

CPSC 540: Machine Learning

Monte Carlo Methods

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Last Time: Markov Chains

- We can use **Markov chains** for density estimation,

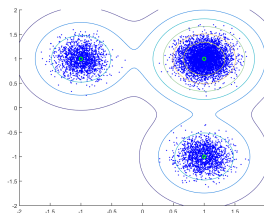
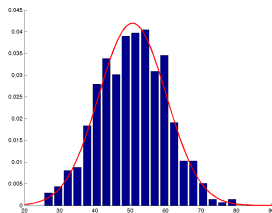
$$p(x) = \underbrace{p(x_1)}_{\text{initial prob.}} \prod_{j=2}^d \underbrace{p(x_j | x_{j-1})}_{\text{transition prob.}},$$

which model **dependency between adjacent features**.

- Different than fmixture models which focus on describe clusters in the data.
- **Homogeneous** chains use same transition probability for all j (**parameter tieing**).
 - Gives more data to estimate transitions, allows examples of different sizes.
- **Inhomogeneous** chains allow different transitions at different times.
- Given a Markov chain model, we overviewed common **computational problems**:
 - Sampling, inference, decoding, conditioning, and stationary distribution.

Fundamental Problem: Sampling from a Density

- A fundamental problem in density estimation is **sampling from the density**.
 - Generating examples x^i that are distributed according to a given density $p(x)$.
 - Basically, the “opposite” of density estimation.



- We've been using pictures of samples to “tell us what the model has learned”.
 - If the samples look like real data, then we have a good density model.
- Samples can also be used in **Monte Carlo** estimation (today):
 - Replace complicated $p(x)$ with **samples** to solve hard problems at test time.

Simplest Case: Sampling from a Bernoulli

- Consider **sampling from a Bernoulli**, for example

$$p(x = 1) = 0.9, \quad p(x = 0) = 0.1.$$

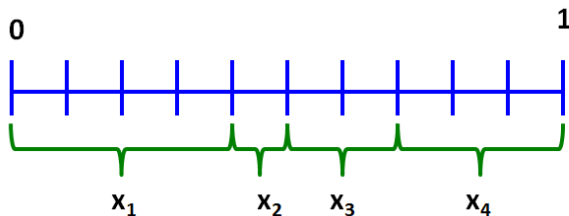
- Sampling methods **assume we can sample uniformly over $[0, 1]$** .
 - Usually, a "pseudo-random" number generator is good enough (like Julia's *rand*).
- How to use a **uniform sample to sample from the Bernoulli** above:
 - 1 Generate a uniform sample $u \sim \mathcal{U}(0, 1)$.
 - 2 If $u \leq 0.9$, set $x = 1$ (otherwise, set $x = 0$).
- If uniform samples are "good enough", then we have $x = 1$ with probability 0.9.

Sampling from a Categorical Distribution

- Consider a more general **categorical density** like

$$p(x = 1) = 0.4, \quad p(x = 2) = 0.1, \quad p(x = 3) = 0.2, \quad p(x = 4) = 0.3,$$

we can divide up the $[0, 1]$ interval based on probability values:



- If $u \sim \mathcal{U}(0, 1)$, 40% of the time it lands in x_1 region, 10% of time in x_2 , and so on.

Sampling from a Categorical Distribution

- Consider a more general **categorical density** like

$$p(x = 1) = 0.4, \quad p(x = 2) = 0.1, \quad p(x = 3) = 0.2, \quad p(x = 4) = 0.3.$$

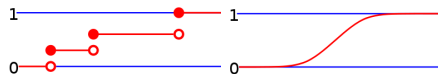
- To **sample from this categorical** density we can use (*sampleDiscrete.jl*):
 - 1 Generate $u \sim \mathcal{U}(0, 1)$.
 - 2 If $u \leq 0.4$, output 1.
 - 3 If $u \leq 0.4 + 0.1$, output 2.
 - 4 If $u \leq 0.4 + 0.1 + 0.2$, output 3.
 - 5 Otherwise, output 4.

