

# Artificial Intelligence

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## *Today program*

Agent in **partially observable environment** maintains a belief state from the percepts observed and a **sensor model** and using a **transition model** the agent can predict how the world might evolve in the next time step.

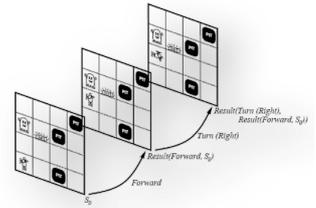
- a **belief state** represents which states of the world are currently possible (by explicit enumeration of states or by logical formulas)
- the probability theory allows to quantify the degree of belief in elements of the believe state
- we can also describe probability of state transitions

### **Probabilistic reasoning over time**

- representation of state transitions
- basic inference tasks
- inference algorithms for temporal models
- specific kinds of models (hidden Markov models, dynamic Bayesian networks)



In **situation calculus**, we view the world as a series of snapshots (**time slices**). A similar approach can be applied in probabilistic reasoning.



Each time slice (**state**) is described as a set of random variables:

- **hidden** (not observable) random **variables**  $X_t$
- **observable** random **variables**  $E_t$  (with observed values  $e_t$ )

$t$  is an identification of the time slice (we assume **discrete time** with uniform time steps)

Notation:

- $X_{a:b}$  denotes a set of variables from  $X_a$  to  $X_b$

### *A model example (umbrella world)*

You are the security guard stationed at a secret underground installation and you want to know whether it is **raining today**:

- hidden random variable  $R_t$

But your only access to the outside world occurs each morning when you see the the director coming in **with, or without, an umbrella**.

- observable random variable  $U_t$



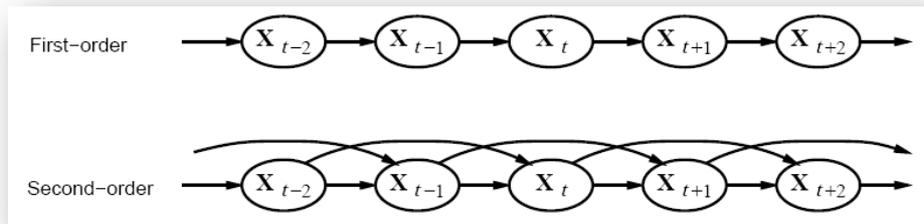
The **transition model** specifies the probability distribution over the latest state variables given the previous values.

This is given by  $P(\mathbf{X}_t | \mathbf{X}_{0:t-1})$ .

**Problem #1:** the set  $\mathbf{X}_{0:t-1}$  is unbounded in size as  $t$  increases

- we can make a **Markov assumption** – the current state depends only on a finite fixed number of previous states; processes satisfying this assumption are called **Markov processes** or **Markov chains**
- **first-order Markov chain** – the current state depends only on the previous state

$$P(\mathbf{X}_t | \mathbf{X}_{0:t-1}) = P(\mathbf{X}_t | \mathbf{X}_{t-1})$$



**Problem #2:** there are infinitely many possible values of  $t$

- We assume that changes in the world state are caused by a **stationary process** (a process of change is governed by laws that do not themselves change very time)
- the conditional probability tables  $P(\mathbf{X}_t | \mathbf{X}_{t-1})$  are identical for all  $t$

A **sensor (observation) model** describes how the evidence (observed) variables  $\mathbf{E}_t$  depend on other variables.

They could depend on previous variables as well as the current state variables.

We make a **sensor Markov assumption** – the evidence variables depend only on the hidden state variables  $\mathbf{X}_t$  from the same time.

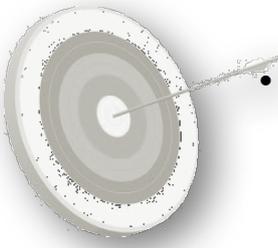
$$P(\mathbf{E}_t | \mathbf{X}_{0:t}, \mathbf{E}_{1:t-1}) = P(\mathbf{E}_t | \mathbf{X}_t)$$



The first-order Markov assumption says that the state variables contain all the information needed to characterize the probability distribution for the next time slice.

What if this assumption is only approximate?

