# CS 188: Artificial Intelligence 

## Review of Probability, Bayes' nets

DISCLAIMER: It is insufficient to simply study these slides, they are merely meant as a quick refresher of the high-level ideas covered. You need to study all materials covered in lecture, section, assignments and projects !

## Probability recap

- Conditional probability $\quad P(x \mid y)=\frac{P(x, y)}{P(y)}$
- Product rule $\quad P(x, y)=P(x \mid y) P(y)$
- Chain rule $P\left(X_{1}, X_{2}, \ldots X_{n}\right)=P\left(X_{1}\right) P\left(X_{2} \mid X_{1}\right) P\left(X_{3} \mid X_{1}, X_{2}\right) \ldots$
- X, Y independent iff: $\forall x, y: P(x, y)=P(x) P(y)$
equivalently, iff: $\quad \forall x, y: P(x \mid y)=P(x)$
equivalently, iff: $\quad \forall x, y: P(y \mid x)=P(y)$
- X and Y are conditionally independent given Z iff:
$\forall x, y, z: P(x, y \mid z)=P(x \mid z) P(y \mid z)$
equivalently, iff: $\forall x, y, z: P(x \mid y, z)=P(x \mid z)$
equivalently, iff: $\forall x, y, z: P(y \mid x, z)=P(y \mid z)$


## Inference by Enumeration

- P(sun)?
- P (sun | winter)?
- $\mathrm{P}($ sun | winter, hot)?

| S | T | W | P |
| :---: | :---: | :---: | :---: |
| summer | hot | sun | 0.30 |
| summer | hot | rain | 0.05 |
| summer | cold | sun | 0.10 |
| summer | cold | rain | 0.05 |
| winter | hot | sun | 0.10 |
| winter | hot | rain | 0.05 |
| winter | cold | sun | 0.15 |
| winter | cold | rain | 0.20 |

## Bayes' Nets Recap

- Representation
- Chain rule -> Bayes' net = DAG + CPTs
- Conditional Independences
- D-separation
- Probabilistic Inference
- Enumeration (exact, exponential complexity)
- Variable elimination (exact, worst-case exponential complexity, often better)
- Probabilistic inference is NP-complete
- Sampling (approximate)


## Chain Rule $\rightarrow$ Bayes net

- Chain rule: can always write any joint distribution as an incremental product of conditional distributions

$$
\begin{aligned}
& P\left(x_{1}, x_{2}, x_{3}\right)=P\left(x_{1}\right) P\left(x_{2} \mid x_{1}\right) P\left(x_{3} \mid x_{1}, x_{2}\right) \\
& P\left(x_{1}, x_{2}, \ldots x_{n}\right)=\prod_{i} P\left(x_{i} \mid x_{1} \ldots x_{i-1}\right)
\end{aligned}
$$

- Bayes nets: make conditional independence assumptions of the form:

$$
P\left(x_{i} \mid x_{1} \cdots x_{i-1}\right)=P\left(x_{i} \mid \operatorname{parents}\left(X_{i}\right)\right)
$$

giving us:


$$
P\left(x_{1}, x_{2}, \ldots x_{n}\right)=\prod_{i=1}^{n} P\left(x_{i} \mid \text { parents }\left(X_{i}\right)\right)
$$

## Probabilities in BNs

- Bayes' nets implicitly encode joint distributions
- As a product of local conditional distributions
- To see what probability a BN gives to a full assignment, multiply all the relevant conditionals together:

$$
P\left(x_{1}, x_{2}, \ldots x_{n}\right)=\prod_{i=1}^{n} P\left(x_{i} \mid \operatorname{parents}\left(X_{i}\right)\right)
$$

- Example:
$P(+$ cavity, +catch, $\neg$ toothache $)$
- This lets us reconstruct any entry of the full joint
- Not every BN can represent every joint distribution
- The topology enforces certain conditional independencies


## Example: Alarm Network



## Size of a Bayes' Net for $P\left(X_{1}, X_{2}, \ldots X_{n}\right)$

- How big is a joint distribution over N Boolean variables? $2^{N}$
- Size of representation if we use the chain rule $2^{N}$
- How big is an N -node net if nodes have up to k parents? $\mathrm{O}\left(\mathrm{N}^{*} 2^{\mathrm{k}+1}\right)$
- Both give you the power to calculate
- BNs:
- Huge space savings!
- Easier to elicit local CPTs
- Faster to answer queries


