

ON THE SINGULAR LIMIT OF SOLUTIONS TO THE COX–INGERSOLL–ROSS INTEREST RATE MODEL WITH STOCHASTIC VOLATILITY

BEÁTA STEHLÍKOVÁ AND DANIEL ŠEVČOVIČ

In this paper we are interested in term structure models for pricing zero coupon bonds under rapidly oscillating stochastic volatility. We analyze solutions to the generalized Cox–Ingersoll–Ross two factors model describing clustering of interest rate volatilities. The main goal is to derive an asymptotic expansion of the bond price with respect to a singular parameter representing the fast scale for the stochastic volatility process. We derive the second order asymptotic expansion of a solution to the two factors generalized CIR model and we show that the first two terms in the expansion are independent of the variable representing stochastic volatility.

Keywords: Cox–Ingersoll–Ross two factors model, rapidly oscillating volatility, singular limit of solution, asymptotic expansion

AMS Subject Classification: 35C20 35B25 62P05 60H10 35K05

1. INTRODUCTION

Term structure models describe a functional dependence between the time to maturity of a discount bond and its present price. Yield of bonds, as a function of maturity, forms the so-called term structure of interest rates. If we denote by $P = P(t, T)$ the price of a bond paying no coupons at time t with maturity at T then the term structure of yields $R(t, T)$ is given by

$$P(t, T) = e^{-R(t, T)(T-t)}, \quad \text{i. e. } R(t, T) = -\frac{\log P(t, T)}{T-t}$$

(cf. Kwok [7]). Continuous interest rate models are often formulated in terms of stochastic differential equations (SDEs) for the instantaneous interest rate (or short rate) as well as SDEs for other relevant quantities like e. g. volatility of the short rate process. In one-factor models there is a single stochastic differential equation for the short rate. The volatility of the short rate process is given in a deterministic way. It is assumed to be constant (the Vašíček model [9]) or it is a function of the short rate itself. In the classical Cox, Ingersoll, and Ross model (CIR) the short rate

is modelled by a solution to the following stochastic differential equation:

$$dr = \kappa(\theta - r) dt + \sigma\sqrt{r} dw, \quad (1)$$

where $\kappa, \theta, \sigma > 0$ are parameters representing the rate of reversion, the long term interest rate and volatility of the interest rate, respectively (see [2]). By dw we have denoted the differential of the Wiener process. Beside these two simple models there is a wide range of other term structure models including, in particular, the Chan–Karolyi–Longstaff–Sanders model [3], the Hull–White model [6] and many others. Based on the assumption made on the form of the short rate process one can derive a linear scalar parabolic equation for the bond price as a function of the current short rate and time to maturity.

In the two-factor models there are two sources of uncertainty yielding different term structures for the same short rate as they may depend on the value of the other factor. Moreover, two-factor models have a richer variety of possible shapes of term structures including, in particular, nonmonotone yield curves. The reader is referred to the books by Kwok [7] and Brigo and Mercurio [1] for detailed discussion on two-factor interest rate modeling.

There are several ways how to incorporate the second stochastic factor. It is reasonable to conjecture that in a financial market the volatility of a fluctuating underlying process for the short rate can be fluctuating as well. In the so-called two-factor models with a stochastic volatility we allow the volatility to have a stochastic behavior driven by another stochastic differential equation. As a consequence of the multidimensional Itô's lemma the corresponding equation for the bond price is a linear parabolic equation in two space dimensions. These spatial dimensions correspond to the short rate and volatility. It is well known that the density distribution of a stochastic process is a solution to the Fokker–Planck partial differential equation and can be expressed analytically in the case the volatility undergoes the Bessel square root process (see e. g. [5]). The exact value for the stochastic volatility is not known in the market. We can just however estimate its density distribution function and, consequently, its statistical moments like e. g. the mean value, volatility, skewness of the volatility etc. Knowing the density distribution of the stochastic volatility we are able to perform averaging of the bond price and the term structure with respect to the stochastic volatility. Unlike the short rate which is known from the market data on daily basis, as it was already mentioned, the volatility of the short rate process is unknown. Therefore such a volatility averaging is of special importance for practitioners.

The main goal of the paper is to derive an asymptotic expansion of the bond price with respect to a singular parameter representing the fast scale for the stochastic volatility process. We derive the second order asymptotic expansion of a solution to the two factors generalized CIR model and we show that the first two terms in the expansion are independent of the stochastic volatility term.

