

Laboratory: Stochastic Differential Equations

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Quantum Computing in Finance

Stochastic Differential Equations (SDEs) are extensively used in financial engineering. They are used as models to describe the future behaviour of financial assets, such as the price of a stock, interest rates, or the likelihood of bankruptcy. In this lecture we will study Geometric Brownian Motion (GBM), which is one of the most common types of SDEs.

In particular, we will:

- Describe GBM and its solution via the lognormal distribution
- Look at how GBM the classical approaches to GBM in terms of analytical and numerical methods
- Understand how GBM can be modelled using the tools of quantum computing
- Finally, we will present a number of Python programs implementing the classical and quantum versions of GBM

DIS US \$

Market

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1x2

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DIS US Equity

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96 Actions

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98 Table

Line Chart

12/03/2012 -

12/02/2013

Total Retur

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11 Compare

Mov. Avgs

Volume

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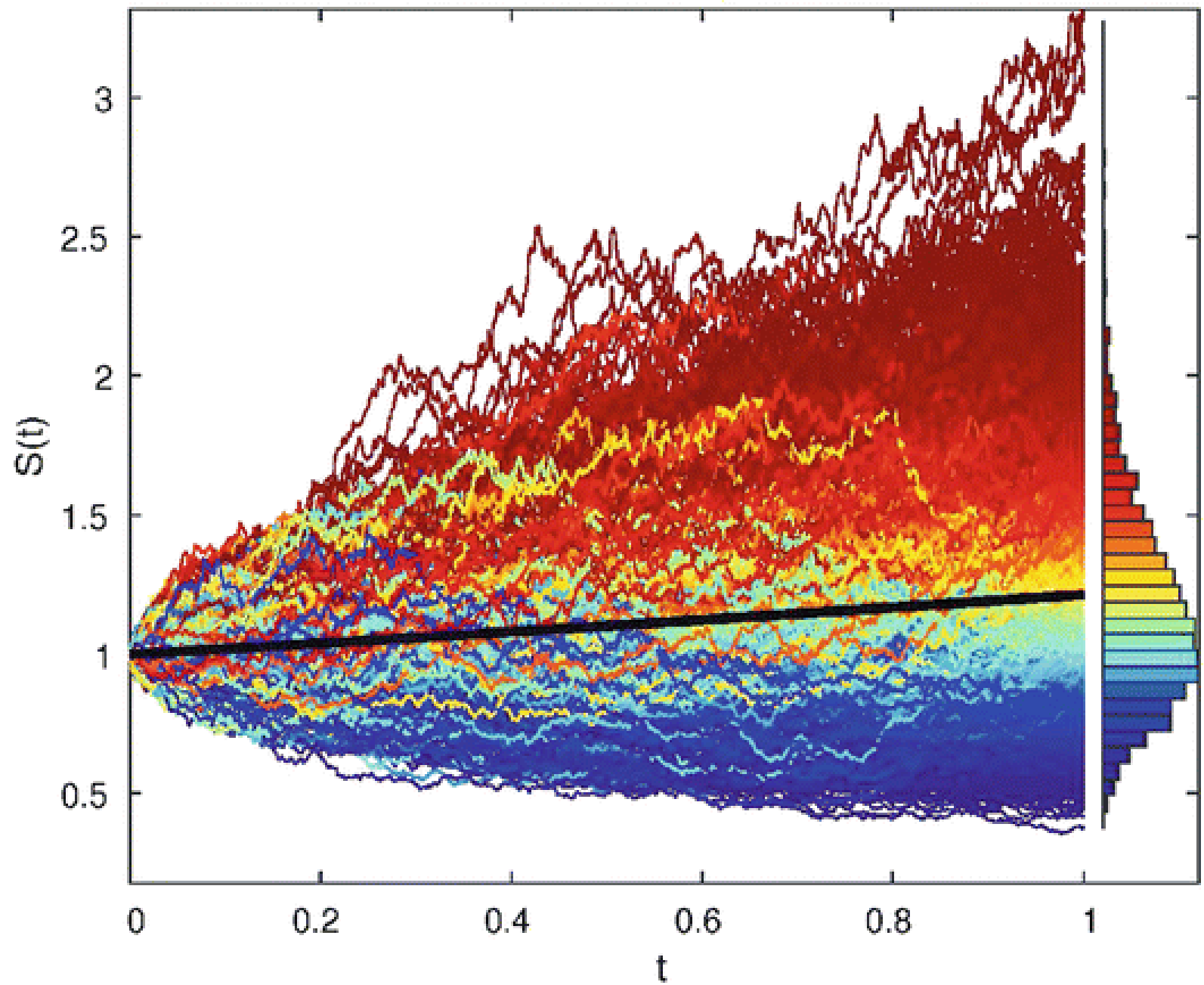
☐ Total Return Index (Gross Dividends) 71.6184
☐ High on 11/26/13 72.2682
☐ Average 61.9866
☐ Low on 12/03/12 49.29



Australia 61 2 9777 8600 Brazil 5511 3048 4500 Europe 44 20 7330 7500 Germany 49 69 9204 1210 Hong Kong 852 2977 6000
 Japan 81 3 3201 8900 Singapore 65 6212 1000 U.S. 1 212 318 2000

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Geometric Brownian Motion: 5000 sample paths, $\mu = 0.15$, $\sigma = 0.3$



The Problem: Geometric Brownian Motion

Geometric Brownian motion is a type of stochastic process.

Stochastic processes are ideal mathematical tools to describe the behavior of financial variables. Remember watching the graphs of a stock index such as the Dow Jones or the EUR/GBP exchange rates.

What do they have in common?

