

Scattering mean free path in continuous complex media: Beyond the Helmholtz equation

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We present theoretical calculations of the ensemble-averaged (or effective or coherent) wave field propagating in a heterogeneous medium considered as one realization of a random process. In the literature, it is usually assumed that heterogeneity can be accounted for by a random scalar function of the space coordinates, termed the potential. Physically, this amounts to replacing the constant wave speed in Helmholtz' equation by a space-dependent speed. In the case of acoustic waves, we show that this approach leads to incorrect results for the scattering mean free path, no matter how weak the fluctuations. The detailed calculation of the coherent wave field must take into account both a scalar and an operator part in the random potential. When both terms have identical amplitudes, the correct value for the scattering mean free paths is shown to be more than 4 times smaller ($13/3$, precisely) in the low-frequency limit, whatever the shape of the correlation function. Based on the diagrammatic approach of multiple scattering, theoretical results are obtained for the self-energy and mean free path within Bourret's and on-shell approximations. They are confirmed by numerical experiments.

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I. INTRODUCTION

Whether quantum or classical, electromagnetic or acoustic, wave phenomena share a common theoretical ground. Hence the universality of fundamental concepts such as coherence, ballistic to diffuse transition, localization, and the observation of related experimental manifestations in all fields of mesoscopic wave physics [1–4].

In this paper, we are interested in the coherent field (i.e., the statistical average of the wave field propagating in an inhomogeneous medium whose characteristic are treated as random variables). In mesoscopic physics, determining the coherent field is the very basis of multiple scattering theory. It allows us to define a scattering mean free path ℓ_s , which is the key parameter in any multiple-scattering problem. ℓ_s is the typical decay length for the intensity of the coherent wave. In the case of a dilute suspension of discrete scatterers embedded in an homogeneous fluid, $1/\ell_s$ is equal to the scattering cross section of a single scatterer multiplied by the number of scatterers per unit volume [5].

Unlike discrete media, what we consider here is an inhomogeneous medium which varies continuously in space. In that case, heterogeneity can be characterized by a random function of the spatial coordinates, called the potential. In this paper, we show that the classical approach to express ℓ_s as a function of the fluctuations σ and correlation length ℓ_c of the potential is incorrect. This is due to an additional term in the acoustic wave equation, which is usually overlooked. When both terms have identical amplitudes, the correct value for the scattering mean free paths is 4 times smaller ($13/3$, precisely) in the low-frequency limit, no matter how weak the fluctuations are and whatever the shape of the correlation function (as long as its second-order moment is finite). As a result, even in the most simple cases (e.g., exponentially correlated disorder) the scattering mean free path can be severely underestimated.

The theoretical framework of the present paper is the diagrammatic approach of multiple scattering [4,6,7]. It yields

an exact equation for the coherent field known as Dyson's equation, the essential ingredient of which is the self-energy Σ . Unfortunately, as often in real life, one has to resort to some degree of approximation to evaluate Σ and obtain tractable expressions for the coherent field. The coherent wave has been extensively studied in the literature with various kinds of waves, both theoretically and experimentally [8–15].

From a theoretical point of view, in the case of continuous heterogeneous media the starting point is usually a wave equation in which disorder is introduced by a space-dependent wave velocity $c(\vec{r})$:

$$\Delta\phi - \frac{1}{c^2} \frac{\partial^2\phi}{\partial t^2} = s(\vec{r}, t). \quad (1)$$

Assuming linearity and time invariance, in order to determine the wave field $\phi(\vec{r}, t)$ generated by any distribution of sources $s(\vec{r}, t)$ it suffices to know the Green's function $g(\vec{r}, \vec{r}_s, t)$ i.e., the solution of Eq. (1) when the source term is $s(\vec{r}, t) = \delta(\vec{r} - \vec{r}_s)\delta(t)$, with appropriate boundary conditions.

An essential point is that in Eq. (1) heterogeneity is fully characterized by a random *scalar* $c(\vec{r})$. The average Green's function $\langle g \rangle$, hence the coherent field $\langle \phi \rangle$, can be calculated, provided the statistical properties of $c(\vec{r})$, particularly its correlation function, are known.

Here we are interested in a very simple case in which the usual approach fails. We consider an acoustic wave propagating in a lossless heterogenous fluid. It is well known that the wave equation for the acoustic pressure does not take the form of Eq. (1): It entails an additional term with a random *operator*, instead of a simple scalar [16–18]. The operator term is usually neglected when dealing with multiple scattering of waves. This implies an important error in the calculation of the mean free path, especially at low frequencies. To our knowledge, this point has been overlooked so far. Let us mention, however, Ref. [19], in the context of acoustic Cerenkov radiation by a moving source in a turbulent medium. Turner *et al.* [20] derived a Dyson equation in the case of isotropic solids with weak fluctuations of density and Lamé coefficients; yet the liquid limit (no shear modulus) of this

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model does not exactly yield the correct mean free path for a fluid, as will be discussed later.

The paper is organized as follows. In the next section we briefly recall the basics of the diagrammatic approach of multiple scattering, applied to the standard wave equation and show why this is inappropriate in the case of acoustic waves. Section III gives a complete solution of the wave equation for the average acoustic pressure. Under the Bourret approximation, we show that the self-energy Σ exhibits three additional terms compared to the standard scalar case. Therefore Sec. III, particularly Eq. (50), is the core of the paper. In the final section, we discuss the importance of these additional terms in the simple situation of an exponentially correlated disorder. Numerical experiments based on averages of numerous simulations of the complete wave equation are found in very good agreement with the theoretical results. Supplementary calculations and numerical details are presented in the Appendix.

II. THE USUAL WAY AND ITS LIMITATIONS

Let us consider first a homogenous and lossless medium, with a constant sound speed c_0 . The corresponding Green's function and its Fourier transforms will be denoted g_0 , G_0 , and \tilde{G}_0 , respectively. In the monochromatic regime, G_0 is the solution of Helmholtz' equation:

$$\Delta G_0 + \frac{\omega^2}{c_0^2} G_0 = \delta(\vec{r} - \vec{r}_s), \quad (2)$$

where ω is the angular frequency and $\omega/c_0 = k_0$ is a reference wave number. With the condition of causality, in unbounded 3D space we have:

$$g_0(\vec{r}, \vec{r}_s, t) = -\frac{1}{4\pi|\vec{r} - \vec{r}_s|} \delta(t - |\vec{r} - \vec{r}_s|/c_0), \quad (3)$$

$$G_0(\vec{r}, \vec{r}_s, \omega) = -\frac{1}{4\pi|\vec{r} - \vec{r}_s|} e^{jk_0|\vec{r} - \vec{r}_s|}, \quad (4)$$

$$\tilde{G}_0(\vec{k}, \omega) = \frac{1}{k_0^2 - k^2}, \quad (5)$$

with \vec{k} the dual variable for $\vec{r} - \vec{r}_s$ and $j = \sqrt{-1}$. The tilde $\tilde{\cdot}$ denotes a spatial Fourier transform. In the sequel, the analysis is performed in the frequency domain and the ω dependence is dropped for brevity.

A. Scalar potential

Assuming that a heterogeneous medium can be simply characterized by a space dependence of the wave speed amounts to replacing c_0 by $c(\vec{r})$ in the wave equation. Then it is convenient to define the scalar potential α :

$$\alpha(\vec{r}) = 1 - c_0^2/c^2(\vec{r}). \quad (6)$$

The monochromatic Green's function $G(\vec{r}, \vec{r}_s)$ in the heterogeneous medium is such that

$$\Delta G + k_0^2 G = k_0^2 \alpha G + \delta(\vec{r} - \vec{r}_s). \quad (7)$$

Note that in Eq. (7) $c_0 = \omega/k_0$ is a reference speed which could be chosen arbitrarily. It is often convenient to set

$$\frac{1}{c_0^2} = \left\langle \frac{1}{c^2(\vec{r})} \right\rangle, \quad (8)$$

so $\langle \alpha \rangle = 0$; the brackets denote an ensemble average.

For media such that the typical speed fluctuation δc is much smaller than the average speed $\langle c \rangle$, Eq. (8) amounts to choosing $c_0 = \langle c \rangle$, hence the reference speed is actually the average sound speed, but this is not true in the general case.

The potential α fully characterizes the heterogeneity, in that it measures the gap between the reference and the actual wave speed, at any point in space. The term *potential* comes from quantum mechanics, where the relevant wave field is the complex amplitude of probability and obeys Schrödinger equation, in which case the heterogeneity of the medium is an actual potential energy [4]. Here α is simply a dimensionless function of the space coordinate \vec{r} .

