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# Determination of geometrical parameters of the microstructure of a porous medium: Application to cementitious materials



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## ABSTRACT

Mass transfer in porous media is governed, besides the species diffusion coefficient, by the porosity and the geometrical parameters of the materials microstructure. This paper aims at developing a methodology for the determination of tortuosity and constrictivity factor based on a diffusion modeling in a porous medium of neutral species. Tortuosity was obtained by the computation of particles trajectory. After which, constrictivity was deduced from the formation factor, represented by the relative diffusion coefficient. Results show that tortuosity evolves in the opposite direction that the porosity while constrictivity evolves in the same direction. In addition, both parameters further slowdown the transfer for lower porosities.

# 1. Introduction

The diffusion of aggressive ions, moisture, gases and other aggressive agents in porous cementitious materials is central for predicting the durability of reinforced concrete (RC) structures. In the case of marine or deicing salt environments, chloride penetration induces the corrosion of reinforcements in RC [1]. Sulphates induces the concrete degradation [2], and the reaction of carbon dioxide with dissolved hydrates coupled with moisture transfer [3] contributes to both concrete damage and steel reinforcement corrosion [4]. These phenomena affect the durability of the materials and consequently reduces the service life of structures. The macroscopic properties of cementitious materials are linked directly to their microstructure and, more precisely, to the morphology of their pores. There are several parameters used to characterize the microstructure of cement-based materials. Porosity is one of the most used. Another essential parameter describing porous media is pore connectivity. This property directly affects the durability of cementitious materials. In fact, a high degree of pore connectivity means an interconnected porosity and consequently the aggressive agents easily penetrate into the porous network of the material. Different methods are used for studying the microstructure of cementitious materials, such as mercury intrusion porosimetry (MIP) [5], nitrogen adsorption BET [6], measurement of electrical resistivity [7], nuclear magnetic resonance [8,9], Small-Angle scattering [10], etc.

Moreover, due to the complexity of cementitious materials, the porosity and the connectivity of the pores are not sufficient to characterize their microstructure finely. Another morphological property represents a common characteristic of transport phenomena in cementitious materials, it is associated with the geometry of the pore structure. This geometry is described by two parameters: tortuosity, usually defined as the ratio of the mean effective path length of the fluid through a porous medium  $(L_e)$  and the material length (L) [11,12] and constrictivity, usually defined as the ratio between two successive different sections of pores [5]. In recent studies [13-16], the fractal theory is used to characterize the transport properties of porous media and to establish a link with geometrical parameters of microstructure such as tortuosity. These works are not interested particularly in cementitious materials, but are focused on porous media that by nature have fractal characteristics. They contribute to provide clearer answers about the general problem treated in this study in terms of identification of the geometrical parameters affecting the transport phenomena in a porous medium. Regarding cementitious materials, electrical resistivity measurements are also proposed to establish a link with the morphology of the microstructure network of cementitious materials [7,17]. These approaches involve an empirical relationship expressing the tortuosity as a function of the volume fraction of pore entrapment. Indirect methods, such as ultrasonic methods [18], were used to estimate the tortuosity from the measurements of reflected waves at two

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Nomenclature		$\mathbf{D}^{hom}$	Homogenized diffusion tensor			
		Ι	Identity tensor			
δ	Constrictivity of the porous medium	q	Position vector of particle			
Γ	Microscopic boundary of $\Omega$	v	Velocity of particle			
$\Gamma_{ff}$	Fluid-fluid interface	х	Macroscopic coordinates			
$\Gamma_{sf}$	Solid-fluid interface	У	Microscopic coordinates			
χ	Local variable	F	Geometrical factor			
Ω	Microscopic domain	D	Local diffusion coefficient in fluid phase			
$\Omega_f$	Fluid phase of $\Omega$	L	Macroscopic characteristic length			
$\Omega_s$	Solid phase of $\Omega$	1	Microscopic characteristic length			
σ	Standard deviation	$l_i$	Length path travelled by particle <i>i</i>			
τ	Tortuosity of the porous medium	Ν	Number of paths			
$\varphi$	Porosity of the porous medium	Nb <sub>c</sub>	Number of circular inclusions			
$\Delta_y$	Laplacien operator according to the microscopic co-	R	Radius of circular inclusions			
	ordinates	R <sub>max</sub>	Maximum radius of circular inclusions			
$\langle \cdot \rangle$	Average operator upon $\Omega$	R <sub>min</sub>	Minimum radius of circular inclusions			
$(\cdot)^T$	Transposition operator	$S_p$	Specific surface			
:	Arithmetic mean of ·	t	Time			
<i>div</i> <sub>y</sub>	Divergence operator according to the microscopic co- ordinates	и	Dimensionless concentration of tracer			

oblique incidence angles. However, different factors such as the geometry and the heterogeneity of the cementitious materials, the presence of steel reinforcement and moisture, make these methods less precise. Boukhatem et al. [19] used a soft computing approach to predict the transport tortuosity of the pore system of fly ash concrete by building an intelligent hybrid system. In their system, a genetic algorithm was used to optimize the structure and the hyper- parameters of the network. The system presented in their study takes into account the effect of the porous network concrete in terms of porosity, and it does not consider the transport properties (permeation or diffusion), which are governed by the tortuosity and the constrictivity of the porous network. 3D reconstructions obtained by X-ray tomography were used to quantify the tortuosity and connectivity of the pores using a specific 3D-image analysis and a random walk simulation [20-22]. Authors showed that this technique provides a promising non-destructive alternative for the pore-related characterization of cementitious materials. However, cementitious materials present a very fine micro porosity at nano-metric scale and the use of such approaches is currently limited by the spatial resolution and the precision that can be obtained. Several theoretical models were also used to assess the tortuosity of cementitious porous media [23-26], based on a certain model of the structure of a porous medium. However, they are very idealized. Empirical approaches have been also proposed to estimate a combination of the two parameters tortuosity and constirctivity of porous media as a function of porosity [27-30]. They contain parameters determined by adjusting experimental data. In contrast, these adjustable parameters vary considerably depending on the materials used and on of their pores geometry [31], and their values reported in literature vary widely.

