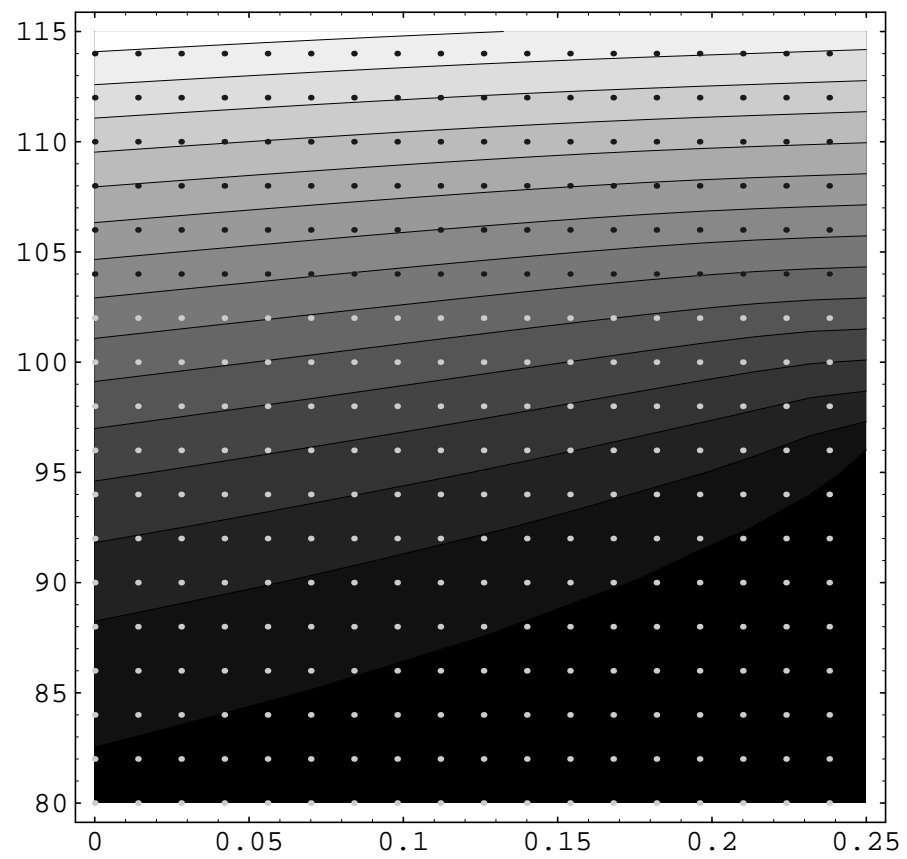


Numerical Methods

All science is dominated
by the idea of approximation.
— Bertrand Russell

Finite-Difference Methods

- Place a grid of points on the space over which the desired function takes value.
- Then approximate the function value at each of these points (p. 667).
- Solve the equation numerically by introducing difference equations in place of derivatives.



Example: Poisson's Equation

- It is $\partial^2\theta/\partial x^2 + \partial^2\theta/\partial y^2 = -\rho(x, y)$.
- Replace second derivatives with finite differences through central difference.
- Introduce evenly spaced grid points with distance of Δx along the x axis and Δy along the y axis.
- The finite difference form is

$$-\rho(x_i, y_j) = \frac{\theta(x_{i+1}, y_j) - 2\theta(x_i, y_j) + \theta(x_{i-1}, y_j)}{(\Delta x)^2} + \frac{\theta(x_i, y_{j+1}) - 2\theta(x_i, y_j) + \theta(x_i, y_{j-1}))}{(\Delta y)^2}.$$

Example: Poisson's Equation (concluded)

- In the above, $\Delta x \equiv x_i - x_{i-1}$ and $\Delta y \equiv y_j - y_{j-1}$ for $i, j = 1, 2, \dots$
- When the grid points are evenly spaced in both axes so that $\Delta x = \Delta y = h$, the difference equation becomes

$$\begin{aligned} -h^2 \rho(x_i, y_j) &= \theta(x_{i+1}, y_j) + \theta(x_{i-1}, y_j) \\ &\quad + \theta(x_i, y_{j+1}) + \theta(x_i, y_{j-1}) - 4\theta(x_i, y_j). \end{aligned}$$

- Given boundary values, we can solve for the x_i s and the y_j s within the square $[\pm L, \pm L]$.
- From now on, $\theta_{i,j}$ will denote the finite-difference approximation to the exact $\theta(x_i, y_j)$.

Explicit Methods

- Consider the diffusion equation
 $D(\partial^2\theta/\partial x^2) - (\partial\theta/\partial t) = 0$.
- Use evenly spaced grid points (x_i, t_j) with distances Δx and Δt , where $\Delta x \equiv x_{i+1} - x_i$ and $\Delta t \equiv t_{j+1} - t_j$.
- Employ central difference for the second derivative and forward difference for the time derivative to obtain

$$\left. \frac{\partial\theta(x, t)}{\partial t} \right|_{t=t_j} = \frac{\theta(x, t_{j+1}) - \theta(x, t_j)}{\Delta t} + \dots, \quad (78)$$

$$\left. \frac{\partial^2\theta(x, t)}{\partial x^2} \right|_{x=x_i} = \frac{\theta(x_{i+1}, t) - 2\theta(x_i, t) + \theta(x_{i-1}, t)}{(\Delta x)^2} + \dots. \quad (79)$$

Explicit Methods (continued)

