

CPSC 540: Machine Learning

Message Passing

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Last Time: Monte Carlo Methods

- If we want to approximate expectations of random functions,

$$\mathbb{E}[g(x)] = \underbrace{\sum_{x \in \mathcal{X}} g(x)p(x)}_{\text{discrete } x} \quad \text{or} \quad \mathbb{E}[g(x)] = \underbrace{\int_{x \in \mathcal{X}} g(x)p(x)dx}_{\text{continuous } x},$$

the Monte Carlo estimate is

$$\mathbb{E}[g(x)] \approx \frac{1}{n} \sum_{i=1}^n g(x^i),$$

where the x^i are independent samples from $p(x)$.

- We can use this to approximate marginals,

$$p(x_j = c) = \frac{1}{n} \sum_{i=1}^n \mathcal{I}[x_j^i = c].$$

Exact Marginal Calculation

- In typical settings Monte Carlo has **sublinear convergence** like stochastic gradient.
 - $O(1/t)$ convergence rate where constant is **variance** of samples.
 - If all samples look the same, it converges quickly.
 - If samples look very different, it can be **painfully slow**.
- For **discrete-state** Markov chains, we can actually **compute marginals directly**:
 - We're given **initial probabilities** $p(x_1 = s)$ for all s as part of the definition.
 - We can use **transition probabilities** to **compute** $p(x_2 = s)$ for all s :

$$p(x_2) = \underbrace{\sum_{x_1=1}^k p(x_2, x_1)}_{\text{marginalization rule}} = \sum_{x_1=1}^k \underbrace{p(x_2 | x_1)p(x_1)}_{\text{product rule}}.$$

- We can repeat this calculation to obtain $p(x_3 = s)$ and subsequent marginals.

Exact Marginal Calculation

- Recursive formula for marginals at time j :

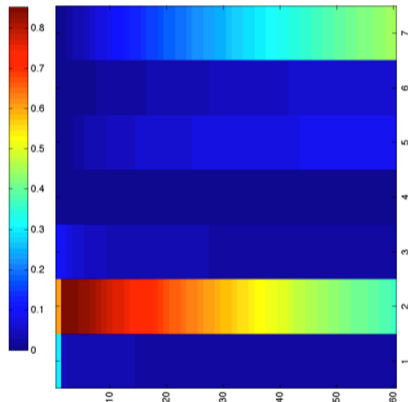
$$p(x_j) = \sum_{x_{j-1}=1}^k p(x_j | x_{j-1})p(x_{j-1}),$$

called the **Chapman-Kolmogorov (CK) equations**.

- Cost:
 - Given previous time, CK equations for one x_j costs $O(k)$.
 - Given previous time, to compute $p(x_j)$ for all k states costs $O(k^2)$.
 - Can be written as matrix-vector product with $k \times k$ transition probabilities matrix.
 - So **cost to compute marginals up to time d is $O(dk^2)$** .
 - An example of **dynamic programming**: efficiently sums over k^d paths.

Marginals in CS Grad Career

- CK equations can give all marginals $p(x_j = c)$ from CS grad Markov chain:



- Each row j is a state and each column c is a year.

Continuous-State Markov Chains

- The CK equations also apply if we have **continuous states**:

$$p(x_j) = \int_{x_{j-1}} p(x_j | x_{j-1})p(x_{j-1}),$$

but this integral **may not have a closed-form solution**.

- **Gaussian probabilities** are an important special case:
 - If $p(x_{j-1})$ and $p(x_j | x_{j-1})$ are Gaussian, then $p(x_j)$ is Gaussian.
 - So we can write $p(x_j)$ in closed-form in terms of mean and variance.
- If the probabilities are non-Gaussian, usually **can't represent $p(x_j)$ distribution**.
 - You are stuck using Monte Carlo or other approximations.