The effect of pore geometry on the transport properties of cementbased materials makes the problem of predicting tortuosity and constrictivity extremely complex. Indeed, it is known that cementitious materials are very heterogeneous porous media and have a complex microstructure at the microscopic and nanoscopic scales [32]. The global behavior of these materials is then governed by local mechanisms acting at the pore scale. The major challenge consists of a better comprehension of physical phenomena that intervene at the scale of the pore, and their relation with the global behavior, all this taking into account the effect of pore geometry. The possible solution is the use of homogenization methods to describe finely these very heterogeneous media at the pore scale. One of the widely used homogenization methods in literature is the periodic homogenization method [33-37]. It is mainly based on the method of asymptotic developments and it assumes that the microstructure of the medium is constituted of a periodic repetition of a certain basic elementary cell. In [36,37], the periodic homogenization method is

based on a dimensional analysis of the transport equations written on the microscopic scale, making naturally appear dimensionless numbers characterizing the problem. Recent contributions in the literature have presented asymptotic periodic homogenization studies on cementitious materials, applied to heat and moisture transfers [38–40], prediction of effective mechanical, diffusive, and chemo-expansive properties [41], diffusion-reaction problems [42], alkaline - silica reaction (ASR) [43], thermal conduction [44,45], modeling of calcium leaching [46], and the transfers of chloride ions taking into account ionic electrocapillary interactions with the solid matrix [37,47–49], etc. Finally, let us cite also recent works using homogenization by two scales convergence methods [50,51].

The aim of this paper is to develop a pertinent approach to evaluate the tortuosity and constrictivity of porous media, mainly cementitious materials, by asymptotic periodic homogenization from transport equations written at the local scale for each phase of the cementitious material. The limitations mentioned above have been overcome and the effect of the pore geometry on the tortuosity and constrictivity of the material has been considered by rigorously incorporating the geometric characteristics of the microstructure. The equations of the model, developed by the upscaling micro-macro method, depend directly on the geometry of the porous medium through the homogenized parameters and the boundary value problem calculated numerically on microstructures of the material considered. The article is organized as follows. Section 2 presents a review of the homogenized model developed by asymptotic periodic homogenization. Then, the properties of the material are calculated by numerical simulations in section 3 by solving the boundary value problem which depends on the microstructure of the material. Different geometries of the porous medium were considered for the tortuosity calculation and the constrictivity deduction. Finally, final conclusions are given in section 4.

#### 2. Double-scale asymptotic method

The periodic homogenization method is one of many other upscaling techniques used to describe the physical phenomena at the global scale starting from the local one. Let us consider that the porous material studied occupies the domain *S* of the three-dimensional space  $\mathbb{IR}^3$ , whose characteristic length is denoted *L* (Fig. 1). The macroscopic domain *S* is limited by the boundary  $\partial S$ . A point of the macroscopic domain *S* will be denoted  $\mathbf{x} = (x_1, x_2, x_3)$ . The microstructure of the considered porous material is assumed to be periodic and constituted of the repetition of the elementary cell  $\Omega = \Omega_s \cup \Omega_f$  composed of a solid phase  $\Omega_s$  and a fluid phase  $\Omega_f$  (Fig. 1). The characteristic length of the elementary cell is denoted *l*.



Fig. 1. Description of the porous material.

The boundary of the domain  $\Omega$  is noted  $\Gamma$ . It is composed of the solidfluid interface  $\Gamma_{sf}$  between the solid and the fluid phases and of the fluidfluid interface  $\Gamma_{ff}$  separating two neighbouring elementary cells at the micro-scale. A point of the domain  $\Omega$  at the elementary cell will be noted  $\mathbf{y} = (y_1, y_2, y_3)$ . The characteristic length of the elementary cell *l* must be very small compared to the macroscopic characteristic length *L*.

If we consider classical molecular diffusion in cementitious materials or ionic diffusion without electrical double layer effects [48], the diffusion is governed by Fick equation or Nernst-Planck equation (without charge effects) and the homogenized diffusion tensor is given by<sup>1</sup>

$$\mathbf{D}^{\text{hom}} = \frac{1}{|\Omega|} \prod_{\Omega_f} D\left( \mathbf{I} + \left( \frac{\partial \chi}{\partial \mathbf{y}} \right)^T \right) d\Omega$$
(1)

where vector  $\boldsymbol{\chi}$  is periodic and of zero average on  $\Omega_{f_2}$  solution of the following boundary value problem:

$$\begin{cases} \operatorname{div}_{y}\left(D\left(\mathbf{I} + \left(\frac{\partial \chi}{\partial \mathbf{y}}\right)^{T}\right)\right) = 0 & \text{in } \Omega_{f} \\ \left(D\left(\mathbf{I} + \frac{\partial \chi}{\partial \mathbf{y}}\right)\right) \cdot \mathbf{n} = 0 & \text{on } \Gamma_{sf} \end{cases}$$

$$\tag{2}$$

In (2), **I** denotes the identity tensor and  $()^T$  is the transposition operator. *D* is the local diffusion coefficient in the fluid phase  $\Omega_f$ .