## Bayes Nets: Assumptions

- Assumptions made by specifying the graph:

$$
P\left(x_{i} \mid x_{1} \cdots x_{i-1}\right)=P\left(x_{i} \mid \text { parents }\left(X_{i}\right)\right)
$$

- Given a Bayes net graph additional conditional independences can be read off directly from the graph
- Question: Are two nodes guaranteed to be independent given certain evidence?
- If no, can prove with a counter example
- I.e., pick a set of CPT's, and show that the independence assumption is violated by the resulting distribution
- If yes, can prove with
- Algebra (tedious)
- D-separation (analyzes graph)


## D-Separation

- Question: Are $X$ and $Y$ conditionally independent given evidence vars $\{Z\}$ ?
- Yes, if $X$ and $Y$ "separated" by $Z$
- Consider all (undirected) paths from $X$ to $Y$
- No active paths = independence!
- A path is active if each triple is active:
- Causal chain $\mathrm{A} \rightarrow \mathrm{B} \rightarrow \mathrm{C}$ where B is unobserved (either direction)
- Common cause $A \leftarrow B \rightarrow C$ where $B$ is unobserved
- Common effect (aka v-structure) $A \rightarrow B \leftarrow C$ where $B$ or one of its descendents is observed
- All it takes to block a path is a single inactive segment

Active Triples






Inactive Triples





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## D-Separation

- Given query $X_{i} \xrightarrow{R} X_{j} \mid\left\{X_{k_{1}}, \ldots, X_{k_{n}}\right\}$
- Shade all evidence nodes
- For all (undirected!) paths between and
- Check whether path is active
- If active return $X_{i} \not \nmid X_{j} \mid\left\{X_{k_{1}}, \ldots, X_{k_{n}}\right\}$
- (If reaching this point all paths have been checked and shown inactive)
- Return $X_{i} \Perp X_{j} \mid\left\{X_{k_{1}}, \ldots, X_{k_{n}}\right\}$



## All Conditional Independences

- Given a Bayes net structure, can run dseparation to build a complete list of conditional independences that are necessarily true of the form

$$
X_{i} \Perp X_{j} \mid\left\{X_{k_{1}}, \ldots, X_{k_{n}}\right\}
$$

- This list determines the set of probability distributions that can be represented by Bayes' nets with this graph structure


## Topology Limits Distributions

- Given some graph topology G, only certain joint distributions can be encoded
- The graph structure guarantees certain (conditional) independences
- (There might be more independence)
- Adding arcs increases the set of distributions, but has several costs


Full conditioning can encode any distribution

$\{X \Perp Z \mid Y$






## Inference by Enumeration

- Given unlimited time, inference in BNs is easy
- Recipe:
- State the marginal probabilities you need
- Figure out ALL the atomic probabilities you need
- Calculate and combine them
- Example:

$$
\begin{aligned}
& P(+b \mid+j,+m)= \\
& \quad \frac{P(+b,+j,+m)}{P(+j,+m)}
\end{aligned}
$$



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## Example: Enumeration

- In this simple method, we only need the BN to synthesize the joint entries

$$
\begin{aligned}
& P(+b,+j,+m)= \\
& P(+b) P(+e) P(+a \mid+b,+e) P(+j \mid+a) P(+m \mid+a)+ \\
& P(+b) P(+e) P(-a \mid+b,+e) P(+j \mid-a) P(+m \mid-a)+ \\
& P(+b) P(-e) P(+a \mid+b,-e) P(+j \mid+a) P(+m \mid+a)+ \\
& P(+b) P(-e) P(-a \mid+b,-e) P(+j \mid-a) P(+m \mid-a)
\end{aligned}
$$

## Variable Elimination

- Why is inference by enumeration so slow?
- You join up the whole joint distribution before you sum out the hidden variables
- You end up repeating a lot of work!
- Idea: interleave joining and marginalizing!
- Called "Variable Elimination"
- Still NP-hard, but usually much faster than inference by enumeration


## Variable Elimination Outline

- Track objects called factors
- Initial factors are local CPTs (one per node)

