# COMP9418: Advanced Topics in Statistical Machine Learning

## Markov Chains and Hidden Markov Models

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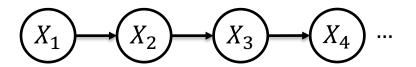
University of New South Wales

## Introduction

- This lecture discusses two classes of Graphical Models
  - Markov chains
  - Hidden Markov Models (HMM)
- Both models are instances of Dynamic Bayesian Networks (DBN)
  - They have a repeating structure that grows with time or space
  - Such structure is simple and uses the Markov property
- The Markov property states that future states are independent of past ones given the current state
  - In Markov chains, all states are observable
  - HMM extend the chains by allowing hidden states
- We will discuss specialised inference algorithms for both classes
  - Applications of these graphical models in domains such as robot localisation

# Time and Space

- Several problems require reasoning about sequences
  - Such sequences may represent the problem dynamics in time or space
  - Examples of applications are speech recognition and robot localization
- Dynamic Bayesian Networks (DBN) allow to incorporate time or space in our models
  - The two simplest instances of DBNs are Markov chains and Hidden Markov models



## **Markov Chains**

#### Markov chain is a state machine

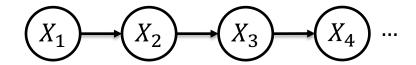
- X is a discrete variable and each value is called a state
- Transitions between states are nondeterministic

#### Parameters

- Prior probabilities  $P(X_1)$
- Transition probabilities or dynamics  $P(X_t|X_{t-1})$

#### Stationary assumption

- lacktriangle Transition probabilities are the same for all values of t
- Also known as a time-homogeneous chain



## Markov Chains: Weather

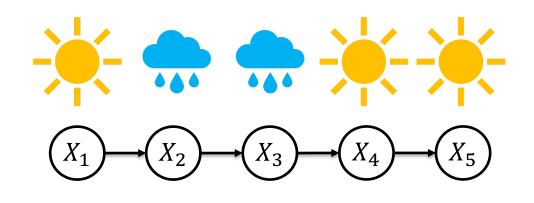
#### States

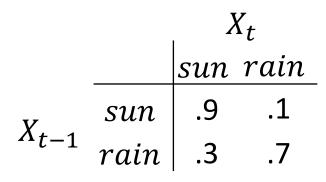
- $\bullet$   $X = \{sun, rain\}$
- Initial distribution

$$X_1 = \begin{pmatrix} 1 & sun \\ 0 & rain \end{pmatrix}$$

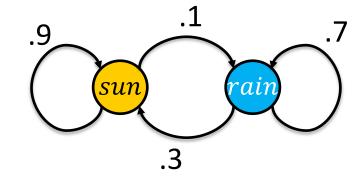
## Transition probabilities

$X_{t-1}$	$X_t$	$P(X_t X_{t-1})$
sun	sun	.9
sun	rain	.1
rain	sun	.3
rain	rain	.7





Matrix of transition probabilities



Transition or state diagram

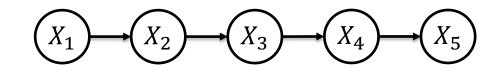
## Markov Chains: Independencies

- An relevant question is which independencies are implied in this chain
  - We can use d-separation to visually infer the independencies
  - This independence assumption is known as (first order) Markov property
- Independences are also apparent when we look at the chain rule
  - Chain rule if Bayesian networks for this example

$$P(X_1, X_2, X_3, X_4, X_5) = P(X_1)P(X_2|X_1)P(X_3|X_2)P(X_4|X_3)P(X_5|X_4)$$

• Chain rule in general  $P(X_1, X_2, X_3, X_4, X_5) = P(X_1)P(X_2|X_1)P(X_3|X_2, X_1)P(X_4|X_3, X_2, X_1)P(X_5|X_4, X_3, X_2, X_1)$ 

$$X_3 \perp X_1 \mid X_2 \qquad X_4 \perp X_1, X_2 \mid X_3 \qquad X_5 \perp X_1, X_2, X_3 \mid X_4$$



$$X_1 \perp X_3 | X_2$$

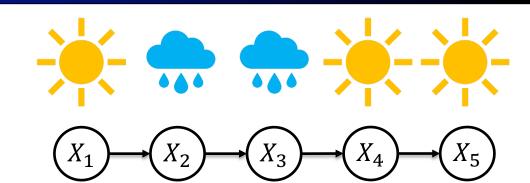
$$X_2 \perp X_4 | X_3$$

More generally,  $X_{t+1} \perp X_{t-1} | X_t$ 

# Markov Chains: Independencies

### In general

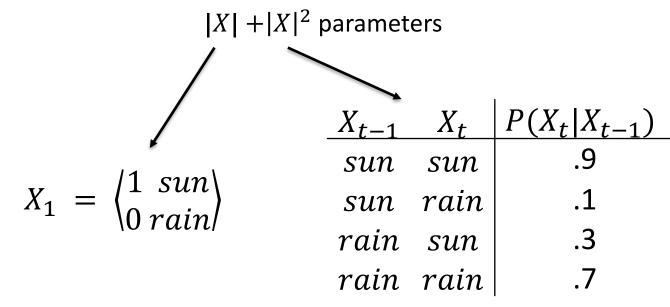
$$P(X_1, ..., X_n) = P(X_1) \prod_{i=2}^{n} P(X_t | X_{t-1})$$



$$|X|^n$$
 parameters

$$|X| + (n-1)|X|^2$$
 parameters  $\longrightarrow$ 

We also assume that  $P(X_t|X_{t-1})$  is the same for all t (stationarity)

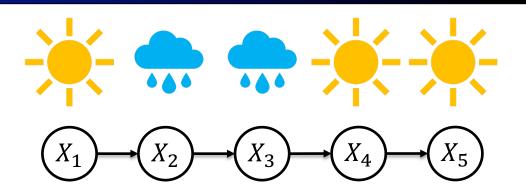


# Probability of a State Sequence

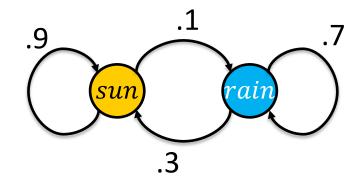
- The probability of a sequence is the product of the transition probabilities
  - This comes directly from the chain rule

$$P(X_1, X_2, X_3, X_4, X_5) = P(X_1)P(X_2|X_1)P(X_3|X_2)P(X_4|X_3)P(X_5|X_4)$$

$$P(sum, rain, rain, sun, sun) = 1(.1)(.7)(.3)(.9) = .189$$



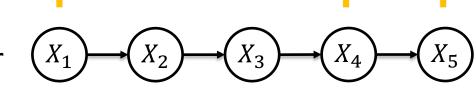
For example, what is the probability of the sequence: sun, rain, rain, sun, sun?



$$X_1 = \begin{pmatrix} 1 & sun \\ 0 & rain \end{pmatrix}$$

# Probability of Staying in a Certain State

- The probability of staying in a certain state for d steps
  - It is the probability of a sequence in this state for d-1 steps then going to a different state



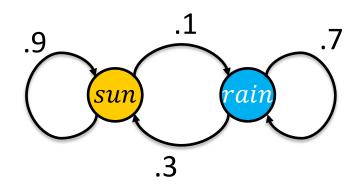
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$$S_S^d = \{X_i = s : 1 \le i \le d\}$$

$$P(\mathbf{S}_s^d) = P(s|s)^{d-1}(1 - P(s|s))$$

$$P(S_{rain}^3) = P(rain|rain)^{3-1}(1 - P(rain|rain)) = (.7^2)(1 - .7) = .147$$

For example, what is the probability of three raining days?



