Level 9 HW: Part C

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Part C: MC 101

Question C.a

Study the source code in the file TestMC.cpp and relate it to the theory that we have just discussed.

```
namespace SDEDefinition
{ // Defines drift + diffusion + data
  OptionData* data;
                            // The data for the option MC
  double drift(double t, double X)
  { // Drift term
    return (data → r)*X; // r - D
  }
  double diffusion(double t, double X)
  { // Diffusion term
    double betaCEV = 1.0;
    return data→sig * pow(X, betaCEV);
  }
  double diffusionDerivative(double t, double X)
  { // Diffusion term, needed for the Milstein method
    double betaCEV = 1.0;
    return 0.5 * (data→sig) * (betaCEV) * pow(X, 2.0 * betaCEV - 1.0);
  }
} // End of namespace
```

The above code parameterizes the SDE in Equation (6) and Equation (9) which shows that that the drift and diffusion coefficients are proportional to the current value, \bar{x} .

$$dX = \alpha X dt + b X dW$$

Now the code moves on to the main function.

```
std::cout << "1 factor MC with explicit Euler\n";
OptionData myOption;
myOption.K = 65.0;
myOption.T = 0.25;
myOption.r = 0.08;
myOption.sig = 0.3;
myOption.type = -1; // Put -1, Call +1
double S_0 = 60;

long N = 100;
std::cout << "Number of subintervals in time: ";
std::cin >> N;
```

First, parameters are set, including the option contract and the quantity of subintervals. N, needed to be created.

```
// Create the basic SDE (Context class)
Range<double> range (0.0, myOption.T);
double VOId = S_0;
double VNew;
std::vector<double> x = range.mesh(N);
```

Then, the grid over time is represented by range, which is an object of class $\frac{Range < double>}{Range < double>}$. Then N+1 grid points are generated by code, $\frac{Range < double>}{Range < double>}$. Which results in a vector of doubles.

$$0 = t_0 < t_1, \dots < t_n < t_{n+1} < \dots < t_N = T$$

```
// V2 mediator stuff
long NSim = 50000;
std::cout << "Number of simulations: ";
std::cin >> NSim;

double k = myOption.T / double (N);
double sqrk = sqrt(k);

// Normal random number
double dW;
double price = 0.0; // Option price

// NormalGenerator is a base class
NormalGenerator* myNormal = new BoostNormal();

using namespace SDEDefinition;
SDEDefinition::data = &myOption;
```

```
std::vector<double> res;
int coun = 0; // Number of times S hits origin
```

Next, parameters are defined for the Monte Carlo simulation and useful expressions are precomputed such as double sqrk = sqrt(k). The user is asked to input the number of simulations. Variables that will be used inside the scope of the MC loop are initialized, such as std::vector < double > res, which stores the results of the MC simulates (the value that is to the averaged over), and a counter variable for the number of times the underlying stock price hits 0. If I remember correctly, it's possible for S to hit 0, but only once. It is an observing state. If we instead had the CIR model the variance would go to S as S making this less likely to occur (maybe impossible).

```
for (long i = 1; i \le NSim; ++i)
{ // Calculate a path at each iteration
  if ((i/10000) * 10000 == i)
  {// Give status after each 1000th iteration
       std::cout << i << std::endl;
  }
  VOId = S_0;
  for (unsigned long index = 1; index < x.size(); ++index)
    // Create a random number
    dW = myNormal → getNormal();
    // The FDM (in this case explicit Euler)
    VNew = VOld + (k * drift(x[index-1], VOld))
            + (sqrk * diffusion(x[index-1], VOId) * dW);
    VOId = VNew;
    // Spurious values
    if (VNew <= 0.0) coun++;
  }
  double tmp = myOption.myPayOffFunction(VNew);
  price += (tmp)/double(NSim);
}
// D. Finally, discounting the average price
price *= exp(-myOption.r * myOption.T);
// Cleanup; V2 use scoped pointer
delete myNormal;
```

Above is the main loop for implementing the Monte Carlo (MC) pricing of a European Call. The outer loop, for (long i = 1; i <= NSim; ++i), computes the MC price NSIM times. Each simulated results in accumulating the MC price divided

by a normalizing factor NSIM so the resulting sum at the end of the loop is the average price, price += (tmp)/double(NSim) .

The inner-loop is more interesting as it mimics the structure of the option. The idea is to simulate the stock price forward to period T. Because there is no early-exercise we only need to consider the payout of the option in period T.

To increment forward by one time-step we make a draw from the normal distribution to match ΔW_n in Equation (10). We substitute the integral of the continuous path of the Wiener process with a draw from the Normal distribution, which exactly captures its distribution qualities, $dW = myNormal \rightarrow getNormal()$.