In this study, the local diffusion coefficient D is assumed to be constant. The boundary value problem (2) then simplified to:

$$\begin{cases}
\Delta \chi_i = 0 \text{ in } \Omega_f \\
\frac{\partial \chi_i}{\partial y} \cdot n = -n_i \text{ on } \Gamma_{sf} \\
\chi_i(y) = \chi_i(y+l) \text{ on } \Gamma_{ff} \\
\langle \chi_i \rangle = \frac{1}{|\Omega|} \int_{\Omega f} \chi_i d\Omega \text{ in } \Omega_f
\end{cases}$$
(3a,b,c,d)

where  $\chi_i$  represents the components of the vector  $\boldsymbol{\chi}$  and  $\langle \chi_i \rangle$  denotes the average of  $\chi_i$  over  $\Omega$ .

In numerous references in the literature for non reactive molecular diffusion, authors define the ratio  $D^{hom}/D$  as a geometrical factor  $\mathcal{F}$  including porosity, tortuosity and constrictivity of the pore network, when the diffusion is electrically neutral, i.e. without interactions between species or at the solid-fluid interface  $\Gamma_{sf}$ .

We propose in this work to focus on the most common definition of the geometrical factor used in literature [53-56] that is

$$\frac{D^{hom}}{D} = \frac{\varphi\delta}{\tau^2} \tag{3}$$

where  $\varphi$  denotes the porosity and  $\delta$  the constrictivity. The relationship

(4) highlights that for the particular case of a cylindrical porous network, i.e. when  $\delta$  and  $\tau$  are equal to 1, the ratio between  $D^{hom}$  and D is directly given by the porosity, i.e. by the reduction of the section through which the species diffuses. The tortuosity  $\tau$  of the porous medium is generally defined as:

$$\tau = \frac{\left(\frac{1}{N}\sum_{i=1}^{N}l_{i}\right)}{l}$$
(5)

where  $l_i$  denotes the *i*-th tortuous path, *N* is the number of paths and *l* is the size of the elementary cell.

To determine the length of possible paths  $l_i$ , we propose to solve the following auxiliary (dimensionless) diffusion problem posed on the elementary cell with associated homogeneous Neumann boundary conditions:

$$\begin{cases}
\Delta_y u = 0 \text{ in } \Omega_f \\
\frac{\partial u}{\partial y} \cdot n = 0 \text{ on } \Gamma_{sf} \\
u(0, y_1) = 1 \text{ on } \Gamma_{ff} (Inlet) \\
u(l, y_2) = 0 \text{ on } \Gamma_{ff} (Outlet)
\end{cases}$$
(6a,b,c,d)

where u presents the tracer dimensionless concentration. For this steady-problem, the paths (trajectories) are superimposed to the streamlines of problem (6). To determine trajectories, we solve the following equation of motion of massless particles:

$$\frac{d\mathbf{q}}{dt} = \mathbf{v} \tag{7}$$

where *t* denotes the time, **q** is the particle position vector of components  $\mathbf{q} = (q_{y_1}q_{y_2})$  and  $\mathbf{v}$  is the velocity function of diffusion flux streamlines obtained from the solution of problem (6) as:

$$\mathbf{v} = \left(-\frac{\partial u}{\partial y_1} - \frac{\partial u}{\partial y_2}\right)^T \tag{6}$$

After the determination of the trajectories of particles, we calculate the distance travelled by each particle. Let us notice that the path number is equal to the particle number injected in  $\Omega_f$  along the inlet interface  $\Gamma_{ff}$  where the boundary condition (6c) is imposed. Indeed, the trajectory of each particle allows to determine one value of tortuosity  $\tau_i$ . The tortuosity in the considered direction is the arithmetic average of all tortuosity particles according to eq. (5).

In parallel, we will solve closure problem (3) to determine vector  $\chi$  on the same elementary cell characterizing the porous network and compute the homogenized diffusion tensor **D**<sup>hom</sup> from eq. (1). Then, from (4), we can deduce the value of the constrictivity  $\delta$  for each considered porosity  $\varphi$ . This procedure will be applied in the next section for different elementary cell geometries.

#### 3. Numerical simulations and parametric study

The aim of this section is the determination of the tortuosity and constrictivity parameters on 2D elementary cells.

## 3.1. Case of one pore with variable size

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We consider the periodic elementary cell of size  $(l \times l)$ , which is constituted of a solid domain  $\Omega_s$  and a fluid domain  $\Omega_f$ . The interface solid-fluid  $\Gamma_{sf}$  is defined quadratic Bézier curve of three points  $(P_1(0,0); P_2(l/2, d); P_3(l, 0))$  for the bottom interface and  $(P_4(0, l); P_5(l/2, l - d); P_6(l, l))$  for the top interface (Fig. 2). The positions of points  $P_2$ and  $P_5$  varies according to the variation of parameter  $d \in [0.1 \ 0.9]$  with a step of about 0.1, leading to porosities  $\varphi \in [0.486 \ 0.943]$ . The finite element method is used to solve problems (3), (6) and (7).

Let us apply the numerical procedure described previously to

<sup>&</sup>lt;sup>1</sup> This is a classical result well established in periodic homogenization techniques whose details may be found in [48,52]. In this example, we have focused on a clearer picture of the results.

compute the tortuosity  $\tau$ . It is important to underline that according to the geometry of the elementary cell, only  $\chi_1$  is non null. Thus, we solve problem (3) for the variable  $\chi_1$  only. The periodic boundary conditions (3c) are imposed on boundaries  $\Gamma_{ff}^{-1}$  and  $\Gamma_{ff}^{-2}$ .

