

Microscopic analysis of sound attenuation in low-temperature amorphous solids reveals quantitative importance of non-affine effects

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Sound attenuation in low temperature amorphous solids originates from their disordered structure. However, its detailed mechanism is still being debated. Here we analyze sound attenuation starting directly from the microscopic equations of motion. We derive an exact expression for the zero-temperature sound damping coefficient. We verify that the sound damping coefficients calculated from our expression agree very well with results from independent simulations of sound attenuation. The small wavevector analysis of our expression shows that sound attenuation is primarily determined by the non-affine displacements' contribution to the sound wave propagation coefficient coming from the frequency shell of the sound wave. Our expression involves only quantities that pertain to solids' static configurations. It can be used to evaluate the low temperature sound damping coefficients without directly simulating sound attenuation.

I. INTRODUCTION

The physics of sound attenuation in amorphous solids is drastically different than in crystalline solids. At low temperatures, when thermal effects can be neglected, sound is attenuated due to the inherent disorder of amorphous solids, whereas the attenuation is absent in crystalline solids. To understand the physical mechanism behind sound attenuation one can examine its wavevector k dependence. Sound attenuation in amorphous solids has a complicated dependence on the wavevector¹, but small wavevector k^4 scaling of sound damping coefficients has long been conjectured on an experimental basis^{2,3}. An initial interpretation was that this small wavevector behavior originates from Rayleigh scattering of sound waves from the solid's inhomogeneities. Recent computer simulations⁴⁻⁶ verified that in classical three-dimensional zero-temperature amorphous solids at the smallest wavevectors sound damping coefficients scale as k^4 , although a logarithmic correction to this scaling was also claimed⁷.

The specific physical mechanism of sound attenuation in low temperature amorphous solids is still debated. Zeller and Pohl² obtained the Rayleigh scattering law using an "isotopic scattering"³ model in which every atom of the glass is an independent source of scattering. Several recent experimental and simulational results were analyzed within the framework of the fluctuating elasticity theory of Schirmacher⁸⁻¹⁰. This theory posits that an amorphous solid can be modeled as a continuous medium with spatially varying elastic constants. The inhomogeneity of the elastic constants causes sound scattering and attenuation. In the limit of the wavelength being much larger than the characteristic spatial scale of the inhomogeneity this mechanism is equivalent to Rayleigh scattering and the theory predicts that sound damping

coefficients scale with the wavevector as k^4 . If the elastic constant variations have slowly decaying, power-law-like correlations, the theory predicts a logarithmic correction to Rayleigh scattering^{7,15}. Other physical approaches, *e.g.* local oscillator^{9,11-14} and random matrix¹⁶⁻¹⁸ models, can also be used to derive the Rayleigh scattering law. For this reason, Rayleigh scaling cannot serve to distinguish between different models⁹, and other model predictions must be used to determine the mechanism behind sound attenuation.

Three recent studies came to very different conclusions regarding the applicability of the fluctuating elasticity theory for sound attenuation. First, Caroli and Lemaître¹⁹ analyzed a version of the theory derived from microscopic equations of motion. They obtained *all* the parameters needed to calculate sound attenuation from the theory from the same simulations that were used to test the theoretical predictions. Caroli and Lemaître showed that this version of the theory underestimates sound damping coefficients by about two orders of magnitude.

Second, Kapteijns *et al.*²⁰ analyzed the dependence of sound attenuation in a two-dimensional glass on a parameter δ , which "resembles" changing the stability of the amorphous solid. To calculate the disorder parameter⁸ of the fluctuating elasticity theory they replaced fluctuations of local elastic constants (which are used in the theoretical description) by the sample-to-sample fluctuations of bulk elastic constants. In this way they were able to sidestep the issue of the definition of local elastic constants²¹ and of the correlation volume. While Kapteijns *et al.* showed that the disorder parameter and the sound damping coefficient have the same dependence on δ , they left the calculation of the pre-factor for the scaling for further research.

Finally, Mahajan and Pica Ciamarra²² argued that sound attenuation is proportional to the square of the disorder parameter γ according to a version of fluctuating elasticity theory that incorporates an elastic correlation length^{9,23}. They relied upon a relation between the bo-

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son peak, the speed of sound, and an elastic correlation length to show that the speed of sound and the boson peak frequency can be used to infer the change of the sound damping coefficient. Again, the magnitude of the sound damping coefficient was not addressed.

The results described above show that it is difficult to distinguish between and to validate different semi-phenomenological models invoked to describe sound attenuation in zero-temperature amorphous solids. One of the reasons is that most of these approaches involve an adjustable parameter (or parameters) and therefore are able to predict trends rather than absolute values of sound damping coefficients. For example, neither Kapteijns *et al.* nor Mahajan and Pica Ciamarra calculated the values of sound damping coefficients (in contrast to Caroli and Lemaître), but rather investigated the variation of the sound attenuation between different glasses. Limited range of glasses that can be created *in silico* makes it difficult to distinguish between trends predicted by different models or different versions of a model.

Our goal is to understand the microscopic origin of the sound attenuation. We derive an exact expression for the sound damping coefficient in terms of quantities that can be calculated from static configurations of amorphous solids, without the need to directly simulate sound attenuation. Our expression is analogous to well-known Green-Kubo formulae for transport coefficients²⁴. The latter expressions allow one to calculate transport coefficients without explicitly simulating transport processes. While both our expression and Green-Kubo formulae need to be evaluated numerically, they can also serve as starting points for approximate analyses and treatments that can shed light on the validity of semi-phenomenological models. We hope that the results of one such analysis, which we present at the end of the paper, can inspire new models or be incorporated into the existing ones.

In Sec. II we start from the microscopic equations of motion for harmonic vibrations. We derive an exact equation of motion for an auto-correlation function that has been used to determine sound attenuation. We identify the self-energy and show that its real part reproduces the non-Born contribution to the zero-temperature wave propagation coefficients. The imaginary part of the self-energy is the origin of sound attenuation. We show that sound damping coefficients calculated this way agree very well with those obtained from direct simulations of sound attenuation in zero-temperature glasses with different stability. In Sec. III we present the small wavevector expansion of our expression for the sound damping coefficient. It shows that the limiting k^4 sound attenuation originates from the same physics as the non-Born contribution to the elastic constants and wave propagation coefficients, *i.e.* from the forces inducing non-affine displacements, which appear due to the amorphous solids' disordered structure. More precisely, attenuation of the sound wave is primarily determined by the contribution to the non-Born part of the wave propagation coefficient

from a shell of frequencies around the frequency of the sound wave. We thus show the common origin of the renormalization of the elastic constants and of sound attenuation. In Sec. IV we discuss the results of an approximate evaluation of our expression for the sound damping coefficient which assumes that the exact eigenvectors of the Hessian matrix can be replaced by plane waves. These results allow us to critically evaluate the relation between our exact expression and the fluctuating elasticity theory. We end the paper with a discussion of our results and related descriptions of the sound attenuation.