- Any known values are selected
- E.g. if we know $L=+\ell$, the initial factors are

| $P(R)$ | $P(T \mid R)$ | $P(+\ell \mid T)$ |
| :---: | :---: | :---: |
| $+r$ 0.1 <br> $-r$ 0.9 | $+r$ $+t$ 0.8 <br> $+r$ $-t$ 0.2 <br> $-r$ $+t$ 0.1 <br> $-r$ $-t$ 0.9 | $+t$ +1 0.3 <br> $-t$ +1 0.1 |

- VE: Alternately join factors and eliminate variables



## Example

$P(B \mid j, m) \propto P(B, j, m)$

$$
P(B) \quad P(E) \quad P(A \mid B, E) \quad P(j \mid A) \quad P(m \mid A)
$$

Choose A
$\left.\begin{array}{l}P(A \mid B, E) \\ P(j \mid A) \\ P(m \mid A)\end{array} \quad \boxed{\times} P(j, m, A \mid B, E) \quad \sum\right\rangle P(j, m \mid B, E)$

$$
P(B) \quad P(E) \quad P(j, m \mid B, E)
$$

## Example

| $P(B)$ | $P(E)$ | $P(j, m \mid B, E)$ |
| :--- | :--- | :--- |

Choose E

$$
\begin{gathered}
\begin{array}{ccc}
P(E) & \boxed{x} & P(j, m, E \mid B) \\
P(j, m \mid B, E) & \boxed{\sum} & P(j, m \mid B) \\
& P(B) & P(j, m \mid B) \\
\hline
\end{array}
\end{gathered}
$$

Finish with B
$P(B)$
$P(j, m \mid B)$$\stackrel{\times>}{ } \quad P(j, m, B) \xrightarrow{\text { Normalize }} \boldsymbol{P}$

## General Variable Elimination

- Query: $P\left(Q \mid E_{1}=e_{1}, \ldots E_{k}=e_{k}\right)$
- Start with initial factors:
- Local CPTs (but instantiated by evidence)
- While there are still hidden variables (not Q or evidence):
- Pick a hidden variable H
- Join all factors mentioning H
- Eliminate (sum out) H
- Join all remaining factors and normalize


## Another (bit more abstractly worked out) Variable Elimination Example

Query: $P\left(X_{3} \mid Y_{1}=y_{1}, Y_{2}=y_{2}, Y_{3}=y_{3}\right)$
Start by inserting evidence, which gives the following initial factors:
$p(Z) p\left(X_{1} \mid Z\right) p\left(X_{2} \mid Z\right) p\left(X_{3} \mid Z\right) p\left(y_{1} \mid X_{1}\right) p\left(y_{2} \mid X_{2}\right) p\left(y_{3} \mid X_{3}\right)$
Eliminate $X_{1}$, this introduces the factor $f_{1}\left(Z, y_{1}\right)=\sum_{x_{1}} p\left(x_{1} \mid Z\right) p\left(y_{1} \mid x_{1}\right)$, and:

$p(Z) f_{1}\left(Z, y_{1}\right) p\left(X_{2} \mid Z\right) p\left(X_{3} \mid Z\right) p\left(y_{2} \mid X_{2}\right) p\left(y_{3} \mid X_{3}\right)$
Eliminate $X_{2}$, this introduces the factor $f_{2}\left(Z, y_{2}\right)=\sum_{x_{2}} p\left(x_{2} \mid Z\right) p\left(y_{2} \mid x_{2}\right)$, and:
$p(Z) f_{1}\left(Z, y_{1}\right) f_{2}\left(Z, y_{2}\right) p\left(X_{3} \mid Z\right) p\left(y_{3} \mid X_{3}\right)$
Eliminate $Z$, this introduces the factor
$f_{3}\left(y_{1}, y_{2}, y_{3}, X_{3}\right)=\sum_{z} p(z) f_{1}\left(z, y_{1}\right) f_{2}\left(z, y_{2}\right) p\left(X_{3} \mid z\right) p\left(y_{3} \mid X_{3}\right)$, and:
$f_{3}\left(y_{1}, y_{2}, y_{3}, X_{3}\right)$
Normalizing over $X_{3}$ gives $P\left(X_{3} \mid y_{1}, y_{2}, y_{3}\right)$.

Computational complexity critically depends on the largest factor being generated in this process. Size of factor = number of entries in table. In example above (assuming binary) all factors generated are of size 2 --- as they all only have one variable ( $Z, Z$, and $X 3$ respectively).