The problem (6) is then solved in direction  $y_1$  by imposing Dirichlet boundary conditions (6c)-(6d) on  $\Gamma_{ff}^{-1}$  and  $\Gamma_{ff}^{-2}$  with u = 1 at  $y_1 = 0$  and u = 0 at  $y_2 = l$ , respectively. The Neumann boundary conditions (3b) and (6b) of problems (3) and (6) are applied on the interface solid-fluid  $\Gamma_{sf}$ . Fig. 3(a) shows the distribution of variable  $\chi_1$ .

After solving problem (3), the homogenized diffusion tensor **D**<sup>hom</sup> is determined by using expression (1) for different values of the porosity  $\varphi$ . Fig. 3(b) shows the variation of  $D^{hom}/D$  versus porosity  $\varphi$  of the elementary cell of Fig. 2. As expected, we observe that the relative diffusion coefficient  $D^{hom}/D$  increases with the increase of the porosity.

The second step of this numerical procedure consists in solving problem (6) in order to determine *u* in fluid domain  $\Omega_f$  (see Fig. 4(a)). From *u*, we determine the velocity vector given by expression (8). The equation of motion (7) is then solved by injecting particles on the boundary  $\Gamma_{ff}^{1}$ , which represents the inlet of particles ( $\Gamma_{ff}^{2}$  is the outlet). The number of particles injected is N = 100 uniformly distributed on the inlet  $\Gamma_{ff}^{1}$ . The time step is about  $\Delta t = 5 \times 10^{-3}s$ , the simulation stops when all particles reach the outlet boundary  $\Gamma_{ff}^{2}$ . Figs. 4(b) and 4(c) show the trajectories and velocity magnitude of particles injected for two time values.

The determination of the elementary cell tortuosity of Fig. 2 in the direction  $y_1$ , requires the calculation of the length of each particle trajectory. For any particle *i* (*i* = 1, *N*), the length of path  $l_i$  is given by:

$$\Delta l_i(t) = ((q_{y_1}(t + \Delta t) - q_{y_1}(t))^2 + (q_{y_2}(t + \Delta t) - q_{y_2}(t))^2)^{\frac{1}{2}}$$
(9)

$$l_i = t_{t=0}^{l_f} \Delta l_i(t) \tag{10}$$

where  $\Delta l_i(t)$  is the increment of distance travelled at time *t* and  $t_f$  is the time required for particle *i* to reach the outlet  $\Gamma_{ff}^{2}$ .

Fig. 4(d) shows the distance travelled by each particle. We remark that in the vicinity of the solid-fluid interface  $\Gamma_{sf}$  the paths are longer. Far from  $\Gamma_{sf}$  the paths length becomes closer to the size of the elementary cell *l*. The tortuosity is finally determined from eq. (5) with N = 100 and for different values of porosity  $\varphi$  (see Fig. 5(a)).

We remark that the squared tortuosity  $\tau^2$  obtained from eq. (5) is lower than the ratio  $(D\varphi/D^{hom})$  derived from homogenization procedure. This means that the geometrical tortuosity effects should be completed by the constrictivity  $\delta$ , accounting for shrinkage or enlargement between obstacles. All these geometrical effects are taken into account in the expression of the homogenized diffusion tensor  $D^{hom}$ . The knowledge of tortuosity  $\tau$  and ratio  $(D^{hom}/D\varphi)$  enables to determinate the constrictivity  $\delta$  from eq. (4). Fig. 5(b) shows the variation of  $\delta$  versus the porosity  $\varphi$ . We notice that  $\delta$  increases with the increase of the porosity  $\varphi$ , even if its value is contained in the interval [0.9, 1] in the configuration studied.

#### 3.2. Case of a circular inclusion

Let us consider an elementary cell with a circular inclusion of radius R located in the center of the periodic cell of size  $1 \times 1$  (Fig. 6). The radius R varies in the interval  $R \in [0.05 \ 0.45]$  leading to a variation of the porosity in the range  $\varphi \in [0.3, \ 0.99]$ .

The same procedure detailed in section 3.1 is used again for the new elementary cell considered here. Because of the geometrical symmetry of the considered unit cell, we solve problems (3), (6) and (7) only in the direction  $y_1$ . In that case the homogenized diffusion tensor is isotropic. The periodic boundary conditions are imposed on the external boundary (fluid-fluid interface  $\Gamma_{ff}$ ) of the elementary cell. The boundary of the circular inclusion represents the fluid-solid interface  $\Gamma_{sf}$  where boundary conditions (3b) and (6b) are imposed. For each value of porosity  $\varphi$ , we solve the Neumann problem (3) and we determine the homogenized diffusion coefficient  $D^{hom}$  by using eq. (1).



Fig. 2. Description of the elementary cell.

coefficient  $D^{hom}/D$  versus the porosity  $\varphi$ . We notice that the homogenized diffusion coefficient increases with the increase of the porosity as observed in the example of section 3.1. The same result, which is intuitive, is also obtained in [48,57].

Then, we solve problem (6) and (7) to determine the length of all possible paths in direction  $y_1$  for different values of porosity. Fig. 7(b) shows the paths of 100 particles. We observe that the path length increases in the vicinity of the circular inclusion. Far from the inclusion, the paths are closer to an ideal path equal to the unit cell size l = 1.