Equation (7) is similar to Eq. (2), with a source term $\alpha(\vec{r})G(\vec{r}, \vec{r}_s) + \delta(\vec{r} - \vec{r}_s)$ that entails the Green's function itself. Hence G can be expressed implicitly in a recursive manner (Lippmann-Schwinger form) as:

$$G(\vec{r}, \vec{r}_s) = G_0(\vec{r}, \vec{r}_s) + k_0^2 \int G_0(\vec{r}, \vec{r}_1) \alpha(\vec{r}_1) G(\vec{r}_1, \vec{r}_s) d^3 r_1. \quad (9)$$

For an arbitrary source distribution, the resulting wave field at \vec{r} would be

$$\phi(\vec{r}) = \int G(\vec{r}, \vec{r}_s) s(\vec{r}_s) d^3 r_s. \quad (10)$$

Substituting G under the integral by the right-hand side of Eq. (9) and reiterating the process yields an exact expression for G as an infinite sum of multiple integrals, known as Born's expansion:

$$\begin{aligned} G(\vec{r}, \vec{r}_s) &= G_0(\vec{r}, \vec{r}_s) + k_0^2 \int G_0(\vec{r}, \vec{r}_1) \alpha(\vec{r}_1) G_0(\vec{r}_1, \vec{r}_s) d^3 r_1 \\ &+ k_0^4 \iint G_0(\vec{r}, \vec{r}_1) \alpha(\vec{r}_1) G_0(\vec{r}_1, \vec{r}_2) \\ &\times \alpha(\vec{r}_2) G_0(\vec{r}_2, \vec{r}_s) d^3 r_1 d^3 r_2 + \dots \end{aligned} \quad (11)$$

The single-scattering approximation (which is commonly made in imaging of weakly heterogeneous media) consists in neglecting all terms beyond the first integral on the right-hand side of Eq. (11). In that case the Green's function G can be easily computed for any function α , and the inverse problem (i.e., reconstructing α from G) may be solved. Naturally, this approach completely fails as soon as multiple scattering is not negligible.

In this paper, we consider multiple scattering of waves but we do not aim at solving Eq. (9). Considering α as a random variable with known statistical parameters, we are interested in the statistical average $\langle \phi \rangle$ of the wave field, which amounts to determine $\langle G \rangle$. We will refer to it as the *coherent* field. This approach implies that we consider a given medium as one particular realization, among the infinity of possible outcomes, of the same random process. From a physical point of view, what an experimentalist would measure with a source distribution $s(\vec{r}_s)$ and a point receiver at \vec{r} is ϕ , not $\langle \phi \rangle$. But ϕ can be formally written as $\langle \phi \rangle + \delta\phi$ with $\langle \delta\phi \rangle = 0$ (i.e., a mean

term plus statistical fluctuations, changing from one realization to an other). How $\langle \phi \rangle$ can be estimated in an actual experiment, and how robust the estimation is, is not our concern in the present paper. We focus on theoretical calculations for $\langle \phi \rangle$, derived from the statistical properties of α . By taking the average of the Born expansion, this will obviously require us to know the statistical moments of α with any order n [i.e., quantities such as $\langle \alpha(\vec{r}_1)\alpha(\vec{r}_2) \dots \alpha(\vec{r}_n) \rangle$].

The diagrammatic theory of multiple scattering shows that $\langle G \rangle$ obeys an exact equation, known as Dyson's equation [6,7]. The basic ingredient in Dyson's equation is a quantity Σ referred to as the self-energy or the mass operator in the literature. Σ can be fully determined by the statistical properties of α and takes into account all orders of multiple scattering. In a nutshell, the basic idea is to rewrite the statistical average of Eq. (11) as an implicit, recursive expression for $\langle G \rangle$. The resulting Dyson's equation reads:

$$\langle G(\vec{r}, \vec{r}_s) \rangle = G_0(\vec{r}, \vec{r}_s) + \iint G_0(\vec{r}, \vec{r}_1) \Sigma(\vec{r}_1, \vec{r}_2) \times \langle G(\vec{r}_2, \vec{r}_s) \rangle d^3r_1 d^3r_2. \quad (12)$$

Assuming that the medium is statistically homogeneous (i.e., its statistical parameters are invariant under translation) $\Sigma(\vec{r}_1, \vec{r}_2)$ only depends on $\vec{x} = \vec{r}_1 - \vec{r}_2$, and since so does G_0 , Eq. (12) is a double convolution product on the variable \vec{x} . Hence it can be simply solved by a spatial Fourier transform, which yields

$$\langle \tilde{G}(\vec{k}) \rangle = \frac{1}{k_0^2 - k^2 - \tilde{\Sigma}(\vec{k})}, \quad (13)$$

where \vec{k} is the dual variable for \vec{x} . Assuming further that the medium is statistically isotropic (i.e., its statistical parameters are also invariant under rotation), both $\tilde{\Sigma}$ and \tilde{G} only depend on $k = |\vec{k}|$.

The key issue is naturally to determine Σ . Mathematically, Σ can be written as a perturbative development in $1/(k_0 \ell_s)$, an infinite series of integrals involving statistical moments of α at all orders, which can be represented by the following diagrams:

$$\Sigma = \text{O} + \text{O} \text{---} \text{O} + \text{O} \text{---} \text{O} \text{---} \text{O} + \dots \quad (14)$$

Following the usual conventions, a continuous line joining two points represents the free-space Green's function G_0 between these points; a dashed line linking n points is the n -order moment of α multiplied by k_0^{2n} . The inner points of a diagram are dummy variables. For the establishment of Eq. (14), see, for instance, Refs. [6,7].

The Bourret approximation (or first-order smoothing approximation) only keeps the first two terms in the perturbative development of Σ . This yields a simple analytical expression for Σ as a function of the first- and second-order moments [i.e., the mean $\langle \alpha \rangle$ and correlation function $C_{\alpha\alpha}(\vec{r}_1 - \vec{r}_2) = \langle \alpha(\vec{r}_1)\alpha(\vec{r}_2) \rangle$]. Note that the Bourret approximation does not imply at all that multiple scattering terms are neglected beyond second-order scattering but rather than all multiple scattering events are assumed to be similar to a succession of uncorrelated single or double-scattering sequences.

Under the Bourret approximation, and having chosen c_0 such that $\langle \alpha \rangle = 0$, the expression for the self-energy is [6,7]:

$$\Sigma(\vec{x}) = k_0^4 G_0(\vec{x}) C_{\alpha\alpha}(\vec{x}). \quad (15)$$

The last step is to determine the average Green's function from Eq. (13). In the most general case, it is an arbitrary function of \vec{k} . To determine the average Green's function $\langle G(\vec{x}) \rangle$ in real space, one has to perform an inverse Fourier transform. This is not always possible analytically and does not always lead to a simple effective medium; Σ is said to be nonlocal. We will not discuss these issues in the present paper. Instead, we make a further approximation referred to as the on-shell approximation. It usually requires $\tilde{\Sigma}(\vec{k})$ to be sufficiently smaller than k_0^2 . Indeed, if $\tilde{\Sigma}(\vec{k})$ is weak enough, then we can reasonably assume that the effect it will have on the homogeneous wave number k_0 is small, so when performing the inverse three-dimensional Fourier transform, the volume that essentially contributes to $\langle G(\vec{r} - \vec{r}_s) \rangle$ is the vicinity of the shell defined by $|k| = k_0$. In other words, this amounts to performing a zero-order development of the self-energy around k_0 , hence replacing $\tilde{\Sigma}(\vec{k})$ by $\tilde{\Sigma}(\vec{k}_0)$ in Eq. (13). In that case the expression of the average Green's function in real space is straightforward:

$$\langle G(\vec{r} - \vec{r}_s) \rangle = -\frac{1}{4\pi |\vec{r} - \vec{r}_s|} e^{jk_{\text{eff}}|\vec{r} - \vec{r}_s|}. \quad (16)$$

This means that the average Green's function is that of a fictitious homogeneous absorbing medium with a complex-valued wave number k_{eff} such that

$$k_{\text{eff}} = \sqrt{k_0^2 - \tilde{\Sigma}(\vec{k}_0)}. \quad (17)$$

Equation (17) is a dispersion relation from which phase and group velocities for the coherent field can be determined. Most importantly, the intensity of the coherent wave is found to decay exponentially with the distance (Beer-Lambert's law). Since there is no absorption, the losses are entirely due to scattering and the scattering mean-free path is $\ell_s = 1/[2 \text{Im}(k_{\text{eff}})]$.

ℓ_s is an essential parameter in all multiple scattering problems. Particularly, the range of validity of the Bourret approximation can be shown to be $k_0 \ell_s \gg 1$ [7]. It should also be mentioned that, as a refinement of the on-shell approximation, k_{eff} can be determined more accurately with an iterative algorithm using k_0 as a first guess.

So, within the Bourret approximation, as long as the correlation function for the potential α is known, the effective wave speed and scattering mean free path can be determined quite easily and sometimes analytically.

