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Multiscale Stochastic Modeling of Random Anisotropic Elastic Media With a Complex Microstructure

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Abstract: A generic microscale stochastic model is presented for a large class of random anisotropical elastic microstructures allowing a parametric analysis of the Representative Volume Element (RVE) size to be performed. This new approach can be useful for a direct experimental identification of random anisotropic elastic microstructures when the standard method cannot easily be applied. Such a RVE is used to construct the macroscopic properties in the context of stochastic homogenization. The probability analysis is not performed as usual for a given particular random microstructure defined in terms of its constituents, but is performed for a large class of random anisotropic elastic microstructures. For this class, the probability distribution of the random effective stiffness tensor is explicitly constructed and allows a full probability analysis of the RVE size to be carried out with respect to the spatial correlation length of the random microstructure.

Keywords: Microstructures, probability and statistics, anisotopic material, elastic material, inhomogeneous material

Introduction

The probabilistic model of a random microstructure (such as a composite constituted of several constituents) can directly be constructed from the geometry and mechanical properties of its constituents. This is the case for the class of random heterogeneous materials whose microstructures can be modeled as a distribution of inclusions or cavities of well-defined geometry in a given matrix, see for instance (Willis, 1982), (Torquato and Stell, 1985), (Drugan and Willis, 1996), (Torquato, 1997), (Quintanilla and Torquato, 1997), (Nemat-Nasser and Hori, 1999), (Quintanilla, 1999), (Ostoja-Starzewski, 1998), (Kachanov et al., 2001), (Milton, 2002), (Roberts and Garboczi, 2002), (Torquato, 2002), (Drugan, 2003), (Monetto and Drugan, 2004). Generally, the statistics-based bounding techniques only use the lower-order statistics (first- and second-order moments) and the probability distributions which give the detailed probabilistic information are not taken into account. Recently, a global stochastic model of the local stiffness tensor of the random microstructure has been constructed using chaos decomposition (Jardak and Ghanem, 2004). The random microstructure can be homogenized if there is a representative volume element (RVE) size such that the random fluctuations of the random effective stiffness tensor around the statistical mean value of the random effective stiffness tensor is "negligible". RVE size has received a particular attention; see for instance (Cailletaud et al., 1994), (Drugan and Willis, 1996), (Gusev, 1997), (Nemat-Nasser and Hori, 1999), (Kanit et al., 2003), (Monetto and Drugan, 2004), (Sab and Nedjar, 2005).

The prime objective of this paper is not to analyse a particular random isotropic or anisotropic microstructure described in terms of its constituents, but is to propose (when the standard method cannot easily be applied to anisotropic elastic microstructures) a new way which could be useful for a direct experimental identification of random anisotropic elastic microstructures introducing a microscale stochastic model. The main idea of this paper is then to directly introduce a microscale stochastic model of the random anisotropic elastic microstructure, which is not deduced from the stochastic models of its constituents. Such a microscale stochastic model must verify fundamental mathematical properties to obtain a physical model of any anisotropic elastic microstructure. The random anisotropic elastic microstructure (for instance a mortar constituted of a cement paste with embedded sand

particles, some porous media such as plaster boards, some cortical bones, some biological membranes and more generally, some living tissues, etc) is then modeled by an equivalent random continuous anisotropic elastic medium which is completely defined by its local stiffness fourth-order tensor-valued random field $\mathbf{x} \mapsto \mathbb{C}(\mathbf{x}) = \{\mathbb{C}_{ijkh}(\mathbf{x})\}_{ijkh}$. The random field \mathbb{C} is then constituted of 21 mutually dependent real-valued random fields modeling the anisotropic microstructure at the microscale level. The theory proposed allows strong anisotropic random fluctuations to be taken into account. Such an equivalent random anisotropic elastic medium can also be viewed (but it is not necessary) as the stochastic homogenization of the random anisotropic microstructure on a microscale RVE. The great interest of such a direct construction of a microscale stochastic model of the random anisotropic elastic microstructure is the capability to identify the parameters of the random field $\mathbf{x} \mapsto \mathbb{C}(\mathbf{x})$ using strain measurements on the boundary of tested specimens at the RVE- or macro-scale and solving an inverse stochastic problem. It is then necessary to choose a stochastic representation of $\mathbf{x} \mapsto \mathbb{C}(\mathbf{x})$ in a class of random fields for which only a few parameters are required to define its system of marginal probability distributions. Therefore, the inverse problem related to such an experimental identification of $\mathbf{x} \mapsto \mathbb{C}(\mathbf{x})$ is more feasible. In addition, such a class of random fields $\mathbf{x} \mapsto \mathbb{C}(\mathbf{x})$ must be constructed using only the available information and not "hypothetical" information for which no statistics are available or for which the number of experimental specimens is too small to obtain a good convergence of the statistical estimators. For the tensor-valued random field $\mathbf{x} \mapsto \mathbb{C}(\mathbf{x})$, the largest class can be constructed using as available information: the symmetry properties, the mean value $\mathbf{x} \mapsto \underline{\mathbb{C}}(\mathbf{x})$ which is assumed to be known and a stochastic nonuniform ellipticity condition for the corresponding linear elasticity stochastic differential operator. Clearly, any random anisotropic elastic microstructure belongs to this class. The results presented in this paper could allow the microscale stochastic model to be identified from RVE- or macro-scale measurements solving an inverse stochastic problem.

1. Macroscopic Properties of a Random Anisotropic Microstructure

Consider a random microstructure constituted of a random heterogeneous anisotropic elastic linear medium. The random local (or microscopic) constitutive equation is written as $\sigma(\mathbf{x}) = \mathbb{C}(\mathbf{x}) : \varepsilon(\mathbf{x})$ which means $\sigma_{jk}(\mathbf{x}) = \mathbb{C}_{jk\ell m}(\mathbf{x}) \varepsilon_{\ell m}(\mathbf{x})$, in which $\mathbf{x} \mapsto \sigma_{jk}(\mathbf{x})$ is the random local stress tensor field, $\varepsilon_{\ell m}(\mathbf{D}(\mathbf{x})) = \frac{1}{2}(\partial D_{\ell}(\mathbf{x})/\partial x_m + \partial D_m(\mathbf{x})/\partial x_{\ell})$ is the random local strain tensor field, $\mathbf{x} \mapsto \mathbf{D}(\mathbf{x}) = (D_1(\mathbf{x}), D_2(\mathbf{x}), D_3(\mathbf{x}))$ is the random local displacement field, $\mathbf{x} \mapsto \mathbb{C}(\mathbf{x})$ is the fourth-order tensor-valued random field allowing the elastic properties of the random microstructure to be characterized and where $\mathbf{x} = (x_1, x_2, x_3)$ is a point of the RVE which is a 3D bounded open domain Ω in \mathbb{R}^3 . The random effective (or macroscopic) stress and strain tensors are usually defined as the average in the RVE of the random local stress and strain tensor fields,

$$<\sigma> = \frac{1}{|\Omega|} \int_{\Omega} \sigma(\mathbf{x}) d\mathbf{x} , <\varepsilon> = \frac{1}{|\Omega|} \int_{\Omega} \varepsilon(\mathbf{x}) d\mathbf{x} .$$
 (1)

