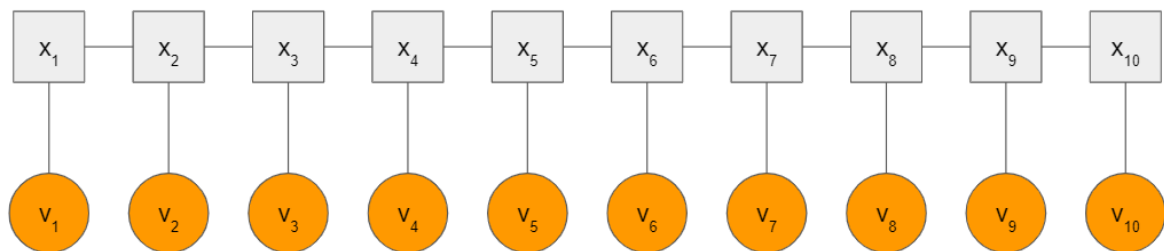


Ruijie_Rao_HW7_HMM

Consider a variable x with domain $\{1, 2, 3 \dots 10\}$. Let v_t be the value of x at timestep t . v_{t+1} is equal to $v_t - 1$ or $v_t + 1$ with probability 0.5 each, except when $v_t = 1$ or $v_t = 10$, in which case $v_{t+1} = 2$ or $v_{t+1} = 9$, respectively. At each timestep t , we also get noisy measurements of v_t . That is, $v_t - 1$, v_t or $v_t + 1$ can be returned with equal probabilities. Your task is to use a Hidden Markov Model to figure out the most likely sequence of values $v_1 v_2 \dots v_{10}$ when the observation sequence is 8, 6, 4, 6, 5, 4, 5, 5, 7, 9. At timestep $t = 1$, v_1 can be any value in $\{1, 2, 3 \dots 10\}$ with equal prior probabilities.



Viterbi Algorithm

```
In [1]: import numpy as np
import pandas as pd
```

Viterbi Algorithm is composed of 2 recursions.

First recursion constructs 2 matrices in a forward direction. S records the previous state, in other words records the routes starting from all possible initiations. P records the possibilities of the most probable state. At the end of the recursion, P gathers all the probability information in the last column and ready to find the optimal solution.

Second recursion goes backward and simply follows the trail of S , starting from the last state that has the highest probability in P .

```
In [13]: class Viterbi:
    def __init__(self, A, B, pi):
        self.A = A
        self.B = B
        self.pi = pi

    def fit(self, y):
        T = y.shape[0] # Number of total t
        N = self.A.shape[0] # Number of Possibilities
        P = np.zeros((T, N)) # Possibility Records
        P[0] = self.pi * self.B[:, y[0]] # Initiate with first overservation
        S = np.zeros((T - 1, N)) # State Records without initiation state

        for t in range(T - 1):
            for i in range(N):
                p = P[t] + self.A[:, i] + self.B[i, y[t + 1]] # calculation of possibilit
                S[t, i] = np.argmax(p) # pick the most propable state
                P[t + 1, i] = np.max(p) # record the best probability

        result = []
```

```

best_state = np.argmax(P[-1, :])
result.append(best_state)
for t in range(T-2, -1, -1):
    best_state = int(S[t, best_state])
    result.append(best_state)

result = result[::-1]
return result

```

In order to initiate the algorithm, 3 matrices needs to be inputted:

1. A is the *State Transition Matrix*. a_{ij} records the probability of x_i becomes x_j .
2. B is the *Emission Matrix*. b_{ij} records the probability of v_j is observed given x_i .
3. Pi is the *Initiation Probabilities* that records the probabilities of each initiating states.

```

In [16]: y = np.array([8, 6, 4, 6, 5, 4, 5, 5, 7, 9])
pi = 1/np.linspace(1,10,10)

A = np.zeros((10,10),float)
for i in range(10):
    for j in range(10):
        if (i+1==j) or (i-1==j):
            A[i,j] = 0.5
A[0,1] = 1.
A[9,8] = 1.

B = np.zeros((10,10),float)
for i in range(10):
    for j in range(10):
        if (i+1==j) or (i-1==j) or (i==j):
            B[i,j] = 1/3

```

```

In [14]: myviterbi = Viterbi(A,B,pi)

```

Outcome

```

In [15]: myviterbi.fit(y)

```

```

Out[15]: [7, 6, 5, 6, 5, 4, 5, 6, 7, 8]

```

Data Structure

Transfer Matrix is a NxN matrix. It has its diagonal surrounded by 0.5, because x_i can only transfer to +1 or -1 at the next t . An exception is for $i=1$ and $i=10$, where they only have one option with possibility 1.

```

In [17]: A

