

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

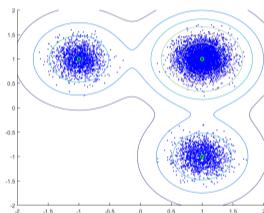
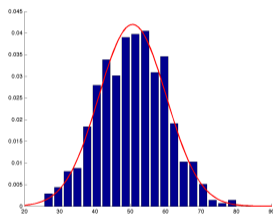
Fundamental Problem: Sampling from a Density

- A common inference task is **sampling from a density**.
 - **Generating examples x^i that are distributed according to a given density $p(x)$.**
 - Basically, the “opposite” of density estimation: **going from a model to data.**

$$p(x) = \begin{cases} 1 & \text{w.p. } 0.5 \\ 2 & \text{w.p. } 0.25 \\ 3 & \text{w.p. } 0.25 \end{cases} \Rightarrow X = \begin{bmatrix} 1 \\ 2 \\ 1 \\ 1 \\ 3 \\ 2 \\ 1 \\ 3 \end{bmatrix} .$$

Fundamental Problem: Sampling from a Density

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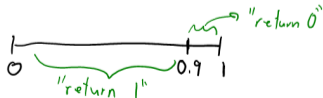
- We've been using pictures of samples to “tell us what the model has learned”.
 - If the samples look like real data, then we have a good density model.
- Samples can also be used in **Monte Carlo** estimation (today):
 - **Replace complicated $p(x)$ with samples** to solve hard problems at test time.

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:
 - Generate a uniform sample $u \sim \mathcal{U}(0, 1)$.
 - If $u \leq 0.9$, set $x = 1$ (otherwise, set $x = 0$).



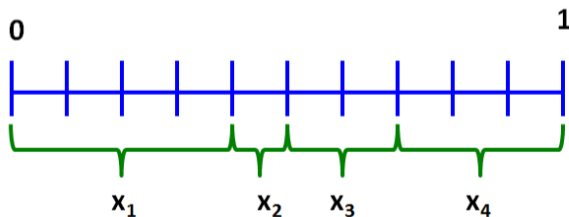
- If uniform samples are “good enough”, then we have $x = 1$ with probability 0.9.

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

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

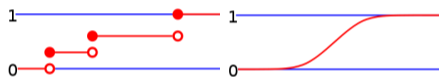
- To **sample from this categorical** density we can use (*sampleDiscrete* function):
 - 1 Generate $u \sim \mathcal{U}(0, 1)$.
 - 2 If $u \leq 0.4$, output 1.
 - 3 If $u \leq 0.4 + 0.1$, output 2.
 - 4 If $u \leq 0.4 + 0.1 + 0.2$, output 3.
 - 5 Otherwise, output 4.

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 $O(k + t \log k)$ instead of $O(tk)$:
 - 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 $u \leq p(x \leq c)$.

Cumulative Distribution Function (CDF)

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



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

- The “binary search for smallest c ” method finds **smallest c such that $u \leq F(c)$** .
 - This same approach works for continuous and general densities.
- General approach uses the **inverse CDF** (or “quantile”) function:
 - $F^{-1}(u) = \inf\{c \mid F(c) \geq u\}$.
 - Given a number u between 0 and 1, returns smallest c with $p(x \leq c) = u$.
 - If F is invertible, then F^{-1} is the usual inverse.

Inverse Transform Method (Exact 1D Sampling)

- Inverse transform method for exact sampling in 1D:

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

- Why this works (invertible case);

$$\begin{aligned} p(F^{-1}(u) \leq c) &= p(u \leq F(c)) && \text{(apply monotonic } F \text{ to both sides)} \\ &= F(c) && \text{(since } p(u \leq y) = y \text{ for uniform } u) \end{aligned}$$

- So this algorithm has the same CDF as the distribution we want to sample.
- Video on pseudo-random numbers and inverse-transform sampling:
 - <https://www.youtube.com/watch?v=C82JyCmtKWg>

Example: Sampling from a 1D Gaussian

- Consider a Gaussian distribution,

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

- CDF has the form

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

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

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).