The paper is organized as follows. In the next section we present an empirical evidence of a short rate process for which the volatility is fluctuating and it has two concentration values. Next we discuss a model for statistical distribution capturing such a volatility clustering. Section 3 is devoted to the asymptotic analysis of

solutions to the bond pricing equation in the case when the fluctuating volatility is rapidly oscillating. We derive explicit formulae for the first three terms in the asymptotic expansion of a solution with respect to a small parameter representing the fast time scale for rapidly oscillating volatility.

2. EMPIRICAL EVIDENCE OF EXISTENCE OF VOLATILITY CLUSTERS AND THEIR MODELING

The key feature of the CIR modeling consists of the assumption made on constant volatility of the stochastic process (1) driving the short rate r . However, in real financial markets we can observe a substantial deviation from this assumption. To provide an empirical evidence for such a volatility process, we computed maximum Nowman's type of Gaussian estimates of the dispersion for the CIR model for 20-day-long intervals using three months treasury bills data. Figure 1 (left) shows the estimated dispersion as a function of time. Higher and lower volatility periods can be distinguished. They can be seen also on the kernel density estimates of the values in Figure 1 (right).

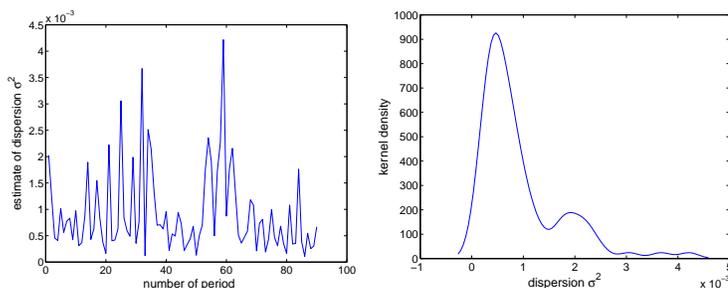


Fig. 1. Left: Estimates of CIR model's dispersion σ^2 from 20-day intervals (3-months Treasury bills, 90 intervals starting in January 1990). Right: the density distribution of estimates of the dispersion σ^2 .

In order to capture such a behavior of the dispersion σ^2 we shall consider a model in which the limiting density of the dispersion (as $t \rightarrow \infty$) has two local maxima. It corresponds to the so-called volatility clustering phenomenon where the dispersion can be observed in the vicinity of two local maxima of the density distribution (see [8]). The desired behavior of the process and its limiting density are shown in the Figure 2. A natural candidate for such a volatility process is

$$dy = \alpha(y) dt + \omega(y) dw \quad (2)$$

having a drift function $\alpha(y)$ such that the differential equation $\frac{dy}{dt} = \alpha(y)$ has two stable stationary solutions. With added stochastic part $\omega(y) dw$ of the process, these stationary solutions become values, around which the volatility concentrates. Recall that the cumulative distribution function $\tilde{G} = \tilde{G}(y, t) = \text{Prob}(y(t) < y | y(0) = y_0)$ of the process $y = y(t)$ satisfying (2) and starting almost surely from the initial datum

y_0 can be obtained from a solution $\tilde{g} = \partial\tilde{G}/\partial y$ to the so-called Fokker–Planck equation for the density function:

$$\frac{\partial\tilde{g}}{\partial t} = \frac{1}{2} \frac{\partial^2}{\partial y^2} (\omega(y)^2 \tilde{g}) - \frac{\partial}{\partial y} (\alpha(y)\tilde{g}), \quad \tilde{g}(y, 0) = \delta(y - y_0) \tag{3}$$

(cf. Goodman [5]). Here $\delta(y - y_0)$ denotes the Dirac delta function located at y_0 . The limiting density $g(y) = \lim_{t \rightarrow \infty} \tilde{g}(y, t)$ of the process is therefore a stationary solution to the Fokker–Planck equation (3) and it forgets any information about the initial datum y_0 , i. e.

$$L_0^* g \equiv \frac{1}{2} \frac{\partial^2}{\partial y^2} (\omega(y)^2 g) - \frac{\partial}{\partial y} (\alpha(y)g) = 0. \tag{4}$$

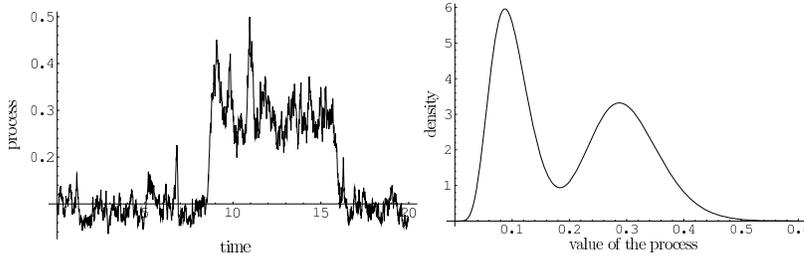


Fig. 2. Simulation of a process (left) and its asymptotic distribution (right).

