

Consider a distribution defined over many variables

$$\pi(\underline{x}) \quad \text{with} \quad \underline{x} = (x_0, x_1, \dots, x_N)$$

N large (eg. $N=100$)

Suppose each x_i can take k possible values so

$$x_i \in S = \{s_1, \dots, s_k\}$$

Total number of states \underline{x} is k^N — very big — this makes it hard to do computations.

E.g. Suppose $\pi(\underline{x}) = \frac{1}{Z} e^{-E(\underline{x})}$ Gibbs distribution

Then computing $Z = \sum_{\underline{x}} e^{-E(\underline{x})}$ may require k^N computations — evaluate all possible states — too many!

computing $\hat{\underline{x}} = \underset{\underline{x}}{\text{ARG MAX}} \pi(\underline{x})$ — $\pi(\hat{\underline{x}}) \geq \pi(\underline{x})$ for all \underline{x} may require k^N computations — too many! (unless N & k small)

Also, how to sample from $\pi(\underline{x})$ to get i.i.d

Samples $\underline{x}^1 = (x_0^1, \dots, x_N^1)$

$$\underline{x}^2 = (x_0^2, \dots, x_N^2)$$

$$\underline{x}^3, \underline{x}^4, \dots$$

?

How to estimate quantities like $\sum_{\underline{x}} \pi(\underline{x}) h(\underline{x})$ for some $h(\underline{x})$?

(2) Today we discuss a special class of probability distributions for which we can solve these problems efficiently by Dynamic Programming (DP).

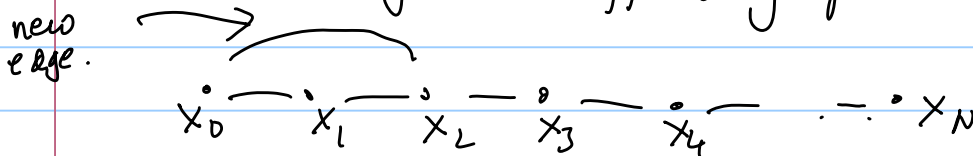
Background probability distributions $\pi(x)$ can be represented graphically, by assigning a node to each variable x_i , and an edge linking nodes which are directly related (Markov condition).

$$x_0 \quad x_1 \quad x_2 \quad x_3 \quad x_4 \quad \dots \quad x_N$$

$$\pi(x) = \frac{1}{Z} e^{-\sum_{i=1}^N \varphi_i(x_{i-1}, x_i)} \quad \varphi(\cdot, \cdot) \text{ is a potential.}$$

The potentials $\varphi_i(x_{i-1}, x_i)$ 'relate' variables x_{i-1} to x_i - so have 'edges' in the graph.

If we added a new potential $\varphi(x_0, x_2)$
 - so $\pi(x) = \frac{1}{\tilde{Z}} e^{-\varphi(x_0, x_2) - \sum_{i=1}^N \varphi_i(x_{i-1}, x_i)}$ -
 (note $\tilde{Z} \neq Z$, different normalization constant.)
 then we would get a different graph.



Today's lecture will stick to a distribution

$$\pi(x) = \frac{1}{Z} e^{-\sum_{i=1}^N \varphi_i(x_{i-1}, x_i)} \quad \left(\text{graph has no closed loops} \right)$$

(3)

Probabilities like $\pi(x) = \frac{1}{Z} e^{-\sum_{i=1}^N \phi_i(x_{i-1}, x_i)}$

Note Title

can be used to model events over time - eg. 4/2/2006

x_i is state of system at time i

x_{i+1} is state of system at time $i+1$

state of system at time i is specified (probabilistically) by state of system at time $i-1$ (more on this next lecture.)

Or, more generally, any 1-D structure like the sequences of DNA

E.g. GATTACA...

A - adenine

C - cytosine

G - guanine

T - thymine.

And many others...

Dynamic Programming can be used to find the global maximum of $\pi(x)$, \bar{x} , and $\pi(\bar{x})$ in $O(Nk^2)$ operations. (not k^N)

DP can also compute $Z = \sum_x e^{-\sum_{i=1}^N \phi_i(x_{i-1}, x_i)}$ in $O(Nk^2)$ operations

DP can also find the marginal distribution $\pi_i(x_i)$ and draw exact random samples from $\pi(x)$ efficiently.

