Numerical Methods II

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If x is a random variable uniformly distributed on [0, 1] then the expectation of a function f(x) is equal to its integral:

$$\overline{f} = \mathbb{E}[f(x)] = I[f] = \int_0^1 f(x) \, \mathrm{d}x.$$

The generalisation to a d-dimensional "cube" $I^d = [0, 1]^d$, is

$$\overline{f} = \mathbb{E}[f(x)] = I[f] = \int_{I^d} f(x) \, \mathrm{d}x.$$

Thus the problem of finding expectations in finance is directly connected to the problem of numerical quadrature (integration), often in very large dimensions.

Suppose we have a sequence x_n of independent samples from the uniform distribution.

An approximation to the expectation/integral is given by

$$I_N[f] = N^{-1} \sum_{n=1}^N f(x_n).$$

Two key features:

• Unbiased: $\mathbb{E}\left[I_N[f]\right] = I[f]$

Convergent:

 $\lim_{N \to \infty} I_N[f] = I[f]$

In general, define

- error $\varepsilon_N(f) = I[f] I_N[f]$
- bias = $\mathbb{E}[\varepsilon_N(f)]$
- RMSE, "root-mean-square-error" = $\sqrt{\mathbb{E}[(\varepsilon_N(f))^2]}$

The Central Limit Theorem proves that for large ${\cal N}$

$$\varepsilon_N(f) \sim \sigma N^{-1/2} Z$$

with Z a N(0,1) random variable and σ^2 the variance of f:

$$\sigma^2 = \mathbb{E}[(f - \overline{f})^2] = \int_{I^d} \left(f(x) - \overline{f} \right)^2 \, \mathrm{d}x.$$

More precisely, provided σ is finite, then as $N \longrightarrow \infty$,

$$\mathsf{CDF}(N^{1/2}\sigma^{-1}\varepsilon_N) \longrightarrow \mathsf{CDF}(Z)$$

so that

$$\mathbb{P}\left[N^{1/2}\sigma^{-1}\varepsilon_N < s\right] \longrightarrow \mathbb{P}\left[Z < s\right] = \Phi(s)$$

and

$$\mathbb{P}\left[\left|N^{1/2}\sigma^{-1}\varepsilon_{N}\right| > s\right] \longrightarrow \mathbb{P}\left[\left|Z\right| > s\right] = 2 \Phi(-s)$$
$$\mathbb{P}\left[\left|N^{1/2}\sigma^{-1}\varepsilon_{N}\right| < s\right] \longrightarrow \mathbb{P}\left[\left|Z\right| < s\right] = 1 - 2 \Phi(-s)$$

Given N samples, the empirical variance is

$$\widetilde{\sigma}^2 = N^{-1} \sum_{n=1}^{N} \left(f(x_n) - I_N \right)^2 = I_N^{(2)} - (I_N)^2$$

where

$$I_N = N^{-1} \sum_{n=1}^N f(x_n), \qquad I_N^{(2)} = N^{-1} \sum_{n=1}^N (f(x_n))^2$$

 $\widetilde{\sigma}^2$ is a slightly biased estimator for σ^2 ; an unbiased estimator is

$$\widehat{\sigma}^2 = (N-1)^{-1} \sum_{n=1}^{N} \left(f(x_n) - I_N \right)^2 = \frac{N}{N-1} \left(I_N^{(2)} - (I_N)^2 \right)$$

Objective: want an accuracy of $\overline{\varepsilon}$ with confidence c. i.e. $|\varepsilon| < \overline{\varepsilon}$ with probability c.

How many samples do we need to use?

Recall,

$$\mathbb{P}\left[N^{1/2}\sigma^{-1}|\varepsilon| < s\right] \approx 1 - 2 \ \Phi(-s),$$

so define function s(c) such that

$$1 - 2 \Phi(-s) = c \iff s = -\Phi^{-1}((1-c)/2)$$

С	0.683	0.9545	0.9973	0.99994
S	1.0	2.0	3.0	4.0

Then $|\varepsilon| < N^{-1/2} \sigma s(c)$ with probability c, so to get $|\varepsilon| < \overline{\varepsilon}$ we can put

$$N^{-1/2} \widehat{\sigma} s(c) = \overline{\varepsilon} \implies N = \left(\frac{\widehat{\sigma} s(c)}{\overline{\varepsilon}}\right)^2$$

Note: twice as much accuracy requires 4 times as many samples.

How does Monte Carlo integration compare to grid based methods for *d*-dimensional integration?

MC error is proportional to $N^{-1/2}$ independent of the dimension.