The we compute the new value of the underlying vnew based the Euler-Maruyama method, void,

```
VNew = VOId + (k * drift(x[index-1], VOId)) + (sqrk * diffusion(x[index-1], VOId) * dW) .
```

Finally, we set void to vNew for the next iteration of the for loop, and increment the counter that tracks times the underlying hit 0.

Once the inner-loop is finished we compute the payoff and increment price as discussed previously.

At the end of the simulation, we need to discount the price back to time t_0 to get the final price of the call, price *= exp(-myOption.r * myOption.T) .

```
// Cleanup; V2 use scoped pointer delete myNormal;

std::cout << "Price, after discounting: " << price << ", " << std::endl; std::cout << "Number of times origin is hit: " << coun << endl;
```

Lastly, we delete the object myNormal from memory, and display the results of the MC simulation to the user.

Question C.b

Run the MC program again with data from Batches 1 and 2. Experiment with different values of NT (time steps) and NSIM (simulations or draws). In particular, how many time steps and draws do you need in order to get the same accuracy as the exact solution? How is the accuracy affected by different values for NT / NSIM?

Batch 1: Accuracy of European Call: $T=0.25,\,K=65,\,\sigma=0.30,\,r=0.08,\,S=60$

Reference Call Price = 2.13337

Table values are absolute errors.

NT , NSIM	NSIM =10	NIM =100	NSIM = 1_000	NSIM =10,000	NIM =100_000	NSIM =1_000_000
NT = 5	2.024499e-01	6.718065e-01	1.043850e-02	8.194448e-03	3.693774e-02	3.394879e-02
NT = 10	5.047594e-02	2.518777e-01	5.335607e-02	1.040406e-02	2.220257e-02	1.662575e-02
NT = 25	8.388749e-01	4.280889e-02	1.820541e-01	6.395586e-02	3.201728e-02	3.427768e-03
NT = 50	1.077734e+00	1.086020e-01	8.705175e-02	6.530664e-02	2.685270e-02	1.377353e-04
NT = 75	6.487481e-02	2.116941e-01	5.879898e-02	1.245244e-02	5.169615e-03	1.694676e-03

NT = 100	6.434635e-01	3.844738e-02	5.927214e-02	4.429913e-03	2.944562e-03	6.577634e-04
NT = 1000	8.123001e-01	4.187805e-01	3.704392e-02	1.746732e-03	1.128285e-02	1.189853e-03

Batch 2: Accuracy of European Call: T=1.0, K=100, $\sigma=0.30$, r=0.08, S=100

Reference Call Price = 7.96557

Table values are absolute errors.

NT , NSIM	NSIM =10	NIM =100	NSIM = 1_000	NSIM =10,000	NIM =100_000	NSIM =1_000_000
NT = 5	1.144529e-01	1.987845e+00	1.403461e-02	7.290750e-02	8.822713e-03	1.128689e-02
NT = 10	6.253628e-01	4.771328e-01	9.580447e-02	1.511568e-02	4.269981e-03	3.653877e-03
NT = 25	2.818773e+00	1.705325e-01	3.679201e-01	9.967205e-02	7.558557e-02	9.635184e-03
NT = 50	4.200774e+00	9.950853e-02	8.227195e-02	1.805510e-01	6.259688e-02	1.120600e-02
NT = 75	3.492576e-02	1.886736e-01	2.200147e-01	8.501766e-03	1.018476e-03	1.100765e-03
NT = 100	1.589411e+00	1.904438e-01	2.299962e-01	2.460466e-02	2.195357e-02	6.577634e-04
NT = 1000	2.613153e+00	9.024101e-01	9.392745e-02	4.331281e-02	1.627487e-02	6.111349e-01

Conclusion: I need something like NT=75 and NSIM=100000 to get close to the accuracy of the closed-form solution

Question C.c.

Now we do some stress-testing of the MC method. Take Batch 4. What values do we need to assign to NT and NSIM in order to get an accuracy to two places behind the decimal point? How is the accuracy affected by different values for NT/NSIM?

Ok, so I ran NT=1.000 and and NSIM=100_000_000 , and the resulting pricing error was 3.23e-01, so I think it suffice to say this batch would need an extreme number of simulations to reach 1e-02.

```
Simulation number: 99930000
Simulation number: 99940000
Simulation number: 99960000
Simulation number: 99970000
Simulation number: 99980000
Simulation number: 99980000
Simulation number: 100000000
Pricing Error: 3.230219e-01
Number of times origin is hit: 0
```

I would think based on what I observed NT = 1_000 and NSIM = 1_000_000_000 should get close 1e-02 , as $10^{-\frac{1}{2}}\approx 3.23\times 10^{-1}$.