Sampling from a Categorical Distribution

- General case for sampling from categorical.
 - 1 Generate $u \sim \mathcal{U}(0, 1)$.
 - 2 If $u \leq p(x = 1)$, output 1.
 - 3 If $u \leq p(x \leq 2)$, output 2.
 - 4 If $u \leq p(x \leq 3)$, output 3.
 - 5 ...
- The value $p(x \leq c) = p(x = 1) + p(x = 2) + \dots + p(x = c)$ is the **CDF**.
 - “**Cumulative distribution function**”.
- Worst case cost with k possible states is $O(k)$ by incrementally computing CDFs.
- But to generate t samples only **costs** $O(k + t \log k)$:
 - One-time $O(k)$ cost to store the CDF $p(x \leq c)$ for each c .
 - Per-sample $O(\log k)$ cost to do **binary search** for smallest c with $u \leq p(x \leq c)$.

Inverse Transform Method (Exact 1D Sampling)

- We often use $F(c) = p(x \leq c)$ to denote the CDF.
 - $F(c)$ is between 0 and 1 and gives proportion of times x is below c .
 - F can be used for discrete and continuous variables:



https://en.wikipedia.org/wiki/Cumulative_distribution_function

- The **inverse CDF** (or “**quantile**” function) F^{-1} is its inverse:
 - Given a number u between 0 and 1, returns c such that $p(x \leq c) = u$.
- **Inverse transform** method for exact sampling in 1D:
 - 1 Sample $u \sim \mathcal{U}(0, 1)$.
 - 2 Return $F^{-1}(u)$.
- Video on pseudo-random numbers and inverse-transform sampling:
 - <https://www.youtube.com/watch?v=C82JyCmtKWg>

Sampling from a 1D Gaussian

- Consider a Gaussian distribution,

$$x \sim \mathcal{N}(\mu, \sigma^2).$$

- CDF has the form

$$F(x) = p(x \leq c) = \frac{1}{2} \left[1 + \operatorname{erf} \left(\frac{c - \mu}{\sigma\sqrt{2}} \right) \right],$$

where “erf” is the CDF of $\mathcal{N}(0, 1)$.

- Inverse CDF has the form

$$F^{-1}(u) = \mu + \sigma\sqrt{2}\operatorname{erf}^{-1}(2u - 1).$$

- To sample from a Gaussian:

- 1 Generate $u \sim \mathcal{U}(0, 1)$.
- 2 Return $F^{-1}(u)$.

Sampling from a Product Distribution

- Consider a **product distribution**,

$$p(x_1, x_2, \dots, x_d) = p(x_1)p(x_2) \cdots p(x_d).$$

- Because variables are **independent**, we can **sample independently**:
 - Sample x_1 from $p(x_1)$.
 - Sample x_2 from $p(x_2)$.
 - ...
 - Sample x_d from $p(x_d)$.
- Example: sampling from a **multivariate Gaussian with diagonal covariance**.
 - Sample each variable independently based on μ_j and σ_j^2 .

Digression: Sampling from a Multivariate Gaussian

- In some cases we can sample from multivariate distributions by transformation.
- Recall the affine property of multivariate Gaussian:
 - If $x \sim \mathcal{N}(\mu, \Sigma)$, then $Ax + b \sim \mathcal{N}(A\mu + b, A\Sigma A^T)$.
- To sample from a general multivariate Gaussian $\mathcal{N}(\mu, \Sigma)$:
 - 1 Sample x from a $\mathcal{N}(0, I)$ (each x_j coming independently from $\mathcal{N}(0, 1)$).
 - 2 Transform to a sample from the right Gaussian using the affine property:

$$Ax + \mu \sim \mathcal{N}(\mu, AA^T),$$

where we choose A so that $AA^T = \Sigma$ (e.g., by Cholesky factorization).

Ancestral Sampling

- Another way to sample non-product distributions is using **chain rule**,

$$p(x_1, x_2, x_3, \dots, x_d) = p(x_1)p(x_2 \mid x_1)p(x_3 \mid x_2, x_1) \cdots p(x_d \mid x_{d-1}, x_{d-2}, \dots, x_1),$$

which comes from repeated application of the **product rule**
($p(a, b) = p(a)p(b \mid a)$).

- The chain rule suggests the following sampling strategy:
 - Sample x_1 from $p(x_1)$.
 - Given x_1 , sample x_2 from $p(x_2 \mid x_1)$.
 - Given x_1 and x_2 , sample x_3 from $p(x_3 \mid x_2, x_1)$.
 - ...
 - Given x_1 through x_{d-1} , sample x_d from $p(x_d \mid x_{d-1}, x_{d-2}, \dots, x_1)$.
- This is called **ancestral sampling**.
 - It's easy if (conditional) probabilities are simple, since sampling in 1D is usually easy.

