COMPSCI 614: Randomized Algorithms with Applications to Data Science

Prof. Cameron Musco University of Massachusetts Amherst. Spring 2024. Lecture 22

Logistics

- Optional Problem Set 5 due 5/13 at 11:59pm.
- Final exam will be Tuesday 5/14, 10:30-12:30pm in the classroom. Study materials to be posted soon.
- Final project due the last day of finals: Friday 5/17.

Summary

Last Time:

- Finish up coupling. Example applications to shuffling, random walks on hypercubes, and exponential convergence of TV distance.
- Markov Chain Monte Carlo example of sampling random independent sets.
- Start on Metropolis Hastings algorithms and application to sampling from the hardcore model.

Today:

- · Finish the Metropolis Hastings algorithm.
- · Sampling to counting reduction for independent sets.

Mixing Time and Eigenvalues

A Markov chain is reversible if $\pi(i)P_{ij} = \pi(j)P_{ji}$ for all i,j. I.e., if the probability of transitioning from state i to state j is equal to the probability of transitioning from state j to state i in the steady state distribution. 'Detailed balance' condition.

- If the chain is irreducible and reversible, P has all real eigenvalues, $1 = \lambda_1 > \lambda_2 \dots > \lambda_n$.
- The eigenvalue gap is $\gamma = \lambda_1 \max\{|\lambda_2|, |\lambda_n|\}$.
- The mixing time is equal to $\tau(\epsilon) = \tilde{O}(\frac{1}{\gamma})$.

Mixing Time and Eigenvalues

Claim: If a Markov chain is reversible (i.e., $\pi(i)P_{ij} = \pi(j)P_{ji}$ for all i, j), then P has all real eigenvalues.

Proof:

- Let $D = diag(\pi)$. Then $D^{-1/2}PD^{1/2}$ is symmetric (and thus has real eigenvalues)
- The above is a similarity transform. The eigenvalues of P are identical to the eigenvalues of $D^{-1/2}PD^{1/2}$ and are thus real.

MCMC Methods Continued

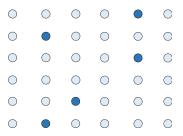
Achieving a Non-Uniform Stationary Distribution

Suppose we want to sample an independent set *X* from our graph with probability:

$$\pi(X) = \frac{\lambda^{|X|}}{\sum_{Y \text{ independent }} \lambda^{|Y|}},$$

for some 'fugacity' parameter $\lambda > 0$.

Known as the 'hard-core model' in statistical physics.

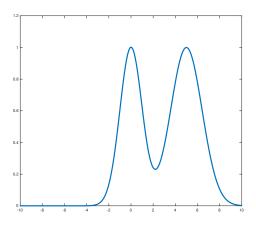


Metropolis-Hastings Algorithm

A very generic way of designing a Markov chain over state space [m] with stationary distribution $\pi \in [0,1]^m$.

- Assume the ability to efficiently compute a density $p(X) \propto \pi(X)$.
- Assume access to some symmetric transition function with transition probability matrix $Q \in [0, 1]^{m \times m}$.
- At step t, generate a 'candidate' state X_{t+1} from X_t according to Q.
- With probability min $\left(1, \frac{\rho(X_{t+1})}{\rho(X_t)}\right)$, 'accept' the candidate. Else 'reject' the candidate, setting $X_{t+1} = X_t$.

Metropolis-Hastings Intuition



Metropolis-Hastings Analysis

Need to check that for the Metropolis-Hastings algorithm, $\pi P = \pi$.

Suffices to show that pP = p where $p(i) \propto \pi(i)$ is our efficiently computable density.

$$[pP](i) = \underbrace{\sum_{j} p(j) \cdot Q_{j,i} \cdot \min\left(1, \frac{p(i)}{p(j)}\right)}_{aceptances} + \underbrace{p(i) \cdot \sum_{j} Q_{i,j} \left(1 - \min\left(1, \frac{p(j)}{p(i)}\right)\right)}_{rejections}$$

$$= \underbrace{\sum_{j} Q_{i,j} \cdot \min\left(p(j), p(i)\right) + p(i) \cdot \sum_{j} Q_{i,j} - \sum_{j} Q_{i,j} \cdot \min(p(i), p(j))}_{p(i)}$$

$$= p(i) \cdot \sum_{j} Q_{i,j} = p(i).$$

Metropolis-Hastings for the Hard-Core Model

Want to sample an independent set X with probability $\pi(X) = \frac{\lambda^{|X|}}{\sum_{X \text{ independent }} \lambda^{|Y|}}.$

- Let $p(X) = \lambda^{|X|}$ and let the transition function Q be given by:
 - · Pick a random vertex v.
 - If $v \in X_t$, set $X_{t+1} = X_t \setminus \{v\}$ with probability min(1, 1/ λ).
 - If $v \notin X_t$ and $X_t \cup \{v\}$ is independent, set $X_{t+1} = X_t \cup \{v\}$.
 - Else set $X_{t+1} = X_t$ with probability min(1, λ).
- Need to accept the transition with probability min $\left(1, \frac{p(X_{t+1})}{p(X_t)}\right)$.

The key challenge then becomes to analyze the mixing time.