If the integrand is sufficiently smooth, trapezoidal integration with $M = N^{1/d}$ points in each direction has

Error
$$\propto M^{-2} = N^{-2/d}$$

This scales better than MC for d < 4, but worse for d > 4. i.e. MC is better at handling high dimensional problems.

Geometric Brownian motion for single asset:

$$S(T) = S_0 \exp\left(\left(r - \frac{1}{2}\sigma^2\right)T + \sigma W(T)\right)$$

W(T) has a Normal distribution with mean 0, variance T; from this we will calculate the risk-neutral expectation for

$$V = \mathbb{E}\left[f(S(T))\right]$$

We can put

$$W(T) = \sqrt{T} Y = \sqrt{T} \Phi^{-1}(U)$$

where Y is a N(0,1) random variable, and U is uniformly distributed on [0,1].

Thus

$$V = \mathbb{E}\left[f(S(T))\right] = \int_0^1 f(S(T)) \, \mathrm{d}U,$$

with

$$S(T) = S_0 \exp\left(\left(r - \frac{1}{2}\sigma^2\right)T + \sigma\sqrt{T}Y\right)$$
$$= S_0 \exp\left(\left(r - \frac{1}{2}\sigma^2\right)T + \sigma\sqrt{T}\Phi^{-1}(U)\right)$$

For the European call option,

$$f(S) = \exp(-rT) (S - K)^+$$

while for the European put option

$$f(S) = \exp(-rT) (K-S)^+$$

where K is the strike price.

For numerical experiments we will consider a European call with r = 0.05, $\sigma = 0.2$, T = 1, $S_0 = 110$, K = 100.

The analytic value is known for comparison.



MC Lecture 2 – p. 13

MC calculation with up to 10^6 paths; true value = 17.663



MC Lecture 2 – p. 14

The upper and lower bounds are given by

Mean
$$\pm \frac{3 \widetilde{\sigma}}{\sqrt{N}}$$
,

so more than a 99.7% probability that the true value lies within these bounds.

MATLAB code:

r=0.05; sig=0.2; T=1; S0=110; K=100;

N = 1:100000;

- U = rand(1,max(N)); % uniform random variable
- Y = norminv(U); % inverts Normal cum. fn.
- $S = S0 + exp((r-sig^2/2) + T + sig + sqrt(T) + Y);$

$$F = \exp(-r * T) * \max(0, S - K);$$

sum1 = cumsum(F); % cumulative summation of sum2 = cumsum(F.^2); % payoff and its square val = sum1./N; rms = sqrt(sum2./N - val.^2);

err = european_call(r,sig,T,S0,K,'value') - val;

plot(N,err, ... N,err-3*rms./sqrt(N), ... N,err+3*rms./sqrt(N)) axis([0 length(N) -1 1]) xlabel('N'); ylabel('Error') legend('MC error','lower bound','upper bound')

New application: basket option

European call for arithmetic average of M stocks which are correlated so that

 $\mathrm{d}S_i = r\,S_i\,\mathrm{d}t + \sigma_i S_i \mathrm{d}W_i$

with the different dW_i <u>not</u> independent.

As before, get

$$S_i(T) = S_i(0) \exp\left(\left(r - \frac{1}{2}\sigma_i^2\right)T + \sigma_i W_i(T)\right)$$

If $\sigma_i W_i(T)$ have covariance matrix Σ , then use Cholesky factorisation $LL^T = \Sigma$ to get

$$S_i(T) = S_i(0) \exp\left(\left(r - \frac{1}{2}\sigma_i^2\right)T + \sum_j L_{ij}Y_j\right)$$

where Y_j are independent N(0, 1) random variables.

Each Y_i can in turn be expressed as $\Phi^{-1}(U_i)$ where the U_i are uniformly, and independently, distributed on [0, 1].

The payoff is

$$f = \exp(-rT) \left(\frac{1}{M}\sum_{i}S_{i} - K\right)^{+}$$

and so the expectation can be written as the *M*-dimensional integral

$$\int_{I^M} f(U) \, \mathrm{d}U.$$

This is a good example for Monte Carlo simulation – cost scales linearly with the number of stocks, whereas it would be exponential for grid-based numerical integration.

Final Words

- Monte Carlo quadrature is straightforward and robust
- Confidence bounds can be obtained as part of the calculation
- Can calculate the number of samples N needed for chosen accuracy
- Much more efficient than grid-based methods for high dimensions
- Accuracy = $O(N^{-1/2})$, CPU time = O(N)
 - \implies accuracy = $O(CPU time^{-1/2})$
 - \implies CPU time = $O(\text{accuracy}^{-2})$