each time step.
    alphas = self._forward_iter()
    betas = self._backward_iter()
    # recast as np.arrays to perform element-wise multiplication
    probs = np.array(alphas)[:, 0:] * np.array(betas)[:, :self.n_obs + 1]
    probs = probs / np.sum(probs, 0)
    # for i in range(self.n_obs):
           print(np.sum(alphas[:, i] * betas[:, i]))
    return probs, alphas, betas
def forward_backward_eln(self):
    logalphas = self._forward_iter_eln()
    logbetas = self._backward_iter_eln()
    probs = np.zeros((self.n_states, self.n_obs + 1))
    for k in range (0, self.n_obs + 1):
        norm = -np.inf
        for i in range (self.n_states):
             probs[i, k] = elnproduct(logalphas[i, k], logbetas[i, k])
             norm = elnsum(norm, probs[i, k])
         for i in range (self.n_states):
             probs\left[\,i\;,\;\;k\,\right]\;=\;eexp\left(\,elnproduct\left(\,probs\left[\,i\;,\;\;k\,\right]\,,\;\;-norm\right)\,\right)
    # for i in range(self.n_obs):
    # print(sum([p for p in list(probs[:, i]) if p != -np.inf]))
return probs, logalphas, logbetas
def _forward(self):
    The forward (filtering) pass starting from the initial time step to the ending
        time step.
    alphas stores the calculated alpha value at each iteration normalized across
        states to avoid
    vanishing probabilities. The initial time step t_0 is initialized with the
        initial state
    probabilities.
    alphas = np.zeros((self.n_states, self.n_obs + 1))
    alphas[:, 0] = self.init
    for k in range(0, self.n_obs):
        alphas\,[:\,,\ k\,+\,1]\,=\,alphas\,[:\,,\ k\,]\,.\,dot\,(\,self\,.\,trans\,.\,transpose\,(\,)\,)\ *\ self\,.\,emis\,[
             self.obs[k], :]
    return alphas
def _backward(self):
    The backward (smoothing) pass starting from the an arbitrary future time step to the beginning time step. The matrix 'betas' stores the calculated beta value at
        each
```

```
iteration normalized across states to avoid vanishing probabilities.
          betas = np.zeros((self.n_states, self.n_obs))
          betas [:, -1] = 1
          for k in range (self.n_obs, 0, -1):
              beta_vec = np.matrix(betas[:, k]).transpose()
              betas [:, k - 1] = (np.matrix(self.trans.transpose()) * np.matrix(np.diag(
                   self.emis.transpose()[:, self.obs[k-1]])) * beta\_vec).transpose()
          return betas
     def _forward_iter(self):
          alphas = np.zeros((self.n_states, self.n_obs + 1))
         # base case
         alphas[:, 0] = self.init
         # recursive case
         for k in range (0, self.n_obs):
              for i in range (self.n_states):
                   for j in range(self.n_states):
                        alphas\left[\,i\,\,,\,\,k\,+\,\,1\,\right] \;+\!\!=\; alphas\left[\,j\,\,,\,\,k\,\right] \;*\; self.trans\left[\,i\,\,,\,\,k\,\right]
                                                                               j] * self.emis[self.
                                                                                    obs[k], i]
         return alphas
     def _backward_iter(self):
         betas = np.zeros((self.n_states, self.n_obs + 1))
         # base case
         betas[:, -1] = 1
         # recursive case
         for k in range(self.n_obs, -1, -1):
              for i in range(self.n_states):
                   for j in range(self.n_states):
                        betas\left[\,i\;,\;k\;-\;1\,\right]\;+\!\!=\;s\,e\,l\,f\;.\,t\,r\,a\,n\,s\,\left[\,j\;,\;\;i\;\right]\;*\;\backslash
                             self.emis[self.obs[
                                 k-1, j * betas[j, k]
          return betas
     def _forward_iter_eln(self):
         logalphas = np.zeros((self.n_states, self.n_obs + 1))
         # base case
         logalphas[:, 0] = [eln(x) for x in self.init]
         # recursive case
         for k in range(1, self.n_obs + 1):
              for j in range(self.n_states):
                   logalpha = -np.inf
                   for i in range(self.n_states):
                       logalpha = elnsum(logalpha, elnproduct(
                   logalphas[i, k-1], eln(self.trans.transpose()[i, j]))) \\ logalphas[j, k] = elnproduct(logalpha, eln(
                        self.emis[self.obs[k-1], j])
         return logalphas
     def _backward_iter_eln(self):
         logbetas = np.zeros((self.n_states, self.n_obs + 1))
         # base case
         logbetas[:, -1] = 0
         # recursive case
          for k in range (self.n_obs, -1, -1):
              for i in range(self.n_states):
                   logbeta = -np.inf
                   for j in range(self.n_states):
                        logbeta = elnsum(logbeta, elnproduct(eln(self.trans.transpose()[i, j
                            ]), elnproduct(
                   \begin{array}{c} \text{self.emis} \left[ \text{ self.obs} \left[ k-1 \right], \text{ } j \right], \text{ } logbetas \left[ j, \text{ } k \right] \right) \right) \\ logbetas \left[ i, \text{ } k-1 \right] = logbeta \end{array}
          return logbetas
# likelihood_sampling_infererence.py
import numpy as np
from numpy.random import choice
from observations import y_obs_long, y_obs_short
```