```

```

Out[17]: array([[0. , 1. , 0. , 0. , 0. , 0. , 0. , 0. , 0. , 0. ],
        [0.5, 0. , 0.5, 0. , 0. , 0. , 0. , 0. , 0. , 0. ],
        [0. , 0.5, 0. , 0.5, 0. , 0. , 0. , 0. , 0. , 0. ],
        [0. , 0. , 0.5, 0. , 0.5, 0. , 0. , 0. , 0. , 0. ],
        [0. , 0. , 0. , 0.5, 0. , 0.5, 0. , 0. , 0. , 0. ],
        [0. , 0. , 0. , 0. , 0.5, 0. , 0.5, 0. , 0. , 0. ],
        [0. , 0. , 0. , 0. , 0. , 0.5, 0. , 0.5, 0. , 0. ],
        [0. , 0. , 0. , 0. , 0. , 0. , 0.5, 0. , 0.5, 0. ],
        [0. , 0. , 0. , 0. , 0. , 0. , 0. , 0.5, 0. , 0. ],
        [0.5, 0. , 0.5, 0. , 0. , 0. , 0. , 0. , 0. , 0. ]])

```

```
[0. , 0. , 0. , 0. , 0. , 0. , 0.5, 0. , 0.5, 0. ],
[0. , 0. , 0. , 0. , 0. , 0. , 0. , 0.5, 0. , 0.5],
[0. , 0. , 0. , 0. , 0. , 0. , 0. , 0. , 0. , 1. , 0. ]])
```

Emission Matrix is a NxN matrix. It has a 3 width diagonal filled with 1/3, because for every x_i , its observed v_j can only be -1, +1 or 0 of its own value.

In [20]:

```
B
```

Out[20]:

```
array([[0.33333333, 0.33333333, 0.          , 0.          , 0.          ,
        0.          , 0.          , 0.          , 0.          , 0.          ],
       [0.33333333, 0.33333333, 0.33333333, 0.          , 0.          ,
        0.          , 0.          , 0.          , 0.          , 0.          ],
       [0.          , 0.33333333, 0.33333333, 0.33333333, 0.          ,
        0.          , 0.          , 0.          , 0.          , 0.          ],
       [0.          , 0.33333333, 0.33333333, 0.33333333, 0.33333333,
        0.          , 0.          , 0.          , 0.          , 0.          ],
       [0.          , 0.          , 0.33333333, 0.33333333, 0.33333333,
        0.          , 0.          , 0.          , 0.          , 0.          ],
       [0.          , 0.          , 0.          , 0.          , 0.          ,
        0.33333333, 0.33333333, 0.33333333, 0.          , 0.          ],
       [0.          , 0.          , 0.          , 0.          , 0.33333333,
        0.33333333, 0.          , 0.          , 0.          , 0.          ],
       [0.33333333, 0.          , 0.          , 0.          , 0.          ,
        0.          , 0.          , 0.          , 0.33333333, 0.          ],
       [0.          , 0.          , 0.          , 0.          , 0.          ,
        0.33333333, 0.33333333, 0.          , 0.          , 0.33333333],
       [0.33333333, 0.33333333, 0.          , 0.          , 0.          ,
        0.          , 0.          , 0.          , 0.          , 0.          ],
       [0.          , 0.          , 0.          , 0.          , 0.          ,
        0.33333333, 0.33333333, 0.33333333, 0.          , 0.          ],
       [0.          , 0.          , 0.          , 0.          , 0.          ,
        0.33333333, 0.33333333, 0.33333333, 0.33333333, 0.          ],
       [0.          , 0.          , 0.          , 0.          , 0.          ,
        0.          , 0.          , 0.          , 0.          , 0.          ],
       [0.          , 0.          , 0.          , 0.          , 0.          ,
        0.          , 0.          , 0.33333333, 0.33333333, 0.33333333],
       [0.          , 0.          , 0.          , 0.          , 0.          ,
        0.          , 0.          , 0.          , 0.          , 0.          ],
       [0.          , 0.          , 0.          , 0.          , 0.33333333,
        0.33333333, 0.33333333, 0.33333333, 0.33333333, 0.33333333]])
```

Initiating Possibilities are simply 1/10.

In [21]:

```
pi
```

Out[21]:

```
array([1.          , 0.5          , 0.33333333, 0.25          , 0.2          ,
        0.16666667, 0.14285714, 0.125          , 0.11111111, 0.1          ])
```

Challenges and Optimizations

Challenge: I was very challenged on how to convert Value Elimination concept to codes at the beginning, until I searched for Viterbi Algorithm and found out that it has the same concept of aggregating information to its last node.

Optimization: Created a algorithm object and has made the process more efficient in implementation.