## Variable Elimination Ordering

- For the query $\mathrm{P}\left(\mathrm{X}_{\mathrm{n}} \mid \mathrm{y}_{1}, \ldots, \mathrm{y}_{\mathrm{n}}\right)$ work through the following two different orderings as done in previous slide: $Z, X_{1}, \ldots, X_{n-1}$ and $X_{1}, \ldots, X_{n-1}$, Z. What is the size of the maximum factor generated for each of the orderings?

- Answer: $2^{n}$ versus 2 (assuming binary)
- In general: the ordering can greatly affect efficiency.


## Computational and Space Complexity of Variable Elimination

- The computational and space complexity of variable elimination is determined by the largest factor
- The elimination ordering can greatly affect the size of the largest factor.
- E.g., previous slide's example $2^{\text {n }}$ vs. 2
- Does there always exist an ordering that only results in small factors?
- No!


## Worst Case Complexity?

- Consider the 3-SAT clause:
$\left(x_{1} \vee x_{2} \vee \neg x_{3}\right) \wedge\left(\neg x_{1} \vee x_{3} \vee \neg x_{4}\right) \wedge\left(x_{2} \vee \neg x_{2} \vee x_{4}\right) \wedge\left(\neg x_{3} \vee \neg x_{4} \vee \neg x_{5}\right) \wedge\left(x_{2} \vee x_{5} \vee x_{7}\right) \wedge\left(x_{4} \vee x_{5} \vee x_{6}\right) \wedge\left(\neg x_{5} \vee x_{6} \vee \neg x_{7}\right) \wedge\left(\neg x_{5} \vee \neg x_{6} \vee x_{7}\right)$ which can be encoded by the following Bayes' net:

- If we can answer $P(z)$ equal to zero or not, we answered whether the 3-SAT problem has a solution.
- Subtlety: why the cascaded version of the AND rather than feeding all OR clauses into a single AND? Answer: a single AND would have an exponentially large CPT, whereas with representation above the Bayes' net has small CPTs only.
- Hence inference in Bayes' nets is NP-hard. No known efficient probabilistic inference in general. ${ }^{27}$


## Polytrees

- A polytree is a directed graph with no undirected cycles
- For poly-trees you can always find an ordering that is efficient
- Try it!!
- Cut-set conditioning for Bayes' net inference
- Choose set of variables such that if removed only a polytree remains
- Think about how the specifics would work out? ${ }^{\text {B }}$


## Approximate Inference: Sampling

- Basic idea:
- Draw N samples from a sampling distribution S
- Compute an approximate posterior probability
- Show this converges to the true probability P
- Why? Faster than computing the exact answer
- Prior sampling:
- Sample ALL variables in topological order as this can be done quickly
- Rejection sampling for query $P\left(Q \mid E_{1}=e_{1}, \ldots E_{k}=e_{k}\right)$
- = like prior sampling, but reject when a variable is sampled inconsistent with the query, in this case when a variable $\mathrm{E}_{\mathrm{i}}$ is sampled differently from $e_{i}$
- Likelihood weighting for query $P\left(Q \mid E_{1}=e_{1}, \ldots E_{k}=e_{k}\right)$
- = like prior sampling but variables $\mathrm{E}_{\mathrm{i}}$ are not sampled, when it's their turn, they get set to $\mathrm{e}_{\mathrm{i}}$, and the sample gets weighted by $P\left(e_{i} \mid\right.$ value of parents $\left(e_{i}\right)$ in current sample)
- Gibbs sampling: repeatedly samples each non-evidence variable



## Example

- We'll get a bunch of samples from the BN:

$$
\begin{aligned}
& +c,-s,+r,+w \\
& +c,+s,+r,+w \\
& -c,+s,+r,-w \\
& +c,-s,+r,+w \\
& -c,-s,-r,+w
\end{aligned}
$$



- If we want to know $\mathrm{P}(\mathrm{W})$
- We have counts <+w:4, -w:1>
- Normalize to get $\mathrm{P}(\mathrm{W})=<+w: 0.8,-w: 0.2>$
- This will get closer to the true distribution with more samples
- Can estimate anything else, too
- What about $\mathrm{P}(\mathrm{C} \mid+\mathrm{w})$ ? $\quad \mathrm{P}(\mathrm{C} \mid+\mathrm{r},+\mathrm{w})$ ? $P(\mathrm{C} \mid-\mathrm{r},-\mathrm{w})$ ?
- Fast: can use fewer samples if less time