The tortuosity  $\tau$  is determined from eq. (5) for different values of the porosity  $\varphi$ . Fig. 8(a) shows the variations of the tortuosity ( $\tau$  and  $\tau^2$ ) and of the ratio  $D\varphi/D^{hom}$  with respect to the porosity  $\varphi$  for the elementary cell of Fig. 6. It is important to underscore that, in the case of the circular inclusion, the tortuosity  $\tau$  follows a quasilinear behavior with respect to the porosity  $\varphi$ . We notice also that the variation of the ratio  $D\varphi/D^{hom}$  is nonlinear for  $\varphi < 0.5$ . The ratio  $D\varphi/D^{hom}$  increases sharply when the porosity is in the vicinity of the threshold of fluid phase connectivity ( $\varphi \simeq 0.33$ ). We underline that the ratio  $D\varphi/D^{hom}$  is higher than the squared tortuosity  $\tau^2$ . The difference between  $D\varphi/D^{hom}$  and  $\tau^2$  is due to the reduction of the distance between neighbouring inclusions. This effect is represented by the constrictivity factor  $\delta$  according to eq. (4), whose variation versus the porosity  $\varphi$  is plotted in Fig. 8(b). We observe that  $\delta$  increases with the increase of the porosity, with non linear variations for porosity  $\varphi < 0.5$ . In this range (low value of  $\varphi$ ), the constrictivity effect is more important because of the decrease of the distance between solid inclusions.

## 3.3. Case of a random porous media

In this section we are interested in the case of a random bi-dimensional porous medium (Fig. 9). The periodic elementary cell, of size  $1 \times 1$ , is composed of discs representing the solid phase  $\Omega_s$ , the domain between discs is the fluid phase  $\Omega_f$ . The disc boundaries represent the solid-fluid interface  $\Gamma_{sf}$ . The external boundary of the elementary cell defines the fluid-fluid interface  $\Gamma_{ff}$ .

The circular inclusions are generated randomly, the radius *R* of each disc is chosen in the interval  $R_{min} \leq R \leq R_{max}$ . In addition, positions of particle centers are also determined arbitrarily in order to avoid the overlapping between discs. For that purpose, we impose that the separation distance between the centers of disc to all neighbouring ones to be smaller than a minimum distance. With this criterion, the connectivity of the fluid phase is guaranteed. The number of discs depends on the value of the fixed porosity. As a first step, we consider four values of porosity (40 % 50 % 60 % 70%).<sup>2</sup>

 $<sup>^2</sup>$  Other numerical simulations with a global porosity of the sample between 8% and 20% could be performed to be more representative of porosities of classical concretes or mortars. In this example, we have focused on a clearer picture of the results.



Fig. 3. (a) Variable  $\chi_1$ . (b) Variation of relative homogenized diffusion coefficient versus porosity  $\varphi$ .



Fig. 4. (a) Contours of variable u. (b) Particle trajectories at t = 0.2s and (c) at t = 1.01s. (d) Path travelled by each particle.

For each porosity value, six samples are generated randomly. The chose of 6 samples seems to be sufficient to take into account a geometric variability. Table 1 summarizes the geometrical properties considered for the corresponding samples for all porosities. The minimum and maximum radii are

fixed to be  $R_{min} = R_{max}/20$  and  $R_{max} = 0.15 \times l$  where l = 1 is the size of the unit cell. We determine for each sample the mean radius  $\overline{R}$  and the standard deviation  $\sigma$ . Finally  $Nb_c$  denotes the number of circles necessary to reach the fixed value of the porosity.



Fig. 5. (a) Variation of tortuosity versus porosity  $\varphi$ . (b) Variation of constrictivity  $\delta$  versus porosity  $\varphi$ .



Fig. 6. Elementary cell with circular inclusion.

The disc size distribution curve is presented in Fig. 10, for four values of the porosity. We remark for several samples that when the porosity increases, the dispersion of the size disc distribution becomes important. We observe that  $\overline{R}$  and  $\sigma$  are closer for lowest porosity (see Table 1). Progressively, the gap increases with the increase of the porosity.

As previously, we adopt the procedure outlined in section 3.1. Firstly, we solve the boundary value problem (3) in direction  $y_1$  and  $y_2$ 

to determine the components  $\chi_1$  and  $\chi_2$  of vector  $\chi$ . After that, we calculate the components of the homogenized diffusion tensor using (1). In that case, **D**<sup>hom</sup> is an anisotropic tensor.

In a second time, we solve the diffusion problem (6) to determine scalar variables  $u_1$  and  $u_2$  in direction  $y_1$  and  $y_2$  respectively. The periodic boundary conditions (3c) of problem (3) are applied on the external fluid-fluid interfaces  $\Gamma_{ff}$ . Neumann boundary conditions (3b) and (6b) of problems (4) and (6) are applied on the solid-fluid interface  $\Gamma_{sf}$ . In addition, boundary conditions (6b) are applied on  $\Gamma_{ff}^3$  and  $\Gamma_{ff}^4$  ( $\Gamma_{ff}^1$ and  $\Gamma_{ff}^{-2}$ ), when problem (6) is solved in direction  $y_1$  (in direction  $y_2$ respectively). The Dirichlet boundary conditions (6c)-(6d) are applied on  $\Gamma_{ff}^{-1}$  with  $u_1 = 1$  and on  $\Gamma_{ff}^{-2}$  with  $u_1 = 0$  in direction  $y_1$  (on  $\Gamma_{ff}^{-3}$  with  $u_2 = 1$  and on  $\Gamma_{ff}^{-4}$  with  $u_2 = 0$  in direction  $y_2$ ).