A typical example is that of a disordered random medium with an exponentially decaying correlation function $C_{\alpha\alpha}(\vec{x}) = \sigma_\alpha^2 \exp(-x/\ell_c)$. σ_α^2 denotes the variance of α and ℓ_c its correlation length. In that case, under the Bourret approximation the self-energy is

$$\tilde{\Sigma}(\vec{k}) = -\frac{\sigma_\alpha^2 k_0^4}{k^2 + (jk_0 - 1/\ell_c)^2}. \quad (18)$$

The on-shell approximation yields

$$k_{\text{eff}} = k_0 \sqrt{1 + \frac{(\sigma_\alpha k_0 \ell_c)^2}{1 - 2jk_0 \ell_c}}, \quad (19)$$

whose imaginary part determines the scattering mean free path. Furthermore, if $\Sigma(k_0) \ll k_0^2$, a first-order development of the square root gives a simple analytic expression for the scattering mean free path ℓ_s , as a function of frequency, correlation length, and variance of the potential:

$$\ell_s = \frac{1}{\sigma_\alpha^2 k_0} \frac{1 + 4k_0^2 \ell_c^2}{2k_0^3 \ell_c^3}, \quad (20)$$

allowing us to work with practical dimensionless quantities and express $k_0 \ell_s \sigma_\alpha^2$ as a function of $k_0 \ell_c$. Beyond this simple example, whatever the shape of the correlation function and whatever the nature of the wave (acoustic, electromagnetic, etc.), the same formalism will hold as soon as we deal with a wave field propagating in a heterogeneous medium in which heterogeneity is fully described by a scalar function such as α . Such is the case when the local wave speed $c(\vec{r})$ suffices to capture the heterogeneity, which is usually assumed as a starting point in multiple scattering theories. However, this description of heterogeneity may sometimes be completely misleading, even in very simple situations.

B. Operator potential

Let us consider the case of acoustic waves in a lossless fluid. The medium is characterized by its mass density ρ and compressibility χ at rest. With no sources, the linearized equations of elastodynamics read

$$\rho \frac{\partial \vec{v}}{\partial t} = -\vec{\nabla} p, \quad (21)$$

$$\vec{\nabla} \cdot \vec{u} = -\chi p. \quad (22)$$

$\vec{u}(\vec{r}, t)$ is the displacement undergone by the particle initially at rest at point \vec{r} , $\vec{v} = \partial \vec{u} / \partial t$ is the particle velocity, and $p(\vec{r}, t)$ is the acoustic pressure. \vec{u} , \vec{v} , and p are first-order infinitesimal quantities. To establish Eqs. (21) and (22) all second-order nonlinear terms have been neglected whatever their physical origin (convective or thermodynamic). From a physical point of view, Eqs. (21) and (22) are an expression of Newton's second law and Hooke's law (the relative dilation $\vec{\nabla} \cdot \vec{u}$ undergone by an infinitesimal volume of fluid is proportional and opposed to the acoustic pressure). If neither ρ nor χ depend on space coordinate \vec{r} , then Eqs. (21) and (22) lead to the usual wave equation with a constant sound velocity $c_0 = 1/\sqrt{\rho\chi}$, which applies to all quantities describing the sound wave (acoustic pressure, velocity, displacement, dilation, etc.).

In a heterogeneous fluid, the local sound velocity naturally depends on the space coordinate \vec{r} as $c(\vec{r}) = 1/\sqrt{\rho(\vec{r})\chi(\vec{r})}$. It is therefore tempting to replace c_0 by $c(\vec{r})$ in Helmholtz's equation for a homogeneous fluid [Eq. (2)], but this is not always correct.

Combining Eqs. (21) and (22) yields the following equations for the acoustic pressure and velocity:

$$\Delta p - \frac{1}{c^2} \frac{\partial^2 p}{\partial t^2} - \frac{\vec{\nabla} \rho \cdot \vec{\nabla} p}{\rho} = 0, \quad (23)$$

$$\vec{\Delta} \vec{v} - \frac{1}{c^2} \frac{\partial^2 \vec{v}}{\partial t^2} - \frac{\vec{\nabla} \chi \cdot \vec{\nabla} \cdot \vec{v}}{\chi} = 0. \quad (24)$$

If ρ does not vary in space, then Eq. (23) yields the usual wave equation for the acoustic pressure p , with a space-dependent velocity c . And if χ does not vary in space, Eq. (24) yields the usual wave equation for the velocity \vec{v} . But in the general case where *both* ρ and χ vary with \vec{r} , none of the acoustic variables satisfy the usual wave equation [16,17]. However, the resulting equation for the acoustic pressure [Eq. (23)] can be Fourier transformed over time and then rearranged in order to define a potential, as we did before. We obtain an equation similar to Helmholtz's:

$$\Delta P + k_0^2 P = k_0^2 \gamma P. \quad (25)$$

The potential γ is such that:

$$\gamma(\vec{r}) = \alpha(\vec{r}) + \frac{1}{k_0^2} \vec{\nabla} \beta(\vec{r}) \cdot \vec{\nabla}. \quad (26)$$

α is defined in Eq. (6) and

$$\beta(\vec{r}) = \ln \left[\frac{\rho}{\rho_0} \right]. \quad (27)$$

ρ_0 is an arbitrary constant with the dimensions of a mass density.

The first term in the expression of γ is the usual potential α , related to the spatial fluctuations of sound speed. In addition, there is an other term, which unlike α is not a simple scalar but an *operator* acting on the field it is applied to. It should be noticed that the same problem arises for different kinds of waves, e.g., electromagnetic waves propagating in a heterogeneous medium showing fluctuations of both relative permeability $\mu(\vec{r}, \omega)$ and permittivity $\epsilon(\vec{r}, \omega)$. In that case, Maxwell's equations yield the following wave equation for the monochromatic electric field $\vec{E}(\vec{r}, \omega)$, with a potential that contains an operator part [21]:

$$\vec{\nabla} \times \vec{\nabla} \times \vec{E} - k_0^2 \epsilon \mu \vec{E} - \vec{\nabla} [\ln(\mu)] \times \vec{\nabla} \times \vec{E} = \vec{0}. \quad (28)$$

Note that the real issue is not to determine when fluctuations in permeability (or density, in the acoustic case) can be neglected compared to fluctuations in permittivity (respectively, compressibility): Both kinds of fluctuations, whether separated or combined, will result in a space-dependent wave speed $c(\vec{r})$. The question should rather be set in terms of comparing the operator part and the scalar part in the potential describing the heterogeneity.

Whatever the physical nature of the wave, the applicability of the diagrammatic techniques when the relevant potential γ has both a scalar part and an operator part as well as the impact of the operator part on the final result are not trivial. In the next section, we deal with this problem and obtain the average Green's function, in the case of acoustic waves.

III. SELF-ENERGY CALCULATION

In this section, the acoustic pressure P is chosen as the relevant variable for the wave field. The issue is to determine the average Green's function of Eq. (25). The reference velocity is still chosen according to Eq. (8). Since the medium is assumed to be statistically invariant under translation $\langle \beta(\vec{r}) \rangle$

does not depend on the space coordinate \vec{r} . Hence, despite the presence of β , the average value of the potential γ will still be zero, since

$$\langle \vec{\nabla} \beta \rangle = \vec{\nabla} \langle \beta \rangle = \vec{0}. \quad (29)$$

As usual, the Green's function for Eq. (25) can be written as a Lippmann-Schwinger equation,

$$G(\vec{r}, \vec{r}_s) = G_0(\vec{r}, \vec{r}_s) + k_0^2 \int G_0(\vec{r}, \vec{r}_1) \gamma(\vec{r}_1) G(\vec{r}_1, \vec{r}_s) d^3 r_1. \quad (30)$$

γ is not a simple scalar function, which precludes the usual definition of its autocorrelation function and that of the self-energy. To overcome this difficulty, we start by introducing a two-variable potential V such that

$$\begin{aligned} V(\vec{r}_1, \vec{r}_2) &= \gamma(\vec{r}_1) \delta(\vec{r}_1 - \vec{r}_2) \\ &= \alpha(\vec{r}_1) \delta(\vec{r}_1 - \vec{r}_2) + \frac{1}{k_0^2} \vec{\nabla}_{\vec{r}_1} \beta(\vec{r}_1) \cdot \vec{\nabla}_{\vec{r}_1} \delta(\vec{r}_1 - \vec{r}_2). \end{aligned} \quad (31)$$

Equation (30) becomes

$$\begin{aligned} G(\vec{r}, \vec{r}_s) &= G_0(\vec{r}, \vec{r}_s) + k_0^2 \iint G_0(\vec{r}, \vec{r}_1) V(\vec{r}_1, \vec{r}_2) \\ &\quad \times G(\vec{r}_2, \vec{r}_s) d^3 r_1 d^3 r_2. \end{aligned} \quad (32)$$