II. MICROSCOPIC ANALYSIS OF SOUND ATTENUATION

We start from microscopic equations of motion for small displacements of N spherically symmetric particles of unit mass comprising our model amorphous solid,

$$\partial_t^2 \mathbf{u}_i = - \sum_j \mathcal{H}_{ij} \cdot \mathbf{u}_j. \quad (1)$$

Here \mathbf{u}_i is the displacement of the i th particle from its inherent structure (potential energy minimum) position \mathbf{R}_i and \mathcal{H} is the Hessian calculated at the inherent structure,

$$\mathcal{H}_{ij} = \sum_{l \neq i} \frac{\partial^2 V(R_{il})}{\partial \mathbf{R}_i \partial \mathbf{R}_j} \quad (2)$$

where $V(r)$ is the pair potential and \mathcal{H}_{ij} is a 3x3 tensor.

To derive an expression for the sound damping coefficient we use a slightly modified procedure proposed by Gelin *et al.*⁷. We assume that at $t = 0$ the particles are displaced from their equilibrium positions according to $\mathbf{u}_i(t = 0) = \hat{\mathbf{e}} \exp(-i\mathbf{k} \cdot \mathbf{R}_i)$, $\dot{\mathbf{u}}_i(t = 0) = 0$, where $\hat{\mathbf{e}}$ is a unit vector and wavevector \mathbf{k} is one of the wavevectors allowed by periodic boundary conditions. We then analyze the time dependence of the auto-correlation function of the single-particle displacement averaged over the whole system, $C(t) = N^{-1} \sum_i \mathbf{u}_i^*(t = 0) \cdot \mathbf{u}_i(t)$. We anticipate that in the limit of small wavevectors \mathbf{k} the auto-correlation function will exhibit damped oscillations, $C(t) \propto \cos(vkt) \exp(-\Gamma(k)t/2)$, and we will identify v as the speed of sound and $\Gamma(k)$ as the damping coefficient.

Solving Eqs. (1) with our initial conditions is equivalent to solving the following equations

$$\partial_t^2 \mathbf{a}_i = - \sum_j \mathcal{H}_{ij}(\mathbf{k}) \cdot \mathbf{a}_j, \quad (3)$$

where $\mathcal{H}(\mathbf{k})$ is the wavevector-dependent Hessian, $\mathcal{H}_{il}(\mathbf{k}) = \mathcal{H}_{il} \exp[i\mathbf{k} \cdot (\mathbf{R}_i - \mathbf{R}_l)]$, with initial conditions $\mathbf{a}_i(t = 0) = \hat{\mathbf{e}}$, $\dot{\mathbf{a}}_i(t = 0) = 0$. In terms of the new variables, $C(t) = N^{-1} \sum_i \mathbf{a}_i^*(t = 0) \cdot \mathbf{a}_i(t)$.

To analyze $C(t)$ we use the standard projection operator approach²⁵. First, we define a scalar product of two displacement vectors, \mathbf{a}_i and \mathbf{b}_j , $i, j = 1, \dots, N$, $\langle a|b \rangle = \sum_i \mathbf{a}_i^* \cdot \mathbf{b}_i$. Next, we define a unit vector $|1\rangle$ with

components $\mathbf{1}_i = N^{-1/2}\hat{\mathbf{e}}_i$, and projection operator \mathcal{P} on the unit vector, $\mathcal{P} = |\mathbf{1}\rangle\langle\mathbf{1}|$ and orthogonal projection \mathcal{Q} , $\mathcal{Q} = \mathcal{I} - |\mathbf{1}\rangle\langle\mathbf{1}|$ where \mathcal{I} is the identity matrix.

Using the projection operator approach we obtain the following expression for the Fourier transform $C(\omega) = \int_0^\infty C(t) \exp(i(\omega + i\epsilon)t) dt$ of the displacement auto-correlation function,

$$C(\omega) = \frac{i(\omega + i\epsilon)}{(\omega + i\epsilon)^2 - \langle\mathbf{1}|\mathcal{H}(\mathbf{k})|\mathbf{1}\rangle + \Sigma(\mathbf{k};\omega)}, \quad (4)$$

where the self-energy $\Sigma(\mathbf{k};\omega)$ reads

$$\Sigma(\mathbf{k};\omega) = \left\langle\mathbf{1}\left|\mathcal{H}(\mathbf{k})\mathcal{Q}\frac{1}{-\omega + i\epsilon + \mathcal{Q}\mathcal{H}(\mathbf{k})\mathcal{Q}}\mathcal{Q}\mathcal{H}(\mathbf{k})\right|\mathbf{1}\right\rangle. \quad (5)$$

Equations (4-5) are exact. While it is straightforward to calculate $\langle\mathbf{1}|\mathcal{H}(\mathbf{k})|\mathbf{1}\rangle$, evaluation of the self-energy requires inversion of a large-dimensional matrix for each allowed wavevector. To make the numerical effort manageable, in the denominator in Eq. (5) we approximate $\mathcal{H}(\mathbf{k})$ by \mathcal{H} . As argued in Appendix A, this approximation does not influence the small wavevector dependence of the sound damping coefficients.