In [8] one of the authors proposed a model with a property that the limiting density is a combination of two Gamma densities. Indeed, let us consider the following two mean reverting Bessel square root stochastic processes:

$$dy_i = \kappa_y (\theta_i - y_i) dt + v\sqrt{y_i} dw_i, \quad i = 1, 2, \tag{5}$$

where $\theta_i > 0$, $2\kappa_y \theta_i > v^2 > 0$ for $i = 1, 2$, and dw_1, dw_2 are uncorrelated differentials of the Wiener processes. Solving the stationary Fokker–Planck equation (4) it turns out that the limiting distributions of the processes y_1, y_2 are the Gamma distributions with shape parameters $2\kappa_y/v^2$ and $2\kappa_y\theta_i/v^2$. Denote their densities by g_1 and g_2 . Then $g_i(y) = C_i y^{\frac{2\kappa_y\theta_i}{v^2}-1} \exp(-\frac{2\kappa_y}{v^2}y)$ for $y > 0$ and $g_i(y) = 0$ otherwise. Here $C_i > 0$ is a normalization constant such that $\int_{\mathbb{R}} g_i(y) dy = 1$. Choose a parameter $k \in (0, 1)$. Our aim is to construct a process with asymptotic density

$$g(y) = kg_1(y) + (1 - k)g_2(y), \tag{6}$$

corresponding to a convex mixture of densities g_1 and g_2 . In the following theorem we see that for the same square root volatility function of the form $v\sqrt{y}$ it is possible to achieve this goal. Drift of the process $\alpha(y)$ can be written as a weighted sum of drifts $\alpha_i(y) = \kappa(\theta_i - y)$, $i = 1, 2$, with the weights depending on y .

Theorem 1. (Stehlíková [8, Section 5]) Suppose that the drift term α has the form: $\alpha(y) = w(y)\alpha_1(y) + (1 - w(y))\alpha_2(y)$ where $w(y) = kg_1(y)/(kg_1(y) + (1 - k)g_2(y))$ and $\alpha_i(y) = \kappa(\theta_i - y)$. Then the stochastic process driven by the SDE: $dy = \alpha(y)dt + v\sqrt{y}dw$ has the limiting distribution g given by the convex combination (6) of densities g_1, g_2 .

3. GENERALIZED CIR MODEL WITH RAPIDLY OSCILLATING STOCHASTIC VOLATILITY AND ITS ASYMPTOTIC ANALYSIS

The aim of this section is to provide a tool for modeling the effects of rapidly oscillating stochastic volatility that can be observed in real markets. If the length of the time scale for dispersion y is denoted by ε , the equation for the variable y reads as follows:

$$dy = \frac{\alpha(y)}{\varepsilon} dt + \frac{v\sqrt{y}}{\sqrt{\varepsilon}} dw_y. \quad (7)$$

In what follows we will assume that $0 < \varepsilon \ll 1$ is a small singular parameter. Notice that the limiting density function g of the stochastic process driven by SDE (7) is independent of the scaling parameter $\varepsilon > 0$. The statement follows directly from the stationary Fokker–Planck equation (4). Concerning structural assumption made on the drift function $\alpha : R \rightarrow R$ we shall henceforth assume the following hypothesis:

$$(A) \quad \alpha \text{ is a } C^1 \text{ function on } [0, \infty), \quad \frac{2\alpha(0)}{v^2} > 1, \quad \limsup_{y \rightarrow \infty} \frac{\alpha(y)}{y} < 0.$$

Now it is straightforward computation to verify the following auxiliary lemma.

Lemma 1. Let the drift function $\alpha(y)$ be defined as a mixture of two Gamma limiting distributions as in Theorem 1. Then the function α satisfies the hypothesis (A) with $\alpha(0) = \kappa \min(\theta_1, \theta_2)$ and $\limsup_{y \rightarrow \infty} \frac{\alpha(y)}{y} = -\kappa < 0$.