Clearly their evolution is not smooth.

Financial graphs are full of ups and downs, and broken features, dramatic changes. In short they are what mathematicians call **fractals**.

And how we can model fractals?

Well, using stochastic differential equations. It was the **French mathematician Bachelier** who first introduced fractals via Brownian motion as a model to capture this “broken” behavior observed in the financial markets.

This was in 1901 and largely forgotten. It was not until the late 20th century that the model was revised and it became the basis of such important financial theories such as the pricing of financial derivatives. In fact, GBM is at the heart of the **Black Scholes** option pricing model of 1973.

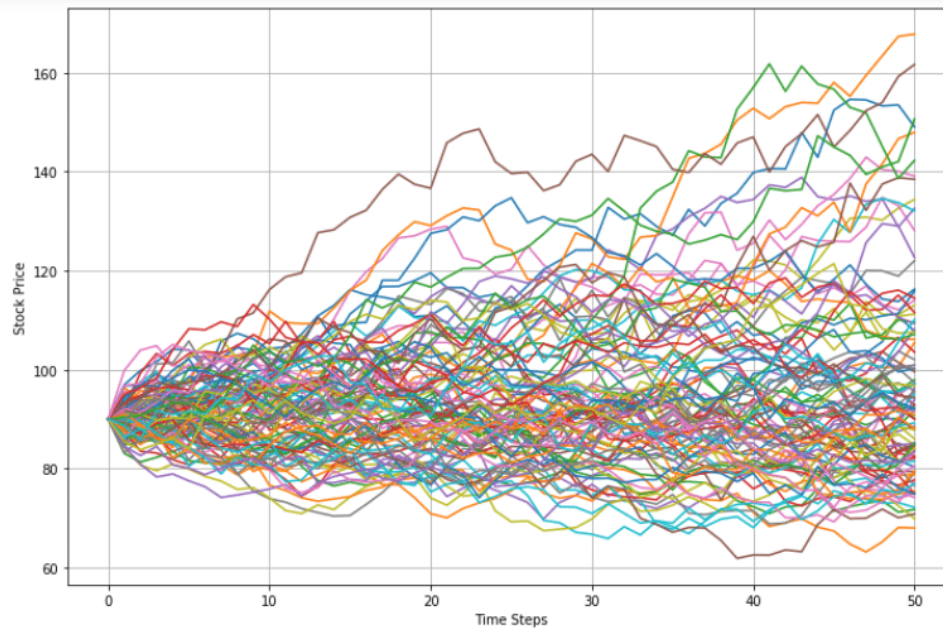


Even though stochastic differential equations in general, or GBM in particular, may seem esoteric this is very far from the truth.

As we will showing this lecture, it is very easy to construct a simple stochastic process in a computer program using random numbers generators and Monte Carlo simulation.

Our point of departure is then the GBM stochastic process of the form:

$$dS = rSdt + vSdW$$



We are going to use GBM as a way to **project into the future** information we have of today.

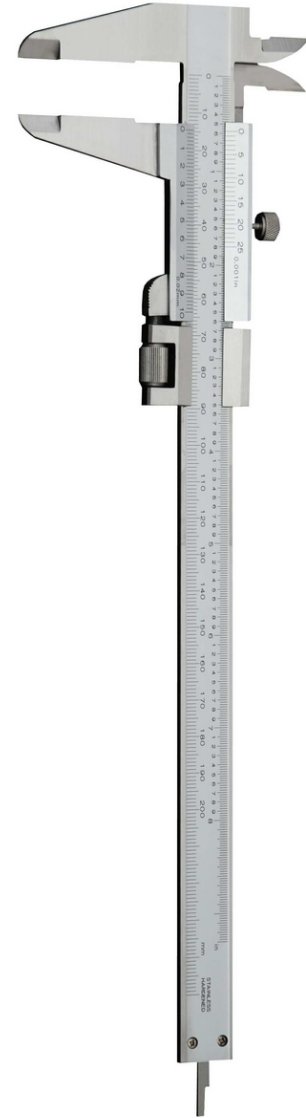
For example if we have the value of a stock price today, **we will use GBM** to generate many possible ways in which this stock price can evolve into the future. We can say that these are different possible futures or scenarios that the stock can take in the next week, month, or years.

How do we use GBM in practice?

Well we would need to solve it. In the case of the standard plain GBM, we are fortunate to have an analytical solution. This is very important for us at this stage because we will use it as a benchmark.

However in practice, we rarely have the luxury to have analytical solutions. The vast majority of financial calculations are done using numerical approximations.

This could include Monte Carlo simulation, like we will describe in this book, but also other types of numerical techniques such as binomial trees, partial differential equations, machine learning, and others.





Important: in financial engineering, we may use different models to describe stocks, interest rates or forex rates, but what is common is that our main interest is to find the expected value of a statistical distribution. This is what we are going to do in this chapter.

In the classical approach we will do this with Monte Carlo simulation which converges slowly to this expected value.

Quantum computing solution that we will show does the same thing, i.e. find the expected value of a statistical distribution, but it does it much faster – in fact with quadratic improvement.