- Next, assemble Eqs. (78) and (79) into a single equation at (x_i, t_j) .
- But we need to decide how to evaluate x in the first equation and t in the second.
- Since central difference around x_i is used in Eq. (79), we might as well use x_i for x in Eq. (78).
- Two choices are possible for t in Eq. (79).
- The first choice uses $t = t_j$ to yield the following finite-difference equation,

$$\frac{\theta_{i,j+1} - \theta_{i,j}}{\Delta t} = D \frac{\theta_{i+1,j} - 2\theta_{i,j} + \theta_{i-1,j}}{(\Delta x)^2}. \quad (80)$$

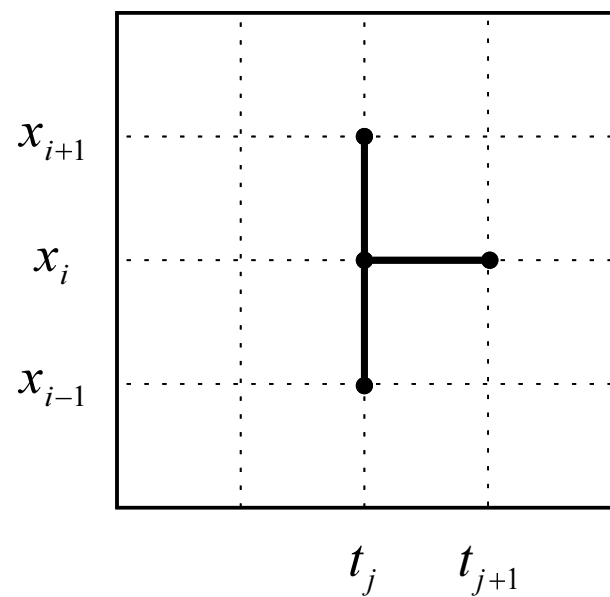
Explicit Methods (continued)

- The stencil of grid points involves four values, $\theta_{i,j+1}$, $\theta_{i,j}$, $\theta_{i+1,j}$, and $\theta_{i-1,j}$.
- Rearrange Eq. (80) on p. 671 as

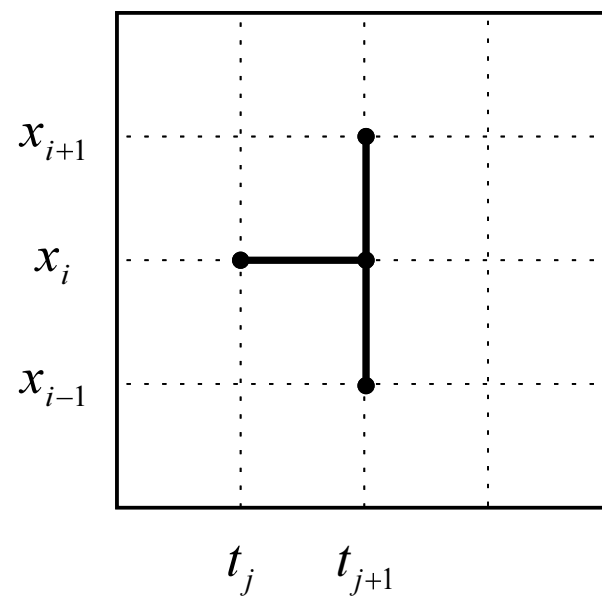
$$\theta_{i,j+1} = \frac{D\Delta t}{(\Delta x)^2} \theta_{i+1,j} + \left(1 - \frac{2D\Delta t}{(\Delta x)^2}\right) \theta_{i,j} + \frac{D\Delta t}{(\Delta x)^2} \theta_{i-1,j}.$$

- We can calculate $\theta_{i,j+1}$ from $\theta_{i,j}$, $\theta_{i+1,j}$, $\theta_{i-1,j}$, at the previous time t_j (see exhibit (a) on next page).

Stencils



(a)



(b)

Explicit Methods (concluded)

- Starting from the initial conditions at t_0 , that is, $\theta_{i,0} = \theta(x_i, t_0)$, $i = 1, 2, \dots$, we calculate

$$\theta_{i,1}, \quad i = 1, 2, \dots .$$

- And then

$$\theta_{i,2}, \quad i = 1, 2, \dots .$$

- And so on.

Stability

- The explicit method is numerically unstable unless

$$\Delta t \leq (\Delta x)^2 / (2D).$$

- A numerical method is unstable if the solution is highly sensitive to changes in initial conditions.
- The stability condition may lead to high running times and memory requirements.
- For instance, halving Δx would imply quadrupling $(\Delta t)^{-1}$, resulting in a running time 8 times as much.

Explicit Method and Trinomial Tree

- Recall that

$$\theta_{i,j+1} = \frac{D\Delta t}{(\Delta x)^2} \theta_{i+1,j} + \left(1 - \frac{2D\Delta t}{(\Delta x)^2}\right) \theta_{i,j} + \frac{D\Delta t}{(\Delta x)^2} \theta_{i-1,j}.$$

- When the stability condition is satisfied, the three coefficients for $\theta_{i+1,j}$, $\theta_{i,j}$, and $\theta_{i-1,j}$ all lie between zero and one and sum to one.
- They can be interpreted as probabilities.
- So the finite-difference equation becomes identical to backward induction on trinomial trees!
- The freedom in choosing Δx corresponds to similar freedom in the construction of trinomial trees.

Implicit Methods

- Suppose we use $t = t_{j+1}$ in Eq. (79) on p. 670 instead.
- The finite-difference equation becomes

$$\frac{\theta_{i,j+1} - \theta_{i,j}}{\Delta t} = D \frac{\theta_{i+1,j+1} - 2\theta_{i,j+1} + \theta_{i-1,j+1}}{(\Delta x)^2}. \quad (81)$$

- The stencil involves $\theta_{i,j}$, $\theta_{i,j+1}$, $\theta_{i+1,j+1}$, and $\theta_{i-1,j+1}$.
- This method is implicit:
 - The value of any one of the three quantities at t_{j+1} cannot be calculated unless the other two are known.
 - See exhibit (b) on p. 673.

