# Big Data Analysis (MA60306) 

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Lecture 23
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## Sampling methods

Proposition The probability that $X \sim g$ is accepted is $\frac{1}{c}$, and is maximized when $c=\sup _{x} \frac{f(x)}{g(x)}$

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Proof

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\begin{aligned}
P\left(U \leq \frac{f(X)}{c g(X)}\right) & =\int_{-\infty}^{\infty} \int_{0}^{\frac{f(t)}{c g(t)}} g(t) d u d t \\
& =\int_{-\infty}^{\infty} \frac{f(t)}{c g(t)} g(t) d t \\
& =\int_{-\infty}^{\infty} \frac{f(t)}{c} d t=\frac{1}{c}
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Efficiency Since $\sup _{x} \frac{f(x)}{g(x)}=\sqrt{\frac{2 e}{\pi}}$ is the acceptance rate, by the above result, it would be $\sqrt{\frac{\pi}{2 e}}=0.7602$. Thus if we generate $100 x$-values from $g$, we can expect 75 of them would be retained, and the others discarded.

## Sampling methods

Question What is the acceptance rate when the standard normal is sampled using Accept-Reject method through the standard Cauchy density

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g(x)=\frac{1}{\pi\left(1+x^{2}\right)}
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Gamma with parameters $n$ and $\lambda$ generate $X \sim G(n, \lambda)$, first generate $n$ independent values $X_{1}, \ldots, X_{n}$ from a standard exponential, and use $X=\lambda\left(X_{1}+\ldots+X_{n}\right)$

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Beta with parameters $m, n$ To generate $X \sim \operatorname{Be}(m, n)$, generate $U \sim G(m, 1), V \sim G(n, 1)$ independently, and use $X=\frac{U}{U+V}$

## Sampling methods

$t$ distribution with $n$ degrees of freedom To generate $X \sim t(n)$, first generate $Z_{1}, Z_{2}, \ldots, Z_{n+1} \sim \mathcal{N}(0,1)$ independently, and use

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Wishart with general parameters To generate $S \sim W_{p}(n, \Sigma)$, generate $X_{1}, X_{2}, \ldots, X_{n} \sim \mathcal{N}_{p}(0, \Sigma)$ independently, and use

$$
S=\sum_{i=1}^{n} X_{i} X_{i}^{T}
$$

Uniform on the surface of the unit ball Generate $Z_{1}, \ldots, Z_{d} \sim \mathcal{N}(0,1)$ independently, and use

$$
X_{i}=\frac{Z_{i}}{\sqrt{Z_{1}^{2}+\ldots+Z_{d}^{2}}}, i=1, \ldots, d
$$

## Sampling methods

Markov Chain Monte Carlo (MCMC)
$\rightarrow$ The standard simulation techniques are usually difficult to apply, for example, when the target distribution is an unconventional one, or even worse, it is known only up to a normalizing constant such as:

$$
f(x)=\frac{h(x)}{c}
$$

for some explicit function $h$, but only $c$ an implicit normalizing constant $c$ because it cannot be computed exactly, or even to a high degree of accuracy

## Sampling methods

$\rightarrow$ For example, the problem of simulating from posterior densities of a parameter(s)

$$
\pi(\theta \mid x)=\frac{f(x \mid \theta) \pi(\theta)}{m(x)}
$$

where $f(x \mid \theta)$ is the likelihood function, $\pi(\theta)$ is the prior density, and $m(x)$ is the marginal density density of the observable $X$ induced by $(f, \pi)$. Thus

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If the parameter $\theta$ is high-dimensional, and the prior density $\pi(\theta)$ is not a very conveniently chosen one, then $m(x)$ usually cannot be calculated in closed-form, or even to a high degree of numerical approximation.
All the simulation methods discussed in the previous section are useless in such a situation