$$X_1 = \begin{pmatrix} 1 & sun \\ 0 & rain \end{pmatrix}$$

## Expected Time in a State

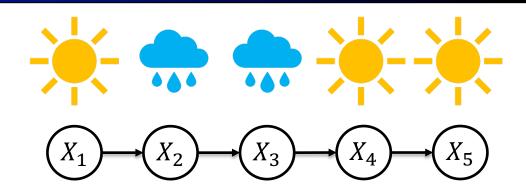
- The average duration of a sequence is a certain state
  - It is the expected number of time steps in that state

$$\mathbb{E}[S_s] = \sum_{i}^{\infty} P(S_s^i)i$$

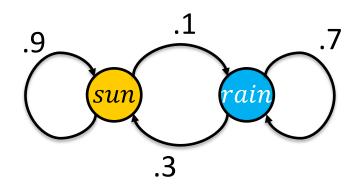
$$= \sum_{i}^{\infty} i P(s|s)^{i-1} (1 - P(s|s))$$

$$= \frac{1}{1 - P(s|s)}$$

$$\mathbb{E}(S_{rain}) = \frac{1}{1 - .7} = 3.33$$



For example, what is the expected number of raining days?



$$X_1 = \begin{pmatrix} 1 & sun \\ 0 & rain \end{pmatrix}$$

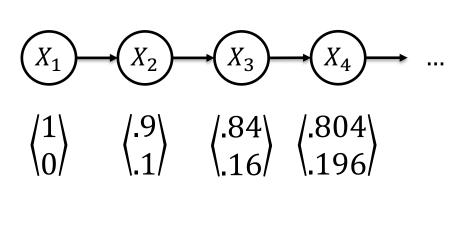
# Mini-Forward Algorithm

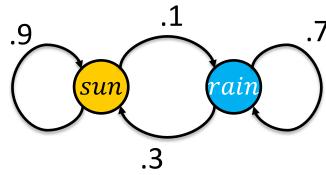
- What is P(X) on some day t?
  - We can obtain an answer by simulating the chain

$$P(x_1)$$
 is known

$$P(x_t) = \sum_{x_{t-1}} P(x_t, x_{t-1})$$

$$= \sum_{x_{t-1}} P(x_t | x_{t-1}) P(x_{t-1})$$



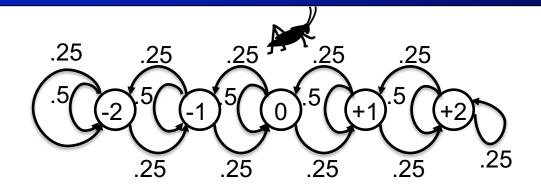


# Mini-Forward Algorithm

```
Input: time n, transition probability P(X_t|X_{t-1}), prior probability of states P(X_1) Output: P(X_t) for each state x do p[x,1] \leftarrow P(X_1 = x) for t \leftarrow 2 to n do p[x_t,t] = 0 for each state x_t do p[x_t,t] \leftarrow p[x_t,t] + p[x_{t-1},t-1]P(x_t|x_{t-1}) return p[x,n]
```

$$O(n|X|^2)$$

## Grasshopper Example



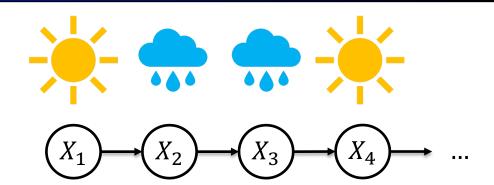
	<b>-</b> 2	-1	0	1	2
$P(X_1)$	0	0	1	0	0
$P(X_2)$	0	.25	.5	.25	0
$P(X_3)$	$.25^2 = .0625$	2(.5)(25) = .25	$.5^2 + 2(.25)^2 = .375$	2(.5)(25) = .25	$.25^2 = .0625$

$$P(X_t) = P(X_{t-1})T \qquad \begin{bmatrix} 0 & 0 & 1 & 0 & 0 \end{bmatrix} \begin{bmatrix} .75 & .25 & & & \\ .25 & .5 & .25 & & \\ & .25 & .5 & .25 & \\ & .25 & .75 \end{bmatrix} = \begin{bmatrix} 0 & .25 & .5 & .25 & \\ .25 & .75 & .25 & \\ & .25 & .75 \end{bmatrix}$$

# **Stationary Distributions**

Starting with a sunny day

$$\begin{pmatrix} 1 \\ 0 \end{pmatrix}$$
  $\begin{pmatrix} .9 \\ .1 \end{pmatrix}$   $\begin{pmatrix} .84 \\ .16 \end{pmatrix}$   $\begin{pmatrix} .804 \\ .196 \end{pmatrix}$  ...  $\begin{pmatrix} .75 \\ .25 \end{pmatrix}$ 

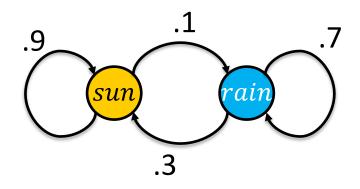


Starting with a rainy day

$$\begin{pmatrix} 0 \\ 1 \end{pmatrix}$$
  $\begin{pmatrix} .3 \\ .7 \end{pmatrix}$   $\begin{pmatrix} .48 \\ .52 \end{pmatrix}$   $\begin{pmatrix} .588 \\ .412 \end{pmatrix}$  ...  $\begin{pmatrix} .75 \\ .25 \end{pmatrix}$ 

Starting with an unknown day

$$\binom{p}{1-p}$$
 ...  $\binom{.75}{.25}$ 



# Stationary Distributions

#### For most chains

- Influence of the initial distribution gets less and less over time
- The distribution we end up in is independent of the initial distribution

## Stationary distribution

- The *stationary distribution*  $\pi$  of the chain is the distribution we obtain if the chain converges
- The stationary distribution satisfies

$$P_{\infty}(X) = P_{\infty+1}(X) = \sum_{x} P(X|x)P_{\infty}(x)$$
$$\pi(X) = \sum_{x} P(X|x)\pi(x)$$

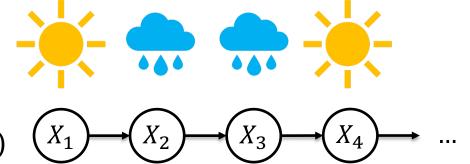
$$\pi(X) = \sum_{x} P(X|x)\pi(x)$$

$$\pi = \pi T$$

# **Stationary Distributions**

## • Question: What is $P(X_{\infty})$ ?

$$\pi(sun) = P(sun|sun)\pi(sun) + P(sun|rain)\pi(rain)$$
  
$$\pi(rain) = P(rain|sun)\pi(sun) + P(rain|rain)\pi(rain)$$



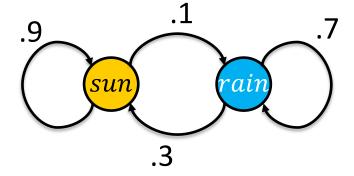
$$\pi(sun) = 0.9 \pi(sun) + 0.3 \pi(rain)$$
  
 $\pi(rain) = 0.1 \pi(sun) + 0.7 \pi(rain)$ 

$$\pi(sun) = 3\pi(rain)$$
  
 $\pi(rain) = 1/3\pi(sun)$ 

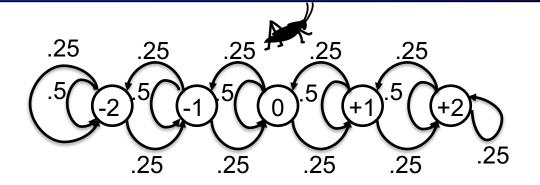
$$\pi(sun) + \pi(rain) = 1$$

$$\pi(sun) = 3/4$$

$$\pi(rain) = 1/4$$



# Stationary Distributions: Grasshopper

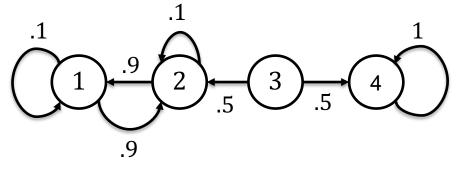


What is the stationary distribution?