Implicit Methods (continued)

- Equation (81) can be rearranged as

$$\theta_{i-1,j+1} - (2 + \gamma) \theta_{i,j+1} + \theta_{i+1,j+1} = -\gamma \theta_{i,j},$$

where $\gamma \equiv (\Delta x)^2 / (D \Delta t)$.

- This equation is unconditionally stable.
- Suppose the boundary conditions are given at $x = x_0$ and $x = x_{N+1}$.
- After $\theta_{i,j}$ has been calculated for $i = 1, 2, \dots, N$, the values of $\theta_{i,j+1}$ at time t_{j+1} can be computed as the solution to the following tridiagonal linear system,

Implicit Methods (continued)

$$\begin{bmatrix} a & 1 & 0 & \cdots & \cdots & 0 \\ 1 & a & 1 & 0 & \cdots & 0 \\ 0 & 1 & a & 1 & \cdots & 0 \\ \vdots & \ddots & \ddots & \ddots & \ddots & \vdots \\ \vdots & \ddots & \ddots & \ddots & \ddots & \vdots \\ 0 & \cdots & \cdots & 1 & a & 1 \\ 0 & \cdots & \cdots & \cdots & 1 & a \end{bmatrix} \begin{bmatrix} \theta_{1,j+1} \\ \theta_{2,j+1} \\ \theta_{3,j+1} \\ \vdots \\ \vdots \\ \vdots \\ \vdots \\ \vdots \\ \theta_{N,j+1} \end{bmatrix} = \begin{bmatrix} -\gamma\theta_{1,j} - \theta_{0,j+1} \\ -\gamma\theta_{2,j} \\ -\gamma\theta_{3,j} \\ \vdots \\ \vdots \\ \vdots \\ \vdots \\ -\gamma\theta_{N-1,j} \\ -\gamma\theta_{N,j} - \theta_{N+1,j+1} \end{bmatrix},$$

where $a \equiv -2 - \gamma$.

Implicit Methods (concluded)

- Tridiagonal systems can be solved in $O(N)$ time and $O(N)$ space.
 - Never invert a matrix to solve a tridiagonal system.
- The matrix above is nonsingular when $\gamma \geq 0$.
 - A square matrix is nonsingular if its inverse exists.

Crank-Nicolson Method

- Take the average of explicit method (80) on p. 671 and implicit method (81) on p. 677:

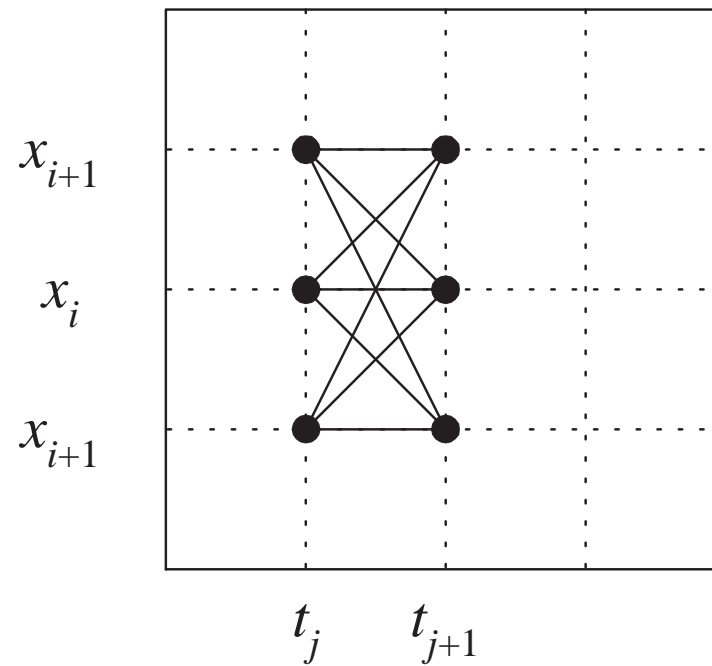
$$\begin{aligned} & \frac{\theta_{i,j+1} - \theta_{i,j}}{\Delta t} \\ = & \frac{1}{2} \left(D \frac{\theta_{i+1,j} - 2\theta_{i,j} + \theta_{i-1,j}}{(\Delta x)^2} + D \frac{\theta_{i+1,j+1} - 2\theta_{i,j+1} + \theta_{i-1,j+1}}{(\Delta x)^2} \right). \end{aligned}$$

- After rearrangement,

$$\gamma\theta_{i,j+1} - \frac{\theta_{i+1,j+1} - 2\theta_{i,j+1} + \theta_{i-1,j+1}}{2} = \gamma\theta_{i,j} + \frac{\theta_{i+1,j} - 2\theta_{i,j} + \theta_{i-1,j}}{2}.$$

- This is an unconditionally stable implicit method with excellent rates of convergence.

Stencil



Numerically Solving the Black-Scholes PDE

- See text.
- Brennan and Schwartz (1978) analyze the stability of the implicit method.

Monte Carlo Simulation^a

- Monte Carlo simulation is a sampling scheme.
- In many important applications within finance and without, Monte Carlo is one of the few feasible tools.
- When the time evolution of a stochastic process is not easy to describe analytically, Monte Carlo may very well be the only strategy that succeeds consistently.

^aA top 10 algorithm according to Dongarra and Sullivan (2000).

The Big Idea

- Assume X_1, X_2, \dots, X_n have a joint distribution.
- $\theta \equiv E[g(X_1, X_2, \dots, X_n)]$ for some function g is desired.
- We generate

$$\left(x_1^{(i)}, x_2^{(i)}, \dots, x_n^{(i)}\right), \quad 1 \leq i \leq N$$

independently with the same joint distribution as (X_1, X_2, \dots, X_n) .