Next we shall show the limiting density g of the process driven by SDE (7) is uniquely given by the following lemma:

Lemma 2. Under the hypothesis (A) made on the drift function α the stationary Fokker–Planck equation $L_0^*g \equiv \frac{v^2}{2} \frac{\partial^2}{\partial y^2}(yg) - \frac{\partial}{\partial y}(\alpha(y)g) = 0$ has a unique solution g such that $g(0) = 0$ for $y \leq 0$. It can be explicitly expressed as:

$$g(y) = Cy^{-1} \exp\left(\frac{2}{v^2} \int_1^y \frac{\alpha(\xi)}{\xi} d\xi\right) = Cy^{\frac{2\alpha(0)}{v^2}-1} \exp\left(\frac{2}{v^2} \int_1^y \hat{\alpha}(\xi) d\xi\right)$$

for $y > 0$ and $g(y) = 0$ for $y \leq 0$. Here $\hat{\alpha}(y) = (\alpha(y) - \alpha(0))/y$ and $C > 0$ is a normalization constant such that $\int_0^\infty g(y) dy = 1$.

Proof. It follows by direct verification of the equation. The other linearly independent solution g_2 to the equation (4) has a nontrivial limit $g_2(0+) > 0$. \square

In what follows, we denote by σ^2 , $D > 0$, and S the limiting mean value, dispersion and skewness of the stochastic process for the y -variable representing stochastic dispersion, i. e.

$$\sigma^2 = \int_0^\infty yg(y) dy, \quad D = \int_0^\infty (y - \sigma^2)^2 g(y) dy, \quad S = \frac{1}{D^{\frac{3}{2}}} \int_0^\infty (y - \sigma^2)^3 g(y) dy. \tag{8}$$

Notice that $D = -\int_0^\infty \int_0^y (\xi - \sigma^2)g(\xi) d\xi dy$. In the generalized CIR model with a stochastic volatility, the instantaneous interest rate (short rate) r will be modelled by the mean reverting process of the form (1) where the volatility of is replaced by a square root of a stochastic dispersion y , i. e.

$$dr = \kappa(\theta - r) dt + \sqrt{y}\sqrt{r} dw_r. \tag{9}$$

The differentials of the Wiener processes dw_y and dw_r are assumed to be uncorrelated throughout the paper, i. e. $E(dw_y dw_r) = 0$. Then the corresponding partial differential equation for the bond price $P^\varepsilon = P^\varepsilon(t, r, y)$ has the following form:

$$\begin{aligned} \frac{\partial P^\varepsilon}{\partial t} &+ (\kappa(\theta - r) - \tilde{\lambda}_1(y, r)r^{\frac{1}{2}}\sqrt{y})\frac{\partial P^\varepsilon}{\partial r} + \frac{1}{2}ry\frac{\partial^2 P^\varepsilon}{\partial r^2} - rP^\varepsilon \\ &+ \frac{1}{\sqrt{\varepsilon}}\left(-\tilde{\lambda}_2(y, r)v\sqrt{y}\frac{\partial P^\varepsilon}{\partial y}\right) + \frac{1}{\varepsilon}\left(\alpha(y)\frac{\partial P^\varepsilon}{\partial y} + \frac{v^2y}{2}\frac{\partial^2 P^\varepsilon}{\partial y^2}\right) = 0, \end{aligned} \tag{10}$$

$(t, r, y) \in Q_T \equiv (0, T) \times \mathbb{R}^+ \times \mathbb{R}^+$, where $\tilde{\lambda}_1, \tilde{\lambda}_2$ are the so-called market prices of risk (cf. [7, Chapter 7]). By a solution P^ε to (11) we mean a bounded function $P^\varepsilon \in C^{1,2}(Q_T) \cap C(\bar{Q}_T)$ satisfying equation (11) on \bar{Q}_T . Concerning the structural form of market prices of risk functions $\tilde{\lambda}_1, \tilde{\lambda}_2$ we shall suppose that

$$\tilde{\lambda}_1(t, r, y) = \lambda_1\sqrt{r}\sqrt{y}, \quad \tilde{\lambda}_2(t, r, y) = \lambda_2\sqrt{y},$$

where $\lambda_1, \lambda_2 \in \mathbb{R}$ are constants. It is worthwhile noting that the latter assumption is not restrictive as the original one-factor CIR model assumes such a form of the market price of risk (cf. Kwok [7]). We shall rewrite PDE (11) in the operator form:

$$(\varepsilon^{-1}\mathcal{L}_0 + \varepsilon^{-1/2}\mathcal{L}_1 + \mathcal{L}_2)P^\varepsilon = 0, \tag{11}$$

where the linear differential operators $\mathcal{L}_0, \mathcal{L}_1, \mathcal{L}_2$ are defined as follows:

$$\begin{aligned} \mathcal{L}_0 &= \alpha(y)\frac{\partial}{\partial y} + \frac{v^2y}{2}\frac{\partial^2}{\partial y^2}, \quad \mathcal{L}_1 = -\lambda_2vy\frac{\partial}{\partial y}, \\ \mathcal{L}_2 &= \frac{\partial}{\partial t} + (\kappa(\theta - r) - \lambda_1ry)\frac{\partial}{\partial r} + \frac{1}{2}ry\frac{\partial^2}{\partial r^2} - r. \end{aligned} \tag{12}$$