The Classical Solution

Solution 1: analytical solution

GBM (equation 1) can be shown to have the solution

$$f(S; \mu, \sigma) = \frac{1}{S\sigma\sqrt{2\pi}} \exp \left[-\frac{(\log(S) - \mu)^2}{2\sigma^2} \right]$$

Which depends on parameters μ and σ , whilst S is the independent variable (which can represent stock price for example). The above equation can be shown to be the probability density function of the lognormal distribution,

$$f(S; \mu, \sigma) = PDF_{LOGNORM}(S; \mu, \sigma)$$

The mapping between the variables used in the GBM (r, v) and lognormal distribution (μ, σ) is

$$\mu = \left[\left(r - \frac{1}{2}v^2 \right) T + \log(S_0) \right]$$

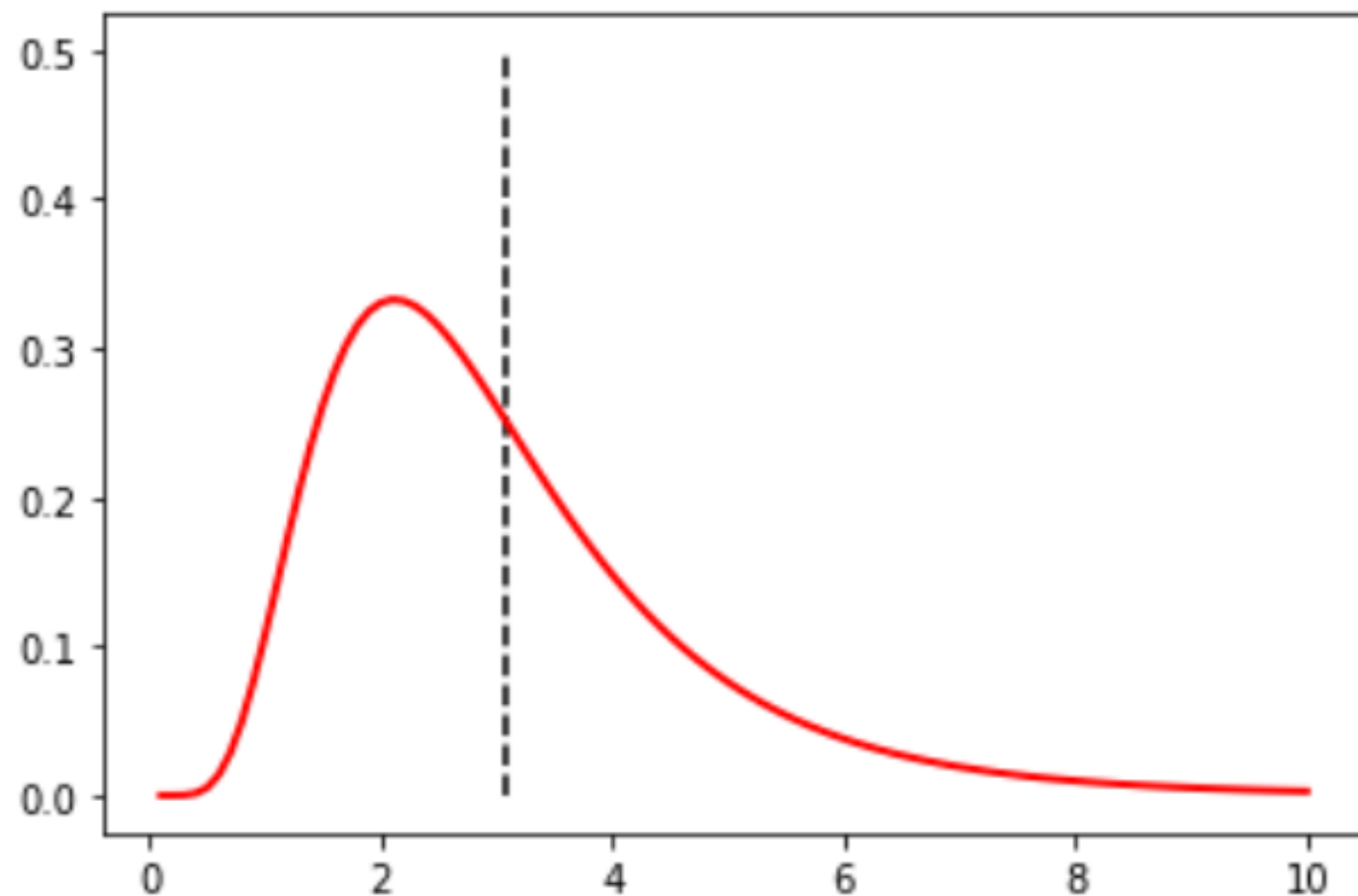
$$\sigma = v\sqrt{T}$$

Finally, the expected value, or the mean of the lognormal distribution can be computed as:

$$\bar{S} = E[S] = \exp\left(\mu + \frac{1}{2}\sigma^2\right)$$



```
# CODE 4.1 Lognormal distribution PDF graph
import numpy as np
import matplotlib.pyplot as plt
```

```
_____ parameters _____  
mu = 1.0  sigma = 0.5
```

```
_____ results _____  
exact_mean = 3.080216848918031
```

Solution 2: numerical solution

The numerical solution of GBM can be obtained by discretizing GBM as:

$$dS = rSdt + vSdW$$

$$\Delta S = rS\Delta t + vS\Delta W$$

$$S^{i+1} - S^i = rS^i\Delta t + vS^i\varepsilon_i\sqrt{\Delta t}$$

$$S^{i+1} = S^i + rS^i\Delta t + vS^i\varepsilon_i\sqrt{\Delta t}$$

$$S^{i+1} = S^i(1 + r\Delta t + v\varepsilon_i\sqrt{\Delta t})$$

Which can then be used step by step to move forward from time zero to time T. Once this process is repeated many times, we can determine the final value as the average of the ST generated.

$$S_T^j = S^N$$

$$\bar{S} = E[S] = \frac{1}{M} \sum_{j=1}^M S_T^j$$

The Composition

In both cases, analytical and numerical, we obtain an estimate of the underlying variable of interest at some future date T . As mentioned before, we are also interested to know what is the expected value of the contract and that future date. This means that using each individual estimate of the underlying, we will then evaluate the contract using a payoff function H , and then take the average. This will result in the expected price of the contract and not only the expected price of the underlying.

