# **Sampling Methods**

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Rejection Sampling Importance Sampling

- 1. Markov Chains
- 2. Importance Sampling
- 3. Gibbs Sampling
- 4. Markov Chain Monte Carlo

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- Let q(x) be a proposal distribution from which we can sample.
- Let *M* be a constant such that  $M \ge \frac{\tilde{p(x)}}{q(x)}$  for all *x*.
- Then, we can sample from p(x) by sampling from q(x) and accepting the sample with probability  $\frac{p(x)}{Mq(x)}$ .

Notebook: rejection-sampling.ipynb







#### 7



#### 8









### **Rejection Sampling (Rejected Sample)**



#### **Rejection Sampling (Accepted Sample)**



# Rejection Sampling (10 samples)



# Rejection Sampling (10 samples) (KDE)



### **Rejection Sampling (1000 samples)**



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### **Rejection Sampling (10000 samples)**



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- Acceptance Probability α(x): Probability that we accept a sample x<sub>s</sub> generated from q(x).

$$\alpha(x_s) = \frac{\tilde{p}(x_s)}{Mq(x_s)} = P(Accept|x_s)$$
(1)
$$P(x_{s}|Accept) = \frac{P(Accept|x_{s})P(x_{s})}{P(Accept)}$$
(2)

where P(x<sub>s</sub>|Accept) is the density of accepted sample x<sub>s</sub>. We want to evaluate this and show this is p(x<sub>s</sub>).

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$$P(Accept) = \frac{Z}{M}$$
(7)

where Z is the normalization constant of  $\tilde{p}(x)$ .

$$P(x_{s}|Accept) = \frac{P(Accept|x_{s})P(x_{s})}{P(Accept)}$$
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$$P(x_s|Accept) = p(x_s) \tag{12}$$

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- How to choose multiplier M?
- Match the densities at the peak of  $\tilde{p}(x)$  and q(x), i.e. at  $x = \vec{0}$ .

• 
$$\tilde{p}(x) = \frac{1}{(2\pi)^{D/2} \sigma_p^D} \exp(-\frac{1}{2\sigma_p^2} x^T x)$$

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• 
$$q(x) = \frac{1}{(2\pi)^{D/2}\sigma_q^D} \exp(-\frac{1}{2\sigma_q^2}x^Tx)$$

• At 
$$x = \vec{0}$$
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• 
$$M = \frac{\tilde{p}(x)}{q(x)} = \frac{\sigma_q^D}{\sigma_p^D} = (\frac{\sigma_q}{\sigma_p})^D$$

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• Let us assume 
$$\sigma_p = 1$$
 and  $\sigma_q = 1.1$ 





• Acceptance probability is very low as D increases.

- Rejection sampling is inefficient when the target distribution is very different from the proposal distribution. In this case, we will reject a lot of samples.
- This is a problem when sampling from high-dimensional distributions. Acceptance probability α(x) is very low.

- We want to compute posterior predictive distribution (or something similar)
- We would typically use Monte Carlo methods to do this.
- $I = \int f(x)p(x)dx$  where p(x) is the posterior distribution.
- We can approximate I by  $\frac{1}{N} \sum_{i=1}^{N} f(x_i)$ , where  $x_i \sim p(x)$ .
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- But, we do not have access to p(x). Rather, we have access to p̃(x), which is proportional to p(x).
- In rejection sampling, we took a sample  $x_i$  from q(x) and accepted it with probability  $\frac{\tilde{p}(x_i)}{Mq(x_i)}$ .
- Can we use all samples  $x_i$  from q(x) without rejection?

- In rejection sampling, we took a sample  $x_i$  from q(x) and accepted it with probability  $\frac{\tilde{p}(x_i)}{Mq(x_i)}$ .
- Can we use all samples  $x_i$  from q(x) without rejection?
- $I = \int f(x)p(x)dx \approx \frac{1}{N}\sum_{i=1}^{N} f(x_i)$ , where  $x_i \sim p(x)$ .
- Let us choose a proposal distribution q(x) which has support over the entire domain of p(x).

• 
$$I = \int f(x)p(x)dx = \int f(x)\frac{p(x)}{q(x)}q(x)dx$$

•  $I = \int f(x)w(x)q(x)dx$ , where  $w(x) = \frac{p(x)}{q(x)}$ . w(x) is called the importance weight.

• 
$$I = \mathbb{E}_q[f(x)w(x)] = \sum_{i=1}^N f(x_i)w(x_i)$$
, where  $x_i \sim q(x)$ .