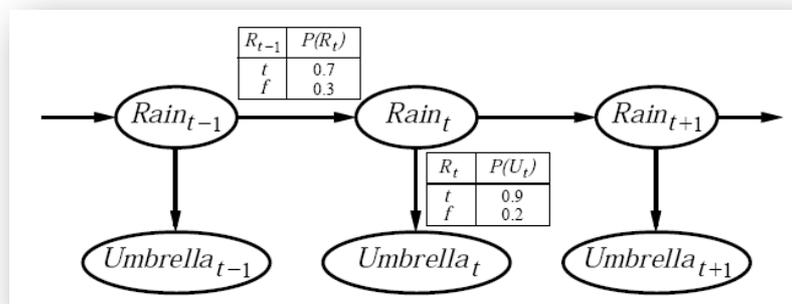
- **increase the order** of the Markov process model
- **increase the set** of state variables
  - For example we could add  $\text{Season}_t$  to incorporate historical records or we could add  $\text{Temperature}_t$ ,  $\text{Humidity}_t$ ,  $\text{Pressure}_t$  to use a physical model of rainy conditions.
  - The first solution (increasing the order) can always be reformulated as an increase in set of state variables.



*A Bayesian network view*

The transition and sensor models can be described using a **Bayesian network**.

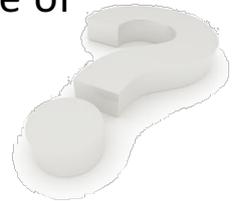
In addition to  $\mathbf{P}(\mathbf{X}_t | \mathbf{X}_{t-1})$  and  $\mathbf{P}(\mathbf{E}_t | \mathbf{X}_t)$  we need to say how everything gets started  $\mathbf{P}(\mathbf{X}_0)$ .



We have a specification of the complete joint distribution:

$$\mathbf{P}(\mathbf{X}_{0:t}, \mathbf{E}_{1:t}) = \mathbf{P}(\mathbf{X}_0) \prod_i \mathbf{P}(\mathbf{X}_i | \mathbf{X}_{i-1}) \mathbf{P}(\mathbf{E}_i | \mathbf{X}_i)$$

- **Filtering:** the task of computing the posterior distribution over the *most recent state*, given all evidence to date  $P(\mathbf{X}_t | \mathbf{e}_{1:t})$
- **Prediction:** the task of computing the posterior distribution over the *future state*, given all evidence to date  $P(\mathbf{X}_{t+k} | \mathbf{e}_{1:t})$  for  $k > 0$
- **Smoothing:** the task of computing posterior distribution over a *past state*, given all evidence up to the present  $P(\mathbf{X}_k | \mathbf{e}_{1:t})$  for  $k: 0 \leq k < t$
- **Most likely explanation:** the task to find the sequence of states that is most likely generated a given sequence of observations  $\text{argmax}_{\mathbf{x}_{1:t}} P(\mathbf{x}_{1:t} | \mathbf{e}_{1:t})$



## Filtering

The task of computing the posterior distribution over the *most recent state*, given all evidence to date –  $P(\mathbf{X}_t | \mathbf{e}_{1:t})$ .

A useful filtering algorithm needs to maintain a current state estimate and update it, rather than going back over (**recursive estimation**):

$$P(\mathbf{X}_{t+1} | \mathbf{e}_{1:t+1}) = f(\mathbf{e}_{t+1}, P(\mathbf{X}_t | \mathbf{e}_{1:t}))$$

How to define the function  $f$ ?

$$\begin{aligned} P(\mathbf{X}_{t+1} | \mathbf{e}_{1:t+1}) &= P(\mathbf{X}_{t+1} | \mathbf{e}_{1:t}, \mathbf{e}_{t+1}) \\ &= \alpha P(\mathbf{e}_{t+1} | \mathbf{X}_{t+1}, \mathbf{e}_{1:t}) P(\mathbf{X}_{t+1} | \mathbf{e}_{1:t}) \\ &= \alpha P(\mathbf{e}_{t+1} | \mathbf{X}_{t+1}) P(\mathbf{X}_{t+1} | \mathbf{e}_{1:t}) \\ &= \alpha P(\mathbf{e}_{t+1} | \mathbf{X}_{t+1}) \sum_{\mathbf{x}_t} P(\mathbf{X}_{t+1} | \mathbf{x}_t, \mathbf{e}_{1:t}) P(\mathbf{x}_t | \mathbf{e}_{1:t}) \\ &= \alpha P(\mathbf{e}_{t+1} | \mathbf{X}_{t+1}) \sum_{\mathbf{x}_t} P(\mathbf{X}_{t+1} | \mathbf{x}_t) P(\mathbf{x}_t | \mathbf{e}_{1:t}) \end{aligned}$$