1.1. Localization and Random Effective Stiffness Tensor

The localization is done with a given random effective strain $\underline{\varepsilon}$ on the boundary $\partial\Omega$ of the RVE which is independent of \mathbf{x} . Consequently, the random local displacement field \mathbf{D} in the microstructure Ω can be constructed by solving the following stochastic boundary value problem (BVP) in Ω ,

$$-\operatorname{div} \mathbf{\sigma} = 0 \quad \text{in} \quad \Omega \quad , \quad \mathbf{D}(\mathbf{x}) = \underline{\boldsymbol{\varepsilon}} \mathbf{x} \quad \text{on} \quad \partial \Omega \quad . \tag{2}$$

Since the solution **D** of Eq. (2) depends linearly on $\underline{\varepsilon}$, the random local strain tensor can be written as

$$\varepsilon(\mathbf{D}(\mathbf{x})) = \mathbb{H}(\mathbf{x}) : \underline{\varepsilon} \quad , \tag{3}$$

in which the symmetric fourth-order tensor-valued random field $\mathbf{x} \mapsto \mathbb{H}(\mathbf{x})$ corresponds to the strain localization associated with the stochastic BVP defined by Eq. (2). In order to construct the random field \mathbb{H} , for all ℓ and m in $\{1,2,3\}$, the second-order tensors $g^{\ell m}$ are introduced such that

$$g_{jk}^{\ell m} = \frac{1}{2} (\delta_{j\ell} \delta_{km} + \delta_{jm} \delta_{k\ell}) \quad , \tag{4}$$

in which $\delta_{j\ell}$ is the Kronecker symbol. For all ℓ and m in $\{1,2,3\}$, let $\mathbf{D}^{\ell m}$ be the random local displacement field which is the solution of the following stochastic BVP in Ω ,

$$-\operatorname{div} \mathbf{\sigma}^{\ell m} = 0 \text{ in } \Omega \quad , \quad \mathbf{D}^{\ell m}(\mathbf{x}) = g^{\ell m} \mathbf{x} \text{ on } \partial \Omega \quad , \tag{5}$$

in which $\sigma^{\ell m}(\mathbf{x}) = \mathbb{C}(\mathbf{x}) : \varepsilon(\mathbf{D}^{\ell m}(\mathbf{x}))$. We then have

$$\mathbb{H}_{jk\ell m}(\mathbf{x}) = \varepsilon_{jk}(\mathbf{D}^{\ell m}(\mathbf{x})) \quad , \tag{6}$$

with the property $<\mathbb{H}_{jk\ell m}>=g_{jk}^{\ell m}$. The symmetric fourth-order random effective stiffness tensor $\mathbb{C}^{\mathrm{eff}}$ which is defined by $<\sigma>=\mathbb{C}^{\mathrm{eff}}:<\varepsilon>$ can then be calculated by

$$\mathbb{C}^{\text{eff}} = \langle \mathbb{C} : \mathbb{H} \rangle \quad . \tag{7}$$

1.2. Microscale Stochastic Model for Random Anisotropic Elastic Microstructures

For all \mathbf{x} fixed in Ω , the random fourth-order tensor $\mathbb{C}(\mathbf{x})$ has a given mean value, must verify the symmetry property and a stochastic nonuniform ellipticity condition related to positive-definiteness properties. The random field \mathbb{C} is constituted of 21 mutually dependent real-valued random fields and the system of marginal probability distributions of \mathbb{C} is required because the unknown random solution $\mathbf{x} \mapsto \mathbf{D}^{\ell m}(\mathbf{x})$ of the stochastic BVP defined by Eq. (5) is a nonlinear mapping of the random field \mathbb{C} .

The mean value of the random field $\mathbb C$ is a deterministic tensor-valued field $\mathbf x \mapsto \{\underline{\mathbb C}_{ijkh}(\mathbf x)\}_{ijkh}$. The microscale stochastic model has to be such that $E\{\mathbb C_{ijkh}(\mathbf x)\} = \underline{\mathbb C}_{ijkh}(\mathbf x)$ for all $\mathbf x$, where E is the mathematical expectation. The known symmetries can be taken into account with the mean model represented by the tensor $\{\underline{\mathbb C}_{ijkh}(\mathbf x)\}_{ijkh}$. This paper deals with the case for which the random fluctuation tensor $\{\mathbb C_{ijkh}(\mathbf x)\}_{ijkh}$ around the mean tensor is purely anisotropic, without any symmetries.

We present an extension of the probability model proposed in (Soize, 2004 and 2006) which only uses the available information. The usual uniform elliptic condition is not introduced to construct this probability model because such a condition does not correspond to available information (objective data). For the proposed stochastic model, a non uniform ellipticity condition is introduced for $\mathbf{x} \mapsto \mathbb{C}(\mathbf{x})$ which corresponds to the available information and which, for all ℓ and m fixed in $\{1,2,3\}$, allows the random weak formulation of the stochastic BVP defined by Eq. (5) to have a unique second-order random solution $\mathbf{x} \mapsto \mathbf{D}^{\ell m}(\mathbf{x})$.

In order to define the stochastic model of the tensor-valued random field \mathbb{C} , the (6×6) matrix representation $[\mathbf{A}(\mathbf{x})]$ of the fourth-order tensor $\mathbb{C}(\mathbf{x})$ is introduced. Therefore, let I and J be the new indices belonging to $\{1,\ldots,6\}$ such that I=(j,k) and $J=(\ell,m)$ with the following correspondence: 1=(1,1), 2=(2,2), 3=(3,3), 4=(1,2), 5=(1,3) and 6=(2,3). Thus, for all \mathbf{x} in Ω , the random (6×6) real matrix $[\mathbf{A}(\mathbf{x})]$ is such that

$$[\mathbf{A}(\mathbf{x})]_{IJ} = \mathbb{C}_{jk\ell m}(\mathbf{x}) \quad . \tag{8}$$

For all \mathbf{x} fixed in Ω , due to the symmetry and positive-definiteness properties of the random fourth-order tensor $\mathbb{C}(\mathbf{x})$, it can be deduced that $[\mathbf{A}(\mathbf{x})]$ is a random variable with values in the set $\mathbb{M}_6^+(\mathbb{R})$ of all the (6×6) real symmetric positive-definite matrices. The $\mathbb{M}_6^+(\mathbb{R})$ -valued random field $\{[\mathbf{A}(\mathbf{x})], \mathbf{x} \in \Omega\}$, indexed by Ω , defined on the probability space (Θ, \mathcal{T}, P) , is constituted of $6\times (6+1)/2=21$ mutually dependent real-valued random fields defining the fourth-order tensor-valued random field \mathbb{C} indexed by Ω .

