# Uncertain Darcy's problem and the stochastic particle transport

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## 1 Introduction

In this project we study a model for describing the trajectory of particles of pollutants in ground-water. A successful model for describing underground flows is given by the uncertain Darcy's problem. Given a bounded two-dimensional domain D, the pressure and velocity fields p and u are given by the solution of the following Partial Differential Equation (PDE)

$$\begin{cases}
 u = -A\nabla p, & \text{in } D, \\
 \nabla \cdot u = f, & \text{in } D, \\
 p = p_0, & \text{on } \Gamma_{in}, \\
 p = 0, & \text{on } \Gamma_{out}, \\
 \nabla p \cdot n = 0, & \text{on } \Gamma_N,
\end{cases}$$
(1)

where  $\Gamma_{in}$ ,  $\Gamma_{out}$  are the inlet and outlet portions of the boundary of D, and an impermeability condition is imposed on  $\Gamma_N$ . The coefficient A is the permeability constant of the material. In practical applications a measure of the composition of the ground in every point is not available, therefore it is necessary to model A using a random field. The problem (1) is then a random PDE, where p, u and A depend on the space variable and on the event  $\omega$ . Moreover, the smoothness of the random field A is not granted, which makes the problem harder to solve and its solution rough. One can be interested in studying the features of the trajectory of a pollutant which is dispersed in the groundwater. For example, an accurate knowledge of the characteristic time it takes for a particle to percur a fixed distance within D could help estimating security zones around extraction wells. The trajectory of pollutant particles in a transport field is intrinsically stochastic and is modeled by a Stochastic Differential Equation (SDE). In this frame, the transport field is given by the solution of the Darcy's problem and the model chosen for the particles' trajectory is the Ornstein-Uhlenbeck process. Given a starting point  $X_0$  in the domain, the trajectory X(t) is given by the solution of

$$\begin{cases} dX(t) = u(X(t))dt + \sigma dW(t), & 0 \le t \le T, \\ X(0) = X_0 \in D, \end{cases}$$
 (2)

where W(t) is a standard two-dimensional Wiener process and the scalar value  $\sigma$  represents the stochastic diffusion. Since the transport field u is defined only on the domain D, the equation (2) has to be properly equipped with boundary conditions. Both the Darcy problem and the solution of the SDE are stochastic, thus Montecarlo simulations have to be performed to estimate the trajectories of the solution of (2). Therefore, in this project we search an efficient numerical scheme in order to compute the parameters of interest of the solution.

This outline of the work is the following.

In Section 2 we investigate the performances of three numerical schemes used to approximate the mean exit time of the solution of (2) from a domain. Theoretical results regarding the weak convergence of the schemes are tested and verified.

In Section 3 we present a brief theoretical investigation regarding the convergence of the analytical and numerical solution of an SDE subject to a perturbation in the transport field.

In Section 4 we apply the techniques presented in Section 2 and 3 to the uncertain Darcy problem and we present results for the mean exit time in this case.

Finally, in Section 4 we conclude reporting some considerations about possible future developments.

# 2 Expected exit time from a domain

We aim to estimate the exit time of a particle driven by a deterministic transport field and a stochastic diffusion from a domain  $D \subset \mathbb{R}^d$ . Given a vector W(t) of m independent Brownian motions and two functions  $f \colon \mathbb{R}^d \to \mathbb{R}^d, g \colon \mathbb{R}^d \to \mathbb{R}^{d \times m}$ , we consider the following stochastic

differential equation (SDE)

$$\begin{cases} dX(t) = f(X(t))dt + g(X(t))dW(t), & 0 < t \le T, \\ X(0) = X_0, & X_0 \in D. \end{cases}$$
 (3)

The problem is completed with two different types of boundary conditions, namely

- i. killing boundaries: if the particle exits D the process is stopped,
- ii. reflecting boundaries: the particle trajectory is reflected normally inside D when it touches the boundary  $\partial D$ .

Our aim is to estimate numerically the first exit time of the solution X(t) from D up to the final time T, *i.e.*, the quantity

$$\tau = \min\{\tau_d, T\}, \text{ where } \tau_d = \min\{t \colon X(t) \notin D\}. \tag{4}$$

Let us remark that estimating the value of  $\tau$  is meaningful only if there exists a portion of the boundary  $\Gamma_k \subset \partial D$  that is endowed with killing boundary conditions. Otherwise, the process X(t) will stay in D for the whole time interval, giving as a result  $\tau = T$  for each realisation of X(t). Another quantity of interest is defined as follows

$$\phi = \phi(T, X_0, F) = \mathbb{1}_{\{\tau < T\}} F(X(T)), \tag{5}$$

where  $F: \mathbb{R}^d \to \mathbb{R}$  is a smooth function. An interesting choice of F could be the function mapping every x of  $\mathbb{R}^d$  to one. In this case, the expectation of  $\phi$  is equal to the probability for X(t) to exit the domain before the final time T. Let us choose the notation F = 1 in this case, getting

$$\Phi(T, X_0) := \mathbb{E}(\phi(T, X_0, 1)) = \Pr(\tau < T | X(0) = X_0) \tag{6}$$

In the case of general f, g and for a d-dimensional SDE, there is no closed form for  $\tau$  and  $\Phi$ . Therefore, we approximate the value of  $\tau$  by means of two numerical schemes, briefly presented in the following.

#### 2.1 Numerical Methods

#### 2.1.1 Discrete Euler-Maruyama

Given  $N \in \mathbb{N}$  let us define a partition of [0,T] as  $P_h = \{t_i\}_{i=0}^N, t_i = ih, h = T/N$ . The Discrete Euler-Maruyama method (DEM) for problem (3) is defined as follows

$$\begin{cases}
X_h^d(t_{i+1}) = f(X(t_i))h + g(X(t_i))(W(t_{i+1}) - W(t_i)), \\
X_h^d(0) = X_0.
\end{cases}$$
(7)

The exit time  $\tau$  is approximated with the quantity  $\tau_h^d$  defined as

$$\tau_h^d = \min\{\min\{t_i \colon X_h^d(t_i) \notin D\}, T\}. \tag{8}$$

We approximate analogously  $\phi$  as

$$\phi_h^d = \mathbb{1}_{\{\tau_h^d < T\}} F(X_h^d(T)). \tag{9}$$

Let us state two results concerning the weak error of this method.

**Proposition 2.1.** Under appropriate assumptions of smoothness of  $f, g, D, \partial D, F$ ,

$$|\mathbb{E}(\tau_h^d) - \mathbb{E}(\tau)| = O(\sqrt{h}). \tag{10}$$

**Proposition 2.2.** Under appropriate assumptions of smoothness of  $f, g, D, \partial D, F$ ,

$$|\mathbb{E}(\phi_h^d) - \mathbb{E}(\phi)| = O(\sqrt{h}). \tag{11}$$

An discussion of result 2.1 can be found in [6], its proof in [4]. A proof of 2.2 can be found in [2].

#### 2.1.2 Continuous Euler-Maruyama.

Let us consider the partition  $P_h$  of [0,T] as above. The Continuous Euler-Maruyama (CEM) method is defined as

$$\begin{cases}
X_h^c(t) = f(X(t_i))(t - t_i) + g(X(t_i))(W(t) - W(t_i)), & t_i < t \le t_{i+1}, \\
X_h^c(0) = X_0.
\end{cases}$$
(12)

Let us remark that in case the particle does not exit the domain,  $X_h^c(t_i) = X_h^d(t_i)$  for all  $t_i \in P_h$ . It is possible to compute the probability that a particle has exited the domain at a time t between two consecutive timesteps  $t_i, t_{i+1}$  when D is an half-space with the following formula [3]

$$\Pr(\exists t \in [t_i, t_{i+1}] \quad X_h^d(t) \notin D|X_h^d(t_i) = x_i, X_h^d(t_{i+1}) = x_{i+1}) = p(x_i, x_{i+1}, h), \tag{13}$$

with  $p(x_i, x_{i+1}, h)$  given by

$$p(x_i, x_{i+1}, h) = \exp\left(-2\frac{[n \cdot (x_i - z_i)][n \cdot (x_{i+1} - z_i)]}{hn \cdot (qq^T(x_i)n)}\right), \tag{14}$$

where  $z_i$  is the projection of  $x_i$  on  $\partial D$  and n is the normal to  $\partial D$  in  $z_i$ . At each timestep  $t_{i+1}$  we compute the probability  $p(x_i, x_{i+1}, h)$ , and then simulate a variable U distributed uniformly in the interval [0, 1], thus obtaining a realization u. Hence, we councild that the particle has left the domain for a time t in  $(t_i, t_{i+1})$  if u is smaller than p. Therefore, we approximate the exit time as

$$\tau_h^c = \min\{T, \min\{t_i = hi : X_h(t_i) \notin D\}, \min\{t_i = hi : u < p(x_{i-1}, x_i, h)\}\},$$
(15)

In the same way as in DEM, we can approximate  $\phi$  as

$$\phi_h^c = \mathbb{1}_{\{\tau_h^c < T\}} F(X_h^c(T)). \tag{16}$$

We show the pseudocode for the implementation of CEM in Algorithm 1. The weak error of this method has been studied exhaustively in previous work.