Next we expand the solution P^ε into Taylor power series:

$$P^\varepsilon(t, r, y) = \sum_{j=0}^\infty \varepsilon^{\frac{j}{2}} P_j(t, r, y), \tag{13}$$

with the terminal conditions $P_0(T, r, y) = 1, P_j(T, r, y) = 0$ for $j \geq 1$ at expiry $t = T$. The main goal of this paper is to examine the singular limiting behavior of a solution P^ε as $\varepsilon \rightarrow 0^+$. More precisely, we shall determine the first three terms P_0, P_1, P_2 of the asymptotic expansion (13). We shall henceforth denote by $\langle \psi \rangle$ the averaged value of the function $\psi \in C([0, \infty))$ with respect to the density g , i. e. $\langle \psi \rangle = \int_0^\infty \psi(y)g(y) dy$. We shall also use the notation $\langle \mathcal{L}_2 \rangle$ standing for the averaged linear operator \mathcal{L}_2 , i. e.

$$\langle \mathcal{L}_2 \rangle \equiv \frac{\partial}{\partial t} + (\kappa(\theta - r) - \lambda_1 r \sigma^2) \frac{\partial}{\partial r} + \frac{1}{2} \sigma^2 r \frac{\partial^2}{\partial r^2} - r. \quad (14)$$

Lemma 3. Let $\psi \in C^1([0, \infty))$ be such that $\mathcal{L}_0 \psi$ is bounded. Then $\langle \mathcal{L}_0 \psi \rangle = 0$.

Proof. Notice that the operator \mathcal{L}_0^* is the adjoint operator to the linear operator \mathcal{L}_0 with respect to the L^2 -inner product $(\psi, \phi) = \int_0^\infty \psi(y)\phi(y) dy$. It means that $\langle \mathcal{L}_0 \psi \rangle = (\mathcal{L}_0 \psi, g) = (\psi, \mathcal{L}_0^* g) = 0$ because the density g is a solution to Eq. (4). \square

The following lemma will be useful when computing higher order term in series expansion (13).

Lemma 4. Let $F \in C([0, \infty))$ be such that $\langle F \rangle = 0$. Then, up to an additive constant, there exists a unique solution $\psi \in C^2((0, \infty)) \cap C([0, \infty))$ to the non-homogeneous equation $\mathcal{L}_0 \psi = \frac{v^2}{2} F$. Its derivative $\frac{\partial \psi}{\partial y}$ is given by

$$\frac{\partial \psi}{\partial y}(y) = \frac{1}{yg(y)} \int_0^y F(\xi)g(\xi) d\xi.$$

Moreover, $\langle \mathcal{L}_1 \psi \rangle = \lambda_2 v \int_0^\infty F(y) yg(y) dy$. In particular, if ψ is a solution to the equation $\mathcal{L}_0 \psi = 0$ then ψ is a constant function with respect to the y -variable.

Proof. Using equation (4) for the limiting density g and inserting $\frac{\partial \psi}{\partial y}$ into the operator \mathcal{L}_0 we obtain that ψ is a solution to the equation $\mathcal{L}_0 \psi = \frac{v^2}{2} F$. Other independent solutions are not continuous at $y = 0$. The formula for $\langle \mathcal{L}_1 \psi \rangle$ follows from the definition of the operator \mathcal{L}_1 by applying integration by parts formula. \square

Now we proceed with collecting the terms of the power series expansion of (11).