## Sampling methods

Markov chain Monte Carlo
Graph: A pair of sets $G=(V, E), E \subseteq E \times E, V \neq \emptyset$. Two vertices $v_{i}, v_{j} \in V$ are adjacent if $\left(v_{i}, v_{j}\right) \in E$.
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Walk: A sequence of vertices and edges

$$
v_{1}, e_{1}, v_{2}, e_{2}, \ldots, v_{k}, e_{k}, v_{k+1}
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such that end points of $e_{i}$ are $v_{i}$ and $v_{i+1}, 1 \leq i \leq k$
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The problem of a random walk ${ }^{1}$ was posed by Louis Bachelier in his thesis devoted to the theory of financial speculations in 1900. The term random walks was first introduced by Karl Pearson in 1905.

[^0]
## Sampling methods

For each $t$, let $X_{t}$ denote the index of the vertex at which the walker resides. Hence $\left\{X_{0}, X_{1}, \ldots\right\}$ is a stochastic process (Markov Chain) taking values in $\{1,2,3,4\}$ stands at time $t$


At time 0 , the random walker stands at $v_{1}: p\left(X_{0}=1\right)=1$
At time 1, flips a fair coin and moves immediately to $v_{2}$ or $v_{4}$ according to whether the coin comes up heads or tails:
$p\left(X_{1}=v_{2}\right)=\frac{1}{2}=p\left(X_{1}=v_{4}\right)$.

## Sampling methods



Markov property

$$
\begin{aligned}
& p\left(X_{t+1}=v_{1} \mid X_{0}=i_{0}, X_{1}=i_{1}, \ldots, X_{t-1}=i_{t-1}, X_{t}=v_{2}\right)=\frac{1}{2} \\
& p\left(X_{t+1}=v_{3} \mid X_{0}=i_{0}, X_{1}=i_{1}, \ldots, X_{t-1}=i_{t-1}, X_{t}=v_{2}\right)=\frac{1}{2}
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for any choice of $i_{0}, \ldots, i_{t-1}$.

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Time homogeneity

$$
p\left(X_{t+1} \mid X_{t}=v\right)=c
$$

for all $t$, for any $v \in V$

## Sampling methods



Markov chain A sequence of random variables $\left\{X_{n}\right\}, n \geq 0$, is said to be a Markov chain if for some countable set $S \subset \mathbb{R}$, and any $n \geq 1$, $s_{n+1}, s_{n}, \ldots, s_{0} \in S$,

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P\left(X_{n+1}=s_{n+1} \mid X_{0}=s_{0}, \ldots, X_{n}=s_{n}\right)=P\left(X_{n+1}=s_{n+1} \mid X_{n}=s_{n}\right)
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The set $S$ is called the state space of the chain. If $S$ is a finite set, the chain is called a finite state Markov chain. $X_{0}$ is called the initial state.

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Example? the indices can be treated as time and the rvs as the observation of a process: Surfing Webpages, Weather prediction

## Random walks

initial distribution the distribution of the initial state $X_{0}$, which tells us how the Markov chain starts

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Distribution of the Markov Chain Let $\mu^{(0)}$ denote the initial distribution of the Markov chain, defined as

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\mu^{(0)}=\left(p\left(X_{0}=s_{1}\right), p\left(X_{0}=s_{2}\right), \ldots, p\left(X_{0}=s_{k}\right)\right)
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with state space $S=\left\{s_{1}, \ldots, s_{k}\right\}$. Similarly,

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\mu^{(t)}=\left(p\left(X_{t}=s_{1}\right), p\left(X_{t}=s_{2}\right), \ldots, p\left(X_{t}=s_{k}\right)\right)
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For example, $\mu^{(0)}=(1,0,0,0), \mu^{(1)}=\left(0, \frac{1}{2}, 0, \frac{1}{2}\right)$ in the previous example.


[^0]:    ${ }^{1}$ Klafter, J. and Sokolov, I.M., 2011. First steps in random walks: from tools to applications. Oxford University Press.