In the small wavevector limit, the first non-trivial term in the denominator in Eq. (4), $\langle\mathbf{1}|\mathcal{H}(\mathbf{k})|\mathbf{1}\rangle$, can be expressed in terms of the Born contributions to the zero-temperature wave propagation coefficients²⁶,

$$\langle\mathbf{1}|\mathcal{H}(\mathbf{k})|\mathbf{1}\rangle = \rho^{-1}\hat{\mathbf{e}}_\alpha A_{\alpha\beta\gamma\delta}^{\text{Born}} k_\beta \hat{\mathbf{e}}_\gamma k_\delta + o(k^2) \quad (6)$$

where $\rho = N/V$ is the number density, Greek indices denote Cartesian components, the Einstein summation convention for Greek indices is hereafter adopted, and $A_{\alpha\beta\gamma\delta}^{\text{Born}}$ is the Born wave propagation coefficient, which can be expressed as the average of the local Born wave propagation coefficients $A_{j,\alpha\beta\gamma\delta}^{\text{Born}}$,

$$A_{j,\alpha\beta\gamma\delta}^{\text{Born}} = \frac{\rho}{2} \sum_{l \neq j} \frac{\partial^2 V(R_{jl})}{\partial R_{j,\alpha} \partial R_{j,\gamma}} R_{jl,\beta} R_{jl,\delta} \quad (7)$$

over the whole system,

$$A_{\alpha\beta\gamma\delta}^{\text{Born}} = N^{-1} \sum_j A_{j,\alpha\beta\gamma\delta}^{\text{Born}}. \quad (8)$$

For example, if the coordinate system is chosen such that $\hat{\mathbf{e}}$ is in the x direction, and we are interested in a transverse wave and choose \mathbf{k} in the y direction, then the right hand side of (6) becomes $\rho^{-1} A_{xyxy}^{\text{Born}} k^2$.

In the absence of the self-energy term, (4) predicts the Born value of the speed of sound and no sound damping. Both the renormalization of the sound speed and the sound attenuation originate from the self-energy.

The self-energy can be calculated using the eigenvalues and eigenvectors of the Hessian. In the thermodynamic limit²⁷, when the spectrum of the Hessian becomes continuous, we can use the Plemelj identity to identify the

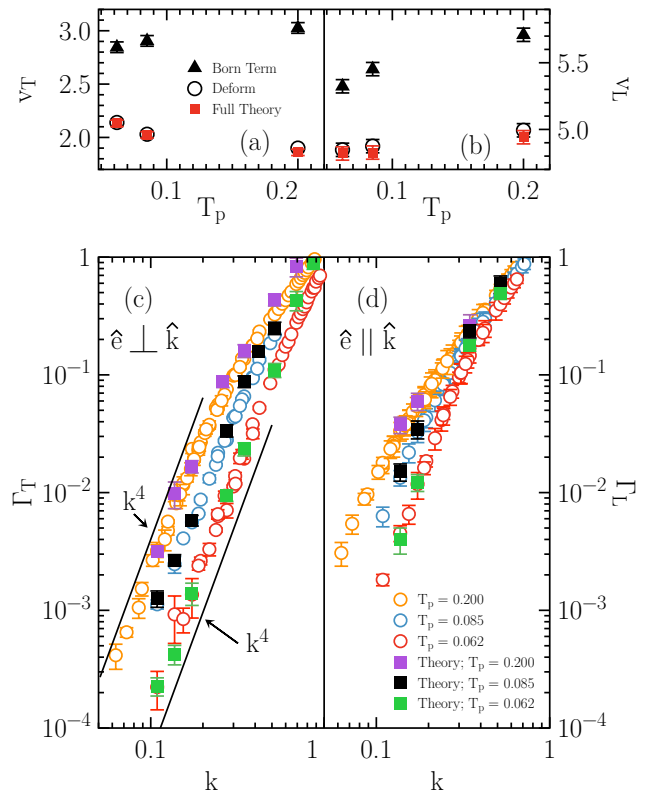


FIG. 1. Upper panels: the transverse (a) and longitudinal (b) speed of sound obtained from the theory, Eq. (13), (red squares) and calculated from the elastic constants (black circles) as a function of parent temperature T_p . The black triangles are the Born values of the speed of sound. Lower panels: the transverse (c) and longitudinal (d) damping coefficients obtained from the theory, Eq. (14), (squares) and obtained from sound attenuation simulations⁵ (circles) for different parent temperatures. Rayleigh scaling $\Gamma \propto k^4$ is recovered at small wavevectors. The error bars for the theoretical calculation represent the uncertainty due to using different bin sizes.

imaginary component of the self-energy, which is responsible for sound attenuation. The real $\Sigma'(\mathbf{k};\omega)$ and imaginary $\Sigma''(\mathbf{k};\omega)$ parts of the self-energy read,

$$\Sigma'(\mathbf{k};\omega) = \mathcal{P} \int d\Omega \Upsilon(\mathbf{k},\Omega) (\Omega^2 - \omega^2)^{-1}, \quad (9)$$

$$\Sigma''(\mathbf{k};\omega) = \frac{\pi}{2\omega} \Upsilon(\mathbf{k},|\omega|), \quad (10)$$

where \mathcal{P} denotes the Cauchy principal value. The function $\Upsilon(\mathbf{k},\Omega)$ is defined through the sum over eigenvectors \mathcal{E}^p of the Hessian matrix with non-zero²⁸ eigenvalues Ω_p^2 such that $\Omega_p \in [\Omega, \Omega + d\Omega]$, where $d\Omega$ is the bin size,

$$\Upsilon(\mathbf{k},\Omega) = (1/d\Omega) \sum_{\Omega_p \in [\Omega, \Omega + d\Omega]} |\langle\mathbf{1}|\mathcal{H}(\mathbf{k})\mathcal{Q}|\mathcal{E}^p\rangle|^2. \quad (11)$$

The key conceptual issue in writing Eq. (11) (and closely related equations (18,21)) is that the thermodynamic

limit is implied for the expression at the right-hand-side. In this limit the spectrum becomes dense and phonon bands are not distinguishable. Thus, to calculate $\Upsilon(\mathbf{k}, \Omega)$ from the analysis of finite-size simulations we need to choose bin size $d\Omega$ such that phonon bands are not resolved. In the numerical calculations described below we tried a few bin sizes between 0.1 and 0.2 and found that within this range the results were not very sensitive to the bin size.