**Proposition 2.3.** Under appropriate smoothness assumptions,

$$|\mathbb{E}(\phi_h^c) - \mathbb{E}(\phi)| = O(h). \tag{17}$$

A proof of result 2.3 can be found in [3].

#### 2.1.3 Reflecting boundaries

The reflecting boundaries are treated in the same way for both DEM and CEM. Let us denote by  $\Gamma_k$  and  $\Gamma_r$  the killing and reflecting subsets of  $\partial D$ , *i.e.* 

$$\Gamma_r \cup \Gamma_k = \partial D, \quad \Gamma_r \cap \Gamma_k = \emptyset$$
 (18)

In case the particle approaches  $\Gamma_k$  the exit is treated as above. If for a timestep of  $t_i \in P_h$ ,  $X(t_i)$  is not in D and has crossed  $\Gamma_r$  at a time  $t_{i-1} < t < t_i$ , we update the solution to be the normal reflection inside D of  $X(t_i)$ .

#### 2.1.4 An adaptive procedure

Let us consider the function g in (3) to be fixed to a real constant  $\sigma$ . The numerical methods we presented can be completed with an adaptive procedure for the step size h [1]. In this frame, adaptivity is performed with respect to the position of the numerical solution  $X_h(t_i)$ , *i.e.*, the step size is reduced if the solution is near to the boundary of the domain D. If we denote by  $d_i$  the distance between  $X_h(t_i)$  and its normal projection on  $\partial D$ , by  $l \in \mathbb{N}$  a fixed parameter and by  $h_0$  a maximum bound for h, the step size is chosen using the following formula

$$h = \max \left\{ 2^{-2l} h_0, \min \left\{ 2^{-l} h_0, \left( \frac{d}{(l+3)\sigma} \right)^2 \right\} \right\}.$$
 (19)

This formula is equivalent to dividing the domain D in three zones, in particular

- an interior zone where  $h = h_{int} = 2^{-l}h_0$ ,
- a boundary zone where  $h = h_{bound} = 2^{-2l}h_0$ ,
- an intermediate zone where  $h = \left(\frac{d}{(l+3)\sigma}\right)^2$ .

Since the loss of 0.5 in the weak order of convergence of Euler-Maruyama is due to the non-detected exits from the domain D, refining the step size only near to the boundary could lead to an error equal to the one obtained using  $h_{bound}$  on the whole domain.

#### 2.2 A PDE approach

It is possible to express the mean exit time and the probability of exit from a domain in terms of the solution of partial differential equations (PDE's). Let us denote by  $\Gamma_k$ ,  $\Gamma_r$  the killing and reflecting subsets of  $\partial D$ . We consider then the expectation of the exit time from the domain D for a trajectory that at t = 0 is at position x, i.e.,

$$\bar{\tau}(x) = \mathbb{E}(\tau | X(0) = x). \tag{20}$$

Let us define the operator  $\mathcal{L}$  induced by (3) acts on a function  $u \colon \mathbb{R}^d \to \mathbb{R}$  as follows

$$\mathcal{L}u = f \cdot \nabla u + \frac{1}{2}gg^T : \nabla \nabla u, \tag{21}$$

where the : operator between two matrices A, B in  $\mathbb{R}^{d \times d}$  is defined as follows

$$A: B = \sum_{i,j=1}^{d} \{A\}_{ij} \{B\}_{ij} = \operatorname{tr}(A^{T}B).$$
 (22)

The following result allows computing the mean exit time as the solution of an appropriate PDE.

**Proposition 2.4.** Let  $\mathcal{L}$  be the differential operator defined as (21). Then, if  $\Gamma_k$  and  $\Gamma_r$  are respectively the killing and reflecting subsets of  $\partial D$ , such that  $\Gamma_k \cup \Gamma_r = \partial D$ ,  $\Gamma_k \cap \Gamma_r = \emptyset$ , the mean

exit time  $\bar{\tau}(x)$  for the solution X(t) of (3) with  $X_0 = x$  is the solution of the following boundary value problem

$$\begin{cases}
\mathcal{L}\bar{\tau}(x) = -1, & \text{in } D, \\
\bar{\tau}(x) = 0, & \text{on } \Gamma_k, \\
\nabla \bar{\tau}(x) \cdot n = 0, & \text{on } \Gamma_r,
\end{cases}$$
(23)

where n is the normal to  $\Gamma_r$ .

Further analytical treatment of the mean exit time can be found in [8, 10]. We now consider the probability of exit from D for a solution X(t) that is equal to x for t = s < T. This probability is the solution of a boundary value problem.

**Proposition 2.5.** Let  $\mathcal{L}$  be the differential operator defined as (21). Then, if  $\Gamma_k$  and  $\Gamma_r$  are respectively the killing and reflecting subsets of  $\partial D$ , such that  $\Gamma_k \cup \Gamma_r = \partial D$ ,  $\Gamma_k \cap \Gamma_r = \emptyset$ 

$$\Pr(\tau < T | X(s) = x) = \Phi(x, s, T) \tag{24}$$

where  $\Phi(x,t,T)$  is the solution of the following backwards PDE

$$\begin{cases} \frac{\partial}{\partial t} \Phi(x, t, T) + \mathcal{L}\Phi(x, t, T) = 0 & in \ D, s \le t < T, \\ \Phi(x, t, T) = 1 & on \ \Gamma_k, s \le t \le T, \\ \nabla \Phi(x, t, T) \cdot n = 0, & on \ \Gamma_r, s \le t \le T, \\ \Phi(x, T, T) = 0 & in \ D, \end{cases}$$

$$(25)$$

where n is the normal to  $\Gamma_r$ .

The proof in case  $\Gamma_k = \partial D$  of this result can be found in [13]. Further treatment in case of mixed boundary conditions and the closed form of the solution for some particular geometries of  $D \subset \mathbb{R}^2$  can be found in [5]. It is therefore possible to approximate  $\bar{\tau}$  and  $\Phi$  by means of classical methods for solving PDE's numerically, such as finite differences or the Finite Elements Method.

#### 2.3 One-Dimensional Case

We consider problem (3) in case d = 1. Given  $f, g : \mathbb{R} \to \mathbb{R}$ , an interval D = [l, r] and a Brownian motion W(t), let us consider the following one dimensional SDE

$$\begin{cases} dX(t) = f(X(t))dt + g(X(t))dW(t), & 0 < t \le T, \\ X(0) = X_0, & X_0 \in D. \end{cases}$$
 (26)

In this case, the boundary of D consists of the two points  $\{l, r\}$ . In order for the problem of the determination of  $\tau$  to be meaningful, at least one of the two points should be endowed with a killing boundary condition.

