

# On the Numerical simulation of the Heston model

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School of Applied Mathematics  
A Thesis  
Submitted

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In Partial Fulfillment of the  
Requirements for the Degree  
of M.Sc. of Applied Mathematics  
at the FGV

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setembro 22, 2017



Approved for the Division  
(Mathematics)

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# Acknowledgements

You get pseudo-order when you seek order; you only get a measure of order and control when you embrace randomness. — Nassim Nicholas Taleb

# Table of Contents

<b>Chapter 1: Introduction</b>	<b>1</b>
<b>Chapter 2: Theoretical Framework</b>	<b>3</b>
2.1 Stochastic Calculus	3
2.1.1 Stochastic differential equation - SDE	3
2.1.2 Brownian Motion	4
2.1.3 Itô's Integral	6
2.2 Black-Scholes Model	7
2.2.1 Basics	7
2.2.2 The model	9
2.2.3 Limitations	10
2.3 Stochastic Volatility models	10
2.3.1 Cox-Ingersoll-Ross model	11
2.3.2 Heston Model	11
2.3.3 Other Models	12
2.4 Numerical Methods	12
2.4.1 Convergence	13
2.4.2 Discretization	14
2.4.3 Stability	16
<b>Chapter 3: The Heston Model Implementation</b>	<b>21</b>
3.1 Characteristic Function	21
3.2 Euler Scheme	22
3.3 Kahl-Jackel	23
3.4 Exact Algorithm	24
3.4.1 Generate a sample of $V_t$ given $V_0$	24
3.4.2 Generate a sample of $\int_0^t V_s ds$ given $V_t, V_0$	25
3.4.3 Compute $\int_0^t \sqrt{V_s} dB_s$ given $V_t, V_0$ and $\int_0^t V_s ds$	27
3.4.4 Limitations	28
<b>Chapter 4: Results</b>	<b>29</b>
<b>Chapter 5: Conclusion</b>	<b>33</b>
<b>Appendix A: Black-Scholes formula</b>	<b>35</b>

<b>Appendix B: Bessel Function</b> . . . . .	<b>37</b>
<b>Appendix C: Implementations</b> . . . . .	<b>39</b>
C.1 Euler Scheme . . . . .	39
C.2 Kahl-Jackel Implementation . . . . .	40
C.3 Exact Algorithm . . . . .	40
C.4 Results with 20 steps . . . . .	43
<b>References</b> . . . . .	<b>45</b>



# List of Tables

2.1	Box calculus . . . . .	7
3.1	Truncation schemes . . . . .	22
4.1	Model Parameters . . . . .	29
4.2	Option Value . . . . .	30
4.3	Estimators . . . . .	32



# List of Figures

2.1	A Wiener process trajectory path example . . . . .	5
2.2	A GBM trajectory path example . . . . .	9
2.3	Volatility Smile . . . . .	11
2.4	Analytical x Euler solutions . . . . .	14
2.5	Graphical representation of Euler's formula . . . . .	15
2.6	Euler's stability whith different timesteps . . . . .	17
2.7	Stability domains . . . . .	19
4.1	Comparison between models, 20 steps. . . . .	30
4.2	Comparison between models, 100 steps. . . . .	31
B.1	Modified Bessel Functions of the First Kind . . . . .	37
C.1	Comparison between models, 20 steps. . . . .	43



# Abstract

In this thesis we revisit some of the numerical methods for solving the Heston model's European call. Specifically, we approach the Euler, the Kahl-Jackel and two versions of the exact algorithm schemes. To perform this task, firstly we present a literature review which brings stochastic calculus, the Black-Scholes (BS) model and its limitations, the stochastic volatility methods and why they resolve the issues of the BS model, and the peculiarities of the numerical methods - convergence, discretization and stability. Since it is impossible to have a deep approach to all these topics, we provide recommendations when we acknowledge that the reader might need more specifics. We introduce the methods aforementioned providing all our implementations in R language. Also, we deliver an R package with these functions and others.

**Keywords:** Heston, Stochastic, Volatility, Black-Scholes, European call, R



# Chapter 1

## Introduction

The french mathematician Louis Bachelier was the trail-blazer that brought Brownian motion, previously restricted to the field of botanics where it was firstly observed, to the financial framework. He modeled the stock prices as a Brownian motion with drift. In 1973, Black and Scholes [1] designed a model based on the Geometric Brownian Motion to price options.

Options are derivatives that give their bearers the rights to buy or sell a specific asset in a future date and with a predetermined price. They are, by design, affected by small variations in the underlying assets' components, for example, the variance.

The Black-Scholes model was once the standard way of option pricing, but was replaced by more modern models that are now prevalent. One of the main drawbacks of the Black-Scholes model is the strong assumption that the stock returns' volatility is constant. Thus, the implied model's volatility results in a flat surface when plotted against the option's strike price and maturity. Real world implied volatility varies with the strike price and maturity, forming what is called the 'volatility smile'.

The Heston model is an extension of the Black-Scholes model that tackles this volatility issue replacing the constant volatility with a stochastic process. There are many models that approach volatility stochastically, but the Heston has valuable characteristics such as presenting an analytical solution to the option pricing and also having a computationally simple implementation when compared to more sophisticated competitors.

We aim with this thesis to better understand the Heston model, its quirkinesses, associated numerical methods and their implementations, and eventually discuss its pros and cons. For such, we will not only discuss the challenges in a vaguely way, but show code snippets of our implementations in R programming language [2] to help clarify possible associated difficulties. Furthermore, since we didn't find code for the exact algorithm implementation we intend to release a CRAN package with the codes as a didactic tool.

This thesis is divided into five chapters, the first being this introduction. Following, we have a literature review that mainly addresses stochastic calculus, the Black-Scholes models and the Heston model. Thereafter we present the different Heston model implementations. Chapter 4 brings the results of the tests we perform and finally, chapter 5 presents the conclusion and work limitations.





# Chapter 2

## Theoretical Framework

This chapter presents the concepts of stochastic calculus, from the historic conception of how it first arose through the basic principles and applications in finance. We address with more care the classical Black-Scholes model and its limitations and the Heston model. This model is also well known, it brings the concept of stochastic volatility in it, which brings its results closer to reality.

### 2.1 Stochastic Calculus

Stochastic calculus arises from stochastic processes and allows the creation of a theory of integration where both the integrand and integrator terms are stochastic processes. Stochastic calculus was created by the Japanese mathematician Kiyosi Itô <sup>1</sup> in the 1940s and 1950s and is used for modeling financial options and in another wide variety of fields [3]. In this chapter we present the historical contexts in which the tools and models are used, but our focus is introducing the concepts and notations that will be further used in our work.

#### 2.1.1 Stochastic differential equation - SDE

At first, before introducing stochastic differential equation, it is helpful to start with ordinary differential equation. Let  $x_t$  denote an asset at time  $t$  so that the change in the asset at time  $t$  is given by the following deterministic differential equation:

$$\begin{aligned} dx_t &= f(t, x_t)dt \\ x(0) &= x_0 \end{aligned} \tag{2.1}$$

---

<sup>1</sup>There is another important stochastic integral, called the *Stratonovich Integral* that unlike the Itô's integral, respects the conventional calculus chain rule. Also, the integral is evaluated at the interval's midpoint, instead of its left extreme. A Stratonovich integral can be expressed as an Itô integral and vice versa.

We now add a “noise” to this equation:

$$\begin{aligned} dx_t &= \underbrace{f(t, x_t) dt}_{\text{drift}} + \underbrace{g(t, x_t) dW_t}_{\text{diffusion}} \\ x(0) &= x_0 \end{aligned} \tag{2.2}$$

This “noise”  $dW_t$  is a *random* Wiener process (which will be clarified below) and  $x_0$  is our initial value.

The  $g(t, x_t)$  part of the SDE is often referred as *diffusion coefficient*, these usually have a continuous paths. Before moving on, we must carefully define what the term *random* means and the best way to begin doing so is to precisely define a probability space:

**Definition 2.1.1** (Probability Space). A triple  $(\Omega, \mathcal{U}, \mathcal{P})$  is called a *probability space* provided  $\Omega$  is any set,  $\mathcal{U}$  is a  $\sigma$ -algebra of subsets of  $\Omega$  and  $\mathcal{P}$  is a probability measure on  $\mathcal{U}$ .

### 2.1.2 Brownian Motion

The Brownian motion is the name given to the irregular motion observed in the motion of pollen particles suspended in fluid resulting from particle collision with atoms or molecules. It is named after Robert Brown, the first to have observed the movement in 1828. He noted two characteristic in the pollen movement [3]:

- the path of a given particle is very irregular, having a tangent at no point
- the motion of two distinct particles appear to be independent

The first quantitative works in Brownian motion come from an interest in stock price fluctuation by Bachelier in 1900. Albert Einstein also leaned over the subject and in 1905 derived the transition density for Brownian motion from molecular-kinetic theory of heat [3,4].

In 1923, the Wiener process was coined in honor of Norbert Wiener mathematical proof of existence of the Brownian motion and stating its properties.<sup>2</sup>

**Definition 2.1.2** (Wiener Process). Given a probability space  $(\Omega, \mathcal{U}, \mathcal{P})$ , a stochastic process  $W_t$  defined in this space is a *Wiener process* if it satisfies the following properties:

- $W_0 = 0$
- The change in  $W$ , given by  $\Delta W = W_{t+\Delta t} - W_t$ , is normally distributed with mean zero and standard deviation  $\sqrt{\Delta t}$ , meaning that  $\Delta W = \epsilon \sqrt{\Delta t}$ , where  $\epsilon$  is  $N(0, 1)$ .

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<sup>2</sup>More can be found on [5–7].

- If the increment  $\Delta t_1$  does not overlap with the time increment  $\Delta t_2$ , then  $\Delta W_1$  and  $\Delta W_2$  are independent.
- The process is continuous, meaning that there are no jumps in the path.
- The process is Markovian. This means that the conditional expectation of  $W_{t+1}$  given its entire history is equal to the conditional expectation of  $W_{t+1}$  given today's information. This can be written as:  $E[W_{t+1}|W_1, \dots, W_t] = E[W_{t+1}|W_t]$ .
- Consider the time interval  $[0, t]$  with  $n$  equally spaced intervals given by  $t_i = \frac{it}{n}$ . Then the paths of the Brownian motion have unbounded variation, this means that they are not differentiable and go towards infinity as  $n$  increases. The quadratic variation is given by  $\sum_{i=1}^n (Z_{t_i} - Z_{t_{i-1}})^2 \rightarrow t$ , meaning that when  $n$  increases it stays constant at  $t$ .

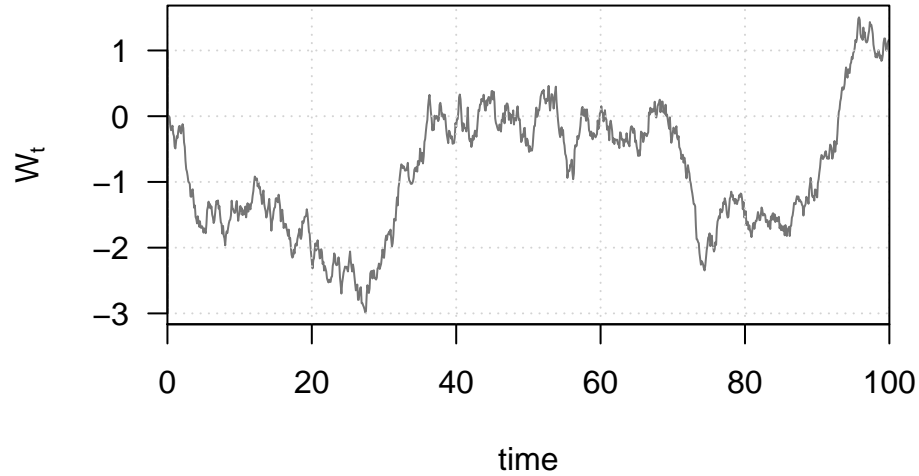


Figure 2.1: A Wiener process trajectory path example

### Correlated Brownian Motions

Two independent Brownian motions that are correlated can describe a new process  $Z_t$ . Let  $W_1$  and  $W_2$  be these two *independent* Brownian motions and let  $-1 \leq \rho \leq 1$  be a given number. For  $0 \leq t \leq T$  define the new process  $Z_t$  as [3]:

$$Z_t = \rho W_{1,t} + \sqrt{1 - \rho^2} W_{2,t} \quad (2.3)$$

This equation is a linear combination of independent normals at each timestep  $t$ , so  $Z_t$  is normally distributed. It is proven that  $Z$  is a Brownian motion and that  $Z$  and  $W_{1,t}$  have correlation  $\rho$  [3].

### 2.1.3 Itô's Integral

Formally, the SDE presented in equation (2.2) is interpreted as [6–11]:

$$x_t = x_0 + \int_0^t f(s, x_s) ds + \int_0^t g(s, x_s) dW_s \quad (2.4)$$

for some  $f(s, x_s)$ ,  $g(s, x_s)$  and  $s \in [0, t]$ .

The Itô integral can, as the Riemann integral, be approximated by a finite sum. Also, it has a definition as a certain limit. Itô's lemma 2.1.1 plays the same role as the fundamental theorem of calculus in allowing to evaluate integrals. It is the formal definition and presents an extra term not encountered in the conventional calculus theorem that is due to the non-smoothness characteristics of Brownian motion paths. It is possible, though, to define the integral in a less rigorous way:

$$Y_{\Delta t}(t) \approx \sum_{t_k < t} g(t_k) \Delta W_k \quad (2.5)$$

with the usual notions  $t_k = k\Delta t$ , and  $\Delta W_k = W(t_{k+1}) - W(t_k)$ . And in a more rigorous form, if the limit exists, then the Ito integral is:

$$Y(t) = \lim_{\Delta t \rightarrow 0} Y_{\Delta t}(t) \quad (2.6)$$

It is essential that the *forward difference* is used rather than the backward difference, which would be wrong.

**Lemma 2.1.1** (Itô's Lemma). *Assume that  $S_t$  has a stochastic differential given by:*

$$dS_t = \mu_t dt + \sigma_t dW_t \quad (2.7)$$

*for  $\mu_t$ ,  $\sigma_t$  and  $t \in [0, T]$ . Assume  $u : \mathbb{R} \times [0, T] \rightarrow \mathbb{R}$  is continuous and that  $\frac{\partial u}{\partial t}$ ,  $\frac{\partial u}{\partial x}$ ,  $\frac{\partial^2 u}{\partial x^2}$  exist and are continuous.*

$$Y_t := u(S_t, t)$$

*Then  $Y$  has the following stochastic differential:*

$$\begin{aligned} dY_t &= \frac{\partial u}{\partial t} dt + \frac{\partial u}{\partial x} dS_t + \frac{1}{2} \frac{\partial^2 u}{\partial x^2} \sigma_t^2 dt \\ &= \left( \frac{\partial u}{\partial t} + \mu_t \frac{\partial u}{\partial x} + \frac{1}{2} \frac{\partial^2 u}{\partial x^2} \sigma_t^2 \right) dt + \sigma_t \frac{\partial u}{\partial x} dW_t \end{aligned} \quad (2.8)$$

*where the argument of  $u$ ,  $\frac{\partial u}{\partial x}$  and  $\frac{\partial^2 u}{\partial x^2}$  above is  $(S_t, t)$ .*

Equation (2.8) is the stochastic equivalent to the chain rule, also known as Itô's formula or Itô's chain rule. The proof to this theorem is based on the Taylor expansion of the function  $f(S_t, t)$  [6,8]. For practical use you should write out a second-order Taylor expansion for the function to be analyzed and apply the multiplication table [3] presented in Table 2.1.

	$dt$	$dW_t$
$dt$	0	0
$dW_t$	0	$dt$

Table 2.1: Box calculus

### Itô's Integral Properties

Let  $f, g \in \mathcal{V}$  and let  $0 \leq t_0 < u < T$ . Then

$$(i) \int_{t_0}^T f dB_t = \int_{t_0}^u f dB_t + \int_u^T f dB_t$$

$$(ii) \int_{t_0}^T (\alpha f + \beta g) dB_t = \alpha \int_{t_0}^T f dB_t + \int_{t_0}^T \beta g dB_t$$

$$(iii) \mathbb{E} \left[ \int_{t_0}^T f dB_t \right] = 0$$

$$(iv) \mathbb{E} \left[ \left( \int_0^t H_s dB_s \right)^2 \right] = \mathbb{E} \left[ \int_0^t H_s^2 ds \right] \text{ (Isometry)}$$

$$(v) \mathbb{E} \left[ \int_{t_0}^T f dB_t \mid \mathcal{F}_s \right] = \int_{t_0}^s f dB_t, \quad \text{for } s < T. \text{ (Martingale)}^3$$

## 2.2 Black-Scholes Model

### 2.2.1 Basics

The Black-Scholes (B-S) model arises from the need to price european options in the derivative markets. Derivatives are financial instruments traded in the market, stock exchange or over-the-counter (OTC) market, whose values depend on the values of an underlying asset. [1,12,13]

- A call option is a derivative that gives its bearer the right, but not the obligation, to purchase a specific asset by a fixed price before or on a given date.
- A put option is a derivative that gives its bearer the right, but not the obligation, to sell a specific asset by a fixed price before or on a given date.

The trading price of the option is called the option *premium* and the asset from which the option derives is called the *underlying asset*. This asset may be interest rates, exchange rates, stock exchanges indices, commodities or stocks. The fixed price in

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<sup>3</sup>A martingale is a stochastic process with certain characteristics. The main one is that the expected value in time  $t + 1$  for  $X$  is the  $X$  value in  $t$ . This means there are no winning strategies when we are dealing with martingales (unlike when we play poker, for example). A Wiener process is a martingale.

contract in which the underlying asset might to be bought or sold is the *strick price*. The option expiration date is called the *maturity*. [1,13]

There are two major different option types: European and American. The difference between these two is that the bearer of the first may exercise it only at the end of its life, at its maturity while the latter can be exercised at any given time until its maturity. [1,14]

**Definition 2.2.1** (Intrinsic value). The intrinsic value of a call is the difference between the underlying asset price and the strike price. The put's intrinsic value operates the other way around, being the difference between the strike and the underling asset prices.

## Geometric Brownian Motion

A stochastic process  $S_t$  is a Geometric Brownian Motion<sup>4</sup> if it is described by the solution of the following stochastic differential equation [3,8,15].

$$dS_t = \mu S_t dt + \sigma S_t dW_t \quad (2.9)$$

for given constants  $\mu \in \mathbb{R}$  and  $\sigma > 0$ . Also, the assumed initial value is positive,  $S_0 > 0$ .

Figure 2.2 shows the GBM,<sup>5</sup> which is quite often applied to model the dynamics of some financial assets because of its properties [16]. Equation (2.10) shows the formula to generate a GBM and we provide proof of this solution in appendix A<sup>6</sup>

$$S_t = S_0 \times \exp\left(\left(\mu - \frac{\sigma^2}{2}\right)t + \sigma W_t\right), \quad t > 0 \quad (2.10)$$

---

<sup>4</sup>There is an Arithmetic Brownian Motion:  $dS_t = \mu dt + \sigma dB_t$ . More information can be obtained at [3].

<sup>5</sup>Also known as exponential Brownian motion.

<sup>6</sup>An intuitive proof can be found at [17].

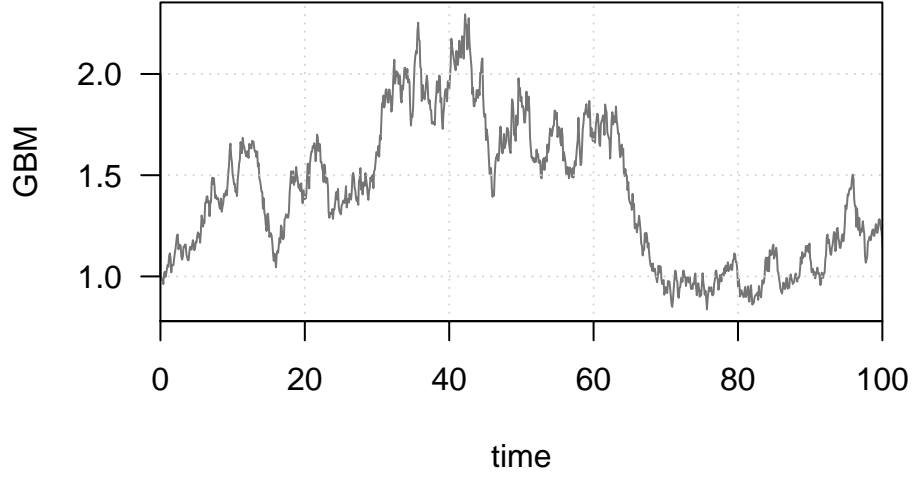


Figure 2.2: A GBM trajectory path example

### 2.2.2 The model

The Black-Scholes model provides analytical solution to the price of a European call at time  $t$  and can be described as follows [1,5,12]:

$$C(S_t, t) = N(d_1)S_t - N(d_2)Ke^{-r(T-t)} \quad (2.11)$$

$$d_1 = \frac{1}{\sigma\sqrt{T-t}} \left[ \ln\left(\frac{S_t}{K}\right) + \left(r + \frac{\sigma^2}{2}\right)(T-t) \right] \quad (2.12)$$

$$d_2 = d_1 - \sigma\sqrt{T-t} \quad (2.13)$$

Where:

- $S_t$  is the spot price of the underlying asset at time  $t$
- $r$  is the risk free rate (generally an annual rate)<sup>7</sup>
- $\sigma$  is the volatility of returns of the underlying asset <sup>8</sup>
- $N(\cdot)$  is the cumulative distribution function of the standard Gaussian distribution
- $K$  is the strike price
- $T - t$  is the time to maturity

Also, the stock price path is a Geometric Brownian Motion as previously stated, and is under the risk-neutral measure with the following dynamics [5,18]:

$$dS_t = (r - q)S_t dt + \sigma S_t dW_t \quad (2.14)$$

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<sup>7</sup>Assumed to be constant.

<sup>8</sup>See footnote 1.

Where  $dW_t$  is a Wiener process [1,18],  $r$  is the risk free rate and  $q$  is the dividend yield<sup>9</sup> and  $t$  denotes the current point in time.

### 2.2.3 Limitations

Although the Black-Scholes is very popular and the *de facto* standard in the market there are implications to the B-S model assumptions that affect the results and that are unrealistic. The main assumption that does not hold up is the deterministic (constant) volatility, that can more accurately be described as a stochastic process since we observe that small moves usually are followed by small moves and large moves by large moves. [5,12]

Other assumptions that are critical to the B-S model and are not always observed in practice refer to the asset's continuity through time (no jumps), being allowed to perform continuous hedge without transactions costs and normal (Gaussian) returns.

Most models focus on the volatility problem because transaction costs often translate to rises in volatility and fat-tails (abnormal) returns can be simulated by stochastic volatility and market or volatility jumps.

## 2.3 Stochastic Volatility models

Introducing stochastic volatility to models brings complexity, but enables modeling some features observed in reality that are crucial, like the random market volatility effects, skewness (market returns are more realistically modeled) and volatility smile<sup>10</sup> (see Figure 2.3). This kind of model is applied highly successfully in foreign exchange and credit markets.

**Definition 2.3.1** (Volatility Smile). Volatility smiles are implied volatility patterns that arise in pricing financial options. In particular for a given expiration, options whose strike price differs substantially from the underlying asset's price command higher prices (and thus implied volatilities) than what is suggested by standard option pricing models. These options are said to be either deep in-the-money or out-of-the-money.

Furthermore, stochastic volatility models use statistical methods as foundations to price and forecast options' behaviors and the underlying's security volatility is arbitrary. The Heston, the 3/2 and other models, like the GARCH<sup>11</sup> and SABR,<sup>12</sup> are considered standard smile models.

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<sup>9</sup> $r$  and  $q$  are assumed to be constant.

<sup>10</sup>The name derives from the concave shape of the graph, which resembles a smile.

<sup>11</sup>generalized autoregressive conditional heteroscedasticity.

<sup>12</sup>stochastic alpha, beta, rho.



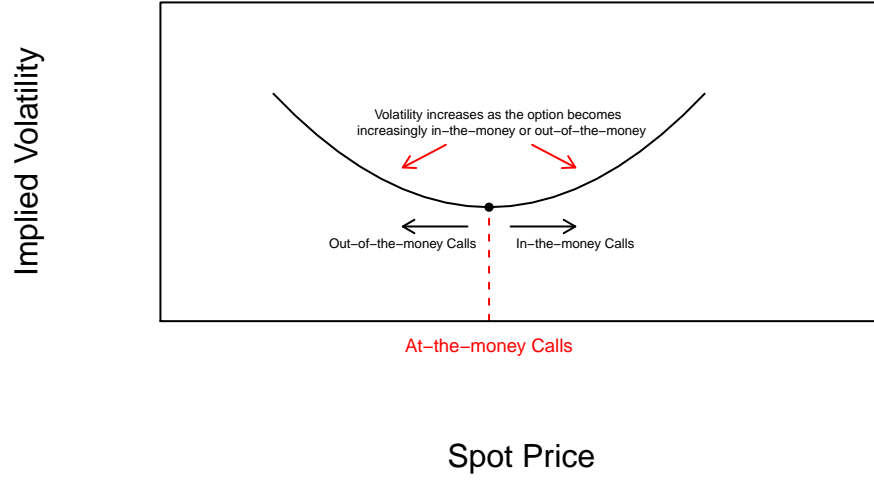


Figure 2.3: Volatility Smile

### 2.3.1 Cox-Ingersoll-Ross model

The Cox-Ingersoll-Ross (CIR) model is a well-known short-rate model that describes the interest rate movements driven by one source of market risk. The dynamics are described as follows[19,20]:

$$dr_t = k(\theta - r_t)dt + \sigma\sqrt{r_t}dB_t \quad (2.15)$$

Where,  $r_t$  is the short rate interest described by parameters  $\kappa$  the speed of mean reversion,  $\theta$  the long-run mean variance and  $\sigma$  the volatility of the variance process.

This model has been widely used to describe the dynamics of the short rate interest because it has some fundamental features like intuitive parametrization, nonnegativity and pricing formulas. Besides, it takes account of anticipations, risk aversion, investment alternatives and preferences about consumption timing and allows for detailed predictions about how changes in a wide range of underlying variables affect the term structure[19]. Furthermore, this equation constitutes one of the two Heston model equations with the volatility taking the short rate interest place.

### 2.3.2 Heston Model

Heston model was introduced in 1993 by Steven Heston to solve the deterministic volatility problems. It was designed to analyze bond and currency options and it introduced the following equations, which represent the dynamics of the stock price

and the variance processes under the risk-neutral measure [20,21]:

$$dS_t = \mu S_t dt + \sqrt{V_t} S_t dW_t^* \quad (2.16)$$

$$dV_t = \kappa(\theta - V_t)dt + \sigma\sqrt{V_t}dB_t \quad (2.17)$$

The second equation, as described in Section 2.3.1, is the CIR model equation. The first equation states the asset price process.  $\mu$  is the asset's rate of return,  $dW_{t,1}$  and  $dW_{t,2}$  are two correlated wiener processes with correlation coefficient of  $\rho$ .

### 2.3.3 Other Models

#### Ornstein-Uhlenbeck

The Ornstein-Uhlenbeck is the earliest recorded SDE. Named after Leonard Ornstein and George Eugene Uhlenbeck, it is a stochastic process that describes the acceleration of a pollen particle in a liquid subject to bombardments by molecules [3]. As we can observe in equation (2.18),  $x_t$  represents the one dimension velocity of the particle, thus  $dx_t$  is the *change* in velocity, in other words, its acceleration. The  $-\theta x_t$  component slows down the acceleration and is to be understood as frictional force. Besides, we add a noise  $W_t$  with intensity  $\sigma$  that models the random bombardment by the molecules.

$$dx_t = -\theta x_t dt + \sigma dW_t \quad (2.18)$$

With  $\theta$  and  $\sigma$  being positive constants. Expressing in terms of  $x_t$  we get:

$$x_t = e^{-\theta t} \times \left[ x_0 + \sigma \int_{t=0}^T e^{\theta t} dW_s \right]. \quad (2.19)$$

#### Langevin

The Langevin equation describes a system that consists of the molecular bombardment of a speck of dust on a water surface. We know that the intensity of the bombardement does not depend on the state variables [22,23].

$$m \frac{dv}{dt} = -\zeta v + \delta F(t) \quad (2.20)$$

$m$  is the mass of the particle,  $v$  it's velocity,  $-\zeta v$  is the frictional force, which is proportional to the velocity, and  $\delta F(t)$  is a *fluctuating* force (random) to the frictional force.

## 2.4 Numerical Methods

Numerical methods are tools that are often applied to solve stochastic differential equations because most of these do not have explicit solution. This means that we are not able to solve these equations using symbolic computation. Although we are

unable to find an analytical solution, when facing real problems, the approximation given by a numerical method is often sufficient. Alongside the analytical issue, the need to calculate the SDE's trajectory through time is the main reason why studying numerical methods is so important. An implementation of a numerical method is called a numerical algorithm.

We will simulate sample paths of time discrete approximations implemented in the R programming language [2] that we base on a finite discretization of a time interval  $[t_0, T]$ . We shall generate approximate values of the sample path for each step contained in the discretized interval [22].

In the fixed step methods, the distance between two contiguous points is the distance  $d_i = t_i - t_{i-1} = \frac{T-t_0}{N} \quad \forall i \mid 1 \leq i \leq N \in \mathbb{N}$ .  $N$  being the time interval partition number.

According to Kloeden [22], in the stochastic realm, simulated sample paths can be statistically analysed to find how good an approximation is compared to an exact solution. Moreover, the computational costs such as time and memory increases polynomially with the problem's dimension, which is good, and it is possible to apply variance reduction methods that allow a considerable decrease in the required sample size.

### 2.4.1 Convergence

As soon as we talk about numerical methods we are required to approach the topic of approximations and how to handle them. Methods efficiency receive the name of *convergence order*. In the SDE domain there are two main methods of convergency, that are classified according to their criteria. Firstly, we present the *strong order of convergence*. A method is said to have strong convergence  $\delta$  to  $Y$  if a time discretized  $Y_\delta$  of a continuous-time process  $Y$ , with  $\delta$  being the maximum time increment of the discretization, and for any fixed time horizon  $T$  holds true that [16]:

$$\mathbb{E} | Y_\delta(T) - Y(T) | \leq C\delta^\gamma, \quad \forall \delta < \delta_0$$

with  $\delta_0 > 0$  and  $C$  a constant not depending on  $\delta$ . Strong convergence addresses the problem of solutions' trajectories. For specific conditions, the Euler method has strong convergence order  $\gamma = \frac{1}{2}$ . Furthermore, there is the *weak order of convergence*. The weak convergence

$$| \mathbb{E} p(Y_n) - \mathbb{E} p(Y(\tau)) | \leq C\Delta t^\gamma$$

Strong and weak convergence are not mutually exclusive [16]. That means that a method with a given strong order of convergence might have a higher weak order of convergence too. This is the case for the Euler scheme, with a strong order of convergence of  $1/2$  and a weak order of  $1$  (under some conditions). For a more detailed and rigorous explanation of convergence we recommend consulting [24].

It is worth noting that, although schemes have a given convergency order, it is not unusual that they behave better than their order for some SDEs specifications.

### 2.4.2 Discretization

We know that convergence is an important feature to a numerical method and studies have found not all time discrete possible approximations of an SDE converge in a useful sense to the solution process as the step size adopted tends toward zero [25,26]. Moreover, particularly for SDEs, some of the more rapidly convergent methods available for ordinary differential equations (ODE) do not work, such as higher order Runge-Kutta methods.<sup>13</sup>

One of the methods that do work for ODEs and SDEs is the Euler method, named after the Swiss mathematician Leonhard Euler. Figure 2.4 shows an example of an implementation for the Newton's cooling law with timestep of 2 seconds compared to its analytical solution. This method (*a.k.a.* forward Euler method) is a first-order numerical procedure. It is the most basic explicit method<sup>14</sup> for numerical integration.

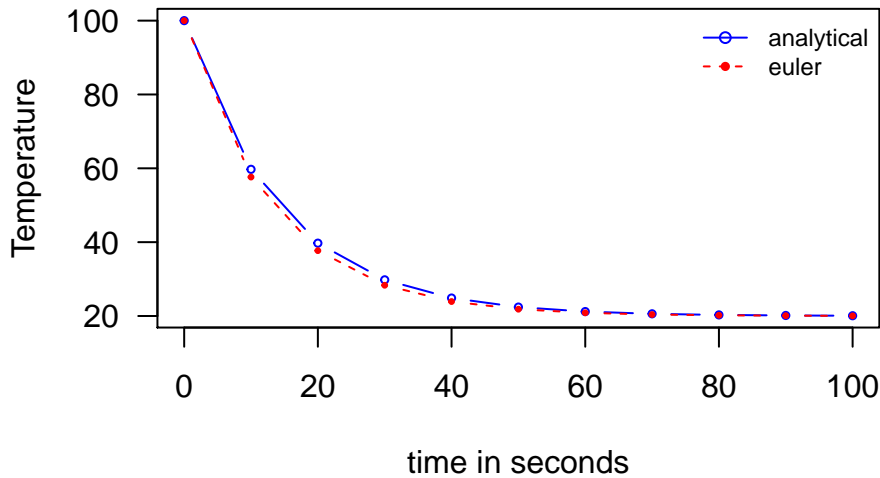


Figure 2.4: Analytical x Euler solutions

The method is first-order, as stated above, this means that the error in each step is a proportion of the square of the step size. Also, the global error at a given time is a function of the step size. We proceed to apply the Euler method to SDEs. Consider the equation:

$$dS_t = \mu(S_t, t)dt + \sigma(S_t, t)dW_t \quad (2.21)$$

$dW_t$  is the Brownian motion,  $\mu$  and  $\sigma$  are functions depending on  $S_t$  and  $t$ , over an

<sup>13</sup>The euler method is the simplest Runge-Kutta method.

<sup>14</sup>Explicit methods calculate the state of a system at a later time from the state of the system at the current time. Mathematically we have something like  $Y(t + \Delta t) = F(Y(t))$ .

interval  $[0, T]$ , and we want to discretize it as  $0 = t_1 < t_2 < \dots < t_m = T$  with increments equally spaced  $dt$ .

Integrating it from  $t$  to  $t+dt$  we have the starting point for our (and any) discretization scheme:

$$S_{t+dt} = S_t + \int_t^{t+dt} \mu(S_u, u) du + \int_t^{t+dt} \sigma(S_u, u) dW_u \quad (2.22)$$

To use the Euler discretization is the equivalent of approximating integrals using the left-point rule as in Figure 2.5<sup>15</sup>, we then have:

$$\begin{aligned} \int_t^{t+dt} \mu(S_u, u) dW_u &\approx \mu(S_t, t) \int_t^{t+dt} dW_u \\ &= \mu(S_t, t)(W_{t+dt} - W_t) \\ \int_t^{t+dt} \sigma(S_u, u) dW_u &\approx \sigma(S_t, t) \int_t^{t+dt} dW_u \\ &= \sigma(S_t, t)(W_{t+dt} - W_t) \\ &= \sigma(S_t, t)\sqrt{dt}Z \end{aligned}$$

$W_{t+dt} - W_t$  and  $\sqrt{dt}Z$  have identical distribution,  $Z$  being a standard gaussian variable. The Euler discretization of equation (2.22) is then:

$$S_{t+dt} = S_t + \mu(S_t, t)dt + \sigma(S_t, t)\sqrt{dt}Z \quad (2.23)$$

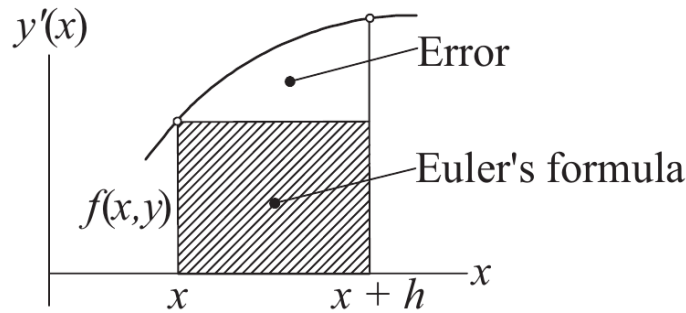


Figure 2.5: Graphical representation of Euler's formula

Source: Numerical methods in Engineering with Python 3.

<sup>15</sup>See Kiusalaas [27]

### Euler method - Heston model

We now proceed to apply the method to our model of interest. We retake the equations (2.16) and (2.17). We begin showing how to discretize the latter [16,24]:

$$V_{t+dt} = V_t + \int_t^{t+dt} \kappa(\theta - V_u)du + \int_t^{t+dt} \sigma\sqrt{V_u}dB_u \quad (2.24)$$

Which discretized turns out as:

$$\begin{aligned} \int_t^{t+dt} \kappa(\theta - V_u)du &\approx \kappa(\theta - V_t)dt \\ \int_t^{t+dt} \sigma\sqrt{V_u}dB_u &\approx \sigma\sqrt{V_t}(W_{t+dt} - W_t) \\ &= \sigma\sqrt{V_t}dtZ_v \end{aligned}$$

And leaves us with:

$$V_{t+dt} = V_t + \kappa(\theta - V_t)dt + \sigma\sqrt{V_t}dtZ_v \quad (2.25)$$

$Z_v$  is a standard normal variable. To avoid problems with negative values in  $\sqrt{V_t}$  we apply the *full truncation* scheme, which substitutes  $V_t$  with  $V_t^+ = \max(0, V_t)$ .<sup>16</sup>

For the  $S_t$  SDE we proceed similarly:

$$S_{t+dt} = S_t + \mu \int_t^{t+dt} S_u du + \int_t^{t+dt} \sqrt{V_u}S_u dW_u \quad (2.26)$$

Discretizing we have:

$$\begin{aligned} \int_t^{t+dt} S_u du &\approx S_t dt \\ \int_t^{t+dt} \sqrt{V_u}S_u dW_u &\approx \sqrt{V_t}S_t(W_{t+dt} - W_t) \\ &= \sqrt{V_t}dtS_tZ_s \end{aligned}$$

$Z_s$  is a standard normal variable with correlation  $\rho$  with  $Z_v$ . We have:

$$S_{t+dt} = S_t + \mu S_t dt + \sqrt{V_t}dtS_tZ_s \quad (2.27)$$

### 2.4.3 Stability

Most differential equations, deterministic or stochastic, cannot be solved explicitly [22]. Hence, stability studies begin with computers and is associated with numerical methods and approximations. Convergent methods were resulting in bigger errors than what was expected that could not be only due to discretization error. Eventually, scientists discovered that this unexpected problem was caused by accumulation of successive

---

<sup>16</sup>Another possible scheme (not used in this work) is the *reflection* scheme where we replace  $V_t$  with  $|V_t|$

truncation errors. Figure 2.6 retakes the cooling example previously approached to show instability due to an increase in size of the timestep and extending to 600 seconds. The top plot, with a step of 25 still converges to the real solution, but we already observe an odd behaviour since the numerical method doesn't follow the analytical solution, but instead revolves around it until convergence. The bottom plot, presents a small increase in the step  $h = 29$ , and this small increment is enough results in numbers completely off target.

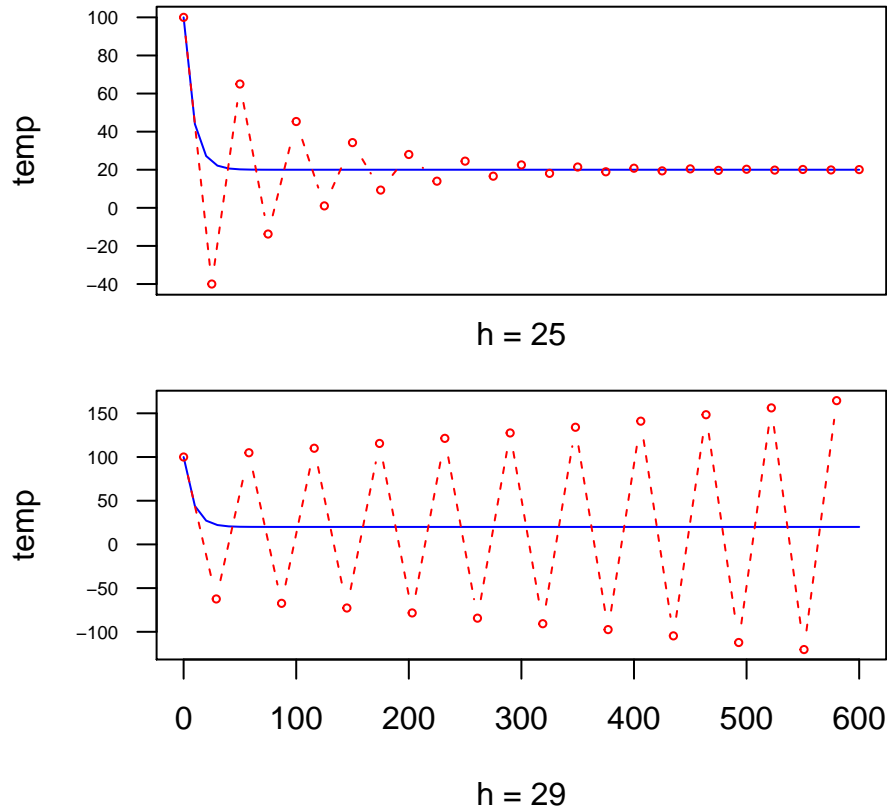


Figure 2.6: Euler's stability with different timesteps

We know that binary machines like computers are not able to represent all the real numbers, but only a subset of them. Thus, solving these errors is not straightforward since it's not possible to eliminate *all* truncation error when using a computer and dealing with numerical solutions. When faced to an incorrect (not acceptable) solution, we have to evaluate and distinguish between two distinct situations:

- i Rounding errors are considerably amplified by the algorithm. This situation is called numerical instability.
- ii Small perturbations of data generate large changes in the solution. This is

termed an ill-conditioned (or sensitive) problem.

Examples of these two classes of problem can be found in [21].

Stiff equations appear very often in mathematical problems and refer to differential equations for which a numerical methods might be unstable for not small enough stepsizes [28]. A differential equation of the form  $y' = f(t, y)$ , if its exact solution  $y(t)$  includes a term that decays exponentially to zero as  $t$  increases, but whose derivatives are greater in magnitude than the term itself. In other words, if it requires a significant depression of the stepsize to avoid stability lost. This is a loose definition but, since we are dealing with numerical methods a proper mathematical definition isn't required. Typically, these equations are of the form  $e^{-ct}$ , where  $c$  is a large positive constant [29]. A practical example of the stiff behavior is the following differential equation [30]:

$$y' = -100y, \quad t > 0, \quad y_0 = 1$$

Whose exact solution is  $y_t = e^{-100t}$  and goes to zero as  $t$  increases. Applying Euler's method to this equation with  $h = 0.1$  we stumble in the following equation

$$y_{n+1} = y_n - 100hy_n = -9y_n$$

which is wrong, since it yields an exponentially growing solution  $y_n = (-9)^n$ . On the other hand, if our timestep is smaller  $h = 10^{-3}$ , our solution using Euler's method becomes  $y_n = (0.9)^n$ . This solution leads to an accurate behavior regarding the exact solution, it rapidly decays to zero.

Sometimes, it is interesting to rank differential equations that are more or less stiff. Thus, people compute the quotient of the largest and the smallest eigenvalues of a linear system. They call it the equations' *stiffness ratio* and, usually the bigger the stiffness ratio, the more likely is to be stiff [28].

## Stability Domain

Let's take the equation:

$$y' = \lambda y, \quad t \geq 0, \quad y_0 = 1 \tag{2.28}$$

where  $\lambda \in \mathbb{C}$  or in other terms  $\lambda = \lambda_r + i\lambda_i$  and whose solution is  $y + t = e^{\lambda t}$ . We can rewrite this equation as a system:

$$\frac{d}{dt} \begin{bmatrix} y^1 \\ y^2 \end{bmatrix} = \begin{pmatrix} \lambda_r & -\lambda_i \\ \lambda_i & -\lambda_r \end{pmatrix} \begin{bmatrix} y^1 \\ y^2 \end{bmatrix} \tag{2.29}$$

The  $\lim_{t \rightarrow \infty} y_t = 0$  if and only if  $\mathbb{R}\lambda < 0$ . The *linear stability domain*  $\mathcal{D}$  is defined as the set of all numbers  $\Delta\lambda \in \mathbb{C}$  such that  $\lim_{n \rightarrow \infty} y_{n=0} = 0$ , with  $\Delta > 0$  being the stepsize. Or, as stated in Kloeden [22], the suitable values of the stepsize are expressed in terms of *region of absolute stability*, consisting of the complex numbers  $\lambda\Delta$  for which an error in  $y_0$  at  $t_0$  will not grow in subsequent iterations of the method.



Without entering all the details, for these we recommend [22,28,30], the euler's stability domain is:

$$\mathcal{D}_{Euler} = \{z \in \mathbb{C} : |1 + z| < 1\}$$

which represents the interior of a complex disc of unit radius and centre  $z = -1$  as can be seen in Figure 2.7, on the left. The right side of the Figure 2.7 shows the stability region called A-stability.<sup>17</sup> If a method is A-stable, the stepsize  $\Delta$  is only constraint by accuracy.

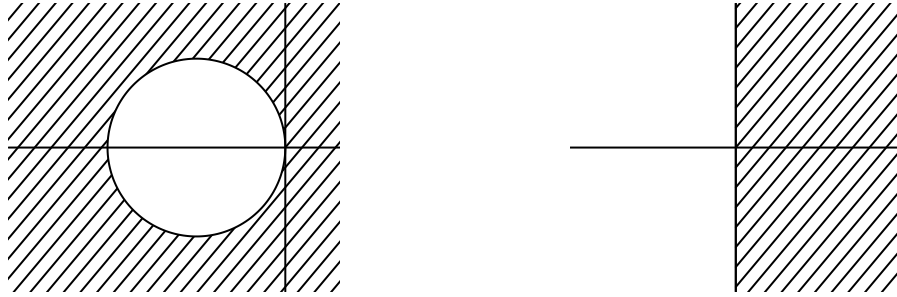


Figure 2.7: Stability domains

Thereby, we claim that stability method study is an important topic, since it enables achieving solutions that are good to stiff equations without having to overly reduce our timesteps which can be very computationally costly.

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<sup>17</sup>Mathematically:  $\mathcal{D} \subseteq \{z \in \mathbb{C} : \operatorname{Re} z < 0\}$ .



## Chapter 3

# The Heston Model Implementation

In section 2.3.2 we presented Heston's SDE system in one of its structures. Another common way [31–33] to write down the system is using the property presented in subsection 2.1.2 as in equation (3.1).

$$\begin{aligned}dS_t &= \mu S_t dt + \rho \sqrt{V_t} dB_t + \sqrt{1 - \rho^2} \sqrt{V_t} S_t dW_t \\dV_t &= k(\theta - V_t) dt + \sigma \sqrt{V_t} dB_t\end{aligned}\tag{3.1}$$

### 3.1 Characteristic Function

The Heston model characteristic function is firstly presented in the 1993 Steven Heston's paper [20] and is described below [34]:

$$f(S_t, V_t, t) = e^{A(T-t) + B(T-t)S_t + C(T-t)V_t + i\phi S_t}\tag{3.2}$$

If we let  $\tau = T - t$ , then the explicit form of the Heston characteristic function is:

$$\begin{aligned}f(i\phi) &= e^{A(\tau) + B(\tau)S_t + C(\tau)V_t + i\phi S_t} \\A(\tau) &= ri\phi\tau + \frac{\kappa\theta}{\sigma^2} \left[ -(\rho\sigma i\phi - \kappa - M)\tau - 2 \ln \left( \frac{1 - Ne^{M\tau}}{1 - N} \right) \right] \\B(\tau) &= 0 \\C(\tau) &= \frac{(e^{M\tau} - 1)(\rho\sigma i\phi - \kappa - M)}{\sigma^2(1 - Ne^{M\tau})}\end{aligned}$$

Where:

$$\begin{aligned}M &= \sqrt{(\rho\sigma i\phi - \kappa)^2 + \sigma^2(i\phi + \phi^2)} \\N &= \frac{\rho\sigma i\phi - \kappa - M}{\rho\sigma i\phi - \kappa + M}\end{aligned}$$

This function is the driving force behind the following formula, that calculates the fair value of a European call option at time  $t$ , given a strike price  $K$ , that expires at

time  $T$  [34]:

$$C = \frac{1}{2}S(t) + \frac{e^{-r(T-t)}}{\pi} \int_0^\infty \Re \left[ \frac{K^{-i\phi} f(i\phi + 1)}{i\phi} \right] d\phi - Ke^{-r(T-t)} \left( \frac{1}{2} + \frac{1}{\pi} \int_0^\infty \Re \left[ \frac{K^{-i\phi} f(i\phi)}{i\phi} \right] d\phi \right) \quad (3.3)$$

## 3.2 Euler Scheme

Given the fact that the underlying asset is temporal dependent upon the solution of the SDE's volatility, we simulate the volatility's path before the asset's. If the Black-Scholes model enabled using Ito's Lemma directly for solving  $S_t$ , this equation system requires numerical methods. We present here the Euler Scheme - Full Truncation algorithm (and compare to other similar schemes) [31] along with some insights on how it was implemented in R. The Euler discretization brings approximation paths to stock prices and variance processes. If we set  $t_0 = 0 < t_1 < \dots < t_M = T$  as partitions of a time interval of  $M$  equal segments of length  $\delta t$ , we have the following discretization for the stock price:

$$S_{t+1} = S_t + rS_t + \sqrt{V_t}S_tZ_s \quad (3.4)$$

And for the variance process:

$$V_{t+1} = f_1(V_t) + \kappa(\theta - f_2(V_t)) + \sigma\sqrt{f_3(V_t)}Z_v \quad (3.5)$$

$Z_s$  being a standard normal random variable, i.e.  $N \sim (0, 1)$ , we set  $Z_t$  and  $Z_v$  as two independent standard normal random variables and  $Z_s$  and  $Z_v$  having correlation  $\rho$ . This means we can write  $Z_s = \rho Z_v + \sqrt{1 - \rho^2}Z_t$ .

The immediate observable problem in the proposed discretization scheme is that  $V$  can become negative with non-zero probability making the computation of  $\sqrt{V_t}$  impossible [32]. There are several proposed fixes that can be used as you can see below: Where  $V^+ = \max(V, 0)$  and  $|V|$  is the absolute value of  $V$ .

Scheme	$f_1(V_t)$	$f_2(V_t)$	$f_3(V_t)$
Reflection	$ V $	$ V $	$ V $
Partial Truncation	$V$	$V$	$V^+$
Full Truncation	$V$	$V^+$	$V^+$

Table 3.1: Truncation schemes

We chose to fix our discretization using the Full-Truncation (FT) scheme and thus, rewrite the equations as follows:

$$S_{t+1} = S_t + rS_t + \sqrt{V_t^+}S_tZ_s \quad (3.6)$$

$$V_{t+1} = V_t + \kappa(\theta - V_t^+) + \sigma\sqrt{V_t^+}Z_v \quad (3.7)$$

Our R implementation follows the euler's scheme with hardly any modifications. It draws two Gaussian random variables ( $Z_v$  and  $Z_t$ ) using the function *rnorm* to create  $Z_s$  with correlation  $\rho$  with  $Z_v$ .

```
Zv <- stats::rnorm(N)
Zt <- stats::rnorm(N)
Zs <- rho * Zv + (sqrt(1 - (rho^2))) * Zt)
```

And use it to compute  $S$  and  $V$  as in the code snippet bellow. The two modifications we apply are previously computing the square roots of the stepsize  $dt$  and the *aux* variable as to improve speed. The *aux* variable is an help variable created to impose positivity to  $V$  as we are operating the full-truncation euler scheme.

```
S <- S * (1 + r * dt + sqrt_aux * Zs * sqrt_dt)
v <- v + k * dt * (theta - aux) +
      sigma * sqrt_aux * Zv * sqrt_dt
```

We could have used the *pmax* function in R, but this function is slow, therefore we opted to create the *aux* variable and impose positive values using the following R syntax:

```
aux <- v
aux[v < 0] <- 0
```

Depending on the reader's R fluency, other parts of the scheme might present a challenge, and that is why a version of the function is fully presented in appendix C.1.

### 3.3 Kahl-Jackel

Kahl-Jackel propose a discretization method they refer to as the “IJK” method [32,33] that coupled with the implicit Milstein scheme for the variance lands the system of equations (3.8) and (3.9). It is possible to verify that this discretization always results in positive paths for  $V$  if  $4\kappa\theta > \sigma^2$ . Unfortunately, this inequality is rarely satisfied when we plug real market data to calibrate the parameters. This means we must have a defined strategy for when the inequality doesn't hold. We use the scheme proposed in Andersen [32], where a truncation similar to the Euler's is applied. Whenever our volatility  $V_t$  drops below zero we use (3.7), and implement  $\hat{V}(t + \Delta)^+$  and  $\hat{V}(t)^+$  instead of  $\hat{V}(t + \Delta)$  and  $\hat{V}(t)$  of equation (3.8). The code guidance to this method can be found in appendix C.2.

$$\ln \hat{S}(t + \Delta) = \ln \hat{S}(t) - \frac{\Delta}{4} \left( \hat{V}(t + \Delta) + \hat{V}(t) \right) + \rho \sqrt{\hat{V}(t)} Z_v \sqrt{\Delta} \quad (3.8)$$

$$+ \frac{1}{2} \left( \sqrt{\hat{V}(t + \Delta)} + \sqrt{\hat{V}(t)} \right) \left( Z_s \sqrt{\Delta} - \rho Z_v \sqrt{\Delta} \right) + \frac{1}{4} \sigma \rho \Delta \left( Z_V^2 - 1 \right)$$

$$\hat{V}(t + \Delta) = \frac{\hat{V}(t) + \kappa \theta \Delta + \sigma \sqrt{\hat{V}(t)} Z_V \sqrt{\Delta} + \frac{1}{4} \sigma^2 \Delta \left( Z_V^2 - 1 \right)}{1 + \kappa \Delta} \quad (3.9)$$

### 3.4 Exact Algorithm

In 2006, Broadie-Kaya [31] propose a method that has a faster convergence rate,  $\mathcal{O}(s^{-1/2})$  than some of the more famous schemes, such as Euler's and Milstein's discretizations. They build their idea to generate an exact sample from the distribution of the terminal stock price based on numerous papers [20]. The stock price and variance are as follows:

$$S_t = S_0 \exp \left[ \mu t - \frac{1}{2} \int_0^t V_s ds + \rho \int_0^t \sqrt{V_s} dB_s + \sqrt{1 - \rho^2} \int_0^t \sqrt{V_s} dW_s \right] \quad (3.10)$$

The squared volatility of the variance process is:

$$V_t = V_0 + \kappa \theta t - \kappa \int_0^t V_s ds + \sigma \int_0^t \sqrt{V_s} dB_s \quad (3.11)$$

The algorithm used to generate the model consists in four steps as follows:

*Step 1.* Generate a sample of  $V_t$  given  $V_0$

*Step 2.* Generate a sample of  $\int_0^t V_s ds$  given  $V_t, V_0$

*Step 3.* Compute *displaystyle*  $\int_0^t \sqrt{V_s} dB_s$  given  $V_t, V_0$  and  $\int_0^t V_s ds$

*Step 4.* Generate a sample from the probability distribution of  $S_t$ , given *displaystyle*  $\int_0^t \sqrt{V_s} dB_s$  and *displaystyle*  $\int_0^t V_s ds$

#### 3.4.1 Generate a sample of $V_t$ given $V_0$

The distribution of  $V_t$  given  $V_0$  for  $0 < t$  is a non-central chi-squared distribution [19,38]:

$$V_t = \frac{\sigma^2(1 - e^{-\kappa t})}{4\kappa} \chi_\delta^2 \left( \frac{4\kappa e^{-\kappa t}}{\sigma^2(1 - e^{-\kappa t})} \times V_0 \right)$$

where  $\delta = \frac{4\theta\kappa}{\sigma^2}$  and  $\chi_\delta^2(\lambda)$  denotes a non-central chi-squared random variable with  $\delta$  degrees of freedom and  $\lambda$  as its non-centrality parameter.

Broadie and Kaya [31] sample generating Poisson and gamma distributions as in Johnson et al. [39]. We used the built-in function in R *rchisq*, which uses this exact method for sampling, see chunk below.

```
d1 <- (4 * k * theta)/(sigma)^2
c0 <- (sigma^2 * (1 - exp(-k*tau)))/(4*k)
dt <- (tau-t)

# sampling V
lambda <- (4*k*exp(-k*dt)*v)/(sigma^2 * (1-exp(-k*dt)))
vt <- c0 * stats::rchisq(n = 1, df = d1, ncp = lambda)
```

### 3.4.2 Generate a sample of $\int_0^t V_s ds$ given $V_t, V_0$

After generating  $V_t$ , we follow the instructions in [31,39]. We use the characteristic function (3.12) to compute the probability density function  $F(x)$ .

$$\begin{aligned}
\Phi(a) &= \mathbb{E} \left[ \exp \left( ia \int_0^t V_s ds \mid V_0, V_t \right) \right] \\
&= \frac{\gamma(a) e^{(-1/2)(\gamma(a)-\kappa)t} (1 - e^{-\kappa t})}{\kappa (1 - e^{-\gamma(a)t})} \\
&\quad \times \exp \left\{ \frac{V_0 + V_t}{\sigma^2} \left[ \frac{\kappa (1 + e^{-\kappa t})}{1 - e^{-\kappa t}} - \frac{\gamma(a) (1 + e^{-\gamma(a)t})}{1 - e^{-\gamma(a)t}} \right] \right\} \\
&\quad \times \frac{I_{0.5\delta-1} \left[ \sqrt{V_0 V_t} \frac{4\gamma(a) e^{-0.5\gamma(a)t}}{\sigma^2 (1 - e^{-\gamma(a)t})} \right]}{I_{0.5\delta-1} \left[ \sqrt{V_0 V_t} \frac{4\kappa e^{-0.5\kappa t}}{\sigma^2 (1 - e^{-\kappa t})} \right]}
\end{aligned} \tag{3.12}$$

where  $\gamma(a) = \sqrt{\kappa^2 - 2\sigma^2 ia}$ ,  $\delta$  was previously defined and  $I_v(x)$  is the modified Bessel function of the first kind.<sup>1</sup> There are no mysteries implementing the characteristic function as you can observe in the chunk below. Although R has a built-in Bessel function, it only accounts for real numbers. Thus, we were obliged to use the Bessel package [40] that accounts for complex numbers. Once again, we pre compute some of the operations that are repeated through the function as to reduce computational time. This is specially important for this method since it involves operations of high complexity.

```

phi_heston <- function(a, v0, v_t, d){

  gamma_a <- sqrt(k^2 - 2 * sigma^2 * 1i*a)
  gammadt <- gamma_a * (tau-t)
  sqrtv0vt <- sqrt(v0*v_t)
  delta <- -k * (tau-t)

  part1 <- (gamma_a * exp(-(gamma_a - k)/2 * (tau-t)) * (1 - exp(delta))) /
    (k * (1 - exp(- gammadt)))
  part2 <- exp((v0+v_t)/(sigma^2) * ( (k * (1 + exp(delta)))/(1-exp(delta)) -
    (gamma_a * (1 + exp(- gammadt)))/(1-exp(- gammadt))))
  part3 <- Bessel::BesselI(z = ((4 * gamma_a * sqrtv0vt)/(sigma^2) *
    exp(- gammadt/2) /
    (1 - exp(- gammadt))), nu = 0.5*d - 1) /
    Bessel::BesselI(z = ((4 * k * sqrtv0vt)/(sigma^2) *
    (exp(delta/2))/(1-exp(delta))), nu = 0.5*d - 1)
  return (part1 * part2 * part3)
}

```

The probability distribution function is obtained in [31,38] by Fourier inversions using Feller [41]. We use the approach in Gil-Pelaez [42], equation (3.13). We

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<sup>1</sup>See *Appendix*  
*refbessel* for more information.

define  $V(u, t)$  the random variable with the same distribution as the integral  $\int_u^t V_s ds$ , conditional on  $V_u$  and  $V_t$ :

$$F(x) \equiv Pr \{V(u, t) \leq x\} = F_X(x) = \frac{1}{2} - \frac{1}{\pi} \int_0^\infty \frac{\text{Im}[e^{-iux} \text{phi}(u)]}{u} du \quad (3.13)$$

$\text{Im}$  denotes the imaginary part of  $e^{-iux} \text{phi}(u)$ . Equation (3.13) is computed numerically and we then sample it by inversion. The integral function of the volatility is composed by five other functions inside of it. Since it acts only as a wrapper to these functions we are going to omit it below, but the complete code can be found in appendix C.3. The first represents our integrand. The following function takes our integrand and actually performs the integration, returning the value. For speed purposes, we limited the integral upper bound to 1000 and increased the tolerance to  $10^{-3}$ . After these calculations we have now the function's cumulative function.

```
integrand <- function(x, phi = cf){

  f2 <- function(u){
    Im(phi(u) * exp(-1i * u * x)) / u
  }
  return(f2)
}

## integrate to "cdf"
F_x <- function(x) {
  y <- 0.5 - 1/pi * stats::integrate(integrand(x), lower= 0, upper= 1000,
                                     rel.tol = 0.001, stop.on.error = FALSE)$value
  return(y)
}

## endsign
endsign <- function(f, sign = 1) {
  b <- sign
  while (sign * f(b) < 0) b <- 10 * b
  return(b)
}
```

After computing the integral, we need to set the inversion function to place. To do that we generate the *endsign* function above. It is a simple function that is used to guarantee that when we call *uniroot*, the function to find roots built-in in R, our bounds have different signs. We set our lower and upper bounds as  $-\infty$  and  $\infty$ , respectively and define an auxiliary function *subcdf* that will perform the inversion subtracting an uniform from it.

```
## inversion
low_bound = -Inf
upp_bound = Inf
invcdf <- function(u) {
  subcdf <- function(t) F_x(t) - u
```



```

if (low_bound == -Inf)
  low_bound <- endsign(subcdf, -1)
if (upp_bound == Inf)
  upp_bound <- endsign(subcdf)
return(uniroot(subcdf, lower=low_bound, upper=upp_bound,
               tol = 0.001220703)$root)
}
U <- stats::runif(n)
sapply(U, invcdf)

```

Furthermore, we also introduce (not in the same simulations, obviously) a simpler version for this step, that computes this integral approximation, using the solution  $\int_u^t V_s ds = \frac{1}{2} (V_u + V_t)$ . This solution is called *drift interpolation* [43] and its implementation is out of the box but we provide it anyway:

```
int_v <- dt * ((1/2) * v + (1/2) * vt)
```

### 3.4.3 Compute $\int_0^t \sqrt{V_s} dB_s$ given $V_t$ , $V_0$ and $\int_0^t V_s ds$

From equation (3.11) we are now able to compute this integral:

$$\int_0^t \sqrt{V_s} dB_s = \frac{V_t - V_0 - \kappa\theta t + \kappa \int_0^t V_s ds}{\sigma} \quad (3.14)$$

The last step of the algorithm consists of computing the conditional distribution of  $\log S_t$  based on the fact that the process for  $V_t$  is independent from  $dB_t$ , and the distribution of  $\int_0^t \sqrt{V_s} dB_s$  is normal with mean 0 and variance  $\int_0^t V_s ds$ , given  $V_t$ .

$$m(u, t) = \log S_0 + \left[ \mu t - \frac{1}{2} \int_0^t V_s ds + \rho \int_0^t \sqrt{V_s} dB_s + \sqrt{1 - \rho^2} \int_0^t \sqrt{V_s} dW_s \right]$$

and variance

$$\sigma^2(0, t) = (1 - \rho^2) \int_0^t V_s ds$$

We generate the  $S_t$  sample using a standard normal random variable  $Z$  and set:

$$S_t = e^{m(0,t) + \sigma(0,t)Z}$$

In R code, we have the following:

```

int_vdw <- (1/sigma) * (vt - v - k * theta * dt + k * int_v)

m <- log(S) + (r * (tau - t) - (1/2) * int_v + rho * int_vdw)
std <- sqrt((1 - rho^2)) * sqrt(int_v)
S <- exp(m + std * stats::rnorm(1))

```

### 3.4.4 Limitations

The biggest limitation this scheme presents is that the second step is computationally costly. It demands the inversion of the distribution function of  $\left(\int_0^t V_s ds \mid V_t, V_0\right)$  numerically. We must perform a root search of  $F(x_i) - U = 0$  testing for different  $x_i$ . Notwithstanding, we do not know the cumulative form of  $F(x_i)$  distribution and have to perform our root finding strategy starting from the characteristic function, which contains two modified Bessel functions inside, in a structure that is rerun until a given tolerance  $\epsilon$  is reached. Mathematically:  $F(x_i) - U = \epsilon$ .

# Chapter 4

## Results

We present here the results of all the implementations that were disclosed in the previous section. We perform numerical comparisons between all the methods, setting out differences accross number of simulations and timesteps. These simulations were performed as a way to validate our implementations, and thus do not have the intent to exhaust all kinds of possibilities or provide sharp comparisons.

Heston [20] gives a closed form used as benchmark and as the ‘true’ option value and enabling the results to be exposed in terms of bias<sup>1</sup> and RMSE (root mean square error).<sup>2</sup>

The simulaton experiments were performed on a notebook with an Intel(R) Core(TM) i7-4500U CPU @ 1.80GHz processor and 8GB of RAM running on a linux x86\_64 based OS, Fedora 25. Codes were all written in R an run on version 3.4.1 “Single Candle” [2].

First of all, we chose a parametrizations based on what we saw in other works, made some adjustments, like reducing the options’ time to maturity due to the slower nature of R language and compared initial the results with the true option price given by the function *callHestoncf* belonging to the package NMOF [44]. Parameters can be seen in Table 4.1 below. As we said above, these parameter weren’t chosen at

Variables	Values
dt	0.05
k	2.00
r	0.05
rho	-0.30
S	100.00
sigma	0.20
t	0.00
tau	1.00
theta	0.09
v	0.09
X	100.00

Table 4.1: Model Parameters

---

<sup>1</sup>Defined as  $\mathbb{E}[\hat{\alpha} - \alpha]$

<sup>2</sup>Defined as  $\sqrt{\mathbb{E}((\hat{\theta} - \theta)^2)}$

random, we took them from Table 2 of the Broadie-Kaya [31] paper, with two minor modifications performed. We reduced the time to maturity from 5 to 1 year and we reduced the  $\sigma$  from 1 to 0.2 following the “rule” they acknowledge in their paper that  $\frac{2\kappa\theta}{\sigma} \leq 1$ . The setup where this inequality isn’t followed produces bigger biases, especially in Euler’s and Kahl-Jackel’s discretizations. In this situations, the EA often outperforms the Euler and KJ solutions in terms of RMSE.

To perform our simulations, we first fixed a seed and saved the results in Table 4.2. Since the value given by the *callHestoncf* function with the parameters in Table 4.1 is 14.176, the method that best approached the “true” value was the modified (drift interpolated) exact algorithm with 100,000 simulations. Although the Euler scheme gives the same result (14.16) as the modified EA with 10,000 simulations, it moves away from the closed form value when we run 100,000 simulations. Results are

Simulations	Euler	KJ	EA-BK	EA-DI
1,000	13.28	13.05	14.71	14.74
10,000	14.16	13.86	14.45	14.38
100,000	14.11	13.83	14.21	14.16

Note: Simulations performed with 20 steps, except the EA BK

Table 4.2: Option Value

observable in Figure 4.1 also. The plot gives a good sense of possible biases associated with each method. It also show that the methods tend to converge quickly to the real result when the number of steps is increased.

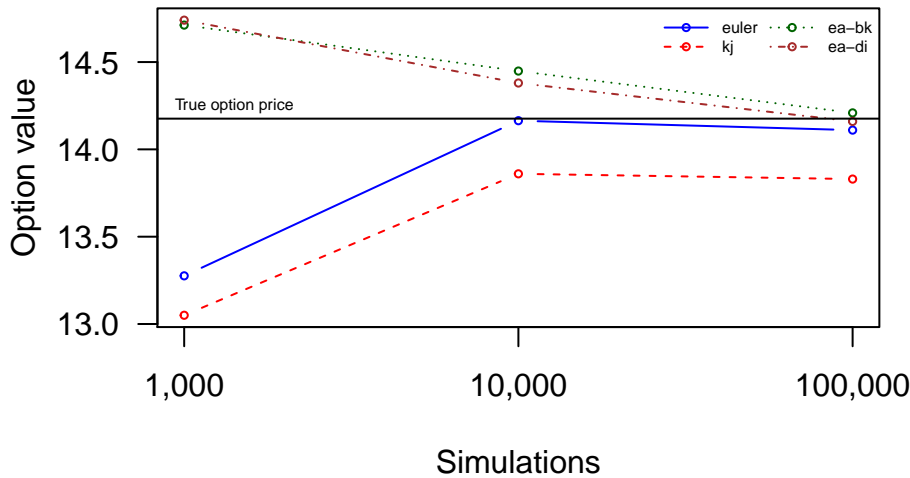


Figure 4.1: Comparison between models, 20 steps.

To verify if in fact these methods present bias, we performed ten thousand simulations of ten thousand paths to the Euler, Kahl-Jackel (KJ) and drift interpolated exact algorithm (EA-DI) with 20 steps ( $dt = 0.05$ ) first. We present these results in appendix C.4 because we present the more precise simulations we made, with 100 steps in Figure 4.2. Results seem to show a little bias, thus we present in Table 4.3 how precise models are.

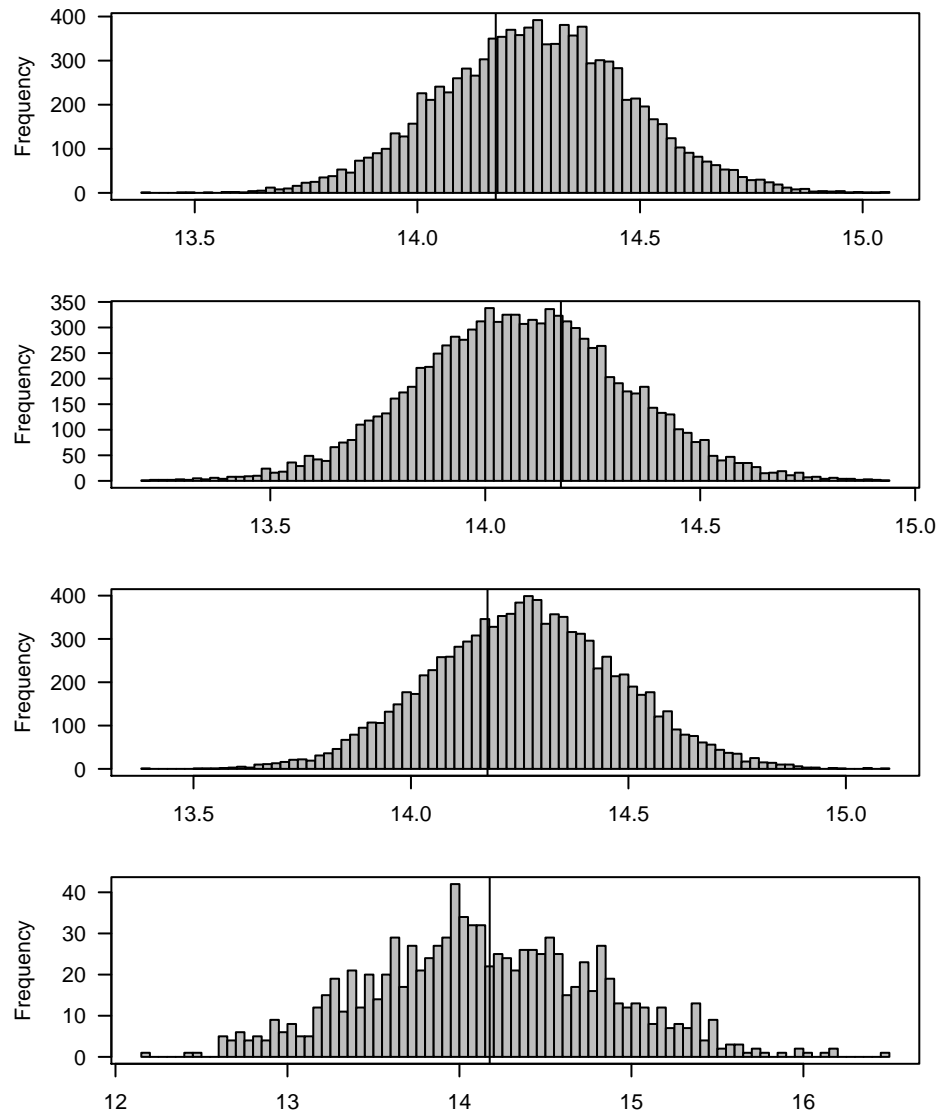


Figure 4.2: Comparison between models, 100 steps.

The exact algorithm performs much slower than the other three we only performed 1000 simulations of one thousand paths each. That is why its RMSE is comparably bigger than the other three implementations. We are aware that, by construction the Broadie-Kaya version of the algorithm has no bias, but we tested the bias of our implementation (even though we didn't put it in the Table 4.3) and the reason is

	Euler	KJ	EA-BK	EA-DI
bias	0.09	-0.09	-	0.09
sd	0.21	0.24	-	0.22
RMSE	0.23	0.26	0.68	0.23

Note: Simulations performed with 100 steps, except the EA BK

Table 4.3: Estimators

twofold: again, we wanted to check if we correctly implemented the algorithm as a whole, but more importantly we wanted to check the outcomes of the adaptations we made (in terms of truncation, relative error and others) were playing into our model. The computed, even with far less simulations was far smaller than the other three methods, with only  $-0.02$ . We understand that this result confirms that our adaptations did not change the nature of the model and with a greater number of simulations the bias would be zero as it should. We observe bias level in the three classical numerical models, the KJ model undervaluing in the same order (0.09) as the other two overvalue the option. In terms of standard deviation the models seem to be equivalent, with the KJ model presenting a slightly bigger standard deviation than the other two (0.24 against 0.21 and 0.22). According to the tests performed, we are confident that our models were implemented correctly.

# Chapter 5

## Conclusion

As we revisit what was presented through this thesis, we highlight that we were unable find codes of the heston's exact algorithm in any shape or form. That means we searched if the algorithm implementation was available in any programming language, or at least a reference to where to find it and we couldn't.

Firstly, we provided a succinct, but rigorous, theoretical framework. Since this topic is very embracing, we provided an abundant amount of literature not only to the core topics approached, but also to the marginal ones. The first topic introduced in our review of literature is the stochastic calculus subject. Following the basics, we present the Black-Scholes model since this is the basic model from which the stochastic volatility ones derive from. We show how the BS formula is computed, and its limitations. The next logical step are the stochastic volatility models since they directly attack and solve the biggest BS model limitation: fixed volatility. We present the CIR before the Heston model as the first concocts the latter. Lastly, approach the challenges brought by the adoption of numerical methods, to better understand if a method is theoretically good, how to perform a discretization, and why a supposedly good discretization scheme can fail, nevertheless.

Thereafter, we thoroughly introduced the algorithms we were going to implement in R, and how we were going to do it. Four different algorithms were presented: the Euler scheme, the Kahl-Jackel algorithm, and two versions of the "exact algorithm". The first trying to be as similar as possible to the one presented in Broadie-Kaya's paper and the modified version bringing an approximation to the algorithm's second step (the costlier step) as we replace the integral with a drift interpolation.

We tested the implementations and presented our results in Chapter @ref(#resultados). We used a specification provided in Broadie-Kaya's paper [31], Table 2 with two slight changes. We reduced the time to maturity from 5 to 1 year and we reduced the  $\sigma$  from 1 to 0.2 following the rules he acknowledges and previously cited. We verified that, in fact, the EA algorithm does not present bias, but because of the nature of the algorithm and of the implementation in R, we doubt it can be used practically without finding a way around the algorithm's second step.

As we performed the EA's implementation we faced some challenges we weren't able to resolve yet. We couldn't parallelize the root inversion, which was a big bottleneck for the algorithm that we will solve. Also, we only created methods for the European

*call* and for four different schemes. We intend to widen the scope to embrace the *put* computation and introduce the binomial tree approach along with the method of moments and maximum likelihood. Finally, the package with all its functions is available at this **link**<sup>1</sup> and in the near future on **CRAN**.

---

<sup>1</sup><https://github.com/fernate7/rnmethods>



# Appendix A

## Black-Scholes formula

In this appendix, we start from the following Geometric Brownian Motion process:

$$dS_t = \mu S_t dt + \sigma S_t dB_t$$

And we claim that the solution to this stochastic differential equation applying Itô's calculus is the following formula:

$$S_T = S_0 \times \exp \left( \left( \mu - \frac{\sigma^2}{2} \right) T + \sigma B_T \right)$$

*Proof.* If  $S$  were deterministic,  $dS_t/S_t$  would be the derivative of  $\ln(S_t)$  with respect to  $S$ . This suggests to find an expression for the stochastic differential of  $\ln(S_t)$ , a function of the single random variable  $S_t$ .

$$f(t, S) = \ln(S)$$

$$df(t, S) = \frac{\partial f}{\partial S} dS + \frac{1}{2} \frac{\partial^2 f}{\partial S^2} (dS)^2$$

$$d \ln(S) = \frac{d \ln(S)}{dS} dS + \frac{1}{2} \frac{d^2 \ln(S)}{dS^2} (dS)^2$$

$$(dS)^2 = \int_0^t (\sigma \times S)^2 ds = \sigma^2 S^2 dt$$

$$d \ln(S) = \frac{1}{S} (\mu S dt + \sigma S dB) + \frac{1}{2} \frac{-1}{S^2} \sigma^2 S^2 dt$$

$$d \ln(S) = \left( \mu - \frac{1}{2} \sigma^2 \right) dt + \sigma dB$$

$$\int_{t=0}^T d \ln(S_T) = \int_{t=0}^T \left( \mu - \frac{\sigma^2}{2} \right) dt + \int_{t=0}^T \sigma dB_t$$

$$\ln(S_T) - \ln(S_0) = \left(\mu - \frac{\sigma^2}{2}\right)T + \sigma B_T$$

$$\ln\left(\frac{S_T}{S_0}\right) = \left(\mu - \frac{\sigma^2}{2}\right)T + \sigma B_T$$

$$S_T = S_0 \times \exp\left(\left(\mu - \frac{\sigma^2}{2}\right)T + \sigma B_T\right)$$

□

# Appendix B

## Bessel Function

The modified Bessel function of the first kind can be described in the shape of a contour integral (below) and is plotted for three different  $\nu$  in Figure B.1.

$$I_n(z) = \frac{1}{2\pi i} \oint e^{(z/2)(t+1/t)} t^{-n-1} dt$$

However, the Broadie-Kaya paper [31] presents it as a power series:

$$I_\nu(z) = \left(\frac{1}{2}z\right)^\nu \sum_{k=0}^{\infty} \frac{\left(\frac{1}{4}z^2\right)^k}{(k!\Gamma(\nu + k + 1))}$$

With  $\Gamma(x)$  is the gamma function and  $\nu$  is a complex number.