## Importance Sampling (with unnormlized $\tilde{p}(x)$ instead of p(x))

$$I = \int f(x)p(x)dx \approx \frac{1}{Z}\frac{1}{S}\sum_{s}f(x_{S})\frac{\tilde{p}(x_{S})}{q(x_{S})}$$
(13)

Now, we need to estimate Z.

$$Z = \int \tilde{\rho}(x) dx = \int \frac{\tilde{\rho}(x)}{q(x)} q(x) dx$$
  
=  $\mathbb{E}_q \left[ \frac{\tilde{\rho}(x)}{q(x)} \right] = \frac{1}{5} \sum_s \frac{\tilde{\rho}(x_s)}{q(x_s)}$  (14)

Thus, we can write I as:

$$I \approx \frac{1}{S} \sum_{s} f(x_{s}) \frac{\tilde{p}(x_{s})/q(x_{s})}{\frac{1}{s} \sum_{t} \tilde{p}(x_{t})/q(x_{t})} =: \sum_{s} f(x_{s}) \tilde{w}_{s} \qquad (15)$$

# **Markov Chains**

https://nipunbatra.github.io/hmm/

Notebook: mcmc=optimization.ipynb

# **Importance Sampling**

In rejection sampling, we saw that due to less acceptance probability, a lot of samples were wasted leading to more time and higher complexity to approximate a distribution.

Computing p(x), q(x) thus seems wasteful. Let us rewrite the equation as:

$$\phi = \int f(x)p(x)dx = \int f(x)\frac{p(x)}{q(x)}q(x)dx$$
$$\sim \frac{1}{N}\sum_{i=1}^{N}f(x_i)\frac{p(x_i)}{q(x_i)} = \frac{1}{N}\sum_{i=1}^{N}f(x_i)w_i$$

Here,  $x_i \sim q(x)$ .  $w_i$  is known as the importance(weight) of sample i.

However the normalization constant  ${\cal Z}$  is generally not known to us. Thus writing:

$$p(x) = \frac{\tilde{p}(x)}{Z} \tag{16}$$

Now inserting this in earlier equations, we get:

$$\phi = \frac{1}{Z} \int f(x)\tilde{p}(x)dx = \frac{1}{Z} \int f(x)\frac{\tilde{p}(x)}{q(x)}q(x)dx$$
$$\sim \frac{1}{NZ} \sum_{i=1}^{N} f(x_i)\frac{\tilde{p}(x_i)}{q(x_i)} = \frac{1}{NZ} \sum_{i=1}^{N} f(x_i)w_i$$

We know that:

$$Z = \int_{\infty}^{\infty} \tilde{p}(x) dx = \int_{\infty}^{\infty} \frac{\tilde{p}(x)}{q(x)} q(x) dx$$
$$= \frac{1}{N} \sum_{i=1}^{N} w_i$$
Substuting this value of Z in the equation above, we get:

$$\phi = \frac{1}{N} \sum_{i=1}^{N} f(x_i) w_i = \frac{\sum_{i=1}^{N} f(x_i) w_i}{\sum_{i=1}^{N} w_i}$$
$$= \sum_{i=1}^{N} f(x_i) W_i$$

Here  $W_i = \frac{w_i}{\sum_{i=1}^N w_i}$  are the normalized weights.

### Limitations

Recall that Var φ̂ = var(f)/N. Importance sampling replaces var(f) with var(f p/q). At positions where p >>> q, the weight can tend to ∞!



# **Gibbs Sampling**

Suppose we wish to sample  $\theta_1, \theta_2 \sim p(\theta_1, \theta_2)$ , but cannot use:

- direct simulation
- accept-reject method
- Metropolis-Hasting

But we can sample using the conditionals i.e.:

- $p(\theta_1|\theta_2)$  and
- $p(\theta_2|\theta_1)$ ,

then we can use Gibbs sampling.

Suppose  $\theta_1, \theta_2 \sim p(\theta_1, \theta_2)$  and we can sample from  $p(\theta_1, \theta_2)$ . We begin with an initial value  $(\theta_1^0, \theta_2^0)$ , the workflow for Gibbs algorithm is:

- 1. sample  $heta_1^j \sim p( heta_1| heta_2^{j-1})$  and then
- 2. sample  $\theta_2^j \sim p(\theta_2|\theta_1^j)$ .

One thing to note here is that the sequence in which the theta's are sampled are not independent!