$$T = \begin{bmatrix} .75 & .25 \\ .25 & .5 & .25 \\ & .25 & .5 & .25 \\ & & .25 & .5 & .25 \\ & & .25 & .75 \end{bmatrix}$$

## Irreducible Markov Chains

- A Markov chain is *irreducible* if every state x' is reachable from every other state x
  - That is, for every pair of states x and x', there is some time t such that the  $P(X_t = x' | X_1 = x) > 0$
  - Also known as regular or ergodic chain
- In this case, the states of the Markov chain are said to be recurrent
  - Each state is guaranteed to be visited an infinite number of times when we simulate the chain



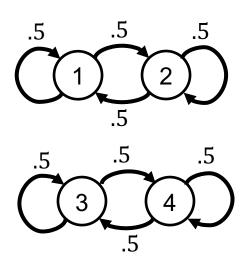
A reducible Markov chain

# **Stationary Distribution**

- Every (finite state) Markov chain has at least one stationary distribution
  - Yet an irreducible Markov chain is guaranteed to have a unique stationary distribution

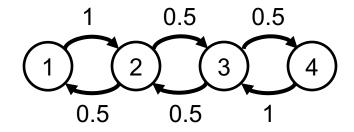


 To guarantee convergence, we need an additional property: Aperiodicity



## **Aperiodic Markov Chains**

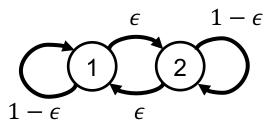
- A Markov chain is aperiodic if it is possible to return to any state at any time
  - There exists an t such that for all state x and all  $t' \ge t$ ,  $P(X_{t'} = x \mid X_1 = x) > 0$
- An irreducible and aperiodic Markov chain converges to a unique stationary distribution
  - Irreducible: we can go from any state to any state
  - Aperiodic: avoids chains that alternates forever between states without ever settling in a stationary distribution



An irreducible but periodic Markov chain

# Markov chains Convergence

- Although an irreducible and aperiodic Markov chain converges to a single stationary distribution, the convergence can be slow
  - In this example, the stationary distribution is close to (0.5, 0.5)
  - For a small  $\epsilon$  it will take a very long time to reach the stationary distribution
  - We stay in the same state with high probability and rarely transition to another state
  - The average of these states will converge to (0.5, 0.5), but the convergence will be very slow



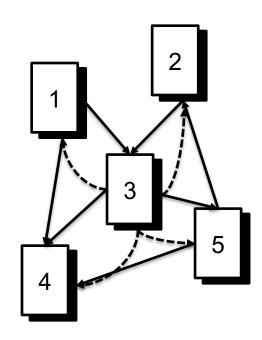
# **Applications of Markov Chains**

#### Markov chains have several well-know applications

- Markov chain Monte Carlo (MCMC) is a powerful approximate inference algorithm used in statistical software such as Stan
- MCs are part of the (LZMA) Lempel-Ziv-Markov compression algorithm used in 7zip
- PageRank algorithm used by Google 1.0 is a direct application of MCs

## PageRank

- Model the web as a state graph: pages are states and hyperlinks are transitions
- Each transition from state i has a probability  $\frac{\alpha}{k_i}$ , where  $\alpha$  is a constant parameter and  $k_i$  is the outgoing degree of node i
- Compute a stationary distribution. But it is not unique. Why?
- Augment the graph with phantom transitions of weight  $\frac{1-\alpha}{N}$ , where N is the number of nodes in the graph



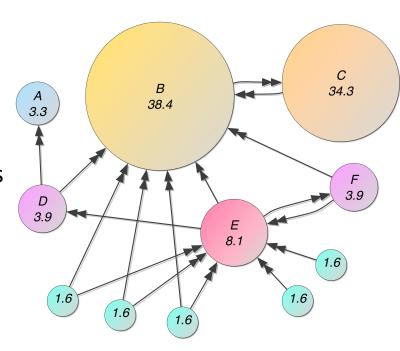
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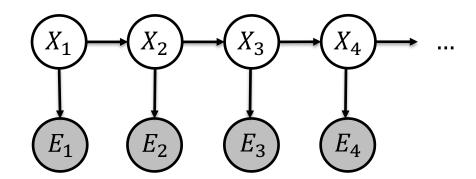
#### PageRank

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# Hidden Markov Models (HMM)

- Hidden Markov Models (HMM) are Markov chains where the states are not directly observable
  - In the weather example, the weather may not be directly observable
  - Instead, we use sensors, such as temperature, air pressure, humidity, wind speed, etc.

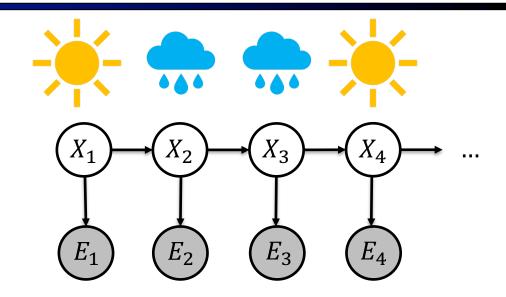


- HMM has two components
  - Underlying Markov chain over states X
  - Observable outputs (effects of the states) at each time step
  - These outputs are often called *emissions*

# HMM Weather Example

## HMM parameters

- Initial distribution  $P(X_1)$
- Transition probabilities  $P(X_t|X_{t-1})$
- Emission probabilities  $P(E_t|X_t)$



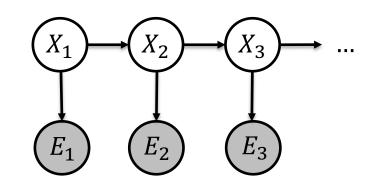
$X_1$	$P(X_1)$	_	$X_{t-1}$	$X_t$	$P(X_t X_{t-1})$
sun	.5		sun	sun	.7
rain	.5		sun	rain	.3
			rain	sun	.3
			rain	rain	.7

$X_t$	$E_{t}$	$P(E_t X_t)$
sun	umb	.2
sun	$\overline{umb}$	.8
rain	umb	.9
rain	$\overline{umb}$	.1

# HMM: Independencies

The chain rule of Bayesian networks for HMMs

$$P(X_1, E_1, \dots, X_n, E_n) = P(X_1)P(E_1|X_1) \prod_{t=1}^n P(X_t|X_{t-1})P(E_t|X_t)$$



- Independences are also apparent when we look at the chain rule
  - Chain rule for Bayesian networks for this example

$$P(X_1, E_1, X_2, E_2, X_3, E_3) = P(X_1)P(E_1|X_1)P(X_2|X_1)P(E_2|X_2)P(X_3|X_2)P(E_3|X_3)$$

Chain rule in general

$$P(X_1, E_1, X_2, E_2, X_3, E_3) = P(X_1)P(E_1|X_1)P(X_2|X_1, E_1)P(E_2|X_2, X_1, E_1)P(X_3|X_2, X_1, E_2, E_1)P(E_3|X_3, X_2, X_1, E_2, E_1)$$

$$X_2 \perp E_1 | X_1$$
  $X_3 \perp X_1, E_1, E_2 | X_2$   $E_2 \perp X_1, E_1 | X_2$   $E_3 \perp X_1, X_2, E_1, E_2 | X_3$ 