Bayes rule

sensor Markov assumption

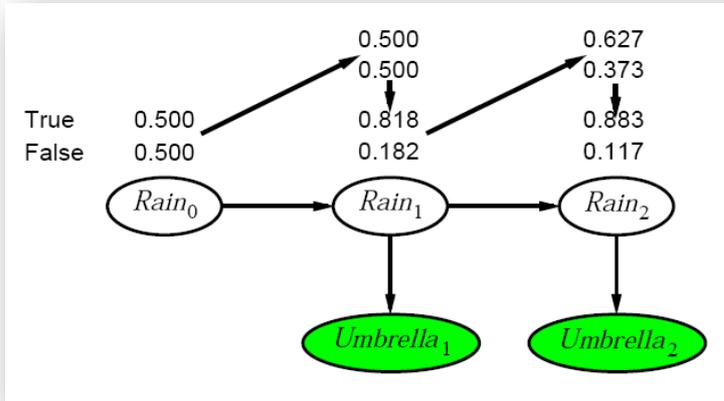
conditioning  
 $P(\mathbf{Y}) = \sum_{\mathbf{z}} P(\mathbf{Y} | \mathbf{z}) P(\mathbf{z})$

A message  $\mathbf{f}_{1:t}$  is propagated forward over the sequence:

$$\begin{aligned} P(\mathbf{X}_t | \mathbf{e}_{1:t}) &= \mathbf{f}_{1:t} \\ \mathbf{f}_{1:t+1} &= \alpha \text{FORWARD}(\mathbf{f}_{1:t}, \mathbf{e}_{t+1}) \\ \mathbf{f}_{1:0} &= P(\mathbf{X}_0) \end{aligned}$$



$$P(R_{t+1} | u_{1:t+1}) = \alpha P(u_{t+1} | R_{t+1}) P(R_{t+1} | u_{1:t}) = \alpha P(u_{t+1} | R_{t+1}) \sum_{r_t} P(R_{t+1} | r_t) P(r_t | u_{1:t})$$



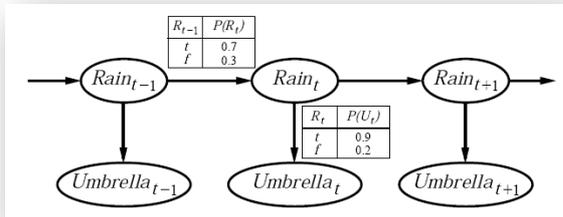
$$P(R_0) = \langle 0.5, 0.5 \rangle$$

$$P(R_1) = \sum_{r_0} P(R_1 | r_0) P(r_0) = \langle 0.5, 0.5 \rangle$$

$$P(R_1 | u_1) = \alpha P(u_1 | R_1) P(R_1) = \alpha \langle 0.9, 0.2 \rangle \langle 0.5, 0.5 \rangle \approx \langle 0.818, 0.182 \rangle$$

$$P(R_2 | u_1) = \sum_{r_1} P(R_2 | r_1) P(r_1 | u_1) = \langle 0.7, 0.3 \rangle \times 0.818 + \langle 0.3, 0.7 \rangle \times 0.182 \approx \langle 0.627, 0.372 \rangle$$

$$P(R_2 | u_1, u_2) = \alpha P(u_2 | R_2) P(R_2 | u_1) = \alpha \langle 0.9, 0.2 \rangle \langle 0.627, 0.372 \rangle = \langle 0.883, 0.117 \rangle$$



## Prediction

The task of computing the posterior distribution over the *future state*, given all evidence to date –  $P(X_{t+k} | e_{1:t})$  for some  $k > 0$ .

We can see this task as filtering without the addition of new evidence:

$$P(X_{t+k+1} | e_{1:t}) = \sum_{x_{t+k}} P(X_{t+k+1} | x_{t+k}) P(x_{t+k} | e_{1:t})$$

After some time (**mixing time**) the predicted distribution converges to the **stationary distribution** of the Markov process and remains constant.



The task of computing posterior distribution over a *past state*, given all evidence up to the present –  $P(\mathbf{X}_k | \mathbf{e}_{1:t})$  for  $k: 0 \leq k < t$ .