To evaluate the displacement auto-correlation function we need to find complex poles of the denominator at the right-hand-side of Eq. (4). In the small wavevector limit this can be done perturbatively, using k as the small parameter. This leads to the following pair of poles,

$$\omega_{\pm} = \pm vk - i\Sigma''(\mathbf{k}; vk)/(2vk) \quad (12)$$

where the renormalized speed of sound v is given by

$$v^2 = \lim_{k \rightarrow 0} k^{-2} [\langle 1|\mathcal{H}(\mathbf{k})|1\rangle - \Sigma'(\mathbf{k}; 0)]. \quad (13)$$

The last term in Eq. (12) is our main result. It says that the sound damping in zero-temperature amorphous solids is determined by $\Upsilon(\mathbf{k}, \Omega)/\Omega^2$ calculated at the wave's frequency, $\Omega = vk$,

$$\Gamma(k) = \frac{\Sigma''(\mathbf{k}; vk)}{vk} = \frac{\pi}{2} \frac{\Upsilon(\mathbf{k}, vk)}{(vk)^2}. \quad (14)$$

We emphasize that $\Upsilon(\mathbf{k}, \Omega)/\Omega^2$ is the same function that, after integration over the whole frequency spectrum, determines the renormalization of the wave propagation coefficients. Note that v , $\Gamma(k)$, and related quantities defined below depend on the angle between the polarization of the initial condition $\hat{\mathbf{e}}$ and the direction of the wavevector $\hat{\mathbf{k}}$.

To verify Eqs. (13-14) we calculated v and $\Gamma(k)$ for model zero-temperature glasses analyzed in Ref.⁵. These glasses were obtained by instantaneously quenching supercooled liquids equilibrated using the swap Monte Carlo algorithm²⁹ at different parent temperatures T_p to their inherent structures using the fast inertial relaxation engine minimization³⁰. The glasses consist of spherically symmetric, polydisperse particles which interact via a potential $\propto r^{-12}$, with a smooth cutoff, see Appendix B and Refs.^{5,29} for details. The parent temperature controls the glass's stability and thus its properties^{5,31,32}.

We calculated eigenvalues and eigenvectors of the Hessian using ARPACK³³ and Intel Math Kernel Library³⁴. Then, using Eqs. (13-14) we evaluated v and $\Gamma(k)$ for the longitudinal, $\hat{\mathbf{e}} \parallel \hat{\mathbf{k}}$, and the transverse, $\hat{\mathbf{e}} \perp \hat{\mathbf{k}}$, sound. To calculate $\Upsilon(\mathbf{k}, \Omega)$ one chooses a \mathbf{k}_n compatible with periodic boundary conditions. Then one calculates $|\langle 1|\mathcal{H}(\mathbf{k}_n)\mathcal{Q}|\mathcal{E}^p\rangle|$ and bins the results according to the square root of the eigenvalue of $|\mathcal{E}^p\rangle$ to determine $\Upsilon(\mathbf{k}_n, \Omega)$. Note that Ω_p^2 is the eigenvalue corresponding to $|\mathcal{E}^p\rangle$. The damping is given by (14) where $\Upsilon(\mathbf{k}_n, \Omega)$ is evaluated at $\Omega = |\mathbf{k}_n|v$.

Fig. 1 shows results for v and Γ for three parent temperatures. $T_p = 0.2$; glasses obtained by quenching liquid samples equilibrated at 0.2 are much less stable than

typical laboratory glasses. $T_p = 0.085$, which is between the mode-coupling temperature $T_c \approx .108$ and the estimated laboratory glass transition temperature $T_g \approx 0.072$; glasses obtained by quenching samples equilibrated at 0.085 are about as stable as typical laboratory glasses. $T_p = 0.062$, which is well below estimated T_g ; glasses obtained by quenching liquid samples equilibrated at 0.062 are as stable as laboratory ultrastable glasses obtained by the vapor deposition method^{35,36}. We previously showed⁵ that sound damping coefficients decrease by more than an order of magnitude over this range of stability.

For all three parent temperatures there is excellent agreement between results of Eqs. (13-14) and transverse and longitudinal sound speeds, v_T and v_L , and transverse and longitudinal sound damping coefficients, Γ_T and Γ_L obtained previously⁵ from direct simulations of sound attenuation. At small wavevectors we recover Rayleigh scaling, $\Gamma \propto k^4$, but the theory also accurately predicts sound damping for wavevectors outside the Rayleigh scaling regime. The predicted damping coefficients depart from the simulation results for larger wavevectors, but at larger wavevectors the assumptions used to find the poles, Eq. (12), become invalid.

III. THE ORIGIN OF SOUND ATTENUATION: NON-AFFINE EFFECTS

To get some physical insight into the origin of sound attenuation in zero-temperature amorphous solids we examine the small wavevector expansion of $\langle 1|\mathcal{H}(\mathbf{k})\mathcal{Q}|\mathcal{E}^p\rangle$,

$$\begin{aligned} \langle 1|\mathcal{H}(\mathbf{k})\mathcal{Q}|\mathcal{E}^p\rangle &= -iN^{-1/2} \sum_j \Xi_{j,\gamma\delta} \hat{\mathbf{e}}_\gamma k_\delta \cdot \mathcal{E}_j^p \quad (15) \\ &+ \rho^{-1} N^{-1/2} \sum_j [A_{j,\alpha\beta\gamma\delta}^{\text{Born}} - \hat{\mathbf{e}}_\alpha \hat{\mathbf{e}}_\mu A_{\mu\beta\gamma\delta}^{\text{Born}}] \hat{\mathbf{e}}_\gamma k_\beta k_\delta \mathcal{E}_{j,\alpha}^p + o(k^2). \end{aligned}$$

In Eq. (15) \mathcal{E}_j^p denotes the component of the p th eigenvector of the Hessian corresponding to particle j and $\mathcal{E}_{j,\alpha}^p$ denotes its Cartesian component α . Furthermore, $\Xi_{j,\beta\gamma}$ denotes the vector field describing forces due affine deformations,

$$\Xi_{j,\gamma\delta} = - \sum_{l \neq j} \frac{\partial^2 V(R_{jl})}{\partial R_{j,\gamma} \partial R_j} R_{jl,\delta}. \quad (16)$$

Specifically, $\Xi_{j,\gamma\delta}$ is proportional to the force on particle j resulting from a deformation along the γ direction that linearly depends on the δ coordinates. Finally, the 2nd term at the right-hand-side of Eq. (15) accounts for the spatial variation of the local Born wave propagation coefficients.