# HMM: Independencies

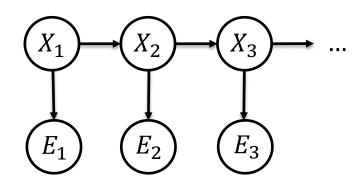
- In general, HMM have the following independency assumptions
  - A state is independent of all past states and all past evidence given the previous state (Markov property)

$$X_t \perp X_1, \dots, X_{t-2}, E_1, \dots, E_{t-2} | X_{t-1} |$$

 Evidence is independent of all past evidence and all past states given the current state (independence of observations)

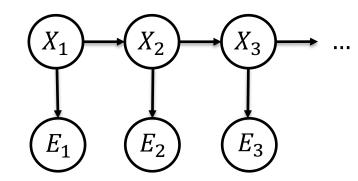
$$E_t \perp X_1, \dots, X_{t-1}, E_1, \dots, E_{t-1} | X_t$$

 Transition and emission probabilities are the same for all values of t (stationary process)



## HMM: Inference

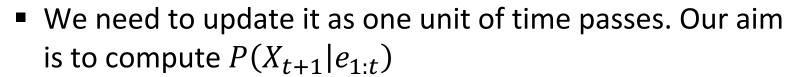
- We start with a first task of tracking the distribution  $P(X_t)$  over time
  - This task is known as filtering or monitoring
  - We use  $B(X_t) = P(X_t|e_1, ..., e_t)$  to denote the *belief of state*
  - We start with  $B(X_1)$ , usually using a uniform distribution
  - Update  $B(X_t)$  as time passes and we get new observations
- The inference has two main steps
  - Passage of time
  - Observation



## Passage of Time

• Suppose we know the current state of belief  $B(X_t)$ 

$$B(X_t) = P(X_t|e_{1:t})$$

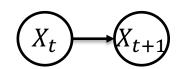


$$P(X_{t+1}|e_{1:t}) = \sum_{x_t} P(X_{t+1}, x_t|e_{1:t})$$

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

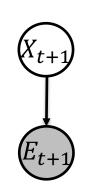
$$= \sum_{x_t} P(X_{t+1}|x_t) P(x_t|e_{1:t})$$

$$= \sum_{x_t} P(X_{t+1}|x_t) B(x_t)$$



## Observation

- Given we updated the belief with passage of time
  - We know  $P(X_{t+1}|e_{1:t})$  and we need to update it to  $B(X_{t+1}) = P(X_{t+1}|e_{1:t+1})$



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

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

$$\propto P(X_{t+1}, e_{t+1}|e_{1:t})$$

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

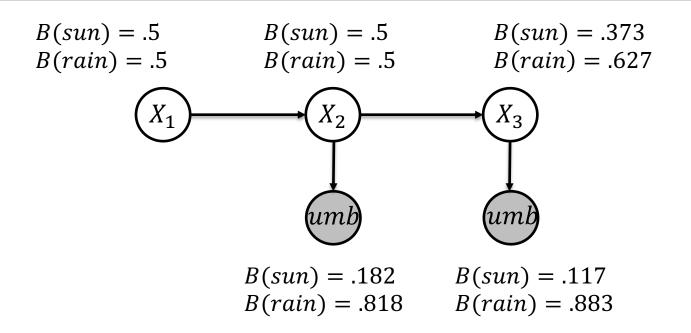
$$= P(e_{t+1}|X_{t+1}, e_{1:t}) P(X_{t+1}|e_{1:t})$$

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

$$B(X_{t+1}) \propto P(e_{t+1}|X_{t+1}) P(X_{t+1}|e_{1:t})$$

We must renormalise the results by  $\sum B(X_{t+1})$ 

# HMM Weather Example



$X_1$	$P(X_1)$
sun	.5
rain	.5

$X_{t-1}$	$X_t$	$P(X_t X_{t-1})$
sun	sun	.7
sun	rain	.3
rain	sun	.3
rain	rain	.7

$X_t$	$E_t$	$P(E_t X_t)$
sun	umb	.2
sun	$\overline{umb}$	.8
rain	umb	.9
rain	$\overline{umb}$	.1

## Forward Algorithm

 Suppose we have a sequence of evidence observations and we want to know the state belief at the end of the sequence

$$B(X_t) = P(X_t | e_{1:t})$$

$$\begin{split} P(X_{t}|e_{1:t}) &\propto P(X_{t},e_{1:t}) \\ &= \sum_{x_{t-1}} P(X_{t},x_{t-1},e_{1:t}) \\ &= \sum_{x_{t-1}} P(X_{t},x_{t-1},e_{t},e_{1:t-1}) \\ &= \sum_{x_{t-1}} P(x_{t-1}) P(X_{t}|x_{t-1}) P(e_{t}|x_{t-1},X_{t}) P(e_{1:t-1}|e_{t},x_{t-1},X_{t}) \\ &= \sum_{x_{t-1}} P(x_{t-1}) P(X_{t}|X_{t-1}) P(e_{t}|X_{t}) P(e_{1:t-1}|x_{t-1}) \\ &= \sum_{x_{t-1}} P(X_{t}|x_{t-1}) P(e_{t}|X_{t}) P(e_{1:t-1},x_{t-1}) \\ &= P(e_{t}|X_{t}) \sum_{x_{t-1}} P(X_{t}|x_{t-1}) P(x_{t-1},e_{1:t-1}) \end{split}$$

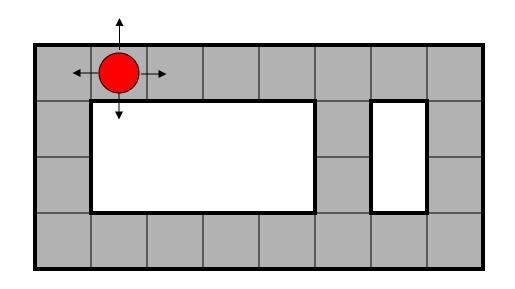
You can renormalise every step, but this algorithm often renormalised only the final one

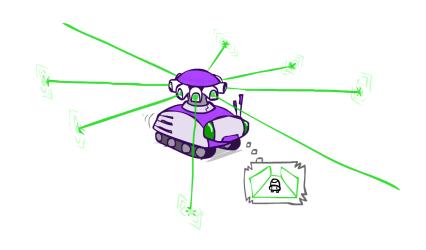
# Forward Algorithm

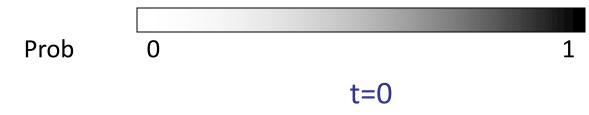
```
Input: time n, transition probability T, emission probability E, prior probability of
states P(X_1), sequence of observations \{e_2, \dots, e_t\}
Output: B(X_t)
for each state x do
    p[x,1] \leftarrow P(X_1 = x)
for t \leftarrow 2 to n do
     for each state x_t do
          p[x_t, t] = 0
           for each state x_{t-1} do
                 p[x_t, t] \leftarrow p[x_t, t] + p[x_{t-1}, t-1]T(x_t|x_{t-1})
           p[x_t, t] \leftarrow p[x_t, t] E(e_t | x_t)
return p[x, n] for all states x
                                     O(n|X|^2)
```