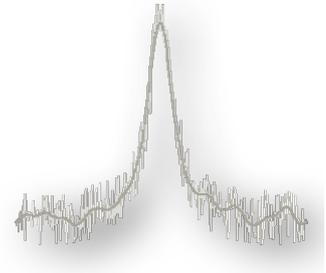
We again exploit a recursive message-passing approach, now in two parts.

$$\begin{aligned}
 P(\mathbf{X}_k | \mathbf{e}_{1:t}) &= P(\mathbf{X}_k | \mathbf{e}_{1:k}, \mathbf{e}_{k+1:t}) && \text{Bayes rule} \\
 &= \alpha P(\mathbf{X}_k | \mathbf{e}_{1:k}) P(\mathbf{e}_{k+1:t} | \mathbf{X}_k, \mathbf{e}_{1:k}) && \text{conditional independence} \\
 &= \alpha P(\mathbf{X}_k | \mathbf{e}_{1:k}) P(\mathbf{e}_{k+1:t} | \mathbf{X}_k) \\
 &= \alpha \mathbf{f}_{1:k} \times \mathbf{b}_{k+1:t}
 \end{aligned}$$

$$\begin{aligned}
 P(\mathbf{e}_{k+1:t} | \mathbf{X}_k) &= \sum_{\mathbf{x}_{k+1}} P(\mathbf{e}_{k+1:t} | \mathbf{X}_k, \mathbf{x}_{k+1}) P(\mathbf{x}_{k+1} | \mathbf{X}_k) && \text{conditioning} \\
 &= \sum_{\mathbf{x}_{k+1}} P(\mathbf{e}_{k+1:t} | \mathbf{x}_{k+1}) P(\mathbf{x}_{k+1} | \mathbf{X}_k) && \text{conditional independence} \\
 &= \sum_{\mathbf{x}_{k+1}} P(\mathbf{e}_{k+1}, \mathbf{e}_{k+2:t} | \mathbf{x}_{k+1}) P(\mathbf{x}_{k+1} | \mathbf{X}_k) \\
 &= \sum_{\mathbf{x}_{k+1}} P(\mathbf{e}_{k+1} | \mathbf{x}_{k+1}) P(\mathbf{e}_{k+2:t} | \mathbf{x}_{k+1}) P(\mathbf{x}_{k+1} | \mathbf{X}_k) && \text{conditional independence}
 \end{aligned}$$

Using the backward message-passing notation:

$$\begin{aligned}
 P(\mathbf{e}_{k+1:t} | \mathbf{X}_k) &= \mathbf{b}_{k+1:t} \\
 \mathbf{b}_{k+1:t} &= \text{BACKWARD}(\mathbf{b}_{k+2:t}, \mathbf{e}_{k+1}) \\
 \mathbf{b}_{t+1:t} &= P(\mathbf{e}_{t+1:t} | \mathbf{X}_t) = P(| \mathbf{X}_t) = \mathbf{1}
 \end{aligned}$$



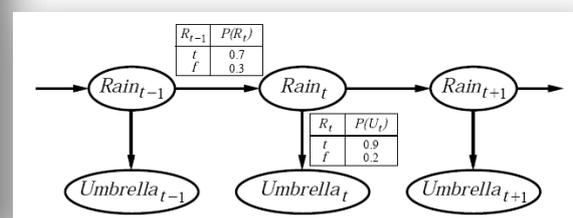
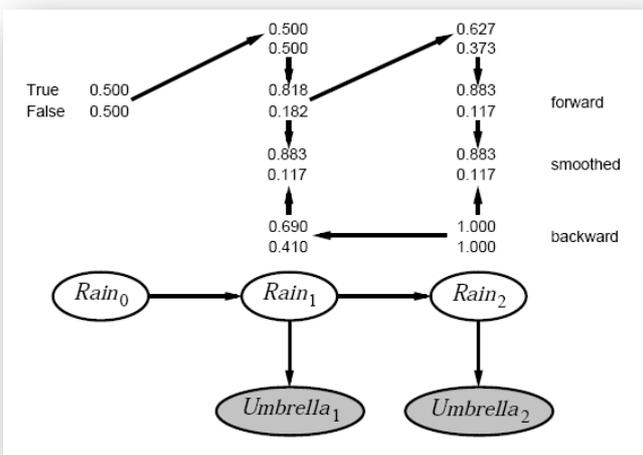
## Smoothing (example)

$$P(R_k | u_{1:t+1}) = \alpha P(R_k | u_{1:k}) P(u_{k+1:t} | R_k)$$

$$P(u_{k+1:t} | R_k) = \sum_{r_{k+1}} P(u_{k+1} | r_{k+1}) P(u_{k+2:t} | r_{k+1}) P(r_{k+1} | R_k)$$

$$P(| R_2) = \mathbf{1}$$

$$\begin{aligned}
 P(u_2 | R_1) &= \sum_{r_2} P(u_2 | r_2) P(| r_2) P(r_2 | R_1) \\
 &= 0.9 \times 1 \times \langle 0.7, 0.3 \rangle + 0.2 \times 1 \times \langle 0.3, 0.7 \rangle = \langle 0.69, 0.41 \rangle
 \end{aligned}$$