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 .

Ancestral Sampling

- To **sample dependent** random variables we can use the **chain rule of probability**,

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

- The chain rule suggests the following sampling strategy:
 - Sample x_1 from $p(x_1)$.
 - Given x_1 , sample x_2 from $p(x_2 | x_1)$.
 - Given x_1 and x_2 , sample x_3 from $p(x_3 | x_2, x_1)$.
 - ...
 - Given x_1 through x_{d-1} , sample x_d from $p(x_d | x_{d-1}, x_{d-2}, \dots, x_1)$.
- This is called **ancestral sampling**.
 - It's easy if (conditional) probabilities are simple, since sampling in 1D is usually easy.
 - But may not be simple, binary **conditional j has 2^j values** of $\{x_1, x_2, \dots, x_j\}$.

Ancestral Sampling Examples

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

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

- So **ancestral sampling simplifies** too:

- ① Sample x_1 from initial probabilities $p(x_1)$.
- ② Given x_1 , sample x_2 from transition probabilities $p(x_2 | x_1)$.
- ③ Given x_2 , sample x_3 from transition probabilities $p(x_3 | x_2)$.
- ④ ...
- ⑤ Given x_{d-1} , sample x_d from transition probabilities $p(x_d | x_{d-1})$.

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

$$p(x, z) = p(z)p(x | z),$$

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

- If you want samples of x , sample (x, z) pairs and **ignore the z values**.

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 row to column):

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

- So $p(x_t = \text{“Grad School”} \mid x_{t-1} = \text{“Industry”}) = 0.01$.

Example of Sampling x_1

- Initial probabilities are:
 - 0.1 (Video Games)
 - 0.6 (Industry)
 - 0.3 (Grad School)
 - 0 (Video Games with PhD)
 - 0 (Academia)
 - 0 (Deceased)
- So initial CDF is:
 - 0.1 (Video Games)
 - 0.7 (Industry)
 - 1 (Grad School)
 - 1 (Video Games with PhD)
 - 1 (Academia)
 - 1 (Deceased)
- To sample the initial state x_1 :
 - First generate a uniform number u , for example $u = 0.724$.
 - Now find the first CDF value bigger than u , which in this case is "Grad School".

Example of Sampling x_2 , Given $x_1 = \text{"Grad School"}$

- So we sampled $x_1 = \text{"Grad School"}$.
 - To sample x_2 , we'll use the **"Grad School"** row in transition probabilities:

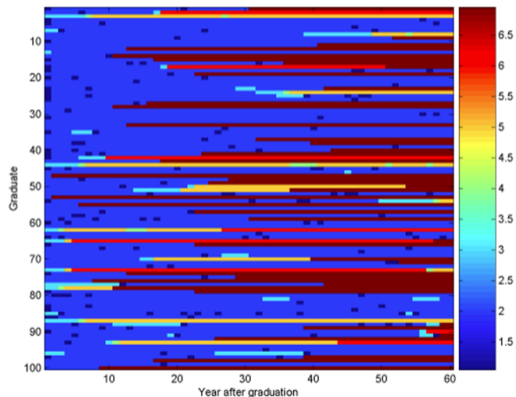
From\to	Video Games	Industry	Grad School	Video Games (with PhD)	Industry (with PhD)	Academia	Deceased
Video Games	0.08	0.90	0.01	0	0	0	0.01
Industry	0.03	0.95	0.01	0	0	0	0.01
Grad School	0.06	0.06	0.75	0.05	0.05	0.02	0.01
Video Games (with PhD)	0	0	0	0.30	0.60	0.09	0.01
Industry (with PhD)	0	0	0	0.02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

Example of Sampling x_2 , Given $x_1 = \text{"Grad School"}$

- Transition probabilities:
 - 0.06 (Video Games)
 - 0.06 (Industry)
 - 0.75 (Grad School)
 - 0.05 (Video Games with PhD)
 - 0.02 (Academia)
 - 0.01 (Deceased)
- So transition CDF is:
 - 0.06 (Video Games)
 - 0.12 (Industry)
 - 0.87 (Grad School)
 - 0.97 (Video Games with PhD)
 - 0.99 (Academia)
 - 1 (Deceased)
- To sample the second state x_2 :
 - First generate a uniform number u , for example $u = 0.113$.
 - Now find the first CDF value bigger than u , which in this case is "Industry".