- Set

$$Y_i \equiv g\left(x_1^{(i)}, x_2^{(i)}, \dots, x_n^{(i)}\right).$$

The Big Idea (concluded)

- Y_1, Y_2, \dots, Y_N are independent and identically distributed random variables.
- Each Y_i has the same distribution as

$$Y \equiv g(X_1, X_2, \dots, X_n).$$

- Since the average of these N random variables, \bar{Y} , satisfies $E[\bar{Y}] = \theta$, it can be used to estimate θ .
- The strong law of large numbers says that this procedure converges almost surely.
- The number of replications (or independent trials), N , is called the sample size.

Accuracy

- The Monte Carlo estimate and true value may differ owing to two reasons:
 1. Sampling variation.
 2. The discreteness of the sample paths.^a
- The first can be controlled by the number of replications.
- The second can be controlled by the number of observations along the sample path.

^aThis may not be an issue if the financial derivative only requires discrete sampling along the time dimension, such as the discrete barrier option.

Accuracy and Number of Replications

- The statistical error of the sample mean \bar{Y} of the random variable Y grows as $1/\sqrt{N}$.
 - Because $\text{Var}[\bar{Y}] = \text{Var}[Y]/N$.
- In fact, this convergence rate is asymptotically optimal.^a
- So the variance of the estimator \bar{Y} can be reduced by a factor of $1/N$ by doing N times as much work.
- This is amazing because the same order of convergence holds independently of the dimension n .

^aThe Berry-Esseen theorem.

Accuracy and Number of Replications (concluded)

- In contrast, classic numerical integration schemes have an error bound of $O(N^{-c/n})$ for some constant $c > 0$.
 - n is the dimension.
- The required number of evaluations thus grows exponentially in n to achieve a given level of accuracy.
 - The curse of dimensionality.
- The Monte Carlo method, for example, is more efficient than alternative procedures for multivariate derivatives.

Monte Carlo Option Pricing

- For the pricing of European options on a dividend-paying stock, we may proceed as follows.
- Assume $dS/S = \mu dt + \sigma dW$.
- Stock prices S_1, S_2, S_3, \dots at times $\Delta t, 2\Delta t, 3\Delta t, \dots$ can be generated via

$$S_{i+1} = S_i e^{(\mu - \sigma^2/2) \Delta t + \sigma \sqrt{\Delta t} \xi}, \quad \xi \sim N(0, 1). \quad (82)$$

Monte Carlo Option Pricing (continued)

- If we discretize $dS/S = \mu dt + \sigma dW$ directly, we will obtain

$$S_{i+1} = S_i + S_i \mu \Delta t + S_i \sigma \sqrt{\Delta t} \xi.$$

- But this is locally normally distributed, not lognormally, hence biased.^a
- In practice, this is not expected to be a major problem as long as Δt is sufficiently small.

^aContributed by Mr. Tai, Hui-Chin (R97723028) on April 22, 2009.

Monte Carlo Option Pricing (continued)

- Non-dividend-paying stock prices in a risk-neutral economy can be generated by setting $\mu = r$ and $\Delta t = T$.
 - 1: $C := 0$; {Accumulated terminal option value.}
 - 2: **for** $i = 1, 2, 3, \dots, m$ **do**
 - 3: $P := S \times e^{(r-\sigma^2/2)T + \sigma\sqrt{T}\xi}$, $\xi \sim N(0, 1)$;
 - 4: $C := C + \max(P - X, 0)$;
 - 5: **end for**
 - 6: **return** Ce^{-rT}/m ;

Monte Carlo Option Pricing (concluded)

- Pricing Asian options is also easy.
 - 1: $C := 0$;
 - 2: **for** $i = 1, 2, 3, \dots, m$ **do**
 - 3: $P := S$; $M := S$;
 - 4: **for** $j = 1, 2, 3, \dots, n$ **do**
 - 5: $P := P \times e^{(r - \sigma^2/2)(T/n) + \sigma\sqrt{T/n} \xi}$;
 - 6: $M := M + P$;
 - 7: **end for**
 - 8: $C := C + \max(M/(n + 1) - X, 0)$;
 - 9: **end for**
 - 10: **return** Ce^{-rT}/m ;

How about American Options?

- Standard Monte Carlo simulation is inappropriate for American options because of early exercise (why?).
- It is difficult to determine the early-exercise point based on one single path.
- But Monte Carlo simulation can be modified to price American options with small biases (pp. 743ff).^a

^aLongstaff and Schwartz (2001).

Delta and Common Random Numbers

- In estimating delta, it is natural to start with the finite-difference estimate

$$e^{-r\tau} \frac{E[P(S + \epsilon)] - E[P(S - \epsilon)]}{2\epsilon}.$$

- $P(x)$ is the terminal payoff of the derivative security when the underlying asset's initial price equals x .
- Use simulation to estimate $E[P(S + \epsilon)]$ first.
- Use another simulation to estimate $E[P(S - \epsilon)]$.
- Finally, apply the formula to approximate the delta.
- This is also called the bump-and-revalue method.

Delta and Common Random Numbers (concluded)

- This method is not recommended because of its high variance.
- A much better approach is to use common random numbers to lower the variance:

$$e^{-r\tau} E \left[\frac{P(S + \epsilon) - P(S - \epsilon)}{2\epsilon} \right].$$

- Here, the *same* random numbers are used for $P(S + \epsilon)$ and $P(S - \epsilon)$.
- This holds for gamma and cross gammas (for multivariate derivatives).

