Introduction to Monte Carlo Simulation With **Applications**

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Introduction (1)

Suppose our computer could hand us, upon demand, numbers,

$$
U_1, U_2, \ldots, U_n,
$$

that were independent and identically distributed (iid) samples from the continuous uniform distribution on the unit interval (0, 1):

$$
P(U\leq x)=x,\;x\in (0,1).
$$

The density function is given by

$$
f(x) = 1, x \in (0,1), \quad f(x) = 0, x \notin (0,1).
$$

QUESTION: What useful/practicle things can we do with these numbers?

ANSWER: As we shall see, by using the U_i to construct random variables X with any desired distribution, $F(x) = P(X \le x)$, we can then

- construct various stochastic processes (random walks, point processes, Markov chains, Brownian motion, etc.) so that we can
- simulate complicated models (queueing models in telecommunications, insurance risk models, asset price models, etc.) and numerically estimate quantities of interest (average delay at a call center, probability of ruin, price of an exotic option or derivative).

Outline

- Simple integration example; strong law of large numbers, the central limit theorem, confidence intervals for justifying the method.
- Basics techniques of stochastic simulation: Inverse transform method, acceptance rejection, polar method for generating random numbers X via using uniforms U .
- Simulating a Poisson process, and more generally a renewal process, Markov chains
- The binomial lattice model (BLM) for risky assets; some option pricing
- Simulating Brownian motion, geometric Brownian motion
- Markov Chain Monte Carlo (MCMC) method

Integration (1)

Suppose we wish to compute an integral

$$
\alpha=\int_0^1g(x)dx,
$$

for a function q that has no known antiderivative.

Observe that if $U \sim \text{unif}(0, 1)$, then the expected value of the random variable $g(U)$ is in fact the integral:

$$
E(g(U))=\int_0^1g(x)dx.
$$

This follows from the more general basic fact in probability that if X is a rv with density function $f(x)$, then the expected value of the random variable $g(X)$, for a function g is given by

$$
E(g(X))=\int g(x)f(x)dx.
$$

Integration (2)

In the following we shall be applying $X_i = g(U_i)$, where the U_i hence the X_i are iid rvs with mean $E(X) = \alpha = \int_0^1 g(x) dx$.

The famous and fundamental Strong Law of Large Numbers (SLLN) in probability theory asserts that:

Theorem (SLLN)

For any iid sequence of random variables X_1, X_2, \ldots with finite mean $E(X)$, it holds that with probability 1:

$$
\lim_{n\to\infty}\frac{1}{n}\sum_{j=1}^n X_j=E(X).
$$

Thus

$$
E(X) \approx \frac{1}{n} \sum_{j=1}^{n} X_j
$$
, for n large

Thus, using our computer's iid numbers U_1, U_2, \ldots, U_n , for large n, and defining $X_1 = g(U_1), \ldots, X_n = g(U_n)$, yields an approximation for our integral:

$$
\int_0^1 g(x)dx \approx \frac{1}{n}\sum_{j=1}^n g(U_j).
$$

Integration (4)

This is an example of Monte Carlo Simulation: We can numerically estimate an integral by first expressing it as an expected value, and then applying the SLLN.

This method works for multidimensional integrals as well: For example the $3 - d$ integral

$$
\int_0^1 \int_0^1 \int_0^1 g(x, y, z) dx dy dz = E(g(U_1, U_2, U_3)).
$$

We then define $X_1 = g(U_1, U_2, U_3)$, $X_2 = g(U_4, U_5, U_6)$ and so on and once again use the approximation

$$
\frac{1}{n}\sum_{j=1}^n X_j.
$$

The precision of this method does not depend on the dimension of the integral!!

The integration limits \int_0^1 can be more generally \int_a^b by using the change of variables $y = (x - a)/(b - a)$, $x \in (a, b)$. For example

$$
\int_a^b g(x)dx = (b-a)\int_0^1 g(a+(b-a)y)dy.
$$

The general theme: We need to compute something which equivalently can be expressed as an expected value, $E(X)$, for some (perhaps very complicated) rv X . We ask the computer to simulate iid copies of X , then use the SLLN to estimate $E(X)$ as an empirical average. We will give other examples soon, but first let us recall the notion and usefulness of confidence intervals.