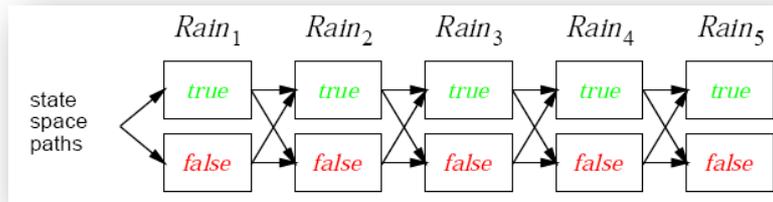
## Most likely explanation/sequence

The task to find the sequence of states that is most likely generated a given sequence of observations

$$\operatorname{argmax}_{x_{1:t}} P(x_{1:t} \mid e_{1:t}).$$

This is different from smoothing for each past state and taking the sequence of most probable states!

We can see each sequence as a **path through a graph** whose nodes are possible states at each time step:



Because of the Markov property the most likely path to a given state consists of the most likely path to some previous state followed by a transition to that state.

This can be described using a **recursive formula**.

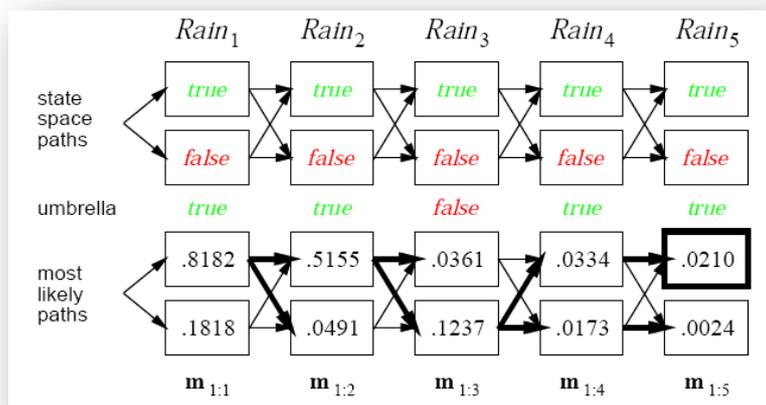
## Viterbi algorithm

The most likely path to a given state consists of the most likely path to some previous state followed by a transition to that state.

$$\begin{aligned} & \max_{x_1, \dots, x_t} P(x_1, \dots, x_t, X_{t+1} \mid e_{1:t+1}) \\ & = \alpha P(e_{t+1} \mid X_{t+1}) \max_{x_t} (P(X_{t+1} \mid x_t) \max_{x_1, \dots, x_{t-1}} P(x_1, \dots, x_t \mid e_{1:t})) \end{aligned}$$

Again, we use an approach of forward message passing:

$$\begin{aligned} m_{1:t} &= \max_{x_1, \dots, x_{t-1}} P(x_1, \dots, x_{t-1}, X_t \mid e_{1:t}) \\ m_{1:t+1} &= P(e_{t+1} \mid X_{t+1}) \max_{x_t} (P(X_{t+1} \mid x_t) m_{1:t}) \end{aligned}$$



Assume that the state of process is described by a single discrete random variable  $X_t$  (there is also a single evidence variable  $E_t$ ).

This is called a **hidden Markov model** (HMM).

This restricted model allows for a simple and elegant **matrix implementation** of all the basic algorithms.

Assume that variable  $X_t$  takes values from the set  $\{1, \dots, S\}$ , where  $S$  is the number of possible states.