Afterwards, we solve the motion eq. (7) of massless particles by injecting N = 200 particles on boundary  $\Gamma_{ff}^{1}$  and on boundary  $\Gamma_{ff}^{3}$  corresponding to the inlet, according to  $y_1$  and  $y_2$  directions. The N particles injected are uniformly distributed on the inlet boundaries and they are collected on boundaries  $\Gamma_{ff}^{2}$  and  $\Gamma_{ff}^{4}$ . The velocity of particles is then computed from the flux diffusion streamlines obtained using the gradient of variables  $u_1$  and  $u_2$  as follows:

$$\mathbf{v}_{1} = \begin{pmatrix} -\frac{\partial u_{1}}{\partial y_{1}} \\ -\frac{\partial u_{1}}{\partial y_{2}} \end{pmatrix}; \qquad \mathbf{v}_{2} = \begin{pmatrix} -\frac{\partial u_{2}}{\partial y_{1}} \\ -\frac{\partial u_{2}}{\partial y_{2}} \end{pmatrix}$$
(11)



Fig. 7. (a) Variation of  $D^{hom}/D$  versus porosity. (b) Path travelled by each particle for  $\varphi = 49.7\%$ .



Fig. 8. (a) Variation of tortuosity versus  $\varphi$ . (b) Variation of constrictivity versus  $\varphi$ .



Fig. 9. Periodic elementary cell of the random porous medium (bi-dimensional case).

mesh is refined in the vicinity of solid-fluid interface  $\Gamma_{sf}$  (Fig. 11), where the code color represents the normalized size of triangular element of the mesh.

Fig. 12(a) and 12(b) show the average<sup>3</sup> of components  $\chi_1$  and  $\chi_2$  of vector  $\chi$  in direction  $y_1$  and  $y_2$ , respectively for sample number 5 with porosity of 40%. From vector  $\chi$ , we calculate the components of homogenized diffusion tensor  $\mathbf{D}^{\text{hom}}/D$ . The numerical values of the relative diagonals values  $D_{11}^{hom}/D$  and  $D_{22}^{hom}/D$  are summarized in Table 2. We notice that the values of  $D_{11}^{hom}/D$  and  $D_{22}^{hom}/D$  present some dispersion for fixed porosity (see Fig. 13) due to the difference on the grain size distribution presented in Fig. 10. This dispersion becomes more important with the increase of the porosity.

The variations of the mean value and the standard deviation (error bar) of  $D_{11}^{hom}/D$  and of  $D_{22}^{hom}/D$  as function of porosity are presented in Fig. 13. In that case, the mean value is the arithmetic average on the samples with fixed porosity. As expected, the relative homogenized diffusion coefficients increase with  $\varphi$ . We observe that the mean values of  $D_{11}^{hom}/D$  and of  $D_{22}^{hom}/D$  are close due to the relatively large number of samples considered for each porosity and to the random

6 109					
109					
7.65					
1.41					
2.86					
2.58					
$\varphi = 70\%$					
6					
18					
9.45					
1.43					
6.13					
4.04					
-					

 Table 1

 Geometrical parameters of the six samples considered for each value of porosity.

The determination of the vector position from eq. (7) enables to calculate the distance travelled by each particle in the two-dimensional porous medium. Finally, we determine the tortuosity using eq. (5) and we deduce the constrictivity form eq. (4).

process used. In this case, we can consider the homogenized diffusion tensor and the microstructure as (nearly) isotropic.

We use finite element method to solve problems (3), (6) and (7). The

<sup>&</sup>lt;sup>3</sup> The average is calculated on the six samples tested and listed in Table 1.



Fig. 10. Grain size distribution curve for different porosities.



Fig. 11. Mesh size distribution (the code color corresponds to the size of elements).

Figs. 14(a) and 14(b) show the trajectories in the direction  $y_1$  and  $y_2$  respectively with a normalized magnitude velocity distribution.

Tortuosity  $\tau_{11}$  and  $\tau_{22}$  are determined from eq. (5) in the direction  $y_1$  and  $y_2$  respectively. These values are presented in Table 2 for each sample.

Fig. 15 shows the variation of the mean value of the tortuosity  $\overline{\tau}$  and  $\overline{\tau}^2$  with respect to the porosity  $\varphi$  in direction  $y_1$  and  $y_2$ . In the same figure, the variation of the mean values of the ratios  $D\varphi/D_{11}^{hom}$  and  $D\varphi/D_{22}^{hom}$  versus porosity are plotted. The difference between mean value of  $\tau_{ii}^2$  and the ratio  $D\varphi/D_{i1}^{hom}$  can be quantified by the constrictivity  $\delta_{11}$  and  $\delta_{22}$  in direction  $y_1$  and  $y_2$  as in previous cases (Table 2).

Fig. 16(a) shows the variation of the average constrictivity versus the porosity in directions  $y_1$  and  $y_2$ . We note that the constrictivities  $\overline{\delta_{11}}$  and  $\overline{\delta_{22}}$  increase with porosity and tend to unity for high porosity

values. In addition, we observe that the constrictivity  $\overline{\delta_{11}} < \overline{\delta_{22}}$ , except for porosity of 50%. This is due to the variation of homogenized diffusion tensor components, which present the same behavior in Fig. 13 (even if it is less marked). The specific surface  $S_p = \frac{|\Gamma_{sf}|}{|\Omega|}$  takes into account the size and number of inclusions composing the numerical porous medium. For a fixed porosity, we note that the specific surface is different for different samples having the same porosity (Table 2). For a fixed value of the porosity, the specific surface presents a dispersion linked to the random way of generating the random porous medium. This dispersion is more important for smallest porosities and decreases strongly for higher porosities (Fig. 16(b)). We notice that the tortuosity increases and constricitivity decreases with the increase of the specific surface  $S_p$ . Indeed, when the discs are closer, the paths length required to bypass the solid obstacles increases. If constrictivity is considered to be the ratio of smallest to highest pore sizes, the increase of  $S_p$  leads to a pore size decrease, which corresponds to a decrease of constricivity.