The next steps are as usual to develop Eq. (32) into a Born expansion by iteration and then to take its statistical average and write it under a recursive form (Dyson's equation). Under the Bourret approximation, only the first two terms in the self-energy are kept. The first one vanishes since c_0 is set so $\langle \gamma \rangle = 0$. The second term reads:

$$\Sigma(\vec{r}_a, \vec{r}_b) = k_0^4 \iint \langle V(\vec{r}_a, \vec{r}_1) G_0(\vec{r}_1, \vec{r}_2) V(\vec{r}_2, \vec{r}_b) \rangle d^3 r_1 d^3 r_2. \quad (33)$$

As a consequence, the self-energy Eq. (33) gives rise to four terms, involving the following dimensionless correlation functions and their derivatives:

$$\begin{aligned} C_{\alpha\alpha}(\vec{r}_1, \vec{r}_2) &= \langle \alpha(\vec{r}_1) \alpha(\vec{r}_2) \rangle, \\ C_{\beta\beta}(\vec{r}_1, \vec{r}_2) &= \langle \beta(\vec{r}_1) \beta(\vec{r}_2) \rangle, \\ C_{\alpha\beta}(\vec{r}_1, \vec{r}_2) &= \langle \alpha(\vec{r}_1) \beta(\vec{r}_2) \rangle, \\ C_{\beta\alpha}(\vec{r}_1, \vec{r}_2) &= \langle \beta(\vec{r}_1) \alpha(\vec{r}_2) \rangle. \end{aligned} \quad (34)$$

We assume that the random processes α and β are *jointly* stationary and invariant under rotation. Then all correlation functions only depend on $x = |\vec{r}_1 - \vec{r}_2|$ and $C_{\alpha\beta}(x) = C_{\beta\alpha}(x)$. Only three correlation functions suffice to characterize the disorder. They can be rewritten as

$$\begin{aligned} C_{\alpha\alpha}(x) &= \sigma_\alpha^2 c_{\alpha\alpha}(x), \\ C_{\beta\beta}(x) &= \sigma_\beta^2 c_{\beta\beta}(x), \\ C_{\alpha\beta}(x) &= C_{\beta\alpha}(x) = \sigma_\alpha \sigma_\beta c_{\alpha\beta}(x). \end{aligned} \quad (35)$$

σ_α^2 , σ_β^2 are the variances of α and β , respectively. $c_{\alpha\alpha}(x)$ is the correlation coefficient between $\alpha(\vec{r})$ and $\alpha(\vec{r} + \vec{x})$, $c_{\beta\beta}(x)$ is the correlation coefficient between $\beta(\vec{r})$ and $\beta(\vec{r} + \vec{x})$, $c_{\alpha\beta}(x)$ is the correlation coefficient between $\alpha(\vec{r})$ and $\beta(\vec{r} + \vec{x})$. Replacing

V in Eq. (33) by Eq. (31), we can write the self-energy Σ as a sum of four contributions:

$$\Sigma = \Sigma_{\alpha\alpha} + \Sigma_{\beta\alpha} + \Sigma_{\alpha\beta} + \Sigma_{\beta\beta}. \quad (36)$$

A. First term

The first term is

$$\Sigma_{\alpha\alpha}(\vec{r}_a - \vec{r}_b) = k_0^4 \langle \alpha(\vec{r}_a) G_0(\vec{r}_a - \vec{r}_b) \alpha(\vec{r}_b) \rangle \quad (37)$$

$$= k_0^4 G_0(\vec{r}_a - \vec{r}_b) C_{\alpha\alpha}(\vec{r}_a - \vec{r}_b). \quad (38)$$

With $\vec{x} = \vec{r}_a - \vec{r}_b$ we find the usual contribution to the self-energy Eq. (15), involving only the scalar potential α .

The three additional terms are more complicated, since they entail combinations of α and β as well as spatial derivatives.

B. Second term

The second term mixes contributions from β and α :

$$\begin{aligned} \Sigma_{\beta\alpha}(\vec{r}_a - \vec{r}_b) &= k_0^2 \int \langle \vec{\nabla}_{\vec{r}_a} \beta(\vec{r}_a) \cdot \vec{\nabla}_{\vec{r}_a} [\delta(\vec{r}_a - \vec{r}_1)] \\ &\quad \times G_0(\vec{r}_1 - \vec{r}_b) \alpha(\vec{r}_b) \rangle d^3 r_1. \end{aligned} \quad (39)$$

Performing two integrations by parts and using the properties of the Dirac distribution yields:

$$\Sigma_{\beta\alpha}(\vec{r}_a - \vec{r}_b) = k_0^2 \vec{\nabla}_{\vec{r}_a} G_0(\vec{r}_a - \vec{r}_b) \cdot \vec{\nabla}_{\vec{r}_a} C_{\beta\alpha}(\vec{r}_a - \vec{r}_b). \quad (40)$$

Taking advantage of the assumed radial symmetry, we obtain:

$$\Sigma_{\beta\alpha}(x) = k_0^2 \frac{\partial G_0}{\partial x} \frac{\partial C_{\beta\alpha}}{\partial x}. \quad (41)$$

C. Third term

Similarly to the second term, the third one implies both β and α :

$$\begin{aligned} \Sigma_{\alpha\beta}(\vec{r}_a - \vec{r}_b) &= k_0^2 \int \langle \alpha(\vec{r}_a) G_0(\vec{r}_a - \vec{r}_2) \\ &\quad \vec{\nabla}_{\vec{r}_2} [\beta(\vec{r}_2)] \cdot \vec{\nabla}_{\vec{r}_2} [\delta(\vec{r}_2 - \vec{r}_b)] \rangle d^3 r_2. \end{aligned} \quad (42)$$

Again, performing two integrations by part and using the properties of the Dirac distribution, we obtain:

$$\Sigma_{\alpha\beta}(\vec{r}_a - \vec{r}_b) = -k_0^2 \vec{\nabla}_{\vec{r}_b} \cdot [G_0(\vec{r}_a - \vec{r}_b) \vec{\nabla}_{\vec{r}_b} \{C_{\alpha\beta}(\vec{r}_a - \vec{r}_b)\}]. \quad (43)$$

Since all functions involved here have radial symmetry, the expression above simplifies into

$$\Sigma_{\alpha\beta}(x) = -\frac{k_0^2}{x^2} \frac{\partial}{\partial x} \left[x^2 G_0(x) \frac{\partial C_{\alpha\beta}}{\partial x} \right]. \quad (44)$$

D. Fourth term

The last term is the most complicated one:

$$\begin{aligned} \Sigma_{\beta\beta}(\vec{r}_a - \vec{r}_b) &= \iint \langle \vec{\nabla}_{\vec{r}_a} \beta(\vec{r}_a) \cdot \vec{\nabla}_{\vec{r}_a} [\delta(\vec{r}_a - \vec{r}_1)] \\ &\quad \times G_0(\vec{r}_1 - \vec{r}_2) \vec{\nabla}_{\vec{r}_2} \beta(\vec{r}_2) \cdot \vec{\nabla}_{\vec{r}_2} [\delta(\vec{r}_2 - \vec{r}_b)] \rangle \\ &\quad \times d^3 r_1 d^3 r_2. \end{aligned} \quad (45)$$

Equation (45) contains a product of two scalar products, which can be written as a tensorial product. Again, two integrations by parts and the properties of the Dirac distribution yield:

$$\Sigma_{\beta\beta}(\vec{r}_a - \vec{r}_b) = -\vec{\nabla}_{\vec{r}_b} \cdot [\vec{\nabla}_{\vec{r}_b} \otimes \vec{\nabla}_{\vec{r}_a} \{C_{\beta\beta}(\vec{r}_a - \vec{r}_b)\} \vec{\nabla}_{\vec{r}_a} G_0(\vec{r}_a - \vec{r}_b)]. \quad (46)$$

Given the radial symmetry, in three dimensions we have:

$$\Sigma_{\beta\beta}(x) = -\vec{\nabla} \cdot [\vec{\nabla} \otimes \vec{\nabla} \{C_{\beta\beta}(x)\} \vec{\nabla} G_0(x)]. \quad (47)$$

The tensorial product between the two gradients is a Hessian matrix. In spherical coordinates and for a function with radial symmetry, we have [22]:

$$\vec{\nabla} \otimes \vec{\nabla} C_{\beta\beta} = \begin{bmatrix} \frac{\partial^2 C_{\beta\beta}}{\partial x^2} & 0 & 0 \\ 0 & \frac{1}{x} \frac{\partial C_{\beta\beta}}{\partial x} & 0 \\ 0 & 0 & \frac{1}{x} \frac{\partial C_{\beta\beta}}{\partial x} \end{bmatrix}. \quad (48)$$

Hence:

$$\Sigma_{\beta\beta}(x) = -\frac{1}{x^2} \frac{\partial}{\partial x} \left[x^2 \frac{\partial^2 C_{\beta\beta}}{\partial x^2} \frac{\partial G_0}{\partial x} \right]. \quad (49)$$

IV. RESULTS AND DISCUSSION

In real 3D space, the four contributions to the self-energy add up to give:

$$\Sigma(x) = k_0^4 \sigma_\alpha^2 G_0 c_{\alpha\alpha} - k_0^2 \sigma_\alpha \sigma_\beta G_0 [c''_{\alpha\beta} + 2c'_{\alpha\beta}/x] - \sigma_\beta^2 [c''_{\beta\beta} G_0'' + c'''_{\beta\beta} G_0' + 2c''_{\beta\beta} G_0'/x]. \quad (50)$$

For simplicity the x dependence of G_0 and of the correlation coefficients have been omitted, and the prime means derivation with respect to x .