## Likelihood Weighting

- Sampling distribution if $z$ sampled and e fixed evidence

$$
S_{W S}(\mathbf{z}, \mathbf{e})=\prod_{i=1}^{l} P\left(z_{i} \mid \operatorname{Parents}\left(Z_{i}\right)\right)
$$

- Now, samples have weights

$$
w(\mathbf{z}, \mathbf{e})=\prod_{i=1}^{m} P\left(e_{i} \mid \operatorname{Parents}\left(E_{i}\right)\right)
$$



- Together, weighted sampling distribution is consistent
$S_{\mathrm{WS}}(z, e) \cdot w(z, e)=\prod_{i=1}^{l} P\left(z_{i} \mid \operatorname{Parents}\left(z_{i}\right)\right) \prod_{i=1}^{m} P\left(e_{i} \mid \operatorname{Parents}\left(e_{i}\right)\right)$
$=P(\mathbf{z}, \mathbf{e})$


## Gibbs Sampling

- Idea: instead of sampling from scratch, create samples that are each like the last one.
- Procedure: resample one variable at a time, conditioned on all the rest, but keep evidence fixed.
- Properties: Now samples are not independent (in fact they' re nearly identical), but sample averages are still consistent estimators!
- What's the point: both upstream and downstream variables condition on evidence.


## Markov Models

- A Markov model is a chain-structured BN
- Each node is identically distributed (stationarity)
- Value of $X$ at a given time is called the state
- As a BN:

- The chain is just a (growing) BN
- We can always use generic BN reasoning on it if we truncate the chain at a fixed length
- Stationary distributions
- For most chains, the distribution we end up in is independent of the initial distribution
- Called the stationary distribution of the chain
- Example applications: Web link analysis (Page Rank) and Gibbs Sampling


## Hidden Markov Models

- Underlying Markov chain over states S
- You observe outputs (effects) at each time step

- Speech recognition HMMs:
- $\mathrm{X}_{\mathrm{i}}$ : specific positions in specific words; $\mathrm{E}_{\mathrm{i}}$ : acoustic signals
- Machine translation HMMs:
- $\mathrm{X}_{\mathrm{i}}$ : translation options; $\mathrm{E}_{\mathrm{i}}$ : Observations are words
- Robot tracking:
- $\mathrm{X}_{\mathrm{i}}$ : positions on a map; $\mathrm{E}_{\mathrm{i}}$ : range readings


## Online Belief Updates

- Every time step, we start with current $P(X \mid$ evidence $)$
- We update for time:

$$
\begin{equation*}
P\left(x_{t} \mid e_{1: t-1}\right)=\sum_{x_{t-1}} P\left(x_{t-1} \mid e_{1: t-1}\right) \cdot P\left(x_{t} \mid x_{t-1}\right) \tag{X}
\end{equation*}
$$

- We update for evidence:

$$
P\left(x_{t} \mid e_{1: t}\right) \propto_{X} P\left(x_{t} \mid e_{1: t-1}\right) \cdot P\left(e_{t} \mid x_{t}\right)
$$

- The forward algorithm does both at once (and doesn' t normalize)


## Recap: Particle Filtering

- Particles: track samples of states rather than an explicit distribution

Elapse
Weight
Resample


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


Particles:
articles
$(3,2)$
$(2,3)$
$(3,2)$
$(2,3)$
$(2,3)$
$(3,2)$
$(3,2)$
$(3,1)$
$(3,1)$
$(3,3)$
$(3,2)$
$(13)$
$(1,3)$
$(2)$
$(2,3)$
$(3,2)$
$(3,2)$
$(2,2)$
$(2,2)$


Particles $(3,2) w=.9$ $(2,3) w=.2$ $(3,2) w=.9$
$(3,1) w=.4$
3,1) $w=.4$
$(3,3) w=.4$
$(3,2) \quad w=.9$
$(1,3) \quad w=.1$
$(2,3) w=.2$
$(3,2) w=.9$
$(2,2) w=.4$