(4)

Maximizing $\Pi(\underline{x})$ is equivalent to
minimizing $E(\underline{x}) = \varphi_1(x_0, x_1) + \dots + \varphi_N(x_{N-1}, x_N)$

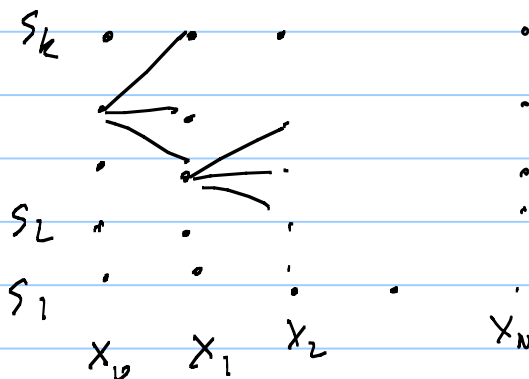
Forward Pass of DP

DP acts recursively: $\left(\begin{array}{l} \text{ie. } x_0 = s_1 \text{ or} \\ s_2 \text{ or } s_3 \text{ or} \\ \dots \text{ or } s_k \end{array} \right)$

• Define $m_1(x_1) = \min_{s_i \in S} \varphi_1(s_i, x_1)$ for $x_1 = s_1, \dots, s_k$

• Recursively compute $m_t(x_t) = \min_{s_i \in S} \{ m_{t-1}(s_i) + \varphi_t(s_i, x_t) \}$ for $x_t = s_1, \dots, s_k$.

Claim: Optimal value $E(\underline{x})$ is obtained by $\min_{s \in \{s_1, \dots, s_k\}} m_N(x_N)$



To compute $m_1(x_1)$ for $x_1 = s_1, \dots, s_k$

needs $O(k^2)$ operations.

Computing all $m_t(x_t) \dots$ requires $O(kN^2)$

Justify Claim: minimum of $m_t(x_t)$ is the minimum of $\varphi_1(x_0, x_1)$

proof by induction: $\min_{x_t \in S} m_t(x_t) = \min_{x_0, \dots, x_1 \in S} \{ \varphi_1(x_0, x_1) + \dots + \varphi_t(x_{t-1}, x_t) \}$.

(5)

To find the optimal path \hat{x} we trace backwards
Backward Pass of DP

• Let \hat{x}_N be the minimizer of $m_N(x_N)$
$$\hat{x}_N = \arg \min_{s_i \in S} m_N(s_i)$$

(Break ties arbitrarily - ties mean that there are several \underline{x} 's st. $E(\underline{x}) = E(\underline{x}')$)

• For $t = N-1, N-2, \dots, 0$

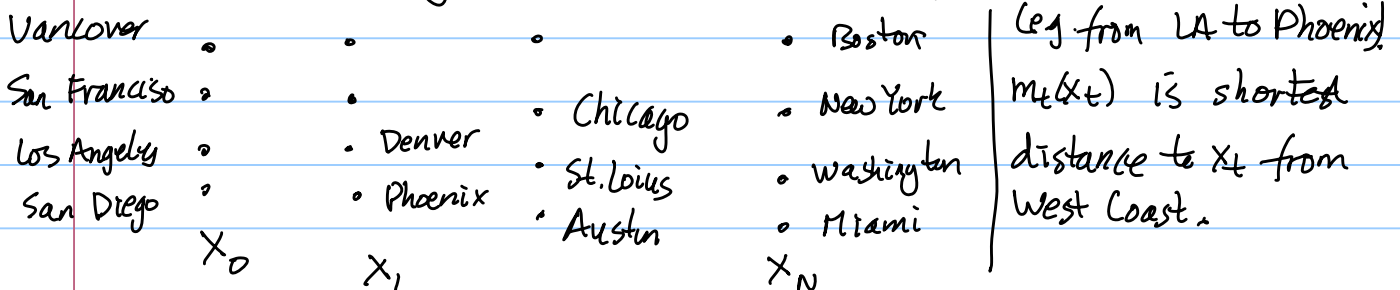
Let $\hat{x}_t = \arg \min_{s_i \in S} \{m_t(s_i) + \varphi_{t+1}(s_i, \hat{x}_{t+1})\}$
(Break ties arbitrarily)

$\hat{x} = (\hat{x}_0, \dots, \hat{x}_N)$ obtained in

this way is the minimizer. If ties, then several minimizers \underline{x}
(doesn't matter which you pick)

Example: Task - find shortest path from West Coast of US to East Coast.