#### 2.3.1 Analytical expression of the mean exit time

In this simple frame, it is possible to deduce an analytical solution  $\bar{\tau}$  of (23). Let us consider the boundary condition at x = l fixed as *killing* and vary the boundary condition at x = r. Since the scope is deducing the exit time of a particle from D, this assumption is plausible. In this frame, it is possible to rewrite (23) as

$$\begin{cases}
f(x)\bar{\tau}'(x) + \frac{1}{2}g^2(x)\bar{\tau}''(x) = -1, & l < x < r, \\
\bar{\tau}(l) = 0, & \\
\bar{\tau}(r) = 0, & \text{if for } x = r \text{ the boundary is } killing, \\
\bar{\tau}'(r) = 0, & \text{if for } x = r \text{ the boundary is } reflecting.
\end{cases} \tag{27}$$

It is possible to show [8, 10] that  $\bar{\tau}$  is in the one-dimensional case given by

$$\bar{\tau}(x) = -2 \int_{l}^{x} \exp(-\psi(z)) \int_{l}^{z} \frac{\exp(\psi(y))}{g^{2}(y)} dy + c_{1} \int_{l}^{x} \exp(-\psi(y)) dy + c_{2}, \tag{28}$$

where the function  $\psi$  is defined as

$$\psi(x) = \int_{1}^{x} \frac{2f(y)}{g^{2}(y)} dy, \tag{29}$$

and the constants  $c_1, c_2 \in \mathbb{R}$  depend on the boundary conditions as follows

$$c_{1} = 2 \frac{\int_{l}^{r} \exp(-\psi(z)) \int_{l}^{z} \frac{\exp(\psi(y))}{g^{2}(y)} dy}{\int_{l}^{r} \exp(-\psi(y)) dy}, \text{ if for } x = r \text{ the boundary is } killing,$$

$$c_{1} = 2 \int_{l}^{r} \frac{\exp(-\psi(y))}{g(y)^{2}} dy, \text{ if for } x = r \text{ the boundary is } reflecting,$$

$$c_{2} = 0.$$
(30)

Let us remark that in case f = -V' for some smooth function V and  $g = \sigma \in \mathbb{R}$ , the expression of  $\psi$  semplifies to

$$\psi(x) = 2\frac{V(l) - V(x)}{\sigma^2}. (31)$$

The value for the expected exit time given by (28) will be used as a reference for verifying the order of convergence of the numerical methods.

#### 2.3.2 Numerical experiments - Estimation of $\tau$

We consider as a domain for (26) the interval D = [-1, 1], final time T = 5 and the following functions

$$f(x) = -V'(x)$$
, where  $V(x) = 0.1(8x^4 - 8x^2 + x + 2)$ ,  
 $q(x) = \sigma = 3$ . (32)

We approximate the value of  $\tau$  with a Montecarlo simulation of  $\tau_h^d$  and  $\tau_h^c$  computed as in (8) and (15) from the solutions provided by DEM and CEM respectively. In order to verify the order of convergence of the methods, we let N vary in the set  $2^i, i=3,\ldots,12$  and we fix the number of trajectories M to 10000. In this way, the error caused by the Montecarlo estimation should not spoil the order of convergence. In Figure 1 we show the errors obtained fixing  $X_0=0$  in both the cases of killing and reflecting boundary conditions in x=1. Moreover, in Figure 2 we show an approximation of  $\tau$  obtained with the two methods with h=T/128 and M=1000 for a set of 10 initial values equispaced along D. It is possible to remark that computing the probability of exit between two consecutive timesteps as in (14) allows correcting the overestimation of  $\tau$  obtained simply using DEM. We want to estimate the computational time for both the method. We consider M=10000, killing boundary conditions and  $N=2^i, i=3,\ldots,12$ . We remark that the computational time required by CEM is higher than for DEM if the same value of h is employed. On the other hand, fixing the error, CEM is faster than DEM in this case.

## 2.3.3 Numerical approximation of $\Phi$ with the PDE approach

Let us consider D as the interval [l,r], the boundary condition in l to be fixed to killing and in r to be either killing or reflecting. In this case and for f independent of t and  $g = \sigma \in \mathbb{R}$  (25) can be written as the following initial value PDE

$$\begin{cases} -\frac{\partial}{\partial t} \Phi(t,x) + f \frac{\partial}{\partial x} \Phi(t,x) + \frac{1}{2} \sigma^2 \frac{\partial^2}{\partial x^2} \Phi(t,x) = 0, & l < x < r \\ & \Phi(t,l) = 1, \\ & \Phi(t,r) = 1, & \text{if for } x = r \text{ the boundary is } killing \\ & \frac{\partial}{\partial x} \Phi(t,r) = 0, & \text{if for } x = r \text{ the boundary is } reflecting \\ & \Phi(0,x) = 0. \end{cases}$$

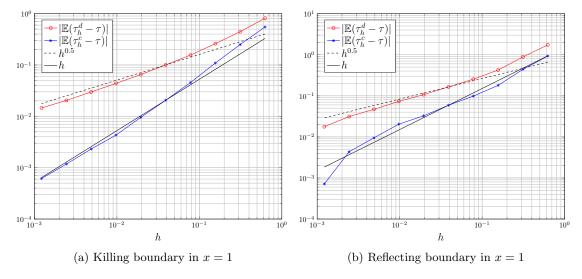


Figure 1: Approximation of  $\tau$ . Orders of convergence for DEM and CEM in the one-dimensional case.

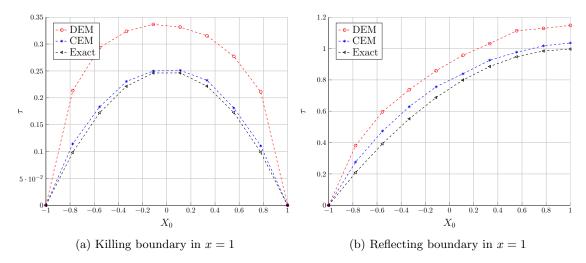


Figure 2: Approximation of  $\tau$  as a function of the initial value  $X_0$ .

This equation can be solved, e.g., using finite differences. We employ the theta method for solving (33). Let us consider the case in which r is a killing boundary, i.e., the PDE is endowed with Dirichlet boundary conditions. Given a step size  $\Delta_t$  for time integration and an uniform grid  $x_i = l + i\Delta_x, i = 0, \ldots, N+1, x_{N+1} = r$ , at each timestep k one has to find the solution of the linear system

$$(I - \Delta_t \theta A)u^{k+1} = (I + \Delta_t (1 - \theta)A)u^k + hF, \ 0 \le \theta \le 1, \tag{34}$$

where I is the identity matrix of  $\mathbb{R}^{N\times N}$ . The matrix A of  $\mathbb{R}^{N\times N}$  and the vector F of  $\mathbb{R}^N$  define the space discretization and the boundary conditions and are defined by

$$A = \frac{1}{2\Delta_x} \begin{pmatrix} \alpha_1 & \beta_1 \\ \gamma_1 & \alpha_2 & \beta_2 \\ & \ddots & \ddots & \ddots \end{pmatrix}, \quad F = \frac{1}{2\Delta_x} \left( F_1 \cdots F_N \right)^T$$
 (35)

and the coefficients are given by

$$\alpha_{i} = -\frac{2\sigma^{2}}{\Delta_{x}}, i = 1, \dots, N,$$

$$\beta_{i} = \frac{\sigma^{2}}{\Delta_{x}} + f(x_{i}), i = 1, \dots, N - 1,$$

$$\gamma_{i} = \frac{\sigma^{2}}{\Delta_{x}} - f(x_{i+1}), i = 1, \dots, N - 1,$$

$$F_{1} = \frac{\sigma^{2}}{\Delta_{x}} - f(x_{1}),$$

$$F_{N} = \frac{\sigma^{2}}{\Delta_{x}} - f(x_{N-1}).$$
(36)

The case of reflecting boundary condtion in x = r is similar and affects only the computation of the matrix A and the vector F. In particular, we introduce a ghost node at position  $x = r + \Delta_x$ , compute the derivative using a centrate approximation and impose that it is equal to 0, which leads to the condition that the value in the node in x = r is equal to the value in  $x = r - \Delta_x$ . Since the matrix defining the system (34) is tridiagonal, one can choose  $\Delta_t$ ,  $\Delta_x$  to be small and obtain a precise solution of (33) in a reasonable computational time. In the following, we will compare the values given by Montecarlo simulations using DEM and CEM with the solution of the theta method with  $\theta = 0.5$ .

#### 2.3.4 Numerical experiments - Estimation of $\Phi$

We consider (26) with D = [-1, 1], the final time T = 1, initial condition  $X_0 = 0$  and we define

$$f(x) = -V'(x)$$
, where  $V(x) = 0.1(8x^4 - 8x^2 + x + 2)$ ,  
 $g(x) = \sigma = 2$ . (37)

In order to approximate  $\Phi$ , we perform a Montecarlo simulation using both DEM and CEM, with  $M=10^5$  trajectories. We consider the number of timesteps for the time integration to be  $N=2^i, i=3,\ldots,9$ . Numerical results (Figure 3a) confirm that the weak error for DEM is of order 0.5. For CEM we have some difficulties in showing a clear order of convergence one. This is due to the fact that the results are accurate already with a big step size, therefore the statistical error is not negligible with respect to the error due to time integration. In order to avoid the noise on the order of convergence, it would be necessary to increase the number of trajectories dramatically, leading to unaffordable computational times. Let us remark that CEM approximates the exit probability better than DEM for any initial value  $X_0$ , as shown in Figure 4.