## **Example: Robot Localization**

Example from Michael Pfeiffer





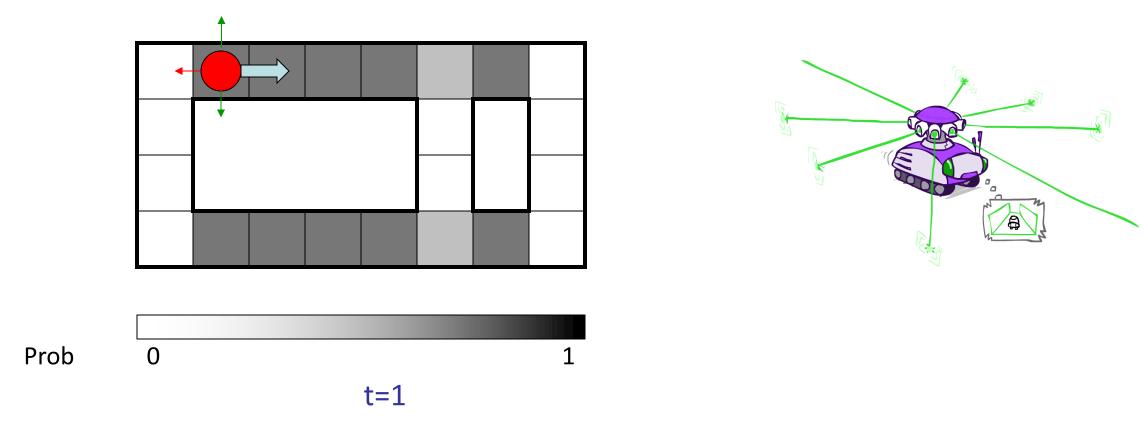


Sensor model: can read in which directions there is a wall, never more than 1 mistake

Motion model: may not execute action with small prob.

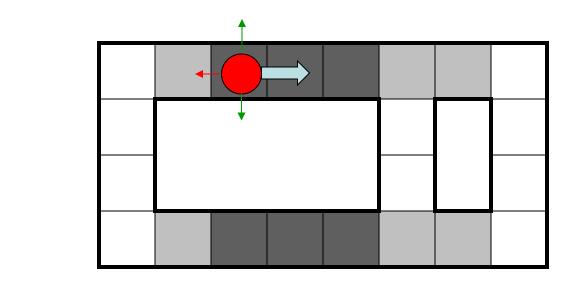
Slide from Berkeley Al course

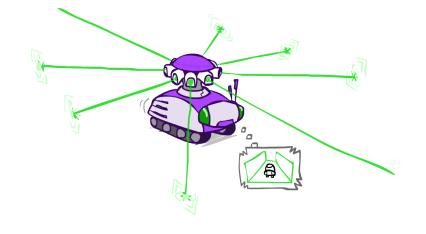
## **Example: Robot Localization**



Lighter grey: was possible to get the reading, but less likely b/c required 1 mistake

# **Example: Robot Localization**



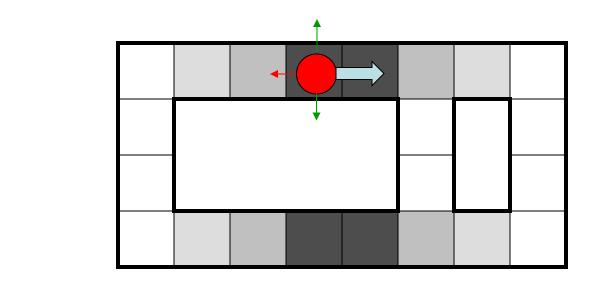


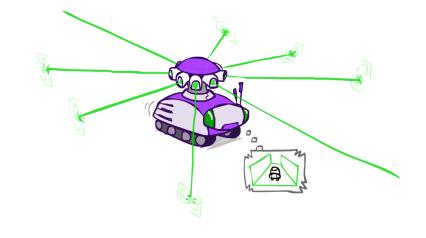
Prob 0 1

t=2

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## **Example: Robot Localization**



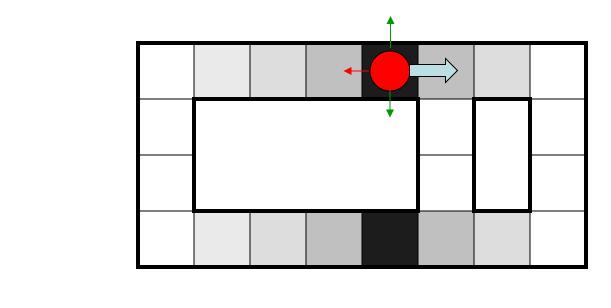


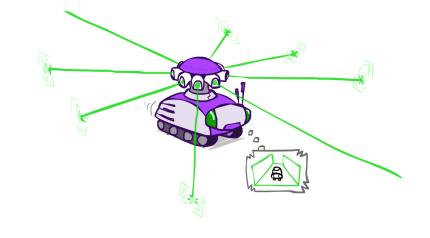
Prob 0 1

t=3

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## **Example: Robot Localization**



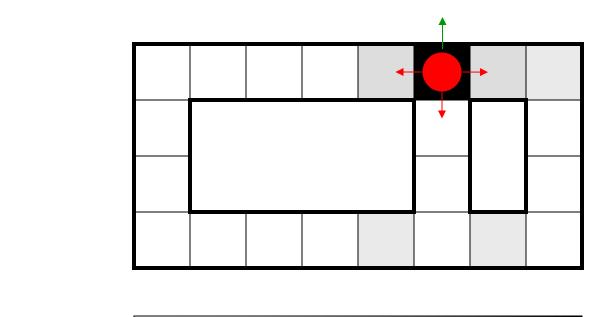




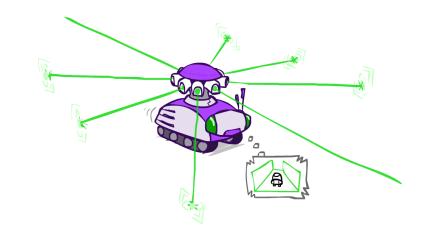


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## **Example: Robot Localization**



Prob

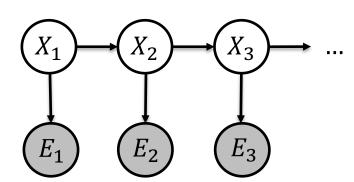


t=5

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## Most Probable Explanation (MPE)

- The forward algorithm tracks the probability of the states
  - These probabilities are updates with as time passes and we observe evidence
- A different task is to provide the most likely explanation
  - Considering all possible state combinations, which one has the highest probability considering the evidence
  - Therefore, we want to compute  $argmax_{x_{1:t}}P(x_{1:t}|e_{1:t})$

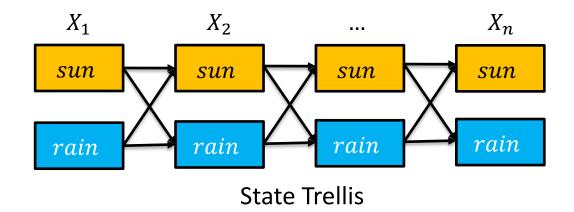


#### State Trellis

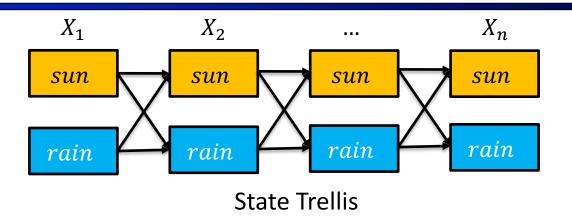
- A state trellis is a graph that illustrates the state transition over time
  - Each arc represents a time passage/evidence observation with weight

$$P(x_t|x_{t-1})P(e_t|x_t)$$

- A path is a sequence of states
  - The product of weights on a path is the sequence probability according to the evidence
  - The forward algorithm computes sums of paths probabilities that end in a same state, such as  $X_n = sun$
  - We will see now the Viterbi algorithm that computes the path with highest probability