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**

Marginalization and Conditioning

- 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 directly model marginals $p(x_j)$, and conditionals are marginals: $p(x_j | x_{j'}) = p(x_j)$.

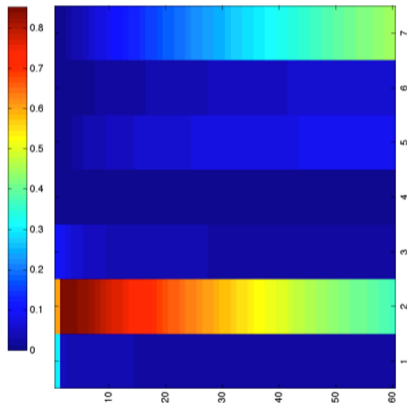
- This is also easy for mixtures of simple independent models.
 - Do inference for each mixture, add results using mixture probabilities:

$$p(x_j) = \sum_z p(z, x_j) = \sum_z p(z) \underbrace{p(x_j | z)}_{\text{inference within cluster}}$$

- 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 state and each column c is a year.

Monte Carlo: Marginalization 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 :(
 - Run physics simulator to “sample”, then see if it leads to a chain reaction.

Monte Carlo Method for Rolling Di

- Monte Carlo estimate of the probability of an event A :

$$\frac{\text{number of samples where } A \text{ happened}}{\text{number of samples}}.$$

- Computing probability of a pair of dice rolling a sum of 7:
 - Roll two dice, check if the sum is 7.
 - Roll two dice, check if the sum is 7.
 - Roll two dice, check if the sum is 7.
 - Roll two dice, check if the sum is 7.
 - Roll two dice, check if the sum is 7.
 - ...
- Monte Carlo estimate: fraction of samples where sum is 7.

Monte Carlo Method for Inequalities

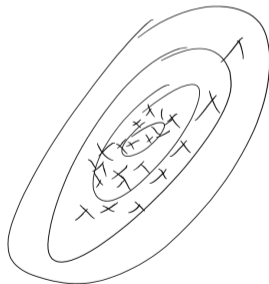
- Monte Carlo estimate of **probability that variable is above threshold**:
 - Compute fraction of examples where sample is above threshold.



Monte Carlo Method for Mean

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

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



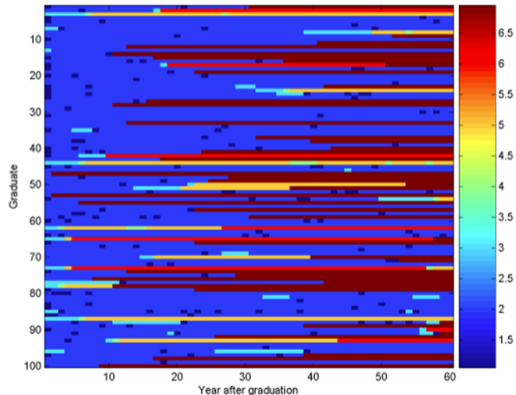
- Visual demo of Monte Carlo approximation of mean and variance:
 - <http://students.brown.edu/seeing-theory/basic-probability/index.html>

Monte Carlo for Non-Uniform Coupon Collecting

- Recall the **coupon collecting** problem:
 - You have a set of n possible objects.
 - **How many IID samples** do you need before you see all n ?
- With uniform probabilities, expected value is $O(n \log n)$.
- With non-uniform probabilities, you can approximate it with Monte Carlo:
 - Take IID samples until you have seen all objects.
 - Repeat many times and take the average time.
 - **Don't even need to necessarily know the probabilities.**

Monte Carlo for Markov Chains

- Our samples from the CS grad student Markov chain:



- We can estimate probabilities by looking at frequencies in samples.
 - In how many out of the 100 chains did we have $x_{10} = \text{“industry”}$?
- This works for continuous states too (for inequalities and expectations).

