# Artificial Intelligence

#### Roman Barták

Department of Theoretical Computer Science and Mathematical Logic

## 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.

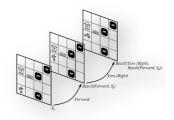
- a belief state represents which states of the world are currently possible (by explicit enumeration of states or by logical formulas)
- the probability theory allows to quantify the degree of belief in elements of the believe state
- we can also describe probability of state transitions

#### 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<sub>t</sub>
- observable random variables  $\mathbf{E}_{t}$  (with observed values  $\mathbf{e}_{t}$ )
- 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<sub>t</sub>

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

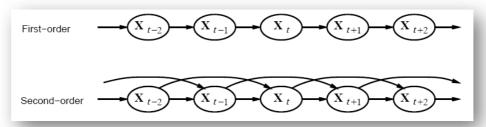
observable random variable U<sub>t</sub>

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

This is given by  $P(X_t | X_{0:t-1})$ .

**Problem #1**: the set  $X_{0:t-1}$  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  $P(X_t \mid X_{0:t-1}) = P(X_t \mid X_{t-1})$



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  $P(X_t | X_{t-1})$  are identical for all t

Sensor model

A **sensor (observation) model** describes how the evidence (observed) variables **E**<sub>t</sub> depend on other variables.

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

We make a **sensor Markov assumption** – the evidence variables depend only on the hidden state variables **X**<sub>t</sub> from the same time.

$$P(E_{t} | X_{0:t}, E_{1:t-1}) = P(E_{t} | X_{t})$$



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.

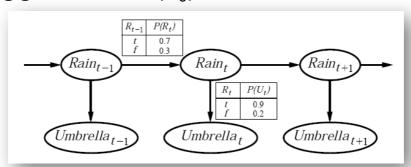
What if this assumption is only approximate?

- increase the order of the Markov process model
- increase the set of state variables
  - For example we could add Season<sub>t</sub> to incorporate historical records or we could add Temperature<sub>t</sub>, Humidity<sub>t</sub>, Pressure<sub>t</sub> 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  $P(X_t \mid X_{t-1})$  and  $P(E_t \mid X_t)$  we need to say how everything gets started  $P(X_0)$ .



We have a specification of the complete joint distribution:

$$P(X_{0:t}, E_{1:t}) = P(X_0) \prod_i P(X_i \mid X_{i-1}) P(E_i \mid X_i)$$

- Filtering: the task of computing the posterior distribution over the most recent state, given all evidence to date
   P(X<sub>t</sub> | e<sub>1:t</sub>)
- Prediction: the task of computing the posterior distribution over the future state, given all evidence to date
   P(X<sub>t+k</sub> | e<sub>1:t</sub>) for k>0
- Smoothing: the task of computing posterior distribution over a past state, given all evidence up to the present
   P(X<sub>k</sub> | e<sub>1:t</sub>) for k: 0 ≤ k < t</li>
- Most likely explanation: the task to find the sequence of states that is most likely generated a given sequence of observations

$$\operatorname{argmax}_{\mathbf{x}_{1:t}} P(\mathbf{x}_{1:t} \mid \mathbf{e}_{1:t})$$

#### Filtering

The task of computing the posterior distribution over the most recent state, given all evidence to date  $-P(X_t|e_{1:t})$ .

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

$$P(X_{t+1} | e_{1:t+1}) = f(e_{t+1}, P(X_t | e_{1:t}))$$

How to define the function f?

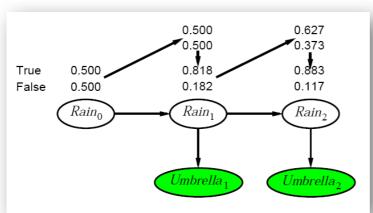
$$\begin{split} & \textbf{P}(\textbf{X}_{t+1} | \textbf{e}_{1:t+1}) = \textbf{P}(\textbf{X}_{t+1} | \textbf{e}_{1:t}, \textbf{e}_{t+1}) \\ & = \alpha \ \textbf{P}(\textbf{e}_{t+1} | \textbf{X}_{t+1}, \textbf{e}_{1:t}) \ \textbf{P}(\textbf{X}_{t+1} | \textbf{e}_{1:t}) \\ & = \alpha \ \textbf{P}(\textbf{e}_{t+1} | \textbf{X}_{t+1}) \ \textbf{P}(\textbf{X}_{t+1} | \textbf{e}_{1:t}) \\ & = \alpha \ \textbf{P}(\textbf{e}_{t+1} | \textbf{X}_{t+1}) \ \boldsymbol{\Sigma}_{\textbf{X}_t} \ \textbf{P}(\textbf{X}_{t+1} | \textbf{x}_t, \textbf{e}_{1:t}) \ \textbf{P}(\textbf{X}_t | \textbf{e}_{1:t}) \\ & = \alpha \ \textbf{P}(\textbf{e}_{t+1} | \textbf{X}_{t+1}) \ \boldsymbol{\Sigma}_{\textbf{X}_t} \ \textbf{P}(\textbf{X}_{t+1} | \textbf{x}_t) \ \textbf{P}(\textbf{X}_t | \textbf{e}_{1:t}) \end{split}$$

