HITTING TIME FOR BESSEL PROCESSES—WALK ON MOVING SPHERES ALGORITHM (WOMS)¹

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In this article we investigate the hitting time of some given boundaries for Bessel processes. The main motivation comes from mathematical finance when dealing with volatility models, but the results can also be used in optimal control problems. The aim here is to construct a new and efficient algorithm in order to approach this hitting time. As an application we will consider the hitting time of a given level for the Cox–Ingersoll–Ross process. The main tools we use are on one side, an adaptation of the method of images to this particular situation and on the other side, the connection that exists between Cox–Ingersoll–Ross processes and Bessel processes.

1. Introduction. The aim of this paper is to study the hitting time of some curved boundaries for the Bessel process. Our main motivations come from mathematical finance, optimal control and neuroscience. In finance Cox-Ingersoll-Ross processes are widely used to model interest rates. As an application, in this article we will consider the simulation of the first hitting time of a given level for the CIR by using its relation with the Bessel process. In neuroscience the firing time of a neuron is usually modelled as the hitting time of a stochastic process associated with the membrane potential behavior; for introduction of noise in neuron systems, see Part I Chapter 5 in [7]. The literature proposes different continuous stochastic models like, for instance, the family of integrate-and-fire models; see Chapter 10 in [6]. Most of them are related to the Ornstein–Uhlenbeck process which appears in a natural way as extension of Stein's model, a classical discrete model. In Feller's model, generalized Bessel processes appear as a more realistic alternative to the Ornstein–Uhlenbeck process; see, for instance, [10] for a comparison of these models. Therefore the interspike interval, which is interpreted as the first passage time of the membrane potential through a given threshold is closely related to the first hitting time of a curved boundary for some Bessel processes.

Our main results and the main algorithm are obtained for the case of Bessel processes. We use in our numerical method the particular formula that we obtain for

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the hitting time of some curved boundaries for the Bessel process and the connection that exists between a Bessel process and the Euclidean norm of a Brownian motion when calculating the hitting position. As an application we consider the hitting time of a given level for the Cox–Ingersoll–Ross process. In order to obtain this, we use first of all the connections that exist between CIR processes and Bessel processes and second, the method of images for this particular situation.

The study of Bessel processes and their hitting times occupies a huge portion of mathematical literature. Let us only mention few of them: Göing-Jaeschke and Yor [8] consider a particular case of CIR processes which are connected with radial Ornstein–Uhlenbeck processes and their hitting times; L. Alili and P. Patie [1] investigate as a special situation the Bessel processes via some boundary crossing identities for diffusions having the time inversion property; recently, Salminen and Yor considered the hitting time of affine boundaries for the 3-dimensional Bessel process [16].

In a recent paper Hamana and Matsumoto [9] gave explicit expressions for the distribution functions and the densities of the first hitting time of a given level for the Bessel process. Their results cover all the cases. Let us also mention a recent work of Byczkowski, Malecki and Ryznar [2]. By using an integral formula for the density of the first hitting time of a given level of the Bessel process, they are able to obtain uniform estimates of the hitting time density function.

In all these papers the formulae are explicit and are hard to use for a numerical purposes as they exhibit Bessel functions. The main idea of this paper is to get rid of this difficulty by using two important tools: first of all the method of images that allow us to obtain, for some particular boundaries, an explicit form for the density of the hitting time, and second, the connection between δ -dimensional Bessel processes and the Euclidean norm of a δ -dimensional Brownian motion in order to get the simulated exit position. By coupling these ingredients we are able to construct a numerical algorithm that is easy to implement and very efficient and which approaches the desired hitting time.

We will use here a modified version of the *random walk on spheres* method which was first introduced by Muller [13] in 1956. This procedure allows us to solve a Dirichlet boundary value problem. The idea is to simulate iteratively, for the Brownian motion, the exit position from the largest sphere included in the domain and centered in the starting point. This exit position becomes the new starting point and the procedure goes on until the exit point is close enough to the boundary. Let us notice that the simulation of the exit time from a sphere is numerically costly.

The method of images was introduced in 1969 by Daniels [4] as a tool to construct nonlinear boundaries for which explicit calculations for the exit distribution for the Brownian motion are possible. The method was developed also in Lerche [11]. We adapt this method for the Bessel process by using the explicit form of its density. For some particular curved boundaries we can explicitly evaluate the density of the Bessel hitting time.

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The paper is organized as follows. First we present some new results on hitting times for Bessel processes. Second, we construct the new algorithm for approaching the hitting time, the so called *walk on moving spheres algorithm*. Finally we present some numerical results and as a particular application the evaluation of the hitting time of a given level for the Cox–Ingersoll–Ross process.

2. Hitting time for Bessel processes. Bessel processes play an important role both in the study of stochastic processes like Brownian motion and in various theoretical and practical applications as, for example, in finance.

Let us consider the δ -dimensional Bessel process starting from *y*, the solution of the following stochastic differential equation:

(2.1)
$$\begin{cases} Z_t^{\delta, y} = Z_0^{\delta, y} + \frac{\delta - 1}{2} \int_0^t (Z_s^{\delta, y})^{-1} \, \mathrm{d}s + B_t, \\ Z_0^{\delta, y} = y, \qquad y \ge 0, \end{cases}$$

where $(B_t)_{t\geq 0}$ is a one-dimensional Brownian motion. We denote

(2.2)
$$v = \frac{\delta}{2} - 1,$$

the *index* of this process. We call δ the *dimension* of the process. This terminology is coming from the fact that, in the case of positive integer $\delta \in \mathbb{N}$, a δ -dimensional Bessel process can be represented as the Euclidean norm of a δ -dimensional Brownian motion. This will be a key point in our numerical method.

The density of this process starting from y is given by

(2.3)
$$p_y(t,x) = \frac{x}{t} \left(\frac{x}{y}\right)^v \exp\left(-\frac{x^2 + y^2}{2t}\right) I_v\left(\frac{xy}{t}\right) \quad \text{for } t > 0, y > 0, x \ge 0,$$

where $I_{\nu}(z)$ is the Bessel function whose expression gives

(2.4)
$$I_{\nu}(z) = \sum_{n=0}^{\infty} \left(\frac{z}{2}\right)^{\nu+2n} \frac{1}{n! \Gamma(\nu+n+1)}.$$

When starting from y = 0, the density of $Z_t^{\delta,0}$ is

(2.5)
$$p_0(t,x) = \frac{1}{2^{\nu}} \frac{1}{t^{\nu+1}} \frac{1}{\Gamma(\nu+1)} x^{\delta-1} \exp\left(-\frac{x^2}{2t}\right) \quad \text{for } t > 0, x \ge 0.$$

2.1. The method of images for Bessel processes. In this section, we investigate the first hitting time of a curved boundary for the Bessel process starting from the origin. Let $\psi(t)$ denote the boundary, and introduce the following stopping time:

(2.6)
$$\tau_{\psi} = \inf\{t > 0; Z_t^{\delta, 0} \ge \psi(t)\}.$$

For some suitable choice of the boundary, the distribution of τ_{ψ} can be explicitly computed. The idea is given by the following remark on the method of images (see, e.g., [4] for the origin of this method and [11] for a complete presentation):

Fundamental idea. Suppose that F is a positive σ -finite measure satisfying some integrability conditions (to be specified later on), and define

(2.7)
$$u(t,x) = p_0(t,x) - \frac{1}{a} \int_{\mathbb{R}_+} p_y(t,x) F(\mathrm{d}y)$$

for some real constant a > 0. Let

$$\psi(t) = \inf \{ x \in \mathbb{R}; u(t, x) < 0 \} \qquad \text{for all } t > 0$$

Then u(t, x) is solution of the partial differential equation

(2.8)
$$\begin{cases} \frac{\partial u}{\partial t}(t,x) = \frac{1}{2} \frac{\partial^2 u}{\partial x^2}(t,x) - \frac{\delta - 1}{2} \frac{\partial}{\partial x} \left(\frac{1}{x} u(t,x)\right), & \text{on } \mathbb{R}_+ \times \mathbb{R}, \\ u(t,\psi(t)) = 0, & \text{for all } t > 0, \\ u(0,\cdot) = \delta_0(\cdot), & \text{on } (-\infty,\psi(0+)]. \end{cases}$$

From this remark we deduce an interesting expression for the hitting time. We can prove that

$$\tau_{\psi} = \inf\{t > 0; u(t, Z_t^{\delta, 0}) = 0\}.$$

This means simply that in order to obtain information on the hitting time it suffices to look for $u(t, Z_t^{\delta, 0}) = 0$. Let us express this in a general result.

THEOREM 2.1. Let F(dy) be a positive σ -finite measure such that $\int_0^\infty p_0(t, t)$ $\sqrt{\varepsilon} v$) $F(dv) < \infty$ for all $\varepsilon > 0$. Let a > 0 and define the function

(2.9)
$$u(t,x) = p_0(t,x) - \frac{1}{a} \int_{\mathbb{R}_+} p_y(t,x) F(\mathrm{d}y)$$

Consider $\psi(t)$ such that $u(t, \psi(t)) = 0$. Then the probability density function of τ_{ψ} is given by

(2.10)
$$\mathbb{P}_{0}(\tau_{\psi} \in dt) = \left[-\frac{1}{2} \frac{\partial u}{\partial x}(t, x) \Big|_{x=\psi(t)} + \frac{1}{2} \frac{\partial u}{\partial x}(t, x) \Big|_{x=0} - \frac{\delta - 1}{2x} u(t, x) \Big|_{x=0} \right] dt.$$

PROOF. We will only point out the main ideas for the proof in this case as it follows mainly the ideas introduced in [11]. A complete description of the method and this result for the Brownian motion case can be found in [11].