- In the order ε^{-1} we have $\mathcal{L}_0 P_0 = 0$. According to Lemma 4 we have $P_0 = P_0(t, r)$, i. e. P_0 is independent of the y -variable.
- In the order $\varepsilon^{-1/2}$ we have $\mathcal{L}_0 P_1 + \mathcal{L}_1 P_0 = 0$. Since $P_0 = P_0(t, r)$ we deduce $\mathcal{L}_1 P_0 = 0$ and so $\mathcal{L}_0 P_1 = 0$. By Lemma 4, $P_1 = P_1(t, r)$ is independent of y .
- In the order ε^0 we have $\mathcal{L}_0 P_2 + \mathcal{L}_1 P_1 + \mathcal{L}_2 P_0 = 0$. Since $P_1 = P_1(t, r)$ we have $\mathcal{L}_1 P_0 = 0$. Hence $\mathcal{L}_0 P_2 + \mathcal{L}_2 P_0 = 0$. Taking the average $\langle \cdot \rangle$ of both sides of the latter equation we obtain $\langle \mathcal{L}_0 P_2 \rangle + \langle \mathcal{L}_2 P_0 \rangle = 0$. By Lemma 3 and the fact that P_0 is independent of y -variable we conclude $\langle \mathcal{L}_2 \rangle P_0 = \langle \mathcal{L}_2 P_0 \rangle = 0$. Therefore P_0 is a

solution to the classical one-factor PDE equation for the CIR model satisfying the terminal condition $P_0(T, r) = 1$ for any $r \geq 0$. It is well known that the solution $P_0 = P_0(t, r)$ to the equation $\langle \mathcal{L}_2 \rangle P_0 = 0$ is given by the explicit formula:

$$P_0(t, r) = A_0(t)e^{-B(t)r}, \quad (15)$$

where $A'_0 = \kappa\theta A_0 B$ and $B' = (\kappa + \lambda_1\sigma^2)B + \frac{\sigma^2}{2}B^2 - 1$, $A_0(T) = 1$, $B(T) = 0$, i. e.

$$A_0(t) = \left(\frac{2\phi e^{(\phi+\psi)(T-t)/2}}{(\phi+\psi)(e^{\phi(T-t)} - 1) + 2\phi} \right)^{\frac{2\kappa\theta}{\sigma^2}}, \quad B(t) = \frac{2(e^{\phi(T-t)} - 1)}{(\phi+\psi)(e^{\phi(T-t)} - 1) + 2\phi},$$

$\psi = \kappa + \lambda_1\sigma^2$, $\phi = \sqrt{\psi^2 + 2\sigma^2}$ (cf. Kwok [7, Chapter 7]). Since $\langle \mathcal{L}_2 \rangle P_0 = 0$ we have $-\mathcal{L}_2 P_0 = (\langle \mathcal{L}_2 \rangle - \mathcal{L}_2) P_0 = (\sigma^2 - y)f(t)re^{-B(t)r}$ where

$$f(t) = \left(\lambda_1 B(t) + \frac{1}{2}B(t)^2 \right) A_0(t).$$

Hence $\mathcal{L}_0 P_2 = -\mathcal{L}_2 P_0 = (\sigma^2 - y)f(t)re^{-B(t)r}$. According to Lemma 4 we have

$$\frac{\partial P_2}{\partial y} = -\frac{2}{v^2}f(t)re^{-B(t)r}H(y), \quad H(y) = \frac{1}{yg(y)} \int_0^y (\xi - \sigma^2)g(\xi) d\xi. \quad (16)$$

• In the order $\varepsilon^{1/2}$ we have $\mathcal{L}_0 P_3 + \mathcal{L}_1 P_2 + \mathcal{L}_2 P_1 = 0$. Since $\langle \mathcal{L}_0 P_3 \rangle = 0$ we have $\langle \mathcal{L}_1 P_2 \rangle + \langle \mathcal{L}_2 P_1 \rangle = 0$. The function $P_1 = P_1(t, r)$ is independent of the y -variable and therefore $\langle \mathcal{L}_2 \rangle P_1 = \langle \mathcal{L}_2 P_1 \rangle = -\langle \mathcal{L}_1 P_2 \rangle$. By Lemma 4 we have

$$\mathcal{L}_1 P_2 = \frac{2\lambda_2}{v}f(t)re^{-B(t)r}yH(y), \quad -\langle \mathcal{L}_1 P_2 \rangle = K_1 f(t)re^{-B(t)r},$$

where $K_1 = -\frac{2\lambda_2}{v} \int_0^\infty \int_0^y (\xi - \sigma^2)g(\xi) d\xi dy = \frac{2\lambda_2}{v}D$ is a constant (see (8)). Notice that the constant K_1 and the function $f(t)$ depend on the first two moments σ^2 and D of the stochastic dispersion only. Equation $\langle \mathcal{L}_2 \rangle P_1 = \langle \mathcal{L}_2 P_1 \rangle = -\langle \mathcal{L}_1 P_2 \rangle$ reads as:

$$\frac{\partial P_1}{\partial t} + (\kappa(\theta - r) - \lambda_1 r \sigma^2) \frac{\partial P_1}{\partial r} + \frac{1}{2}r\sigma^2 \frac{\partial^2 P_1}{\partial r^2} - rP_1 = K_1 f(t)re^{-B(t)r}. \quad (17)$$