In the example we use in this chapter, the payoff H will be represented by a simple linear function. This has the form α times X plus β .

$$H = H(x; \text{parameters})$$

$$H(x; \alpha, \beta) = \alpha x + \beta$$

$$H(x; 1, 0) = x$$

As before we can compute the expected value as the average of the terminal values for the estimates of the underlying

$$\bar{S} = E[S_T] = \frac{1}{M} \sum_{j=1}^M S_T^j$$

And we can do the same thing for the expected payoff:

$$\bar{H} = E[H(S_T)] = \frac{1}{M} \sum_{j=1}^M H(S_T^j)$$

Python Lab

LABORATORY 1: Classical Analytical Solution

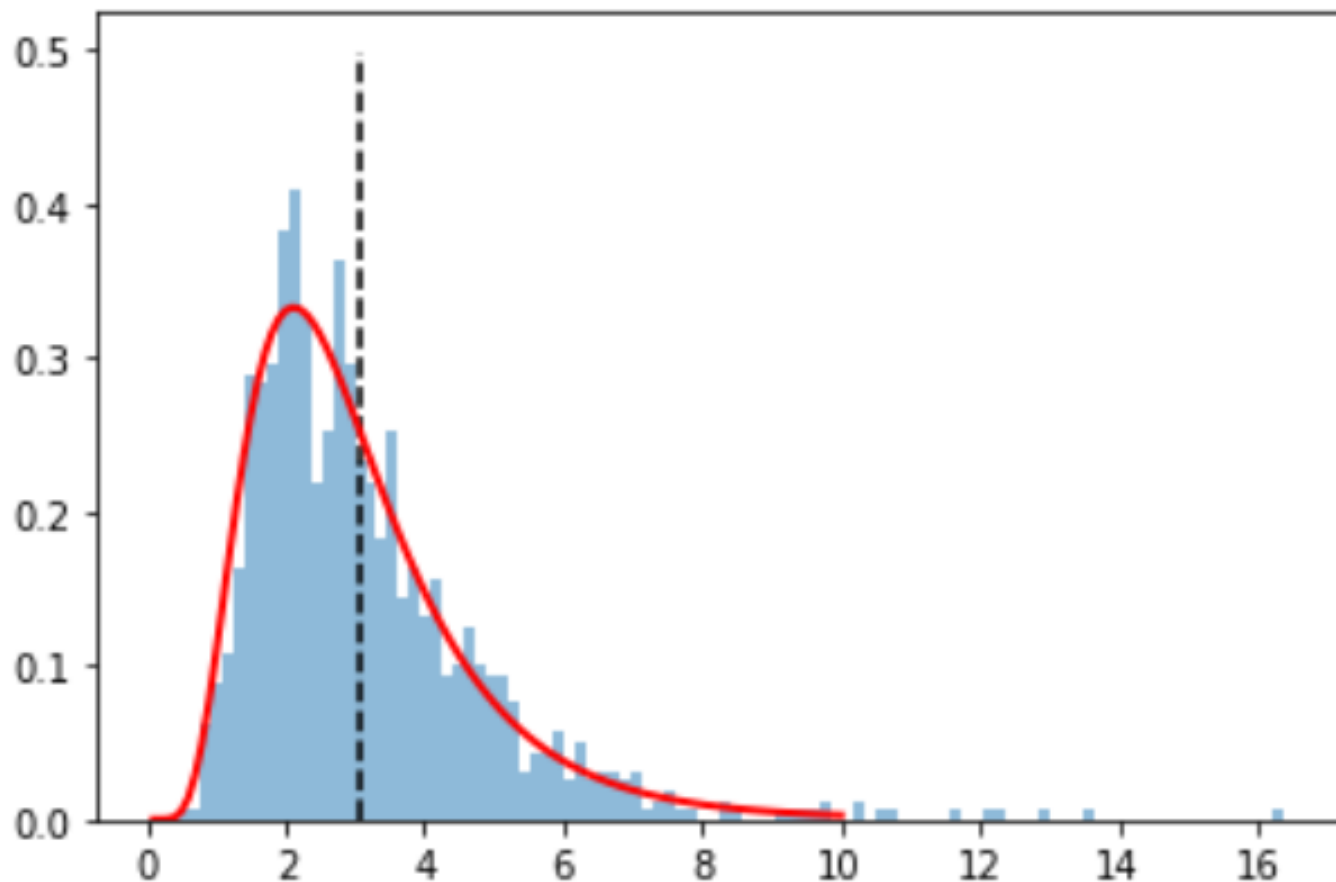
This code illustrates the analytical methods to obtain the estimated mean value of the Lognormal distribution. We have used parameters $\mu=1$ and $\sigma=0.5$. The technique is based on the sampling of distribution based on the Python function `xxxx`. We have sampled the distribution $M=100$ times and then simply taking the average of samples. As can be seen, the exact meaning of this tuition is 3.0802, which compares with the estimated value of the Monte Carlo sampling of 3.1801.

The graph illustrates an overlay of the lognormal probability density function in red, the histogram of the samples, and the exact meaning of distribution indicated by the dashed vertical line.