Problems with the Bump-and-Revalue Method

- Consider the binary option with payoff

$$\begin{cases} 1, & \text{if } S(T) > X, \\ 0, & \text{otherwise.} \end{cases}$$

- Then

$$P(S + \epsilon) - P(S - \epsilon) = \begin{cases} 1, & \text{if } P(S + \epsilon) > X \text{ and} \\ & P(S - \epsilon) < X, \\ 0, & \text{otherwise.} \end{cases}$$

- So the finite-difference estimate per run for the (undiscounted) delta is 0 or $O(1/\epsilon)$.
- This means high variance.

Gamma

- The finite-difference formula for gamma is

$$e^{-r\tau} E \left[\frac{P(S + \epsilon) - 2 \times P(S) + P(S - \epsilon)}{\epsilon^2} \right].$$

- For a correlation option with multiple underlying assets, the finite-difference formula for the cross gamma $\partial^2 P(S_1, S_2, \dots) / (\partial S_1 \partial S_2)$ is:

$$e^{-r\tau} E \left[\frac{P(S_1 + \epsilon_1, S_2 + \epsilon_2) - P(S_1 - \epsilon_1, S_2 + \epsilon_2)}{4\epsilon_1 \epsilon_2} - \frac{P(S_1 + \epsilon_1, S_2 - \epsilon_2) + P(S_1 - \epsilon_1, S_2 - \epsilon_2)}{4\epsilon_1 \epsilon_2} \right].$$

Gamma (continued)

- Choosing an ϵ of the right magnitude can be challenging.
 - If ϵ is too large, inaccurate Greeks result.
 - If ϵ is too small, unstable Greeks result.
- This phenomenon is sometimes called the curse of differentiation.

Gamma (continued)

- In general, suppose

$$\frac{\partial^i}{\partial \theta^i} e^{-r\tau} E[P(S)] = e^{-r\tau} E \left[\frac{\partial^i P(S)}{\partial \theta^i} \right]$$

holds for all $i > 0$, where θ is a parameter of interest.

- Then formulas for the Greeks become integrals.
- As a result, we avoid ϵ , finite differences, and resimulation.

Gamma (concluded)

- This is indeed possible for a broad class of payoff functions.^a
 - Roughly speaking, any payoff function that is equal to a sum of products of differentiable functions and indicator functions with the right kind of support.
 - For example, the payoff of a call is

$$\max(S(T) - X, 0) = (S(T) - X)I_{\{S(T) - X \geq 0\}}.$$

- The results are too technical to cover here.

^aTeng (R91723054) (2004) and Lyuu and Teng (R91723054) (2011).

Biases in Pricing Continuously Monitored Options with Monte Carlo

- We are asked to price a continuously monitored up-and-out call with barrier H .
- The Monte Carlo method samples the stock price at n discrete time points t_1, t_2, \dots, t_n .
- A sample path $S(t_0), S(t_1), \dots, S(t_n)$ is produced.
 - Here, $t_0 = 0$ is the current time, and $t_n = T$ is the expiration time of the option.

Biases in Pricing Continuously Monitored Options with Monte Carlo (continued)

- If all of the sampled prices are below the barrier, this sample path pays $\max(S(t_n) - X, 0)$.
- Repeating these steps and averaging the payoffs yield a Monte Carlo estimate.