## 2.4 Two-dimensional case

We are interested in estimating the exit time of a particle from a domain  $D \subset \mathbb{R}^2$ . Given W(t) a vector of two independent Brownian motions, we consider the equation (3). In this case,  $f \colon \mathbb{R}^2 \to \mathbb{R}^2$ ,  $g \colon \mathbb{R}^2 \to \mathbb{R}^{2 \times 2}$ . We compute the mean exit time and the exit probability using DEM and CEM and compare results with the numerical solution of the PDE's presented in 2.2.

#### 2.4.1 Numerical experiments - Estimation of $\tau$

**Killing boundary conditions.** We consider a simple case of (3) in  $D = [-1, 1] \times [-1, 1]$ , where

$$f = 0 \in \mathbb{R}^2, \ g = \sigma I \in \mathbb{R}^{2 \times 2}, \sigma \in \mathbb{R}.$$

Moreover, we consider  $\partial D$  to be a killing boundary. The solution in this case is a Brownian motion. In this case, the partial differential equation (23) reduces to

$$\begin{cases}
-\sigma^2 \Delta \bar{\tau} = 2, & \text{in } D, \\
\bar{\tau} = 0, & \text{on } \partial D.
\end{cases}$$
(38)

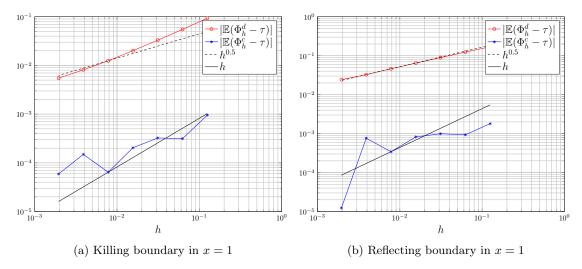


Figure 3: Approximation of  $\Phi$ . Orders of convergence of DEM and CEM in the one-dimensional case.

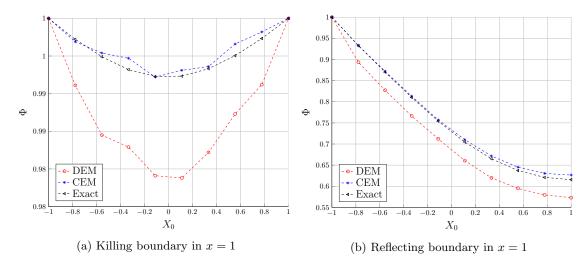


Figure 4: Profile of the probability of exit for various values of the initial condition  $X_0$  along the domain D.

This is the Poisson equation, hence it is possible to solve it numerically with the Finite Elements Method or the finite differences avoiding a high computational cost. We use the Finite Elements Method adopting a regular mesh with equal constant spacing in the x and y directions, obtaining a solution as in Figure 5b. In order to verify the orders of convergence of DEM and CEM, we set T=3,  $\sigma=1$ ,  $X_0=(0,0)^T$ , with M=10000 and  $N=2^i, i=3,\ldots,9$ . We then compare the Montecarlo estimation we obtain with the value of  $\bar{\tau}$  in (0,0), obtained by interpolation on the Finite Elements solution. The orders of convergence for this numerical experiment are shown in Figure 5a. The results confirm the theoretical orders of convergence for DEM and CEM, with an average order of 0.55 for DEM and 0.93 for CEM, which corrects to 0.98 if the last point is not taken into account.

Mixed boundary conditions. We consider the same problem as above with mixed killing and reflecting boundary conditions. f and g are the same as above, so the SDE model does not change, but we consider the two left and right boundaries of D, defined by  $x = \pm 1$ , to be reflecting. We denote this portion of the boundary as  $\Gamma_r$ , and the rest as  $\Gamma_k$ . In this case, the equation for  $\bar{\tau}$ 

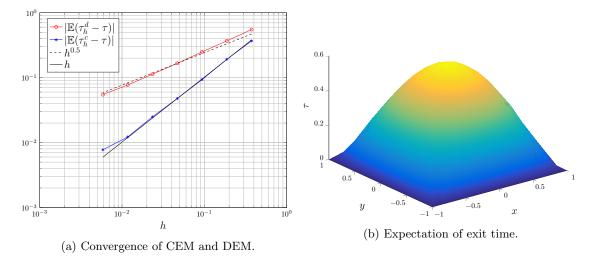


Figure 5: Summary of the results for  $\tau$  in the two-dimensional case with pure killing boundary conditions.

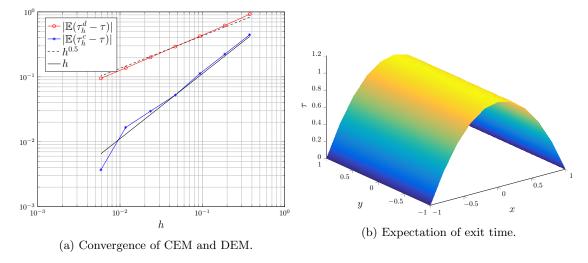


Figure 6: Summary of the results for  $\tau$  in the two-dimensional case with mixed boundary conditions.

becomes 
$$\begin{cases}
-\sigma^2 \Delta \bar{\tau} = 2, & \text{in } D, \\
\bar{\tau} = 0, & \text{on } \Gamma_k, \\
\partial \bar{\tau} \cdot n = 0, & \text{on } \Gamma_r.
\end{cases}$$
(39)

The solution of this equation is shown in Figure 6b. We compute the expectation of  $\tau$  with DEM and CEM with the same parameters as above. The results (Figure 6a), show that the theoretical orders of convergence are not spoilt by this choice of boundary conditions. The mean order for DEM in this case is 0.55, while for CEM it is 1.15.

## 2.4.2 Numerical experiments - Estimation of $\Phi$

Killing boundary conditions. We consider the same simple case as in section 2.4.1 We consider  $\partial D$  to be a killing boundary. The solution of (3) is in this case a Brownian motion. In this case,

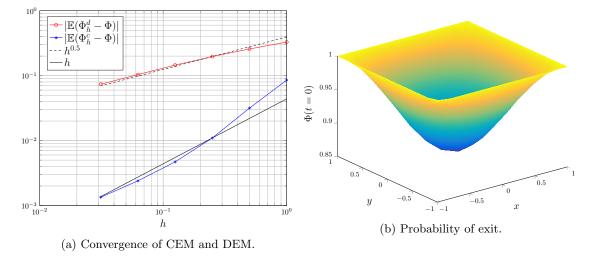


Figure 7: Summary of the results for  $\Phi$  in the two-dimensional case with pure killing boundary conditions.

the partial differential equation (25) reduces to

$$\begin{cases}
\frac{\partial}{\partial t}\Phi(x,t,T) + \frac{1}{2}\sigma^2\Delta\Phi(x,t,T) = 0, & \text{in } D,0 \le t < T, \\
\Phi(x,t,T) = 1, & \text{on } \partial D,0 \le t < T, \\
\Phi(x,T,T) = 0, & \text{in } D.
\end{cases} \tag{40}$$

We solve this problem numerically with the Finite Elements Method as for (39). The solution at t=0 is shown in Figure 7b. We verify the orders of convergence of DEM and CEM setting  $X_0=(0,0)^T$ ,  $\sigma=1,T=1$ . We consider M=100000 trajectories and  $N=2^i,i=0,\ldots,5$ . We then compare the Montecarlo estimation with the value of  $\Phi$  in (0,0), obtained by interpolation on the Finite Elements solution. The orders of convergence for this numerical experiment are shown in Figure 7a. The theoretical orders of convergence are confirmed in this case as well, with an average order of 0.43 for DEM and 1.19 for CEM.

Mixed boundary conditions. We consider the same values for the parameters, the time integration and the Montecarlo estimation as in the pure killing case. In this case, we set the boundary conditions to be reflecting on the subset of the boundary of D defined by  $x = \pm 1$  and killing for the other boundaries. Therefore, in this case the exit probability  $\Phi$  is the solution of the following PDE

$$\begin{cases}
\frac{\partial}{\partial t}\Phi(x,t,T) + \frac{1}{2}\sigma^2\Delta\Phi(x,t,T) = 0, & \text{in } D,0 \le t < T, \\
\Phi(x,t,T) = 1, & \text{on } \Gamma_k, 0 \le t < T, \\
\nabla\Phi(x,t,T) \cdot n = 0, & \text{on } \Gamma_r, 0 \le t < T, \\
\Phi(x,T,T) = 0, & \text{in } D.
\end{cases} \tag{41}$$

The solution of this equation computed with Finite Elements is shown in Figure 8b. The convergence results for DEM and CEM are shown in Figure 8a. The mean orders in this case are 0.37 for DEM and 0.87 for CEM, which is less than the prediction given by theoretical results. This decrease in the convergence rate is remarkable for small values of h. This could mean that the error caused by the Finite Element approximation of the solution of (41) is not negligible with respect to the error of CEM.