```
from \ parameters \ import \ pxk\_xkm1 \,, \ pyk\_xk \,, \ px0
from exponential_scaling import eexp, eln, elnsum, elnproduct
import colored_traceback
colored_traceback.add_hook()
class LikelihoodSamplingInference():
    def __init__(self, transition_probs, evidence_probs, initial_probs, observations):
        self.n_states = len(initial_probs)
        self.init_probs = initial_probs
        self.ev_probs = evidence_probs
        self.trans\_probs = transition\_probs
        self.n_observations = len(observations)
        self.observations = observations
    def max_likelihood_state_estimate(self, probs):
        probs = probs
        states = [np.argmax(np.array(prob_vec)) + 1 for prob_vec in probs]
        return states
    def run_inference(self, n_samples=100):
        samples, weights = self.run_sampling(n_samples)
probs = []
        for k in range(self.n_observations + 1):
            prob = []
            xk\_stack = list(list(zip(*samples))[k])
            for s in range (self.n_states):
                idxs = [idx for idx, value in enumerate(xk_stack) if value == s]
                prob.append(sum([w for idx, w in enumerate(weights) if idx in idxs]) /
                    sum(weights))
            probs.append(prob)
        return probs
    def run_sampling(self, n_samples):
        samples = []
        weights = []
        for n in range(n_samples):
            sample, weight = self.liklihood_sampling()
            samples.append(sample)
            weights.append(weight)
        return samples, weights
    def liklihood_sampling(self):
        w = 1
        X = []
        xkm1 = self.sample_state(self.init_probs)
        X. append (xkm1)
        for k in range(0, self.n_observations):
            xk = self.sample_state(list(self.trans_probs[:, xkm1]))
            ob_idx = self.observations[k]
            py_xk = self.ev_probs[ob_idx, xk]
            w = w * py_xk
            X. append (xk)
            xkm1\,=\,xk
        return (X, w)
    def sample_state(self, probs):
        return choice (4, 1, p=probs)[0]
# exponential_scaling.py
import numpy as np
def eexp(x):
    if np. isinf(x):
        return 0
    else:
        return np.exp(x)
def eln(x):
```

```
if x = 0:
        return -np.inf
    if x > 0:
        return np.log(x)
        raise ValueError ("x_must_be_non-negative_for_eln")
def elnsum(elnx, elny):
    if np.isinf(elnx) or (np.isinf(elny)):
        if (np.isinf(elnx)):
            return elnv
        else:
            return elnx
    else:
        if (elnx > elny):
            return elnx + eln(1 + np.exp(elny - elnx))
        else:
            return elny + eln(1 + np.exp(elnx - elny))
def elnproduct(elnx, elny):
    if np.isinf(elnx) or np.isinf(elny):
        return -np.inf
        return elnx + elny
# viterbi.py
import numpy as np
from observations import y_obs_long, y_obs_short
from parameters import pxk_xkm1, pyk_xk, px0
from \ exponential\_scaling \ import \ eexp\,, \ eln\,, \ elnsum\,, \ elnproduct
import colored_traceback
colored_traceback.add_hook()
class ViterbiHMM():
    def __init__(self, transition_probs, evidence_probs, initial_probs, observations):
        self.n_states = len(initial_probs)
        self.init = initial_probs
        self.emis = evidence_probs
        self.trans = transition_probs
        self.n_obs = len(observations)
        self.obs = observations
    def viterbi_path(self):
        init = np.array([eln(val) for val in np.nditer(self.init)])
        trans = np.array([[eln(v) for v in np.nditer(axis)] for axis in np.nditer(self.
            trans)]).reshape(self.n_states, self.n_states)
        print(np.array([[eln(v) for v in np.nditer(axis)] for axis in np.nditer(self.
            emis)]))
        emis = np. array([[eln(v) for v in np.nditer(axis)] for axis in np.nditer(self.
            emis)]).reshape(self.emis.shape[1], self.emis.shape[0])
        best\_states = []
        logprob = init
        for k in range (0, self.n_obs):
            trans_p = np.zeros([self.n_states, self.n_states])
            for i in range(self.n_states):
                for j in range(self.n_states):
                    trans_p[i, j] = elnsum(logprob[i], elnproduct(trans[i, j], emis[i,
                        self.obs[k-1]))
            # Get the indices of the max probs that give the best prior states.
            best_states.append(np.argmax(trans_p, axis=0))
            logprob = np.max(trans_p, axis=0)
        print(len(best_states))
        # Most likely final state.
        final_state = np.argmax(logprob)
        print(final_state)
```

```
# Reconstruct path by backtracking through likeliest states.
prior_state = final_state
best_path = [prior_state + 1]
for best in reversed(best_states):
    prior_state = best[prior_state]
    best_path.append(prior_state + 1)
return list(reversed(best_path)), logprob[final_state]
```