Ancestral Sampling Examples

- For **Markov chains** the **chain rule simplifies** to

$$p(x_1, x_2, x_3, \dots, x_d) = p(x_1)p(x_2 \mid x_1)p(x_3 \mid x_2) \cdots p(x_d \mid x_{d-1}),$$

- So **ancestral sampling simplifies** too:

- ➊ Sample x_1 from initial probabilities $p(x_1)$.
- ➋ Given x_1 , sample x_2 from transition probabilities $p(x_2 \mid x_1)$.
- ➌ Given x_2 , sample x_3 from transition probabilities $p(x_3 \mid x_2)$.
- ➍ ...
- ➎ Given x_{d-1} , sample x_d from transition probabilities $p(x_d \mid x_{d-1})$.

- For **mixture models** with cluster variables z we could write

$$p(x, z) = p(z)p(x \mid z),$$

so we can **first sample cluster z** and then **sample x given cluster z** .

- You can just ignore the z values to get samples of x .

Markov Chain Toy Example: CS Grad Career

- “Computer science grad career” Markov chain:

- Initial probabilities:

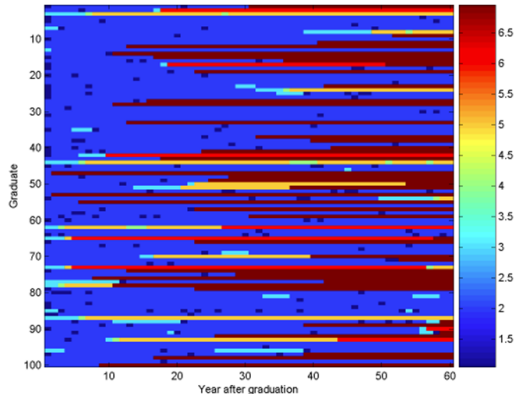
State	Probability	Description
Industry	0.60	They work for a company or own their own company.
Grad School	0.30	They are trying to get a Masters or PhD degree.
Video Games	0.10	They mostly play video games.

- Transition probabilities:

From\to	Video Games	Industry	Grad School	Video Games (with PhD)	Industry (with PhD)	Academia	Deceased
Video Games	0.08	0.90	0.01	0	0	0	0.01
Industry	0.03	0.95	0.01	0	0	0	0.01
Grad School	0.06	0.06	0.75	0.05	0.05	0.02	0.01
Video Games (with PhD)	0	0	0	0.30	0.60	0.09	0.01
Industry (with PhD)	0	0	0	0.02	0.95	0.02	0.01
Academia	0	0	0	0.01	0.01	0.97	0.01
Deceased	0	0	0	0	0	0	1

Markov Chain Toy Example: CS Grad Career

- Samples from “computer science grad career” Markov chain:



- State 7 (“deceased”) is called an **absorbing state** (no probability of leaving).
- Samples often give you an idea of what model knows (and what should be fixed).

Outline

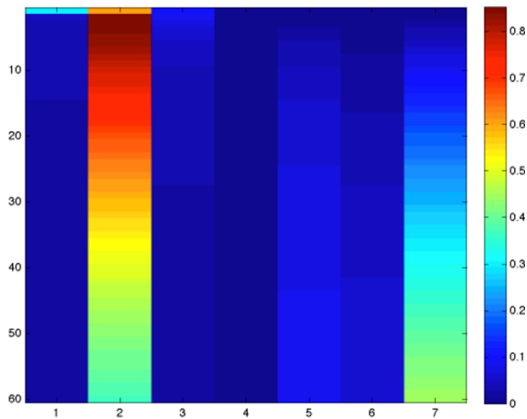
- 1 Introduction to Sampling
- 2 Monte Carlo Approximation

Marginal and Conditional Inference

- Given density estimator, we often want to make **probabilistic inferences**:
 - **Marginals**: what is the probability that $x_j = c$?
 - What is the probability we're in industry 10 years after graduation?
 - **Conditionals**: what is the probability that $x_j = c$ given $x_{j'} = c'$?
 - What is the probability of industry after 10 years, if we immediately go to grad school?
- This is easy for simple independent models:
 - We are directly modeling marginals $p(x_j)$.
 - By independence, conditionals are marginals: $p(x_j \mid x_{j'}) = p(x_j)$.
- This is also easy for mixtures of simple independent models.
 - Do inference for each mixture.
- For Markov chains, it's more complicated...