Let us consider

(2.11)
$$u(t,x) = p_0(t,x) - \frac{1}{a} \int_{\mathbb{R}_+} p_y(t,x) F(\mathrm{d}y),$$

where F(dy) is a measure on \mathbb{R}_+ . We consider $\psi(t)$ the solution of $u(t, \psi(t)) = 0$. Let us define as before $\tau_{\psi} = \inf\{t \ge 0; Z_t^{\delta,0} \ge \psi(t)\}$. Then $u(t, x) dx = \mathbb{P}(\tau_{\psi} > t, Z_t^{\delta,0} \in dx)$ and

(2.12)
$$\mathbb{P}_0(\tau_{\psi} > t) = \int_0^{\psi(t)} u(t, x) \, \mathrm{d}x.$$

In order to get the distribution of τ_{ψ} we have to evaluate the derivative of $\mathbb{P}_0(\tau_{\psi} > t)$. By using equality (2.12) we obtain

$$\mathbb{P}_{0}(\tau_{\psi} \in dt)$$

$$(2.13) = \left(-\psi'(t)u(t,\psi(t)) - \int_{0}^{\psi(t)} \frac{\partial u}{\partial t}(t,x) dx\right) dt$$

$$= \left(-\frac{1}{2} \int_{0}^{\psi(t)} \frac{\partial^{2} u}{\partial x^{2}}(t,x) dx + \frac{\delta - 1}{2} \int_{0}^{\psi(t)} \frac{\partial}{\partial x} \left(\frac{1}{x}u(t,x)\right) dx\right) dt,$$

as u(t, x) is solution of partial differential equation (2.8). We thus obtain

(2.14)

$$\mathbb{P}_{0}(\tau_{\psi} \in dt) = \left(-\frac{1}{2}\frac{\partial u}{\partial x}(t,x)\Big|_{x=\psi(t)} + \frac{\delta - 1}{2\psi(t)}u(t,\psi(t)) + \left(\frac{1}{2}\frac{\partial u}{\partial x}(t,x) - \frac{\delta - 1}{2x}u(t,x)\right)\Big|_{x=0}\right)dt$$

As $\frac{\delta - 1}{2\psi(t)}u(t, \psi(t)) = 0$, and this ends the proof of the theorem. \Box

The idea behind the method of images is that for some particular forms of F(dy), we can derive explicit formulae of the hitting time distribution. More precisely:

PROPOSITION 2.2. Let us denote, for $\delta = 2\nu + 2 > 0$ and a > 0 by

$$\operatorname{Supp}(\tau_{\psi}) = \left[0, \left(\frac{a}{\Gamma(\nu+1)2^{\nu}}\right)^{1/(\nu+1)}\right].$$

We define, for all $t \in \text{Supp}(\tau_{\psi})$ *, the function*

(2.15)
$$\psi_a(t) = \sqrt{2t \log \frac{a}{\Gamma(\nu+1)t^{\nu+1}2^{\nu}}}.$$

Then the probability density of τ_{ψ} has its support in $\text{Supp}(\tau_{\psi})$ and is given by

(2.16)
$$\mathbb{P}_0(\tau_{\psi} \in \mathrm{d}t) = \frac{1}{2at} \left(2t \log \frac{a}{\Gamma(\nu+1)t^{\nu+1}2^{\nu}} \right)^{\nu+1} \mathbb{1}_{\mathrm{Supp}(\tau_{\psi})}(t) \,\mathrm{d}t.$$

PROOF. By using the expression in (2.3) we remark first that

(2.17)
$$y^{2\nu+1}p_y(t,x) = x^{2\nu+1}p_x(t,y)$$

Let us consider, as in Theorem 2.1,

(2.18)
$$u(t,x) = p_0(t,x) - \frac{1}{a} \int_{\mathbb{R}_+} p_y(t,x) F(\mathrm{d}y),$$

with $F(dy) = y^{2\nu+1} \mathbb{1}_{\{y>0\}} dy$. In this situation the function *u* defined in (2.18) gives

(2.19)
$$u(t,x) = p_0(t,x) - \frac{1}{a}x^{2\nu+1} \\ = \left(\frac{1}{2^{\nu}}\frac{1}{t^{\nu+1}}\frac{1}{\Gamma(\nu+1)}\exp\left(-\frac{x^2}{2t}\right) - \frac{1}{a}\right)x^{2\nu+1}.$$

For simplicity we will write ψ instead of ψ_a . Following the result in the Theorem 2.1, we are looking for $\psi(t)$ such that $u(t, \psi(t)) = 0$. This yields

(2.20)
$$x = \psi(t) = \sqrt{2t \log \frac{a}{\Gamma(\nu+1)t^{\nu+1}2^{\nu}}}$$

under the obvious condition $t^{\nu+1} \leq \frac{a}{\Gamma(\nu+1)2^{\nu}}$.

We can now notice that

$$p_0(t, \psi(t)) = \frac{1}{a} (\psi(t))^{2\nu+1}$$

and we can prove easily that

$$\frac{\partial u}{\partial x}(t,x) = (\delta - 1)\frac{u(t,x)}{x} - \frac{x}{t}p_0(t,x).$$

We obtain, after replacing in (2.10) and after applying the Theorem 2.1, for this particular case,

$$\mathbb{P}_0(\tau_{\psi} \in \mathrm{d}t) = \frac{1}{2t}\psi(t)p_0(t,\psi(t))\,\mathrm{d}t$$
$$= \frac{1}{2at}\psi^{2\nu+2}(t)\,\mathrm{d}t$$
$$= \frac{1}{2at}\left(2t\log\frac{a}{\Gamma(\nu+1)t^{\nu+1}2^{\nu}}\right)^{\nu+1}\,\mathrm{d}t,$$

which gives the desired result. \Box

The second boundary which allows us to express explicit results is obtained by using the Markov property for the Bessel process.

PROPOSITION 2.3. Let us, for $\delta = 2\nu + 2 > 0$, s > 0 and a > 0 fixed, denote by

$$\operatorname{Supp}(\tau_{\psi}) = \begin{cases} [0, +\infty), & \text{for } a \ge 1, \\ \left[0, \frac{s}{(1/a)^{1/(\nu+1)} - 1}\right], & \text{for } 0 < a < 1. \end{cases}$$

For $t \in \text{Supp}(\tau_{\psi})$ *we define the function*

 $\mathbb{P}_0(\tau_{\psi} \in \mathrm{d}t)$

(2.21)
$$\psi_a(t) = \sqrt{\frac{2t(t+s)}{s}} \left[(\nu+1) \log\left(1+\frac{s}{t}\right) + \log a \right]$$

Then the probability density function of the hitting time τ_{ψ} is given by

(2.22)
$$= \frac{1}{\Gamma(\nu+1)} \frac{1}{t} \left(\frac{t+s}{s}\right)^{\nu} \left[\log\left(a\left(\frac{t+s}{t}\right)^{\nu+1}\right) \right]^{\nu+1} \\ \times \exp\left[-\frac{t+s}{s}\log\left(a\left(\frac{t+s}{t}\right)^{\nu+1}\right)\right] \mathbb{1}_{\operatorname{Supp}(\tau_{\psi})}(t) \, \mathrm{d}t.$$

PROOF. We will only sketch the proof as it follows the same ideas as the one of the Theorem 2.1. Let us consider the measure $F(dy) = p_0(s, y) dy$ for s > 0 fixed. Then, when evaluating the corresponding u(t, x), we have

$$\begin{aligned} u(t,x) &= p_0(t,x) - \frac{1}{a} \int_{\mathbb{R}_+} p_0(s,y) p_y(t,x) \, \mathrm{d}y \\ &= p_0(t,x) - \frac{1}{a} p_0(t+s,x) \\ &= \frac{1}{2^{\nu}} \frac{1}{\Gamma(\nu+1)} x^{2\nu+1} \Big[\frac{1}{t^{\nu+1}} \exp\left(-\frac{x^2}{2t}\right) - \frac{1}{a} \frac{1}{(t+s)^{\nu+1}} \exp\left(-\frac{x^2}{2(t+s)}\right) \Big], \end{aligned}$$
by using the Markov property. We obtain the form of $u(t)$ by the condition

by using the Markov property. We obtain the form of $\psi(t)$ by the condition $u(t, \psi(t)) = 0$ which gives

(2.23)
$$\psi(t) = \sqrt{\frac{2t(t+s)}{s}} \Big[(v+1) \log \Big(1 + \frac{s}{t} \Big) + \log a \Big],$$
$$\begin{cases} \text{for } t \ge 0 \text{ if } a \ge 1, \\ \text{for } t \le \frac{s}{(1/a)^{1/(v+1)} - 1} \text{ if } a < 1. \end{cases}$$

In order to obtain the distribution of τ_{ψ} , one has only to evaluate

(2.24)
$$\frac{\partial u}{\partial x}(t,x) = (\delta-1)\frac{u(t,x)}{x} - \frac{s}{t(t+s)}xp_0(t,x),$$

and $\frac{u(t,x)}{x}$ for x = 0 and $x = \psi(t)$ and replace the values in the general form (2.14). The expression (2.22) follows. \Box

REMARK 2.4. We can notice that the function $\psi_a(t)$ defined by (2.21) satisfies, for large times,

$$\begin{cases} \psi_a(t) \simeq \sqrt{t}, & \text{for } a = 1, \\ \psi_a(t) \simeq t, & \text{for all } a > 1 \end{cases}$$

In particular, we can approach large times by considering this kind of boundary.

A new boundary can be obtained by using the Laplace transform of the square of the δ -dimensional Bessel process starting from *x*. More precisely:

PROPOSITION 2.5. Let, for $\delta = 2\nu + 2 > 0$, $\lambda > 0$ and a > 0 fixed,

(2.25)
$$\psi_a(t) = \frac{\lambda t}{1 + 2\lambda t} + t \sqrt{\left(\frac{\lambda}{1 + 2\lambda t}\right)^2 + \frac{2}{t} \log \frac{a(1 + 2\lambda t)^{\nu+1}}{2^{\nu} t^{\nu+1} \Gamma(\nu+1)}}$$

for all $t \in \text{Supp}(\tau_{\psi})$, where $\text{Supp}(\tau_{\psi})$ is defined by

(2.26)
$$\operatorname{Supp}(\tau_{\psi}) = \begin{cases} \left[0, \frac{1}{(2^{\nu}\Gamma(\nu+1)/a)^{1/(\nu+1)} - 2\lambda}\right], \\ if \lambda < \frac{1}{2} \left(\frac{2^{\nu}\Gamma(\nu+1)}{a}\right)^{1/(\nu+1)}, \\ [0, +\infty), \\ if \lambda \ge \frac{1}{2} \left(\frac{2^{\nu}\Gamma(\nu+1)}{a}\right)^{1/(\nu+1)}. \end{cases}$$

Then the probability density function of the hitting time is given by

(2.27)
$$\mathbb{P}_{0}(\tau_{\psi} \in dt) = \sqrt{\left(\frac{\lambda}{1+2\lambda t}\right)^{2} + \frac{2}{t}\log\frac{a(1+2\lambda t)^{\nu+1}}{2^{\nu}t^{\nu+1}\Gamma(\nu+1)}}p_{0}(t,\psi(t))\mathbb{1}_{\mathrm{Supp}(\tau_{\psi})}(t)\,dt.$$

PROOF. We present only the main ideas as the result follows as above from the general method in Theorem 2.1 applied to the measure $F(dy) = y^{2\nu+1}e^{-\lambda y^2} \mathbb{1}_{\{y \ge 0\}} dy$. For this measure u(t, x) takes the form

$$u(t, x) = p_0(t, x) - \frac{1}{a} \int_{\mathbb{R}_+} p_y(t, x) F(dy)$$

= $p_0(t, x) - \frac{1}{a} \int_{\mathbb{R}_+} p_y(t, x) y^{2\nu+1} e^{-\lambda y^2} dy$
= $p_0(t, x) - \frac{1}{a} \int_{\mathbb{R}_+} x^{2\nu+1} p_x(t, y) e^{-\lambda y^2} dy$
= $p_0(t, x) - \frac{x^{2\nu+1}}{a} \mathbb{E}(e^{-\lambda Z_t^{\delta, x}}).$

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By using the expression of the Laplace transform for $Z_t^{\delta, x}$ we obtain

(2.28)
$$u(t,x) = p_0(t,x) - \frac{x^{2\nu+1}}{a} \frac{1}{(1+2\lambda t)^{\delta/2}} \exp\left(-\frac{\lambda x}{1+2\lambda t}\right)$$

We consider first the equality $u(t, \psi(t)) = 0$ in (2.28), and this gives the form of $\psi(t)$ in (2.25). Afterwards, we can evaluate once again in this particular situation

$$\frac{\partial u}{\partial x}(t,x) = (\delta - 1)\frac{u(t,x)}{x} - \left(\frac{x}{t} - \frac{\lambda t}{1 + 2\lambda t}\right)p_0(t,x).$$

