# Sampling Methods

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- Goal: sample from  $p(x)$ .

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- Let  $q(x)$  be a proposal distribution from which we can sample.
- Let M be a constant such that  $M \geq \frac{p(x)}{q(x)}$  $\frac{p(x)}{q(x)}$  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>

















## Rejection Sampling (Rejected Sample)



### Rejection Sampling (Accepted Sample)



## Rejection Sampling (10 samples)



## Rejection Sampling (10 samples) (KDE)



## Rejection Sampling (1000 samples)



## Rejection Sampling (1000 samples) (KDE)



### Rejection Sampling (10000 samples)



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- Aim:
	- Show that the samples we accept are distributed according to  $p(x)$ .
	- Or, the density of an accepted sample (say  $x_s$ ) is  $p(x_s)$  (and not  $\tilde{p}(x_s)$ ).
- Acceptance Probability  $\alpha(x)$ : Probability that we accept a sample  $x_s$  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)

 $\bullet$  where  $P(\mathsf{x}_s | \mathsf{Accept})$  is the density of accepted sample  $\mathsf{x}_s$ . We want to evaluate this and show this is  $p(x_s)$ .

$$
P(x_s|Accept) = \frac{P(Accept|x_s)P(x_s)}{P(Accept)} \tag{2}
$$

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- $P(x_s) = q(x)$  is the density of samples we draw from  $q(x)$ .

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- $P(Accept|x_s)$  is  $\alpha(x_s)$
- $P(x_s) = q(x)$  is the density of samples we draw from  $q(x)$ .
- $P(Accept)$  is the unconditional probability that we accept a sample generated from  $q(x)$ .

$$
P(Accept) = \int P(Accept |xs)P(xs)dxs
$$
 (3)

$$
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$$
 (3)

$$
P(Accept) = \int \alpha(x_s)q(x_s)dx_s \tag{4}
$$

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$$
 (3)

$$
P(Accept) = \int \alpha(x_s)q(x_s)dx_s \qquad (4)
$$

$$
P(Accept) = \int \frac{\tilde{p}(x_s)}{Mq(x_s)} q(x_s) dx_s
$$
 (5)

$$
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$$
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$$

$$
P(Accept) = \int \frac{\tilde{p}(x_s)}{Mq(x_s)} q(x_s) dx_s
$$
 (5)

$$
P(Accept) = \frac{1}{M} \int \tilde{p}(x_s) dx_s
$$
 (6)

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$$
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$$

$$
P(Accept) = \int \frac{\tilde{p}(x_s)}{Mq(x_s)} q(x_s) dx_s
$$
 (5)

$$
P(Accept) = \frac{1}{M} \int \tilde{p}(x_s) dx_s
$$
 (6)

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

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

$$
P(x_s|Accept) = \frac{P(Accept|x_s)P(x_s)}{P(Accept)}
$$
(8)

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(8)

$$
P(x_s|Accept) = \frac{\alpha(x_s)q(x_s)}{P(Accept)}
$$
(9)

$$
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$$
(8)

$$
P(x_s|Accept) = \frac{\alpha(x_s)q(x_s)}{P(Accept)}
$$
(9)

$$
P(x_s|Accept) = \frac{\frac{\tilde{p}(x_s)}{Mq(x_s)}q(x_s)}{\frac{Z}{M}}
$$
(10)

$$
P(x_s|Accept) = \frac{P(Accept|x_s)P(x_s)}{P(Accept)}
$$
(8)

$$
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$$
(9)

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$$
(10)

$$
P(x_s|Accept) = \frac{\tilde{p}(x_s)}{Z}
$$
 (11)

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P(x_s|Accept) = p(x_s)
$$
 (12)

• Let us assume  $\tilde{p}(x)$  is D dimensional Gaussian  $\mathcal{N}_D(0, \sigma_p^2 I)$ 

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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  $x = \vec{0}$ .

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

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• 
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• At 
$$
x = \vec{0}
$$
,  $\tilde{p}(x) = \frac{1}{(2\pi)^{D/2} \sigma_p^D}$  and  $q(x) = \frac{1}{(2\pi)^{D/2} \sigma_q^D}$ 

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• 
$$
M = \frac{\tilde{\rho}(x)}{q(x)} = \frac{\sigma_q^D}{\sigma_p^D} = (\frac{\sigma_q}{\sigma_p})^D
$$

• 
$$
M = \frac{\tilde{\rho}(x)}{q(x)} = \frac{\sigma_q^D}{\sigma_p^D} = \left(\frac{\sigma_q}{\sigma_p} \right)D
$$
  
\n• Let us assume  $\sigma_p = 1$  and  $\sigma_q = 1.1$   
\n
$$
\begin{bmatrix}\n0^q \\
\vdots \\
0^{10^q} \\
0^{10^q}\n\end{bmatrix}
$$
\n
$$
\begin{bmatrix}\n0^q \\
\vdots \\
0^{10^n}\n\end{bmatrix}
$$
\n
$$
\begin{bmatrix}\n0^q \\
\vdots \\
0^{10^n}\n\end{bmatrix}
$$
\n
$$
\begin{bmatrix}\n0 \\
\vdots \\
0\n\end{bmatrix}
$$