If $\sigma_\beta = 0$, then the self-energy Eq. (50) is reduced to the usual (i.e., scalar only) term $\Sigma_{\alpha\alpha}$. In the general case where heterogeneity is such that a scalar (α) and an operator (β) term coexist, it is not obvious to determine the orders of magnitude of the additional terms in the self-energy, since they involve five physical parameters: two variances and three correlation lengths. In order to highlight the importance of the additional terms relatively to the first one, we focus on the most simple case where $\sigma_\alpha = \sigma_\beta = \sigma$ and $c_{\alpha\alpha} = c_{\alpha\beta} = c_{\beta\beta}$. The variance σ^2 appears as a mere multiplicative term, and from a physical point of view everything will depend on the typical correlation length ℓ_c . Exponential and Gaussian correlation functions were tested, and similar trends were obtained. We only give the result for the exponential case in 3D, which entails simpler analytical expressions. The 2D case is dealt with in the Appendix (Sec. A 2).

In the case of an exponentially correlated disorder, we have

$$c_{\alpha\alpha}(x) = \exp\left[-\frac{x}{\ell_c}\right]. \quad (51)$$

In 3D, the calculation of $\tilde{\Sigma}$ yields:

$$\tilde{\Sigma}(k_0) = -\frac{\sigma^2}{\ell_c^2} \left[\frac{(k_0 \ell_c)^4 + j k_0 \ell_c}{1 - 2j k_0 \ell_c} - \frac{1 - 2(k_0 \ell_c)^2}{k_0 \ell_c} \arctan\left(\frac{k_0 \ell_c}{1 - j k_0 \ell_c}\right) \right]. \quad (52)$$

As long as $k_0 \ell_c \gg 1$ (i.e., the correlation length is much larger than the wavelength), the first term dominates:

$$\tilde{\Sigma}(k_0) \sim \tilde{\Sigma}_{\alpha\alpha}(k_0) = -\frac{\sigma^2}{\ell_c^2} \frac{(k_0 \ell_c)^4}{1 - 2j k_0 \ell_c}. \quad (53)$$

Therefore at high frequencies, even though the scalar and operator parts have equal importance in the random potential ($\sigma_\alpha = \sigma_\beta$), considering the usual wave equation with a space-dependent wave speed $c(\vec{r})$ instead of c_0 is legitimate to determine the coherent pressure field. However, it becomes completely wrong as soon as $k_0 \ell_c$ is comparable to unity. In that case, the impact of the three additional terms ($\tilde{\Sigma}_{\alpha\beta}$, $\tilde{\Sigma}_{\beta\alpha}$, and $\tilde{\Sigma}_{\beta\beta}$) on the effective wave number k_{eff} and particularly the scattering mean free path $\ell_s = 1/[2 \text{Im}(k_{\text{eff}})]$ can be far from negligible. More precisely, the difference is less than 6% for $k_0 \ell_c > 10$; but below $k_0 \ell_c \sim 1.5$, the three additional terms in the self-energy are larger than the first one. As a result, at low frequencies the actual mean free path can be more than 4 times smaller than expected. The exact ratio is 13/3; the same behavior was obtained in the case of a Gaussian-correlated disorder. Interestingly, it can be shown that the 13/3 ratio is independent of the correlation function (as long as its second-order moment is finite, see the Appendix, paragraph A 1).

As an illustration, Figs. 1 and 2 compare the scattering mean free paths obtained with (ℓ_s) and without [$\ell_s^{(\alpha\alpha)}$] the additional terms.

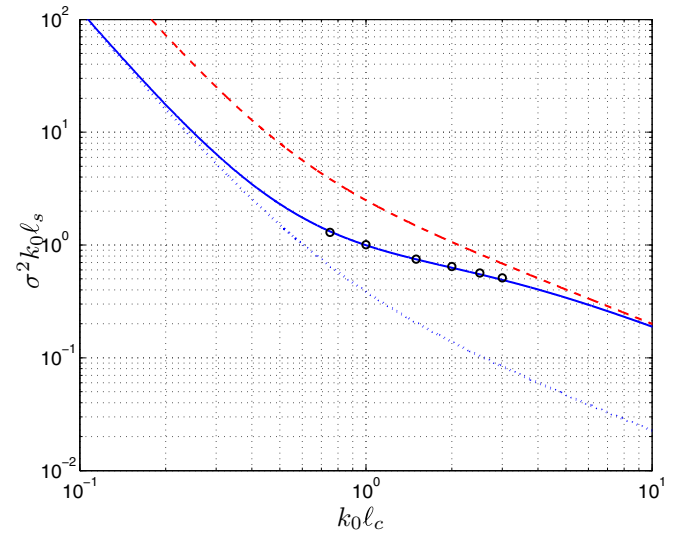


FIG. 1. (Color online) Scattering mean free path as a function of $k_0 \ell_c$ with (solid line) and without (dashed line) the additional terms in the self-energy for an exponentially correlated disorder. Results from the numerical experiments are indicated by black circles. For comparison, the fluid limit of the model from Ref. [20] is also plotted (dotted line).

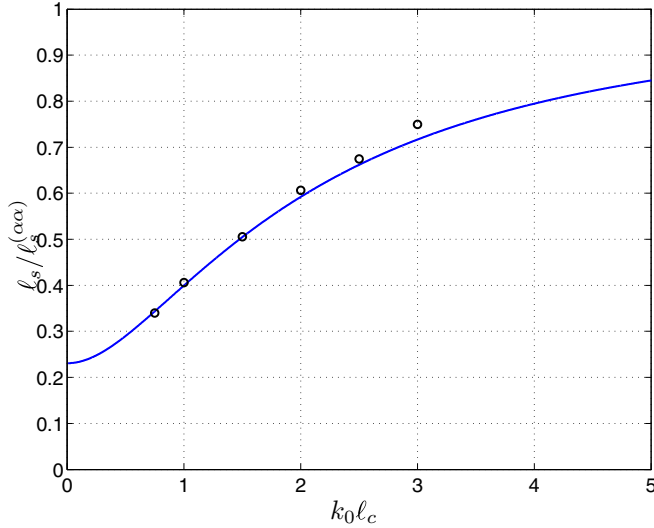


FIG. 2. (Color online) Ratio of the scattering mean free paths obtained with and without the additional terms in the self-energy as a function of $k_0\ell_c$ for an exponentially correlated disorder. Results from the numerical experiments are indicated by black circles. The value at zero frequency is $3/13$.

Note that care should be taken when taking the low-frequency limit; the on-shell approximation usually requires $\tilde{\Sigma}(k_0)$ to be much smaller than k_0^2 , hence [from Eq. (52)] when $k_0\ell_c \rightarrow 0$ the results are consistent only if the variance is kept such that $\sigma^2 \ll (k_0\ell_c)^2$. Interestingly, in the standard (scalar) case at low frequency the same condition implies that $\sigma^2 \ll 1/(k_0\ell_c)^2$. This means that whatever the fluctuations σ , a weak disorder approximation [$\tilde{\Sigma}(k_0) \ll k_0^2$] is always fulfilled at zero frequency if the random operator is purely scalar. This is no longer true when the operator term cannot be neglected: For a finite σ , there is a cut-off frequency (typically $k_0\ell_c \sim \sigma$) below which $\tilde{\Sigma}(k_0)$ is not small compared to k_0^2 .