The mean function $\mathbf{x} \mapsto [\underline{a}(\mathbf{x})]$ of the random field $[\mathbf{A}]$ is assumed to be a given function from Ω into $\mathbb{M}_{6}^{+}(\mathbb{R})$ such that, for all \mathbf{x} fixed in Ω ,

$$E\{[\mathbf{A}(\mathbf{x})]\} = [\underline{a}(\mathbf{x})] \quad . \tag{9}$$

Since $[\underline{a}(\mathbf{x})]$ belongs to $\mathbb{M}_6^+(\mathbb{R})$, there is an upper triangular invertible matrix $[\underline{L}(\mathbf{x})]$ in $\mathbb{M}_6(\mathbb{R})$ (the set of all the (6×6) real matrices) such that

$$[\underline{a}(\mathbf{x})] = [\underline{L}(\mathbf{x})]^T [\underline{L}(\mathbf{x})] . \tag{10}$$

It is assumed that $\mathbf{x} \mapsto [\underline{L}(\mathbf{x})]$ is bounded on Ω and that $\mathbf{x} \mapsto [\underline{a}(\mathbf{x})]$ satisfies the usual uniform ellipticity condition on Ω . For all \mathbf{x} fixed in Ω , the random matrix $[\mathbf{A}(\mathbf{x})]$ is written as

$$[\mathbf{A}(\mathbf{x})] = [\underline{L}(\mathbf{x})]^T [\mathbf{G}(\mathbf{x})] [\underline{L}(\mathbf{x})] , \qquad (11)$$

in which $\mathbf{x} \mapsto [\mathbf{G}(\mathbf{x})]$ is a random field defined on (Θ, \mathcal{T}, P) , indexed by \mathbb{R}^3 , with values in $\mathbb{M}_6^+(\mathbb{R})$, such that for all \mathbf{x} in \mathbb{R}^3 , $E\{[\mathbf{G}(\mathbf{x})]\} = [I]$ in which [I] is the identity matrix. The random field $[\mathbf{G}]$ is completely defined in Section 3.

2. Stochastic Model of the Random Fields [G] and [A]

Let $d \geq 1$ and $n \geq 1$ be two given integers. The random field $\mathbf{x} = (x_1, \dots, x_d) \mapsto [\mathbf{G}(\mathbf{x})]$ is indexed by \mathbb{R}^d with values in $\mathbb{M}_n^+(\mathbb{R})$. In Eq. (11), we have d=3 and n=6. As explained in Section 2.2, the extended probability model presented below is based on the construction and the mathematical analysis of the random field $[\mathbf{G}]$ performed in (soize, 2006). The results which allow the numerical calculation to be performed are summarized below. The random field $\mathbf{x} \mapsto [\mathbf{G}(\mathbf{x})]$ is constructed as a homogeneous and normalized non-Gaussian positive-definite matrix-valued random field, defined on probability space (Θ, \mathcal{T}, P) , indexed by \mathbb{R}^d , with values in $\mathbb{M}_n^+(\mathbb{R})$. This random field is constructed as a nonlinear mapping of stochastic germs.

2.1. Random Fields $U_{jj'}$ as the Stochastic Germs of the Random Field [G]

The stochastic germs are constituted of n(n+1)/2 independent second-order centered homogeneous Gaussian random fields $\mathbf{x}\mapsto U_{jj'}(\mathbf{x}),\ 1\leq j\leq j'\leq n$, defined on the probability space (Θ,\mathcal{T},P) , indexed by \mathbb{R}^d , with values in \mathbb{R} and such that $E\{U_{jj'}(\mathbf{x})\}=0$ and $E\{U_{jj'}(\mathbf{x})^2\}=1$. Consequently, all these random fields are completely and uniquely defined by the n(n+1)/2 autocorrelation functions $R_{U_{jj'}}(\eta)=E\{U_{jj'}(\mathbf{x}+\eta)U_{jj'}(\mathbf{x})\}$ defined for all $\mathbf{\eta}=(\eta_1,\ldots,\eta_d)$ in \mathbb{R}^d and such that $R_{U_{jj'}}(0)=1$. In order to obtain a class having a reasonable number of parameters, these autocorrelation functions are written as $R_{U_{jj'}}(\eta)=\rho_1^{jj'}(\eta_1)\times\ldots\times\rho_d^{jj'}(\eta_d)$ in which, for all $k=1,\ldots,d$, one has $\rho_k^{jj'}(0)=1$ and for all $\eta_k\neq 0$,

$$\rho_k^{jj'}(\eta_k) = 4(L_k^{jj'})^2 / (\pi^2 \eta_k^2) \sin^2 \left(\pi \eta_k / (2L_k^{jj'}) \right) , \qquad (12)$$

in which $L_1^{jj'},\ldots,L_d^{jj'}$ are positive real numbers. Each random field $U_{jj'}$ is then mean-square continuous on \mathbb{R}^d and it power spectral measure has a compact support. Such a model has $d\,n(n+1)/2$ real parameters $L_1^{jj'},\ldots,L_d^{jj'}$ for $1\leq j\leq j'\leq n$ which represent the spatial correlation lengths of the stochastic germs $U_{jj'}$.

2.2. Defining an Adapted Family of Functions

The construction of the random field $[\mathbf{G}]$ requires the introduction of an adapted family of functions $\{u\mapsto h(\alpha,u)\}_{\alpha>0}$. Let α be a positive real number. The function $u\mapsto h(\alpha,u)$ from $\mathbb R$ into $]0\,,+\infty[$ is introduced such that $\Gamma_\alpha=h(\alpha,U)$ is a gamma random variable with parameter α while U is a normalized Gaussian random variable ($E\{U\}=0$ and $E\{U^2\}=1$). Consequently, for all u in $\mathbb R$, we have

$$h(\alpha, u) = F_{\Gamma_{\alpha}}^{-1}(F_U(u))$$
 (13)

in which $u\mapsto F_U(u)=\int_{-\infty}^u\frac{1}{\sqrt{2\pi}}e^{-v}\,dv$ is the cumulative distribution function of the normalized Gaussian random variable U. The function $p\mapsto F_{\Gamma_\alpha}^{-1}(p)$ from $]0\,,1[$ into $]0\,,+\infty[$ is the reciprocal function of the cumulative distribution function $\gamma\mapsto F_{\Gamma_\alpha}(\gamma)=\int_0^\gamma\frac{1}{\Gamma(\alpha)}\,t^{\alpha-1}\,e^{-t}\,dt$ of the gamma random variable Γ_α with parameter α in which $\Gamma(\alpha)$ is the gamma function defined by $\Gamma(\alpha)=\int_0^{+\infty}t^{\alpha-1}\,e^{-t}\,dt$.

2.3. Defining the Random Field [G]

The random field $\mathbf{x} \mapsto [\mathbf{G}(\mathbf{x})]$, defined on the probability space (Θ, \mathcal{T}, P) , indexed by \mathbb{R}^d , with values in $\mathbb{M}_n^+(\mathbb{R})$ is constructed as follows:

(i) Let $\{U_{jj'}(\mathbf{x}), \mathbf{x} \in \mathbb{R}^d\}_{1 \leq j \leq j' \leq n}$ be the n(n+1)/2 independent random fields introduced in Section 2.1. For all \mathbf{x} in \mathbb{R}^d and for all i and j such that $1 \leq j \leq j' \leq n$, we have

$$E\{U_{ij'}(\mathbf{x})\} = 0$$
 , $E\{U_{ij'}(\mathbf{x})^2\} = 1$. (14)

- (ii) Let δ be the real number, independent of \mathbf{x} and n, such that $0 < \delta < \sqrt{(n+1)(n+5)^{-1}} < 1$. This parameter which is assumed to be known (resulting, for instance, from an experimental identification solving an inverse problem) allows the dispersion of the random field $[\mathbf{G}]$ to be controlled.
- (iii) For all \mathbf{x} in \mathbb{R}^d , the random matrix $[\mathbf{G}(\mathbf{x})]$ is written

$$[\mathbf{G}(\mathbf{x})] = [\mathbf{L}(\mathbf{x})]^T [\mathbf{L}(\mathbf{x})] , \qquad (15)$$

in which $[\mathbf{L}(\mathbf{x})]$ is the upper $(n \times n)$ real triangular random matrix defined as follows:

For $1 \le j \le j' \le n$, the n(n+1)/2 random fields $\mathbf{x} \mapsto [\mathbf{L}(\mathbf{x})]_{jj'}$ are independent.