#### 2.4.3 Numerical experiments - Adaptivity

We apply to a test case the adaptivity procedure explained in section 2.1.4. In particular, we fix  $f=0, g=\sigma=1, T=3$  and  $X_0=(0,0)^T$  in the square domain  $D=[0,1]^2$  with pure killing boundary conditions. We vary l in (19) in the range  $l=0,\ldots,7$  and we run three methods

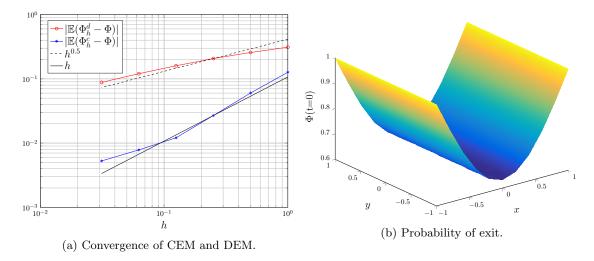


Figure 8: Summary of the results for  $\Phi$  in the two-dimensional case with mixed boundary conditions.

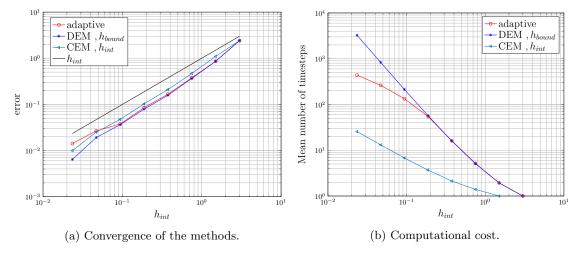


Figure 9: Estimation of  $\tau$  in the square domain with mixed boundary conditions. Comparison between DEM with fixed or adaptive step size.

- 1. Adaptive method with  $h_0 = T$
- 2. DEM with  $h = h_{bound}$  for each timestep
- 3. CEM with  $h = h_{int}$  for each timestep

We expect the three methods to have an error of order  $O(h_{int})$  when estimating the mean exit time. In Figure 9a it is possible to remark that the error is of the same order for the three methods. Moreover, the methods have approximately the same error, *i.e.*, the constant multipling  $h_{int}$  is the same for the three methods. In order to choose which method performs better from the point of view of computational cost, we compute the mean number of timesteps that they perform in order to estimate the exit time. From Figure 9b it is clear that CEM with  $h_{int}$  for every timestep performs better than the other two methods, with the adaptive procedure applied to DEM which implies lower computational time than DEM with constant timestep equal to  $h_{bound}$ . We can conclude that the approach based on Brownian bridge proposed in [3] has better performances when estimating the mean exit time from a domain.

# 3 Theoretical investigation

In this section we analyse the impact of a perturbation on the transport field on the analytical and numerical solution of an SDE. The theoretical investigation we present is needed in order to give a meaning to the results we find in the Darcy case. Since the velocity field is approximated by means of the Finite Elements method, we hope that the solution of an SDE that has the numerical approximation of the velocity field as transport term would converge to the solution of the SDE that uses the exact velocity field. Moreover, it can be costly to evaluate the Finite Element solution at each timestep of DEM or CEM, but an interpolation procedure can be exploited in order to obtain faster simulations. Hence, another source of perturbation is introduced, and a theoretical background is fundamental to strengthen the basis of our method.

## 3.1 Analysis of perturbed SDEs

Let us consider  $\sigma \in \mathbb{R}, f : \mathbb{R}^d \to \mathbb{R}^d$ ,  $W_1(t)$  a standard d-dimensional Wiener process and the following SDE

$$\begin{cases} dX(t) = f(X(t))dt + \sigma I dW_1(t), & 0 < t \le T, \\ X(0) = X_0, \end{cases}$$
(42)

where I is the identity matrix in  $\mathbb{R}^{d \times d}$ . Moreover, let us consider a perturbation of the transport field  $f^{\varepsilon} : \mathbb{R}^d \to \mathbb{R}^d$  and a standard d-dimensional Wiener process  $W_2(t)$  not necessarily independent of  $W_1(t)$ . Then, we consider the perturbed SDE

$$\begin{cases} dX^{\varepsilon}(t) = f^{\varepsilon}(X^{\varepsilon}(t))dt + \sigma I dW_2(t), & 0 < t \le T, \\ X^{\varepsilon}(0) = X_0. \end{cases}$$
(43)

Finally, let us introduce the following notation

$$||f - f^{\varepsilon}||_{\infty} = \sup_{x \in \mathbb{R}} |f(x) - f^{\varepsilon}(x)|.$$

We can state the following preliminary result

**Lemma 3.1.** With the notation above, let us consider d = 1,  $W_1 = W_2 = W$  almost everywhere. Then, given a real constant K and with the following assumptions

1. 
$$|f(x) - f(y)| \le K|x - y|, \forall x, y \in \mathbb{R}$$

2. 
$$||f - f^{\varepsilon}||_{\infty} \stackrel{\varepsilon \to 0}{\longrightarrow} 0$$
,

the solution  $X^{\varepsilon}(t)$  of (43) converges strongly to the solution X(t) of (42) in  $L^{2}(\Omega)$ , i.e.

$$\mathbb{E} \sup_{0 < t < T} |X^{\varepsilon}(t) - X(t)|^2 \xrightarrow{\varepsilon \to 0} 0.$$

*Proof.* For almost all  $\omega$ 

$$\begin{split} |X^{\varepsilon}(t) - X(t)|^2 &= \Big| \int_0^t (f^{\varepsilon}(X^{\varepsilon}(s)) - f(X(s))) ds \Big|^2 \\ &\leq T \int_0^t |f^{\varepsilon}(X^{\varepsilon}(s)) - f(X(s))|^2 ds \\ &\leq 2T \int_0^t |f^{\varepsilon}(X^{\varepsilon}(s)) - f(X^{\varepsilon}(s))|^2 ds + 2T \int_0^t |f(X^{\varepsilon}(s)) - f(X(s))|^2 ds \\ &\leq 2T^2 \|f - f^{\varepsilon}\|_{\infty}^2 + 2T^2 K^2 \int_0^t |X^{\varepsilon}(s) - X(s)|^2 ds. \end{split}$$

We then apply Gronwall's inequality, which gives

$$|X^{\varepsilon}(t) - X(t)|^2 \le 2T^2 ||f - f^{\varepsilon}||_{\infty}^2 e^{2K^2T^2}, \quad a.e.$$
 (44)

Since the right hand side of the inequality is constant, we can then take the supremum over time and the expectation at both sides, *i.e.*,

$$\mathbb{E} \sup_{0 \le t \le T} |X^{\varepsilon}(t) - X(t)|^2 \le 2T^2 ||f - f^{\varepsilon}||_{\infty}^2 e^{2K^2 T^2}.$$

Then, assumption 2. gives the result.

We now consider the case of two independent Wiener processes.

**Lemma 3.2.** With the notation above and the assumptions 1. and 2. of Lemma 3.1, let us consider  $W_1$  independent of  $W_2$  and d = 1. Then

$$\mathbb{E}\sup_{0\leq t\leq T}|X^{\varepsilon}(t)-X(t)|^2\leq 4T(T\|f-f^{\varepsilon}\|_{\infty}^2+4\sigma^2)e^{2K^2T^2}.$$

*Proof.* Let us compute the difference between  $X^{\varepsilon}(t)$  and X(t). Applying Young's inequality, we get

$$\mathbb{E} \sup_{0 \le t \le T} |X^{\varepsilon}(t) - X(t)|^{2} \le 2\mathbb{E} \sup_{0 \le t \le T} \left| \int_{0}^{t} (f^{\varepsilon}(X^{\varepsilon}(s)) - f(X(s))) ds \right|^{2} + 2\sigma^{2}\mathbb{E} \sup_{0 \le t \le T} \left| \int_{0}^{t} dW_{1}(s) - \int_{0}^{t} dW_{2}(s) \right|^{2}.$$

Let us define  $Z(t) := W_1(t) - W_2(t)$ . Z(t) is a standard Wiener process with variance 2t, thus it is a martingale. Hence, we can apply Doob's maximal quadratic inequality (e.g., [11, Page 11]) to the second term, obtaining

$$\mathbb{E} \sup_{0 \le t \le T} \left| \int_0^t dW_1(s) - \int_0^t dW_2(s) \right|^2 = \mathbb{E} \sup_{0 \le t \le T} |Z(t)|^2 \le 4\mathbb{E} |Z(T)|^2 = 8T.$$

For the first term, we can apply the same technique as in Lemma 3.1. Therefore, we get the result

$$\mathbb{E} \sup_{0 < t < T} |X^{\varepsilon}(t) - X(t)|^{2} \le (4T^{2} ||f - f^{\varepsilon}||_{\infty}^{2} + 16T\sigma^{2})e^{2K^{2}T^{2}}.$$

With those preliminary results, we can now consider the general case of a d-dimensional SDE.