For this particular case, there is only one nonvanishing term in expression (2.10) of $\mathbb{P}_0(\tau_{\psi} \in dt)$, that is, the term $-(\frac{x}{t} - \frac{\lambda t}{1+2\lambda t})p_0(t,x)$ of $\frac{\partial u}{\partial x}(t,x)$ for $x = \psi(t)$, and this is exactly given by the right-hand side of formula (2.27). \Box

COROLLARY 2.6. The previous results give, for $\delta = 2$:

(1) for a > 0, $0 \le t \le a$ and $\psi(t) = \sqrt{2t \log \frac{a}{t}}$, the density of the hitting time τ_{ψ} is

$$\mathbb{P}_0(\tau_{\psi} \in \mathrm{d}t) = \frac{1}{2a} \log\left(\frac{a}{t}\right) \mathbb{1}_{\{0 \le t \le a\}}(t) \,\mathrm{d}t;$$

(2) for s > 0, a > 0, $0 \le t \le \frac{sa}{1-a}$ and $\psi(t) = \sqrt{\frac{2t(t+s)}{s}\log(a\frac{t+s}{t})}$, the probability density function of τ_{ψ} is given by

$$\mathbb{P}_{0}(\tau_{\psi} \in \mathrm{d}t) = \frac{t+s}{t} \log\left(a\frac{t+s}{t}\right) \exp\left[-\frac{t+s}{t} \log\left(a\frac{t+s}{t}\right)\right] \mathbb{1}_{\left\{0 \le t \le sa/(1-a)\right\}}(t) \,\mathrm{d}t;$$

(3) for a > 0 and $\psi(t) = \frac{\lambda t}{1+2\lambda t} + t\sqrt{(\frac{\lambda}{1+2\lambda t})^2 + \frac{2}{t}\log\frac{a(1+2\lambda t)}{t}}$, for $t \in \text{Supp}(\tau_{\psi})$, where

(2.29)
$$\operatorname{Supp}(\tau_{\psi}) = \begin{cases} [0, +\infty), & \text{if } \lambda \ge \frac{1}{2a}, \\ \left[0, \frac{a}{1 - 2\lambda a}\right], & \text{if } \lambda < \frac{1}{2a}, \end{cases}$$

the probability density function of τ_{ψ} is

$$\mathbb{P}_{0}(\tau_{\psi} \in \mathrm{d}t) = \sqrt{\left(\frac{\lambda}{1+2\lambda t}\right)^{2} + \frac{2}{t}\log\frac{a(1+2\lambda t)}{t}} p_{0}(t,\psi(t))\mathbb{1}_{\mathrm{Supp}(\tau_{\psi})}(t)\,\mathrm{d}t.$$

2.2. Approximation of the first hitting time for Bessel processes starting from the origin. In this section we will construct a stepwise procedure, the so-called random walk on moving spheres (WoMS) algorithm, which allows us to approach the first time the standard Bessel process hits a given level l > 0. Of course, this stopping time $\tau_l = \inf\{t > 0; Z_t^{\delta,x} = l\}$ can be characterized by its well-known Laplace transform computed by solving an eigenvalue problem. Indeed if $(Z_t^{\delta,x}, t \ge 0)$ is the Bessel process starting from x, of index $\nu = \frac{\delta}{2} - 1$, then for $\nu > 0$ and $x \le l$, we get

$$\mathbb{E}_{x}\left[e^{-\lambda\tau_{l}}\right] = \frac{x^{-\nu}I_{\nu}(x\sqrt{2\lambda})}{l^{-\nu}I_{\nu}(l\sqrt{2\lambda})}, x > 0 \quad \text{and} \quad \mathbb{E}_{0}\left[e^{-\lambda\tau_{l}}\right] = \frac{(l\sqrt{2\lambda})^{\nu}}{2^{\nu}\Gamma(\nu+1)}\frac{1}{I_{\nu}(l\sqrt{2\lambda})}.$$

Here I_{ν} denotes the modified Bessel function. This Laplace transform can be used to describe the following tail distribution: Ciesielski and Taylor [3] proved that, for $\delta = 2\nu + 2 \in \mathbb{N}^*$,

$$\mathbb{P}_{0}(\tau_{l} > t) = \frac{1}{2^{\nu-1}\Gamma(\nu+1)} \sum_{k=1}^{\infty} \frac{j_{\nu,k}^{\nu-1}}{\mathcal{J}_{\nu+1}(j_{\nu,k})} e^{(-j_{\nu,k}^{2}/(2l^{2}))t}$$

where \mathcal{J}_{ν} is the Bessel function of the first kind, and $(j_{\nu,k})_{\nu,k}$ is the associated sequence of its positive zeros.

We are looking for a numerical approach for the hitting time and these formulae are not easy to handle and approach, in particular we cannot compute the inverse cumulative function! The aim of this section is to construct an easy and efficient algorithm without need of inverting the Laplace transform and without using discretization schemes since the hitting times are unbounded. In the next step we will extend this procedure to the hitting time of time-dependent boundaries like straight lines, useful in the description of the hitting time of a given level for the CIR process (the Laplace transform is then unknown).

2.2.1. Hitting time of a given level for the Bessel process with positive integer dimension δ . Let us consider δ independent one-dimensional Brownian motions $(B_t^{(k)}, t \ge 0), 1 \le k \le \delta$. Then the Bessel process of index δ starting from 0, satisfies the following property:

$$(Z_t^{\delta,0}, t \ge 0)$$
 has the same distribution as $(\sqrt{(B_t^{(1)})^2 + \dots + (B_t^{(\delta)})^2}, t \ge 0).$

Let

(2.30)
$$\tau_l = \inf\{t \ge 0; Z_t^{\delta, 0} \ge l\}.$$

In particular, we can express τ_l by using the first time when the δ -dimensional Brownian motion $\mathbf{B} = (B^{(1)}, \dots, B^{(\delta)})$ exits from the Euclidean ball *D* centered in the origin with radius *l*. Approximating the exit time and the exit position for the 2-dimensional Brownian motion of a convex domain was already studied by

Milstein [12]. He used the so-called random walk on spheres algorithm which allows one to approach the exit location and the exit time through an efficient algorithm. The exit position is really simple to obtain (as it is uniformly distributed on the circle) while the exit time is much more difficult to approach. That is why we will construct an adaptation of this initial procedure in order to obtain nice and efficient results concerning the Bessel process exit time.

Let us introduce now our *walk on moving spheres (WoMS)* algorithm. We first define a continuous function $\rho : \mathbb{R}^2 \to \mathbb{R}_+$ which represents the distance to the boundary of D,

(2.31)
$$\rho(x) = \inf\{\|x - y\|; y \in D^c\} = l - \|x\|.$$

For any small enough parameter $\varepsilon > 0$, we will denote by D^{ε} the sphere centered at the origin with radius $l - \varepsilon$,

(2.32)
$$D^{\varepsilon} = \{x \in D; \|x\| \le l - \varepsilon\} = \{x \in D; \rho(x) \ge \varepsilon\}.$$

Algorithm (A1) for $\delta = 2$. Let us fix a parameter $0 < \gamma < 1$. Initialization: Set $X(0) = (X_1(0), X_2(0)) = 0$, $\theta_0 = 0$, $\Theta_0 = 0$, $A_0 = \gamma^2 l^2 e/2$. First step: Let (U_1, V_1, W_1) be a vector of three independent random variables uniformly distributed on [0, 1]. We set

$$\begin{cases} \theta_1 = A_0 U_1 V_1, & \Theta_1 = \Theta_0 + \theta_1, \\ X(1)^{\mathsf{T}} = (X_1(1), X_2(1))^{\mathsf{T}} = X(0)^{\mathsf{T}} + \psi_{A_0}(\theta_1) \begin{pmatrix} \cos(2\pi W_1) \\ \sin(2\pi W_1) \end{pmatrix}, \end{cases}$$

where

(2.33)
$$\psi_a(t) = \sqrt{2t \log \frac{a}{t}}, \qquad t \le a, a > 0.$$

At the end of this step we set $A_1 = \gamma^2 \rho(X(1))^2 e/2$.

The *n*th step: While $X(n-1) \in D^{\varepsilon}$, simulate (U_n, V_n, W_n) a vector of three independent random variables uniformly distributed on [0, 1] and define

(2.34)
$$\begin{cases} \theta_n = A_{n-1}U_n V_n, & \Theta_n = \Theta_{n-1} + \theta_n, \\ X(n)^{\mathsf{T}} = (X_1(n), X_2(n))^{\mathsf{T}} = X(n-1)^{\mathsf{T}} + \psi_{A_{n-1}}(\theta_n) \begin{pmatrix} \cos(2\pi W_n) \\ \sin(2\pi W_n) \end{pmatrix} \end{cases}$$

At the end of this step we set $A_n = \gamma^2 \rho(X(n))^2 e/2$.

When $X(n-1) \notin D^{\varepsilon}$ the algorithm is stopped: we set $\theta_n = 0$, $\Theta_n = \Theta_{n-1}$ and X(n) = X(n-1).

Outcome: The hitting time Θ_n and the exit position X(n).

REMARK 2.7. The WoMS algorithm describes a *D*-valued Markov chain $(X(n), n \ge 0)$. Each step corresponds to an exit problem for the 2-dimensional

Brownian motion. If X(n) = x, then we focus our attention to the exit problem of the ball centered in x and of radius ($\psi_a(t), t \ge 0$): the exit location corresponds to X(n+1) and the exit time to θ_{n+1} . Of course the choice of the parameter *a* plays a crucial role since the moving sphere has to belong to the domain *D* as time elapses. When the Markov chain *X* is close to the boundary ∂D , we stop the algorithm and obtain therefore a good approximation of the exit problem of *D*.

Comparison with the classical (*WoS*) algorithm: at each step, the *n*th step of the classical walk on spheres (*WoS*) is based on the exit location and exit time, which are mutually independent, for the Brownian paths exiting from a ball centered in X(n-1) and with radius $\gamma \rho(X(n-1))$. The exit location is uniformly distributed on the sphere while the exit time is characterized by its Laplace transform. Therefore, if one knows X(n-1), then the diameter of the sphere is deterministic. For the *WoMS* the center of the ball should also be X(n-1), but the radius is random, smaller than $\gamma \rho(X(n-1))$. The exit location will also be uniformly distributed on the sphere, but the exit time will be much easier to simulate: in particular, you do not need to evaluate the Bessel functions.

The stochastic process $(X(n), n \ge 0)$ is a homogeneous Markov chain stopped at the first time it exits from D^{ε} . In the following, we shall denote N^{ε} this stopping time which represents in fact the number of steps in the algorithm:

$$N^{\varepsilon} = \inf\{n \ge 0; X(n) \notin D^{\varepsilon}\}.$$

We just notice that $X(N^{\varepsilon}) \notin D^{\varepsilon}$ by its definition.

Algorithm (A1) is presented in the 2-dimensional case. Of course we can construct a generalization of this procedure for the δ -dimensional Bessel process when $\delta \in \mathbb{N}^*$. For notational simplicity, we use a slightly different method: instead of dealing with a Markov chain ($X(n), n \in \mathbb{N}$) living in \mathbb{R}^{δ} we shall consider its squared norm, which is also (surprisingly) a Markov chain. At each step, we shall construct a couple of random variables ($\xi_n, \chi(n)$) associated to an exit problem, the first coordinate corresponds to an exit time and the second one to the norm of the exit location.