For the related Glauber dynamics, Luby and Vigoda showed that for graphs with maximum degree Δ , when $\lambda < \frac{2}{\Delta - 2}$, the mixing time is $O(n \log n)$. But when $\lambda > \frac{c}{\Delta}$ for large enough constant c, it is NP-hard to approximately sample from the hard-core model.

MCMC for Approximate Counting

Counting to Sampling Reductions

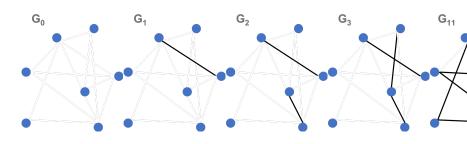
Often if one can efficiently sample from the distribution $\pi(X) = \frac{p(X)}{\sum_{Y} p(Y)}$, one can efficiently approximate the normalizing constant $Z = \sum_{Y} p(Y)$ (often called the partition function).

- If Z is hard to approximate, then this can give a proof that sampling is hard, and thus it is unlikely that any simple MCMC method for sampling from π mixes rapidly.
- This is e.g., how one can show that sampling from the hard-core model is hard when $\lambda = \Omega(1/\Delta)$.
- Let's consider the simple case of $\lambda=1$. I.e., we want to sample a uniformly random independent set.
- In this case, Z = |S(G)|, the number of independent sets in G. It is known that approximating |S(G)| even up to a poly(n) factor is NP-Hard.

Counting Independent Sets

How can we count the number of independent sets |S(G)| in a graph, given an oracle for sampling a uniform random independent set?

Let G_0, G_1, \ldots, G_m be a sequence of graphs with $G_m = G$ and G_i obtained by removing an arbitrary edge from G_{i+1} .



We can write:

$$|S(G)| = \frac{|S(G_m)|}{|S(G_{m-1})|} \cdot \frac{|S(G_{m-1})|}{|S(G_{m-2})|} \cdot \ldots \cdot \frac{|S(G_1)|}{|S(G_0)|} \cdot |S(G_0)|.$$

Counting Independent Sets

$$|S(G)| = \frac{|S(G_m)|}{|S(G_{m-1})|} \cdot \frac{|S(G_{m-1})|}{|S(G_{m-2})|} \cdot \ldots \cdot \frac{|S(G_1)|}{|S(G_0)|} \cdot |S(G_0)| 2^n = 2^n \cdot \prod_{i=1}^m r_i,$$

where $r_i = \frac{|S(G_m)|}{|S(G_{m-i})|}$. If we can estimate each r_i with \tilde{r}_i satisfying

$$\left(1 - \frac{\epsilon}{2m}\right) \cdot r_i \le \tilde{r}_i \le \left(1 + \frac{\epsilon}{2m}\right) \cdot r_i,$$

then:

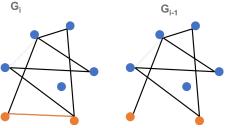
$$(1 - \epsilon) \cdot |S(G)| \le 2^n \cdot \prod_{i=1}^m \tilde{r}_i \le (1 + \epsilon) \cdot |S(G)|$$

since
$$\left(1 + \frac{\epsilon}{2m}\right)^m \le 1 + \epsilon$$
 and $\left(1 - \frac{\epsilon}{2m}\right)^m \ge 1 - \epsilon$.

Independent Set Ratios

Consider the ratio $r_i = \frac{|S(G_i)|}{|S(G_{i-1})|}$. Observe that $r_i \leq 1$.

Further, $r_i \ge 1/2$. Let (u, v) be the edge removed from G_i to obtain G_{i-1} . Then each independent set in $S(G_{i-1}) \setminus S(G_i)$, must contain both u and v.



So, we can map each set in $S(G_{i-1}) \setminus S(G_i)$ to a unique set in $S(G_i)$ by simply removing v.

$$r_i = \frac{|S(G_i)|}{|S(G_{i-1})|} = \frac{|S(G_i)|}{|S(G_i)| + |S(G_{i-1}) \setminus S(G_i)|} \ge \frac{1}{2}.$$

Independent Set Ratios

So Far: We have written $|S(G)| = 2^n \cdot \prod_{i=1}^m r_i$ where $r_i = \frac{|S(G_i)|}{|S(G_{i-1})|}$. Need to get a $1 \pm \epsilon/m$ estimate to each r_i to get a $1 \pm \epsilon$ estimate to |S(G)|.

Let **X** be a random variable generated as follows: pick a random independent set from G_{i-1} and let X = 1 if the set is also independent in G_i . Otherwise let X = 0.

What is $\mathbb{E}[X]$?

How many samples of **X** do we need to take to obtain a $1 \pm \epsilon/m$ approximation to r_i with high probability?

Counting Independent Sets

Upshot: For a graph G with m edges, making $\tilde{O}(m^2/\epsilon^2)$ calls to a uniform random independent set sampler on G or its subgraphs suffices to approximate the number of independent sets in G up to $1 \pm \epsilon$ relative error.

- So a polynomial time algorithm for uniform random independent set sampling, would lead to a polynomial time algorithm for counting independent sets, and hence the collapse of NP to P.
- Observe that near-uniform sampling (as would be obtained e.g., with an MCMC method) would also suffice.