#### Forward and Viterbi Algorithms



The forward algorithm computes the sum of the path probabilities that lead to the same final state

$$s[x_t] = P(x_t|e_{1:t})$$

$$= P(e_t|x_t) \sum_{x_{t-1}} P(x_t|x_{t-1})s[x_{t-1}]$$

 The Viterbi algorithm computes the maximum of the path probabilities that lead to the same final state

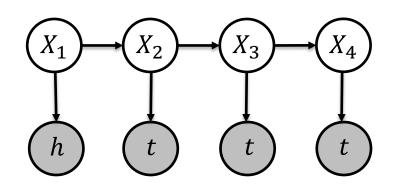
$$m[x_t] = \max_{x_{1:t-1}} P(x_{1:t-1}, x_t | e_{1:t})$$

$$= P(e_t | X_t) \max_{x_{t-1}} P(x_t | x_{t-1}) m[x_{t-1}]$$

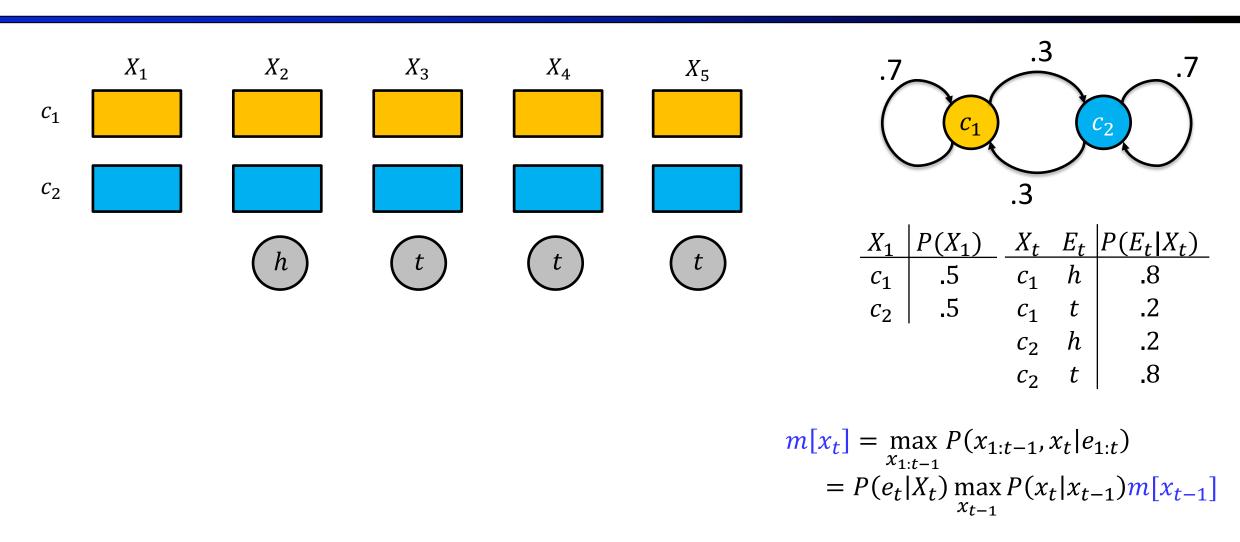
#### • Consider we have two unfair coins, $c_1$ and $c_2$

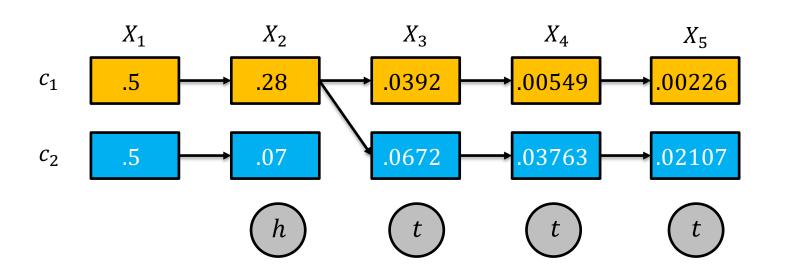
- Someone flips the coins sequentially, but we do not know which one. We only observe the outcomes heads or tails
- But we know that  $c_1$  has a higher probability of *heads* and  $c_2$  of *tails*
- Also, the person has a preference to keep the same coin and he/she starts with a coin chosen randomly with equal probabilities

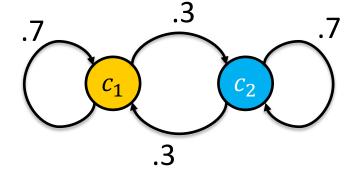
$P(X_t X_t)$	$X_t$	$X_{t-1}$	$P(X_1)$	$X_1$
.7	$c_1$		.5	$c_1$
.3	$c_2$	$c_1$	.5	$c_2$
.3	$c_1$	$c_2$		
_	_	_		



$X_t$	$E_{t}$	$P(E_t X_t)$
$c_1$	h	.8
$c_1$	t	.2
$c_2$	h	.2
$c_2$	t	.8







$$egin{array}{c|ccccc} X_1 & P(X_1) & X_t & E_t & P(E_t|X_t) \\ \hline c_1 & .5 & c_1 & h & .8 \\ c_2 & .5 & c_1 & t & .2 \\ & & c_2 & h & .2 \\ & & c_2 & t & .8 \\ \hline \end{array}$$

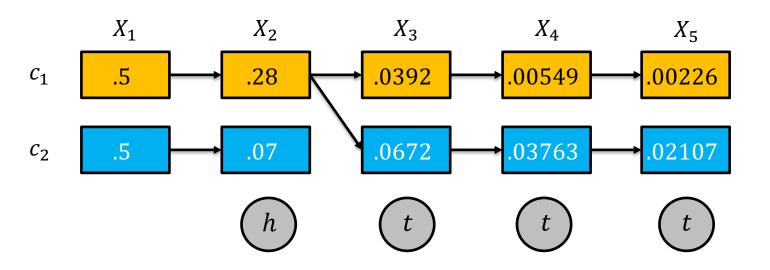
$$m[x_t] = \max_{x_{1:t-1}} P(x_{1:t-1}, x_t | e_{1:t})$$

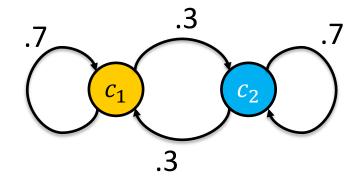
$$= P(e_t | X_t) \max_{x_{t-1}} P(x_t | x_{t-1}) m[x_{t-1}]$$

```
Input: time n, transition probability T, emission probability E, prior probability of
states P(X_1), sequence of observations \{e_2, \dots, e_t\}
Output: max P(x_{1:t-1}, x_t | e_{2:t})
for each state x do
     m[x,1] \leftarrow P(X_1 = x)
for t \leftarrow 2 to n do
     for each state x_t do
           m[x_t, t] = 0
           for each state x_{t-1} do
                 if m[x_{t-1}, t-1]T(x_t|x_{t-1}) > m[x_t, t]
                       m[x_t, t] \leftarrow m[x_{t-1}, t-1]T(x_t|x_{t-1})
           m[x_t, t] \leftarrow m[x_t, t] E(e_t | x_t)
return p[x, n] for all states x
```

 $O(n|X|^2)$ 

- The Viterbi algorithm of the previous slide provides the probability of the most likely sequence
  - However, often we are more interested in the sequence instead of its probability
- There are to common solutions
  - Keep an additional structure pointing to the parent of each node
  - Backtrack the computation from the last node