We introduce some notation: S^{δ} represents the unit ball in \mathbb{R}^{δ} and $\pi_1 : \mathbb{R}^{\delta} \to \mathbb{R}$ the projection on the first coordinate.

Algorithm (A2). Let us fix a parameter $0 < \gamma < 1$. Initialization: Set $\chi(0) = 0$, $\xi_0 = 0$, $\Xi_0 = 0$, $A_0 = (\gamma^2 l^2 e / (\nu + 1))^{\nu+1} \frac{\Gamma(\nu+1)}{2}$. The *n*th step: While $\sqrt{\chi(n-1)} < l - \varepsilon$, we choose U_n a uniform distributed random vector on $[0, 1]^{\lfloor \nu \rfloor + 2}$, G_n a standard Gaussian random variable and V_n an uniformly distributed random vector on S^{δ} . Consider U_n , G_n and V_n independent. We set

(2.35)
$$\begin{cases} \xi_n = \left(\frac{A_{n-1}}{\Gamma(\nu+1)2^{\nu}}U_n(1)\cdots U_n(\lfloor\nu\rfloor+2)\right)^{1/(\nu+1)} \exp\left\{-\frac{\nu-\lfloor\nu\rfloor}{\nu+1}G_n^2\right\},\\ \Xi_n = \Xi_{n-1} + \xi_n,\\ \chi(n) = \chi(n-1) + 2\pi_1(V_n)\sqrt{\chi(n-1)}\psi_{A_{n-1}}(\xi_n) + \psi_{A_{n-1}}^2(\xi_n),\end{cases}$$

where

(2.36)
$$\psi_{a}(t) = \sqrt{2t \log \frac{a}{\Gamma(\nu+1)t^{\nu+1}2^{\nu}}},$$
$$t \le t_{\max}(a) := \left[\frac{a}{\Gamma(\nu+1)2^{\nu}}\right]^{1/(\nu+1)}, \qquad a > 0.$$

At the end of this step we set

$$A_n = (\gamma^2 (l - \sqrt{\chi(n)})^2 e / (\nu + 1))^{\nu + 1} \frac{\Gamma(\nu + 1)}{2}$$

When $\sqrt{\chi(n)} \ge l - \varepsilon$ the algorithm is stopped: we then set $\xi_n = 0$, $\Xi_n = \Xi_{n-1}$ and $\chi(n) = \chi(n-1)$. **Outcome:** The hitting time Ξ_n and the value of the Markov chain $\chi(n)$.

It is obvious that for the particular dimension $\delta = 2$, that is, $\nu = 0$, the stopping times obtained by Algorithms (A1) and (A2) have the same distribution. Moreover, for each n, $\chi(n)$ has the same distribution as $||X(n)||^2$. In other words, if the number of steps of (A1) and (A2) are identical in law, the number of steps will be denoted in both cases N^{ε} .

THEOREM 2.8. Set $\delta \in \mathbb{N}^*$. The number of steps N^{ε} of the Algorithm WoMS (A2) is almost surely finite. Moreover, there exist constants $C_{\delta} > 0$ and $\varepsilon_0(\delta) > 0$, such that

$$\mathbb{E}[N^{\varepsilon}] \le C_{\delta} |\log \varepsilon| \qquad for \ all \ \varepsilon \le \varepsilon_0(\delta).$$

THEOREM 2.9. Set $\delta \in \mathbb{N}^*$. As ε goes to zero, $\Xi_{N^{\varepsilon}}$ converges in distribution toward τ_l , the hitting time of the δ -dimensional Bessel process (with cumulative distribution function F), which is almost surely finite. Moreover, for any $\alpha > 0$ small enough,

(2.37)
$$\left(1 - \frac{\varepsilon}{\sqrt{2\alpha\pi}}\right)F^{\varepsilon}(t-\alpha) \le F(t) \le F^{\varepsilon}(t) \quad \text{for all } t > 0,$$

where $F^{\varepsilon}(t) := \mathbb{P}(\Xi_{N^{\varepsilon}} \leq t)$.

These results and the key ideas of the proofs are adapted from the classical random walk on spheres (*WoS*); see [12].

PROOF OF THEOREM 2.8. Step 1. Let us estimate the number of steps. Since $(\chi(n), n \ge 0)$ is a homogeneous Markov chain, we introduce the operator $P_x f$ defined, for any nonnegative function $f : \mathbb{R}_+ \to \mathbb{R}_+$, by

$$P_x f := \int_{\mathbb{R}_+} f(y) \mathbb{P}(x, \mathrm{d}y),$$

where $\mathbb{P}(x, dy)$ is the transition probability of the Markov chain. By definition, $\chi(n + 1)$ depends only on $\chi(n)$, V_n and ξ_n . Let us note that, by construction, V_n and ξ_n are independent. Moreover using the result developed in the Appendix, the density of $\xi_n (\frac{2^{\nu} \Gamma(\nu+1)}{A_{n-1}})^{1/(\nu+1)}$ is given by

(2.38)
$$\mu(r) = \frac{(\nu+1)^{\nu+2}}{\Gamma(\nu+2)} r^{\nu} (-\log r)^{\nu+1} \mathbb{1}_{[0,1]}(r).$$

If we denote σ^d , the uniform surface measure on the unit sphere in \mathbb{R}^d , we get

(2.39)
$$P_x f = \int_0^1 \int_{\mathcal{S}^{\delta}} f(x + 2\pi_1(u)\sqrt{x}K(x,r) + K^2(x,r))\mu(r) \,\mathrm{d}r \,\sigma^{\delta}(\mathrm{d}u),$$

with K(x, r) defined by

(2.40)
$$K(x,r) = \psi_A \left(\left[\frac{A}{2^{\nu} \Gamma(\nu+1)} \right]^{1/(\nu+1)} r \right),$$

and A depending on x in the following way:

$$A = \left(\frac{\gamma^2 (l - \sqrt{x})^2 e}{\nu + 1}\right)^{\nu + 1} \frac{\Gamma(\nu + 1)}{2}.$$

We can observe the following scaling property: $\psi_A(A^{1/(\nu+1)}t) = A^{1/(2\nu+2)}\psi_1(t)$. Therefore the definition of ψ_1 leads to

(2.41)
$$K(x,r) = \gamma(l-\sqrt{x})\sqrt{\frac{er}{\nu+1}\log\frac{1}{r^{\nu+1}}} = \gamma(l-\sqrt{x})\sqrt{er(-\log r)}.$$

Step 2. Using classical potential theory for discrete time Markov chains (see, e.g., Theorem 4.2.3 in [14]), we know that

$$\phi(x) = \mathbb{E}_x \left(\sum_{n=0}^{N^c - 1} g(\chi(n)) \right)$$

satisfies, for any nonnegative function g,

(2.42)
$$\begin{cases} \phi(x) = P_x \phi + g(x), & 0 \le x < (l - \varepsilon)^2, \\ \phi((l - \varepsilon)^2) = 0. \end{cases}$$

In particular, for g = 1, we obtain that $\phi(x) = \mathbb{E}_x[N^{\varepsilon}]$. In order to get an upperbound for the averaged number of steps, it suffices to apply a comparison result. Let us first define the constant C_{δ} ,

(2.43)
$$C_{\delta} = \left(\frac{\nu+1}{\nu+2}\right)^{\nu+2} \frac{e}{\Gamma(\nu+2)} \frac{1}{2\delta} \sigma^{\delta}(S^{\delta}).$$

We choose the function

(2.44)
$$U^{\varepsilon}(x) = \left\{ \log((l - \sqrt{x})/\varepsilon) - \log(1 - \gamma) \right\} / (C_{\delta} \gamma^2), \qquad 0 \le x < l^2,$$

which satisfies $U^{\varepsilon}(x) \ge P_x U^{\varepsilon} + 1$, for all $0 < x < (l - \varepsilon)^2$ (see Lemma 2.10 for the definition of the constant and for the inequality) and $U^{\varepsilon}(x) \ge 0$ for all $0 < x < (l - \varepsilon)^2$. A classical comparison result related to the potential theory (see, e.g., Theorem 4.2.3 in [14]) implies that $\mathbb{E}_x[N^{\varepsilon}] \le U^{\varepsilon}(x)$ for all $x \in [0, (l - \varepsilon)^2]$ and consequently leads to the announced statement. \Box

LEMMA 2.10. Let us define, for small $\varepsilon > 0$, $U^{\varepsilon}(x) = \{\log((l - \sqrt{x})/\varepsilon) - \log(1 - \gamma)\}/(C_{\delta}\gamma^2)$ for $x \in [0, l^2[$ and where the constant C_{δ} is given by (2.43) and γ is related to the definition of the WoMS. Then, for any $x \in]0, (l - \varepsilon)^2[$, the following inequality yields

$$P_{x}U^{\varepsilon} - U^{\varepsilon}(x) \leq -1.$$

We recall that $P_x U^{\varepsilon}$ is defined by (2.39) and (2.41).

PROOF. We will split the proof into several steps.

Step 1. First of all, we observe that $U^{\varepsilon} \ge -\log(1-\gamma)/(C_{\delta}\gamma^2)$ in the domain $[0, (l-\varepsilon)^2]$. Let us consider now $\chi(0) = x \in [0, (l-\varepsilon)^2]$ and y in the support of the law of $\chi(1)$ and let us prove that $U^{\varepsilon}(y) \ge 0$. By the definition of $\chi(1)$ we obtain

$$\chi(1) \le \sup_{y \in [-1,1], t \in [0, t_{\max}(A)]} (x + 2y\sqrt{x}\psi_A(t) + \psi_A^2(t)),$$

where $A = (\gamma^2 (l - \sqrt{x})^2 e/(\nu + 1))^{\nu+1} \frac{\Gamma(\nu+1)}{2}$ and both ψ_A and t_{max} are defined by (2.36). The right-hand side of the preceding inequality is increasing with respect to *y* so that

$$\chi(1) \leq \left(\sqrt{x} + \sup_{t \in [0, t_{\max}(A)]} \psi_A(t)\right)^2.$$

Furthermore, for a > 0 the maximum of the function ψ_a is reached for $t_{\max}(a) = \frac{1}{e} (\frac{a}{\Gamma(\nu+1)2^{\nu}})^{1/(\nu+1)}$ and is equal to

(2.45)
$$\sup_{t \in [0, t_{\max}(a)]} \psi_a(t) = \left\{ \frac{2(\nu+1)}{e} \left(\frac{a}{\Gamma(\nu+1)2^{\nu}} \right)^{1/(\nu+1)} \right\}^{1/2}.$$

Finally using the definition of A and the inequality $x \le (l - \varepsilon)^2$, we find the following lower bound:

$$l - \sqrt{\chi(1)} \ge (l - \sqrt{x})(1 - \gamma) \ge \varepsilon(1 - \gamma).$$

We can therefore conclude that, for any y in the support of the law of $\chi(1)$ (even for $y \ge (l-\varepsilon)^2$), $U^{\varepsilon}(y) \ge 0$ which ensures that U^{ε} is well defined and nonnegative in the domain of the operator P_x .