Monte Carlo Methods for Markov Chains

- Some Monte Carlo approximations of inference tasks 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 Methods

- Monte Carlo methods approximate expectations of random functions,

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

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

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

Unbiasedness of Monte Carlo Methods

- Let $\mu = \mathbb{E}[g(x)]$ be the value we want to approximate (not necessarily mean).
- The Monte Carlo estimate is an **unbiased** approximation of μ ,

$$\begin{aligned}\mathbb{E}\left[\frac{1}{n}\sum_{i=1}^n g(x^i)\right] &= \frac{1}{n}\mathbb{E}\left[\sum_{i=1}^n g(x^i)\right] && \text{(linearity of } \mathbb{E}\text{)} \\ &= \frac{1}{n}\sum_{i=1}^n \mathbb{E}[g(x^i)] && \text{(linearity of } \mathbb{E}\text{)} \\ &= \frac{1}{n}\sum_{i=1}^n \mu && (x^i \text{ is IID with mean } \mu) \\ &= \mu.\end{aligned}$$

- 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.

Rate of Convergence of Monte Carlo Methods

- Let f be the squared error in a 1D Monte Carlo approximation,

$$f(x^1, x^2, \dots, x^n) = \left(\frac{1}{n} \sum_{i=1}^n g(x^i) - \mu \right)^2.$$

- If variance is bounded, error with n samples is $O(1/n)$,

$$\begin{aligned} \mathbb{E} \left[\left(\frac{1}{n} \sum_{i=1}^n g(x^i) - \mu \right)^2 \right] &= \text{Var} \left[\frac{1}{n} \sum_{i=1}^n g(x^i) \right] && \text{(unbiased and def'n of variance)} \\ &= \frac{1}{n^2} \text{Var} \left[\sum_{i=1}^n g(x^i) \right] && (\text{Var}(\alpha x) = \alpha^2 \text{Var}(x)) \\ &= \frac{1}{n^2} \sum_{i=1}^n \text{Var}[g(x^i)] && \text{(IID)} \\ &= \frac{1}{n^2} \sum_{i=1}^n \sigma^2 = \frac{\sigma^2}{n}. && (x^i \text{ is IID with var } \sigma^2) \end{aligned}$$

- Similar $O(1/n)$ argument holds for $d > 1$ (notice that faster for small σ^2).

Conditional Probabilities with Monte Carlo

- We often want to compute **conditional probabilities** in Markov chains.
 - We can ask “what lead to $x_{10} = 4$?” with queries like $p(x_1 | x_{10} = 4)$.
 - We can ask “where does $x_{10} = 4$ lead?” with queries like $p(x_d | x_{10} = 4)$.
- **Monte Carlo approach** to estimating $p(x_j | x_{j'})$:
 - 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 $p(x_j = c, x_{j'} = c')$ and $p(x_{j'} = c')$ to give

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

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.

Monte Carlo as a Stochastic Gradient Method

- 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 applying stochastic gradient descent on f 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.

Monte Carlo as a Stochastic Gradient Method

- 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 $O(1/n)$.

Law of the Unconscious Statistician

- We use these identities to define the expectation of a function g applied to a random variable x ,

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

- The transformation from expectation to sum/integral is known as the “law of the unconscious statistician”.
 - It’s usually taken as being true, but it’s proof is a bit of a pain.

Accelerated Monte Carlo: Quasi Monte Carlo

- Unlike stochastic gradient, there are some “accelerated” Monte Carlo methods.
- **Quasi Monte Carlo** methods achieve an accelerated rate of $O(1/n^2)$.
 - Key idea: fill the space strategically with a deterministic “low-discrepancy sequence”.
 - Uniform random vs. deterministic low-discrepancy:

