

SPECIAL ISSUE PAPER**Incorporating boundary conditions in a stochastic volatility model for the numerical approximation of bond prices**Lourdes Gómez-Valle¹ | Miguel Ángel López-Marcos² | Julia Martínez-Rodríguez¹¹Departamento de Economía Aplicada e IMUVA, Facultad de Ciencias Económicas y Empresariales, Universidad de Valladolid, Spain²Departamento de Matemática Aplicada e IMUVA, Facultad de Ciencias, Universidad de Valladolid, Spain**Correspondence**

Julia Martínez-Rodríguez, Departamento de Economía Aplicada e IMUVA, Facultad de Ciencias Económicas y Empresariales, Universidad de Valladolid, Spain. Email: julia@eco.uva.es

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Summary

In this paper, we consider a two-factor interest rate model with stochastic volatility and we assume that the instantaneous interest rate follows a jump-diffusion process. In this kind of problems, a two-dimensional partial integro-differential equation is derived for the values of zero-coupon bonds. To apply standard numerical methods to this equation, it is customary to consider a bounded domain and incorporate suitable boundary conditions. However, for these two-dimensional interest rate models there are not well-known boundary conditions, unlike option models. Here, in order to approximate bond prices we propose new artificial boundary conditions which maintain the discount function property of the zero-coupon bond price. Then, we illustrate the numerical approximation of the corresponding boundary value problem by means of an Alternative Direction Implicit method which has been already applied for pricing options. We test these boundary conditions with several interest rate pricing models.

KEYWORDS:

interest rate model, zero-coupon bonds, jump-diffusion stochastic processes, stochastic volatility, risk-neutral measure, boundary conditions, Alternative Direction Implicit methods

1 | INTRODUCTION

The term structure of interest rates is one of the most important topics in the study of economics and finance. On the one hand, it provides useful information concerning market expectations of future real activity and inflation, that is, the conventional final targets of monetary policy. On the other hand, it is a very important subject in pricing models, risk management, hedging and arbitrage. In fact, many researches focus on the dynamic models of the term structure.

Traditionally, interest rates are assumed to move continuously and, in the literature, diffusion processes are used to model their behaviour^{10,27}. However, different market phenomena, such as shocks or surprises, provide unexpected jumps in the interest rates, as a result of this, the most realistic models use jump-diffusion stochastic processes to model interest rates^{11,22}. Moreover, pricing and hedging with jump-diffusion models is crucial, since ignoring jumps can produce hedging and pricing risk²³.

It is widely known that one-factor interest rate models are very attractive for practitioners because of its simplicity and computational convenience. However, these models have also unrealistic properties. Firstly, they cannot generate all the yield curve shapes and changes that we can find in the markets. Secondly, the changes over infinitesimal periods of any two interest-rate dependent prices will be perfectly correlated. Finally, as Hong and Li¹⁷ show, none of their analyzed one-factor models adequately captures the interest rate dynamics. Therefore, at least two factors are necessary in order to properly model the term

⁰**Abbreviations:** ADI method, Alternative Direction Implicit method; PIDE, partial integro-differential equation; HV method, Hundsdorfer-Verwer method; MHV method, Modified Hundsdorfer-Verwer method; MC method, Monte Carlo method

structure of interest rates. In fact, the number of factors must be a compromise between an efficient numerical implementation and the capability of the model to fit the experimental data.

In the literature, the different functions describing the stochastic variables of the model are usually chosen to be affine: the aim is to obtain a pricing problem with a feasible closed-form solution. However, when more realistic functions are considered¹⁵, a feasible expression of the solution is not achieved. Then, numerical methods should be used to approximate the solution. In such a case, we can underline two techniques. On the one hand, as zero-coupon bond prices can be obtained as expectations, the Monte Carlo approach provides a valuable tool: by the law of large numbers, the integrals that describe the expected value of the underlying random variables can be approximated by means of independent generated paths of such random variables. The Monte Carlo method is simple, flexible, and can handle multiple sources of uncertainty. However, this technique provides low accuracy approximations with a very high computational cost. On the other hand, we can also use numerical methods based on the discretization of the differential equation in order to approximate its solution¹². More specifically, if the model involves a jump-diffusion process, we have to deal with a partial integro-differential equation with a final condition.

From a numerical point of view, in order to discretize the problem, first of all a bounded domain for the state variables must be considered. Hence, suitable boundary conditions must be established. In bond pricing literature, only one-factor interest rate models are usually investigated. In this case, the restriction of the equation when the interest rate is zero and an artificial Dirichlet condition based on the behaviour of the solution for the maximum interest rate are incorporated^{5,8,13}. However, there is a lack of investigation when the number of state variables is greater than one.

If we analyse other financial derivatives, such as options, finite difference schemes are usually considered for pricing European and barrier options in jump-diffusion models⁷. For the multifactor case, first, a transformation in the partial integro-differential equation is usually made to eliminate the mixed derivative term. Then, several suitable methods are considered with very well-known and established boundary conditions^{14,19,20}. However, this is not applicable to general interest rate models.

In this paper, we focus on the approximation of zero-coupon bond prices in two-factor jump-diffusion models where the state variables are the instantaneous interest rate and the volatility. The numerical approach requires the discretization of a multidimensional problem involving a partial integro-differential equation: so we have to discretize some differential terms and a nonlocal integral term. As the problem has only a final condition, we introduce a bounded domain for the state variables. The main difficulty is to fix appropriate boundary conditions compatible with the model. Here we propose such restrictions taking into account the discount function property of the zero-coupon bond price. We illustrate the implementation of a numerical method for the proposed boundary value problem by using a well-known ADI method: the Hundsdorfer-Verwer scheme²⁰.

The paper is organized as follows. Section 2 describes a two-factor jump-diffusion model with stochastic volatility to price interest rate derivatives. In Section 3, we propose suitable boundary conditions for the numerical approximation of this kind of models, and we implement an ADI method for the emerging boundary value problem. In Section 4, we validate the numerical technique by means of the approximation of the yield curves in two well-known term structure models. Section 5 collects the conclusions.

2 | THE JUMP-DIFFUSION MODEL

We present a two-factor jump-diffusion model for pricing zero-coupon bonds. Define $(\Omega, \mathcal{F}, \{\mathcal{F}_t\}_{t \geq 0}, \mathcal{P})$ as a complete filtered probability space which satisfies the usual conditions, where $\{\mathcal{F}_t\}_{t \geq 0}$ is a filtration^{2,6,25}. We assume that the state variables are the dynamics of the instantaneous interest rate, r , and the volatility, V . In order to take into account the abrupt changes of the interest rates in the markets, we consider that such interest rate follows a jump-diffusion stochastic process. We suppose that the volatility is a diffusion process^{1,16}: r is right-continuous (cadlag⁶) and we denote the left limit $r(t-) = \lim_{z \uparrow t} r(z)$ (however, for notational clarity the pre-jump values $r(t-)$ will be added only when necessary to avoid confusion and otherwise, they will be assumed implied). Thus, these factors follow this joint stochastic process:

$$r(t) = r(0) + \int_0^t \mu_r(r(z), V(z)) dz + \int_0^t V(z) dW_r(z) + \int_0^t c(r(z-), V(z)) dq(z), \quad (1)$$

$$V(t) = V(0) + \int_0^t \mu_V(r(z), V(z)) dz + \int_0^t \sigma_V(r(z), V(z)) dW_V(z), \quad (2)$$

where μ_r and μ_V are the drifts and σ_V the volatility of the implied volatility process. Moreover, W_r and W_V are Wiener processes, and the impact of the jump is given by function c and the compound Poisson process $q(t) = \sum_{i=1}^{N(t)} X_i$, with jump times $(\tau_i)_{i \geq 1}$. The Poisson process is represented by $N(t)$ with intensity $\lambda(r, V)$; and X_1, X_2, \dots is a sequence of identically distributed random variables with a Normal probability distribution $\Pi, \mathcal{N}(0, \sigma_X)$. We assume that W_r, W_V are correlated with

$$[W_r, W_V](t) = \rho t,$$

but the jump size distribution and both standard Brownian motions are independent of N . We also assume that the jump magnitude and jump arrival times are uncorrelated with the diffusion parts of the processes. Lastly, we suppose that the functions $\mu_r, \mu_V, \sigma_V, \lambda$ and Π satisfy suitable regularity conditions^{3,24}.

As in previous works¹⁶, we assume that the market is arbitrage-free. Then, there exists an equivalent martingale measure, \mathcal{Q} -measure, which is known as the risk-neutral measure (see the extended Girsanov-type measure transformation^{4,26}). The state variables of the model (1)-(2) under the risk-neutral measure are as follows:

$$r(t) = r(0) + \int_0^t \mu_r^{\mathcal{Q}}(r(z), V(z)) dz + \int_0^t V(z) dW_r^{\mathcal{Q}}(z) + \int_0^t c(r(z-), V(z)) d\tilde{q}^{\mathcal{Q}}(z), \quad (3)$$

$$V(t) = V(0) + \int_0^t \mu_V^{\mathcal{Q}}(r(z), V(z)) dz + \int_0^t \sigma_V(r(z), V(z)) dW_V^{\mathcal{Q}}(z), \quad (4)$$

where $\mu_r^{\mathcal{Q}} = \mu_r - V\theta^{W_r}$, $\mu_V^{\mathcal{Q}} = \mu_V - \sigma_V\theta^{W_V}$. The Wiener processes under \mathcal{Q} -measure are $W_r^{\mathcal{Q}}$ and $W_V^{\mathcal{Q}}$, and $[W_r^{\mathcal{Q}}, W_V^{\mathcal{Q}}](t) = \rho t$. The market prices of risk associated to W_r and W_V Wiener processes are $\theta^{W_r}(r, V)$ and $\theta^{W_V}(r, V)$, respectively. The compensated compound Poisson process under \mathcal{Q} -measure is $\tilde{q}^{\mathcal{Q}}(t) = \sum_{i=1}^{N^{\mathcal{Q}}(t)} X_i - \lambda^{\mathcal{Q}} t E_X^{\mathcal{Q}}[X_1]$, and $\lambda^{\mathcal{Q}}(r, V)$ is the intensity of the Poisson process $N^{\mathcal{Q}}(t)$. Finally, we consider the function $c(r, V) = 1$ in (3).

A zero-coupon bond price at time t , with maturity time $T, t \leq T$, under the above assumptions, is

$$p(t, r, V; T) = E^{\mathcal{Q}} \left[\exp \left(- \int_t^T r(u) du \right) \middle| r(t) = r, V(t) = V \right]. \quad (5)$$

Trivially $p(T, r, V; T) = 1$. Moreover, the yield curve can be obtained as

$$R(t, r, V; T) = \frac{-\ln(p(t, r, V; T))}{T - t}. \quad (6)$$

Using the Feynman-Kac Theorem, it can be shown that $p(t, r, V; T)$ in (5) is the solution of the following partial integro-differential equation with final condition:

$$p_t + \mu_r^{\mathcal{Q}} p_r + \mu_V^{\mathcal{Q}} p_V + \frac{1}{2} V^2 p_{rr} + \frac{1}{2} \sigma_V^2 p_{VV} + \rho V \sigma_V p_{rV} - (r + \lambda^{\mathcal{Q}}) p + \lambda^{\mathcal{Q}} \int_{-\infty}^{\infty} p(t, r + y, V; T) f(y) dy = 0, \quad (7)$$

$$p(T, r, V; T) = 1, \quad (8)$$

for $r > 0, V > 0, 0 \leq t \leq T$, where f is the Normal density function

$$f(y) = \frac{1}{\sigma_X \sqrt{2\pi}} \exp \left(-\frac{y^2}{2\sigma_X^2} \right).$$

In order to solve numerically the problem (7)-(8) it is customary to make a change of the time variable, $\tau = T - t$, to obtain the following initial value problem

$$p_{\tau} = \mu_r^{\mathcal{Q}} p_r + \mu_V^{\mathcal{Q}} p_V + \frac{1}{2} V^2 p_{rr} + \frac{1}{2} \sigma_V^2 p_{VV} + \rho V \sigma_V p_{rV} - (r + \lambda^{\mathcal{Q}}) p + \lambda^{\mathcal{Q}} \int_{-\infty}^{\infty} p(\tau, r + y, V) f(y) dy, \quad (9)$$

$$p(0, r, V) = 1, \quad (10)$$

where we have reduced the notation as $p(\tau, r, V) = p(\tau, r, V; T)$.

With respect to the integral term in (9), which we denote as $J(\tau, r, V)$, we also make a change of variable obtaining the following expression

$$J(\tau, r, V) = \int_{-\infty}^{\infty} p(\tau, z, V) f(z - r) dz.$$

In most of the cases, a closed-form solution for the problem (9)-(10) is not available and then we have to consider numerical methods in order to approximate the solution.

3 | ARTIFICIAL BOUNDARY CONDITIONS AND NUMERICAL METHOD

Numerical techniques based on the discretization of the problem usually involve a bounded domain of the state variables. However, the problem (9)-(10) consists of a pure initial value problem. So, first, we need to restrict our attention to an appropriate portion of the domain and incorporate suitable boundary conditions to the model.

From a mathematical point of view, different boundary conditions are usual: Dirichlet, Newmann or Robin data. However, when we restrict our attention to an artificial bounded region, this kind of conditions can be inappropriate or inconsistent with the modelization. Another valuable process consists of discretizing the restriction of the equation to the boundary. However, here we propose a new kind of boundary conditions based on the financial meaning of the model, which are more efficient for this kind of problems.

In the classical Mathematical Finance literature, in a continuous framework, the discount function takes the form of a negative exponential function, and this is the case of the zero coupon bonds. In fact, the known standard solution for this kind of bond pricing problems¹, such as (7)-(8), has the form

$$p(t, r, V; T) = \exp(A(T - t) + B(T - t)r + C(T - t)V),$$

with A , B , and C as suitable functions satisfying $A(T) = B(T) = C(T) = 0$. As a consequence, the logarithm of the zero-coupon bond price should be linear with respect to the state variables. We want to mimic this behaviour in the general case. So, we propose the following boundary conditions for the problem (9)-(10) on the boundary of the truncated domain $[r_{min}, r_{max}] \times [V_{min}, V_{max}]$:

$$\frac{\partial^2 \ln p}{\partial r^2}(\tau, r_{min}, V) = 0, \quad \frac{\partial^2 \ln p}{\partial r^2}(\tau, r_{max}, V) = 0, \quad (11)$$

$$\frac{\partial^2 \ln p}{\partial V^2}(\tau, r, V_{min}) = 0, \quad \frac{\partial^2 \ln p}{\partial V^2}(\tau, r, V_{max}) = 0. \quad (12)$$

Therefore, we have to solve the problem (9)-(10) in the previous bounded domain with the boundary conditions (11)-(12).

In order to obtain the numerical solution of the pricing problem, first, we introduce a uniform meshgrid over the domain of the state variables $[r_{min}, r_{max}] \times [V_{min}, V_{max}]$. For positive integers m_r and m_V , we define the constant step sizes

$$h_r = \frac{r_{max} - r_{min}}{m_r},$$

$$h_V = \frac{V_{max} - V_{min}}{m_V}.$$

Then, the grid points are defined as:

$$r_i = r_{min} + ih_r, \quad i = 0, 1, \dots, m_r,$$

$$V_j = V_{min} + jh_V, \quad j = 0, 1, \dots, m_V.$$

Here, $P_{i,j}(\tau)$ represents an approximation to $p(\tau, r_i, V_j)$. We consider a semidiscretization of the PIDE (9) by means of approximations to the different terms at the inner grid points of the domain $[r_{min}, r_{max}] \times [V_{min}, V_{max}]$. For the partial derivatives of p with respect to r , we apply central difference formulas. For $i = 1, \dots, m_r - 1$, $j = 1, \dots, m_V - 1$,

$$\frac{\partial p_{i,j}(\tau)}{\partial r} \approx \frac{P_{i+1,j}(\tau) - P_{i-1,j}(\tau)}{2h_r},$$

$$\frac{\partial^2 p_{i,j}(\tau)}{\partial r^2} \approx \frac{P_{i+1,j}(\tau) - 2P_{i,j}(\tau) + P_{i-1,j}(\tau)}{h_r^2}.$$

Analogous finite differences are considered for the partial derivatives of p with respect to V . In the case of the mixed derivative $\partial^2 p / \partial r \partial V$, we use the following formula which is obtained by successive application of the standard central finite difference for the first derivative in the r and V variables

$$\frac{\partial^2 p_{i,j}(\tau)}{\partial r \partial V} \approx \frac{P_{i+1,j+1}(\tau) - P_{i-1,j+1}(\tau) + P_{i-1,j-1}(\tau) - P_{i+1,j-1}(\tau)}{4h_r h_V}, \quad i = 1, \dots, m_r - 1, j = 1, \dots, m_V - 1.$$

All these finite difference approximations are of second order.

To compute numerically the integral term, we reduce the region of integration to a bounded interval: here, we consider the same interval $[r_{min}, r_{max}]$. We approximate the integral by means of a quadrature rule⁷ with the step size h_r ,

$$J_{i,j}(\tau) = J(\tau, r_i, V_j) \approx \sum_{l=0}^{m_r} P_{l,j}(\tau) \int_{r_l - h_r/2}^{r_l + h_r/2} f(z - r_i) dz, \quad i = 1, \dots, m_r - 1, j = 1, \dots, m_V - 1.$$

If we collect all these discretizations, we obtain a system of ordinary differential equations for the time derivatives $P'_{i,j}(\tau)$, $i = 1, \dots, m_r - 1, j = 1, \dots, m_V - 1$, involving the values of $P_{i,j}(\tau)$, $i = 0, 1, \dots, m_r, j = 0, 1, \dots, m_V$. More precisely, we collect all the approximations $P_{i,j}(\tau)$, $i = 0, \dots, m_r, j = 0, \dots, m_V$, in a vector $\mathbf{P}(\tau)$ (we can use lexicographical order, for example). Then, we define the functions $\mathbf{F}^{(1)}$, $\mathbf{F}^{(2)}$, $\mathbf{F}^{(D)}$ and $\mathbf{F}^{(J)}$ with the components: for $i = 1, \dots, m_r - 1, j = 1, \dots, m_V - 1$,

$$\begin{aligned} F_{i,j}^{(1)}(\mathbf{P}) &= -\frac{1}{2}(r_i + \lambda^Q(r_i, V_j))P_{i,j} + \mu_r^Q(r_i, V_j) \frac{P_{i+1,j} - P_{i-1,j}}{2h_r} + \frac{1}{2}V_j^2 \frac{P_{i+1,j} - 2P_{i,j} + P_{i-1,j}}{h_r^2}, \\ F_{i,j}^{(2)}(\mathbf{P}) &= -\frac{1}{2}(r_i + \lambda^Q(r_i, V_j))P_{i,j} + \mu_V^Q(r_i, V_j) \frac{P_{i,j+1} - P_{i,j-1}}{2h_V} + \frac{1}{2}\sigma_V^2(r_i, V_j) \frac{P_{i,j+1} - 2P_{i,j} + P_{i,j-1}}{h_V^2}, \\ F_{i,j}^{(D)}(\mathbf{P}) &= \rho(r_i, V_j)V_j\sigma_V(r_i, V_j) \frac{P_{i+1,j+1} - P_{i-1,j+1} + P_{i-1,j-1} - P_{i+1,j-1}}{4h_r h_V}, \\ F_{i,j}^{(J)}(\mathbf{P}) &= \lambda^Q(r_i, V_j) \sum_{l=0}^{m_r} \left(\int_{r_l - h_r/2}^{r_l + h_r/2} f(z - r_i) dz \right) P_{l,j}. \end{aligned}$$

Now, $\mathbf{F} = \mathbf{F}^{(1)} + \mathbf{F}^{(2)} + \mathbf{F}^{(D)} + \mathbf{F}^{(J)}$, and therefore,

$$P'_{i,j}(\tau) = F_{i,j}(\mathbf{P}(\tau)), \quad i = 1, \dots, m_r - 1, j = 1, \dots, m_V - 1. \quad (13)$$

For the boundary points, we discretize the boundary conditions (11)-(12): we apply the forward (respectively, backward) finite difference approximations for the first (respectively, last) grid point. For example, for $r = r_0 = r_{min}$,

$$\frac{2 \ln P_{0,j}(\tau) - 5 \ln P_{1,j}(\tau) + 4 \ln P_{2,j}(\tau) - \ln P_{3,j}(\tau)}{h_r^2} = 0, \quad j = 0, \dots, m_V.$$

Using the exponential function we can explicit $P_{0,j}(\tau)$. Then, if we apply analogous approximations to each boundary, we obtain

$$P_{0,j}(\tau) = \sqrt{\frac{(P_{1,j}(\tau))^5 P_{3,j}(\tau)}{(P_{2,j}(\tau))^4}}, \quad P_{m_r,j}(\tau) = \sqrt{\frac{(P_{m_r-1,j}(\tau))^5 P_{m_r-3,j}(\tau)}{(P_{m_r-2,j}(\tau))^4}}, \quad j = 0, \dots, m_V, \quad (14)$$

$$P_{i,0}(\tau) = \sqrt{\frac{(P_{i,1}(\tau))^5 P_{i,3}(\tau)}{(P_{i,2}(\tau))^4}}, \quad P_{i,m_V}(\tau) = \sqrt{\frac{(P_{i,m_V-1}(\tau))^5 P_{i,m_V-3}(\tau)}{(P_{i,m_V-2}(\tau))^4}}, \quad i = 0, \dots, m_r. \quad (15)$$

Note that for the values $P_{0,0}$, P_{0,m_V} , $P_{m_r,0}$ and P_{m_r,m_V} we have two different expressions. In the practical implementation, we will choose only one.

Finally, we discretize the time variable. We assume a constant time step size $k = T/N$ and the temporal grid points $\tau_n = nk$, for $n = 0, 1, \dots, N$. Now, $P_{i,j}^n$ represents an approximation to $P_{i,j}(\tau_n)$. Starting from the grid restriction of the initial value (10), we obtain a numerical approximation at the new time level τ_{n+1} , from the approximations at the previous time level τ_n . To this end, we couple two different procedures. On the one hand, to compute the approximations at the inner grid points, we apply a numerical time integrator to the previous system of ordinary differential equations (13). On the other hand, to compute the approximations at the boundary grid points, we use the technique of the artificial boundary condition discretization which provides formulas (14)-(15).

For the numerical time integration, we use a suitable modification of the HV method adapted to this boundary problem: the MHV method. Note that the original HV method¹⁸ is well-established in the literature for option valuation²⁰.

Therefore, the general time step provides \mathbf{P}^{n+1} from \mathbf{P}^n for $n = 0, \dots, N - 1$, as follows

- Stage 1

$$Y_{i,j}^0 = P_{i,j}^n + kF_{i,j}(\mathbf{P}^n), \quad i = 1, \dots, m_r - 1, j = 1, \dots, m_V - 1, \quad (16)$$

- Stage 2

$$Y_{i,j}^1 = Y_{i,j}^0 + \theta kF_{i,j}^{(1)}(\mathbf{Y}^1 - \mathbf{P}^n), \quad i = 1, \dots, m_r - 1, j = 1, \dots, m_V - 1, \quad (17)$$

$$Y_{0,j}^1 = \sqrt{\frac{(Y_{1,j}^1)^5 Y_{3,j}^1}{(Y_{2,j}^1)^4}}, \quad Y_{m_r,j}^1 = \sqrt{\frac{(Y_{m_r-1,j}^1)^5 Y_{m_r-3,j}^1}{(Y_{m_r-2,j}^1)^4}}, \quad j = 0, \dots, m_V, \quad (18)$$

$$Y_{i,0}^1 = \sqrt{\frac{(Y_{i,1}^1)^5 Y_{i,3}^1}{(Y_{i,2}^1)^4}}, \quad Y_{i,m_V}^1 = \sqrt{\frac{(Y_{i,m_V-1}^1)^5 Y_{i,m_V-3}^1}{(Y_{i,m_V-2}^1)^4}}, \quad i = 1, \dots, m_r - 1, \quad (19)$$

- Stage 3

$$Y_{i,j}^2 = Y_{i,j}^1 + \theta kF_{i,j}^{(2)}(\mathbf{Y}^2 - \mathbf{P}^n), \quad i = 1, \dots, m_r - 1, j = 1, \dots, m_V - 1, \quad (20)$$

$$Y_{0,j}^2 = \sqrt{\frac{(Y_{1,j}^2)^5 Y_{3,j}^2}{(Y_{2,j}^2)^4}}, \quad Y_{m_r,j}^2 = \sqrt{\frac{(Y_{m_r-1,j}^2)^5 Y_{m_r-3,j}^2}{(Y_{m_r-2,j}^2)^4}}, \quad j = 0, \dots, m_V, \quad (21)$$

$$Y_{i,0}^2 = \sqrt{\frac{(Y_{i,1}^2)^5 Y_{i,3}^2}{(Y_{i,2}^2)^4}}, \quad Y_{i,m_V}^2 = \sqrt{\frac{(Y_{i,m_V-1}^2)^5 Y_{i,m_V-3}^2}{(Y_{i,m_V-2}^2)^4}}, \quad i = 1, \dots, m_r - 1, \quad (22)$$

- Stage 4

$$\tilde{Y}_{i,j}^0 = Y_{i,j}^0 + \frac{1}{2} kF_{i,j}(\mathbf{Y}^2 - \mathbf{P}^n), \quad i = 1, \dots, m_r - 1, j = 1, \dots, m_V - 1, \quad (23)$$

- Stage 5

$$\tilde{Y}_{i,j}^1 = \tilde{Y}_{i,j}^0 + \theta kF_{i,j}^{(1)}(\tilde{\mathbf{Y}}^1 - \mathbf{Y}^2), \quad i = 1, \dots, m_r - 1, j = 1, \dots, m_V - 1, \quad (24)$$

$$\tilde{Y}_{0,j}^1 = \sqrt{\frac{(\tilde{Y}_{1,j}^1)^5 \tilde{Y}_{3,j}^1}{(\tilde{Y}_{2,j}^1)^4}}, \quad \tilde{Y}_{m_r,j}^1 = \sqrt{\frac{(\tilde{Y}_{m_r-1,j}^1)^5 \tilde{Y}_{m_r-3,j}^1}{(\tilde{Y}_{m_r-2,j}^1)^4}}, \quad j = 0, \dots, m_V, \quad (25)$$

$$\tilde{Y}_{i,0}^1 = \sqrt{\frac{(\tilde{Y}_{i,1}^1)^5 \tilde{Y}_{i,3}^1}{(\tilde{Y}_{i,2}^1)^4}}, \quad \tilde{Y}_{i,m_V}^1 = \sqrt{\frac{(\tilde{Y}_{i,m_V-1}^1)^5 \tilde{Y}_{i,m_V-3}^1}{(\tilde{Y}_{i,m_V-2}^1)^4}}, \quad i = 1, \dots, m_r - 1, \quad (26)$$

- Stage 6

$$P_{i,j}^{n+1} = \tilde{Y}_{i,j}^1 + \theta kF_{i,j}^{(2)}(\mathbf{P}^{n+1} - \mathbf{Y}^2), \quad i = 1, \dots, m_r - 1, j = 1, \dots, m_V - 1, \quad (27)$$

$$P_{0,j}^{n+1} = \sqrt{\frac{(P_{1,j}^{n+1})^5 P_{3,j}^{n+1}}{(P_{2,j}^{n+1})^4}}, \quad P_{m_r,j}^{n+1} = \sqrt{\frac{(P_{m_r-1,j}^{n+1})^5 P_{m_r-3,j}^{n+1}}{(P_{m_r-2,j}^{n+1})^4}}, \quad j = 0, \dots, m_V, \quad (28)$$

$$P_{i,0}^{n+1} = \sqrt{\frac{(P_{i,1}^{n+1})^5 P_{i,3}^{n+1}}{(P_{i,2}^{n+1})^4}}, \quad P_{i,m_V}^{n+1} = \sqrt{\frac{(P_{i,m_V-1}^{n+1})^5 P_{i,m_V-3}^{n+1}}{(P_{i,m_V-2}^{n+1})^4}}, \quad i = 1, \dots, m_r - 1, \quad (29)$$

where θ is a positive parameter. In the numerical experiments, in the next section, we assume $\theta = 1 - \frac{\sqrt{2}}{2}$, as in the original work of in't Hout²¹.

Note that stages 2, 3, 5 and 6 are implicit: we have to compute intermediate values \mathbf{Y} by means of a nonlinear system that can be written as $\mathbf{Y} = \mathbf{G}(\mathbf{Y})$. Function \mathbf{G} collects, on the one hand, the equations linking the inner values $Y_{i,j}$, $i = 1, \dots, m_r - 1$, $j = 1, \dots, m_V - 1$ (see, for example, equation (17) in stage 2) with all the components of \mathbf{Y} . On the other hand, the equations related to the boundary values

$$\begin{aligned} Y_{0,j}, Y_{m_r,j}, \quad j = 0, \dots, m_V, \\ Y_{i,0}, Y_{i,m_V}, \quad i = 1, \dots, m_r - 1, \end{aligned}$$

(see, for example, equations (18)-(19) in stage 2). In order to solve it, we consider an iterative procedure: starting from an initial guess \mathbf{Y}^0 (for example, a suitable choice is $\mathbf{Y}^0 = \mathbf{P}^n$) then, we compute, for $\nu = 0, 1, \dots$

$$\mathbf{Y}^{\nu+1} = \mathbf{G}(\mathbf{Y}^\nu).$$

At each iteration $\nu = 0, 1, \dots$, we propose: first, computing the components of $\mathbf{Y}^{\nu+1}$ at the inner grid points

$$Y_{i,j}^{\nu+1} = G_{i,j}(\mathbf{Y}^\nu), \quad i = 1, \dots, m_r - 1, j = 1, \dots, m_V - 1;$$

second, computing the boundary values

$$\begin{aligned} Y_{0,j}^{\nu+1} = G_{0,j}(\mathbf{Y}^\nu), \quad Y_{m_r,j}^{\nu+1} = G_{m_r,j}(\mathbf{Y}^\nu), \quad j = 0, \dots, m_V, \\ Y_{i,0}^{\nu+1} = G_{i,0}(\mathbf{Y}^\nu), \quad Y_{i,m_V}^{\nu+1} = G_{i,m_V}(\mathbf{Y}^\nu), \quad i = 1, \dots, m_r - 1. \end{aligned}$$

This iteration is repeated until the convergence is achieved. In practice, until the norm of the difference between two consecutive iterants is sufficiently small (less than a prescribed tolerance).

4 | NUMERICAL EXPERIMENTS

In this section, we test the artificial boundary conditions and the MHV method proposed in Section 3 on several interest rate models. The first one possesses an explicit expression of its solution that we can use in order to compare the approximations. For the second one, we do not know a closed-form solution.

Test problem 1: We study the term structure model proposed by Andersen et al. ¹ that considers two state variables: the instantaneous interest rate, which follows a jump-diffusion process, and its volatility, which follows a diffusion process. It is the following affine model

$$\begin{aligned} dr &= (k_1(\mu - r) - \xi_1 V)dt + \sqrt{V} dW_r^Q + d \left(\sum_{i=1}^{N^Q(t)} X_i \right), \\ dv &= (k_2(\alpha - V) - \eta \xi_2 V)dt + \eta \sqrt{V} dW_V^Q, \end{aligned}$$

and the zero-coupon bond price is

$$p(t, r, V; T) = e^{A(T-t) + B(T-t)r + C(T-t)V},$$

(see the work of Andersen et al. ¹ for the expressions of A , B and C , and the estimated values of the parameters).

In order to approximate this solution with the MHV method, first, we change the time variable as in (9). Then, we have to apply the numerical method to the following PIDE

$$\begin{aligned} p_\tau &= (k_1(\mu - r) - \xi_1 V)p_r + (k_2(\alpha - V) - \eta \xi_2 V)p_V + \frac{1}{2}V p_{rr} + \frac{1}{2}\eta^2 V p_{VV} + \rho \eta V p_{rV} \\ &\quad - (r + \lambda^Q)p + \lambda^Q \int_{-\infty}^{\infty} p(\tau, r + y, V) f(y) dy, \end{aligned} \quad (30)$$

with the initial condition (10). Second, we introduce the artificial boundary conditions (11)-(12), then, previously we have to consider a truncated domain at the state variables. This domain window should be chosen wide enough to describe faithfully the dynamic of the model. However, if the window is too wide, the computational cost will increase, and the method will lose efficiency. Therefore, in the numerical experiments we make a balance between both difficulties to fix the domain window.

Taking into account the financial meaning of the state variables, we chose $[r_{min}, r_{max}] \times [V_{min}, V_{max}] = [-0.02, 0.48] \times [0, 0.5]$, which is suitable to represent the dynamics of the problem.

In our first experiment we approximate the zero-coupon bond prices for several maturities, then, we compare the computed yield values that we obtain by means of (6), with the exact ones, at the grid points. In the simulations, we take the same step sizes for both state variables and, to show the convergence, we consider successive refinements of the step sizes dividing them by two. In Table 1, we present the numerical errors for the yields at maturity $T = 10$ over the mesh grid: more specifically, we show the maximum error over the grid points (the maximum norm). We observe that there exist values of the step sizes which do not provide satisfactory approximations: as usual in the literature a relation between the step sizes is necessary. Note that in each column the errors do not change practically: we do not obtain an increase in the accuracy under time refinement. However, in each row the errors go to zero as the step in the state variables decrease. Therefore, the error mainly comes from the discretization of the state variables. In order to observe the order of convergence, it is usual to compare the errors in the same diagonal. In this table, we show the experimental order of convergence in brackets, and we observe that it is of the second order.

$k \setminus h_r = h_V$	1.2500e-1	6.2500e-2	3.1250e-2
5.0000e-2	1.0614e-02		
2.5000e-2	1.0594e-02		
1.2500e-2	1.0589e-02	2.8512e-03 (1.9)	
6.2500e-3	1.0588e-02	2.8485e-03 (1.9)	
3.1250e-3	1.0588e-02	2.8478e-03 (1.9)	7.4785e-04 (1.9)
1.5625e-3	1.0587e-02	2.8477e-03 (1.9)	7.4766e-04 (1.9)

TABLE 1 Errors for 10-year maturity yields obtained with the MHV method in test problem 1. In brackets, the experimental order of convergence comparing the diagonal entries.

Our next goal is to prove the efficiency of the MHV method. To this end, we measure this property by comparison of the accuracy offered with respect to the computational cost required. Again, $m_r = m_v$, that we denote as M .

In Figure 1, we present the results through a log-log efficiency chart where the vertical axis measures the errors in yields (in the maximum norm) and the horizontal axis corresponds to the cost (in CPU time). The stars joined by a solid line are the results of the method with $M = 4, 8, 12, 16, 20$ and $N = 100, 200, 300, 400, 500$.

On the other hand, we compare the results with those obtained with another numerical technique very common in this kind of problems. In the finance literature, when a closed-form solution of a PIDE like (30) is unknown, it is usually expressed as the conditional expectation (5) by the Feynman-Kac Theorem. Then, zero-coupon bond prices are usually approximated by means of MC method. This method approximates this expectation with a great number of paths of the state variables r and V . However, this method involves a high computational cost and produces a low accuracy. For the test problem, we have carried out many computations with different numbers of simulations, denoted by S . In Figure 1, the stars joined by the dashed line are the results with the MC method for $S = 2000, 4000, 16000, 52000, 60000$. We observe that for a similar computational cost, the MC method produces higher errors than the MHV method. In fact, in order to obtain 3 decimals of accuracy with this method we need a day of CPU time. However, with the MHV method this accuracy is obtained in less than a second of computation.

In order to analyze the advantages of the previous artificial boundary conditions, we also price the same zero-coupon bonds with the HV method considering other usual boundary conditions: the restriction of the PIDE to the boundary, that we discretize by means of backward and forward finite difference formulas.

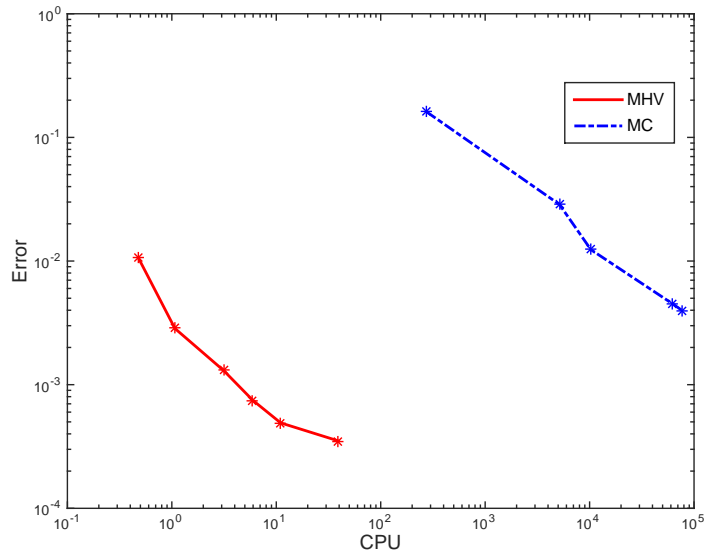


FIGURE 1 Log-log efficiency chart for the 10-year yield approximations with the MHV and MC methods in test problem 1.

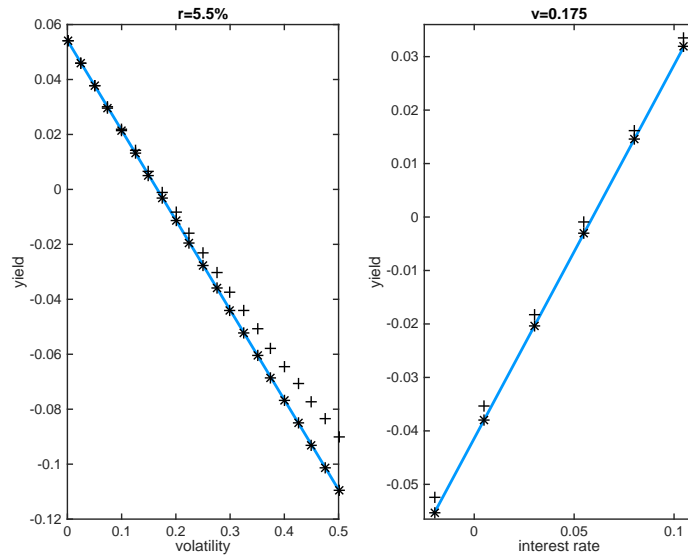


FIGURE 2 3-year yield profiles as function of the volatility when $r = 0.05$ (left), and the interest rate when $V = 0.175$ (right) with the HV and MHV methods in test problem 1.

Figure 2 , depicts some profiles of the 3-year yield. The left picture shows the yield as a function of the volatility along the whole interval, for a fixed value of the interest rate (in this case, $r = 0.05$: a common value of the interest rates in the financial market). The right picture plots the yield as a function of the interest rate only along a truncated interval with financial relevance, for a fixed value of the volatility (in this case, $V = 0.175$: a common value of the volatility in the financial market). The exact solution of the PIDE is represented by a solid line, the approximations with the usual boundary conditions by plus signs and the approximations using the new boundary conditions by asterisk signs. We observe that the approximations obtained with the

(M, N)	$\ \mathbf{R}(2M, 2N) - \mathbf{R}(M, N)\ _\infty$	(γ_h)	$\ \mathbf{R}(M, N) - \tilde{\mathbf{R}}(S)\ _\infty$
(4,40)			1.1244e-04
(8,80)	7.8211e-05		3.8774e-05
(16,160)	2.0571e-05	(1.9)	2.7688e-05
(32,320)	5.5835e-06	(1.9)	2.4761e-05

TABLE 2 Numerical results for the 10-year yields in test problem 2. Second column: comparison of consecutive approximations obtained with the MHV method (experimental order of convergence in brackets). Third column: comparison of the MHV approximations with the MC approximations ($S = 80000$).

MHV method are the closest to the exact yields, they are practically the same. On the one hand, from the left picture the higher the volatility is, the higher the difference. But, from the right picture, the differences between the exact yields and those obtained with the usual boundary conditions increase slightly when the interest rates decrease. We have observed in our computations that, for greater values of the maturity the different behaviour is more remarkable.

Test problem 2: We study the interest rate model proposed by Cotton et al. ⁹. In this case, we have two state variables, the interest rate and another variable which provides the volatility of the interest rates. However, following (3), we add a jump term in the interest rate stochastic process. Then, the stochastic processes in this model are

$$dr = (k_1(\mu - r) - \xi g(s))dt + g(s)dW_r^Q + d\left(\sum_{i=1}^{N(\lambda^Q)} X_i\right),$$

$$ds = k_2(\alpha - s)dt + \eta\sqrt{Y}dW_s^Q.$$

In this case, we do not know a closed-form solution in order to compare the numerical approximations obtained by means of the MHV method. For the numerical experiments, we consider the parameter values estimated in Cotton et al. ⁹, and we assume that the volatility of the interest rate is the positive function $g(s) = e^s$. Then, for pricing the corresponding zero-coupon bonds, we have to solve the PIDE:

$$p_\tau = (k_1(\mu - r) - \xi e^s)p_r + k_2(\alpha - s)p_s + \frac{1}{2}e^{2s}p_{rr} + \frac{1}{2}\eta^2 s p_{ss} + \rho\eta s p_{rs} - (r + \lambda^Q)p + \lambda^Q \int_{-\infty}^{\infty} p(\tau, r + y, s)f(y) dy,$$

with the initial conditions (10).

We apply the proposed MHV method: we take the same truncated domain used in the previous test problem. We calculate the zero-coupon bond prices with a maturity equal to 10 years. In order to analyze the convergence of the approximations, taking into account that the solution is unknown, we compute the experimental order of convergence by means of the comparison between two consecutive prices obtained by the refinement of the step sizes. Denoting by $\mathbf{R}(2M, 2N)$ the approximation at the meshgrid of the yield obtained by the MHV method with $m_r = m_V = M$, for the state variables, and N , for the time variable, we compute the experimental order of convergence in the standard way

$$\gamma_h = \log_2 \frac{\|\mathbf{R}(4M, 4N) - \mathbf{R}(2M, 2N)\|_\infty}{\|\mathbf{R}(2M, 2N) - \mathbf{R}(M, N)\|_\infty}. \quad (31)$$

Second column of Table 2 shows the results, which are compatible with the expected order of convergence observed in the previous test problem. Finally, we compare the approximations provided by the MHV method with those computed with the MC method, denoted by $\tilde{\mathbf{R}}(S)$, where S is the amount of simulations. Last column of Table 2 offers this comparison when $S = 80000$, a sufficiently large number to assure that the approximation is hard to improve with the MC method. We observe the convergence of the MHV approximations to the same values obtained with the MC method.

All the implementations have been done using MATLAB.

5 | CONCLUSIONS

In the finance literature, interest rate models are usually considered to be affine because a closed-form solution is always known. If the functions of the stochastic processes are assumed to be more realistic, an exact solution of the model is, in general, unknown. In such cases, numerical methods are the most important way to approximate the solution. On the one hand, the MC method is one of the most common techniques in the zero-coupon bond pricing. Although, it is very expensive from a computational point of view and not very accurate. On the other hand, other kind of methods, such as finite difference methods, require a bounded region for the state variables. However, this kind of problems typically involves a pure initial value problem.

When the interest rate is the single state variable, different numerical methods have been proposed for solving the pricing equation, jointly with well-established boundary conditions. Nevertheless, this is not the situation when there are two or more state variables: the usual conditions for the one-factor models expand the error from the boundary to all the integration interval.

In this paper, we focus on approximating zero-coupon bond prices in a two-factor jump-diffusion model, where the state variables are the interest rates and the volatility. The problem consists of a partial integro-differential equation with a final condition. Then, for its numerical approximation, we introduce artificial boundaries in the domain that must be discretized. Here, we propose new boundary conditions based on the discount function properties of the bond prices. We discretize these boundary conditions and show how to incorporate them in an ADI method in order to approximate the solution.

An extensive numerical simulation in two different interest rate models shows that this new numerical procedure is a very valuable tool in the approximation of this kind of financial derivatives: it provides approximations of second order, is more efficient than the MC method and fits the profile of the solutions better than other artificial boundary conditions. Besides, this technique can be adapted easily to other numerical methods.

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