In statistics, we estimate an unknown mean $\mu = E(X)$ of a distribution by collecting n iid samples from the distribution, X_1, \ldots, X_n and using the sample mean

$$
\overline{X}(n) = \frac{1}{n} \sum_{j=1}^{n} X_j.
$$
 (1)

This estimate is justified by the SLLN. But how good is our estimate?

Letting $\sigma^2 = \text{Var}(X)$ denote the variance of the distribution, we have

$$
Var(\overline{X}(n)) = \frac{\sigma^2}{n}.
$$
 (2)

The Central Limit Theorem asserts that as $n \to \infty$,

$$
Z_n \stackrel{\text{def}}{=} \frac{\sqrt{n}}{\sigma} (\overline{X}(n) - \mu),
$$

converges in distribution to N(0, 1), the unit normal distribution. Letting Z denote a $N(0, 1)$ rv, we conclude that for n sufficently large, $Z_n \approx Z$ in distribution, meaning that $P(Z_n \le x) \approx \Theta(x)$ for any $x \in \mathbb{R}$.

This allows us to construct confidence intervals for our estimate: (letting $z_{\alpha/2}$ be such that $P(Z > Z_{\alpha/2}) = \alpha/2$):

we say that the interval $\overline{X}(n) \pm z_{\alpha/2} \frac{\sigma}{\sqrt{n}}$ $\frac{1}{n}$ is a 100(1 − α)% confidence interval for the mean μ .

Typically, we would use (say) $\alpha = 0.05$ in which case $z_{\alpha/2} = z_{0.025} = 1.96$, and we thus obtain a 95% confidence interval $\overline{X}(n) \pm (1.96) \frac{\sigma}{\sqrt{n}}$ $\frac{1}{n}$.

We would not know σ in practice, so we instead use an estimate for it, the sample variance $s^2(n)$ defined by

$$
s^{2}(n) = \frac{1}{n-1}\sum_{j=1}^{n}(X_{j}-\overline{X}(n))^{2}.
$$

It can be shown that $s^2(n) \to \sigma^2$, with probability 1, as $n \to \infty$ and that $E(s^2(n)) = \sigma^2$, $n \ge 2$.

So, in practice we would use $s(n)$ is place of σ when constructing our confidence intervals. For example, a 95% confidence interval is given by $\overline{X}(n) \pm (1.96) \frac{s(n)}{\sqrt{n}}$ $\frac{\eta}{\overline{n}}$.

The following recursions can be derived; they are useful when implementing a simulation requiring a confidence interval:

$$
\overline{X}(n+1) = \overline{X}_n + \frac{X_{n+1} - \overline{X}(n)}{n+1},
$$

$$
S(n+1)^2 = \left(1 - \frac{1}{n}\right)S^2(n) + (n+1)(\overline{X}(n+1) - \overline{X}(n))^2.
$$

Inverse transform method

Theorem If $F(x) = P(X \le x)$ is the cumulative distribution of a desired rv, and if

$$
F^{-1}(y) = \min\{x : F(x) \ge y\}, y \in [0, 1],
$$

(the generalized inverse), then setting $X = F^{-1}(U)$ yields a rv distributed exactly as F.

For example, if X has the exponential distribution at rate λ , $F(x) = 1 - e^{-\lambda x}$, then this leads to the algorithm

$$
X=-\lambda^{-1}\ln(U).
$$

This method requires having an explicit closed form for F^{-1} which is not always possible (consider the normal distribution for example). Other methods: acceptance-rejection, polar method, and so on.

Inverse-transform method for discrete random variables

Consider a non-negative discrete rv X with probability mass function (pmf) $p(k) = P(X = k)$, $k \ge 0$. In this case, the construction $X = F^{-1}(U)$ is explicitly given by: $X = 0$ if $U \leq p(0)$,

$$
X = k, \text{ if } \sum_{i=0}^{k-1} p(i) < U \leq \sum_{i=0}^{k} p(i), \ k \geq 1.
$$

This is known as the discrete inverse-transform method. For example, if $P(X = 1) = p$ and $P(X = 0) = 1 - p$ is a Bernoulli (p) rv, then we can generate it via Set $X = 0$ if $U \le 1 - p$, set $X = 1$ if $U > 1 - p$. We also could do this via: Set $X = 1$ if $U \leq p$, set $X = 0$ if $U > p$.