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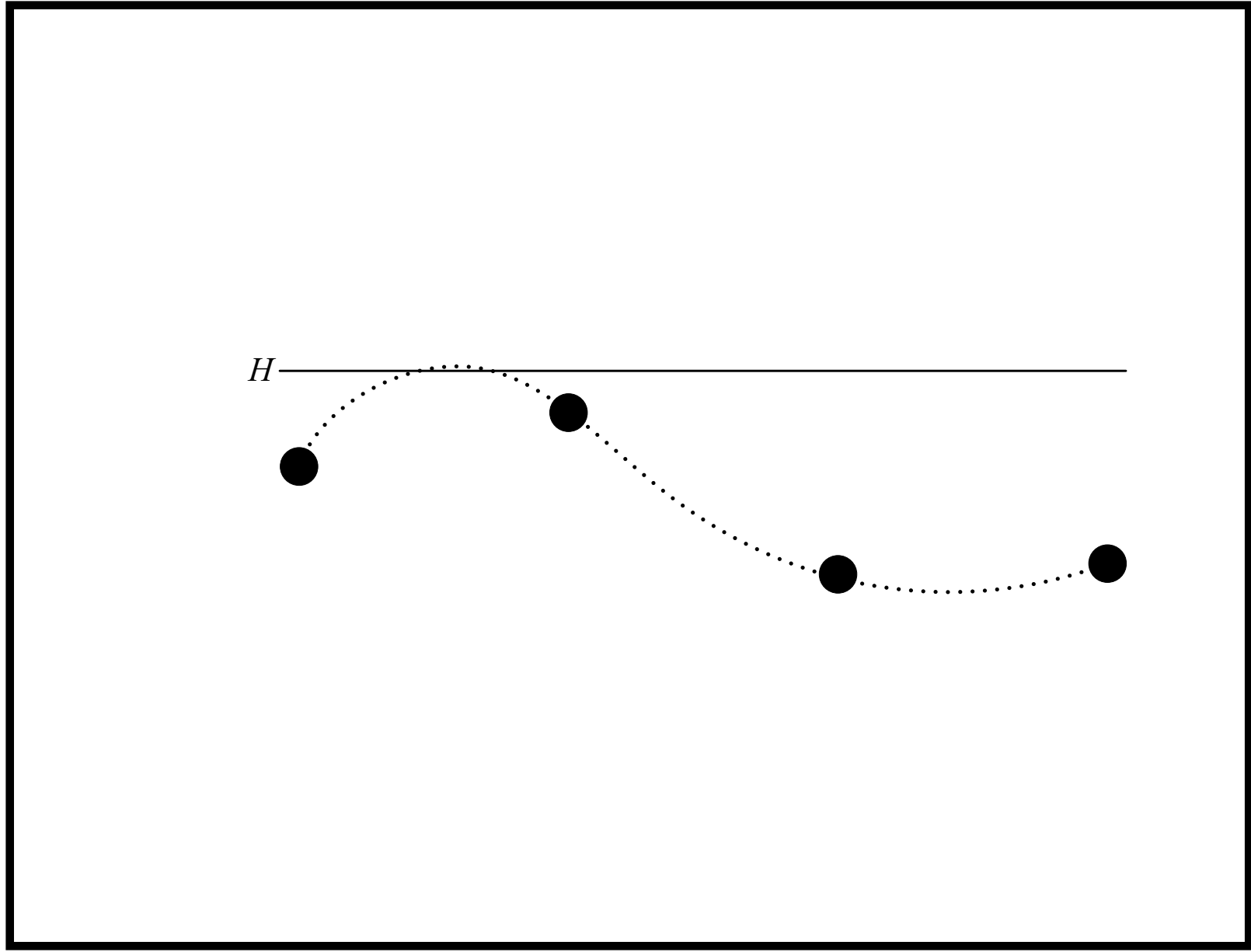
1:  $C := 0$ ;
2: for  $i = 1, 2, 3, \dots, m$  do
3:    $P := S$ ; hit := 0;
4:   for  $j = 1, 2, 3, \dots, n$  do
5:      $P := P \times e^{(r - \sigma^2/2)(T/n) + \sigma\sqrt{(T/n)} \xi}$ ;
6:     if  $P \geq H$  then
7:       hit := 1;
8:       break;
9:     end if
10:  end for
11:  if hit = 0 then
12:     $C := C + \max(P - X, 0)$ ;
13:  end if
14: end for
15: return  $Ce^{-rT}/m$ ;

```

Biases in Pricing Continuously Monitored Options with Monte Carlo (continued)

- This estimate is biased.^a
 - Suppose none of the sampled prices on a sample path equals or exceeds the barrier H .
 - It remains possible for the continuous sample path that passes through them to hit the barrier between sampled time points (see plot on next page).

^aShevchenko (2003).



Biases in Pricing Continuously Monitored Options with Monte Carlo (concluded)

- The bias can certainly be lowered by increasing the number of observations along the sample path.
- However, even daily sampling may not suffice.
- The computational cost also rises as a result.

Brownian Bridge Approach to Pricing Barrier Options

- We desire an unbiased estimate which can be calculated efficiently.
- The above-mentioned payoff should be multiplied by the probability p that a continuous sample path does *not* hit the barrier conditional on the sampled prices.
- This methodology is called the Brownian bridge approach.
- Formally, we have

$$p \equiv \text{Prob}[S(t) < H, 0 \leq t \leq T \mid S(t_0), S(t_1), \dots, S(t_n)].$$

Brownian Bridge Approach to Pricing Barrier Options (continued)

- As a barrier is hit over a time interval if and only if the maximum stock price over that period is at least H ,

$$p = \text{Prob} \left[\max_{0 \leq t \leq T} S(t) < H \mid S(t_0), S(t_1), \dots, S(t_n) \right].$$

- Luckily, the conditional distribution of the maximum over a time interval given the beginning and ending stock prices is known.

Brownian Bridge Approach to Pricing Barrier Options (continued)

Lemma 21 Assume S follows $dS/S = \mu dt + \sigma dW$ and define

$$\zeta(x) \equiv \exp \left[-\frac{2 \ln(x/S(t)) \ln(x/S(t + \Delta t))}{\sigma^2 \Delta t} \right].$$

(1) If $H > \max(S(t), S(t + \Delta t))$, then

$$\text{Prob} \left[\max_{t \leq u \leq t + \Delta t} S(u) < H \mid S(t), S(t + \Delta t) \right] = 1 - \zeta(H).$$

(2) If $h < \min(S(t), S(t + \Delta t))$, then

$$\text{Prob} \left[\min_{t \leq u \leq t + \Delta t} S(u) > h \mid S(t), S(t + \Delta t) \right] = 1 - \zeta(h).$$

Brownian Bridge Approach to Pricing Barrier Options (continued)

- Lemma 21 gives the probability that the barrier is not hit in a time interval, given the starting and ending stock prices.
- For our up-and-out call, choose $n = 1$.
- As a result,

$$p = \begin{cases} 1 - \exp \left[-\frac{2 \ln(H/S(0)) \ln(H/S(T))}{\sigma^2 T} \right], & \text{if } H > \max(S(0), S(T)), \\ 0, & \text{otherwise.} \end{cases}$$

Brownian Bridge Approach to Pricing Barrier Options (continued)

```
1:  $C := 0$ ;  
2: for  $i = 1, 2, 3, \dots, m$  do  
3:    $P := S \times e^{(r-q-\sigma^2/2)T + \sigma\sqrt{T} \xi()};$   
4:   if  $(S < H \text{ and } P < H) \text{ or } (S > H \text{ and } P > H)$  then  
5:      $C := C + \max(P - X, 0) \times \left\{ 1 - \exp \left[ -\frac{2 \ln(H/S) \times \ln(H/P)}{\sigma^2 T} \right] \right\};$   
6:   end if  
7: end for  
8: return  $C e^{-rT} / m$ ;
```

Brownian Bridge Approach to Pricing Barrier Options (concluded)

- The idea can be generalized.
- For example, we can handle more complex barrier options.
- Consider an up-and-out call with barrier H_i for the time interval $(t_i, t_{i+1}]$, $0 \leq i < n$.
- This option thus contains n barriers.
- Multiply the probabilities for the n time intervals to obtain the desired probability adjustment term.

Variance Reduction

- The statistical efficiency of Monte Carlo simulation can be measured by the variance of its output.
- If this variance can be lowered without changing the expected value, fewer replications are needed.
- Methods that improve efficiency in this manner are called variance-reduction techniques.
- Such techniques become practical when the added costs are outweighed by the reduction in sampling.

Variance Reduction: Antithetic Variates

- We are interested in estimating $E[g(X_1, X_2, \dots, X_n)]$, where X_1, X_2, \dots, X_n are independent.
- Let Y_1 and Y_2 be random variables with the same distribution as $g(X_1, X_2, \dots, X_n)$.
- Then

$$\text{Var} \left[\frac{Y_1 + Y_2}{2} \right] = \frac{\text{Var}[Y_1]}{2} + \frac{\text{Cov}[Y_1, Y_2]}{2}.$$

- $\text{Var}[Y_1]/2$ is the variance of the Monte Carlo method with two independent replications.
- The variance $\text{Var}[(Y_1 + Y_2)/2]$ is smaller than $\text{Var}[Y_1]/2$ when Y_1 and Y_2 are negatively correlated.

Variance Reduction: Antithetic Variates (continued)

- For each simulated sample path X , a second one is obtained by *reusing* the random numbers on which the first path is based.
- This yields a second sample path Y .
- Two estimates are then obtained: One based on X and the other on Y .
- If N independent sample paths are generated, the antithetic-variates estimator averages over $2N$ estimates.

Variance Reduction: Antithetic Variates (continued)

- Consider process $dX = a_t dt + b_t \sqrt{dt} \xi$.
- Let g be a function of n samples X_1, X_2, \dots, X_n on the sample path.
- We are interested in $E[g(X_1, X_2, \dots, X_n)]$.
- Suppose one simulation run has realizations $\xi_1, \xi_2, \dots, \xi_n$ for the normally distributed fluctuation term ξ .
- This generates samples x_1, x_2, \dots, x_n .
- The estimate is then $g(\mathbf{x})$, where $\mathbf{x} \equiv (x_1, x_2, \dots, x_n)$.

Variance Reduction: Antithetic Variates (concluded)

- The antithetic-variates method does not sample n more numbers from ξ for the second estimate $g(\mathbf{x}')$.
- Instead, generate the sample path $\mathbf{x}' \equiv (x'_1, x'_2, \dots, x'_n)$ from $-\xi_1, -\xi_2, \dots, -\xi_n$.
- Compute $g(\mathbf{x}')$.
- Output $(g(\mathbf{x}) + g(\mathbf{x}'))/2$.
- Repeat the above steps for as many times as required by accuracy.

Variance Reduction: Conditioning

- We are interested in estimating $E[X]$.
- Suppose here is a random variable Z such that $E[X | Z = z]$ can be efficiently and precisely computed.
- $E[X] = E[E[X | Z]]$ by the law of iterated conditional expectations.
- Hence the random variable $E[X | Z]$ is also an unbiased estimator of $E[X]$.

Variance Reduction: Conditioning (concluded)

- As

$$\text{Var}[E[X | Z]] \leq \text{Var}[X],$$

$E[X | Z]$ has a smaller variance than observing X directly.

- First obtain a random observation z on Z .
- Then calculate $E[X | Z = z]$ as our estimate.
 - There is no need to resort to simulation in computing $E[X | Z = z]$.
- The procedure can be repeated a few times to reduce the variance.

Control Variates

- Use the analytic solution of a similar yet simpler problem to improve the solution.
- Suppose we want to estimate $E[X]$ and there exists a random variable Y with a known mean $\mu \equiv E[Y]$.
- Then $W \equiv X + \beta(Y - \mu)$ can serve as a “controlled” estimator of $E[X]$ for any constant β .
 - However β is chosen, W remains an unbiased estimator of $E[X]$ as

$$E[W] = E[X] + \beta E[Y - \mu] = E[X].$$

Control Variates (continued)

- Note that

$$\text{Var}[W] = \text{Var}[X] + \beta^2 \text{Var}[Y] + 2\beta \text{Cov}[X, Y], \quad (83)$$

- Hence W is less variable than X if and only if

$$\beta^2 \text{Var}[Y] + 2\beta \text{Cov}[X, Y] < 0. \quad (84)$$

Control Variates (concluded)

- The success of the scheme clearly depends on both β and the choice of Y .
 - For example, arithmetic average-rate options can be priced by choosing Y to be the otherwise identical geometric average-rate option's price and $\beta = -1$.
- This approach is much more effective than the antithetic-variates method.

Choice of Y

- In general, the choice of Y is ad hoc, and experiments must be performed to confirm the wisdom of the choice.
- Try to match calls with calls and puts with puts.^a
- On many occasions, Y is a discretized version of the derivative that gives μ .
 - Discretely monitored geometric average-rate option vs. the continuously monitored geometric average-rate option given by formulas (32) on p. 355.
- For some choices, the discrepancy can be significant, such as the lookback option.^b

^aContributed by Ms. Teng, Huei-Wen (R91723054) on May 25, 2004.