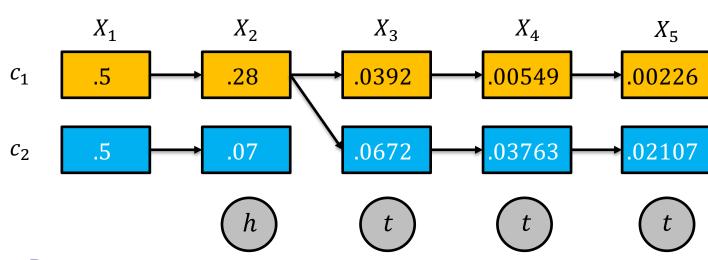


$X_1$	$P(X_1)$	$X_t$	$E_{t}$	$P(E_t X_t)$
$c_1$	.5	$c_1$	h	.8
$c_2$	.5	$c_1$	t	.2
	-	$c_2$	h	.2
		$C_2$	t	.8

$$m[x_t] = \max_{x_{1:t-1}} P(x_{1:t-1}, x_t | e_{1:t})$$

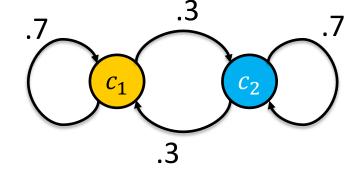
$$= P(e_t | X_t) \max_{x_{t-1}} P(x_t | x_{t-1}) m[x_{t-1}]$$

# Viterbi Algorithm: Backtracking Computation



#### Repeat

- 1. Divide by the probability of evidence
- 2. For each state  $x_{t-1}$  divide by  $P(x_t|x_{t-1})$
- 3. See which value of  $x_{t-1}$  matches the result

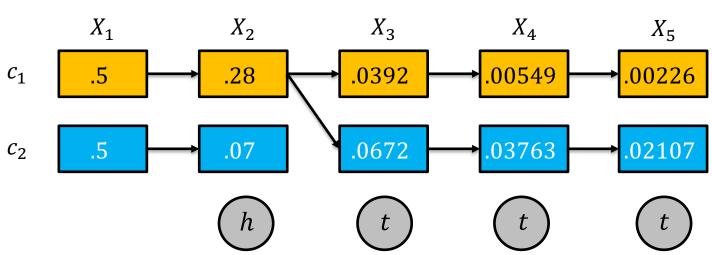


$X_1$	$P(X_1)$	$X_t$	$E_t$	$P(E_t X_t)$
$c_1$	.5	$c_1$	h	.8
$c_2$	.5	$c_1$	t	.2
		$c_2$	h	.2
		$c_2$	t	.8

$$m[x_t] = \max_{x_{1:t-1}} P(x_{1:t-1}, x_t | e_{1:t})$$

$$= P(e_t | X_t) \max_{x_{t-1}} P(x_t | x_{t-1}) m[x_{t-1}]$$

# Viterbi Algorithm: Backtracking Computation



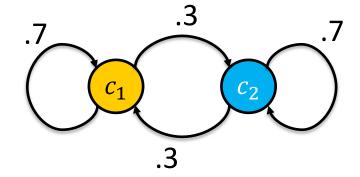
#### Repeat

- 1. Divide by the probability of evidence
- 2. For each state  $x_{t-1}$  divide by  $P(x_t|x_{t-1})$
- 3. See which value of  $x_{t-1}$  matches the result

$$\frac{.02107}{.8} = 0.0263375$$

$$x_4 = c_1 : \frac{0.0263375}{.3} \approx 0.08779$$

$$x_4 = c_2 : \frac{0.0263375}{7} \approx 0.03763$$



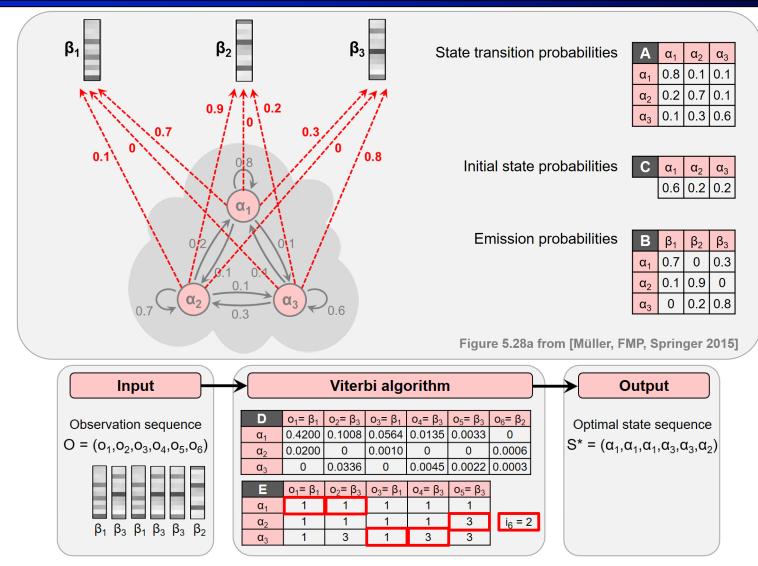
$X_1$	$P(X_1)$	$X_t$	$E_{t}$	$P(E_t X_t)$
$c_1$	<b>.</b> 5	$c_1$	h	.8
$c_2$	.5	$c_1$	t	.2
	•	$c_2$	h	.2
		$c_2$	t	.8

$$m[x_t] = \max_{x_{1:t-1}} P(x_{1:t-1}, x_t | e_{1:t})$$

$$= P(e_t | X_t) \max_{x_{t-1}} P(x_t | x_{t-1}) m[x_{t-1}]$$

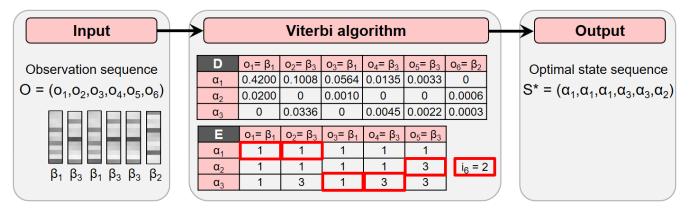
$$x_4 = c_2$$
: 0.03763

# Viterbi Algorithm: Backtracking Computation



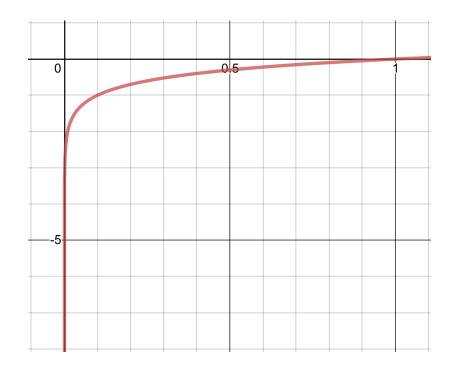
## Viterbi Algorithm: Vanishing Probabilities

- Notice the probabilities decrease as we observe more evidence
  - It is intuitive since the number of paths grows exponentially with the sequence size
  - In this example, the probabilities are around  $10^{-4}$  with just 6 steps
  - Long sequences (such as 100 steps) will cause an underflow and the probabilities will become zero
  - We can fix that using log probabilities, similarly to the Naïve Bayes classifier
  - This approach also replaces multiplications by sums that are more efficiently handled by most computers