For j < j', the real-valued random field $\mathbf{x} \mapsto [\mathbf{L}(\mathbf{x})]_{jj'}$, indexed by \mathbb{R}^d , is defined by $[\mathbf{L}(\mathbf{x})]_{jj'} = \sigma U_{jj'}(\mathbf{x})$ in which σ is such that $\sigma = \delta (n+1)^{-1/2}$.

For j=j', the positive-valued random field $\mathbf{x}\mapsto [\mathbf{L}(\mathbf{x})]_{jj}$, indexed by \mathbb{R}^d , is defined by $[\mathbf{L}(\mathbf{x})]_{jj} = \sigma$ $\sqrt{2\,h(\alpha_j,U_{jj}(\mathbf{x}))}$ in which $\alpha_j=(n+1)/(2\delta^2)+(1-j)/2$.

Let $[\mathbf{B}]$ be a random matrix, defined on (Θ, \mathcal{T}, P) , with values in the set $\mathbb{M}_m(\mathbb{R})$ of all the $(m \times m)$ real matrices. For $\theta \in \Theta$, let $[\mathbf{B}(\theta)] \in \mathbb{M}_m(\mathbb{R})$ be a realization of $[\mathbf{B}]$. The norm $\|\mathbf{B}(\theta)\|$ of $[\mathbf{B}(\theta)]$ induced by the Euclidean norm $\|\mathbf{v}\|$ of \mathbf{v} in \mathbb{R}^m is such that

$$\|\mathbf{B}(\theta)\| = \sup_{\|\mathbf{v}\| \le 1} \|[\mathbf{B}(\theta)]\mathbf{v}\| \quad , \quad \mathbf{v} \in \mathbb{R}^m \quad . \tag{16}$$

It can be proven the following fundamental property (non uniform ellipticity condition replacing the usual uniform ellipticity condition which is not introduced): let Ω be any bounded open domain of \mathbb{R}^d and let $\overline{\Omega} = \Omega \cup \partial \Omega$ be its closure. We then have

$$E\left\{\left(\sup_{\mathbf{x}\in\overline{\Omega}}\|\left[\mathbf{G}(\mathbf{x})\right]^{-1}\|\right)^{2}\right\} = c_{G}^{2} < +\infty , \qquad (17)$$

in which sup is the supremum and where $0 < c_G < +\infty$ is a finite positive constant. Note that the mathematical proof of Eq. (17) can easily be derived from (Soize, 2006) for the extended class introduced in Section 3.1.

2.4. Basic Properties of the Rrandom Field [A] and its Parameters

The random field $\mathbf{x} \mapsto [\mathbf{A}(\mathbf{x})]$ is a second-order random field on Ω : $E\{\|\mathbf{A}(\mathbf{x})\|^2\} \le E\{\|\mathbf{A}(\mathbf{x})\|_F^2\} < +\infty$ in which $\|\cdot\|_F$ is the Frobenius norm. The system of the marginal probability distributions of the random field $\mathbf{x} \mapsto [\mathbf{A}(\mathbf{x})]$ is completely defined, is not Gaussian and is deduced from the system of the marginal probability distributions of the random field $\mathbf{x} \mapsto [\mathbf{G}(\mathbf{x})]$ by using Eq. (11). In general, since $[\underline{a}(\mathbf{x})]$ depends on \mathbf{x} , then the random field $\{[\mathbf{A}(\mathbf{x})], \mathbf{x} \in \Omega\}$ is non homogeneous. For all fixed \mathbf{x} , the dispersion parameter $\delta_A(\mathbf{x})$ of the random matrix $[\mathbf{A}(\mathbf{x})]$, which is defined by

$$\delta_A(\mathbf{x})^2 = E\{\|[\mathbf{A}(\mathbf{x})] - [\underline{a}(\mathbf{x})]\|_F^2\} / \|\underline{a}(\mathbf{x})\|_F^2 \quad , \tag{18}$$

is such that

$$\delta_A(\mathbf{x}) = \frac{\delta}{\sqrt{n+1}} \left\{ 1 + \frac{\left(\text{tr}\left[\underline{a}(\mathbf{x})\right]\right)^2}{\text{tr}\left\{\left[\underline{a}(\mathbf{x})\right]^2\right\}} \right\}^{1/2} . \tag{19}$$

The random field $\mathbf{x}\mapsto [\mathbf{G}(\mathbf{x})]$ almost surely has continuous trajectories. If the function $\mathbf{x}\mapsto [\underline{a}(\mathbf{x})]$ is continuous on $\overline{\Omega}$, then the random field $\mathbf{x}\mapsto [\mathbf{A}(\mathbf{x})]$ almost surely has continuous trajectories on $\overline{\Omega}$. Nevertheless, if the function $\mathbf{x}\mapsto [\underline{a}(\mathbf{x})]$ is not continuous on $\overline{\Omega}$, then the random field $\mathbf{x}\mapsto [\mathbf{A}(\mathbf{x})]$ almost surely does not have continuous trajectories on $\overline{\Omega}$. Then the random field $\mathbf{x}\mapsto [\mathbf{G}(\mathbf{x})]$ is completely and uniquely defined by the following parameters: the $\mathbb{M}_6^+(\mathbb{R})$ -valued mean function $\mathbf{x}\mapsto [\underline{a}(\mathbf{x})]$, the positive real parameter δ and the 63 positive real parameters $L_1^{jj'}, L_2^{jj'}, L_3^{jj'}$ for $1\le j\le j'\le \delta$. The smallest number of parameters corresponds to the following case: $\mathbf{x}\mapsto [\underline{a}(\mathbf{x})]$, δ and $L_d=L_1^{jj'}=L_2^{jj'}=L_3^{jj'}$ for all $1\le j\le j'\le \delta$.

2.5. Spatial Correlation Lengths of the Random Field [A] for the Homogeneous Case

If $[\underline{a}(\mathbf{x})] = [\underline{a}]$ is independent of \mathbf{x} , then the random field $\{[\mathbf{A}(\mathbf{x})] = [\underline{L}]^T [\mathbf{G}(\mathbf{x})] [\underline{L}], \mathbf{x} \in \Omega\}$ can be viewed as the restriction to Ω of a homogeneous random field indexed by \mathbb{R}^3 . Then the dispersion parameter defined by Eq. (18) is independent of \mathbf{x} and then $\delta_A(\mathbf{x}) = \delta_A$. Let $\mathbf{\eta} = (\eta_1, \eta_2, \eta_3) \mapsto r^A(\mathbf{\eta})$ be the function defined from \mathbb{R}^3 into \mathbb{R} by

$$r^{A}(\mathbf{\eta}) = \frac{\operatorname{tr} E\{([\mathbf{A}(\mathbf{x} + \mathbf{\eta})] - [\underline{a}]) ([\mathbf{A}(\mathbf{x})] - [\underline{a}])\}}{E\{\|[\mathbf{A}(\mathbf{x})] - [\underline{a}]\|_{F}^{2}\}} \quad . \tag{20}$$

It can be seen that $r^A(0)=1$ and $r^A(-\eta)=r^A(\eta)$. For k=1,2,3, the spatial correlation length L_k^A of $\mathbf{x}\mapsto [\mathbf{A}(\mathbf{x})]$ and relative to the coordinate x_k can then be defined by

$$L_k^A = \int_0^{+\infty} |r^A(\mathbf{\eta}^k)| d\eta_k \quad , \tag{21}$$

in which $\eta^1 = (\eta_1, 0, 0), \eta^2 = (0, \eta_2, 0)$ and $\eta^3 = (0, 0, \eta_3)$.