Code 4.2 Classical Analytical Solution

In the following code, we present the use of method `lognormal()` from module `random` to generate random numbers from a lognormal distribution.

```
# CODE 4.2 Lognormal distribution MC sampling
import numpy as np
import matplotlib.pyplot as plt
```

_____ parameters _____
mu = 1.0 sigma = 0.5

_____ results _____
Exact_mean = 3.080216848918031
MC_sample_mean = 3.18018716824946

LABORATORY 2: Numerical Solution

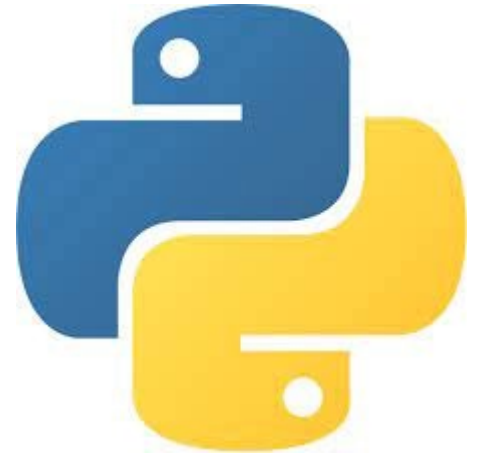
In this laboratory we ought to follow the trajectory of the underlying variable up to some future time T . Each path represents a possible scenario or evolution of the underlying price into the future. The code that we present here implements the iterated process described in the numerical solution section described before.

This is based in both random number generation and Monte Carlo simulation. The code is composed of three parts:

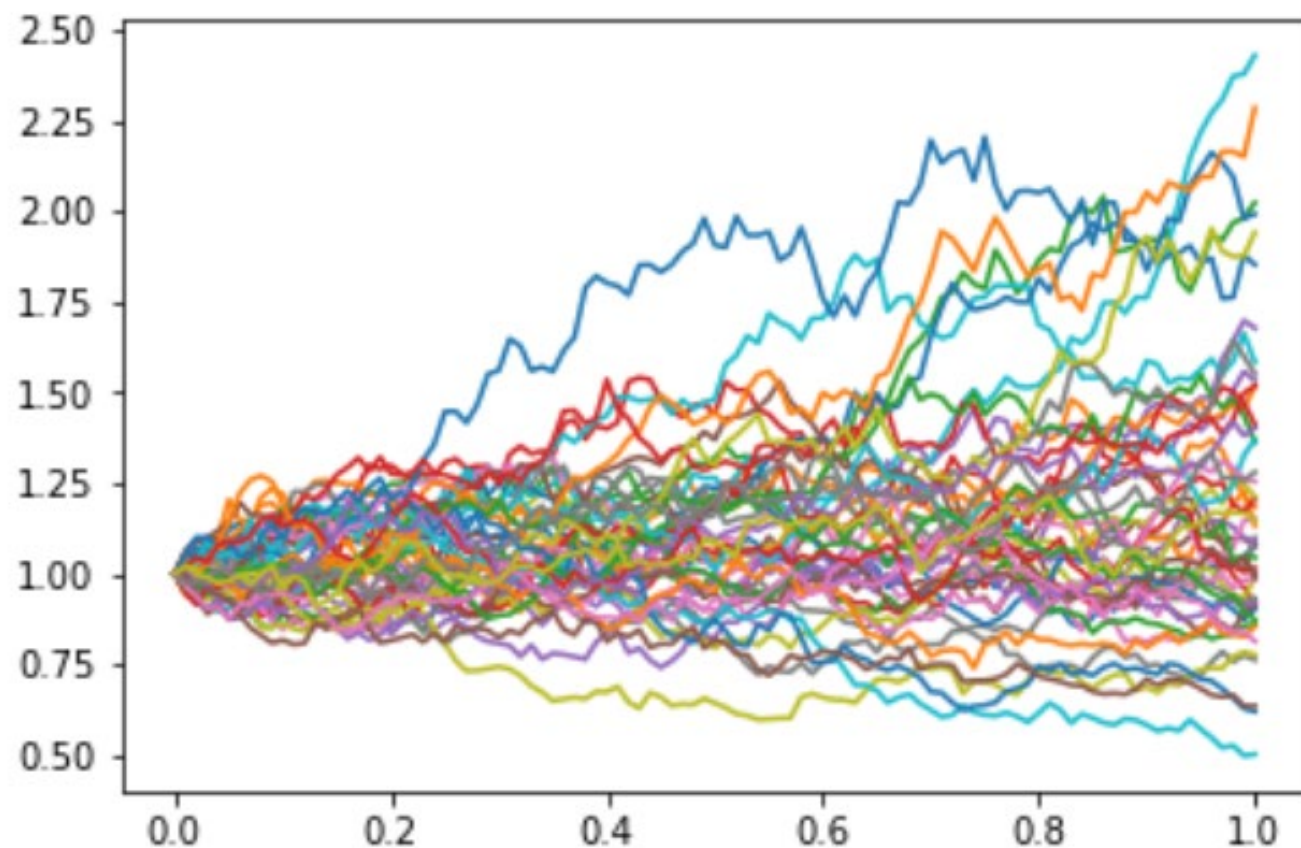
The first Part, the necessary libraries are loaded and the parameters of the problem are specified.

The second part represents the iterated process that is required to advance from time $T = 0$ to T . And the the M repetitions required. These are captured using a double for loop.

The third and final part is where the average value is computed and the results presented in graphical form.



```
# CODE 4.3 Geometric Brownian Motion with Monte Carlo  
import math  
import scipy as sp  
import numpy as np  
import matplotlib.pyplot as plt
```



parameters

$S_0 = 1$ $\text{vol} = 0.3$ $r = 0.1$ $T = 1$
 $N = 100$ $M = 1000$
 $\mu = 0.0550000000000000001$ $\sigma = 0.3$

results

Exact_mean = 1.1051709180756477
MC_ST_mean = 1.1043202409665795

Code 4.4 Classical Numerical Solution: distribution

We can explore in more detail the behavior of the underlying at maturity with plot a histogram of the values at this time. In the next code we again generate multiple trajectories of the evolution of the underlying variable, but this time we collect the values of the underlying at maturity and generates an histogram of them. This is illustrated in the next code and figure.