Stationary Distribution

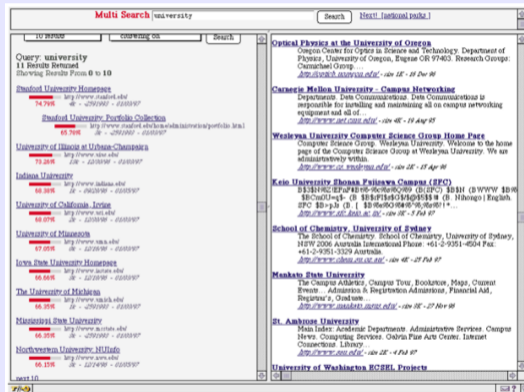
- A **stationary distribution** of a homogeneous Markov chain is a vector π satisfying

$$\pi(c) = \sum_{c'} p(x_j = c \mid x_{j-1} = c')\pi(c').$$

- “The probabilities don’t change across time” (also called **“invariant” distribution**).
- Under certain conditions, **marginals converge to a stationary distribution**.
 - $p(x_j = c) \rightarrow \pi(c)$ as j goes to ∞ .
 - If we fit a Markov chain to the rain example, we have $\pi(\text{“rain”}) = 0.41$.
 - In the CS grad student example, we have $\pi(\text{“dead”}) = 1$.
- Stationary distribution is basis for Google’s **PageRank** algorithm.

Application: PageRank

- Web search before Google:

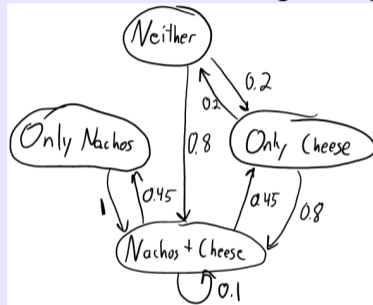


<http://ilpubs.stanford.edu:8090/422/1/1999-66.pdf>

- It was also easy to fool search engines by copying popular websites.

State Transition Diagram

- **State transition diagrams** are common for visualizing homogenous Markov chains:



$$P = \begin{bmatrix} 0 & 0 & 0.2 & 0.8 \\ 0 & 0 & 0 & 1 \\ 0.2 & 0 & 0 & 0.8 \\ 0 & 0.45 & 0.45 & 0.1 \end{bmatrix}$$

- Each **node** is a state, each **edge** is a non-zero transition probability.
 - For web-search, each **node** will be a webpage.
- **Cost of CK equations** is only $O(z)$ if you have only z edges.

Application: PageRank

- Wikipedia's cartoon illustration of Google's PageRank:
 - Large face means higher rank.



<https://en.wikipedia.org/wiki/PageRank>

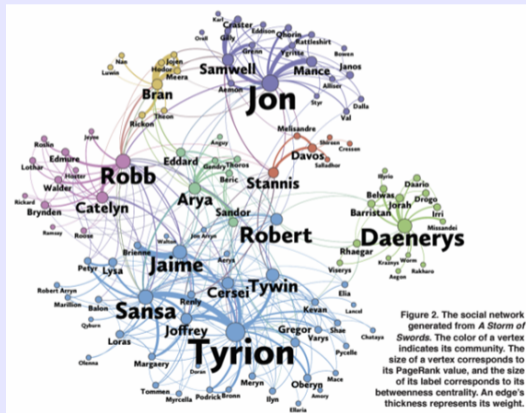
- “Important webpages are linked from other important webpages”.
- “Link is more meaningful if a webpage has few links”.

Application: PageRank

- Google's **PageRank** algorithm for measuring the importance of a website:
 - Stationary probability in “random surfer” Markov chain:
 - With probability α , surfer clicks on a **random link** on the current webpage.
 - Otherwise, surfer goes to a **completely random webpage**.
- To compute the stationary distribution, they use the **power method**:
 - Repeatedly apply the CK equations.
 - Iterations are faster than $O(k^2)$ due to sparsity of links.
 - Can be easily parallelized.
 - Achieves a linear convergence rate.
- More recent works have shown **coordinate optimization can be faster**.

Application: Game of Thrones

- PageRank can be used in other applications.
- “Who is the main character in the Game of Thrones books?”



Existence/Uniqueness of Stationary Distribution

- Does a stationary distribution π **exist** and is it **unique**?
- A sufficient condition for existence/uniqueness is that all $p(x_j = c \mid x_{j'} = c') > 0$.
 - PageRank satisfies this by adding probability α of jumping to a random page.
- Weaker sufficient conditions for existence and uniqueness (“ergodic”):
 - 1 “Irreducible” (doesn’t get stuck in part of the graph).
 - 2 “Aperiodic” (probability of returning to state isn’t on fixed intervals).