**Proposition 3.1.** With the notation above and if  $f: \mathbb{R}^d \to \mathbb{R}^d$ ,  $f^{\varepsilon}: \mathbb{R}^d \to \mathbb{R}^d$ , satisfy

1. 
$$||f(x) - f(y)|| \le K||x - y||, \forall x, y \in \mathbb{R}^d$$

2. 
$$||f_i - f_i^{\varepsilon}||_{\infty} \stackrel{\varepsilon \to 0}{\longrightarrow} 0, i = 1, \dots, d,$$

then it is true for  $X(t), X^{\varepsilon}(t)$  the solutions of (42) and (43) that

$$\mathbb{E}\|X^{\varepsilon}(t) - X(t)\|_{2}^{2} \le \left(4T^{2} \sum_{i=1}^{d} \sup_{x \in \mathbb{R}} |f_{i}(x) - f_{i}^{\varepsilon}(x)|^{2} + 16Td\sigma^{2}\right) e^{2dK^{2}T^{2}}.$$
 (45)

*Proof.* The proof follows from Lemma 3.2. Let us denote by  $X_i(t), X_i^{\varepsilon}(t)$  the *i*-th component of the solution,  $i = 1, \ldots, d$ . Then

$$\mathbb{E}||X^{\varepsilon}(t) - X(t)||^2 = \mathbb{E}\sum_{i=1}^{d} |X_i^{\varepsilon}(t) - X_i(t)|^2$$
$$= \sum_{i=1}^{d} \mathbb{E}|X_i^{\varepsilon}(t) - X_i(t)|^2$$

Then, applying Lemma 3.2 to each component of the solution, one obtains (45).

Remark 3.1. (Convergence considerations.) The result of Proposition 3.1 shows that in case one uses two different Wiener processes  $W_1, W_2$  for the standard and the perturbed SDEs, the strong convergence is not granted. In fact, while the first term in estimation (45) disappears due to the uniform convergence of the perturbed transport field towards the non-perturbed one, the error due to the different Wiener processes is not disappearing with respect to  $\varepsilon$ . However, in underground flow models the variance  $\sigma$  of the Brownian diffusion is often small with respect to the transport field. Therefore, one could consider the solution  $X^{\varepsilon}(t)$  of (43) to be practically converging to the solution X(t) of (42).

Remark 3.2. (Interpolation results.) Let us consider a domain  $D \in \mathbb{R}^d$  where f and  $f^{\varepsilon}$  are defined and  $f^{\varepsilon}$  to be a polynomial interpolation of f on D. Then, the interpolation error of can be estimated if f is regular enough. In particular, the following results are valid [12, Chapter 8] for a grid of size  $\varepsilon$ 

1. f Lipschitz continuous of constant K,  $f^{\varepsilon}$  piecewise constant interpolation,  $C_0$  a real constant

$$\sup_{x \in D} \|f(x) - f^{\varepsilon}(x)\| \le C_0 K \varepsilon.$$

2. f of class  $C^1$ ,  $f^{\varepsilon}$  piecewise constant interpolation,  $C_1$  a real constant

$$\sup_{x \in D} \|f(x) - f^{\varepsilon}(x)\| \le C_1 \varepsilon \sup_{x \in D} \|f'(x)\|.$$

3. f of class  $C^2$ ,  $f^{\varepsilon}$  piecewise linear interpolation,  $C_2$  a real constant

$$\sup_{x \in D} \|f(x) - f^{\varepsilon}(x)\| \le C_2 \varepsilon^2 \sup_{x \in D} \|f''(x)\|.$$

In all these cases, it is true that  $f^{\varepsilon}$  converges uniformly to f with respect to  $\varepsilon$ , therefore  $f^{\varepsilon}$  satisfies the assumptions of Proposition 3.1.

#### 3.2 Analysis of numerical convergence

We now consider the Euler-Maruyama method applied to the perturbed equation (43). We would like to find a balance between the error due to the numerical integration of the SDE with step size h and the approximation of the transport field f with  $f^{\varepsilon}$ . In this way, one can choose wisely the two parameters  $\varepsilon$ , h in order to avoid extra computational time.

Let us consider  $X(t), X^{\varepsilon}(t)$  the solution of (42) and (43) respectively, and let us denote with  $X_n, X_n^{\varepsilon}$  the numerical solution obtained with Euler-Maruyama method applied to the two equations at time  $t_n = hn, t_N = T$ . It is known that if f is Lipschitz continuous with constant K, and if  $\sigma$  is a constant, then the strong error for  $X_n$  approximating X(t) is given by (e.g., [7, Chapter 10])

$$\sup_{n=0,\dots,N} \mathbb{E}||X(nh) - X_n|| \le Ch, \tag{46}$$

for a constant C independent of h. It is possible to obtain a similar estimate for  $X_n^{\varepsilon}$  estimating X(t).

**Proposition 3.2.** Let us consider (42), (43) with  $W_1 = W_2 = W$  almost everywhere. Given h > 0 such that hN = T for  $N \in \mathbb{N}^*$ , let us consider  $X_n^{\varepsilon}$  the numerical solution given by the Euler-Maruyama method applied to (43), i.e.,

$$\begin{cases} X_{n+1}^{\varepsilon} = X_n^{\varepsilon} + f^{\varepsilon}(X_n^{\varepsilon})h + \sigma(W(t_{n+1}) - W(t_n)), & n = 0, \dots, N-1, \\ X_0^{\varepsilon} = X_0. \end{cases}$$

Then, if  $f, f^{\varepsilon}$  satisfy the assumptions 1. and 2. of Lemma 3.1

$$\sup_{n=0,\dots,N} \mathbb{E}||X(nh) - X_n^{\varepsilon}|| \le Ch + ||f^{\varepsilon} - f||_{\infty} \frac{e^{KT} - 1}{K}, \tag{47}$$

with C a real constant independent of h and depending only on the final time T and the Lipschitz constant K of f.

*Proof.* We consider  $X_n$  the numerical approximation of X(t) obtained using Euler-Maruyama with the same initial condition  $X_0$ . If we add and substract  $X_n$  and apply the triangular inequality we obtain

$$\mathbb{E}||X_n^{\varepsilon} - X(nh)|| \le \mathbb{E}||X_n^{\varepsilon} - X_n|| + \mathbb{E}||X_n - X(nh)||.$$

Since f is regular enough and  $\sigma$  is a constant, for the second term it is known that

$$\sup_{n=0,...,N} \mathbb{E}||X_n - X(nh)|| \le Ch.$$
 (48)

We can then make a recursive analysis of the first term. For almost all  $\omega$ 

$$\begin{split} \|X_{n}^{\varepsilon} - X_{n}\| &\leq \|X_{n-1}^{\varepsilon} - X_{n-1}\| + h\|f^{\varepsilon}(X_{n-1}^{\varepsilon}) - f(X_{n-1})\| \\ &\leq \|X_{n-1}^{\varepsilon} - X_{n-1}\| + h\|f^{\varepsilon}(X_{n-1}^{\varepsilon}) - f(X_{n-1}^{\varepsilon})\| + h\|f(X_{n-1}^{\varepsilon}) - f(X_{n-1})\| \\ &\leq h\|f^{\varepsilon} - f\|_{\infty} + (1 + hK)\|X_{n-1}^{\varepsilon} - X_{n-1}\| \\ &\leq h\|f^{\varepsilon} - f\|_{\infty} + (1 + hK)(h\|f^{\varepsilon} - f\|_{\infty} + (1 + hK)\|X_{n-2}^{\varepsilon} - X_{n-2}\|) \\ &\stackrel{(\cdots)}{\leq} h\|f^{\varepsilon} - f\|_{\infty} \sum_{i=0}^{n-1} (1 + hK)^{i} + (1 + hK)^{n}\|X_{0}^{\varepsilon} - X_{0}\| \end{split}$$

Since  $X_0^{\varepsilon} = X_0$  and using the geometric sum

$$||X_n^{\varepsilon} - X_n|| \le h||f^{\varepsilon} - f||_{\infty} \frac{(1 + hK)^n - 1}{hK}$$
$$\le ||f^{\varepsilon} - f||_{\infty} \frac{(1 + hK)^N - 1}{K}$$
$$\le ||f^{\varepsilon} - f||_{\infty} \frac{e^{KT} - 1}{K},$$

where the last inequality is valid since N = T/h and K, T, h are all positive real numbers. Since the bound we found is independent of  $\omega$  and n, we can take the expectation and the supremum, obtaining

$$\sup_{n=0,\dots,N} \mathbb{E}||X_n^{\varepsilon} - X_n|| \le ||f^{\varepsilon} - f||_{\infty} \frac{e^{KT} - 1}{K}.$$
 (49)

This result combined with (48) concludes the proof.