As discussed in the literature^{37,38}, forces encoded in vector field $\Xi_{j,\gamma\delta}$ do not seem to possess any longer-range correlations. In contrast, non-affine displacements given by $\mathcal{H}^{-1} \cdot \Xi$ exhibit characteristic vortex-like structures and correlations extending over many particle

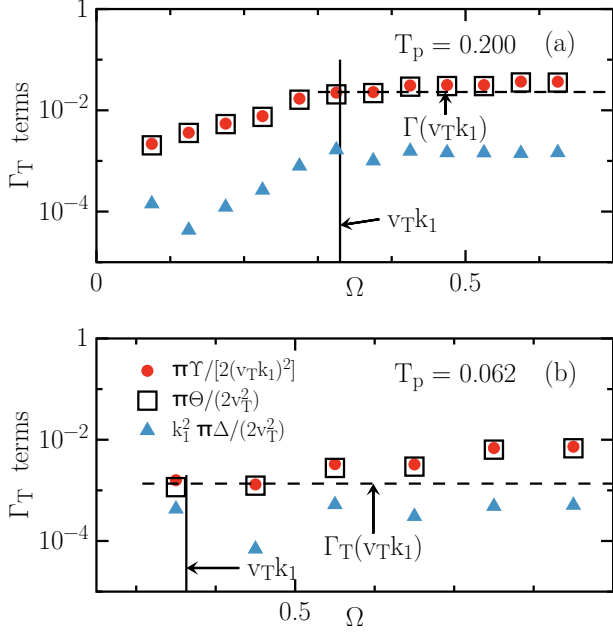


FIG. 2. The terms that contribute to the transverse sound attenuation for $k_1 = 2\pi/L = 0.13722$, where L is the length of the simulation box, given by Eq. 20 for $T_p = 0.200$ (a) and $T_p = 0.062$ (b). The vertical lines marks the frequency $\Omega = v_T k_1$, where v_T is the transverse speed of sound, and the horizontal line is the sound attenuation calculated in simulations from Ref.⁵. The value of $\pi\Upsilon(\mathbf{k}_1, \Omega = v_T k_1)/[2(v_T k_1)^2]$ gives the damping for k_1 .

diameters^{37–40}. The characteristic length of these correlations determines the minimal length scale on which a macroscopic elastic approach can be used to describe the response of amorphous solids³⁹. It follows from the combination of Eqs. (9), (13) and (15) that the renormalization of the wave propagation coefficients originates from the first term in Eq. (15),

$$\lim_{k \rightarrow 0} k^{-2} \Sigma'(\mathbf{k}; 0) = N^{-1} \int d\Omega \Theta(\Omega) \Omega^{-2} \quad (17)$$

where $\Theta(\Omega)$ is defined analogously to $\Upsilon(\mathbf{k}, \Omega)$,

$$\Theta(\Omega) = (1/d\Omega) \sum_{\Omega_p \in [\Omega, \Omega+d\Omega]} \left| \Xi_{\gamma\delta}^p \hat{e}_\gamma \hat{k}_\delta \right|^2 \quad (18)$$

with

$$\Xi_{\gamma\delta}^p = N^{-1/2} \sum_j \Xi_{j,\gamma\delta} \cdot \mathcal{E}_j^p. \quad (19)$$

Equations (17-18) reproduce the exact expression for the non-Born contribution to the wave propagation coefficients derived from the analysis of the non-affine displacements^{37,38}. We note that function $\Theta(\Omega)$ is closely related to function $\Gamma_{\alpha\beta\kappa\chi}(\omega)$ introduced and evaluated by Lemaître and Maloney, see Eq. (32) of Ref.³⁸.

While only the first term in Eq. (15) determines the renormalization of the wave propagation coefficients, both terms contribute to sound attenuation,

$$\Gamma(k) = \frac{\pi}{2v^2} [\Theta(vk) + k^2 \Delta(vk)] \quad (20)$$

where $\Delta(\Omega)$ is defined analogously to $\Upsilon(\mathbf{k}, \Omega)$,

$$\Delta(\Omega) = (1/d\Omega) \sum_{\Omega_p \in [\Omega, \Omega+d\Omega]} \left| \delta A_{\beta\gamma\delta}^p \hat{k}_\beta \hat{e}_\gamma \hat{k}_\delta \right|^2 \quad (21)$$

with

$$\delta A_{\beta\gamma\delta}^p = \rho^{-1} N^{-1/2} \sum_j [A_{j,\alpha\beta\gamma\delta}^{\text{Born}} - \hat{e}_\alpha \hat{e}_\mu A_{\mu\beta\gamma\delta}^{\text{Born}}] \mathcal{E}_{j,\alpha}^p. \quad (22)$$

We note that the second term in Eq. (20) is expressed in term of the fluctuations of the local Born wave propagation coefficients, see Eq. (22). Thus, the physical content of the second term resembles that of the fluctuating elasticity theory. We will discuss this correspondence further in the next section.

It is the first term in Eq. (20) that makes the dominant contribution to the damping coefficient, see Fig. 2. This implies that the sound damping is primarily determined by function $\Theta(\Omega)$, which is the same function that also determines the renormalization of the wave propagation coefficients, Eq. (17). While previous studies suggested⁴¹ and analyzed approximately⁴² the importance the non-affine effects for the sound attenuation, we have presented the first approach that accounts for these effects exactly.

IV. SOUND DAMPING IN THE PLANE-WAVE APPROXIMATION

The most recent version of the fluctuating elasticity theory discussed by Mahajan and Pica Ciammarra²² posits that “amorphous materials can be described as homogeneous isotropic elastic media punctuated by quasiloocalized modes acting as elastic heterogeneities.” This suggests that plane waves should be a reasonable zeroth order approximation for the eigenvectors of the Hessian matrix describing an amorphous solid. To check this supposition we calculated Θ and Δ contributions in Eq. (20) approximating the exact eigenvectors by plane waves, $\mathcal{E}_j^p \propto \hat{e}_\mathbf{q} e^{-i\mathbf{q}\cdot\mathbf{R}_j}$, see Appendix C for details. For the contributions to transverse wave damping coefficient we obtained the following expressions

$$\frac{\pi}{2v^2} \Theta(v_T k) \approx \frac{1}{60\pi} \frac{k^4}{\rho^3} \left[\left(\frac{v_T^2}{v_L^5} + \frac{4}{v_T^3} \right) \langle A_{\alpha\beta xy}^{\text{Born}} A_{\alpha\beta xy}^{\text{Born}} \rangle + \left(\frac{v_T^2}{v_L^5} - \frac{1}{v_T^3} \right) \langle A_{\alpha\alpha xy}^{\text{Born}} A_{\gamma\gamma xy}^{\text{Born}} + A_{\alpha\beta xy}^{\text{Born}} A_{\beta\alpha xy}^{\text{Born}} \rangle \right], \quad (23)$$