#### Viterbi Algorithm: Log Probabilities

```
Input: time n, transition probability T, emission probability E, prior
probability of states P(X_1), sequence of observations \{e_2, \dots, e_t\}
Output: max \log P(x_{1:t-1}, x_t | e_{1:t})
for each state x do
     m[x,1] \leftarrow \log P(X_1 = x)
for t \leftarrow 2 to n do
     for each state x_t do
           m[x_t, t] = -\infty
           for each state x_{t-1} do
                 if m[x_{t-1}, t-1] + \log T(x_t|x_{t-1}) > m[x_t, t]
                       m[x_t, t] \leftarrow m[x_{t-1}, t-1] + \log T(x_t | x_{t-1})
            m[x_t, t] \leftarrow m[x_t, t] + \log E(e_t|x_t)
return p[x, n] for all states x
```



$$m[x_t] = \log \max_{x_{1:t-1}} P(x_{1:t-1}, x_t | e_{1:t-1})$$

$$= \log P(e_t | X_t) + \max_{x_{t-1}} \log P(x_t | x_{t-1}) + m[x_{t-1}]$$

## Particle Filtering

- Filtering: approximate solution
- Sometimes |X| is too big to use exact inference
  - |X| may be too big to even store B(X)
  - E.g. *X* is continuous
- Solution: approximate inference
  - Track samples of X, not all values
  - Samples are called particles
  - Time per step is linear in the number of samples
  - But: number needed may be large
  - In memory: list of particles, not states
- This is how robot localization works in practice
- Particle is just new name for sample

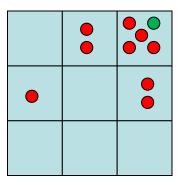
0.0	0.1	0.0
0.0	0.0	0.2
0.0	0.2	0.5



	•
• •	

#### Representation: Particles

- Our representation of P(X) is now a list of N particles (samples)
  - Generally,  $N \ll |X|$
  - Storing map from X to counts would defeat the point



- P(x) approximated by number of particles with value x
  - So, many x may have P(x) = 0!
  - More particles, more accuracy
- For now, all particles have a weight of 1

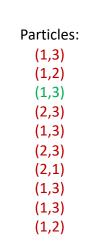
Particles
(1,3)
(1,2)
(1,3)
(2,3)
(1,3)
(2,3)
(2,1)
(1,3)
(1,3)
(1,2)

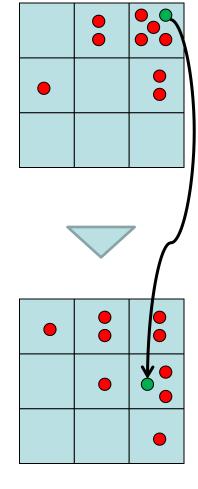
#### Particle Filtering: Elapse Time

 Each particle is moved by sampling its next position from the transition model

$$x' = \text{sample}(P(X'|x))$$

- Here, most samples move clockwise, but some move in another direction or stay in place
- This captures the passage of time
  - If enough samples, close to exact values before and after (consistent)





Particles:

(2,3) (1,1) (1,2)

(1,2)

(2,2)

#### Particle Filtering: Observe

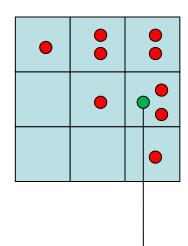
- Slightly trickier:
  - Don't sample observation, fix it
  - Downweigh samples based on the evidence

$$w(x) = P(e|x)$$

$$B(X) \propto P(e|X)B'(X)$$

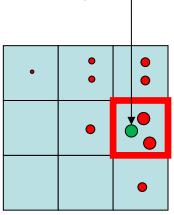
 As before, the probabilities don't sum to one, since all have been downweighed (in fact they now sum to (N times) an approximation of P(e))

Particles:	
(2,3)	
(1,2)	
(2,3)	
(3,3)	
(1,3)	
(2,3)	
(1,1)	
(1,2)	
(1,3)	
(2,2)	



#### Particles: (2,3) w=.9 (1,2) w=.2 (2,3) w=.9 (3,3) w=.4 (1,3) w=.4

- (2,3) w=.9 (1,1) w=.1
- (1,3) w=.4
- (1,2) w=.2 (2,2) w=.4

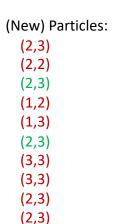


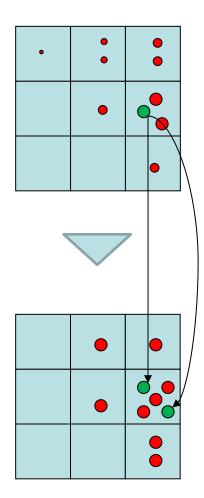
## Particle Filtering: Resample

- Rather than tracking weighted samples, we resample
- N times, we choose from our weighted sample distribution (i.e. draw with replacement)
- This is equivalent to renormalizing the distribution
- Now the update is complete for this time step, continue with the next one

# Particles: (2,3) w=.9 (1,2) w=.2 (2,3) w=.9 (3,3) w=.4 (1,3) w=.4 (2,3) w=.9 (1,1) w=.1 (1,2) w=.2 (1,3) w=.4

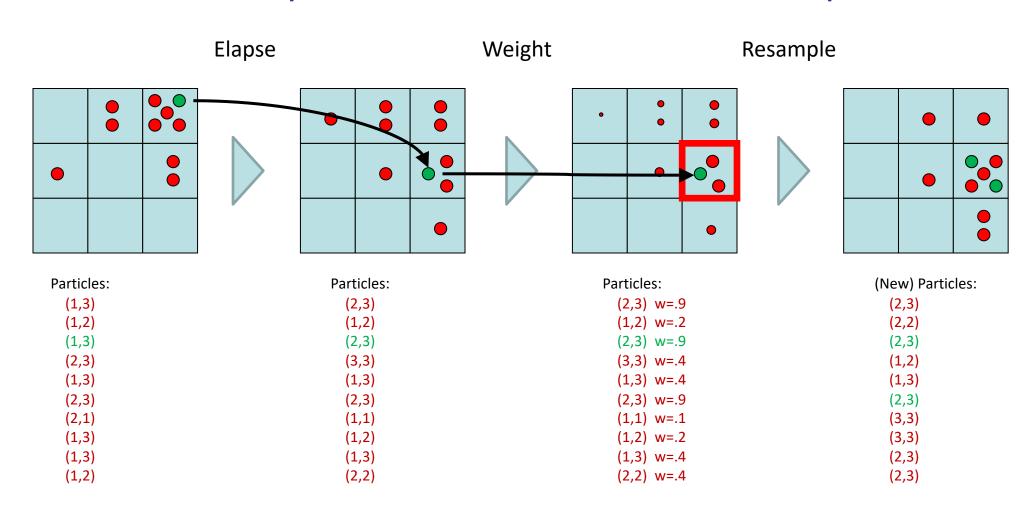
(2,2) w=.4





#### Recap: Particle Filtering

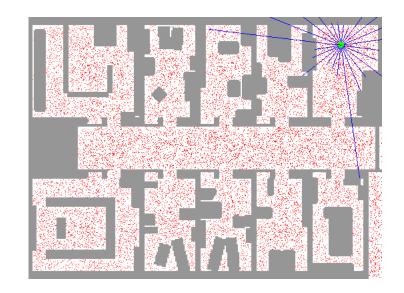
Particles: track samples of states rather than an explicit distribution



#### **Robot Localization**

#### In robot localization:

- We know the map, but not the robot's position
- Observations may be vectors of range finder readings
- State space and readings are typically continuous (works basically like a very fine grid) and so we cannot store B(X)
- Particle filtering is a main technique



#### Conclusion

- Markov chains and Hidden Markov models are simple examples of Dynamic Bayesian networks
  - DBNs are networks that allow us to model changes in time or space
  - Changes in time are specified using transition probabilities
- Markov chains are sequence models
  - It tracks the probability distribution over a series of transitions
  - For many sequences, the probability distribution converges to a stationary distribution
  - The stationary distribution has several applications such as the MCMC algorithms used for approximate inference
- Hidden Markov models are Markov chains with hidden states
  - Those states are never directly observed, but we can make indirect inference through emissions
  - These models are used in several applications such as language and signal processing and robot localisation