As far as the weak convergence is concerned, it is known that the Euler-Maruyama method is of weak order one [7, Chapter 14]. The term due to the perturbation of the transport field can be treated as for the strong error, therefore we get without any further assumption for a constant C independent of h

$$\sup_{n=0,\dots,N} \|\mathbb{E}(X_n^{\varepsilon} - X(nh))\| \le Ch + \|f^{\varepsilon} - f\|_{\infty} \frac{e^{KT} - 1}{K}.$$

Remark 3.3. If  $f^{\varepsilon}$  is the interpolation of f on a regular grid of size  $\varepsilon$ , the result of Proposition 3.2 allows to balance the interpolation error and the error due to numerical integration. In fact, we reported above some results on interpolation of f with piecewise polynomials  $f^{\varepsilon}$ , which we can plug in (47) as follows

1. if f is Lipschitz continuous of constant K or of class  $\mathcal{C}^1$  and  $f^{\varepsilon}$  is piecewise constant, then

$$\sup_{n=0,\dots,N} \mathbb{E}||X_n^{\varepsilon} - X(nh)|| = O(h) + O(\varepsilon).$$

2. if f is of class  $C^2$ , and  $f^{\varepsilon}$  is piecewise linear, then

$$\sup_{n=0,\dots,N} \mathbb{E}||X_n^{\varepsilon} - X(nh)|| = O(h) + O(\varepsilon^2).$$

Therefore, in the first case the step size h should be of the same order of magnitude as  $\varepsilon$ , while in the second case it should scale as  $\varepsilon^2$ .

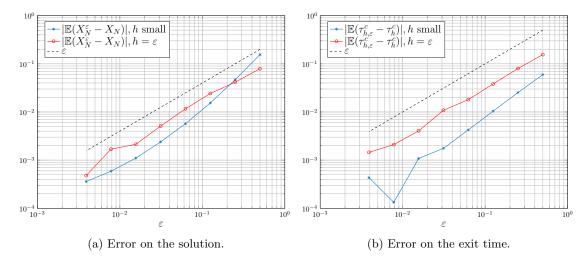


Figure 10: Convergence of the numerical solution with respect to the interpolation characteristic size  $\varepsilon$ .

#### 3.3 Numerical confirmation of the theory

We perform a numerical experiment in order to verify the theoretical bounds presented above. Let us consider the domain  $D = [-1, 1]^2$  and the deterministic transport field given by

$$f(x,y) = \frac{1}{2} (x^2 + y^2, |x - y|)^T.$$

In the domain, the transport field is Lipschitz continuous. Our aim is verifying whether the numerical solution computed using a piecewise constant interpolation of f leads to convergence with respect to the characteristic size of the grid used for interpolation. In the following,  $f^{\varepsilon}$  denotes the piecewise constant interpolation over D of f on a structured grid of equal size  $\varepsilon$  in both directions. Since the theoretical results interest the value of the solution itself and not the mean exit time from a domain, we start by considering the value itself. Then, we verify if the results are valid for the exit time as well.

Numerical solution. We consider the boundary conditions to be reflecting on all the boundary of D. We fix the parameters to be  $T=1, \sigma=1$  and perform a Montecarlo simulation over M=10000 simulations of Euler-Maruyama. Let us remark that since the boundary is completely reflecting, there is no distinction between CEM and DEM. Mantaining the notation of Proposition 3.2, we compute the weak error of  $X_n^{\varepsilon}$  with respect to  $X_n$ , which should be of order  $O(\varepsilon)$ . The reference solution  $X_n$  is computed over  $N=2^{10}$  timesteps over the time span. Then, we vary  $\varepsilon$  in the range  $\varepsilon_i=2^{-i}, i=1,\ldots,8$  and compute the numerical solution  $X_n^{\varepsilon}$  either fixing the number of timesteps  $N=2^9$  or varying it so that  $h_i=\varepsilon_i$ , i.e., the time and space discretizations have the same order of magnitude. In the first case, we wish to overkill the error due to numerical integration, while in the second case we wish to mantain an acceptable computational cost. Moreover, the bound presented in Remark 3.3 implies that theoretically the error should be independent of h, so we would expect similar results for both the approaches. Results (Figure 10a) confirm the theoretical results, with a clear rate of convergence equal to one, and the errors which are similar for the two approaches.

Mean exit time. We perform an experiment with the same values for all parameters as above, but we consider mixed killing and reflecting boundary conditions and we estimate the exit time  $\tau$ . We wish that the theoretical results obtained for the solution apply practically to the exit time itself. We use the same strategies as above, either fixing a small step size h for any value of  $\varepsilon$ , or mantaining h equal to  $\varepsilon$ . Results (Figure 10b) show that the error is  $O(\varepsilon)$  in both cases, with a smaller constant in case a small value of h is chosen. On the other side, we notice that fixing, e.g., the error to 0.01, the computational time is in this case approximately sixteen times smaller in case h and  $\varepsilon$  are balanced.

# 4 The uncertain Darcy problem

The two methods for approximating the mean exit time have been investigated in a general frame. In the following we will consider (3) with  $f: \mathbb{R}^2 \to \mathbb{R}^2$  given by the solution of the uncertain Darcy problem.

#### 4.1 Problem statement

Let us consider a domain  $D \subset \mathbb{R}^2$ . Let us define the Neumann boundaries of D as  $\Gamma_N$ , its inlet boundary as  $\Gamma_{in}$  and its outlet boundary as  $\Gamma_{out}$ . Then, we search the solution of the following problem

$$\begin{cases}
 u = -A\nabla p, & \text{in } D, \\
 \nabla \cdot u = 0, & \text{in } D, \\
 p = p_0, & \text{on } \Gamma_{in}, \\
 p = 0, & \text{on } \Gamma_{out}, \\
 \nabla p = 0, & \text{on } \Gamma_N,
\end{cases}$$
(50)

where A is a random field. The solution u of this equation is used as transport field in equation (3), which can be therefore written as

$$\begin{cases} dX(t) = u(X)dt + \sigma dW(t), & 0 < t \le T, \\ X(0) = X_0, & X_0 \in D, \end{cases}$$

$$(51)$$

where we set the boundary conditions to be reflecting on  $\Gamma_N$  and killing on both  $\Gamma_{in}$ ,  $\Gamma_{out}$ .

## 4.2 Finite Elements solution of the Darcy problem

Let us consider the domain  $D = [-1, 1] \times [-1, 1]$ . The random field A in (50) is chosen to be lognormal, *i.e.*,

$$A = e^{\gamma}, \tag{52}$$

where  $\gamma$  is a normal random field defined by its covariance function  $\operatorname{cov}_{\gamma}(x_1, x_2)$  for any couple of points  $x_1, x_2$  in the domain D. The covariance function is of the Matern family, thus having the following form

$$\operatorname{cov}_{\gamma}(x_1, x_2) = \frac{\sigma_A^2}{\Gamma(\nu) 2^{\nu - 1}} \left( \sqrt{2\nu} \frac{|x_1 - x_2|}{L_c} \right)^{\nu} K_{\nu} \left( \sqrt{2\nu} \frac{|x_1 - x_2|}{L_c} \right), \quad \nu \ge 0.5, \tag{53}$$

where  $\sigma_A^2$  is the variance,  $L_c$  is the correlation length,  $\Gamma$  is the gamma function,  $K_{\nu}$  is the modified Bessel function of the second kind and  $\nu$  is a parameter. Let us remark that the covariance function does not depend on  $x_1, x_2$  but only on their euclidean distance  $|x_1 - x_2|$ . The regularity of the covariance function and of the realizations of A depend on  $\nu$ . In particular, for  $\nu$  equal to 0.5, the covariance is Lipschitz continuous and the field is  $\alpha$ -Hölder continuous for  $\alpha < 0.5$ . Results concerning further regularity properties of A can be found in [9]. The realizations of A are computed using a discrete Fourier transformation on the vertices of a grid of D, equispaced on both the x and y directions with the same spacing  $\Delta_A$ . Then, the numerical solution  $\hat{p}$  of (50) is obtained with linear Finite Elements on a regular mesh  $T_p$  with maximum element size  $\Delta_p$ . Since the vertices of the grid on which we compute A do not coincide with the vertices of  $T_p$ , we interpolate A on  $T_p$  to obtain  $\hat{p}$ . Then, the velocity field  $\hat{u}$  is retrived computing the gradient of  $\hat{p}$ . The results for a realization of A are shown in Figure 11, where the value of the inlet pressure  $p_0$  is equal to 1, and the parameters for the random field are  $\nu = 0.5$ ,  $L_c = 0.05$ .