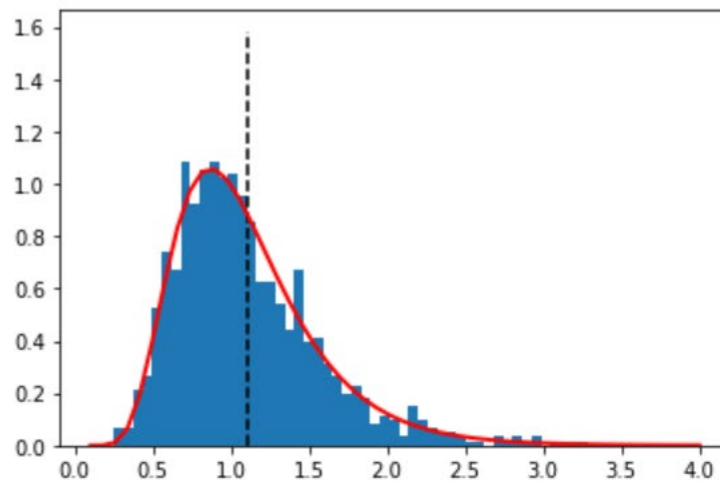
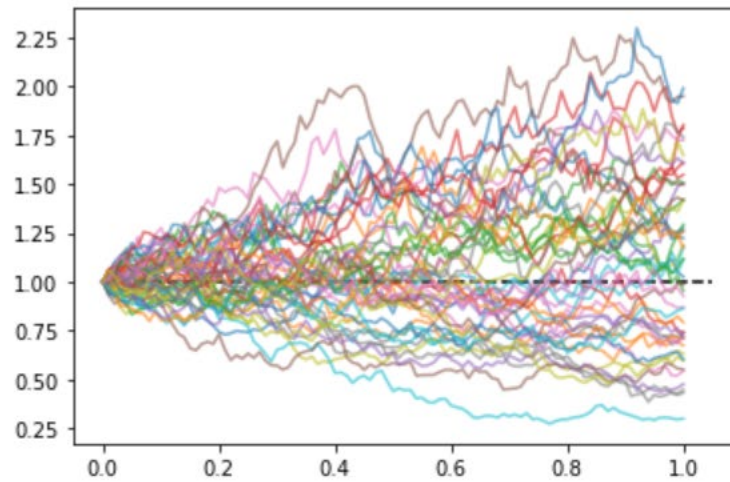
```
# CODE 4.4 GBM PATH MC SAMPLING

import math

import scipy as sp

import numpy as np

import matplotlib.pyplot as plt
```



parameters

$S_0 = 1$ $\text{vol} = 0.4$ $r = 0.1$ $T = 1$
 $N = 100$ $M = 1000$
 $\mu = 0.019999999999999999$ $\sigma = 0.4$

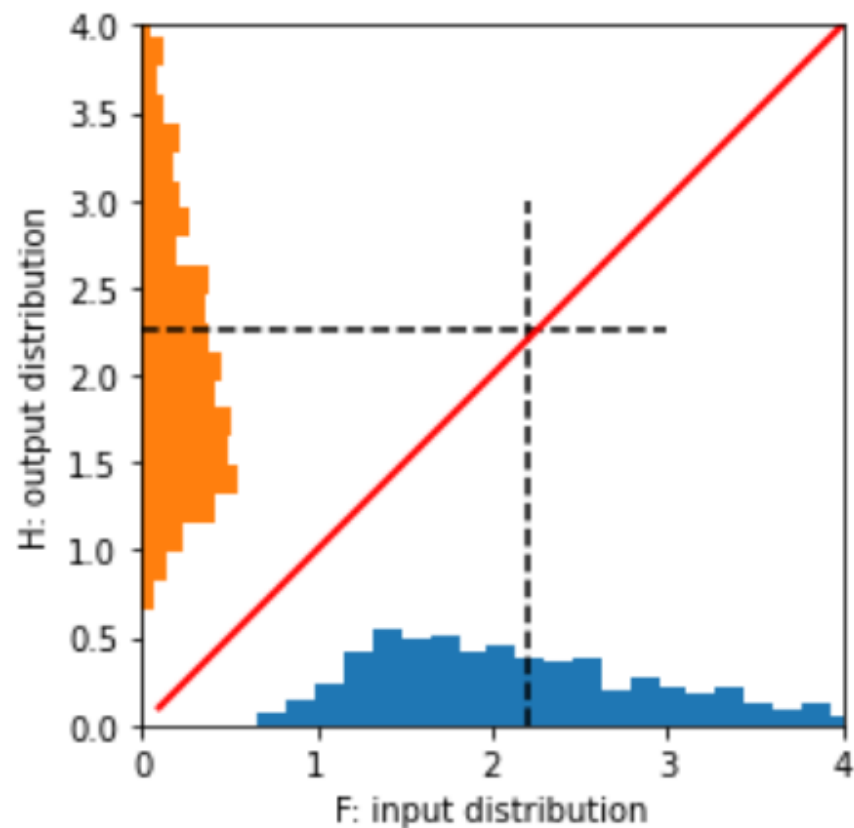
results

Exact_mean = 1.1051709180756477
 MC_sample_mean = 1.1012865878720308



Code 4.5 Classical Numerical Solution: composition

To study the behavior the payoff on the predicted values of the underlying using this technique, we can explore this in the following Python code. In this code we generate as in the previous one distribution of potential values for the underlying, and subsequently for each of them, we apply the payoff H . Once we have obtained M predicted values of the payoff based on M predicted values of the underlying, we can then compute the expected payoff simply by taking an arithmetic average



parameters

$S_0 = 2$ $\text{vol} = 0.4$ $r = 0.1$ $T = 1$
 $N = 100$ $M = 1000$
 $\mu = 0.7131471805599453$ $\sigma = 0.4$

results

Exact mean LOGNORMAL = 2.210341836151295
 MC estimate $E[H(ST)]$ = 2.2559784667535405

The Quantum Solution

The Quantum Computing Solution

How can we approach the problem of computing the expected value of the statistical distribution using quantum computing?