#### 3.4. Discussion and comparison

In this numerical study, we are interested in the calculation of the geometrical parameters of porous media and in the determination of the tortuosity and constrictivity with particular application to cementitious materials. The first case studied concern a unit cell constituted of a single pore with variable size. This simple geometry allowed us to highlight the effect of constrictivity and tortuosity in the diffusion process in a porous media and to extract its effect from the macroscopic ratio  $D\varphi/D^{hom}$  obtained by periodic homogenization and often used in literature. These results were confirmed by the second example considered (a unit cell with a circular inclusion) where the effects of tortuosity and constrictivity are more important. In both cases, we observe the decreasing of tortuosity and the increasing of constrictivity when porosity increases. The last more complex case studied is a random porous media composed of polydisperse circular inclusions. The variation of  $D\varphi/D^{hom}$  is slightly different in the two directions due to a non perfectly isotropic granular assembly. The same



Fig. 12. Numerical results for the sample number 5 with a porosity  $\varphi = 0.4$ . (a) Distribution of  $\chi_1$ . (b) Distribution of  $\chi_2$ .

 Table 2

 Numerical results obtained on the six different samples for each value of porosity.

	$\varphi = 40\%$						$\varphi$ = 50%					
Samples	1	2	3	4	5	6	1	2	3	4	5	6
$S_p$	30.23	27.80	30.39	34.69	42.11	28.55	17.16	15.49	16.93	14.03	25.74	19.59
$D_{11}^{hom}/D$	0.216	0.204	0.211	0.212	0.223	0.212	0.283	0.271	0.302	0.317	0.275	0.294
$D_{22}^{hom}/D$	0.211	0.215	0.216	0.219	0.218	0.215	0.296	0.278	0.275	0.266	0.311	0.288
$\tau_{11}$	1.177	1.173	1.171	1.158	1.163	1.159	1.128	1.117	1.116	1.119	1.141	1.130
$\tau_{22}$	1.173	1.159	1.172	1.149	1.168	1.157	1.131	1.122	1.134	1.128	1.138	1.142
$\delta_{11}$	0.637	0.598	0.618	0.614	0.649	0.615	0.640	0.608	0.675	0.710	0.628	0.667
$\delta_{22}$	0.619	0.625	0.635	0.631	0.636	0.623	0.670	0.626	0.625	0.600	0.710	0.660
	$\varphi = 60\%$						$\omega = 70\%$					
							1					
Samples	1	2	3	4	5	6	1	2	3	4	5	6
Samples	1 8.49	2 14.37	3 12.50	4 11.23	5 11.09	6 14.57	1 8.66	2	3 8.54	4 8.51	5 8.12	6 6.94
Samples $S_p$ $D_{11}^{hom}/D$	1 8.49 0.369	2 14.37 0.400	3 12.50 0.388	4 11.23 0.374	5 11.09 0.353	6 14.57 0.383	1 8.66 0.487	2 7.66 0.461	3 8.54 0.511	4 8.51 0.494	5 8.12 0.493	6 6.94 0.546
Samples $S_p$ $D_{11}^{hom}/D$ $D_{22}^{hom}/D$	1 8.49 0.369 0.375	2 14.37 0.400 0.381	3 12.50 0.388 0.330	4 11.23 0.374 0.388	5 11.09 0.353 0.412	6 14.57 0.383 0.401	1 8.66 0.487 0.506	2 7.66 0.461 0.535	3 8.54 0.511 0.509	4 8.51 0.494 0.504	5 8.12 0.493 0.514	6 6.94 0.546 0.465
Samples $S_p$ $D_{11}^{hom}/D$ $D_{22}^{hom}/D$ $\tau_{11}$	1 8.49 0.369 0.375 1.100	2 14.37 0.400 0.381 1.100	3 12.50 0.388 0.330 1.093	4 11.23 0.374 0.388 1.102	5 11.09 0.353 0.412 1.117	6 14.57 0.383 0.401 1.107	1 8.66 0.487 0.506 1.068	2 7.66 0.461 0.535 1.092	3 8.54 0.511 0.509 1.094	4 8.51 0.494 0.504 1.076	5 8.12 0.493 0.514 1.072	6 6.94 0.546 0.465 1.063
Samples $S_p$ $D_{11}^{hom}/D$ $D_{22}^{hom}/D$ $\tau_{11}$ $\tau_{22}$	1 8.49 0.369 0.375 1.100 1.106	2 14.37 0.400 0.381 1.100 1.112	3 12.50 0.388 0.330 1.093 1.095	4 11.23 0.374 0.388 1.102 1.089	5 11.09 0.353 0.412 1.117 1.109	6 14.57 0.383 0.401 1.107 1.119	1 8.66 0.487 0.506 1.068 1.080	2 7.66 0.461 0.535 1.092 1.077	3 8.54 0.511 0.509 1.094 1.079	4 8.51 0.494 0.504 1.076 1.077	5 8.12 0.493 0.514 1.072 1.090	6 6.94 0.546 0.465 1.063 1.082
$Samples$ $S_{p}$ $D_{11}^{hom}/D$ $D_{22}^{hom}/D$ $\tau_{11}$ $\tau_{22}$ $\delta_{11}$	1 8.49 0.369 0.375 1.100 1.106 0.679	2 14.37 0.400 0.381 1.100 1.112 0.733	3 12.50 0.388 0.330 1.093 1.095 0.712	4 11.23 0.374 0.388 1.102 1.089 0.686	5 11.09 0.353 0.412 1.117 1.109 0.659	6 14.57 0.383 0.401 1.107 1.119 0.708	1 8.66 0.487 0.506 1.068 1.080 0.743	2 7.66 0.461 0.535 1.092 1.077 0.719	3 8.54 0.511 0.509 1.094 1.079 0.799	4 8.51 0.494 0.504 1.076 1.077 0.759	5 8.12 0.493 0.514 1.072 1.090 0.755	6 6.94 0.546 0.465 1.063 1.082 0.829



**Fig. 13.** Variation of the mean value of the relative homogenized diffusion coefficients in direction  $y_1$  and  $y_2$  versus porosity. The error bar represents the dispersion (standard deviation) of the results.

tendency is still observed with an amplification of the variations. In addition, the constrictivity in this case presents some dispersion due to the random way to generate the porous medium. The specific surface which takes into account the variation of the number and size inclusions, seems to have also an important effect on the constrictivity.