Level 9 HW: Part D

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Date: 4/2/2025

Part D: Advanced Monte Carlo

Question D.a

Create generic functions to compute the standard deviation and standard error based on the above formulae. The inputs are a vector of size M (M = NSIM), the interest-free rate and expiry time T. Integrate this newcode into TestMC.cpp. Make sure that the code compiles.

I implemented two functions. One generic function computes the standard deviation, <code>cmptSD()</code>, and one computes the standard error, <code>cmptSE()</code>, which calls <code>cmptSD()</code> using the specified inputs. These functions are generic in that they allow for different numeric types, <code>NumType</code>.

```
// Global function for computing std. dev.
template<typename NumType>
NumType cmptSD(vector<NumType> draws, NumType r, NumType T)
{
    // This function computes the standard deviation of the Monte Carlo
    // pricing algorithm based on a vector of draws, the risk-free rate,
    // and time-to-expiry.

// Extract M
    unsigned M = draws.size();

// Compute sum_C := \sum_j C_{T,j}, and sum_sq_C := \sum_j C_{T,j}^2
NumType sum_sq_draws = 0;
NumType sum_draws = 0;
for (unsigned i = 0; i < M; ++i)
{
    sum_draws += draws[i];
    sum_sq_draws += draws[i] * draws[i];
}</pre>
```

```
return sqrt((sum_sq_draws - (pow(sum_draws, 2) / M)) / (M - 1)) * exp(-r * T);
}

// Global function for computing standard error
template<typename NumType>
NumType cmptSE(vector<NumType> draws, NumType r, NumType T)

{
    // This function computes the standard error of the Monte Carlo
    // pricing algorithm based on a vector of draws, the risk-free rate,
    // and time-to-expiry.

// Extract M
    unsigned M = draws.size();

// Compute standard deviation
    NumType std_dev = cmptSD(draws, r, T);

return std_dev / sqrt(M);
}
```

Additionally I needed to change part of the main loop to first compute and assign the payoff at time T, or equivalently $C_{T,j}$, assign it to variable payoff, store that in a vector of payoffs cs to then pass at the end of the loop to the aforementioned functions.

```
// Initalize the vector of Calls
vector<double> Cs(NSim);

// Main MC loop
for (long i = 1; i <= NSim; ++i) {
    if (i % 10000 == 0) {
        std::cout << "Simulation number: " << i << std::endl;
    }

VOId = S_0;
for (size_t index = 1; index < x.size(); ++index) {
    dW = myNormal → getNormal();

VNew = VOId + (k * drift(x[index - 1], VOId))
        + (sqrk * diffusion(x[index - 1], VOId) * dW);

VOId = VNew;</pre>
```

```
if (VNew <= 0.0) coun++;
}

payoff = myOption.myPayOffFunction(VNew);
Cs.push_back(payoff); // Push C_{T,j} to vector
price += payoff / double(NSim);
}

price *= exp(-myOption.r * myOption.T);</pre>
```

Question D.b

Run the MC program again with data from Batches 1 and 2. Experiment with different values of NT (time steps) and NSIM (simulations or draws). How do SD and SE react for these different run parameters, and is there any pattern in regards to the accuracy of the MC (when compared to the exact method)?

Batch 1: Accuracy of European Call:
$$T=0.25$$
, $K=65$, $\sigma=0.30$, $r=0.08$, $S=60$

Reference Call Price = 2.13337

Table 1: The **absolute error** of the MC simulation across grid points (NT) and number of simulations (NSIM) for Batch 1

NT , NSIM	NSIM =100	NSIM =1,000	NSIM =10,000	NSIM =100,000	NSIM =1,000,000
NT = 2	2.632417e-01	9.203648e-02	1.200790e-01	8.510446e-02	8.340218e-02
NT = 3	5.035559e-01	3.235333e-01	1.224027e-01	7.813884e-02	5.147937e-02
NT = 4	5.987387e-01	9.473616e-02	1.460946e-02	3.346980e-02	3.233861e-02
NT = 5	5.466716e-01	7.456874e-02	9.855942e-03	3.847193e-02	1.890097e-02
NT = 6	1.061755e+00	1.753611e-01	5.521700e-02	1.889628e-02	1.844404e-02
NT = 7	7.457166e-01	2.418215e-01	1.316571e-01	3.636494e-02	1.242936e-02
NT = 8	2.669690e-01	2.127838e-01	1.708850e-02	2.627710e-02	2.159024e-02
NT = 9	6.084229e-02	5.314298e-02	2.094094e-02	5.702243e-03	1.804855e-02
NT = 10	3.212444e-01	1.475987e-01	7.430848e-04	1.402769e-02	2.317850e-02
NT = 100	4.266600e-01	1.503501e-01	5.603340e-03	1.537361e-02	1.758639e-03
NT = 1,000	4.071797e-01	2.320302e-02	3.167530e-02	1.191878e-02	4.454782e-03