On date t , you start in city x_{t+1} and drive distance $\varphi_{t+1}(x_{t+1}, x_t)$ to city x_t



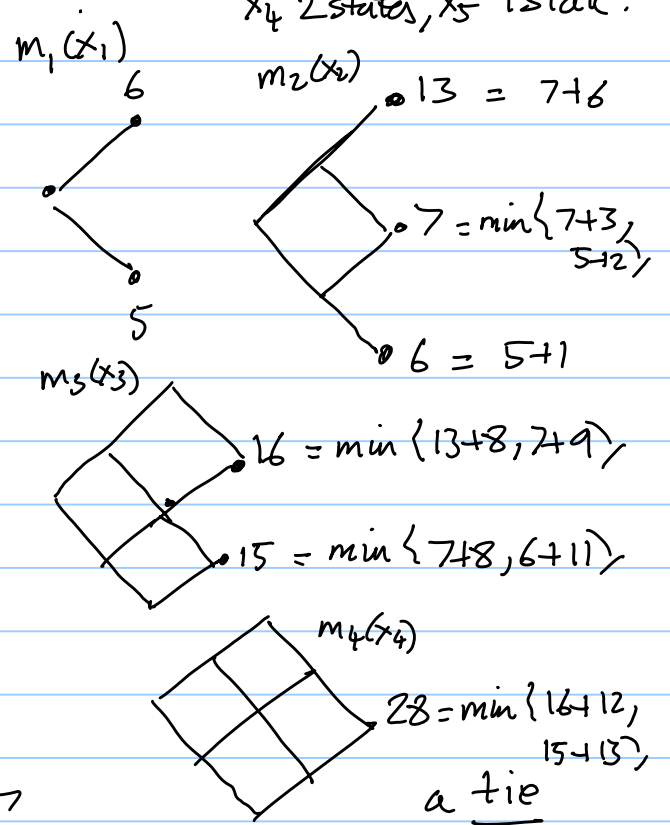
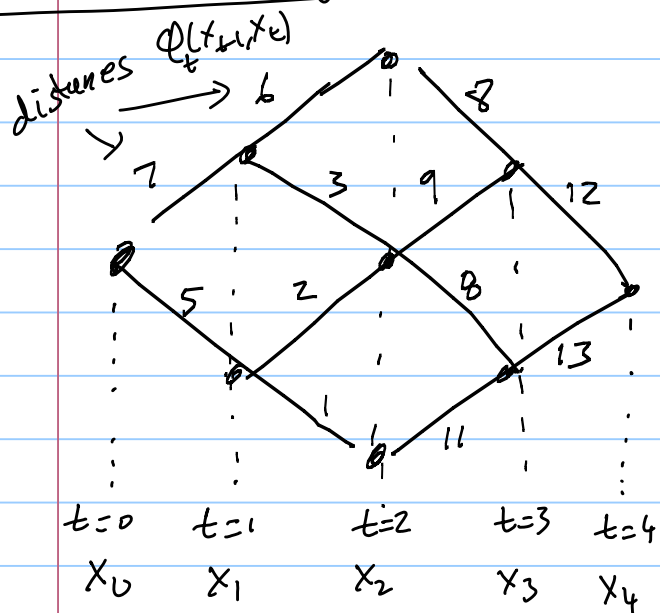
(6) Intuition: DP solves the minimization problem efficiently by breaking it down into separate parts.

E.g. If you want to find the shortest path between the West Coast and the East Coast which goes through Chicago, then you can do this by solving two independent problems:

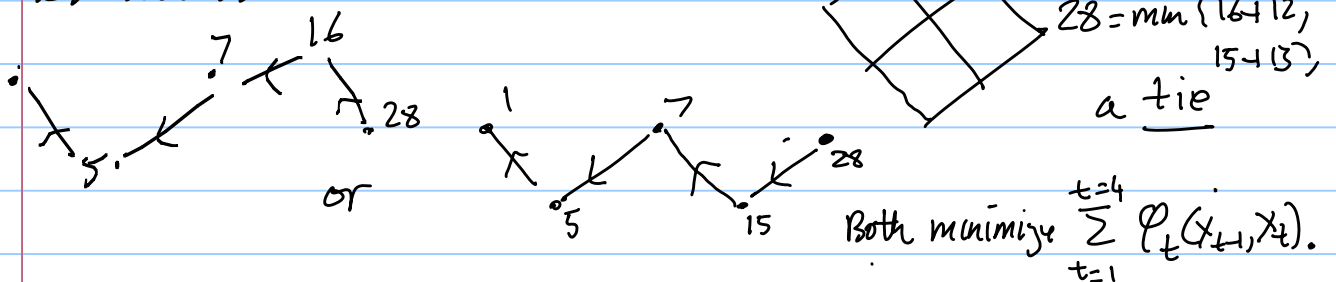
- (i) Find shortest path from West Coast to Chicago
- (ii) Find shortest path from Chicago to East Coast

Note: DP only works if the probability distribution $\pi(x)$ is defined on a graph with no closed loops. Like $x_0 \rightarrow x_1 \rightarrow x_2 \rightarrow x_3$

Simplest Example of DP: x_0 has 1 state only, x_1 2 states, x_2 3 states, x_3 2 states, x_4 1 state.



Two Best Solutions:



Both minimize $\sum_{t=1}^{t=4} \rho_t(x_{t-1}, x_t)$.