In order to test the validity of the theoretical results above, we have performed numerical simulations of the inhomogeneous wave equations [(21) and (22)], using a finite-difference software developed in our lab [23]¹. Simulations were carried out for conditions typical of ultrasonic experiments. The reference (unperturbed) medium was water ($c_0 = 1500$ m/s, $\rho_0 = 1000$ kg/m³) and the incoming waveform was a short pulse with a central frequency ranging from 1 to 2 MHz. Using a random number generator, exponentially correlated 3D maps with zero mean and standard deviation σ were fabricated and used for both α and β . For a given realization of disorder, once $\alpha(\vec{r})$ and $\beta(\vec{r})$ are determined the corresponding density, compressibility, and sound speed are accessible from Eqs. (6) and (27). The typical correlation length was $\ell_c = 0.240$ mm, so $k_0\ell_c = 1$ at 1 MHz. Given the frequency spectrum, $k_0\ell_c$ spans typically between 0.75 and 3 in the numerical experiments. Further details on the simulations are given in the Appendix, Sec. A 3.

A plane wave is launched from one side of the random medium ($z = 0$) in the z direction. The resulting pressure is

measured at every grid point (x, y, z) and time t . A robust estimation of the coherent wave is obtained by a two-step average. For each realization of disorder, the pressure field is averaged in the (x, y) plane, as would do a plane detector perpendicular to the initial direction of propagation. Second, ensemble averaging is performed over at least 25 different realizations of disorder. As a result we obtain an estimate of the coherent pressure field $\langle p(z, t) \rangle$ as a function of depth z and time t . In all the numerical experiments, the total thickness of the map was at least $3\ell_s$ and we always ensured that the measured wave front was an accurate estimator of the coherent field (that is to say, remaining random fluctuations could be considered as negligible). A digital Fourier transform is performed; the coherent field's intensity $I_c = |\langle P(z, \omega) \rangle|^2$ is found to decay exponentially with z . An estimation of the scattering mean free path is obtained by a linear fit of $\log(I_c)$ with z at each frequency. Simulating both types of media separately (either $\sigma_\beta = 0$ hence no operator term in the random potential, or $\sigma_\beta = \sigma_\alpha = \sigma$), we plotted in Fig. 2 the ratio of the corresponding mean free paths. The results are in very good agreement with the analytical results presented earlier and support the validity of the theoretical analysis. The numerical results were also compared to another model derived from acoustics in polycrystals with randomly varying elastic properties but macroscopically isotropic [20]. In the limit where the second Lamé coefficient μ tends to zero (no shear stress), the results should be valid for the case of an inhomogeneous liquid. Interestingly, this model does predict the $13/3$ factor at zero frequency, yet it yields incorrect results at higher frequencies, especially above $k_0\ell_c > 0.1$. The essential reason is that in the solid model, the fluctuations in mass density and elastic constants are assumed to be very weak from the very beginning (i.e., the linearized equations of elastodynamics). In our approach, fluctuations are not necessarily weak initially, what is considered as weak is the second-order terms in the development of the self-energy (Bourret approximation). The weak fluctuation limit, if necessary, is only taken afterwards. Assuming that the fluctuations are weak from the beginning amounts to misestimate some of the additional terms in the self-energy. In the case of a fluid with $\sigma_\alpha = \sigma_\beta$, our results indicate that they cannot be discarded, no matter how weak fluctuations are; hence the results presented here for an inhomogeneous fluid cannot be seen as a particular case of the solid model. Note that we do not claim at all that the solid model in Ref. [20] is wrong: it is very well suited for polycrystals (e.g., coarse-grain steel), in which fluctuations of mass density and elasticity are indeed very weak compared to their mean value.

It is also interesting to plot the exponent $n = -\omega/\ell_s d\omega/d\ell_s$ as a function of frequency (see Fig. 3). Indeed, since a power-law dependence of the attenuation length is often assumed, n commonly serves as an indicator of the scattering regime. In both cases, $1/\ell_s$ is found to be proportional to ω^4 at low frequency and ω^2 at high frequency. These two trends are usually referred to as the Rayleigh and stochastic regimes and are used to characterize scattering media based on the measured dependence of acoustic attenuation with frequency. Figure 3 shows how misleading the omission of the additional terms in the wave equation can be, especially at intermediate frequencies ($k_0\ell_c \sim 1$): The exponent can be 35% lower than expected.

¹All information regarding the software are available on the website www.simsonic.fr

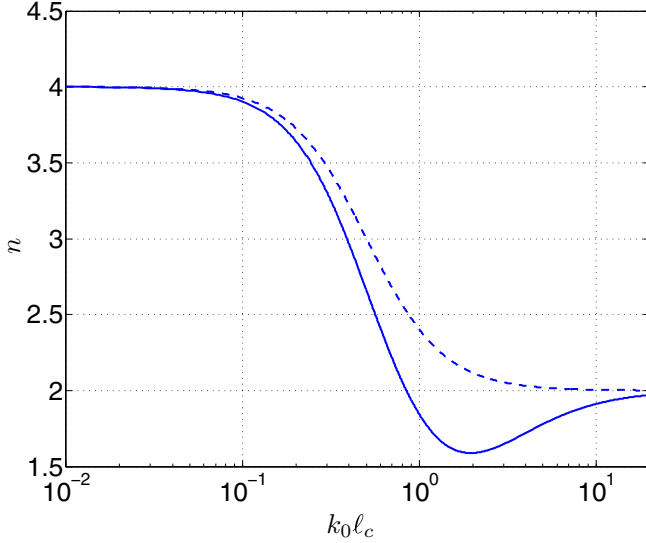


FIG. 3. (Color online) Characteristic exponent n obtained with (solid line) and without (dashed line) the additional terms in the self-energy as a function of dimensionless frequency $k_0 \ell_c$ for an exponentially correlated disorder.

However, as to the velocity of the coherent field, the effect of the additional terms is very limited since the real part of the wave vector is

$$\text{Re}(k_{\text{eff}}) \approx k_0 - \frac{\text{Re}[\tilde{\Sigma}(k_0)]}{2k_0}, \quad (54)$$

which will always remain close to k_0 within the on-shell approximation.

V. CONCLUSION

Starting from the wave equation for the acoustic pressure in an heterogeneous and nondissipative fluid, we have calculated the coherent wave, taking into account spatial variations of both density and compressibility such that the relevant random potential contains both a scalar and an operator part, α and β . The calculation is based on the diagrammatic approach of multiple scattering, within Bourret and on-shell approximations. Interestingly, the results show that discarding the random operator term [as is usually done when treating the problem as Helmholtz' equation with a space-dependent wave speed $c(\vec{r})$] amounts to overestimate the scattering mean free path by up to a factor of 13/3 when the fluctuation of α and β have similar magnitude. The error is particularly large at low frequencies, when the correlation length is comparable to or smaller than the wavelength. The theoretical analysis has been conducted in two and three dimensions and validated by numerical experiments. Though the results presented here are theoretical and rather academic, we believe they are of importance for all practical applications involving multiple scattering of acoustic waves (e.g., characterization of inhomogeneous media). Moreover, from a theoretical point of view, the scattering mean free path is the basic ingredient to describe universal wave phenomena in complex media, such as coherent backscattering, ballistic-to-diffuse transition,

radiative transport of energy, etc. It is therefore crucial to determine it properly.

ACKNOWLEDGMENTS

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APPENDIX A:

1. 3D calculations

Assuming that the correlation functions $C_{\alpha\alpha}$, $C_{\alpha\beta}$, and $C_{\beta\beta}$ are identical, we have

$$\begin{aligned} \Sigma(x) &= k_0^4 G_0 C - k_0^2 G_0 [C'' + 2C'/x] \\ &\quad - [C'' G_0'' + C''' G_0' + 2C'' G_0'/x] \\ &= G_0 [k_0^4 C - 2k_0^2 C'/x + (1/x - jk_0) C''']. \end{aligned} \quad (A1)$$

Given the radial symmetry, the 3D Fourier transform of Σ is

$$\tilde{\Sigma}(k_0) = \frac{4\pi}{k_0} \int_0^\infty \Sigma(x) \sin(k_0 x) x dx. \quad (A2)$$

Hence the imaginary part:

$$\begin{aligned} \text{Im } \tilde{\Sigma}(k_0) &= \frac{1}{k_0} \int_0^\infty [k_0^4 C - 2k_0^2 C'/x + C'''/x] \sin^2(k_0 x) dx \\ &\quad - \frac{1}{2} \int_0^\infty C''' \sin(2k_0 x) dx. \end{aligned} \quad (A3)$$

In order to study its behavior in the low-frequency regime ($k_0 x \rightarrow 0$) a Taylor expansion of the sines up to the sixth order followed by integrations by parts are performed. It yields

$$\text{Im } \tilde{\Sigma}(k_0) \rightarrow k_0^5 \int_0^\infty \frac{13}{3} x^2 C(x) dx. \quad (A4)$$

If the additional terms due to the random operator are neglected, Eq. (A3) reduces to

$$\text{Im } \tilde{\Sigma}(k_0) = \frac{1}{k_0} \int_0^\infty k_0^4 C \sin^2(k_0 x) dx, \quad (A5)$$

$$\rightarrow k_0^5 \int_0^\infty x^2 C(x) dx. \quad (A6)$$

As a consequence, in the low-frequency limit $k_0 \ell_c \rightarrow 0$, the ratio of the mean free path calculated with (ℓ_s) or without [$\ell_s^{(\alpha\alpha)}$] the additional terms is 13/3. This ratio does not depend on the precise shape of the correlation function $C(x)$, as long as its second-order moment is finite.