Outline

- 1 Exact Marginals and PageRank
- 2 Message Passing

Decoding: Maximizing Joint Probability

- **Decoding** in density models: finding x with highest joint probability:

$$\operatorname{argmax}_{x_1, x_2, \dots, x_d} p(x_1, x_2, \dots, x_d).$$

- For CS grad student ($d = 60$) the decoding is “industry” for all years.
 - The decoding often doesn't look like a typical sample.
 - The decoding can change if you increase d .
- **Decoding is easy for independent** models:
 - We can just optimize each x_j independently.
 - For example, with four variables we have

$$\max_{x_1, x_2, x_3, x_4} \{p(x_1)p(x_2)p(x_3)p(x_4)\} = \left(\max_{x_1} p(x_1)\right) \left(\max_{x_2} p(x_2)\right) \left(\max_{x_3} p(x_3)\right) \left(\max_{x_4} p(x_4)\right).$$

- Can we also maximize the marginals to decode a Markov chain?

Example of Decoding vs. Maximizing Marginals

- Consider the “plane of doom” 2-variable Markov chain:

$$X = \begin{bmatrix} \text{“land”} & \text{“alive”} \\ \text{“land”} & \text{“alive”} \\ \text{“crash”} & \text{“dead”} \\ \text{“explode”} & \text{“dead”} \\ \text{“crash”} & \text{“dead”} \\ \text{“land”} & \text{“alive”} \\ \vdots & \vdots \end{bmatrix} .$$

- 40% of the time the plane lands and you live.
- 30% of the time the plane crashes and you die.
- 30% of the time the plane crashes and you die.

Example of Decoding vs. Maximizing Marginals

- Initial probabilities are given by

$$p(x_1 = \text{"land"}) = 0.4, \quad p(x_1 = \text{"crash"}) = 0.3, \quad p(x_1 = \text{"explode"}) = 0.3,$$

and x_2 is "alive" iff x_1 is "land".

- If we apply the **CK equations** we get

$$p(x_2 = \text{"alive"}) = 0.4, \quad p(x_2 = \text{"dead"}) = 0.6,$$

so maximizing the marginals $p(x_j)$ independently gives ("land", "dead").

- This actually has probability 0.
- **Decoding** considers the **joint assignment to x_1 and x_2** maximizing probability.
 - In this case it's ("land", "alive"), which has probability 0.4.

Distributing Max across Product

- Note that decoding **can't be done forward in time** as in CK equations.
 - We need to **optimize over all k^d assignments to all variables**.
 - Even if $p(x_1 = 1) = 0.99$, the most likely sequence could have $x_1 = 2$.
- Fortunately, the **Markov property** makes the max simplify:

$$\begin{aligned}
 \max_{x_1, x_2, x_3, x_4} p(x_1, x_2, x_3, x_4) &= \max_{x_1, x_2, x_3, x_4} p(x_4 | x_3) p(x_3 | x_2) p(x_2 | x_1) p(x_1) \\
 &= \max_{x_4} \max_{x_3} \max_{x_2} \max_{x_1} p(x_4 | x_3) p(x_3 | x_2) p(x_2 | x_1) p(x_1) \\
 &= \max_{x_4} \max_{x_3} \max_{x_2} p(x_4 | x_3) p(x_3 | x_2) \max_{x_1} p(x_2 | x_1) p(x_1) \\
 &= \max_{x_4} \max_{x_3} p(x_4 | x_3) \max_{x_2} p(x_3 | x_2) \max_{x_1} p(x_2 | x_1) p(x_1),
 \end{aligned}$$

where we're using that $\max_i \alpha a_i = \alpha \max_i a_i$ for non-negative α .

Decoding with Memoization

- The Markov property writes decoding as a sequence of max problems:

$$\max_{x_1, x_2, x_3, x_4} p(x_1, x_2, x_3, x_4) = \max_{x_4} \max_{x_3} p(x_4 | x_3) \max_{x_2} p(x_3 | x_2) \max_{x_1} p(x_2 | x_1) p(x_1),$$

but note that we can't just "solve" \max_{x_1} once because it's a function of x_2 .