The **transition model**  $P(X_t | X_{t-1})$  becomes an  $S \times S$  matrix  $\mathbf{T}$ , where:

$$\mathbf{T}_{(i,j)} = P(X_t = j | X_{t-1} = i)$$

We also put the **sensor model** in matrix form. Now we know the value of the evidence variable  $e_t$  so we describe  $P(E_t = e_t | X_t = i)$ , using a diagonal matrix  $\mathbf{O}_t$ , where:

$$\mathbf{O}_{t(i,i)} = P(E_t = e_t | X_t = i)$$

## Matrix formulation of algorithms

### The forward message propagation (from filtering)

$$P(X_t | \mathbf{e}_{1:t}) = \mathbf{f}_{1:t}$$

$$\mathbf{f}_{1:t+1} = \alpha P(\mathbf{e}_{t+1} | \mathbf{X}_{t+1}) \sum_{\mathbf{x}_t} P(\mathbf{X}_{t+1} | \mathbf{x}_t) P(\mathbf{x}_t | \mathbf{e}_{1:t})$$

can be reformulated using matrix operations (message  $\mathbf{f}_{1:t}$  is modelled as a one-column matrix) as follows:

$$\mathbf{T}_{(i,j)} = P(X_t = j | X_{t-1} = i)$$

$$\mathbf{O}_{t(i,i)} = P(E_t = e_t | X_t = i)$$

$$\mathbf{f}_{1:t+1} = \alpha \mathbf{O}_{t+1} \mathbf{T}^T \mathbf{f}_{1:t}$$

### The backward message propagation (from smoothing)

$$P(\mathbf{e}_{k+1:t} | \mathbf{X}_k) = \mathbf{b}_{k+1:t}$$

$$\mathbf{b}_{k+1:t} = \sum_{\mathbf{x}_{k+1}} P(\mathbf{e}_{k+1} | \mathbf{x}_{k+1}) P(\mathbf{e}_{k+2:t} | \mathbf{x}_{k+1}) P(\mathbf{x}_{k+1} | \mathbf{X}_k)$$

can be reformulated using matrix operations (message  $\mathbf{b}_{k:t}$  is modelled as a one-column matrix) as follows:

$$\mathbf{b}_{k+1:t} = \mathbf{T} \mathbf{O}_{k+1} \mathbf{b}_{k+2:t}$$

What if we need to **smooth the whole sequence of states?**

$$P(\mathbf{X}_k | \mathbf{e}_{1:t}) = \alpha \mathbf{f}_{1:k} \times \mathbf{b}_{k+1:t}$$

The time complexity of smoothing with respect to evidence  $\mathbf{e}_{1:t}$  is  $O(t)$

One obvious method to smooth the whole sequence is to run the smoothing algorithm for each time step – this results in time complexity  $O(t^2)$ .

A better approach uses **dynamic programming** (reuse already computed information) reducing the time complexity to  $O(t)$ .

- **forward-backward algorithm**
- the practical drawback of this approach is that its space complexity can be too high – it is  $O(|\mathbf{f}|t)$ .

## Forward-backward algorithm

```

function FORWARD-BACKWARD(ev, prior) returns a vector of probability distributions
  inputs: ev, a vector of evidence values for steps  $1, \dots, t$ 
           prior, the prior distribution on the initial state,  $P(\mathbf{X}_0)$ 
  local variables: fv, a vector of forward messages for steps  $0, \dots, t$ 
                    b, a representation of the backward message, initially all 1s
                    sv, a vector of smoothed estimates for steps  $1, \dots, t$ 

  fv[0]  $\leftarrow$  prior
  for  $i = 1$  to  $t$  do
    fv[ $i$ ]  $\leftarrow$  FORWARD(fv[ $i - 1$ ], ev[ $i$ ])
  for  $i = t$  downto  $1$  do
    sv[ $i$ ]  $\leftarrow$  NORMALIZE(fv[ $i$ ]  $\times$  b)
    b  $\leftarrow$  BACKWARD(b, ev[ $i$ ])
  return sv

```

Can be smoothing the whole sequence of states done with smaller memory consumption while keeping the time complexity  $O(t)$ ?

**Ideas:**

- For message-passing in one direction we need constant space independent of  $t$ .
- Can the message  $\mathbf{f}_{1:t}$  be obtained from the message  $\mathbf{f}_{1:t+1}$ ?
- Then we can pass the forward message in the reverse (backward) direction together with the backward message.