The final results given and plotted in the paper were established for an exponentially correlated disorder. In the Gaussian case where $C(x) = \sigma^2 \exp(-x^2/\ell_c^2)$, we obtain

$$\begin{aligned} \frac{\tilde{\Sigma}(k_0)}{k_0^2} &= \sqrt{\pi} \frac{\sigma^2}{4} \left[j k_0 \ell_c (9E - 1) + \frac{8}{\sqrt{\pi}} + \frac{1}{k_0 \ell_c} 4j(3E - 1) \right. \\ &\quad \left. + \frac{1}{(k_0 \ell_c)^2} \frac{8}{\sqrt{\pi}} + \frac{8j(E - 1)}{(k_0 \ell_c)^3} \right]. \end{aligned} \quad (A7)$$

In the expression above, we have introduced a dimensionless constant E :

$$E = [1 + \operatorname{erf}(jk_0\ell_c)]e^{-k_0^2\ell_c^2}. \quad (\text{A8})$$

If the additional terms are neglected, then we have

$$\frac{\tilde{\Sigma}(k_0)}{k_0^2} = \sqrt{\pi} \frac{\sigma^2}{4} [jk_0\ell_c(E - 1)]. \quad (\text{A9})$$

For the sake of simplicity, the ratio $\ell_s/\ell_s^{(\alpha\alpha)}$ has not been plotted in the Gaussian case, but its general trend is very similar to the exponential case.

2. 2D calculations

In 2D space, we have:

$$G_0(\vec{r} - \vec{r}_s) = \frac{-i}{4} H_0^{(1)}(k_0|\vec{r} - \vec{r}_s|), \quad (\text{A10})$$

$$\langle G(\vec{r} - \vec{r}_s) \rangle = \frac{-i}{4} H_0^{(1)}(k_{\text{eff}}|\vec{r} - \vec{r}_s|). \quad (\text{A11})$$

$H_0^{(1)}(x)$ is the Hankel function of the first kind and of order 0. We still have four contributions to the self-energy [see Eq. (36)]. Assuming circular symmetry, with $x = |\vec{r}_a - \vec{r}_b|$ we have:

$$\begin{aligned} \Sigma_{\alpha\alpha}(x) &= k_0^4 G_0(x) C_{\alpha\alpha}(x) \\ \Sigma_{\beta\alpha}(x) &= k_0^2 \vec{\nabla} G_0(x) \cdot \vec{\nabla} C_{\beta\alpha}(x) \\ \Sigma_{\alpha\beta}(x) &= -k_0^2 \vec{\nabla} \cdot [G_0(x) \vec{\nabla} C_{\alpha\beta}(x)] \\ \Sigma_{\beta\beta}(x) &= -\vec{\nabla} \cdot [\vec{\nabla} \otimes \vec{\nabla} C_{\beta\beta}(x) \vec{\nabla} G_0(x)]. \end{aligned}$$

The difference between the 2D and 3D cases lie in the expressions of the gradient, divergence, and Hessian of a function with circular (or spherical) symmetry. In particular, $\Sigma_{\beta\beta}$ requires the Hessian of a circularly symmetric function in polar coordinates:

$$\vec{\nabla} \otimes \vec{\nabla} C_{\beta\beta} = \begin{bmatrix} \frac{\partial^2 C_{\beta\beta}}{\partial x^2} & 0 \\ 0 & \frac{1}{x} \frac{\partial C_{\beta\beta}}{\partial x} \end{bmatrix}. \quad (\text{A12})$$

As a whole, in 2D the expression of the self-energy [equivalent of Eq. (50) in 3D] reads:

$$\begin{aligned} \Sigma(x) &= k_0^4 G_0(x) \sigma_\alpha^2 c_{\alpha\alpha}(x) \\ &\quad - k_0^2 G_0(x) \sigma_\alpha \sigma_\beta \left[c_{\alpha\beta}''(x) + \frac{1}{x} c_{\alpha\beta}'(x) \right] \\ &\quad - \sigma_\beta^2 \left\{ G_0'(x) c_{\beta\beta}'''(x) + \left[G_0''(x) + \frac{1}{x} G_0'(x) \right] c_{\beta\beta}''(x) \right\}. \end{aligned} \quad (\text{A13})$$

Using Eq. (A10) along with differentiation and recurrence properties for Bessel and Hankel functions (see Ref. [24],

p. 361), it is straightforward to obtain:

$$\begin{aligned} \Sigma(x) &= G_0(x) \left\{ k_0^4 \sigma_\alpha^2 c_{\alpha\alpha}(x) \right. \\ &\quad \left. - k_0^2 \sigma_\beta \left[\sigma_\alpha c_{\alpha\beta}''(x) + \frac{1}{x} \sigma_\alpha c_{\alpha\beta}'(x) - \sigma_\beta c_{\beta\beta}''(x) \right] \right\} \\ &\quad - G_0'(x) \sigma_\beta^2 c_{\beta\beta}'''(x). \end{aligned} \quad (\text{A14})$$

And for identical correlation functions $C_{\alpha\alpha}$, $C_{\alpha\beta}$, and $C_{\beta\beta}$, we have:

$$\Sigma(x) = \sigma^2 G_0(x) \left[k_0^4 c(x) - \frac{k_0^2}{x} c'(x) \right] - \sigma^2 G_0'(x) c'''(x). \quad (\text{A15})$$

Once an analytical expression for $\Sigma(x)$ is obtained, we have to calculate its spatial Fourier transform in order to determine the effective wave number. In 2D, the Fourier transform of a circularly symmetric function is the zero-order Hankel transform:

$$\tilde{\Sigma}(k) = 2\pi \int_0^\infty \Sigma(x) J_0(kx) x dx, \quad (\text{A16})$$

where $J_0(x)$ is the cylindrical Bessel function of order 0. Calculating the mean free path $\ell_s = 1/[2 \operatorname{Im}(k_{\text{eff}})]$ amounts to numerically evaluating three integrals:

$$\tilde{\Sigma}(k) = \frac{-i\pi}{2} (k_0^4 I_1 - k_0^2 I_2 + k_0 I_3), \quad \text{where} \quad (\text{A17})$$

$$I_1 = \int_0^\infty C(x) H_0^{(1)}(k_0 x) J_0(kx) x dx, \quad (\text{A18})$$

$$I_2 = \int_0^\infty C'(x) H_0^{(1)}(k_0 x) J_0(kx) dx, \quad (\text{A19})$$

$$I_3 = \int_0^\infty C'''(x) H_1^{(1)}(k_0 x) J_0(kx) x dx, \quad (\text{A20})$$

whatever the shape of the correlation function $C(x)$.

In the low-frequency regime ($k_0 x \rightarrow 0$), a Taylor expansion of the Bessel functions followed by integrations by parts yield

$$\operatorname{Im} \tilde{\Sigma}(k_0) \rightarrow -\frac{\pi}{2} k_0^4 \int_0^\infty \frac{9}{2} x C(x) dx. \quad (\text{A21})$$

If the additional terms due to the random operator are neglected (i.e., $I_2 = I_3 = 0$), then Eq. (A20) reduces to

$$\operatorname{Im} \tilde{\Sigma}(k_0) = \frac{4\pi}{k_0} \int_0^\infty k_0^4 C \sin^2(k_0 x) dx, \quad (\text{A22})$$

$$\rightarrow -\frac{\pi}{2} k_0^4 \int_0^\infty x C(x) dx. \quad (\text{A23})$$

Hence, in the low-frequency limit $k_0\ell_c \rightarrow 0$, the ratio of the 2D mean free paths calculated with (ℓ_s) or without [$\ell_s^{(\alpha\alpha)}$] the additional terms is 9/2, as opposed to 13/3 in 3D. And again, this ratio does not depend on the precise shape of the correlation function $C(x)$, as long as its second-order moment is finite.

3. Numerical simulations

The acoustic wave propagation in heterogeneous media is numerically simulated with Simsonic, a 3D Cartesian FDTD approach to solve the elastodynamic equations.