$$\frac{\pi k^2}{2v^2} \Delta(v_T k) \approx \frac{1}{12\pi} \frac{k^4}{\rho^3} \left(\frac{1}{v_L^3} + \frac{2}{v_T^3} \right) \left[\langle (\Delta A_{xyxy}^{\text{Born}})^2 \rangle + \langle (A_{yyxy}^{\text{Born}})^2 \rangle + \langle (A_{zyxy}^{\text{Born}})^2 \rangle \right], \quad (24)$$

where we implicitly assumed analyticity of the correlation functions of local wave propagation coefficients at the vanishing wavevector. For example, we assumed that at $\mathbf{q} \rightarrow 0$,

$$\langle (\Delta A_{xyxy}^{\text{Born}})^2 \rangle = \lim_{\mathbf{q} \rightarrow 0} N^{-1} \left| \sum_j e^{i\mathbf{q} \cdot \mathbf{R}_j} (A_{j,xyxy}^{\text{Born}} - A_{xyxy}^{\text{Born}}) \right|^2 \quad (25)$$

and other similar equalities, as discussed in Appendix C.

We note that while the exact formula (18) for Θ contribution involves non-affine forces Ξ , approximate formula (23) is expressed in terms of correlations of local wave propagation coefficients. This follows from the fact that, as shown in Appendix C, for small wavevectors \mathbf{q}

$$\sum_j \Xi_{j,\gamma\delta} \cdot \hat{\mathbf{e}}_{\mathbf{q}} e^{-i\mathbf{q} \cdot \mathbf{R}_j} = \frac{i}{\rho} \sum_j A_{j,\alpha\beta\gamma\delta}^{\text{Born}} \hat{\mathbf{e}}_{\mathbf{q}\alpha} q_\beta e^{-i\mathbf{q} \cdot \mathbf{R}_j} + o(q). \quad (26)$$

Furthermore, we note that formulae (23-24) are reminiscent of Zeller and Pohl's "isotopic scattering" model in that every atom j is a source of scattering of a plane wave, with the amplitude depending on its local wave propagation coefficient $A_{j,\alpha\beta\gamma\delta}^{\text{Born}}$. Importantly, our approximate formulae involve correlations of local wave propagation coefficients that vanish at the macroscopic level and thus do not appear in the semi-phenomenological fluctuating elasticity theory, *e.g.* $A_{j,\gamma\delta}^{\text{Born}}$.

The plane wave approximation recovers analytically the Rayleigh scattering k^4 scaling of the sound damping coefficient. However, it is quantitatively quite inaccurate, see Fig. 3. This implies that at least for the purpose of calculating sound damping, eigenvectors of the Hessian are not well approximated by plane waves. We note that the plane-wave approximation becomes more accurate with decreasing parent temperature or increasing glass stability.

Finally, we note that the first term in square brackets in Eq. (24), which involves correlations of the fluctuations of the local shear modulus, $\Delta A_{j,xyxy}^{\text{Born}}$, represents the result of the microscopic, isotopic scattering-like, version of the fluctuating elasticity theory. As shown in Fig. 3, this term is about 2.5-4 times smaller than the complete plane-wave result, and thus it severely underestimates sound attenuation.

In Fig. 3 we also show the result of a semi-phenomenological fluctuating elasticity theory. To calculate this result we started from the celebrated formula of Rayleigh⁴³ that predicts the attenuation of a transverse wave due to inclusions of volume V_d and number density n , $\Gamma^{\text{R}}(k) = nV_d v_T \gamma k^4 / (6\pi)$, where γ is the disorder parameter. In Rayleigh's calculation γ characterized the variation of "optical density". To adopt his calculation to the present problem we expressed γ in terms of the variation of the square of the transverse speed of sound, $\gamma = (\delta v_T^2)^2 V_d / v_T^4$. Next, we added to

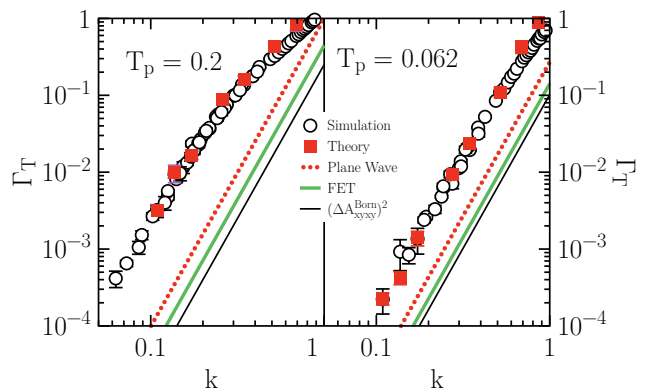


FIG. 3. Comparison of the sound damping coefficients obtained from the theory (squares) and evaluated from sound attenuation simulations (circles) with predictions of the plane-wave approximation, Eqs. (23-24) (red symbols) and microscopic version of the fluctuating elasticity theory, *i.e.* the contribution due to $\langle (\Delta A_{xyxy}^{\text{Born}})^2 \rangle$ term in Eq. (24) (black line). Also shown is the result of a semi-phenomenological fluctuating elasticity theory (FET, green line).

Rayleigh's expression the the contribution of the longitudinal waves excited due to the presence of the inclusions, $\delta\Gamma(k) = nV_d v_T \gamma k^4 (v_T^3 / (2v_L^3)) / (6\pi)$. The complete formula of the semi-phenomenological fluctuating elasticity theory thus reads

$$\Gamma^{\text{FET}}(k) = \frac{k^4 v_T}{6\pi} \left(1 + \frac{1}{2} \frac{v_T^3}{v_L^3} \right) nV_d \gamma. \quad (27)$$

We note that if one makes the identification $\langle (\Delta A_{xyxy}^{\text{Born}})^2 \rangle / (\rho^3 v_T^4) = nV_d \gamma$, the contribution to sound attenuation due to the first term in square brackets in Eq. (24) becomes identical to expression (27). To calculate the value of Γ^{FET} we need the disorder parameter γ and the volume fraction of the inclusions nV_d . For γ we used previously obtained results for the fluctuations of local elastic constants⁴⁴. We recall that disorder parameters calculated this way increase slightly with increasing box size used to define local elastic constants, thus we used the largest box size considered in Ref.⁴⁴. Furthermore, we note that Mahajan and Pica Ciamarra's formulation of fluctuating elasticity theory assumes $nV_d \ll 1$, see the SI of Ref.²². To calculate the upper bound for Γ^{FET} we substituted $nV_d = 1$. Figure 3 shows that the result of this procedure significantly underestimates sound attenuation.