Step 2. Furthermore the Taylor expansion yields

$$U^{\varepsilon}(y) \le U^{\varepsilon}(x) + \frac{\sqrt{x} - \sqrt{y}}{C_{\delta}\gamma^{2}(l - \sqrt{x})} - \frac{(\sqrt{x} - \sqrt{y})^{2}}{2C_{\delta}\gamma^{2}(l - \sqrt{x})^{2}} + \frac{(\sqrt{x} - \sqrt{y})^{3}}{3C_{\delta}\gamma^{2}(l - \sqrt{x})^{3}},$$
(2.46)

$$x, y \in [0, l^{2}[.$$

If $\chi(0) = x$ and y is in the support of the random variable $\chi(1)$, then

$$\sqrt{y} - \sqrt{x} = \sqrt{x + 2\pi_1(u)\sqrt{x}K(x,r) + K^2(x,r)} - \sqrt{x}$$
$$\geq \pi_1(u)K(x,r).$$

By expansion (2.46) and the definition of the operator P_x given by (2.39), the following upper-bound for the operator P_x holds:

$$P_{x}U^{\varepsilon} = \int_{0}^{1} \int_{\mathcal{S}^{\delta}} U^{\varepsilon} (x + 2\pi_{1}(u)\sqrt{x}K(x,r) + K^{2}(x,r))\mu(r) \, \mathrm{d}r \, \sigma^{\delta}(\mathrm{d}u),$$

$$\leq U^{\varepsilon}(x) - \int_{0}^{1} \int_{\mathcal{S}^{\delta}} \frac{\pi_{1}(u)K(x,r)}{C_{\delta}\gamma^{2}(l-\sqrt{x})}\mu(r) \, \mathrm{d}r \, \sigma^{\delta}(\mathrm{d}u)$$

$$- \int_{0}^{1} \int_{\mathcal{S}^{\delta}_{+}} \frac{\pi_{1}^{2}(u)K^{2}(x,r)}{2C_{\delta}\gamma^{2}(l-\sqrt{x})^{2}}\mu(r) \, \mathrm{d}r \, \sigma^{\delta}(\mathrm{d}u)$$

$$- \int_{0}^{1} \int_{\mathcal{S}^{\delta}} \frac{\pi_{1}^{3}(u)K^{3}(x,r)}{3C_{\delta}\gamma^{2}(l-\sqrt{x})^{3}}\mu(r) \, \mathrm{d}r \, \sigma^{\delta}(\mathrm{d}u),$$

where

(2.47)
$$S_{+}^{\delta} := \{ u \in S^{\delta} : \pi_{1}(u) > 0 \}.$$

Due to symmetry properties, the first and the third integral terms vanish. Then (2.41) leads to

$$P_{x}U^{\varepsilon} \leq U^{\varepsilon}(x) - \frac{I}{C_{\delta}} \int_{S^{\delta}_{+}} \pi_{1}^{2}(u) \sigma^{\delta}(\mathrm{d}u)$$

with

$$I = \frac{(\nu+1)^{\nu+2}e}{2\Gamma(\nu+2)} \int_0^1 r^{\nu+1} (-\log r)^{\nu+2} \,\mathrm{d}r.$$

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The description of the probability density function in the Appendix leads to the following explicit value:

$$I = \left(\frac{\nu+1}{\nu+2}\right)^{\nu+2} \frac{e}{\Gamma(\nu+2)}$$

In order to complete the proof, it suffices to choose the particular constant given by (2.43) after noticing that

$$\int_{S^{\delta}_{+}} \pi_{1}^{2}(u) \sigma^{\delta}(\mathrm{d}u) = \frac{1}{2\delta} \sigma^{\delta}(S^{\delta}).$$

PROOF OF THEOREM 2.9. The proof is split in two parts. First, the steps of the algorithm and the hitting time of the Bessel process of index ν shall be related to stopping times of a δ -dimensional Brownian motion ($\nu = \frac{\delta}{2} - 1$). Second, we point out that the corresponding stopping times are close together by evaluating deviations of the Brownian paths.

Step 1. Let $\mathbf{B} = (B^{(1)}, \hat{B}^{(2)}, \dots, B^{(\delta)})$ be a δ -dimensional Brownian motion. Then the norm of **B** has the same distribution as a Bessel process of index ν ; see, for instance, [15]. Hence the first hitting time τ_l is identical in law to the stopping time

$$\mathbb{T}_l = \inf\{t \ge 0; \mathbf{B}_t \notin D\},\$$

where *D* is the Euclidean ball centered at the origin and of radius *l*. We introduce then a procedure in order to come close to \mathbb{T}_l . For the first step we shall focus our attention to the first exit time of a moving sphere centered at the origin and of radius $\psi_a(t)$ defined by (2.36), we denote $\hat{\xi}_1$ this stopping time. Of course this moving sphere should always stay in *D*, so we choose *a* such that the maximum of ψ_a stays smaller than *l*. By (2.45), we get

$$\sup_{t \le a} \psi_a(t) < l \quad \Longleftrightarrow \quad a < \frac{\Gamma(\nu+1)}{2} \left(\frac{el^2}{\nu+1}\right)^{\nu+1}.$$

For $a = A_0 = \frac{\Gamma(\nu+1)}{2} (\frac{e\gamma^2 l^2}{\nu+1})^{\nu+1}$ with a parameter $\gamma < 1$, the condition is satisfied, sup_{t \le a} \psi_a(t) = \gamma^{2\nu+2} l < l. Let us describe the law of $(\hat{\xi}_1, \mathbf{B}_{\hat{\xi}_1})$. The norm of the Brownian motion is identical in law with the Bessel process; therefore Proposition 2.2 implies that the density function of $\hat{\xi}_1$ is given by (2.16) with *a* replaced by A_0 . Using the law described in Proposition A.1, we can prove that $\hat{\xi}_1$ has the same distribution as

$$\left(\frac{A_0}{\Gamma(\nu+1)2^{\nu}}\right)^{1/(\nu+1)}e^{-Z},$$

where Z is Gamma distributed with parameters $\alpha = \nu + 2$ and $\beta = \frac{1}{\nu+1}$. By construction we deduce that $\hat{\xi}_1 \stackrel{(d)}{=} \xi_1$ where ξ_1 is defined in the Algorithm *WoMS* (A2). Knowing the stopping time, we can easily describe the exit location since the Brownian motion is rotationnally invariant: $\mathbf{B}_{\hat{\xi}_1}$ is then uniformly distributed on

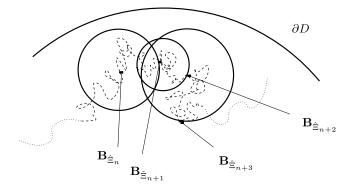


FIG. 1. Walk on moving spheres.

the sphere of radius $\psi_{A_0}(\hat{\xi}_1)$. Hence

$$(\hat{\xi}_1, \|\mathbf{B}_{\hat{\xi}_1}\|) \stackrel{(\mathrm{d})}{=} (\xi_1, \chi(1)) \quad \text{and} \quad \hat{\xi}_1 < \mathbb{T}_l.$$

By this procedure we can construct a sequence of stopping times $(\hat{\xi}_n, n \ge 1)$ and define $\hat{\Xi}_n = \hat{\xi}_1 + \cdots + \hat{\xi}_n$; $\hat{\Xi}_n$ is the first time after $\hat{\Xi}_{n-1}$ such that the Brownian motion exits from a sphere centered in $\mathbf{B}_{\hat{\Xi}_{n-1}}$ of radius ψ_{a_n} initialized at time $\hat{\Xi}_{n-1}$. See Figure 1. The moving sphere should stay in the domain *D*, so we choose

$$a_n = \left(\gamma^2 (l - \sqrt{\mathbf{B}_{\hat{\Xi}_{n-1}}})^2 e / (\nu + 1)\right)^{\nu + 1} \frac{\Gamma(\nu + 1)}{2}.$$

Using the same arguments as before and by the Markov property for the Brownian motion, we obtain the identities in law

$$(a_n, n \ge 1) \stackrel{\text{(d)}}{=} (A_n, n \ge 1), \qquad (\hat{\Xi}_n, \|\mathbf{B}_{\hat{\Xi}_n}\|)_{n \ge 1} \stackrel{\text{(d)}}{=} (\Xi_n, \chi(n))_{n \ge 1}$$

with $\hat{\Xi}_n < \mathbb{T}_l$ and Ξ_n , A_n , $\chi(n)$ defined in the Algorithm *WoMS* (A2). Consequently defining $\hat{N}^{\varepsilon} = \inf\{n \ge 0; \mathbf{B}_{\hat{\Xi}_n} \notin D^{\varepsilon}\}$, the following identity yields

(2.48)
$$(\hat{\Xi}_{\hat{N}^{\varepsilon}}, \|\mathbf{B}_{\hat{\Xi}_{\hat{N}^{\varepsilon}}}\|) \stackrel{(\mathrm{d})}{=} (\Xi_{N^{\varepsilon}}, \chi(N^{\varepsilon})) \text{ and } \hat{\Xi}_{\hat{N}^{\varepsilon}} < \mathbb{T}_{l}.$$

Step 2. Let us now estimate the difference between $\hat{\Xi}_{\hat{N}^{\varepsilon}}$ and \mathbb{T}_{l} . By (2.48) we first deduce

(2.49)
$$F(t) := \mathbb{P}(\tau_l \le t) = \mathbb{P}(\mathbb{T}_l \le t) \le F^{\varepsilon}(t) := \mathbb{P}(\Xi_{N^{\varepsilon}} \le t), \qquad t > 0.$$

Furthermore, for any small $\alpha > 0$,

(2.50)

$$1 - F(t) = \mathbb{P}(\mathbb{T}_{l} > t, \Xi_{\hat{N}^{\varepsilon}} \le t - \alpha) + \mathbb{P}(\mathbb{T}_{l} > t, \Xi_{\hat{N}^{\varepsilon}} > t - \alpha)$$

$$\leq \mathbb{P}(\mathbb{T}_{l} > t, \widehat{\Xi}_{\hat{N}^{\varepsilon}} \le t - \alpha) + \mathbb{P}(\widehat{\Xi}_{\hat{N}^{\varepsilon}} > t - \alpha)$$

$$\leq \mathbb{P}(\mathbb{T}_{l} > t, \widehat{\Xi}_{\hat{N}^{\varepsilon}} \le t - \alpha) + 1 - F^{\varepsilon}(t - \alpha).$$

At time $\hat{\Xi}_{\hat{N}^{\varepsilon}}$ the Brownian motion is in the ε -neighborhood of the boundary ∂D , hence $l - \|\mathbf{B}_{\hat{\Xi}_{\hat{\Omega}^{\varepsilon}}}\| \le \varepsilon$. Using the strong Markov property, we obtain

(2.51)
$$\mathbb{P}(\mathbb{T}_l > t, \, \hat{\Xi}_{\hat{N}^{\varepsilon}} \le t - \alpha) \le F^{\varepsilon}(t - \alpha) \sup_{y \in D \setminus D^{\varepsilon}} \mathbb{P}_y(\mathbb{T}_l > \alpha)$$

Since the Brownian motion is rotationally invariant, it suffices to choose $y = (l - \varepsilon, 0, ..., 0)$. Due to the convexity of *D*, the following upper-bound holds:

(2.52)
$$\mathbb{P}_{y}(\mathbb{T}_{l} > \alpha) \leq \mathbb{P}_{0}\left(\sup_{0 \leq t \leq \alpha} \overline{B}_{t}^{(1)} < \varepsilon\right) = \mathbb{P}_{0}\left(2\left|\overline{B}_{\alpha}^{(1)}\right| < \varepsilon\right) \leq \frac{\varepsilon}{\sqrt{2\alpha\pi}}$$

Combining (2.49) for the upper-bound and (2.50), (2.51) and (2.52) for the lower-bound yields the announced estimation (2.37). \Box

2.2.2. The first time the Bessel process of index v hits a decreasing curved boundary. The algorithm developed in the previous paragraph can be adapted to the problem of hitting a deacreasing curved boundary. Let us define

(2.53) $\tau = \inf\{t \ge 0 : Z_t^{\delta,0} = l(t)\} \quad \text{where } l \text{ is decreasing and } l(0) > 0.$

ASSUMPTION 2.11. There exists a constant $\Delta_{\min} > 0$ which bounds the derivative of *l*

$$l'(t) \ge -\Delta_{\min} \qquad \forall t \ge 0.$$

The procedure then also consists in building a *WoMS* which reaches a neighborhood of the boundary. But instead of dealing with a fixed boundary as in Section 2.2.1, that is a ball of radius l, we shall in this section introduce the following moving boundary: the ball centered in the origin and of radius l(t). The arguments developed in order to prove Theorems 2.8 and 2.9 will be adapted to this new context.