Table 2: The **standard deviation** of the MC simulation across grid points (NT) and number of simulations (NSIM) for Batch 1

NT , NSIM	NSIM =100	NSIM =1,000	NSIM =10,000	NSIM =100,000	NSIM =1,000,000
NT = 2	3.702759e+00	4.205938e+00	4.134713e+00	4.149745e+00	4.170402e+00
NT = 3	4.891625e+00	3.732387e+00	4.185014e+00	4.262213e+00	4.294542e+00
NT = 4	3.517116e+00	4.434011e+00	4.459829e+00	4.364515e+00	4.353021e+00
NT = 5	5.118550e+00	4.516054e+00	4.395748e+00	4.365230e+00	4.393537e+00
NT = 6	2.414266e+00	4.509912e+00	4.366411e+00	4.410379e+00	4.415072e+00
NT = 7	4.981660e+00	4.830431e+00	4.227897e+00	4.394543e+00	4.430881e+00
NT = 8	5.199024e+00	4.446559e+00	4.464114e+00	4.411672e+00	4.424697e+00
NT = 9	4.213768e+00	4.579448e+00	4.453109e+00	4.476104e+00	4.443539e+00
NT = 10	4.057807e+00	4.294339e+00	4.455329e+00	4.449267e+00	4.435539e+00
NT = 100	3.731941e+00	4.548375e+00	4.430099e+00	4.492470e+00	4.515502e+00
NT = 1,000	4.623455e+00	4.741437e+00	4.470627e+00	4.502637e+00	4.516729e+00

Table 3: The **standard errors** of the MC simulation across grid points (NT) and number of simulations (NSIM) for Batch 1.

NT , NSIM	NSIM =100	NSIM =1,000	NSIM =10,000	NSIM =100,000	NSIM =1,000,000
NT = 2	3.702759e-01	1.330034e-01	4.134713e-02	1.312265e-02	4.170402e-03
NT = 3	4.891625e-01	1.180285e-01	4.185014e-02	1.347830e-02	4.294542e-03
NT = 4	3.517116e-01	1.402157e-01	4.459829e-02	1.380181e-02	4.353021e-03
NT = 5	5.118550e-01	1.428102e-01	4.395748e-02	1.380407e-02	4.393537e-03
NT = 6	2.414266e-01	1.426159e-01	4.366411e-02	1.394684e-02	4.415072e-03
NT = 7	4.981660e-01	1.527516e-01	4.227897e-02	1.389677e-02	4.430881e-03
NT = 8	5.199024e-01	1.406125e-01	4.464114e-02	1.395093e-02	4.424697e-03
NT = 9	4.213768e-01	1.448149e-01	4.453109e-02	1.415468e-02	4.443539e-03
NT = 10	4.057807e-01	1.357989e-01	4.455329e-02	1.406982e-02	4.435539e-03
NT = 100	3.731941e-01	1.438322e-01	4.430099e-02	1.420644e-02	4.515502e-03
NT = 1,000	4.623455e-01	1.499374e-01	4.470627e-02	1.423859e-02	4.516729e-03

Batch 2: Accuracy of European Call:
$$T=1.0$$
, $K=100$, $\sigma=0.20$, $r=0$, $S=100$

Reference Call Price = 7.96557

Table 3: The **abs error** of the MC simulation across grid points (NT) and number of simulations (NSIM) for Batch 2.