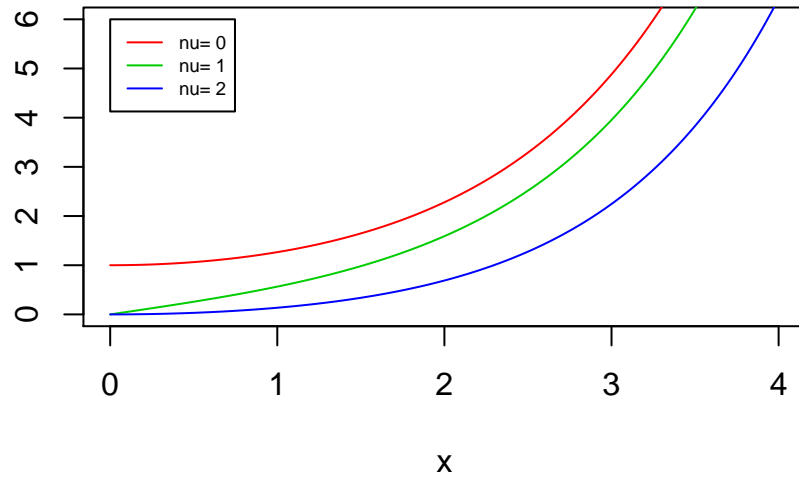


Figure B.1: Modified Bessel Functions of the First Kind



# Appendix C

## Implementations

### C.1 Euler Scheme

```
hestoneuler <- function(S, X, r, v, theta, rho, k,
                        sigma, t = 0, dt = NULL, tau = 1, N){

  if(is.null(dt)){ dt <- (tau-t)/1000}
  sequencia <- seq(t,tau,dt)
  ST <- matrix(NA, length(sequencia), N) #transformar em matrix
  aux <- NULL
  sqrt_dt <- sqrt(dt)

  for(i in sequencia){
    Zv <- stats::rnorm(N)
    Zt <- stats::rnorm(N)
    Zs <- rho * Zv + (sqrt(1 - (rho^2))) * Zt
    aux <- v
    aux[v < 0] <- 0
    sqrt_aux <- sqrt(aux)
    S <- S * (1 + r * dt + sqrt_aux * Zs * sqrt_dt)
    S[S <= 0] = 0
    v <- v + k * dt * (theta - aux) + sigma * sqrt_aux * Zv * sqrt_dt
    ST[j,] <- S
  }
  rm(aux, v, Zv, Zt, Zs, S, j)
  ST <- as.matrix(ST, ncol=N)
  Result <- ST[nrow(ST),] - X
  Result[Result <= 0] = 0
  call = mean(exp(-r*(tau-t))*Result)
  lista = list('call' = call, 'Result' = Result, 'Spot' = ST)
  return(lista)
}
```

## C.2 Kahl-Jackel Implementation

```
Hestoncallkj <- function(S, X, r, q, v, theta, rho, k,
                        sigma, t = 0, dt = NULL, tau = 1, N){

  if(is.null(dt)){ dt <- (T-t)/1000}
  v <- rep(v,N)
  theta<- rep(theta,N)
  sequencia <- seq(t,tau,dt)
  ST <- matrix(NA, length(sequencia), N) #transformar em matrix
  S <- log(S)

  for(i in seq(t,tau,dt)){
    Zv <- stats::rnorm(N)
    Zt <- stats::rnorm(N)
    Zs <- rho * Zv + sqrt(1 - rho^2) * Zt
    vt <- (v + k * theta * dt + sigma * sqrt(v) * Zv * sqrt(dt) +
           (1/4) * sigma^2 * dt * ((Zv)^2 - 1))/(1 + k * dt)
    vt[vt <= 0] <- v[vt <= 0] + k * dt * (theta[vt <= 0] -
                                           v[vt <= 0],0) + sigma * sqrt(v[vt <= 0],0) *
                                           Zv[vt <= 0] * sqrt(dt)

    v <- vt
    v[v<=0] <- 0
    vt[vt<=0] <- 0
    S <- S + (r - (v+vt)/4) * dt + rho * sqrt(v) * Zv * sqrt(dt) +
             (1/2) * (sqrt(v) + sqrt(vt)) * (Zs + rho * Zv) * sqrt(dt) +
             ((rho * sigma * dt)/2) * ((Zv)^2 - 1)
    S[S <= 0] = 0
    ST[j,] <- S
  }
  ST <- as.matrix(ST, ncol=N)
  Result <- exp(ST[nrow(ST),]) - X
  Result[Result <= 0] = 0
  call = mean(exp(-r*tau)*Result)

  lista = list('call' = call, 'Result' = Result, 'Spot' = ST)
  return(lista)
}
```

## C.3 Exact Algorithm

```
phi_heston <- function(a, v0, v_t, d){

  gamma_a <- sqrt(k^2 - 2 * sigma^2 * 1i*a)
  gammadt <- gamma_a * (tau-t)
  sqrtv0vt <- sqrt(v0*v_t)
  delta <- -k * (tau-t)

  part1 <- (gamma_a * exp(-(gamma_a - k)/2 * (tau-t)) * (1 - exp(delta)))/
           (k * (1- exp(- gammadt)))
```

```

part2 <- exp((v0+v_t)/(sigma^2) * ( (k * (1 + exp(delta)))/(1-exp(delta)) -
                                     (gamma_a * (1 + exp(- gammadt)))/(1-exp(- gammadt))))

part3 <- Bessel::BesselI(z = ((4 * gamma_a * sqrtv0vt)/(sigma^2) *
                             exp(- gammadt/2)/
                             (1 - exp(- gammadt))), nu = 0.5*d - 1) /
  Bessel::BesselI(z = ((4 * k * sqrtv0vt)/(sigma^2) * (exp(delta/2))/
                    (1-exp(delta))), nu = 0.5*d - 1)

return (part1 * part2 * part3)
}

intv <- function(n, cf, v_t){
  integrand <- function(x, phi = cf){

    f2 <- function(u){
      Im(phi(u) * exp(-1i * u * x)) /u
    }
    return(f2)
  }

  ## integrate to "cdf"
  F_x <- function (x) {
    y <- 0.5 - 1/pi * integrate(integrand(x), lower= 0, upper= 1000,
                                rel.tol = 0.001, stop.on.error = FALSE)$value
    return(y)
  }

  ## endsign

  endsign <- function(f, sign = 1) {
    b <- sign
    while (sign * f(b) < 0) b <- 10 * b
    return(b)
  }

  ## inversion

  low_bound = -Inf
  upp_bound = Inf
  invcdf <- function(u) {
    subcdf <- function(t) F_x(t) - u
    if (low_bound == -Inf)
      low_bound <- endsign(subcdf, -1)
    if (upp_bound == Inf)
      upp_bound <- endsign(subcdf)
    return(uniroot(subcdf, lower=low_bound, upper=upp_bound,
                  tol = 0.001220703)$root)
  }
  U <- stats::runif(n)
  sapply(U, invcdf)
}

```

```

}

hestonea_mod <- function(S, X, r, v, theta, rho, k, sigma, t = 0, tau = 1){

  d1 <- (4 * k * theta)/(sigma)^2
  c0 <- (sigma^2 * (1 - exp(-k*tau)))/(4*k)
  dt <- (tau-t)
  ST <- NULL

  # sampling V

  lambda <- (4*k*exp(-k*dt)*v)/(sigma^2 * (1-exp(-k*dt)))
  vt <- c0 * stats::rchisq(n = 1, df = d1, ncp = lambda)

  # Sampling int{V}

  phi <- function(a, v0=v, v_t=vt, d=d1){phi_heston(a, v0=v, v_t=vt, d=d1)}

  int_v <- intv(1, cf = phi, v_t=vt)
  # OR if you perform the drift interpolation scheme:
  # int_v <- dt * ((1/2) * v + (1/2) * vt)

  # Sampling int{v}dw
  int_vdw <- (1/sigma) * (vt - v - k * theta * dt + k * int_v)

  # Sampling S
  if( int_v >= 0){
    m <- log(S) + (r * (tau - t) - (1/2) * int_v + rho * int_vdw)
    std <- sqrt((1 - rho^2)) * sqrt(int_v)
    S <- exp(m + std * rnorm(1))
    v <- vt
    ST <- S
  } else {
    v <- vt
    ST <- rbind(ST,NA)}

  Result <- ST - X
  Result[Result <= 0] = 0
  call = exp(-r*tau)*Result
  lista = list('call' = call, 'Result' = Result, 'Spot' = ST)
  return(lista)
}

```



## C.4 Results with 20 steps

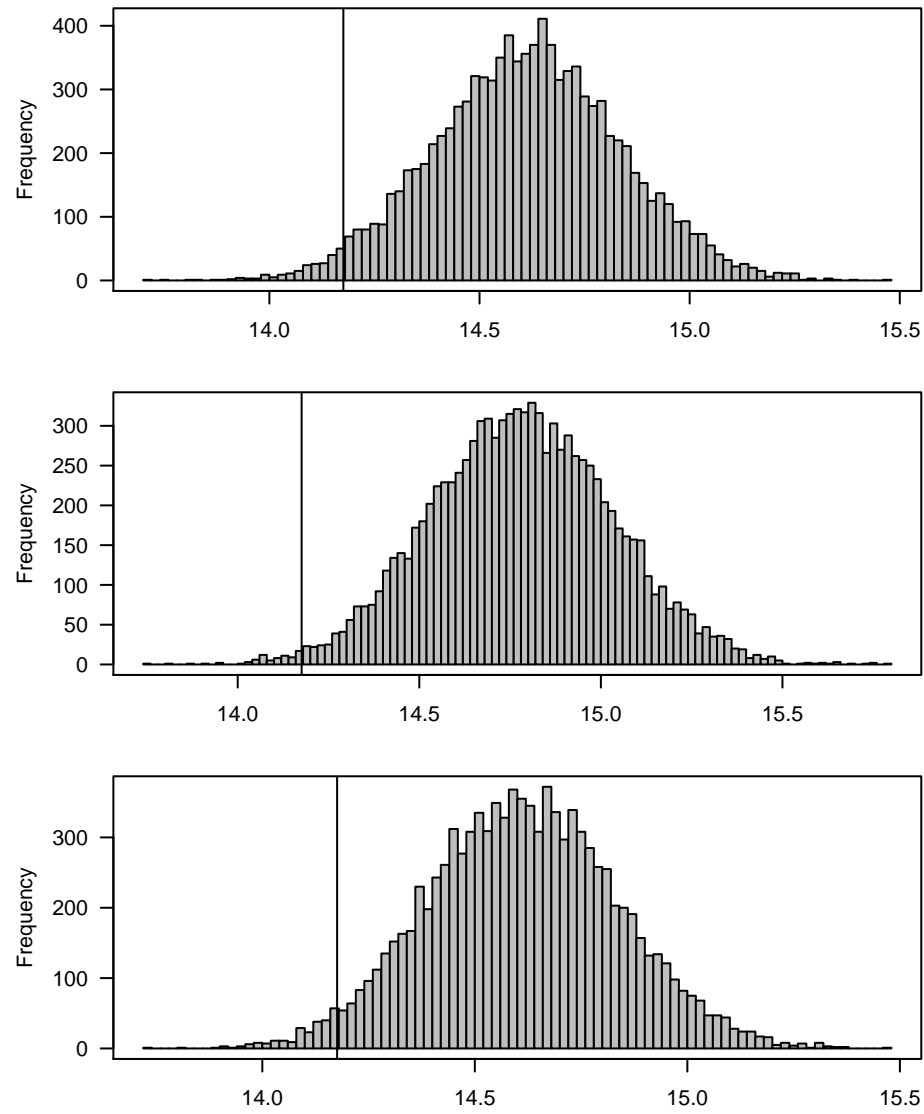


Figure C.1: Comparison between models, 20 steps.



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