- Instead, we'll memoize solution $M_2(x_2) = \max_{x_1} p(x_2 | x_1) p(x_1)$ for all x_2 ,

$$\max_{x_1, x_2, x_3, x_4} p(x_1, x_2, x_3, x_4) = \max_{x_4} \max_{x_3} p(x_4 | x_3) \max_{x_2} p(x_3 | x_2) M_2(x_2).$$

- Now we memoize solution $M_3(x_3) = \max_{x_2} p(x_3 | x_2) M_2(x_2)$ for all x_3 ,

$$\max_{x_1, x_2, x_3, x_4} p(x_1, x_2, x_3, x_4) = \max_{x_4} \max_{x_3} p(x_4 | x_3) M_3(x_3).$$

- And defining $M_4(x_4) = \max_{x_3} p(x_4 | x_3) M_3(x_3)$ the maximum value is given by

$$\max_{x_1, x_2, x_3, x_4} p(x_1, x_2, x_3, x_4) = \max_{x_4} M_4(x_4).$$

Example: Decoding the Plane of Doom

- We have $M_1(x_1) = p(x_1)$ so in “plane of doom” we have

$$M_1(\text{“land”}) = 0.4, \quad M_1(\text{“crash”}) = 0.3, \quad M_1(\text{“explode”}) = 0.3.$$

- We have $M_2(x_2) = \max_{x_1} p(x_2 | x_1)M_1(x_1)$ so we get

$$M_2(\text{“alive”}) = 0.4, \quad M_2(\text{“dead”}) = 0.3.$$

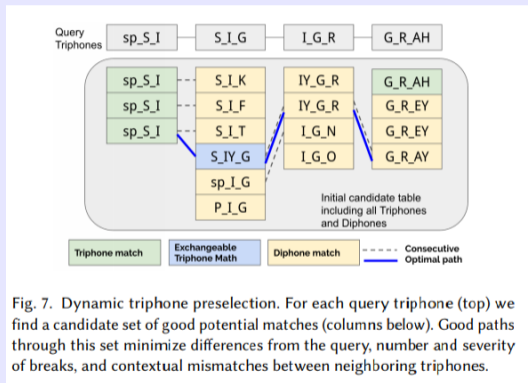
- $M_2(2) \neq p(x_2 = 2)$ because we **needed to choose either “crash” or “explode”**.
- We maximize $M_2(x_2)$ to find that the optimal decoding ends with “alive”.
 - We now need to **backtrack** to find the state that lead to “alive”, giving “land”.

Viterbi Decoding

- What is $M_j(x_j)$ in words?
 - “Probability of most likely length- j sequence that ends in x_j (ignoring future)”.
- The **Viterbi decoding** algorithm (special case of **dynamic programming**):
 - 1 Set $M_1(x_1) = p(x_1)$ for all x_1 .
 - 2 Compute $M_2(x_2)$ for all x_2 , store value of x_1 leading to the best value of each x_2 .
 - 3 Compute $M_3(x_3)$ for all x_3 , store value of x_2 leading to the best value of each x_3 .
 - 4 ...
 - 5 Maximize $M_d(x_d)$ to find value of x_d in a decoding.
 - 6 **Backtrack** to find the value of x_{d-1} that lead to this x_d .
 - 7 Backtrack to find the value of x_{d-2} that lead to this x_{d-1} .
 - 8 ...
- Computing all $M_j(x_j)$ given all $M_{j-1}(x_{j-1})$ costs $O(k^2)$.
 - Total cost is only $O(dk^2)$ to search over all k^d paths.
 - Has numerous applications like decoding digital TV.

Application: Voice Photoshop

- Application: Adobe VoCo uses Viterbi as part of synthesizing voices:



http://gfx.cs.princeton.edu/pubs/Jin_2017_VTI/Jin2017-VoCo-paper.pdf

- <https://www.youtube.com/watch?v=I3l4XLZ59iw>

Summary

- **Chapman-Kolmogorov equations** compute exact univariate marginals.
 - For discrete or Gaussian Markov chains.
- **Stationary distribution** of homogenous Markov chain.
 - Marginals as time goes to ∞ .
 - Basis of Google's PageRank method.
- **Decoding** is task of finding most probable x .
- **Viterbi decoding** allow efficient decoding with Markov chains.
- Next time: measuring defence in the NBA.

Label Propagation as a Markov Chain Problem

- Basic **label propagation** method has a Markov chain interpretation.
 - We have $n + t$ states, one for each [un]labeled example.
- Monte Carlo approach to label propagation (“adsorption”):
 - At time $t = 0$, set the state to the node you want to label.
 - At time $t > 0$ and on a labeled node, output the label.
 - Labeled nodes are absorbing states.
 - At time $t > 0$ and on an unlabeled node i :
 - Move to neighbour j with probability proportional w_{ij} (or \bar{w}_{ij}).
- Final **predictions are probabilities of outputting each label**.
 - Nice if you only need to label one example at a time (slow if labels are rare).
 - Common hack is to limit random walk time to bound runtime.