Marginals in CS Grad Career

- All marginals $p(x_j = c)$ from “computer science grad career” Markov chain:



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

Monte Carlo: Inference by Sampling

- A basic Monte Carlo method for estimating probabilities of events:

- 1 Generate a large number of samples x^i from the model,

$$X = \begin{bmatrix} 0 & 0 & 1 & 0 \\ 1 & 1 & 1 & 0 \\ 0 & 1 & 1 & 1 \\ 1 & 1 & 1 & 1 \end{bmatrix}.$$

- 2 Compute frequency that the event happened in the samples,

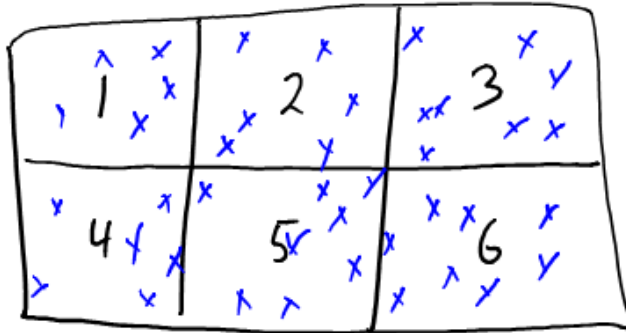
$$p(x_2 = 1) \approx 3/4,$$

$$p(x_3 = 0) \approx 0/4.$$

- Monte Carlo methods are second most important class of ML algorithms.
 - Originally developed to build better atomic bombs :(

Monte Carlo Method for Rolling Di

- Probability of event:
 - $(\text{number of samples where event happend})/(\text{number of samples})$



Monte Carlo Method for Inequalities

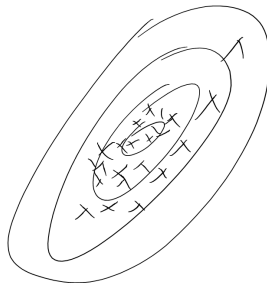
- Monte Carlo estimate of **probability that variable is above threshold**:
 - Compute fraction of examples where sample is above threshold.



Monte Carlo Method for Mean

- A Monte Carlo approximation of the mean:
 - Approximate the mean by average of samples.

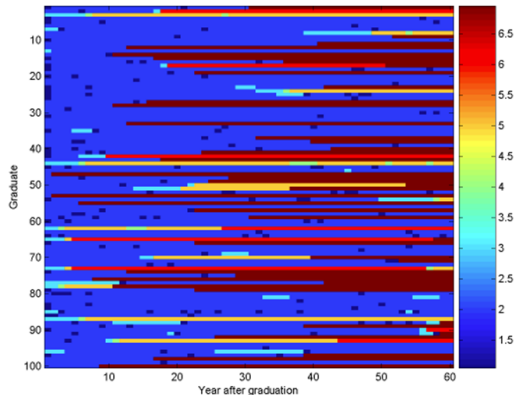
$$E[x] \approx \frac{1}{n} \sum_{i=1}^n x^i.$$



- Visual demo of Monte Carlo approximation of mean and variance:
 - <http://students.brown.edu/seeing-theory/basic-probability/index.html>

Monte Carlo for Markov Chains

- Our samples from the CS grad student Markov chain:



- We can estimate probabilities by looking at frequencies in samples.
- This works for continuous states too.

Monte Carlo Methods

- Monte Carlo methods approximate expectations of random functions,

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

- Computing mean is the special case of $g(x) = x$.
- Computing probability of any event A is also a special case:
 - Set $g(x) = \mathcal{I}[\text{"}A \text{ happened in sample } x^i\text{"}]$.
- To approximate expectation, generate n samples x^i from $p(x)$ and use:

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

Monte Carlo Methods

- Monte Carlo estimate is **unbiased** approximation of expectation,

$$\begin{aligned}\mathbb{E} \left[\frac{1}{n} \sum_{i=1}^n g(x^i) \right] &= \frac{1}{n} \sum_{i=1}^n \mathbb{E}[g(x^i)] \\ &= \frac{1}{n} \sum_{i=1}^n \mathbb{E}[g(x)] = \mathbb{E}[g(x)],\end{aligned}$$

- The **law of large numbers** says that:
 - Unbiased approximators “converge” (probabilistically) to expectation as $n \rightarrow \infty$.
 - So the more samples you get, the closer to the true value you expect to get.
- Convergence rate is $O(1/n)$ (sublinear), similar to stochastic gradient (bonus).