The solution P_1 satisfying the terminal condition $P_1(T, r) = 0$ for $r \geq 0$ can be found in the closed form:

$$P_1(t, r) = (A_{10}(t) + A_{11}(t)r) e^{-B(t)r}, \quad (18)$$

where the functions $A_{10}(t)$, $A_{11}(t)$ are solutions to the system of linear ODEs:

$$\begin{aligned} A'_{11}(t) &= (\kappa\theta B(t) + \kappa + \lambda_1\sigma^2 + \sigma^2 B(t)) A_{11}(t) + K_1 f(t), \\ A'_{10}(t) &= \kappa\theta B(t)A_{10}(t) - \kappa\theta A_{11}(t), \end{aligned} \quad (19)$$

with terminal conditions $A_{10}(T) = 0$, $A_{11}(T) = 0$. We can analytically and also numerically compute A_{10} , A_{11} in a fast and accurate manner. This way we have

obtained the term $P_1(t, r)$. In Figure 3 examples of numerical approximation of the term structure $R^\varepsilon(T - \tau, r) = -\frac{1}{\tau} \log \langle P^\varepsilon(T - \tau, r, \cdot) \rangle$ corresponding to the second order expansion of the averaged value of $\langle P^\varepsilon(t, r, \cdot) \rangle$, $P^\varepsilon(t, r, y) \approx P_0(t, r) + \sqrt{\varepsilon} P_1(t, r)$. We plot term structures starting from the short rate $r = 0.03$ (left) and $r = 0.031$ (right) for parameters $\kappa = 5$, $\theta = 0.03$, $\kappa_y = 100$, $v = 1.1832$, $\theta_1 = 0.025$, $\theta_2 = 0.1$, $k = 1/3$, $\lambda_1 = -1$, $\lambda_2 = -100$ and $\varepsilon = 0, 0.001, 0.01$ (black, red and blue curves).

Having P_1 and $\frac{\partial P_2}{\partial y}$ we can compute the term $\mathcal{L}_1 P_2 + \mathcal{L}_2 P_1$. With regard to Lemma 4 equation $\mathcal{L}_0 P_3 = -\mathcal{L}_1 P_2 - \mathcal{L}_2 P_1$ then yields a formula for $\frac{\partial P_3}{\partial y}$ and

$$\langle \mathcal{L}_1 P_3 \rangle = -\frac{2\lambda_2}{v} \left(\left(\frac{2\lambda_2}{v} K_3 + K_1 \sigma^2 \right) f(t) r e^{-B(t)r} + D \left(-\lambda_1 r \frac{\partial P_1}{\partial r} + \frac{r}{2} \frac{\partial^2 P_1}{\partial r^2} \right) \right),$$

where the constant $K_3 = \int_0^\infty \xi^3 H(\xi) g(\xi) d\xi = -\frac{1}{2} S D^{\frac{3}{2}} - \sigma^2 D$ depends on the first three statistical moments of the stochastic dispersion.

• In the order ε^1 we have $\mathcal{L}_0 P_4 + \mathcal{L}_1 P_3 + \mathcal{L}_2 P_2 = 0$. Proceeding similarly as before we have $\langle \mathcal{L}_0 P_4 \rangle = 0$ and therefore

$$\langle \mathcal{L}_1 P_3 \rangle + \langle \mathcal{L}_2 P_2 \rangle = 0. \quad (20)$$

We decompose the function $P_2(t, r, y)$ in the form

$$P_2(t, r, y) = \bar{P}_2(t, r) + \tilde{P}_2(t, r, y), \quad (21)$$

where \bar{P}_2 is the averaged value of P_2 and \tilde{P}_2 is a zero mean fluctuation, i. e. $\langle \tilde{P}_2 \rangle = 0$. As \bar{P}_2 does not depend on y , we have $\frac{\partial \bar{P}_2}{\partial y} = \frac{\partial P_2}{\partial y}$. Taking into account $\langle \mathcal{L}_2 \tilde{P}_2 \rangle = 0$ we obtain

$$\tilde{P}_2(t, r, y) = -\frac{2}{v^2} f(t) r e^{-B(t)r} \left(\int_0^y H(\xi) d\xi - K_2 \right)$$