### **Bivariate Normal Example**

Suppose

$$heta \sim \mathit{N}_2(0,\Sigma) ext{ and } \Sigma = egin{matrix} 1 & 
ho \ 
ho & 1 \end{pmatrix}$$

Then, we have:

$$\theta_1|\theta_2 \sim N(\rho\theta_2, [1-\rho^2])$$

 $\theta_2|\theta_1 \sim N(\rho\theta_1, [1-\rho^2])$  are the conditional distributions. The Gibbs sampling proceeds as follows:

IterationSample 
$$\theta_1$$
Sample  $\theta_2$ 1 $\theta_1 \sim N(\rho \theta_2^0, [1-\rho^2])$  $\theta_2 \sim N(\rho \theta_1^1, [1-\rho^2])$ 

$$k \qquad \theta_1 \sim N(\rho \theta_2^{k-1}, [1-\rho^2]) \quad \theta_2 \sim N(\rho \theta_1^k, [1-\rho^2])$$



Suppose 
$$\theta = (\theta_1, \theta_2, \dots, \theta_K)$$
, the Gibbs workflow is as follows:  
 $\theta_1^j = p(\theta_1 | \theta_2^{j-1}, \dots, \theta_K^{j-1})$   
 $\theta_2^j = p(\theta_2 | \theta_1^j, \theta_3^{j-1}, \dots, \theta_K^{j-1})$   
.  
 $\theta_k^j = p(\theta_k | \theta_1^j, \dots, \theta_{k-1}^j, \theta_{k+1}^{j-1}, \dots, \theta_K^{j-1})$   
.  
 $\theta_K^j = p(\theta_K | \theta_1^j, \dots, \theta_{K-1}^j)$   
The distributions above are call the full conditional distributions.

Gibbs sampling can be used to draw samples from  $p(\theta)$  when:

- Other methods don't work quite well in higher dimensions.
- Draw samples from the full conditional distributions is easy,  $p(\theta_k | \theta_{-k})$ .

## Markov Chain Monte Carlo

- *Transformation based methods*: Usually limited to drawing from standard distributions.
- *Rejection and Importance sampling*: Require selection of good proposal distirbutions.

In high dimensions, usually most of the density p(x) is concentrated within a tiny subspace of x. Moreover, those subspaces are difficult to be known a priori.

A solution to these are MCMC methods.

Markov Chain: A joint distribution p(X) over a sequence of random variables X = {X<sub>1</sub>, X<sub>2</sub>,..., X<sub>n</sub>} is said to have the Markov property if

$$p(X_i|X_1,...,X_{i-1}) = p(X_i|X_{i-1})$$

The sequence is then called a Markov chain.

• The idea is that the estimates contain information about the shape of the target distribution *p*.

- The basic idea is propose to move to a new state  $x_{i+1}$  from the current state  $x_i$  with probability  $q(x_{i+1}|x_i)$ , where q is called the proposal distribution and our target density of interest is  $p(=\frac{1}{Z}\tilde{p})$ .
- The new state is accepted with probability  $\alpha(x_i, x_{i+1})$ .
  - If  $p(x_{i+1}|x_i) = p(x_i|x_{i+1})$ , then  $\alpha(x_i, x_{i+1}) = \min(1, \frac{p(x_{i+1})}{p(x_i)})$ .
  - If  $p(x_{i+1}|x_i) \neq p(x_i|x_{i+1})$ , then  $\alpha(x_i, x_{i+1}) = \min(1, \frac{p(x_{i+1})q(x_i|x_{i+1})}{p(x_i)q(x_{i+1}|x_i)}) = \min(1, \frac{\tilde{p}(x_{i+1})q(x_i|x_{i+1})}{\tilde{p}(x_i)q(x_{i+1}|x_i)})$
- Evaluating α, we only need to know the target distribution up to a constant of proportionality or without normalization constant.

- 1. Initialize  $x_0$ .
- 2. for i = 1, ..., N do:
- 3. Sample  $x^* \sim q(x^*|x_{i-1})$ .
- 4. Compute  $\alpha = \min(1, \frac{\tilde{p}(x^*)q(x_{i-1}|x^*)}{\tilde{p}(x_{i-1})q(x^*|x_{i-1})})$
- 5. Sample  $u \sim \mathcal{U}(0, 1)$
- 6. if  $u \leq \alpha$ :

$$x_i = x^*$$

else:

$$x_i = x_{i-1}$$

#### How do we choose the initial state $x_0$ ?

How do we choose the initial state  $x_0$ ?

- 1. Start the Markov Chain at an initial  $x_0$ .
- 2. Using the proposal  $q(x|x_i)$ , run the chain long enough, say  $N_1$  steps.
- 3. Discard the first  $N_1 1$  samples (called 'burn-in' samples).
- 4. Treat  $x_{N_1}$  as first sample from p(x).

https://chi-feng.github.io/mcmc-demo/app.html