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

## Dynamic Bayes Nets (DBNs)

- We want to track multiple variables over time, using multiple sources of evidence
- Idea: Repeat a fixed Bayes net structure at each time
- Variables from time $t$ can condition on those from $t-1$

- Discrete valued dynamic Bayes nets are also HMMs


## Best Explanation Queries



- Query: most likely seq:

$$
\underset{x_{1: t}}{\arg \max } P\left(x_{1: t} \mid e_{1: t}\right)
$$

## Best Explanation Query Solution Method 1: Search

$\underset{x_{1: t}}{\arg \max } P\left(x_{1: t} \mid e_{1: t}\right)=\underset{x_{1: t}}{\arg \max } \prod_{i=1}^{t} P\left(x_{t} \mid x_{t-1}\right) P\left(e_{t} \mid x_{t}\right) \quad \begin{gathered}\text { sight abuse of fototion } \\ \text { assuming } P\left(|x| x \mid x_{0}\right)=P\left(x_{1}\right)\end{gathered}$

$$
=\underset{x_{1: t}}{\arg \max } \sum_{i=1}^{t} \log \left(P\left(x_{t} \mid x_{t-1}\right) P\left(e_{t} \mid x_{t}\right)\right)
$$

- States: $\left\{(),+x_{1},-\mathrm{x}_{1},+\mathrm{x}_{2},-\mathrm{x}_{2}, \ldots,+\mathrm{x}_{\mathrm{t}},-\mathrm{x}_{\mathrm{t}}\right\}$
- Start state: ()
- Actions: in state $\mathrm{x}_{\mathrm{k}}$, choose any assignment for state $\mathrm{x}_{\mathrm{k}+1}$
- Cost: $\log \left(P\left(x_{k+1} \mid x_{k}\right) P\left(e_{k+1} \mid x_{k+1}\right)\right)$
- Goal test: goal $\left(\mathrm{x}_{\mathrm{k}}\right)=$ true iff $\mathrm{k}==\mathrm{t}$
$\rightarrow$ Can run uniform cost graph search to find solution
$\rightarrow$ Uniform cost graph search will take $\mathrm{O}\left(\mathrm{t} \mathrm{d}^{2}\right)$. Think about this!

Best Explanation Query Solution Method 2: Viterbi Algorithm (= max-product version of forward algorithm)

$$
\begin{aligned}
x_{1: T}^{*} & =\underset{x_{1: T}}{\arg \max } P\left(x_{1: T} \mid e_{1: T}\right)=\underset{x_{1: T}}{\arg \max } P\left(x_{1: T}, e_{1: T}\right) \\
m_{t}\left[x_{t}\right] & =\max _{x_{1: t-1}} P\left(x_{1: t-1}, x_{t}, e_{1: t}\right) \\
& =\max _{x_{1: t-1}} P\left(x_{1: t-1}, e_{1: t-1}\right) P\left(x_{t} \mid x_{t-1}\right) P\left(e_{t} \mid x_{t}\right) \\
& =P\left(e_{t} \mid x_{t}\right) \max _{x_{t-1}} P\left(x_{t} \mid x_{t-1}\right) \max _{x_{1: t-2}} P\left(x_{1: t-1}, e_{1: t-1}\right) \\
& =P\left(e_{t} \mid x_{t}\right) \max _{x_{t-1}} P\left(x_{t} \mid x_{t-1}\right) m_{t-1}\left[x_{t-1}\right]
\end{aligned}
$$

Viterbi computational complexity: O(t d$\left.{ }^{2}\right)$
Compare to forward algorithm:
$P\left(x_{t}, e_{1: t}\right)=P\left(e_{t} \mid x_{t}\right) \sum_{x_{t-1}} P\left(x_{t} \mid x_{t-1}\right) P\left(x_{t-1}, e_{1: t-1}\right)$

## Parameter Estimation

- Estimating distribution of random variables like X or $\mathrm{X} \mid \mathrm{Y}$
- Empirically: use training data
- For each outcome x , look at the empirical rate of that value:

$$
P_{\mathrm{ML}}(x)=\frac{\operatorname{count}(x)}{\text { total samples }} \quad \begin{aligned}
& P_{\mathrm{ML}}(r)=1 / 3
\end{aligned}
$$