Algorithm (A3):

Let us define the following positive constants:

(2.54)
$$L = \max(l(0), \Delta_{\min}, \sqrt{\nu+1}), \qquad \kappa = \frac{2^{\nu}}{5^{\nu+1}L^{2\nu+2}}\Gamma(\nu+1).$$

Initialization: Set $\chi(0) = 0, \, \xi_0 = 0, \, \Xi_0 = 0, \, A_0 = \kappa(l(0) - \sqrt{\chi(0)})^{2(\nu+1)}.$

The *n*th step: While the condition

$$l(\Xi_{n-1}) - \sqrt{\chi(n-1)} > \varepsilon$$

[denoted by $\mathbb{C}(n-1)$] holds, we simulate U_n an uniform distributed random vector on $[0, 1]^{\lfloor \nu \rfloor + 2}$, G_n a standard Gaussian random variable and V_n a uniformly distributed random vector on S^{δ} . U_n , G_n and V_n have to be independent. We then construct $(\xi_n, \Xi_n, \chi(n))$ using (2.35). At the end of this step we

set $A_n = \kappa (l(\Xi_n) - \sqrt{\chi(n)})^{2(\nu+1)}$. The algorithm stops when $\mathbb{C}(n-1)$ is not longer satisfied: we set $\xi_n = 0$ and so $\Xi_n = \Xi_{n-1}$ and $\chi(n) = \chi(n-1)$. **Outcome** The exit position $\chi(n)$ and the exit time.

Let us note that the stochastic process $(\chi(n), n \ge 0)$ is not a Markov chain since the sequence $(A_n)_{n\ge 0}$ depends on both $(\Xi_n, \chi(n))$. That is why we define the following Markov chain:

$$R_n := (\Xi_n, \chi(n)) \in \mathbb{R}^2_+$$

stopped at the first time the condition $\mathbb{C}(n)$ is not satisfied. In the following, we shall denote N^{ε} this stopping time (number of steps of the algorithm):

$$N^{\varepsilon} = \inf\{n \ge 0; l(\Xi_n) - \sqrt{\chi(n)} \le \varepsilon\}.$$

THEOREM 2.12. The number of steps N^{ε} of the Algorithm WoMS (A3) is almost surely finite. Moreover, there exist a constant $C_{\delta} > 0$ and $\varepsilon_0(\delta) > 0$, such that

$$\mathbb{E}[N^{\varepsilon}] \le C_{\delta} |\log \varepsilon| \qquad for \ all \ \varepsilon \le \varepsilon_0(\delta).$$

THEOREM 2.13. As ε goes to zero, $\Xi_{N^{\varepsilon}}$ converges in distribution toward τ defined by (2.53) (with cumulative distribution function F), which is almost surely finite. Moreover, for any $\alpha > 0$ small enough,

(2.55)
$$\left(1 - \frac{\varepsilon}{\sqrt{2\alpha\pi}}\right)F^{\varepsilon}(t-\alpha) \le F(t) \le F^{\varepsilon}(t) \quad \text{for all } t > 0,$$

where $F^{\varepsilon}(t) := \mathbb{P}(\Xi_{N^{\varepsilon}} \leq t)$.

PROOF OF THEOREM 2.12. The proof is based mainly on arguments already presented in Theorem 2.8. So we let the details of the proof to the reader and focus our attention to the main ideas.

(1) The process $(\Xi_n, \chi(n))$ is a homogeneous Markov chain and the associated operator is given by

(2.56)
$$P_{t,x}f := \int_{(s,y)\in\mathbb{R}^2_+} f(s,y)\mathbb{P}((t,x), (ds, dy)),$$

where *f* is a nonnegative function and $\mathbb{P}((t, x), (ds, dy))$ is the transition probability of the chain. The chain starts with $(\Xi_0, \chi(0)) = (0, 0)$ and is stopped the first time when $l(\Xi_n) - \sqrt{\chi(n)} \le \varepsilon$. Classical potential theory ensures that

$$\phi(t, x) = \mathbb{E}_{t, x} \left(\sum_{n=0}^{N^{\varepsilon} - 1} g\left(\Xi_n, \chi(n) \right) \right)$$

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is solution of the following equation:

(2.57)
$$\begin{cases} \phi(t,x) = P_{t,x}\phi + g(t,x), & (t,x) \in D^{\varepsilon}, \\ \phi(t,x) = 0, & \forall (t,x) \in \partial D^{\varepsilon}, \end{cases}$$

with $D^{\varepsilon} = \{(t, x) \in \mathbb{R}^2_+ : l(t) - \sqrt{x} \le \varepsilon\}$. For the particular choice g = 1, we obtain $\phi(t, x) = \mathbb{E}_{t,x}[N^{\varepsilon}]$, and therefore the averaged number of step is given by $\phi(0, 0)$.

(2) In order to point out an upper-bound for the averaged number of steps, we use a comparison result: we are looking for a function U(t, x) such that

(2.58)
$$\begin{cases} U(t,x) \ge P_{t,x}U+1, & \forall (t,x) \in D^{\varepsilon}, \\ U(t,x) \ge 0, & \forall (t,x) \in \partial D^{\varepsilon}. \end{cases}$$

For such a particular function, we can deduce $\phi(t, x) \leq U(t, x)$. Let us define

$$U(t, x) = c \log\left(\frac{l(t) - \sqrt{x}}{\varepsilon}\right) \mathbb{1}_{\{l(t) - \sqrt{x} \ge 0\}}$$

with some constant c > 0 which shall be specified later on. The positivity assumption on the boundary ∂D^{ε} is trivial. Moreover since *l* is a decreasing function, (2.56) implies

(2.59)

$$P_{t,x}U = \int_{(s,y)\in\mathbb{R}^2_+} U(s,y)\mathbb{P}((t,x), (ds, dy))$$

$$\leq \int_{(s,y)\in\mathbb{R}^2_+} U(t,y)\mathbb{P}((t,x), (ds, dy)).$$

By using the Taylor expansion, we get

(2.60)
$$U(t, y) \le U(t, x) - c \frac{\sqrt{y} - \sqrt{x}}{l(t) - \sqrt{x}} - \frac{c}{2} \frac{(\sqrt{y} - \sqrt{x})^2}{(l(t) - \sqrt{x})^2} - \frac{c}{3} \frac{(\sqrt{y} - \sqrt{x})^3}{(l(t) - \sqrt{x})^3},$$
$$(x, y) \in \mathbb{R}^2_+.$$

Using similar arguments and similar bounds as those presented in Lemma 2.10, the odd powers in the Taylor expansion do not play any role in the integral (2.59). Therefore we obtain

$$P_{t,x}U \le U(t,x) - \frac{c}{2} \int_{(s,y)\in\mathbb{R}^2_+} \frac{(\sqrt{y} - \sqrt{x})^2}{(l(t) - \sqrt{x})^2} \mathbb{P}((t,x), (ds, dy))$$

$$\le U(t,x) - \frac{c}{2} \int_0^1 \int_{S^{\delta}_+} \frac{\pi_1^2(u)K^2(x,r)}{(l(t) - \sqrt{x})^2} \mu(r) \, \mathrm{d}r \, \sigma^{\delta}(\mathrm{d}u),$$

where S_{+}^{δ} is given in (2.47), and K is defined by (2.40) with $A = \kappa (l(s) - \sqrt{x})^{2(\nu+1)}$. We have now

$$P_{t,x}U \leq U(t,x) - \frac{c(\nu+1)}{2} \left(\frac{2K}{\Gamma(\nu+1)}\right)^{1/(\nu+1)} \left(\int_{\mathcal{S}^{\delta}_{+}} \pi_{1}^{2}(u)\sigma^{\delta}(\mathrm{d}u)\right)$$
$$\times \left(\int_{0}^{1} r(-\log r)\mu(r)\,\mathrm{d}r\right).$$

An appropriate choice of the constant c leads to (2.58). Finally we get

$$\mathbb{E}[N^{\varepsilon}] \le U(0,0) = c \log(l(0)/\varepsilon). \qquad \Box$$

PROOF OF THEOREM 2.13. The arguments are similar to those developed for Theorem 2.9, and the extension of the convergence result to curved boundaries is straightforward. That is why we shall not repeat the proof, but just focus our attention on the only point which is quite different. We need to prove that the Markov chain $R_n = (\Xi_n, \chi(n))$ stays in the domain $D^0 = \{(t, x) : 0 \le x \le l^2(t)\}$ so that the hitting time τ defined by (2.53) satisfies $\tau > \Xi_{N^{\varepsilon}}$. In other words, if the Markov chain $R_n = (\Xi_n, \chi(n))$ for the *n*th step is equal to (s, x), then R_{n+1} should belong to $\{(t, x) : t \ge s, x \le l^2(t)\}$. In the *WoMS* setting, for $t \ge s$, this means that the ball centered in *x* and of time-dependent radius $\psi_A(t - s)$ always belongs as time elapses to the ball centered in 0 of radius l(t). We recall that

$$A = \kappa \left(l(s) - \sqrt{x} \right)^{2(\nu+1)}.$$

Therefore we shall prove that

(2.61)
$$\forall t \ge s \qquad \psi_A(t-s) + \sqrt{x} \le l(t).$$

In fact, due to Assumption 2.11 and the definition of ψ_A , it suffices to obtain

(2.62)
$$\psi_A(t-s) \le l(s) - \sqrt{x} - \Delta_{\min}(t-s)$$
 $\forall s \le t \le s + W^2$,

where

$$W = \left(\frac{A}{\Gamma(\nu+1)2^{\nu}}\right)^{1/(2\nu+2)} = \left(\frac{\kappa}{\Gamma(\nu+1)2^{\nu}}\right)^{1/(2\nu+2)} (l(s) - \sqrt{x})$$
$$= \frac{1}{L\sqrt{5}} (l(s) - \sqrt{x}).$$

Due to the definition of the constant L, we have

$$0 \le W \le \frac{1}{2\Delta_{\min}} \frac{2(l(s) - \sqrt{x})\Delta_{\min}}{\sqrt{((2\nu + 2)/e) + 4(l(s) - \sqrt{x})\Delta_{\min}}}$$
$$\le \frac{1}{2\Delta_{\min}} \left\{ \sqrt{\frac{2\nu + 2}{e} + 4(l(s) - \sqrt{x})\Delta_{\min}} - \sqrt{\frac{2\nu + 2}{e}} \right\}.$$

The right-hand side of the preceding inequality is the positive root of the polynomial function $P(X) = \Delta_{\min} X^2 + \sqrt{2(\nu+1)/e} X - (l(s) - \sqrt{x})$. We deduce that $P(W) \le 0$. By (2.45) and $P(W) \le 0$, we obtain

$$\sup_{t \ge s} \psi_A(t-s) = \left(\frac{2(\nu+1)}{e}\right)^{1/2} W$$

$$\leq l(s) - \sqrt{x} - \Delta_{\min} W^2$$

$$\leq l(s) - \sqrt{x} - \Delta_{\min}(t-s) \qquad \forall s \le t \le s + W^2.$$

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Finally we have proved (2.62) and so (2.61). \Box

If Assumption 2.11 is not satisfied, then it is difficult to have a general description of an iterated procedure in order to simulate hitting times. However the particular form of the function ψ_a defined by (2.36) permits us to describe a *WoMS* algorithm for the square root boundaries. Let us therefore consider the following functions:

(2.63)
$$\psi_a(t) = \sqrt{2t \log \frac{a}{\Gamma(\nu+1)t^{\nu+1}2^{\nu}}} \text{ and } f(t) = \sqrt{r-ut},$$

well defined for $t \le t_0 := \min(\alpha^{1/(\nu+1)}, \frac{r}{\nu})$ where $\alpha = a(\Gamma(\nu+1)2^{\nu})^{-1}$.

The algorithm is essentially based on the following result (the constants r and u associated with the hitting problem of a square root boundary for the Bessel process shall be specified in the proof of Proposition 2.15).

LEMMA 2.14. Let us define

(2.64)
$$F_{\nu}(r,u) = \frac{1}{2} \left(\frac{er}{\nu+1}\right)^{\nu+1} \Gamma(\nu+1)e^{-u/2}, \qquad r > 0, u > 0.$$

If $a = F_{\nu}(r, u)$, then

(2.65)
$$\psi_a(t) \le f(t) \quad \text{for all } 0 \le t \le \alpha^{1/(\nu+1)}.$$

PROOF. We are looking for a particular value *a* depending on both *r* and *u* such that the following bound holds: $\psi_a(t) \le f(t)$, for all $0 \le t \le t_0$. Since $t \le t_0$, it suffices to prove that

$$2t\log\frac{\alpha}{t^{\nu+1}} \le r - ut \quad \Longleftrightarrow \quad g(t) := t\left(2\log\frac{\alpha}{t^{\nu+1}} + u\right) \le r.$$

Let us compute the maximum of the function g on the interval $[0, t_0]$, with t_0 fixed,

$$g'(t) = 2\log\frac{\alpha}{t^{\nu+1}} + u - 2(\nu+1).$$

We have

$$g'(t) = 0 \quad \Longleftrightarrow \quad \log \frac{\alpha}{t^{\nu+1}} = \nu + 1 - \frac{u}{2} \quad \Longleftrightarrow \quad t^{\nu+1} = \alpha \exp\left\{\frac{u}{2} - \nu - 1\right\}.$$

In other words the maximum of the function g is reached for

$$t_{\max} = \alpha^{1/(\nu+1)} \exp\left\{\frac{u}{2(\nu+1)} - 1\right\}$$

and is equal to

$$g(t_{\max}) = g_{\max} = 2(\nu + 1)\alpha^{1/(\nu+1)}e^{(u/(2(\nu+1)))-1}.$$

Choosing $g_{\text{max}} \leq r$ we obtain in particular (2.65), which means

$$\alpha \le \left(\frac{er}{2(\nu+1)}\right)^{\nu+1} e^{-u/2} \iff a \le \frac{1}{2} \left(\frac{er}{\nu+1}\right)^{\nu+1} \Gamma(\nu+1) e^{-u/2}.$$

For $a_0 = \frac{1}{2} \left(\frac{er}{\nu+1}\right)^{\nu+1} \Gamma(\nu+1) e^{-u/2}$, we get (2.65) since $t_0 = \alpha^{1/(\nu+1)}$. \Box

The aim is now to construct an algorithm which permits us to approximate the hitting time of the square root boundary. Therefore we consider a Bessel process of dimension δ which hits the decreasing curved boundary f(t) given by (2.63).

Algorithm (A4)—the square root boundary: $l(t) = \sqrt{\beta_0 - \beta_1 t}$ with $\beta_0 > 0$, $\beta_1 > 0$.

Let $\kappa \in]0, 1[$.

Initialization: Set $\chi(0) = 0$, $\xi_0 = 0$, $\Xi_0 = 0$, $A_0 = \kappa F_{\nu}(\beta_0, \beta_1)$. **The** (n + 1)**th step:** While the condition

$$\mathcal{L}(\Xi_n) - \sqrt{\chi(n)} > \varepsilon$$
 (denoted by $\mathbb{C}(n)$)

holds, we define

(2.66)
$$A_n = \kappa F_{\nu} \left(\left(l(\Xi_n) - \sqrt{\chi(n)} \right)^2, \beta_1 \left(1 - \frac{\sqrt{\chi(n)}}{l(\Xi_n)} \right) \right),$$

where F_{ν} is defined by (2.64), and we simulate U_{n+1} , a uniformly distributed random vector on $[0, 1]^{\lfloor\nu\rfloor+2}$, G_{n+1} , a standard Gaussian random variable and V_{n+1} , a uniformly distributed random vector on S^{δ} . U_{n+1} , G_{n+1} and V_{n+1} have to be independent. We then construct $(\xi_{n+1}, \Xi_{n+1}, \chi(n+1))$ using (2.35). The algorithm stops when $\mathbb{C}(n)$ is not longer satisfied: we set $\xi_{n+1} = 0$ and so

The algorithm stops when $\mathbb{C}(n)$ is not longer satisfied: we set $\xi_{n+1} = 0$ and so $\Xi_{n+1} = \Xi_n$ and $\chi(n+1) = \chi(n)$.

PROPOSITION 2.15. The statements of Theorems 2.12 and 2.13 are true for Algorithm (A4) associated with the square root boundary.

PROOF. All the arguments developed for decreasing boundaries with lowerbounded derivatives are easily adapted to the square root boundary. We leave the details to the reader and focus our attention to the following fact: the stochastic process $(\Xi_n, \chi(n), n \ge 0)$ stays in the domain D^0 defined by

$$D^{0} = \{(t, x) \in \mathbb{R}^{2}_{+} : l(t) - \sqrt{x} > 0\}.$$

In the *WoMS* setting, for $t \ge s$, this means that for $(\Xi_n, \chi(n)) = (s, x) \in D^0$ the following step leads to $\sqrt{\chi(n+1)} < l(\Xi_{n+1})$. By (2.35), it suffices to prove that

$$\sqrt{x} + \psi_A(t) < l(s+t)$$

for all $t \in \{u \ge 0 : \min(l(s+u), \psi_A(u)) \ge 0\}$,

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with $A = \kappa F_{\nu}((l(s) - \sqrt{x})^2, \beta_1(1 - \frac{\sqrt{x}}{l(s)}))$, since $\chi(n + 1) \leq (\sqrt{\chi(n)} + \psi_{A_n}(\xi_{n+1}))^2$. By Lemma 2.14 and due to the coefficient κ , we have

$$\psi_A(t) < \sqrt{\left(l(s) - \sqrt{x}\right)^2 - \beta_1 \left(1 - \frac{\sqrt{x}}{l(s)}\right)t}$$

Hence

$$\begin{aligned} (l(s+t) - \sqrt{x})^2 &- \psi_A(t)^2 \\ &> (\sqrt{l(s)^2 - \beta_1 t} - \sqrt{x})^2 - (l(s) - \sqrt{x})^2 + \beta_1 \left(1 - \frac{\sqrt{x}}{l(s)}\right) t \\ &> 2\sqrt{x} (l(s) - \sqrt{l^2(s) - \beta_1 t}) - \frac{\beta_1 \sqrt{x}}{l(s)} t \ge 0. \end{aligned}$$

This leads directly to (2.67). \Box

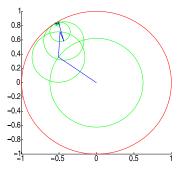
REMARK 2.16. The whole study points out a new efficient algorithm in order to simulate Bessel hitting times for given levels or curved boundaries. We can use this algorithm in two generalized situations:

(1) We have assumed that the Bessel process starts from the origin. Of course the procedure presented here can also be applied to Bessel processes starting from x > 0. It suffices to change the initialization step!

(2) We focused our attention to the Bessel process, but we linked also the initial problem to the exit time of a δ -dimensional Brownian motion from a ball of radius *l*. Algorithm (A1) extended to higher dimensions can also be used in order to evaluate exit times of general compact domains whose boundary is regular.

3. Numerical results. In this part we will illustrate the previous results on some numerical examples. Let us figure first an outcome of our algorithm, the exit position from a sphere with radius depending on time. The figure below is giving this result for an radius l = 1 and a precision $\varepsilon = 10^{-3}$.

Let us compare our algorithm with existing results. Consider the classical Euler scheme for a Brownian motion, and evaluate the first hitting time and hitting position from a disk with given radius.



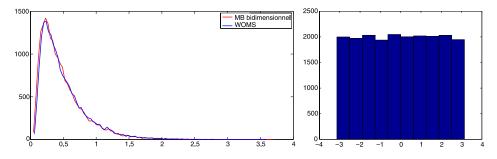


FIG. 2. Distribution of the hitting time [Euler scheme and WoMS Algorithm (A1)]—Histogram of the angle for the exit position.

First of all we can verify that the distribution of the hitting time for the *WoMS* algorithm matches the distribution of the hitting time of a given level for the 2-dimensional Bessel process. Figure 2 gives this result for a starting disk with radius 1, a precision $\varepsilon = 10^{-3}$ and a number of simulations N = 20,000. In the Euler scheme the time step is $\Delta t = 10^{-4}$.

We can also test the fact that the exit position is uniformly distributed on the circle. In order to do this we can evaluate the angle of the exit position in our *WoMS* procedure and show that it is a uniformly distributed random variable with values in $[-\pi, \pi]$. Figure 2 also shows the histogram of the result for a disk of radius 1 an $\varepsilon = 10^{-3}$ and 20,000 simulations.

Let us now present a simulation with Algorithm (A2). We consider the hitting time of the level l = 2 for the Bessel process of index $\nu = 2$, and we illustrate Theorem 2.8 by Figure 3. The curve represents the averaged number of steps versus the precision $\varepsilon = 10^{-k}$, k = 1, ..., 7. We can observe that the number of steps is better than suspected since the curve is sub-linear. We obtain the following values (for $\gamma = 0.9$ and 100,000 simulations in order to evaluate the mean).