• 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  $\alpha(x)$  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  $x_i \sim p(x)$ .
- But, we do not have access to  $p(x)$ . Rather, we have access to  $\tilde{p}(x)$ , which is proportional to  $p(x)$ .
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- But, we do not have access to  $p(x)$ . Rather, we have access to  $\tilde{p}(x)$ , which is proportional to  $p(x)$ .
- $\bullet$  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?
- $\bullet$  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?
- $I = \int f(x)p(x)dx \approx \frac{1}{N}$  $\frac{1}{N}\sum_{i=1}^{N}f(x_i)$ , 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.

$$
Z = \int \tilde{p}(x) dx = \int \frac{\tilde{p}(x)}{q(x)} q(x) dx
$$
  
=  $\mathbb{E}_q \left[ \frac{\tilde{p}(x)}{q(x)} \right] = \frac{1}{S} \sum_{s} \frac{\tilde{p}(x_s)}{q(x_s)}$  (14)

Thus, we can write *I* as:

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

# <span id="page-66-0"></span>[Markov Chains](#page-66-0)

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

Notebook: mcmc=optimization.ipynb

# <span id="page-69-0"></span>[Importance Sampling](#page-69-0)

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  $Z$  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  $\hat{\phi} = \frac{var(f)}{N}$  $\frac{N(t)}{N}$ . Importance sampling replaces var(f) with var(f $\frac{p}{q}$  $\frac{p}{q}$ ). At positions where  $p >> q$ , the weight can tend to  $\infty!$ 



## <span id="page-74-0"></span>[Gibbs Sampling](#page-74-0)

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  $\theta_1^j \sim \rho(\theta_1|\theta_2^{j-1})$  $\binom{J-1}{2}$  and then
- 2. sample  $\theta_2^j \sim \rho(\theta_2 | \theta_1^j)$  $\binom{J}{1}$ .

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

### Bivariate Normal Example

Suppose

$$
\theta \sim N_2(0,\Sigma) \text{ and } \Sigma = \frac{1}{\rho} \frac{\rho}{1}
$$

Then, we have:

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

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

> . .



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



Suppose 
$$
\theta = (\theta_1, \theta_2, ..., \theta_K)
$$
, the Gibbs workflow is as follows:  
\n
$$
\begin{aligned}\n\theta_1^j &= p(\theta_1 | \theta_2^{j-1}, ..., \theta_K^{j-1}) \\
\theta_2^j &= p(\theta_2 | \theta_1^j, \theta_3^{j-1}, ..., \theta_K^{j-1}) \\
\vdots \\
\theta_K^j &= p(\theta_K | \theta_1^j, ..., \theta_{K-1}^j, \theta_{K+1}^{j-1}, ..., \theta_K^{j-1})\n\end{aligned}
$$
\n
$$
\begin{aligned}\n\theta_K^j &= p(\theta_K | \theta_1^j, ..., \theta_{K-1}^j) \\
\text{The distributions above are call the full conditional distributions.}\n\end{aligned}
$$

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,  $p(\theta_k | \theta_{-k}).$

# <span id="page-81-0"></span>[Markov Chain Monte Carlo](#page-81-0)

- 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  $X = \{X_1, X_2, \ldots, X_n\}$  is said to have the Markov property if

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

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  $\alpha(x_i, x_{i+1}) = \min(1, \frac{p(x_{i+1})}{p(x_i)})$  $\frac{p(x_{i+1})}{p(x_i)}$ .
	- If  $p(x_{i+1}|x_i) \neq p(x_i|x_{i+1})$ , then  $\alpha(x_i, x_{i+1}) = \min(1, \frac{p(x_{i+1})q(x_i|x_{i+1})}{p(x_i)q(x_{i+1}|x_i)})$  $\frac{p(\textsf{x}_{i+1})q(\textsf{x}_i|\textsf{x}_{i+1})}{p(\textsf{x}_i)q(\textsf{x}_{i+1}|\textsf{x}_i)}$ ) = min $(1,\frac{\tilde{p}(\textsf{x}_{i+1})q(\textsf{x}_i|\textsf{x}_{i+1})}{\tilde{p}(\textsf{x}_i)q(\textsf{x}_{i+1}|\textsf{x}_i)}$  $\frac{p(x_{i+1})q(x_i|x_{i+1})}{p(x_i)q(x_{i+1}|x_i)}$
- Evaluating  $\alpha$ , 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  $\alpha = \min(1, \frac{\tilde{p}(x^*)q(x_{i-1}|x^*)}{\tilde{p}(x_{i-1})q(x^*)x_{i-1}})$  $\frac{p(x)q(x_{i-1}|x)}{\tilde{p}(x_{i-1})q(x^*|x_{i-1})}$
- 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>