#### 4.3 Solution of the SDE

Once the Finite Element approximation  $\hat{u}$  of the velocity field is available, it is possible to approximate by means of DEM and CEM the solution of (51). The values of the numerical solution  $X_h$  can take any value in D, therefore it is necessary that the velocity field is defined in any point in

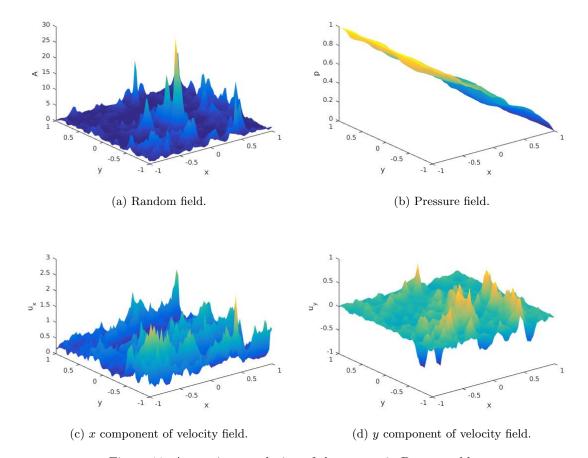


Figure 11: Approximate solution of the uncertain Darcy problem.

D. If an interpolation of  $\hat{u}$  is performed at each step, both DEM and CEM lose in computational efficiency. Hence, an interpolation of  $\hat{u}$  has to be performed before the numerical integration of the SDE. Therefore, we define a grid with spacing  $\Delta_u$ , interpolating the values of  $\hat{u}$  in the center of each square defined by the grid (Figure 12). Let us denote by Q the set of the interpolation points, whose elements are defined by

$$\{Q\}_{ij} = \left(-1 + (i - 0.5)\Delta_u, -1 + (j - 0.5)\Delta_u\right)^T, \quad i, j = 1, \dots, \frac{2}{\Delta_u} =: N_u.$$
 (54)

We compute two matrices  $U_x, U_y$  of  $\mathbb{R}^{N_u \times N_u}$  containing the values of the two components of  $\hat{u}$  interpolated on the points of Q. Then, the velocity field is considered to be piecewise constant in each square of the grid defined by  $\Delta_u$ . Therefore, if we denote by  $\tilde{u}$  the transport field for the SDE, at the *i*-th step of the integration  $\tilde{u}$  is evaluated as follows

$$\tilde{u}(X_h(t_i)) = \begin{pmatrix} U_x(\lceil (X_{h,1}(t_i) + 1)/\Delta_u \rceil, \lceil (X_{h,2}(t_i) + 1)/\Delta_u \rceil) \\ U_y(\lceil (X_{h,1}(t_i) + 1)/\Delta_u \rceil, \lceil (X_{h,2}(t_i) + 1)/\Delta_u \rceil) \end{pmatrix}, \tag{55}$$

where  $X_{h,1}, X_{h,2}$  denote the first and second components of  $X_h$  and  $U_x(i,j)$  represents the element (i,j) of the matrix  $U_x$  (respectively  $U_y$ ). Then, given the step size h, one step of DEM will be defined as

$$X_h(t_{i+1}) = \tilde{u}(X_h(t_i))h + \sigma(W(t_{i+1}) - W(t_i)). \tag{56}$$

Given an input initial value  $X_0$  for (51), we approximate the solution using DEM and CEM using the strategy above. In Figure 13 we display 15 trajectories for  $X_0 = (-0.8, -0.8)^T$  with two different timesteps. The choice of the initial point is made in order to observe reflections on the

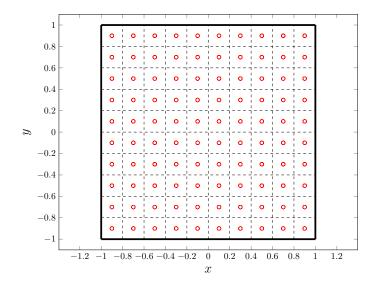


Figure 12: Grid used for interpolation of  $\hat{u}$ . Dots represent the interpolation points.

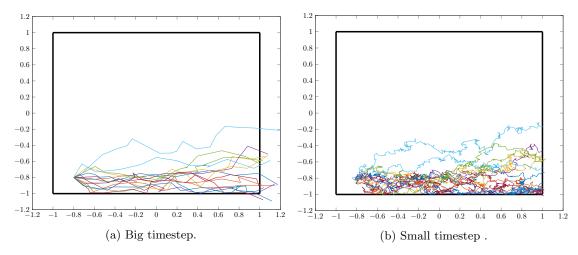


Figure 13: Trajectories of the numerical solution of (51) with DEM.

lower boundary of the domain D on which we compute the solution, as well as the killing boundary at the right side.

#### 4.4 Summary

Let us summarize the procedure we use for approximating the solution of (51).

- 1. Generate A on a grid with spacing  $\Delta_A$  with a Fourier transform method.
- 2. Approximate the solution of (50) with the Finite Element methods on a fine triangulation  $T_p$  with maximum element size  $\Delta_p$ .
- 3. Compute the velocity field from the FEM solution and interpolate it on a grid with spacing  $\Delta_u$  to obtain a piece-wise constant field  $\tilde{u}$ .
- 4. Solve (51) with DEM or CEM to evaluate the mean exit time  $\bar{\tau}$  using  $\tilde{u}$  as the transport field.

#### 4.5 Estimation of the exit time

The numerical and theoretical path we followed in the previous paragraphs enables us to build a Montecarlo simulation using which we estimate the mean exit time from a domain with the

```
Algorithm 2: Estimation of the exit time

Data: number of realizations M_r, number of trajectories M_t

Result: Estimate of the exit time \bar{\tau}

for i = 1, ..., M_r do

Generate the random field A;

Find the solution of the Darcy problem;

Interpolate the velocity field on a structured grid of size \Delta_u;

Estimate \tau_i using CEM with step size \Delta_u over M_t trajectories;

end

\tau = 1/M_r \sum_{i=1}^{M_r} \tau_i;
```

solution of the uncertain Darcy problem as a transport field. All consideration above allow us performing efficiently and accurately the estimation of the exit time for each realization of the Darcy problem. Therefore, it is sufficient to average over different realizations of the Darcy problem in order to obtain an estimation of the exit time, as in Algorithm 2. We consider the square domain  $D = [-1,1]^2$  and the parameters listed in Table 1. We vary the value of  $\sigma$  and generate 100 solutions of the Darcy problem for each value of  $\sigma$ , estimating the exit time with 10000 trajectories for each realization. The results we obtain are in Table 2. We remark that in the transport-dominated case the exit time stabilizes on a value of approximately four seconds.

Random field				Darcy	FEM	Interpolation	Trajectorie	s
ν	$L_c$	$\sigma_A$	$\Delta_A$	$p_0$	$\Delta_p$	$\Delta_u$	$X_0$	T
0.5	0.05	1	0.0039	1	$5\cdot 10^{-3}$	0.0625	$\begin{pmatrix} -0.8, & 0 \end{pmatrix}^T$	20

Table 1: Parameters for the estimation of the exit time in the Darcy case

$\sigma$	1	0.7	0.3	0.1	$10^{-2}$	$10^{-3}$	$10^{-4}$
$\tau$	0.4718	0.9906	3.1318	3.9437	4.2634	3.9767	3.9906

Table 2: Results of exit time estimations in the Darcy case for different values of  $\sigma$ .

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