Monte Carlo Methods for Markov Chain Inference

- Monte Carlo methods allow approximating expectations in Markov chains:
 - Marginal $p(x_j = c)$ is the number of chains that were in state c at time j .
 - Average value at time j , $E[x_j]$, is approximated by average of x_j in the samples.
 - $p(x_j \leq 10)$ is approximate by frequency of x_j being less than 10.
 - $p(x_j \leq 10, x_{j+1} \geq 10)$ is approximated by number of chains where both happen.

Monte Carlo for Conditional Probabilities

- We often want to compute **conditional probabilities** in Markov chains.
 - We can ask “what lead to $x_{10} = 4$?” with queries like $p(x_1 \mid x_{10} = 4)$.
 - We can ask “where does $x_{10} = 4$ lead?” with queries like $p(x_d \mid x_{10} = 4)$.
- **Monte Carlo approach** to estimating $p(x_j \mid x_{j'})$:
 - 1 Generate a large number of samples from the Markov chain, $x^i \sim p(x_1, x_2, \dots, x_d)$.
 - 2 Use Monte Carlo estimates of $p(x_j, x_{j'} = c)$ and $p(x_{j'} = c)$ to give

$$p(x_j = c \mid x_{j'} = c') = \frac{p(x_j = c, x_{j'} = c')}{p(x_{j'} = c')} \approx \frac{\sum_{i=1}^n I[x_j^i = c, x_{j'}^i = c']}{\sum_{i=1}^n I[x_{j'}^i = c']},$$

frequency of first event in samples consistent with second event.

- This is a special case of **rejection sampling** (we'll see general case later).
 - Unfortunately, if $x_{j'} = c'$ is rare then **most samples are “rejected”** (ignored).

Summary

- **Inverse Transform** generates samples from simple 1D distributions.
 - When we can easily invert the CDF.
- **Ancestral sampling** generates samples from multivariate distributions.
 - When conditionals have a nice form.
- **Monte Carlo** methods approximate expectations using samples.
 - Can be used to approximate arbitrary probabilities in Markov chains.
- Next time: the original Google algorithm.

Convergence Rate of Monte Carlo

- Consider case of using Monte Carlo method to estimate mean $\mu = \mathbb{E}[x]$,

$$\mu \approx \frac{1}{n} \sum_{i=1}^n x^i.$$

- We can write this as minimizing the 1-strongly convex

$$f(w) = \frac{1}{2} \|w - \mu\|^2.$$

- The gradient is $\nabla f(w) = (w - \mu)$.
- Consider stochastic gradient using

$$\nabla f_i(w^k) = w^k - x^{k+1},$$

which is unbiased since each x^i is unbiased μ approximation.

- Monte Carlo method is a stochastic gradient method with this approximation.

Convergence Rate of Monte Carlo

- Monte Carlo approximation as a stochastic gradient method with $\alpha_i = 1/(i + 1)$,

$$\begin{aligned}w^n &= w^{n-1} - \alpha_{n-1}(w^{n-1} - x^i) \\&= (1 - \alpha_{n-1})w^{n-1} + \alpha_{n-1}x^i \\&= \frac{n-1}{n}w^{n-1} + \frac{1}{n}x^i \\&= \frac{n-1}{n} \left(\frac{n-2}{n-1}w^{n-2} + \frac{1}{n-1}x^{i-1} \right) + \frac{1}{n}x^i \\&= \frac{n-2}{n}w^{n-2} + \frac{1}{n}(x^{i-1} + x^i) \\&= \frac{n-3}{n}w^{n-3} + \frac{1}{n}(x^{i-2} + x^{i-1} + x^i) \\&= \frac{1}{n} \sum_{i=1}^n x^i.\end{aligned}$$

- We know the rate of stochastic gradient for strongly-convex is $O(1/n)$.