The quantum solution relies on a very different approach than the classical solution. Quantum solution is going to be based on the Grover's search algorithm and its generalization, the so-called quantum amplitude estimation. On the one hand this are advanced tools in quantum computing, but on the other they are easy to use.

Fundamentally the classical and the quantum solutions are similar. Conceptually they are trying to do the same: estimate the expected value of the mean of a statistical distribution. However, the approach the solution in very different ways.

Whilst the classical solution samples statistical distribution using random numbers and therefore constructs a set of samples from which we can then calculate statistical properties, such as the mean. In contrast, the quantum solution relies on a search algorithm, the Grover's search algorithm, which when implemented in quantum computers turns out to be extremely efficient.

In terms of computational efficiency, it has been shown that the classical Monte Carlo sampling of in a statistical solution is proportional to the square root of the number of samples taken. On the contrary, Montanaro has shown that quantum amplitude estimation technique converges according to the inverse of the number of samples done. This represents a quadratic improvement, which is extremely significant.

Qiskit Lab

LABORATORY 3: Quantum Uncertainty Model

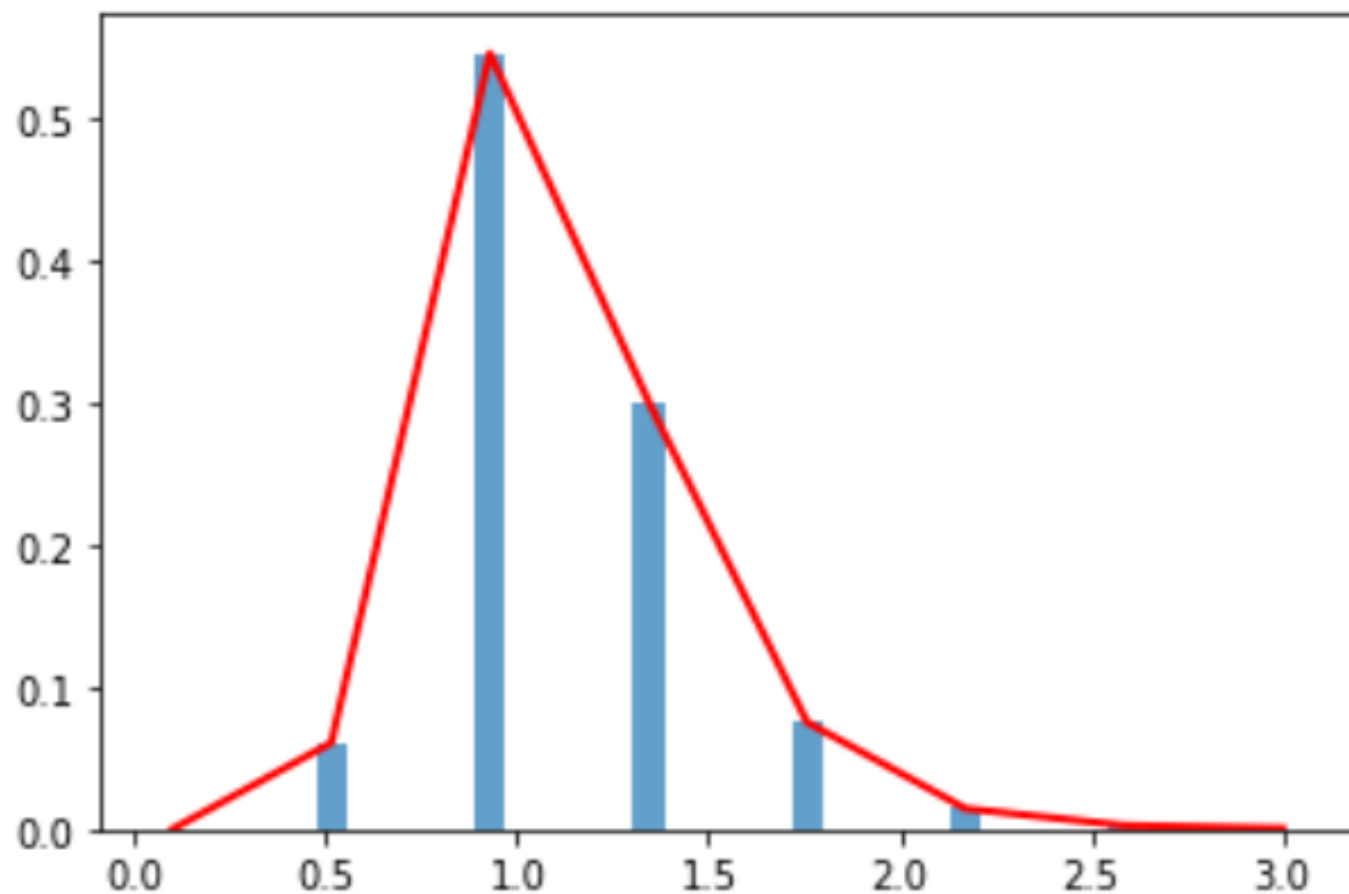
In order to capture the lognormal distribution in qiskit, we will use the function `uncertainty model` with a specified number of qubits. This function allows for the description of not only the lognormal distribution, but also of the normal distribution and the uniform distribution. This flexibility is quite useful in financial calculations. It allows us to describe the statistical behaviour of some underlying variable. Whilst in some cases a live demonstration will be an adequate, like in stock price modelling, in other cases a normal distribution will be more appropriate, such as some types of interest rate models, as will be discussed in a future.