- This is the estimate that maximizes the likelihood of the data
- Laplace smoothing

$$
L(x, \theta)=\prod_{i} P_{\theta}\left(x_{i}\right)
$$

- Pretend saw every outcome k extra times $P_{L A P, k}(x)=\frac{c(x)+k}{N+k|X|}$
- Smooth each condition independently: $P_{L A P, k}(x \mid y)=\frac{c(x, y)+k}{c(y)+k|X|}$


## Decision Networks

- MEU: choose the action which maximizes the expected utility given the evidence
- Can directly operationalize this with decision networks
- Bayes nets with nodes for utility and actions
- Lets us calculate the expected utility for each action
- New node types:
- Chance nodes (just like BNs)


Actions (rectangles, cannot have parents, act as observed evidence)


- Utility node (diamond, depends on action and chance nodes)


## Decision Networks

- Action selection:
- Instantiate all evidence
- Set action node(s) each possible way
- Calculate posterior for all parents of utility node, given the evidence
- Calculate expected utility for each action
- Choose maximizing action



## Example: Decision Networks

Umbrella = leave

$$
\begin{aligned}
& \mathrm{EU}(\text { leave })=\sum_{w} P(w) U(\text { leave }, w) \\
& =0.7 \cdot 100+0.3 \cdot 0=70
\end{aligned}
$$

Umbrella = take

| $\mathrm{EU}($ take $)=\sum_{w} P(w) U($ take,$w)$ |
| :--- |
| $=0.7 \cdot 20+0.3 \cdot 70=35$ |
|  |
| Optimal decision $=$ leave |

$$
\operatorname{MEU}(\varnothing)=\max _{a} \mathrm{EU}(a)=70
$$



## Example: Decision Networks



Optimal decision $=$ take

$$
\operatorname{MEU}(F=\mathrm{bad})=\max _{a} \mathrm{EU}(a \mid \mathrm{bad})=53
$$



## VPI Example: Weather

MEU with no evidence

$$
\operatorname{MEU}(\varnothing)=\max _{a} \mathrm{EU}(a)=70
$$

MEU if forecast is bad

$$
\operatorname{MEU}(F=\mathrm{bad})=\max _{a} \mathrm{EU}(a \mid \mathrm{bad})=53
$$

MEU if forecast is good

$$
\operatorname{MEU}(F=\operatorname{good})=\max _{a} \mathrm{EU}(a \mid \text { good })=95
$$



Forecast distribution

$$
\begin{array}{r}
0.59 \cdot(95)+0.41 \cdot(53)-70 \\
77.8-70=7.8
\end{array}
$$

$$
\operatorname{VPI}\left(E \mid e^{\prime}\right)=\left(\sum_{e^{\prime}} P\left(e^{\prime} \mid e\right) \operatorname{MEU}\left(e, e^{\prime}\right)\right)-\operatorname{MEU}(e)
$$

## Value of Information

- Assume we have evidence $\mathrm{E}=\mathrm{e}$. Value if we act now: $\operatorname{MEU}(e)=\max _{a} \sum_{s} P(s \mid e) U(s, a)$
- Assume we see that $E^{\prime}=e^{\prime}$. Value if we act then:
 $\operatorname{MEU}\left(e, e^{\prime}\right)=\max _{a} \sum_{s} P\left(s \mid e, e^{\prime}\right) U(s, a)$
- BUT E' is a random variable whose value is unknown, so we don't know what e' will be
- Expected value if $\mathrm{E}^{\prime}$ is revealed and then we act: $\operatorname{MEU}\left(e, E^{\prime}\right)=\sum_{e^{\prime}} P\left(e^{\prime} \mid e\right) \operatorname{MEU}\left(e, e^{\prime}\right)$
- Value of information: how much MEU goes up by revealing E' first then acting, over acting now:

$\operatorname{VPI}\left(E^{\prime} \mid e\right)=\operatorname{MEU}\left(e, E^{\prime}\right)-\operatorname{MEU}(e)$


## Example: Ghostbusters

- In (static) Ghostbusters:
- Belief state determined by evidence to date $\{\mathrm{e}\}$
- Tree really over evidence sets
- Probabilistic reasoning needed to predict new evidence given past evidence

- Solving POMDPs
- One way: use truncated expectimax to compute approximate value of actions
- What if you only considered busting or one sense followed by a bust?
- You get a VPI-based agent!