Let us exploit **matrix operations**:

$$\mathbf{f}_{1:t+1} = \alpha \mathbf{O}_{t+1} \mathbf{T}^T \mathbf{f}_{1:t} \quad \rightarrow \quad \mathbf{f}_{1:t} = \alpha' (\mathbf{T}^T)^{-1} (\mathbf{O}_{t+1})^{-1} \mathbf{f}_{1:t+1}$$

**Algorithm:**

- first, run the forward-message propagation to get  $\mathbf{f}_{1:t}$
- then during the backward stage compute both  $\mathbf{f}_{1:k}$  and  $\mathbf{b}_{k+1:t}$

Smoothing with a fixed time lag

Assume smoothing in an on-line setting where smoothed estimates must be computed for a fixed number  $d$  of back time steps –  $\mathbf{P}(\mathbf{X}_{t-d} | \mathbf{e}_{1:t})$ . This is called **fixed-lag smoothing**.

In the ideal case, we want incremental computation in a constant time per update.

we have  $\mathbf{P}(\mathbf{X}_{t-d} | \mathbf{e}_{1:t}) = \alpha \mathbf{f}_{1:t-d} \times \mathbf{b}_{t-d+1:t}$   
 we need  $\mathbf{P}(\mathbf{X}_{t-d+1} | \mathbf{e}_{1:t+1}) = \alpha \mathbf{f}_{1:t-d+1} \times \mathbf{b}_{t-d+2:t+1}$

An incremental approach:

- we can use  $\mathbf{f}_{1:t-d+1} = \alpha \mathbf{O}_{t-d+2} \mathbf{T}^T \mathbf{f}_{1:t-d}$
- we need incremental computation of  $\mathbf{b}_{t-d+2:t+1}$  from  $\mathbf{b}_{t-d+1:t}$

$$\mathbf{b}_{t-d+1:t} = \mathbf{T} \mathbf{O}_{t-d+1} \mathbf{b}_{t-d+2:t} = (\prod_{i=t-d+1, \dots, t} \mathbf{T} \mathbf{O}_i) \mathbf{b}_{t+1:t} = \mathbf{B}_{t-d+1:t} \mathbf{1}$$

$$\mathbf{b}_{t-d+2:t+1} = (\prod_{i=t-d+2, \dots, t+1} \mathbf{T} \mathbf{O}_i) \mathbf{b}_{t+2:t+1} = \mathbf{B}_{t-d+2:t+1} \mathbf{1}$$

$$\mathbf{B}_{t-d+2:t+1} = (\mathbf{O}_{t-d+1})^{-1} \mathbf{T}^{-1} \mathbf{B}_{t-d+1:t} \mathbf{T} \mathbf{O}_{t+1}$$

```

function FIXED-LAG-SMOOTHING( $e_t$ , hmm,  $d$ ) returns a distribution over  $\mathbf{X}_{t-d}$ 
  inputs:  $e_t$ , the current evidence for time step  $t$ 
            hmm, a hidden Markov model with  $S \times S$  transition matrix  $T$ 
             $d$ , the length of the lag for smoothing
  static:  $t$ , the current time, initially 1
             $\mathbf{f}$ , a probability distribution, the forward message  $\mathbf{P}(X_t|e_{1:t})$ , initially  $\text{PRIOR}[hmm]$ 
             $B$ , the  $d$ -step backward transformation matrix, initially the identity matrix
             $e_{t-d:t}$ , double-ended list of evidence from  $t - d$  to  $t$ , initially empty
  local variables:  $\mathbf{O}_{t-d}$ ,  $\mathbf{O}_t$ , diagonal matrices containing the sensor model information

  add  $e_t$  to the end of  $e_{t-d:t}$ 
   $\mathbf{O}_t \leftarrow$  diagonal matrix containing  $\mathbf{P}(e_t|X_t)$ 
  if  $t > d$  then
     $\mathbf{f} \leftarrow$  FORWARD( $\mathbf{f}$ ,  $e_t$ )
    remove  $e_{t-d-1}$  from the beginning of  $e_{t-d:t}$ 
     $\mathbf{O}_{t-d} \leftarrow$  diagonal matrix containing  $\mathbf{P}(e_{t-d}|X_{t-d})$ 
     $B \leftarrow \mathbf{O}_{t-d}^{-1} \mathbf{T}^{-1} \mathbf{B} \mathbf{O}_t$ 
  else  $B \leftarrow \mathbf{B} \mathbf{O}_t$ 
   $t \leftarrow t + 1$ 
  if  $t > d$  then return NORMALIZE( $\mathbf{f} \times \mathbf{B}1$ ) else return null
  
```



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