where $K_2 = \int_0^\infty g(s) \int_0^s H(\xi) d\xi ds$ is a constant and the function H is given by (16). Now we can use decomposition (21) to evaluate $\langle \mathcal{L}_2 P_2 \rangle$. We have $\langle \mathcal{L}_2 P_2 \rangle = \langle \mathcal{L}_2 (\bar{P}_2 + \tilde{P}_2) \rangle = \langle \mathcal{L}_2 \bar{P}_2 \rangle + \langle \mathcal{L}_2 \tilde{P}_2 \rangle$ because \bar{P}_2 is independent of y . Next we can determine $\langle \mathcal{L}_2 \tilde{P}_2 \rangle$ in the following form:

$$\langle \mathcal{L}_2 \tilde{P}_2 \rangle = -\frac{2}{v^2} K_4 f(t) r \left(-\lambda_1 \frac{\partial}{\partial r} + \frac{1}{2} \frac{\partial^2}{\partial r^2} \right) (r e^{-B(t)r})$$

where $K_4 = \int_0^\infty \int_0^y H(\xi) d\xi (y - \sigma^2) g(y) dy$. It is worthwhile noting that both constants K_2, K_4 depend on all nontrivial statistical moments of the stochastic dispersion. Equation (20) then becomes

$$\langle \mathcal{L}_2 \bar{P}_2 \rangle = -\langle \mathcal{L}_2 \tilde{P}_2 \rangle - \langle \mathcal{L}_1 P_3 \rangle = (a(t) + b(t)r + c(t)r^2) e^{-B(t)r}, \quad \bar{P}_2(T, r) = 0,$$

which is a partial differential equation for $\bar{P}_2 = \bar{P}_2(t, r, y)$ with a right hand side which can be explicitly computed from already obtained results in the closed form:

$$\bar{P}_2(t, r) = (A_{20}(t) + A_{21}(t)r + A_{22}(t)r^2) e^{-B(t)r}, \quad (22)$$

where the functions A_{20}, A_{21}, A_{22} are solutions to a linear system of ODEs. We omit details here.

In summary we have shown the following main result of this paper:

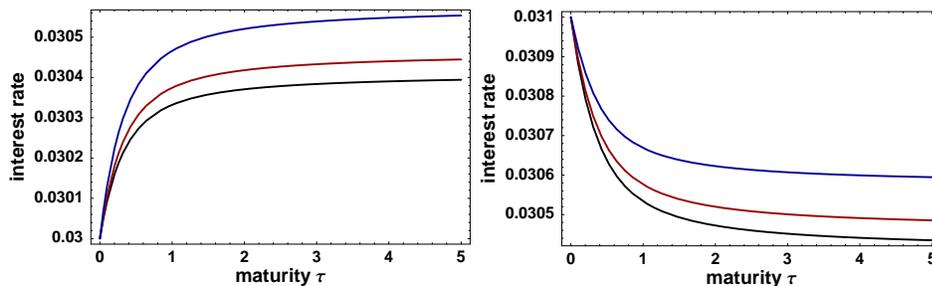


Fig. 3. The approximate term structure $R^\varepsilon = R^\varepsilon(T - \tau, r)$ based on the first two leading terms of the bond price $P^\varepsilon \approx P_0(T - \tau, r) + \sqrt{\varepsilon}P_1(T - \tau, r)$ starting from the short rate $r = 0.03$ (left) and $r = 0.031$ (right) for several values of the singular parameter $\varepsilon = 0, 0.001, 0.01$ (black, red and blue curves), resp.

Theorem 2. The solution $P^\varepsilon = P^\varepsilon(t, r, y)$ of the generalized CIR bond pricing equation (11) with rapidly oscillating dispersion can be approximated, for small values of the singular parameter $0 < \varepsilon \ll 1$, by $P^\varepsilon(t, r, y) \approx P_0(t, r) + \sqrt{\varepsilon}P_1(t, r) + \varepsilon P_2(t, r, y) + O(\varepsilon^{\frac{3}{2}})$.

The first two terms P_0, P_1 are independent of the y -variable representing unobserved stochastic volatility. They depend only on the first two statistical moments (mean value and dispersion) of the stochastic dispersion and other model parameters.

The next term in the expansion P_2 nontrivially depends on the y -variable. P_2 as well as its averaged value $\langle P_2 \rangle$ depends also on all nontrivial statistical moments of the stochastic dispersion.

The terms P_0, P_1, P_2 can be evaluated by closed-form formulae (15), (18), (22).

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*Beáta Stehlíková and Daniel Ševčovič, Department of Applied Mathematics and Statistics, Comenius University, Mlynská dolina, 84248 Bratislava. Slovak Republic.
e-mails: stehlikova,sevcovic@fmph.uniba.sk*