Let us now try to compare<sup>4</sup> the order of magnitude the results obtained here to some existing experimental and numerical models of the literature (Figs. 17 and 18). First, we propose to compare the ratio  $D\varphi/D_{ii}^{hom}$  for random porous medium to experimental data obtained from diffusion test of chloride in cement paste. This experimental data are represented by the fitted evolution laws of the measured values. In addition, numerical models are used for this comparison as the results of Feng et al. [58] and Damrongwiriyanupap et al. [59] which give a good approximation of chloride diffusivity in cement paste. We notice

<sup>&</sup>lt;sup>4</sup> It is important to underscore that the comparison is mainly illustrative as our simulations are in 2D whereas the experimental results are obtained in 3D. Moreover, the latter have been fitted and sometimes extrapolated for a possible comparison.



Fig. 14. Numerical results for the sample number 5 with a porosity  $\varphi = 0.4$ . (a) Particle trajectories with the normalized velocity in direction  $y_1$ . (b) Particle trajectories with the normalized velocity in direction  $y_2$ .



**Fig. 15.** Variation of the mean value of the tortuosity in direction  $y_1$  and  $y_2$  versus porosity.

that our results are of the same order as the experimental data for the high porosities, the gap becomes more important for the low porosities. On Fig. 18, we compare the squared tortuosity obtained for the

random porous media of Fig. 9 to the some empirical expressions obtained from numerical simulations of 2D and 3D geometry [31,60–63].



Fig. 17. Comparison of  $D\varphi/D^{hom}$  with experimental and numerical results in literature.

We remark that there is significant differences between our model and models of random squared overlapping inclusions [60–62]. For high porosities ( $\varphi > 0.5$ ), our model gives squared tortuosity value closer to those obtained in the case of the cubic and tetrahedral packings of monosized spherical inclusions [63]. In the case of a circular inclusion,



Fig. 16. (a) Variation of constrictivity mean value versus porosity. (b) Variation of specific surface mean value versus porosity.



**Fig. 18.** Comparison of  $\overline{\tau^2}$  with numerical results in literature.

the model developed by Sun el al. [31], where the tortuosity is expressed with respect to porosity as  $\tau = 1 - p \ln (\varphi)$ , is very closer to our results with p = 0.409 (see eq. (24) of [31]).

#### 4. Conclusion

In this study, we presented a new approach to calculate numerically the tortuosity and the constrictivity of the pores network of a porous medium, for application to cementitious materials. In this context, the asymptotic periodic homogenization method was used to calculate a homogenized diffusion coefficient of non-reactive and non-charged particles. The latter is directly related to the geometry of the porous medium microstructure considered by means of the formation factor. It has been shown that the diffusion coefficient increases with the increase in the porosity. Then, the monitoring of particles injected in the porous network allowed calculating numerically the tortuosity of the medium. It was noted that the calculated tortuosity value is lower than the relative diffusion coefficient, which derives from the homogenization procedure. The difference is attributed to the constrictivity factor, which is deduced from this difference. In addition, the study presented in this paper aims at considering the effect of various parameters characterizing the microstructure, such as porosity, grain size distribution, specific surface of the pores and geometry of the porous medium, on the tortuosity and the constrictivity of pore network. The results indicate that:

- The values of the constrictivity evolve as the porosity. Regarding the tortuosity, its evolution is inversely proportional to the porosity of the medium.
- The grain size distribution affects the values of the tortuosity and constrictivity of the pore network, due to the variation of the specific surface. In fact, for the same porosity, the tortuosity increases and the constrictivity of pore network decreases with the increase in the specific surface. In this case, the solid inclusions are larger and closer, they induce an increase in the paths necessary to get around the solid phase in the medium.
- For representative cells with generated inclusions, the study proved that the length of the particle path increases in the vicinity of the inclusions which oppose their trajectory. Consequently, a smaller constrictivity of the pore network and a slightly higher tortuosity were observed.
- Finally, concerning the most realistic case where the microstructure of the material is generated in a random way, the calculated tortuosity and constrictivity are different in the main directions of the medium due to the anisotropy of the microstructure.

This work provides fundamental data for a better understanding of the tortuosity and constrictivity of pore networks of complex materials. However, the geometries considered in this study allow obtaining porosities that can correspond to those of usually cement pastes and mortars. On the other hand, the case of concrete, with lower porosity (sometimes < 15%), was not tackled in this study. It is therefore possible subsequently to extend this work to analyze in depth the effect the microstructural parameters governing the durability of concrete. We can consider the implementation of an approach allowing studying geometries corresponding to those of cementitious materials with very low porosities and in 3D dimensions for more realistic microstructure models.

# **Declaration of Competing Interest**

The authors whose names are listed immediately below certify that they have NO a\_liations with or involvement in any organization or entity with any \_nancial interest (such as honoraria; educational grants; participation in speakers' bureaus; membership, employment, consultancies, stock ownership, or other equity interest; and expert testimony or patent-licensing arrangements), or non-\_nancial interest (such as personal or professional relationships, a\_liations, knowledge or beliefs) in the subject matter or materials discussed in this manuscript.

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