We note that in addition to the microscopic version of the fluctuating elasticity theory, originally derived by Caroli and Lemaître and embodied in the first term in square brackets in Eq. (24), and the semi-phenomenological approach resulting in expression (27) one could compare our results to predictions of more sophisticated versions of the fluctuating elasticity theory, *e.g.* the version relying upon the self consistent Born approximation⁸. This comparison is left for future work.

V. DISCUSSION

According to our microscopic analysis, sound attenuation in zero-temperature amorphous solids is primarily determined by internal forces induced by initial affine displacements of the particles, *i.e.* by the physics of non-affine displacement fields. Quantitatively, the damping coefficient is proportional to the non-affine contribution to the wave propagation coefficients from the frequency equal to the frequency of the sound wave. It is not trivial that our exact calculation (as opposed to the plane-wave approximation discussed in the previous section) reproduces the Rayleigh scaling of sound damping coefficients. This fact results from the frequency dependence of Θ and Δ , which deserves further theoretical study.

The mechanism of the attenuation revealed by our microscopic analysis was mentioned by Caroli and Lemaître in Ref.¹⁹. It was investigated in Ref.⁴¹, where Caroli and Lemaître considered separately the effects of the long-wavelength, elastic continuum-like, and small-scale, primarily non-affine, motions with the small-scale motions being the scatterers for the long-wavelength ones.

An earlier study by Wang, Szamel and Flenner⁵ found a strong correlation between the sound attenuation coefficient and the amplitude of the vibrational density of states of quasilocalized modes. The latter modes were defined using a cutoff in the participation ratio, following Mizuno *et al.*⁴⁵ and Wang *et al.*³¹. We attempted to quantify the relative contributions of the extended and quasi-localized modes by separating the contributions of modes with small and large participation ratio. We did not find convincing evidence for the dominance of small participation ratio modes versus larger participation ratio modes.

We note in this context that local oscillator models^{11–14} express the sound attenuation coefficient in terms of the contributions from localized “defects”^{46,47} referred to as “soft modes”. The formulas derived in these approaches are similar to our Eqs. (14) and (20). The details of expressions of Refs.^{11–14} and our Eqs. (14) and (20) differ; in particular we express the self-energy in terms of all the exact eigenvectors and eigenvalues of the Hessian matrix. In order to evaluate the local oscillator based sound attenuation coefficient formulas one needs to characterize the properties of the soft modes. In practical applications one may parametrize the soft modes’ properties and fit the parameters to the experimental results. Such a procedure was used by Schober¹⁴ and resulted in a good agreement between the theory and experiment.

In view of both the previously found⁵ correlation between the sound attenuation coefficient and the amplitude of the vibrational density of states of quasilocalized modes and the success of local oscillator approach¹⁴ we believe that future work should investigate whether dominant contributions to the sound attenuation coefficient formulas (14) and (20) originate from well defined regions that can be identified as “defects”.

Damart *et al.*⁴⁰ demonstrated that the non-affine dis-

placement field was responsible for high-frequency harmonic dissipation in a simulated amorphous SiO₂. Therefore, it appears that non-affine displacements are responsible for dissipation over the full frequency range. Further theoretical development is needed to connect the low-frequency and high-frequency theories.

Recently, Baggioli and Zaccone developed an approximate microscopic theory for the sound attenuation that takes into account non-affine displacements⁴². This theory shares physical insight with our approach but it is quantitatively as inaccurate as the plane-wave version of our exact formula.

As we mentioned in the introduction, Gelin *et al.*⁷ found a logarithmic correction to the Rayleigh scattering scaling of the sound damping coefficients, which within the fluctuating elasticity theory could originate from the slowly decaying correlations between local values of the elastic constants^{7,15}. Within our approach, a logarithmic correction could originate from a logarithmic dependence of $\Theta(\Omega)$ or $\Delta(\Omega)$ on frequency Ω . Our present numerical data are consistent with the absence of such a logarithmic dependence but it would be interesting to investigate this issue farther.

Within the plane-wave approximation a logarithmic correction could result from a logarithmic small wavevector divergence of the correlation functions of local Born wave propagation coefficients. We did not observe such a divergence but we note that our systems were significantly smaller than those discussed in Ref.⁷. We note that if the correlation functions of local wave propagation coefficients are singular, additional terms in the plane-wave approximation will appear. These terms will originate from the anisotropic small wavevector character of the correlation functions of local wave propagation coefficients.

Our approach arrives at the physical picture of sound attenuation different from that postulated in the fluctuating elasticity theory. While the latter theory can predict trends^{20,22}, it is quantitatively very inaccurate, as noted earlier by Caroli and Lemaître¹⁹. Our analysis revealed that the fluctuating elasticity theory misses the dominant non-affine effects. In addition, it does not include the contributions due to fluctuations of local microscopic wave propagation coefficients that vanish at the macroscopic level. Most importantly, the fluctuating elasticity theory uses plane-wave-like picture of sound in low-temperature amorphous solids. The comparison of the results obtained using the full theoretical expression and adopting the plane-wave approximation, shown in Fig. 3, suggests that this leads to large quantitative discrepancies.

Finally, we note that calculating sound attenuation using Eq. (14) or (20) is somewhat numerically demanding but more straightforward than analyzing the time dependence of the velocity or displacement auto-correlation functions. The latter analysis suffers from large finite-size effects^{5,48} that make the evaluation of the sound damping coefficients at the smallest wavevectors allowed by the pe-

periodic boundary conditions difficult. Our approach offers an attractive alternative way to evaluate sound damping coefficients of low temperature elastic solids that does not suffer from finite size effects.

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AUTHOR DECLARATIONS

Conflict of interest

The authors have no conflicts to disclose.

DATA AVAILABILITY

The data that support the findings of this study are available from the corresponding author upon reasonable request.