In the accompanying Python code illustrates the amount distribution in qiskit using three qubits. The resultant number of combinations of grid points is 2 elevated to power of 3, and therefore eight.

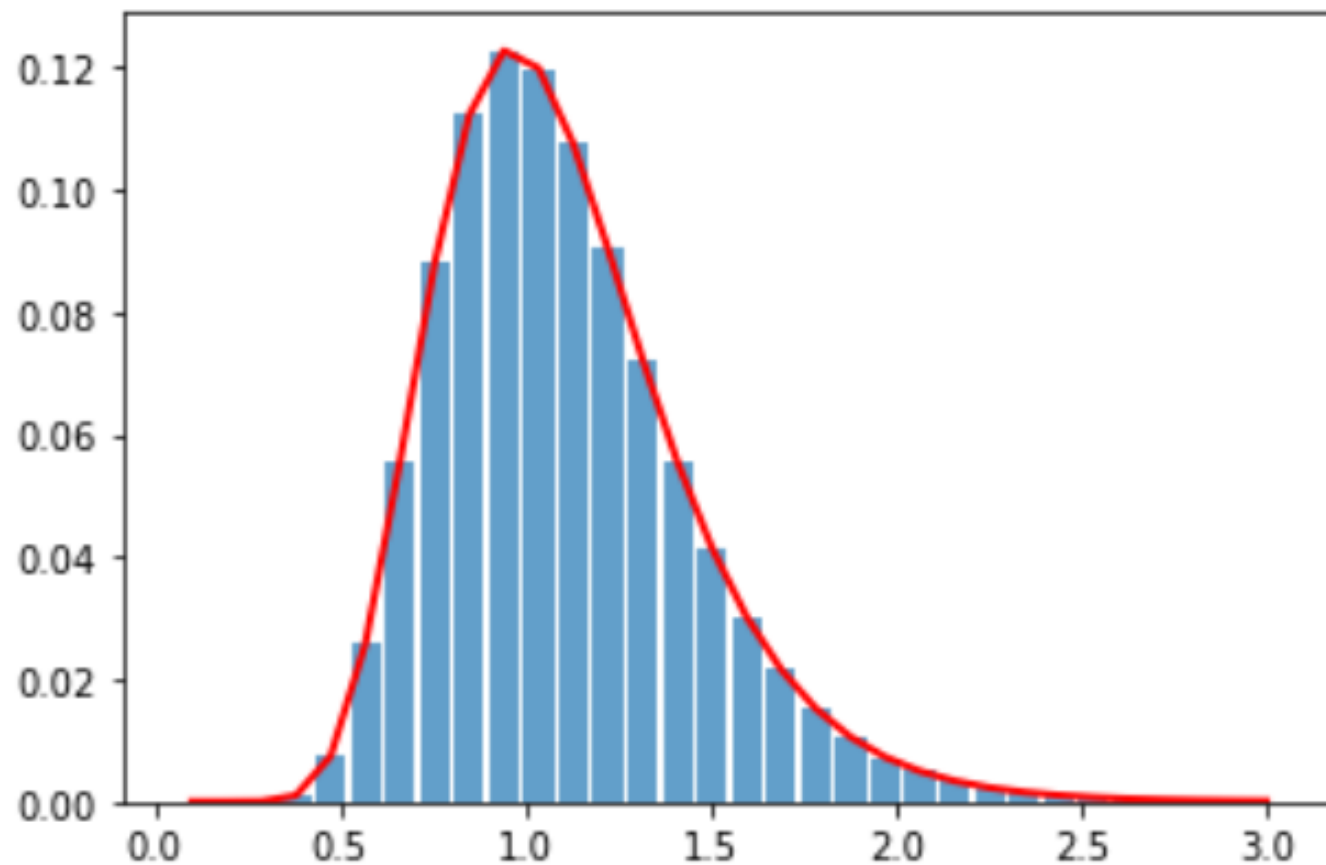


Code 4.6 Quantum LogNormal FEW QUBITS

```
# CODE_4_6_QUANTUM_LOGNORMAL_FEW
import matplotlib.pyplot as plt
import numpy as np
from qiskit import Aer, QuantumCircuit
```



```
_____ parameters _____  
S0= 1  vol= 0.3  r= 0.1  T= 1  
mu = 0.055000000000000001  sigma = 0.3  
_____ quantum LOGN _____  
number qubits   = 3  
resulting grid  = 8
```



```
_____ parameters _____  
S0= 1  vol= 0.3  r= 0.1  T= 1  
mu =  0.055000000000000001  sigma =  0.3  
_____ quantum LOGN _____  
number qubits  =  5  
resulting grid =  32  
_____
```


References

The theory of stochastic processes as used in finance is extremely complex. I invite Grinko, D., Gacon, J., Zoufal, C. et al. Iterative quantum amplitude estimation. npj Quantum Inf 7, 52 (2021). <https://doi.org/10.1038/s41534-021-00379-1>

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