	NSIM =100	NSIM =1,000	NSIM =10,000	NSIM =100,000	NSIM =1,000,000
NT = 2	3.485391e-01	1.202511e-02	6.247254e-02	2.999940e-02	2.521862e-02
NT = 3	1.540124e+00	7.635447e-01	1.544779e-01	4.850612e-02	3.448957e-02
NT = 4	1.722926e+00	4.451495e-01	2.034288e-01	3.689535e-02	4.353653e-02
NT = 5	1.804811e+00	1.569835e-01	8.584167e-02	1.678313e-02	5.679862e-02
NT = 6	3.267576e+00	6.546431e-01	1.049989e-01	3.365332e-02	3.851598e-02
NT = 7	2.364024e+00	6.423566e-01	3.295264e-01	2.332860e-02	4.026738e-02
NT = 8	4.974667e-01	4.276399e-01	1.248284e-01	7.118566e-03	1.174670e-02
NT = 9	5.017132e-01	1.037566e-01	1.213223e-01	6.574340e-02	9.758878e-03
NT = 10	9.372743e-01	2.957527e-01	5.235354e-02	1.579178e-02	1.312259e-02
NT = 100	1.204283e+00	3.692603e-01	1.715665e-02	3.035763e-02	7.707332e-03
NT = 1000	1.055245e+00	9.063880e-02	7.579229e-02	2.864089e-02	1.631538e-02

Table 5: The **standard deviation** of the MC simulation across grid points (NT) and number of simulations (NSIM) for Batch 2.

	NSIM =100	NSIM =1,000	NSIM =10,000	NSIM =100,000	NSIM =1,000,000
NT = 2	1.133303e+01	1.242954e+01	1.226964e+01	1.232857e+01	1.237038e+01
NT = 3	1.416083e+01	1.139988e+01	1.239817e+01	1.257496e+01	1.265163e+01
NT = 4	1.082217e+01	1.300745e+01	1.300898e+01	1.281460e+01	1.278670e+01
NT = 5	1.455594e+01	1.309936e+01	1.288798e+01	1.281749e+01	1.288042e+01
NT = 6	8.381529e+00	1.319100e+01	1.281277e+01	1.291724e+01	1.292694e+01
NT = 7	1.441636e+01	1.394607e+01	1.249895e+01	1.287796e+01	1.296443e+01
NT = 8	1.477896e+01	1.288287e+01	1.303940e+01	1.291955e+01	1.294673e+01
NT = 9	1.248044e+01	1.331490e+01	1.303436e+01	1.307194e+01	1.298788e+01
NT = 10	1.205227e+01	1.261498e+01	1.302500e+01	1.300213e+01	1.297001e+01
NT = 100	1.135694e+01	1.313115e+01	1.295705e+01	1.309141e+01	1.314985e+01
NT = 1000	1.362096e+01	1.362321e+01	1.303816e+01	1.311199e+01	1.315371e+01

Table 6: The **standard errors** of the MC simulation across grid points (NT) and number of simulations (NSIM) for Batch 2.

NT , NSIM	NSIM =100	NSIM =1,000	NSIM =10,000	NSIM =100,000	NSIM =1,000,000
NT = 2	1.133303e+00	3.930564e-01	1.226964e-01	3.898636e-02	1.237038e-02
NT = 3	1.416083e+00	3.604959e-01	1.239817e-01	3.976550e-02	1.265163e-02
NT = 4	1.082217e+00	4.113318e-01	1.300898e-01	4.052333e-02	1.278670e-02
NT = 5	1.455594e+00	4.142380e-01	1.288798e-01	4.053246e-02	1.288042e-02
NT = 6	8.381529e-01	4.171362e-01	1.281277e-01	4.084791e-02	1.292694e-02

NT = 7	1.441636e+00	4.410136e-01	1.249895e-01	4.072369e-02	1.296443e-02
NT = 8	1.477896e+00	4.073921e-01	1.303940e-01	4.085522e-02	1.294673e-02
NT = 9	1.248044e+00	4.210541e-01	1.303436e-01	4.133711e-02	1.298788e-02
NT = 10	1.205227e+00	3.989208e-01	1.302500e-01	4.111634e-02	1.297001e-02
NT = 100	1.135694e+00	4.152434e-01	1.295705e-01	4.139866e-02	1.314985e-02
NT = 1000	1.362096e+00	4.308036e-01	1.303816e-01	4.146374e-02	1.315371e-02

Conclusion

Looking at Table 1 and Table 4, which displays the **pricing error** in absolute terms we see both increases in grid size and number of simulations tends to decrease error. While increasing the number of simulations does not bias the estimate, we are much more likely to get a smaller draw from the distribution of abs. err. when the number of simulations is smaller. Increasing the number of grid points does tend to result in lower bias (smaller error), but the results are minimal. Even with NT=2 we get a good approximation to prices with very large number of simulations.

Looking at Table 2 and Table 5 which shows the **standard deviation** of MC prices we no discernable pattern with the number of grid points or simulations. This is to be expected. The distribution of prices itself should converge to a stable distribution. But as we see in Table 3 and Table 6 which shows the **standard error** (volatility around the mean estimate) we see that the standard error decreases by a factor of roughly \sqrt{NSIM} as we increase the order of magnitude of NSIM.