3. Discretization with Stochastic Finite Elements and Solving the Random Equation

For all ℓ and m in $\{1,2,3\}$, we have to (1) solve the stochastic BVP defined by Eq. (5) with $\sigma^{\ell m}(\mathbf{x}) = \mathbb{C}(\mathbf{x}) : \varepsilon(\mathbf{D}^{\ell m}(\mathbf{x}))$ in order to construct the random local displacement field $\mathbf{D}^{\ell m}$; (2) calculate the fourth-order tensor-valued random field \mathbb{H} defined by Eq. (6) and corresponding to the strain localization; (3) estimate the probabilistic properties of the fourth-order random effective stiffness tensor \mathbb{C}^{eff} defined by Eq. (7).

In order to solve this problem, the following computational stochastic method is used.

The weak formulation of the stochastic BVP is constructed and the existence of a unique second-order stochastic solution can be proven. Note that the usual proof of the existence of a solution can be directly deduced if a uniform ellipticity condition and a uniform boundness condition for the fourth-order tensor-valued random field $\mathbf{x} \mapsto \mathbb{C}(\mathbf{x})$ is used. This usual proof cannot be used for the present case due to the non introduction of these uniformness conditions which are substituted by the more realistic probabilistic hypothesis which has been introduced and which corresponds to the introduction of the non uniformness condition. The adapted proof can be obtained thanks to Eq. (17).

The stochastic finite element method is used for discretizing the weak formulation and Eq. (5). For such a numerical approximation, the random local stiffness tensor field $\mathbf{x} \mapsto \mathbb{C}_{jk\ell m}(\mathbf{x})$ and the random local strain tensor field $\mathbf{x} \mapsto \varepsilon_{jk}(\mathbf{D}^{\ell m}(\mathbf{x}))$ are discretized at all the Gauss-Legendre quadrature points of the finite elements.

The probabilistic quantities are then estimated by using the Monte Carlo simulation method which is made up of 3 main steps: (a) developing a generator for constructing n_s independent realizations $\{\mathbf{x} \mapsto \mathbb{C}(\mathbf{x}, \theta_r), r=1, \dots n_s\}$ of the random field $\mathbf{x} \mapsto \mathbb{C}(\mathbf{x})$ using the probability model presented in Section 3; (b) for each realization $\mathbf{x} \mapsto \mathbb{C}(\mathbf{x}, \theta_r)$, calculating the corresponding realization $\mathbb{C}^{\text{eff}}(\theta_r)$ of the effective tensor (and related quantities such as the random eigenvalues of the random tensor \mathbb{C}^{eff}) by solving a deterministic matrix equation; (c) with the n_s independent realizations, estimating the probabilistic quantities (moments, probability distributions) using the mathematical statistics and studying the convergence with respect to n_s .

3.1. Finite Element Discretization

Let $\mathbb{V}=L^2(\Theta,V)$ be the real Hilbert space of all the second-order random variables $\theta\mapsto\{\mathbf{x}\mapsto\mathbf{D}(\mathbf{x},\theta)\}$ defined on probability space (Θ,\mathcal{T},P) , with values in the Hilbert space $V=(H^1(\Omega))^3$. Let $\mathbf{e}(\mathbf{d})$ be the strain vector defined by $\mathbf{e}(\mathbf{d})=(\varepsilon_{11}(\mathbf{d}),\varepsilon_{22}(\mathbf{d}),\varepsilon_{33}(\mathbf{d}),2\varepsilon_{12}(\mathbf{d}),2\varepsilon_{13}(\mathbf{d}),2\varepsilon_{23}(\mathbf{d}))$ and let $(\mathbf{D},\delta\mathbf{D})\mapsto\mathbf{K}(\mathbf{D},\delta\mathbf{D})$ be the random bilinear form on $\mathbb{V}\times\mathbb{V}$ defined by

$$\mathbf{K}(\mathbf{D}, \delta \mathbf{D}) = \int_{\Omega} \langle [\mathbf{A}(\mathbf{x})] \, \mathbf{e}(\mathbf{D}(\mathbf{x})) \,, \, \mathbf{e}(\delta \mathbf{D}(\mathbf{x})) \rangle \, d\mathbf{x} \quad . \tag{22}$$

The stochastic finite element method is used to discretize the weak formulation. A finite element mesh of domain Ω is carried out using 3D solid finite elements. We then have

 $\Omega = \bigcup_e \Omega_e$ in which Ω_e is the domain of the finite element number e. Any displacement field $\mathbf{x} \mapsto \mathbf{d}(\mathbf{x})$ in V and its associated strain vector field $\mathbf{x} \mapsto \mathbf{e}(\mathbf{x})$ are then approximated by

$$\mathbf{d}(\mathbf{x}) \simeq [B(\mathbf{x})]\mathbf{w} \ , \ \mathbf{x} \in \overline{\Omega} \quad , \quad \mathbf{e}(\mathbf{x}) \simeq [S(\mathbf{x})]\mathbf{w} \ , \ \mathbf{x} \in \Omega \quad ,$$

in which $\mathbf{w}=(w_1,\ldots,w_{\nu})$ is the vector of the ν degrees of freedom corresponding to the values of the components of the field \mathbf{d} at the nodes of the mesh. The $(3\times\nu)$ real matrices $[B(\mathbf{x})]$ and $[S(\mathbf{x})]$ are known matrices usually constructed by using the interpolation functions of the finite elements. For any integrable function $\mathbf{x}\mapsto f(\mathbf{x})$ defined on Ω and continuous on Ω_e , the following usual numerical approximation can be written as $\int_{\Omega_e} f(\mathbf{x}) \, d\mathbf{x} \simeq \sum_{k=1}^{N_e} \omega_{\alpha_k} \, f(\mathbf{x}^{\alpha_k})$, in which $\{\omega_{\alpha_1},\ldots,\omega_{\alpha_{N_e}}\}$ and $\{\mathbf{x}^{\alpha_1},\ldots,\mathbf{x}^{\alpha_{N_e}}\}$ are the sets of all the N_e weights and the N_e Gauss-Legendre quadrature points for the finite element Ω_e . Consequently, it can be written that $\int_{\Omega} f(\mathbf{x}) \, d\mathbf{x} = \sum_e \int_{\Omega_e} f(\mathbf{x}) \, d\mathbf{x} \simeq \sum_{\alpha=1}^{N} \omega_{\alpha} \, f(\mathbf{x}^{\alpha})$, in which $\{\omega_1,\ldots,\omega_N\}$ and $\{\mathbf{x}^1,\ldots,\mathbf{x}^N\}$ are the sets of all the N weights and the N Gauss-Legendre quadrature points for $\Omega = \bigcup_e \Omega_e$ with $N = \sum_e N_e$.