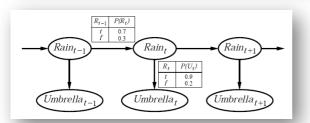
A message  $\mathbf{f}_{1:t}$  is propagated forward over the sequence:

$$P(X_t | e_{1:t}) = f_{1:t}$$
  
 $f_{1:t+1} = \alpha FORWARD(f_{1:t}, e_{t+1})$   
 $f_{1:0} = P(X_0)$ 



$$\begin{split} & \mathsf{P}(\mathsf{R}_{t+1} \,|\, \mathsf{u}_{1:t+1}) \\ & = \alpha \; \mathsf{P}(\mathsf{u}_{t+1} \,|\, \mathsf{R}_{t+1}) \; \mathsf{P}(\mathsf{R}_{t+1} \,|\, \mathsf{u}_{1:t}) = \alpha \; \mathsf{P}(\mathsf{u}_{t+1} \,|\, \mathsf{R}_{t+1}) \; \Sigma_{\mathsf{r}_{\mathsf{t}}} \; \mathsf{P}(\mathsf{R}_{t+1} \,|\, \mathsf{r}_{\mathsf{t}}) \; \mathsf{P}(\mathsf{r}_{\mathsf{t}} \,|\, \mathsf{u}_{1:t}) \end{split}$$





$$\begin{split} \mathbf{P}(\mathbf{R}_0) &= \langle 0.5, 0.5 \rangle \\ \mathbf{P}(\mathbf{R}_1) \\ &= \sum_{\mathbf{r}_0} \mathbf{P}(\mathbf{R}_1 | \mathbf{r}_0) \, \mathbf{P}(\mathbf{r}_0) \\ &= \langle 0.5, 0.5 \rangle \\ \mathbf{P}(\mathbf{R}_1 | \mathbf{u}_1) \\ &= \alpha \, \mathbf{P}(\mathbf{u}_1 | \mathbf{R}_1) \, \mathbf{P}(\mathbf{R}_1) \\ &= \alpha \, \langle 0.9, 0.2 \rangle \langle 0.5, 0.5 \rangle \\ &\approx \langle 0.818, 0.182 \rangle \\ \mathbf{P}(\mathbf{R}_2 | \mathbf{u}_1) \\ &= \sum_{\mathbf{r}_1} \mathbf{P}(\mathbf{R}_2 | \mathbf{r}_1) \, \mathbf{P}(\mathbf{r}_1 | \mathbf{u}_1) \\ &= \langle 0.7, 0.3 \rangle \times 0.818 \\ &+ \langle 0.3, 0.7 \rangle \times 0.182 \\ &\approx \langle 0.627, 0.372 \rangle \\ \mathbf{P}(\mathbf{R}_2 | \mathbf{u}_1, \mathbf{u}_2) \\ &= \alpha \, \mathbf{P}(\mathbf{u}_2 | \mathbf{R}_2) \, \mathbf{P}(\mathbf{R}_2 | \mathbf{u}_1) \\ &= \alpha \, \langle 0.9, 0.2 \rangle \langle 0.627, 0.372 \rangle \\ &= \langle 0.883, 0.117 \rangle \end{split}$$

#### 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.

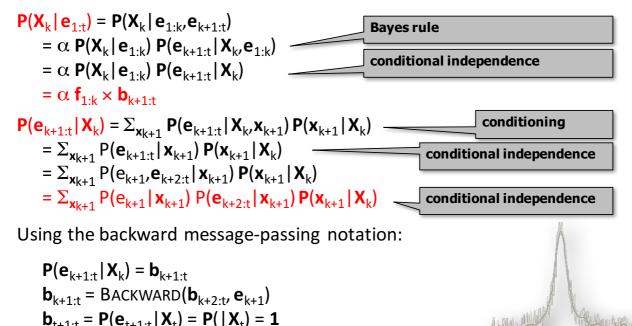
We can see this task as filtering without the addition of new evidence:

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

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

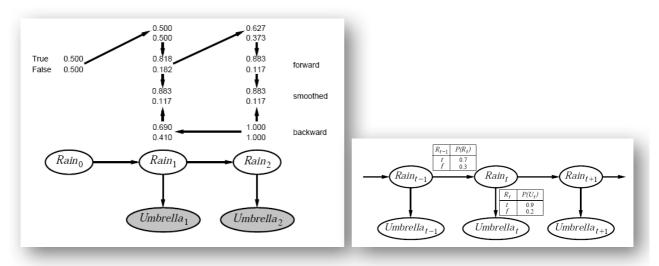
The task of computing posterior distribution over a *past state*, given all evidence up to the present  $-\mathbf{P}(\mathbf{X}_k|\mathbf{e}_{1:t})$  for k:  $0 \le k < t$ .

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



# Smoothing (example)

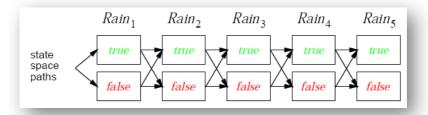
$$\begin{split} \textbf{P}(R_k \,|\, u_{1:t+1}) &= \alpha \; \textbf{P}(R_k \,|\, u_{1:k}) \; \textbf{P}(u_{k+1:t} \,|\, R_k) \\ P(u_{k+1:t} \,|\, R_k) &= \Sigma_{r_{k+1}} \; P(u_{k+1} \,|\, r_{k+1}) \; P(u_{k+2:t} \,|\, r_{k+1}) \; \textbf{P}(r_{k+1} \,|\, R_k) \\ \textbf{P}(\,|\, R_2) &= \textbf{1} \\ \textbf{P}(u_2 \,|\, R_1) &= \Sigma_{r_2} \; P(u_2 \,|\, r_2) \; P(\,|\, r_2) \; \textbf{P}(r_2 \,|\, R_1) \\ &= 0.9 \times 1 \times \langle 0.7, 0.3 \rangle + 0.2 \times 1 \times \langle 0.3, 0.7 \rangle = \langle 0.69, 0.41 \rangle \end{split}$$



The task to find the sequence of states that is most likely generated a given sequence of observations  $argmax_{x_{1:t}} P(x_{1:t} | e_{1:t})$ .

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.

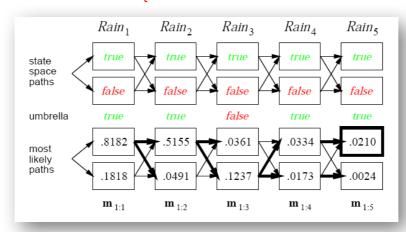
#### Viterbi algorithm

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.

$$\begin{aligned} & \max_{\mathbf{x}_1,\dots,\mathbf{x}_t} \ \mathbf{P}(\mathbf{x}_1,\dots,\mathbf{x}_t,\mathbf{X}_{t+1} | \ \mathbf{e}_{1:t+1}) \\ & = \alpha \ \mathbf{P}(\mathbf{e}_{t+1} | \mathbf{X}_{t+1}) \ \max_{\mathbf{x}_t} \ (\mathbf{P}(\mathbf{X}_{t+1} | \ \mathbf{x}_t) \ \max_{\mathbf{x}_1,\dots,\mathbf{x}_{t-1}} \ \mathbf{P}(\mathbf{x}_1,\dots,\mathbf{x}_t | \ \mathbf{e}_{1:t})) \end{aligned}$$

Again, we use an approach of forward message passing:

$$\begin{aligned} & m_{1:t} = \max_{\mathbf{x}_{1},...,\mathbf{x}_{t-1}} \mathbf{P}(\mathbf{x}_{1},...,\mathbf{x}_{t-1},\mathbf{X}_{t} | \mathbf{e}_{1:t}) \\ & m_{1:t+1} = \mathbf{P}(\mathbf{e}_{t+1} | \mathbf{X}_{t+1}) \max_{\mathbf{x}_{t}} (\mathbf{P}(\mathbf{X}_{t+1} | \mathbf{x}_{t}) m_{1:t}) \end{aligned}$$



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  $\{1,...S\}$ , where S is the number of possible states.