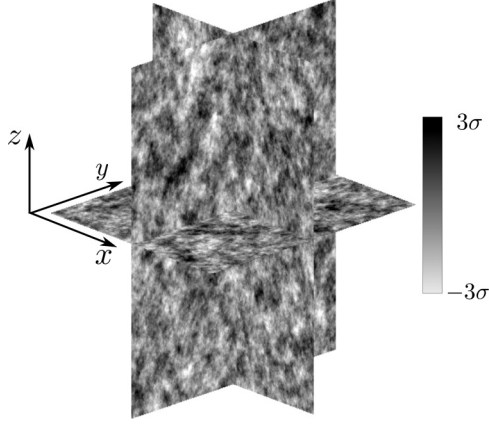


FIG. 4. Example of an exponentially correlated disorder with Gaussian statistics (zero mean, variance σ^2) in 3D space.

Three-dimensional maps of the local wave speed and mass density can be designed by the user. These maps define the propagation medium at each grid point. By properly filtering a 3D white noise of $N_x \times N_y \times N_z$ points with Gaussian statistics, it is possible to build a 3D map exhibiting an exponentially correlated disorder:

$$C_{\alpha\alpha}(r) = \sigma^2 \exp\left[-\frac{r}{\ell_c}\right], \quad (\text{A24})$$

where $r = \sqrt{x^2 + y^2 + z^2}$ is the radial coordinate, ℓ_c the correlation length, and σ^2 the variance. An example of one realization of the media is given in Fig. 4. Values below -3σ or above 3σ are truncated. Various uncorrelated realizations of disorder can be obtained by repeating the procedure. The same map has been employed for α and β , so $\alpha(\vec{r}) = \beta(\vec{r})$ for each realization of disorder. In that case, we have $C_{\alpha\alpha} = C_{\beta\beta} = C_{\alpha\beta}$ which corresponds to the theoretical example detailed in the paper. Using independent, or partially correlated maps, or with different variances or correlation lengths for α and β could also be possible to investigate all possibilities.

The correlation length was set at $\ell_c = 0.240$ mm, so $k_0\ell_c = 1$ for a driving frequency $f = 1$ MHz. The variance σ^2 ranged between 1% and 4%. Various simulations were carried out in order to calculate the scattering mean free paths for frequencies in the range $k_0\ell_c \in [0.75, 3]$ by changing the central frequency of the incoming pulse between 1 and 2 MHz. In order to avoid an additional numerical dissipation of the acoustic energy, it is important to resolve both the correlation length ℓ_c and the wavelength λ with at least 10 grid points in all directions. Furthermore, the CFL (Courant-Friedrichs-Lewy) condition is to be respected based on the maximum propagation speed $c_{\max} = c_0/\sqrt{1 - 3\sigma}$ in the medium. Perfectly matched layers (PML) were implemented outside the scattering region to ensure absorbing boundary conditions. Typically more than 80 Go of RAM were required and a multithreaded parallel version of SIMSONIC3D (OPENMP) was needed to perform these large-scale simulations. As a typical example, a snapshot of the propagating wave front in the (y, z) plane is given in Fig. 5.

Prior to an ensemble average of the acoustic pressure field over N realizations of disorder, p is first spatially averaged along the (x, y) plane (under the hypothesis of spatial ergodicity) to obtain $\bar{p}_i(z, t)$. Then the final mean-field

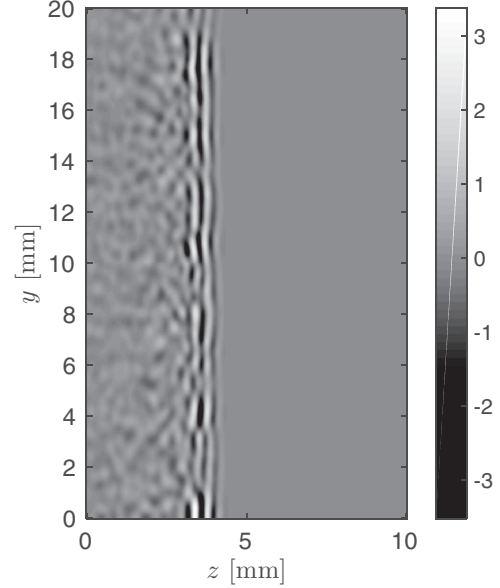


FIG. 5. A pulsed plane wave is generated at $z = 0$ mm and propagates along the z axis. The snapshot is taken at $t = 4 \mu\text{s}$ in the $x = 0$ plane. The resulting pressure is in arbitrary units.

estimator reads:

$$\langle p(z, t) \rangle = \frac{1}{N} \sum_{i=1}^N \bar{p}_i(z, t). \quad (\text{A25})$$

In all simulations we ensured N to be large enough in order that Eq. (A25) represents an accurate estimator of the mean coherent pressure field and that remaining random fluctuations can be neglected.

From the mean pressure field we can calculate the acoustic intensity in the frequency domain, $|\langle P(z, \omega) \rangle|^2$. As seen in Fig. 6, the acoustic intensity decays exponentially; a linear fit of its logarithm gives an estimation of the mean free paths

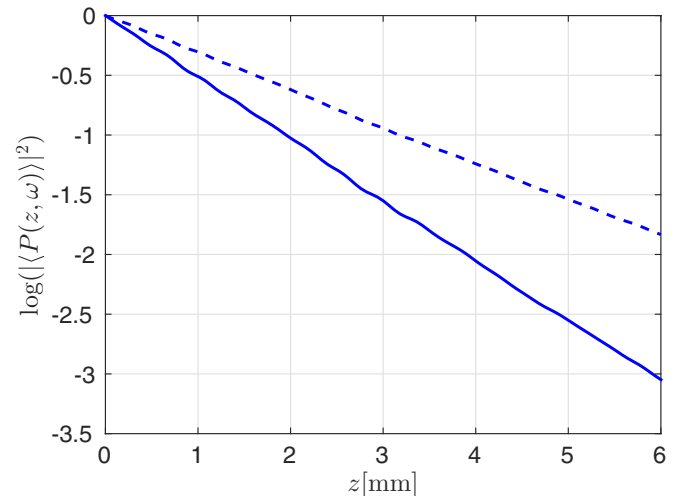


FIG. 6. (Color online) Decay of $\log |\langle P(z, \omega) \rangle|^2$ versus distance z at $f = 2$ MHz and $\sigma = 0.2$. A linear fit of this data gives an estimation of the scattering mean free path in the scalar case (blue dashed line) and in the operator case (blue solid line).

$[\ell_s$ or $\ell_s^{(\alpha\alpha)}$] at a given frequency. In all simulations, we ensured that the propagation distance was at least 3 times larger than the scattering mean free path, so the decay of intensity is significant.

4. Fluid limit of the solid model

In Ref. [20], the self-energy is expressed in terms of fluctuations of mass density ρ and Lamé coefficients λ and

μ . The liquid limit is taken by setting $\mu = 0$ and $\chi = 1/\lambda$. Since the fluctuations of all parameters relative to their mean are assumed to be very small, Eqs. (6) and (27) can be differentiated to obtain linear relations between the two pairs of variables. This leads to $\sigma_\beta = \sigma_\rho$ and $\sigma_\lambda^2 = \langle(\alpha + \beta)^2\rangle$. In the scalar case, $\sigma_\beta = 0$, then $\sigma_\rho = 0$ and $\sigma_\alpha = \sigma_\lambda$. If the operator term is taken into account and $\alpha(\vec{r}) = \beta(\vec{r})$, then $\sigma_\beta = \sigma_\rho = \sigma$ and $\sigma_\alpha = \sigma_\lambda/2 = \sigma$. From Eq. (37) in Ref. [20], we infer

$$\frac{1}{k_0\ell_s^{(\alpha\alpha)}} = \sigma^2(k_0\ell_c)^3 \int_{-1}^1 \frac{dx}{[1 + 2(k_0\ell_c)^2(1-x)]^2} = \sigma^2 \frac{2(k_0\ell_c)^3}{1 + 4(k_0\ell_c)^2} \quad (\text{A26})$$

in the scalar case and

$$\begin{aligned} \frac{1}{k_0\ell_s} &= \sigma^2(k_0\ell_c)^3 \int_{-1}^1 \frac{(x+2)^2}{[1 + 2(k_0\ell_c)^2(1-x)]^2} dx \\ &= \sigma^2 \frac{1 + 8(k_0\ell_c)^2 + 18(k_0\ell_c)^4}{k_0\ell_c(1 + 4(k_0\ell_c)^2)} - \sigma^2 \frac{[1 + 10(k_0\ell_c)^2 + 24(k_0\ell_c)^4] \log[1 + 4(k_0\ell_c)^2]}{4(k_0\ell_c)^3[1 + 4(k_0\ell_c)^2]} \end{aligned} \quad (\text{A27})$$

in the operator case. Equation (A26) is exactly our result in the scalar case [see Eq. (20)]. However, in the operator case,

Eq. (A27) and Eq. (52) disagree, except at zero frequency, as was discussed earlier and shown in Fig. 1.

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