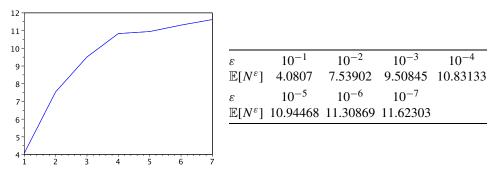


FIG. 3. Averaged number of step of Algorithm (A2) versus ε .

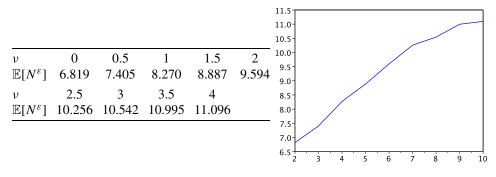


FIG. 4. Averaged number of step of Algorithm (A2) versus $\delta = 2\nu + 2$.

Finally we present the dependence of the averaged number of steps of Algorithm (A2) with respect to the dimension of the Bessel process. See Figure 4. For that purpose, we simulate hitting time of the level l = 2 with $\varepsilon = 10^{-3}$, $\gamma = 0.9$, 50,000 simulations for each estimation of the averaged value, and the dimension of the Bessel process takes value in the set {2, 3, ..., 18}.

4. Application to the Cox–Ingersoll–Ross process. We now aim to estimate the hitting time of a level l > 0 for $(X_t^{\delta}, t \ge 0)$, a Cox–Ingersoll–Ross process. The CIR process is the solution of the following stochastic differential equation:

(4.1)
$$\begin{cases} \mathrm{d}X_t^{\delta} = (a + bX_t^{\delta}) \,\mathrm{d}t + c\sqrt{|X_t^{\delta}|} \,\mathrm{d}B_t, \\ X_0^{\delta} = x_0, \end{cases}$$

where $x_0 \ge 0$, $a \ge 0$, $b \in \mathbb{R}$, c > 0 and $(B_t, t \ge 0)$ is a standard Brownian motion. We denote here $\delta = 4a/c^2$.

We will first recall a connection between this stochastic process and $(Y^{\delta}(t), t \ge 0)$, the square of the Bessel process BESQ(δ), the solution of the equation

(4.2)
$$Y^{\delta}(t) = y_0 + \delta t + 2 \int_0^t \sqrt{|Y^{\delta}(s)|} \, \mathrm{d}B_s, \qquad t \ge 0.$$

LEMMA 4.1. The CIR process has the same distribution as $(\overline{X}_t, t \ge 0)$ which is defined by

(4.3)
$$\begin{cases} \overline{X}_t = e^{bt} Y^{\delta} \left(\frac{c^2}{4b} (1 - e^{-bt}) \right), \\ \overline{X}_0 = Y^{\delta}(0), \end{cases}$$

where Y is the square of a Bessel process in dimension $\delta = 4a/c^2$; see [15].

PROOF. Let us only sketch some ideas of the proof. Let $Y^{\delta}(t)$ be the square of the δ -dimensional Bessel process. By applying Itô's formula, we get the stochastic

differential equation satisfied by the process \overline{X}_t ,

(4.4)

$$d\overline{X}_{t} = b\overline{X}_{t} dt + e^{bt} d\left(Y\left(\frac{c^{2}}{4b}(1-e^{-bt})\right)\right)$$

$$= b\overline{X}_{t} dt + b\delta \frac{c^{2}}{4b} dt + 2e^{bt}\sqrt{|e^{-bt}\overline{X}_{t}|} dB_{(c^{2}/(4b))(1-e^{-bt})}$$

$$= (a+b\overline{X}_{t}) dt + 2e^{bt/2}\sqrt{|\overline{X}_{t}|} dB_{(c^{2}/(4b))(1-e^{-bt})},$$

where $\delta = 4a/c^2$. Let us remark that

$$\frac{c^2}{4b}(1-e^{-bt}) = \int_0^t \rho^2(s) \,\mathrm{d}s \qquad \text{with } \rho(t) = \frac{c}{2}e^{-bt/2}.$$

We can deduce that there exists a Brownian motion $(\beta_t, t \ge 0)$ such that

$$B_{(c^2/(4b))(1-e^{-bt})} = \int_0^t \rho(s) \, \mathrm{d}\beta_s$$

for all $t \ge 0$. With this notation, equation (4.4) gives

$$d\overline{X}_t = (a + b\overline{X}_t) dt + 2e^{bt/2} \sqrt{|\overline{X}_t|} \rho(t) d\beta_t$$
$$= (a + b\overline{X}_t) dt + c \sqrt{|\overline{X}_t|} d\beta_t,$$

and $\overline{X}_0 = Y(0)$. This proves that the process $(\overline{X}_t, t \ge 0)$ has the same distribution as the CIR process given by (4.1). \Box

Let us consider the hitting time of a given level l for the CIR process and denote it by T_l . This time is defined by

$$T_l = \inf\{s \ge 0; X_s^{\delta} = l\}.$$

The previous Lemma 4.1 gives also an equivalence (in distribution) connecting the hitting time of the CIR process and the hitting time of the square of a δ -dimensional Bessel process.

PROPOSITION 4.2. The hitting time T_l of a level l > 0 for a CIR process has the same distribution as $-\frac{1}{b}\log(1-\frac{4b}{c^2}\tau_{\psi})$ where

$$\tau_{\psi} = \inf\left\{t \ge 0; Y^{\delta}(t) = l\left(1 - \frac{4b}{c^2}t\right)\right\},\$$

and Y^{δ} is the square of a Bessel process of dimension $\delta = 4a/c^2$.

PROOF. By using Lemma 4.1, τ_{ψ} has the same distribution as \overline{T}_l given by

(4.5)
$$\overline{T}_l = \inf\left\{s \ge 0; Y^{\delta}\left(\frac{c^2}{4b}(1-e^{-bs})\right) = le^{-bs}\right\}.$$

Define $t = \frac{c^2}{4b}(1 - e^{-bs})$, so we have two situations: *First case*: If b < 0, let $s = \eta(t)$ where

$$\eta(t) = -\frac{1}{b} \log\left(1 - \frac{4b}{c^2}t\right) \quad \text{for } t \ge 0.$$

The map η is a strictly nondecreasing function, and we thus get thus

$$\overline{T}_l = \inf\left\{\eta(t); t \ge 0, Y^{\delta}(t) = l\left(1 - \frac{4b}{c^2}t\right)\right\}$$
$$= \eta\left(\inf\left\{t \ge 0; Y^{\delta}(t) = l\left(1 - \frac{4b}{c^2}t\right)\right\}\right).$$

Second case: If $b \ge 0$, let also $s = \eta(t)$. In this case the variable t takes its values only on the interval $[0, \frac{c^2}{4b})$. So

$$\overline{T}_l = \inf\left\{\eta(t); 0 \le t \le \frac{c^2}{4b}, Y^{\delta}(t) = l\left(1 - \frac{4b}{c^2}t\right)\right\}.$$

The condition $0 \le t \le \frac{c^2}{4b}$ can be omitted in the estimation of the infimum as the boundary to hit: $1 - \frac{4bt}{c^2}$ is negative outside this interval, and the Bessel process is always positive. Furthermore the function η is also nondecreasing for $b \ge 0$, and the result is thus obtained. \Box

Application of Algorithm (A4):

An immediate consequence of Proposition 4.2 is that the hitting time T_l is related to the first time the Bessel process of dimension $\delta = 4a/c^2$ reaches the curved boundary: $f(t) = \sqrt{l(1 - \frac{4b}{c^2}t)}$. We are able to apply Algorithm (A4) if $4a/c^2 \in \mathbb{N}^*$ and b > 0 (the boundary is then decreasing). Let us denote by N^{ε} the number of steps of (A4) and $\Xi_{N^{\varepsilon}}$, the approximated hitting time of the Bessel process associated with the particular curved boundary f. Combining Propositions 2.15 and 4.2 leads to

$$\begin{split} \left(1 - \frac{\varepsilon}{\sqrt{2\alpha\pi}}\right) \mathbb{P}\left(\Xi_{N^{\varepsilon}} \leq \frac{c^2}{4b} (1 - e^{-bt}) - \alpha\right) &\leq \mathbb{P}(T_l \leq t) \\ &\leq \mathbb{P}\left(\Xi_{N^{\varepsilon}} \leq \frac{c^2}{4b} (1 - e^{-bt})\right) \end{split}$$

for α small enough and t > 0.

APPENDIX: SIMULATION OF RANDOM VARIABLES

Let us introduce simulation procedures related to particular probability density functions.

PROPOSITION A.1. Let Z be a random variable with Gamma distribution $Gamma(\alpha, \beta)$, that is,

$$\mathbb{P}(Z \in \mathrm{d}z) = \frac{1}{\Gamma(\alpha)\beta^{\alpha}} z^{\alpha-1} e^{-z/\beta} \mathbb{1}_{\{z>0\}} \mathrm{d}z, \qquad \alpha > 0, \beta > 0.$$

Then $W = \exp(-Z)$ *has the following distribution:*

$$\mathbb{P}(W \in \mathrm{d}r) = \frac{1}{\Gamma(\alpha)\beta^{\alpha}} (-\log r)^{\alpha - 1} r^{1/\beta - 1} \mathbb{1}_{[0,1]}(r) \,\mathrm{d}r.$$

In particular the stopping time τ_{ψ} defined by (2.16) has the same law as $\left[\frac{a}{\Gamma(\nu+1)2^{\nu}}\right]^{1/(\nu+1)}e^{-Z}$. Here Z is a Gamma distributed random variable with parameters $\alpha = \nu + 2$ and $\beta = \frac{1}{\nu+1}$.

PROOF. Let f be a nonnegative function. Using suitable changes of variables, we obtain

$$\mathbb{E}[f(W)] = \frac{1}{\Gamma(\alpha)\beta^{\alpha}} \int_0^\infty f(e^{-z}) z^{\alpha-1} e^{-z/\beta} dz$$
$$= \frac{1}{\Gamma(\alpha)\beta^{\alpha}} \int_0^1 f(r) (-\log r)^{\alpha-1} r^{1/\beta-1} dr$$

In order to end the proof it suffices to multiply W by a constant and use once again a change of variables formula. \Box

We need to simulate Gamma distributed variables. Let us just recall some common facts.

PROPOSITION A.2. (i) If $\alpha \in \mathbb{N}$ (so-called Erlang distributions), then the Gamma distributed variables Z has the same law as

$$-\beta \log(U_1 \cdots U_{\alpha}),$$

where $(U_i)_{1 \le i \le \alpha}$ are independent uniformly distributed random variables. Hence W defined by $W = \exp(-Z)$ can be simulated by

$$(U_1U_2\cdots U_{\alpha})^{\beta}.$$

(ii) If $\alpha - 1/2 \in \mathbb{N}$, then Z has the same law as

$$-\beta \log(U_1 \cdots U_{\lfloor \alpha \rfloor}) + \frac{\beta N^2}{2},$$

where $(U_i)_{1 \le i \le \lfloor \alpha \rfloor}$ are *i.i.d.* uniformly distributed random variables, and N is an independent standard Gaussian r.v.; see, for instance, [5], Chapter IX.3.

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