^bContributed by Mr. Tsai, Hwai (R92723049) on May 12, 2004.

Optimal Choice of β

- Equation (83) on p. 722 is minimized when

$$\beta = -\text{Cov}[X, Y] / \text{Var}[Y].$$

– It is called beta in the book.

- For this specific β ,

$$\text{Var}[W] = \text{Var}[X] - \frac{\text{Cov}[X, Y]^2}{\text{Var}[Y]} = (1 - \rho_{X,Y}^2) \text{Var}[X],$$

where $\rho_{X,Y}$ is the correlation between X and Y .

- Note that the variance can never be increased with the optimal choice.

Optimal Choice of β (continued)

- Furthermore, the stronger X and Y are correlated, the greater the reduction in variance.
- For example, if this correlation is nearly perfect (± 1), we could control X almost exactly.
- Typically, neither $\text{Var}[Y]$ nor $\text{Cov}[X, Y]$ is known.
- Therefore, we cannot obtain the maximum reduction in variance.
- We can guess these values and hope that the resulting W does indeed have a smaller variance than X .

Optimal Choice of β (continued)

- A second possibility is to use the simulated data to estimate these quantities.
 - How to do it efficiently in terms of time and space?
- Observe that $-\beta$ has the same sign as the correlation between X and Y .
- Hence, if X and Y are positively correlated, $\beta < 0$, then X is adjusted downward whenever $Y > \mu$ and upward otherwise.
- The opposite is true when X and Y are negatively correlated, in which case $\beta > 0$.

Optimal Choice of β (concluded)

- Suppose a suboptimal $\beta + \epsilon$ is used instead.
- The variance increases by only $\epsilon^2 \text{Var}[Y]$.^a

^aHan and Lai (2010).

A Pitfall

- A potential pitfall is to sample X and Y independently.
- In this case, $\text{Cov}[X, Y] = 0$.
- Equation (83) on p. 722 becomes

$$\text{Var}[W] = \text{Var}[X] + \beta^2 \text{Var}[Y].$$

- So whatever Y is, the variance is *increased*!
- Lesson: X and Y must be correlated.

Problems with the Monte Carlo Method

- The error bound is only probabilistic.
- The probabilistic error bound of \sqrt{N} does not benefit from regularity of the integrand function.
- The requirement that the points be independent random samples are wasteful because of clustering.
- In reality, pseudorandom numbers generated by completely deterministic means are used.
- Monte Carlo simulation exhibits a great sensitivity on the seed of the pseudorandom-number generator.

Matrix Computation

To set up a philosophy against physics is rash;
philosophers who have done so
have always ended in disaster.
— Bertrand Russell

Definitions and Basic Results

- Let $A \equiv [a_{ij}]_{1 \leq i \leq m, 1 \leq j \leq n}$, or simply $A \in \mathbf{R}^{m \times n}$, denote an $m \times n$ matrix.
- It can also be represented as $[a_1, a_2, \dots, a_n]$ where $a_i \in \mathbf{R}^m$ are vectors.
 - Vectors are column vectors unless stated otherwise.
- A is a square matrix when $m = n$.
- The rank of a matrix is the largest number of linearly independent columns.

Definitions and Basic Results (continued)

- A square matrix A is said to be symmetric if $A^T = A$.
- A real $n \times n$ matrix

$$A \equiv [a_{ij}]_{i,j}$$

is diagonally dominant if $|a_{ii}| > \sum_{j \neq i} |a_{ij}|$ for $1 \leq i \leq n$.

- Such matrices are nonsingular.
- The identity matrix is the square matrix

$$I \equiv \text{diag}[1, 1, \dots, 1].$$

Definitions and Basic Results (concluded)

- A matrix has full column rank if its columns are linearly independent.
- A real symmetric matrix A is positive definite if

$$x^T A x = \sum_{i,j} a_{ij} x_i x_j > 0$$

for any nonzero vector x .

- A matrix A is positive definite if and only if there exists a matrix W such that $A = W^T W$ and W has full column rank.

Cholesky Decomposition

- Positive definite matrices can be factored as

$$A = LL^T,$$

called the Cholesky decomposition.

- Above, L is a lower triangular matrix.

Generation of Multivariate Distribution

- Let $\mathbf{x} \equiv [x_1, x_2, \dots, x_n]^T$ be a vector random variable with a positive definite covariance matrix C .
- As usual, assume $E[\mathbf{x}] = \mathbf{0}$.
- This covariance structure can be matched by $P\mathbf{y}$.
 - $C = PP^T$ is the Cholesky decomposition of C .^a
 - $\mathbf{y} \equiv [y_1, y_2, \dots, y_n]^T$ is a vector random variable with a covariance matrix equal to the identity matrix.

^aWhat if C is not positive definite? See Lai (R93942114) and Lyuu (2007).

Generation of Multivariate Normal Distribution

- Suppose we want to generate the multivariate normal distribution with a covariance matrix $C = PP^T$.
 - First, generate independent standard normal distributions y_1, y_2, \dots, y_n .
 - Then

$$P[y_1, y_2, \dots, y_n]^T$$

has the desired distribution.

- These steps can then be repeated.

Multivariate Derivatives Pricing

- Generating the multivariate normal distribution is essential for the Monte Carlo pricing of multivariate derivatives (pp. 648ff).
- For example, the rainbow option on k assets has payoff

$$\max(\max(S_1, S_2, \dots, S_k) - X, 0)$$

at maturity.

- The closed-form formula is a multi-dimensional integral.^a

^aJohnson (1987); Chen (D95723006) and Lyuu (2009).

Multivariate Derivatives Pricing (concluded)

- Suppose $dS_j/S_j = r dt + \sigma_j dW_j$, $1 \leq j \leq k$, where C is the correlation matrix for dW_1, dW_2, \dots, dW_k .
- Let $C = PP^T$.
- Let ξ consist of k independent random variables from $N(0, 1)$.
- Let $\xi' = P\xi$.
- Similar to Eq. (82) on p. 690,

$$S_{i+1} = S_i e^{(r - \sigma_j^2/2) \Delta t + \sigma_j \sqrt{\Delta t} \xi'_j}, \quad 1 \leq j \leq k.$$