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

$$T_{(i,j)} = P(X_t = j \mid X_{t-1} = i)$$

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

$$O_{t(i,i)} = P(E_t = e_t | X_t = i)$$

### Matrix formulation of algorithms

### The forward message propagation (from filtering)

$$\begin{split} & \mathbf{P}(\mathbf{X}_{t} \,|\, \mathbf{e}_{1:t}) = \mathbf{f}_{1:t} \\ & \mathbf{f}_{1:t+1} = \alpha \, \mathbf{P}(\mathbf{e}_{t+1} \,|\, \mathbf{X}_{t+1}) \, \Sigma_{\mathbf{x}_{t}} \, \mathbf{P}(\mathbf{X}_{t+1} \,|\, \mathbf{x}_{t}) \, \, \mathbf{P}(\mathbf{x}_{t} \,|\, \mathbf{e}_{1:t}) \end{split}$$

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

$$T_{(i,j)} = P(X_t = j \mid X_{t-1} = i)$$
 $O_{t(i,i)} = P(E_t = e_t \mid X_t = i)$ 
 $f_{1:t+1} = \alpha O_{t+1} T^T f_{1:t}$ 

The backward message propagation (from smoothing)

$$\begin{aligned} \mathbf{P}(\mathbf{e}_{k+1:t} | \mathbf{X}_k) &= \mathbf{b}_{k+1:t} \\ \mathbf{b}_{k+1:t} &= \Sigma_{\mathbf{x}_{k+1}} \ \mathsf{P}(\mathbf{e}_{k+1} | \mathbf{x}_{k+1}) \ \mathsf{P}(\mathbf{e}_{k+2:t} | \mathbf{x}_{k+1}) \ \mathsf{P}(\mathbf{x}_{k+1} | \mathbf{X}_k) \\ \text{can be reformulated using matrix operations (message } \mathbf{b}_{k:t} \\ \text{is modelled as a one-column matrix) as follows:} \end{aligned}$$

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

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

$$P(X_k | e_{1:t}) = \alpha f_{1:k} \times b_{k+1:t}$$

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(|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, \ldots, t prior, the prior distribution on the initial state, P(X_0) local variables: fv, a vector of forward messages for steps 0, \ldots, t b, a representation of the backward message, initially all 1s sv, a vector of smoothed estimates for steps 1, \ldots, 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:

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

#### Algorithm:

- first, run the forward-message propagation to get f<sub>1:t</sub>
- 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 –  $P(X_{t-d}|e_{1:t})$ . 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  $\mathbf{f}_{1:t-d+1} = \alpha \mathbf{O}_{t-d+2} \mathbf{T}^\mathsf{T} \mathbf{f}_{1:t-d}$
- we need incremental computation of  $\mathbf{b}_{\text{t-d+2:t+1}}$  from  $\mathbf{b}_{\text{t-d+1:t}}$

$$\begin{aligned} & \mathbf{b}_{t-d+1:t} = \mathbf{T} \ \mathbf{O}_{t-d+1} \ \mathbf{b}_{t-d+2:t} = (\Pi_{i=t-d+1,...,t} \ \mathbf{T} \ \mathbf{O}_{i}) \ \mathbf{b}_{t+1:t} = \mathbf{B}_{t-d+1:t} \ \mathbf{1} \\ & \mathbf{b}_{t-d+2:t+1} = (\Pi_{i=t-d+2,...,t+1} \ \mathbf{T} \ \mathbf{O}_{i}) \ \mathbf{b}_{t+2:t+1} = \mathbf{B}_{t-d+2:t+1} \ \mathbf{1} \end{aligned}$$

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

# Smoothing with a fixed time lag

```
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 x S transition matrix T
             d, the length of the lag for smoothing
  static: t, the current time, initially 1
           f, a probability distribution, the forward message P(X_t|e_{1:t}), initially 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 \text{diagonal matrix containing } \mathbf{P}(e_t|X_t)
  if t \ge d then
       \mathbf{f} \leftarrow \text{FORWARD}(\mathbf{f}, e_t)
       remove e_{t-d-1} from the beginning of e_{t-d:t}
      \mathbf{O}_{t-d} \leftarrow \text{diagonal matrix containing } \mathbf{P}(e_{t-d}|X_{t-d})
\mathbf{B} \leftarrow \mathbf{O}_{t-d}^{-1} \mathbf{T}^{-1} \mathbf{B} \mathbf{T} \mathbf{O}_{t}
  else B \leftarrow \mathbf{BTO}_t
  t \leftarrow t + 1
  if t > d then return NORMALIZE(f x B1) else return null
```



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Department of Theoretical Computer Science and Mathematical Logic bartak@ktiml.mff.cuni.cz