For all \mathbf{x} in $\overline{\Omega}$, the finite element approximation of all \mathbf{D} and $\delta \mathbf{D}$ in \mathbb{V} is then written as $\mathbf{D}(\mathbf{x}) \simeq [B(\mathbf{x})] \mathbf{W}$ and $\delta \mathbf{D}(\mathbf{x}) \simeq [B(\mathbf{x})] \delta \mathbf{W}$, in which \mathbf{W} and $\delta \mathbf{W}$ are \mathbb{R}^{ν} -valued second-order random vectors. Therefore, the corresponding finite element approximation of the random bilinear form defined by Eq. (22) is such that $\langle [\mathbf{K}] \mathbf{W}, \delta \mathbf{W} \rangle = \mathbf{K}([B(\,.\,)] \mathbf{W}, [B(\,.\,)] \delta \mathbf{W})$, defining the $(\nu \times \nu)$ random stiffness matrix $[\mathbf{K}]$ such that

$$[\mathbf{K}] = \int_{\Omega} [S(\mathbf{x})]^T [\mathbf{A}(\mathbf{x})] [S(\mathbf{x})] . \tag{23}$$

The random stiffness matrix $[\mathbf{K}]$ defined by Eq. (23) can then be approximated by the random matrix $[\mathbf{K}_N]$ such that

$$[\mathbf{K}_N] = \sum_{\alpha=1}^N \omega_{\alpha} \left[S(\mathbf{x}^{\alpha}) \right]^T \left[\mathbf{A}(\mathbf{x}^{\alpha}) \right] \left[S(\mathbf{x}^{\alpha}) \right] d\mathbf{x} \quad , \tag{24}$$

in which the $\mathbb{M}_6^+(\mathbb{R})$ -valued random field $\mathbf{x} \mapsto [\mathbf{A}(\mathbf{x})]$ is discretized in the N Gauss-Legendre quadrature points $\mathbf{x}^1,\dots,\mathbf{x}^N$. The integral over Ω in Eq. (24) must be read as a mean-square integral and it can be proven that the convergence is reached when N goes to infinity. For a given fixed N, the random matrix $[\mathbf{K}_N]$ will be a good approximation of the random matrix $[\mathbf{K}]$ in the mean-square sense if the number N of the Gauss-Legendre quadrature points $\mathbf{x}^1,\dots,\mathbf{x}^N$ is sufficiently large and is adapted to the variations of the intercorrelation functions of the real-valued random fields $[\mathbf{A}]_{k\ell}$ and $[\mathbf{A}]_{k'\ell'}$.

functions of the real-valued random fields $[{\bf A}]_{k\ell}$ and $[{\bf A}]_{k'\ell'}$. The finite element discretization of the random field ${\bf D}^{\ell m}$ is then written as ${\bf D}^{\ell m}({\bf x}) \simeq [B({\bf x})] {\bf W}^{\ell m}$, ${\bf x} \in \overline{\Omega}$. The second-order random vector ${\bf W}^{\ell m}$ with values in \mathbb{R}^{ν} can then be written as ${\bf W}^{\ell m} = ({\bf W}_i^{\ell m}, {\bf w}_b^{\ell m})$ in which ${\bf W}_i^{\ell m}$ is the \mathbb{R}^{ν_i} -valued second-order random vector of the ν_i degrees of freedom for the nodes inside the domain Ω and where ${\bf w}_b^{\ell m}$ is the \mathbb{R}^{ν_b} deterministic vector of the ν_b degrees of freedom for the nodes belonging to the boundary $\partial \Omega$. This last vector is such that ${\bf d}^{\ell m}({\bf x}) \simeq [B({\bf x})] {\bf w}_b^{\ell m}$ for all ${\bf x}$ in $\partial \Omega$ and the vector ${\bf w}_b^{\ell m}$ is constituted of the values of the components of the field ${\bf d}^{\ell m}$ at the nodes belonging to boundary $\partial \Omega$. The block writing of the random stiffness matrix $[{\bf K}]$ defined by Eq. (23) is

$$[\mathbf{K}] = \begin{bmatrix} [\mathbf{K}_{ii}] & [\mathbf{K}_{ib}] \\ [\mathbf{K}_{ib}]^T & [\mathbf{K}_{bb}] \end{bmatrix} .$$

The finite element approximation of Eq. (5) is then given by the following random matrix equation allowing the unknown random vector $\mathbf{W}_i^{\ell m}$ to be calculated,

$$\left[\mathbf{K}_{ii}\right]\mathbf{W}_{i}^{\ell m} = -\left[\mathbf{K}_{ib}\right]\mathbf{w}_{b}^{\ell m} \quad . \tag{25}$$

The random values $\mathbb{H}(\mathbf{x}^1),\dots,\mathbb{H}(\mathbf{x}^N)$ of \mathbb{H} are calculated by using Eq. (6). Let $\mathbb{C}(\mathbf{x}^1),\dots,\mathbb{C}(\mathbf{x}^N)$ be the random values of \mathbb{C} . The finite element approximation of the fourth-order random effective stiffness tensor $\mathbb{C}^{\mathrm{eff}}$ defined by Eq. (7) can then be written as

$$\mathbb{C}^{\text{eff}} \simeq \frac{1}{|\Omega|} \sum_{\alpha=1}^{N} \omega_{\alpha} \, \mathbb{C}(\mathbf{x}^{\alpha}) : \mathbb{H}(\mathbf{x}^{\alpha}) \quad . \tag{26}$$

The $\mathbb{M}_6^+(\mathbb{R})$ -valued random matrix $[\mathbf{A}^{\mathrm{eff}}]$ of the fourth-order random effective stiffness tensor $\mathbb{C}^{\mathrm{eff}}$ is such that

$$[\mathbf{A}^{\text{eff}}]_{IJ} = \mathbb{C}^{\text{eff}}_{jk\ell m} \quad . \tag{27}$$

3.2. Defining the Probabilistic Quantities for the Random Effective Stiffness Matrix

It is assumed that the mean model of the microstructure is homogeneous, that is to say $[\underline{a}(\mathbf{x})] = [\underline{a}]$ is independent of \mathbf{x} . Consequently, the random field $[\mathbf{A}]$ describing the fourth-order tensor-valued random field $\mathbb C$ can be viewed as the restriction to Ω of a homogeneous random field indexed by $\mathbb R^3$. Using Eq. (19), the dispersion parameter δ_A which is then independent of \mathbf{x} can easily be deduced from the value of the dispersion parameter δ . The spatial correlation lengths L_1^A , L_2^A and L_3^A defined by Eq. (21) depend on the values of the parameters $L_1^{jj'}$, $L_2^{jj'}$, $L_3^{jj'}$ for $1 \leq j \leq j' \leq 6$ of the stochastic germs. Therefore, every probabilistic analysis of the macroscopic properties of the random anisotropic elastic microstructure is performed for a given value of the parameters $[\underline{a}]$, δ and $L_1^{jj'}$, $L_2^{jj'}$, $L_3^{jj'}$ for $1 \leq j \leq j' \leq 6$. Let $z \mapsto p_Z(z)$ be the probability density function with respect to dz of the random variable Z defined by

$$Z = \|\mathbf{A}^{\text{eff}}\|/E\{\|\mathbf{A}^{\text{eff}}\|\} \quad . \tag{28}$$

The support of p_Z is \mathbb{R}^+ . The cumulative distribution function $z^* \mapsto F_Z(z^*)$ of the random variable Z is then defined by

$$F_Z(z^*) = \int_0^{z^*} p_Z(z) \, dz = E\{\mathbb{1}_{Z \le z^*}\} \quad , \tag{29}$$

in which $\mathbb{1}_{Z \leq z^*} = 1$ if $Z \leq z^*$ and = 0 if not.

3.3. Solving the Random Equation and Computing the Statistical Estimations

The stochastic problem defined by Eqs. (25) to (28) is solved by using the Monte Carlo numerical simulation method with n_s independent realizations $\theta_1,\ldots,\theta_{n_s}$ in Θ . The independent stochastic germs $U_{jj'}$ for $1\leq j\leq j'\leq 6$ are simulated in the N Gauss-Legendre quadrature points $\mathbf{x}^1,\ldots,\mathbf{x}^N$. The mathematical expectation of any random quantity R is estimated by $E\{R\}=1/n_s\sum_{r=1}^{n_s}R(\theta_r)$. The cumulative probability distribution F_Z defined by Eq. (28) is estimated with the usual estimator (Serfling, 1980). Let $[\mathbf{A}_{\nu}^{\mathrm{eff}}]$ be the random effective stiffness matrix calculated with the finite element model having ν degrees of freedom. For a

given value of ν , the convergence of the Monte Carlo numerical simulation with respect to n_s can be controlled by studying the function

$$n_s \mapsto \operatorname{conv}(n_s, \nu) = \frac{1}{\|\underline{\mathbf{A}}_{\nu}^{\text{eff}}\|} \left(\frac{1}{n_s} \sum_{r=1}^{n_s} \|\mathbf{A}_{\nu}^{\text{eff}}(\theta_r)\|^2 \right)^{1/2}$$
(30)

in which $\|\mathbf{A}_{\nu}^{\mathrm{eff}}(\theta_r)\|$ is the realization θ_r of the random variable $\|\mathbf{A}_{\nu}^{\mathrm{eff}}\|$. If the mean model of the microstructure is homogeneous, then the local stiffness matrix $[\underline{a}]$ is independent of \mathbf{x} and then $[\underline{\mathbf{A}}_{\nu}^{\mathrm{eff}}] = [\underline{a}]$. The right-hand side of Eq. (30) corresponds to the statistical estimation of the norm $|||\mathbf{A}_{\nu}^{\mathrm{eff}}||| = (E\{\|\mathbf{A}_{\nu}^{\mathrm{eff}}\|^2\})^{1/2}$ of the random matrix $[\mathbf{A}_{\nu}^{\mathrm{eff}}]$. For a given value of n_s , the convergence with respect to the number ν of degrees of freedom of the finite element model is given by the function $\nu \mapsto \mathrm{conv}(n_s, \nu)$.

4. Application to an Anisotropic Random Microstructure

4.1. Mean Model and Finite Element Discretization of the Mean Model of the Microstructure

The open bounded domain Ω (the RVE) of \mathbb{R}^3 is such that $\Omega=(]0,1[)^3$. The mean model of the microstructure corresponds to a homogeneous anisotropic linear elastic medium whose local stiffness matrix $[\underline{a}]$ belonging to $\mathbb{M}_n^+(\mathbb{R})$ with n=6 is then independent of \mathbf{x} and such that $[\underline{a}]=10^{10}$

$$\times \begin{bmatrix} 3.3617 & 1.7027 & 1.3637 & -0.1049 & -0.2278 & 2.1013 \\ 1.7027 & 1.6092 & 0.7262 & 0.0437 & -0.1197 & 0.8612 \\ 1.3637 & 0.7262 & 1.4653 & -0.1174 & -0.1506 & 1.0587 \\ -0.1049 & 0.0437 & -0.1174 & 0.1319 & 0.0093 & -0.1574 \\ -0.2278 & -0.1197 & -0.1506 & 0.0093 & 0.1530 & -0.1303 \\ 2.1013 & 0.8612 & 1.0587 & -0.1574 & -0.1303 & 1.7446 \end{bmatrix}$$

The finite element model is a regular mesh of $12 \times 12 \times 12 = 1728$ nodes and $11 \times 11 \times 11 = 1331$ finite elements which are 8-nodes solid elements with $2 \times 2 \times 2$ Gauss-Legendre quadrature points. Therefore there are 5184 degrees of freedom, N=10648 Gauss-Legendre quadrature points, $\nu_i=3000$ degrees of freedom for the nodes inside domain Ω and $\nu_b=2184$ degrees of freedom for the nodes belonging to boundary $\partial\Omega$. In this case, the weights ω_α are such that $\omega_\alpha=|\Omega|/N$ and consequently, Eq. (26) can be rewritten as

$$\mathbb{C}^{\mathrm{eff}} \simeq \frac{1}{N} \sum_{\alpha=1}^{N} \mathbb{C}(\mathbf{x}^{\alpha}) : \mathbb{H}(\mathbf{x}^{\alpha}) .$$

4.2. Microscale Stochastic Model, Computational Parameters and Stochastic Response

(i) Microscale stochastic model. The stochastic model of the elasticity tensor of the random anisotropic microstructure is defined in Sections 2.2 and 3. The random field $\mathbf{x} \mapsto [\mathbf{A}(\mathbf{x})]$, indexed by Ω , with values in $\mathbb{M}_n^+(\mathbb{R})$, with n=6, is such that (see Eq. (11)), $[\mathbf{A}(\mathbf{x})] = [\underline{L}]^T[\mathbf{G}(\mathbf{x})][\underline{L}]$ in which the matrix $[\underline{a}]$, defined in Section 5.1 is written (see Eq. (10)) as $[\underline{a}] = [\underline{L}]^T[\underline{L}]$. The stochastic field $\mathbf{x} \mapsto [\mathbf{G}(\mathbf{x})]$, indexed by \mathbb{R}^3 , with values in $\mathbb{M}_n^+(\mathbb{R})$, is defined in Section 3.

(ii) Dispersion parameter. The dispersion parameter δ_A defined by Eq. (18) is then independent of **x** and can be calculated as a function of the parameter δ by using Eq. (19) and yields

 $\delta_A=0.6192 \times \delta$. For $\delta=0.1, 0.2, 0.3$ and 0.4, one then has $\delta_A=0.0619, 0.1238, 0.1858$ and 0.2477 respectively. Below, all the results are given as function of δ instead of δ_A .

(iii) Parameters of the stochastic germs and spatial correlation lengths of the random field [A]. In is assumed that the parameters $L_1^{jj'}, L_2^{jj'}, L_3^{jj'}$ for $1 \leq j \leq j' \leq 6$ of the stochastic germs are such that $L_1^{jj'} = L_2^{jj'} = L_3^{jj'} = L_d$ for all j and j' in which L_d is the unique parameter relative to the length-scales of the stochastic germs. For each given value of δ and for each given value of L_d , the independent realizations of the random matrices $[\mathbf{A}(\mathbf{x}^1)], \ldots, [\mathbf{A}(\mathbf{x}^N)]$ at the N=10648 points $\mathbf{x}^1, \ldots, \mathbf{x}^N$ are calculated. Since $L_1^{jj'} = L_2^{jj'} = L_3^{jj'} = L_d$ for all j and j', it can easily be deduced that the correlation function defined by Eq. (20) is such that $r^A(\eta,0,0) = r^A(0,\eta,0) = r^A(0,\eta,0)$. The following notation $r_d^A(\eta) = r^A(\eta,0,0) = r^A(0,\eta,0) = r^A(0,0,\eta)$ is then used below. It can be proven that the function r_d^A is independent of δ . From Eq. (21), it can then be seen that the spatial correlation lengths L_1^A , L_2^A and L_3^A of the random field $\mathbf{x} \mapsto [\mathbf{A}(\mathbf{x})]$ are then equal to a same value denoted by L_A .

For $L_d=0.1$, Figure 1 displays the graph of the function $\eta\mapsto r_d^A(\eta)$ calculated with Eq. (20) in which the mathematical expectation is estimated by using the Monte Carlo simulation with 2000 independent realizations. For these values, the spatial correlation length L_A is calculated by Eq. (21) and is such that $L_A=0.1113$ for $L_d=0.1$ and $L_A=1.113$ for $L_d=1$. More generally, one has $L_A=1.113$ L_d . Due to this correspondence between L_A and L_d , it is equivalent to present the results in terms of the spatial correlation length L_A or the parameter L_d . Below, the results are presented in function of L_d .

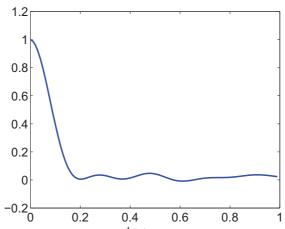


Figure 1. Graph of the correlation function $\eta \mapsto r_d^A(\eta)$ for $L_d = 0.1$. Horizontal axis η . Vertical axis $r_d^A(\eta)$

Let \mathcal{C}_d be the cube $L_A \times L_A \times L_A$. For the smallest value of L_A considered in the numerical calculation ($L_A=0.1113$ corresponding to $L_d=0.1$) there are about 14 Gauss-Legendre quadrature points in \mathcal{C}_d . For $L_A=0.2226$ corresponding to $L_d=0.2$, there are about 112 points. This is sufficient to obtain a good approximation of the random matrix $[\mathbf{K}]$ by $[\mathbf{K}_N]$ in the mean-square sense taking into account the very slow variations (over any interval of length L_A) of the correlation function displays in Figure 1. In addition, Figure 2 shows that the convergence is reached with respect to N for this smallest value of L_A and for the strongest stochastic fluctuations considered ($\delta=0.4$). This is an additional important information to conclude that the convergence is reached in the mean-square sense with a resonable accuracy.

(iv) Stochastic convergence analysis for the random effective stiffness matrix. For each given value of the dispersion parameter δ , for each given value of the parameter L_d and for a given

finite element model having ν degrees of freedom, the probabilistic quantities for the random effective stiffness matrix are estimated by using the Monte Carlo numerical simulation.

The mean-square convergence with respect to n_s has been studied by constructing the function $n_s\mapsto \text{conv}(n_s,\nu)$ defined by Eq. (30). For $\delta=0.4$ (the largest value considered for the dispersion parameter), for $L_d=0.1$ (the smallest value considered for this parameter) and for a finite element model having $\nu=5184$ degrees of freedom and corresponding to $11\times11\times11$ finite elements, convergence is reached for $n_s\geq 500$.

The mean-square convergence with respect to the number ν of degrees of freedom of the finite element model is studied by constructing the function $\nu \mapsto \mathrm{conv}(n_s,\nu)$ defined by Eq. (30). Figure 2 displays the graph of $\nu \mapsto \mathrm{conv}(n_s,\nu)$ for $\delta=0.4$, $L_d=0.1$ and $n_s=900$. Clearly, convergence is reached for $11\times11\times11$ finite elements corresponding to $\nu=5184$.

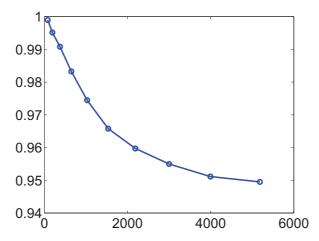


Figure 2. Mean-square convergence of the random effective stiffness matrix with respect to the number ν of degrees of freedom of the finite element model for $\delta=0.4$, $L_d=0.1$ and for $n_s=900$. Graph of function $\nu\mapsto {\rm conv}(n_s,\nu)$. Horizontal axis ν . Vertical axis ${\rm conv}(n_s,\nu)$.

All the results presented below have been computed with $n_s=900$ and $11\times11\times11$ finite elements corresponding to $\nu=5184$. Convergence is reached for the values of L_d and δ which are considered below.

4.3. Probabilistic Analysis of the RVE Size

Let β be a positive real number. Let $\beta \mapsto \mathbb{P}(\beta)$ be the function defined by

$$\mathbb{P}(\beta) = P\{1 - \beta < Z \le 1 + \beta\} = F_Z(1 + \beta) - F_Z(1 - \beta).$$

For $\delta=0.4$ and for L_d belonging to [0.1,0.7], Figure 3 displays the graph of the function \mathbb{P} . For instance, for $L_d=0.2$ and for $\beta=0.02,0.04$ and 0.08, one has $\mathbb{P}=0.36,0.65$ and 0.95, which means that $P\{0.98 < Z \leq 1.02\} = 0.36, P\{0.96 < Z \leq 1.04\} = 0.65$ and $P\{0.92 < Z \leq 1.08\} = 0.95$. For instance, for $\delta=0.4$ (i.e. $\delta_A=0.2477$), if stochastic homogenization is performed with a RVE whose size is five times the spatial correlation length (REV size = 1 and spatial correlation length $L_A=0.2226$), then the probability for which the random fluctuations of the effective stiffness tensor is less than 2%, 4% or 8%, is 0.36, 0.65 or 0.95 respectively.

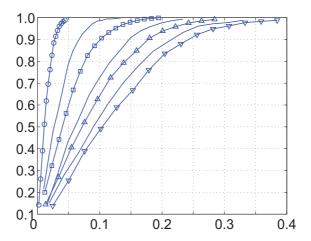


Figure 3. Graph of $\beta \mapsto \mathbb{P}(\beta) = P\{1 - \beta < Z \le 1 + \beta\}$ for $\delta = 0.4$ and for several values of the spatial correlation length $L_d = 0.1$ (circle), 0.2 (no marker), 0.3 (square), 0.4 (no marker), 0.5 (triangle-up), 0.6 (no marker), 0.7 (triangle-down). Horizontal axis β . Vertical axis $\mathbb{P}(\beta)$.

5. Conclusions

We have presented an approach introducing a microscale stochastic model which can be used when the standard method cannot be applied to anisotropic elastic microstructures. This new approach is useful for a direct experimental identification of random anisotropic elastic microstructures. A parametric probabilistic study of the RVE size has been performed with respect to the intensity δ_A of the stochastic fluctuations of the local stiffness tensor-valued random field describing the microscale stochastic model and in function of its correlation lengths L_k^A . Such a study is also useful to get information on the RVE size for which stochastic fluctuations are still significant and consequently, can be measured. Then the results presented in this paper could allow the microscale stochastic model to be identified from RVE- or macro-scale measurements solving an inverse stochastic problem. For fixed spatial correlation lengths and dispersion parameter δ_A , the effective RVE size can effectively be calculated in a probability sense using the cumulative distribution function of the random variable measuring the random fluctuations of the effective stiffness tensor. A large numerical simulation has been carried out and the numerical results obtained allow a probability analysis of the representative volume element size to be performed.

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