Appendix A: Approximation $\mathcal{H}(\mathbf{k}) \approx \mathcal{H}$ in Eq. (5) of the main text

First, we examine the small wavevector expansion of $\mathcal{Q}\mathcal{H}(\mathbf{k})\mathcal{Q}$. The i, j element, which is a 3x3 tensor, reads

$$\begin{aligned} [\mathcal{Q}\mathcal{H}(\mathbf{k})\mathcal{Q}]_{ij} &= \mathcal{H}_{ij}e^{-i\mathbf{k}\cdot(\mathbf{R}_i-\mathbf{R}_j)} - N^{-1} \sum_l \mathcal{H}_{il}e^{-i\mathbf{k}\cdot(\mathbf{R}_i-\mathbf{R}_l)} \cdot \hat{\mathbf{e}}\hat{\mathbf{e}} - N^{-1} \hat{\mathbf{e}} \sum_l \hat{\mathbf{e}} \cdot \mathcal{H}_{lj}e^{-i\mathbf{k}\cdot(\mathbf{R}_l-\mathbf{R}_j)} \\ &+ N^{-2} \hat{\mathbf{e}} \sum_{l,m} \hat{\mathbf{e}} \cdot \mathcal{H}_{lm} \cdot \hat{\mathbf{e}} e^{-i\mathbf{k}\cdot(\mathbf{R}_l-\mathbf{R}_m)} \hat{\mathbf{e}} = [\mathcal{Q}\mathcal{H}(\mathbf{k}=0)\mathcal{Q}]_{ij} + [\delta\mathcal{H}^{\mathcal{Q}1}(\mathbf{k})]_{ij} + [\delta\mathcal{H}^{\mathcal{Q}2}(\mathbf{k})]_{ij} + o(k^2), \end{aligned} \quad (\text{A.1})$$

where the matrix elements of the terms of the first and second order in k , $\delta\mathcal{H}^{\mathcal{Q}1}(\mathbf{k})$ and $\delta\mathcal{H}^{\mathcal{Q}2}(\mathbf{k})$, read

$$[\delta\mathcal{H}^{\mathcal{Q}1}(\mathbf{k})]_{ij} = i(1 - \delta_{ij}) \left\{ \frac{\partial^2 V(R_{ij})}{\partial \mathbf{R}_i^2} \mathbf{k} \cdot (\mathbf{R}_i - \mathbf{R}_j) - N^{-1} \sum_l \hat{\mathbf{e}} \cdot \left[\frac{\partial^2 V(R_{il})}{\partial \mathbf{R}_i^2} \mathbf{k} \cdot (\mathbf{R}_i - \mathbf{R}_l) - \frac{\partial^2 V(R_{lj})}{\partial \mathbf{R}_j^2} \mathbf{k} \cdot (\mathbf{R}_j - \mathbf{R}_l) \right] \hat{\mathbf{e}} \right\}, \quad (\text{A.2})$$

$$\begin{aligned} [\delta\mathcal{H}^{\mathcal{Q}2}(\mathbf{k})]_{ij} &= \frac{1}{2} (1 - \delta_{ij}) \frac{\partial^2 V(R_{ij})}{\partial \mathbf{R}_i^2} (\mathbf{k} \cdot (\mathbf{R}_i - \mathbf{R}_j))^2 - \frac{1}{2} N^{-1} \sum_{l \neq i} \frac{\partial^2 V(R_{il})}{\partial \mathbf{R}_i^2} \cdot \hat{\mathbf{e}} (\mathbf{k} \cdot (\mathbf{R}_i - \mathbf{R}_l))^2 \hat{\mathbf{e}} \\ &- \frac{1}{2} N^{-1} \hat{\mathbf{e}} \sum_{l \neq j} \hat{\mathbf{e}} \cdot \frac{\partial^2 V(R_{lj})}{\partial \mathbf{R}_j^2} (\mathbf{k} \cdot (\mathbf{R}_l - \mathbf{R}_j))^2 + \frac{1}{2} N^{-2} \hat{\mathbf{e}} \sum_{l \neq m} \hat{\mathbf{e}} \cdot \frac{\partial^2 V(R_{lm})}{\partial \mathbf{R}_l^2} \cdot \hat{\mathbf{e}} (\mathbf{k} \cdot (\mathbf{R}_l - \mathbf{R}_m))^2 \hat{\mathbf{e}}. \end{aligned} \quad (\text{A.3})$$

Next, we assume that for small wavevectors k we can treat terms $\delta\mathcal{H}^{\mathcal{Q}1}(\mathbf{k})$ and $\delta\mathcal{H}^{\mathcal{Q}2}(\mathbf{k})$ in the denominator of Eq. (5) of the main text perturbatively. Due to the symmetry, the term of the first order in k , $\delta\mathcal{H}^{\mathcal{Q}1}(\mathbf{k})$, will

contribute in the second order of the perturbation expansion. In contrast, the term of the second order in k , $\delta\mathcal{H}^{\mathcal{Q}2}(\mathbf{k})$, will contribute in the first order. Here we will show the contribution of $\delta\mathcal{H}^{\mathcal{Q}2}(\mathbf{k})$ term. It reads

$$\begin{aligned} &\delta\Sigma^{\mathcal{Q}2}(\mathbf{k}; \omega) \quad (\text{A.4}) \\ = &- \sum_{\substack{\text{eigenvec.} \\ p,q}} \langle 1 | \mathcal{H}(\mathbf{k})\mathcal{Q} | \mathcal{E}_p \rangle \langle \mathcal{E}_p | (-\omega + i\epsilon)^2 + \lambda_p)^{-1} | \mathcal{E}_p \rangle \langle \mathcal{E}_p | \delta\mathcal{H}^{\mathcal{Q}2}(\mathbf{k}) | \mathcal{E}_q \rangle \langle \mathcal{E}_q | (-\omega + i\epsilon)^2 + \lambda_q)^{-1} | \mathcal{E}_q \rangle \langle \mathcal{E}_q | \mathcal{Q}\mathcal{H}(\mathbf{k}) | 1 \rangle. \end{aligned}$$

Counting powers of k in the expression above shows that,

at least perturbatively, term $\delta\mathcal{H}^{\mathcal{Q}2}(\mathbf{k})$ results in a correc-

tion that is higher order in k than the dominant small wavevector result of approximation $\mathcal{H}(\mathbf{k}) \approx \mathcal{H}$ in the denominator of Eq. (5).

Appendix B: Simulation details

We obtained zero-temperature glasses by instantaneously quenching supercooled liquids of unit number density, $\rho = 1.0$, equilibrated through the swap Monte Carlo algorithm²⁹. The constituent particles of these liquids have unit mass and diameters σ selected using distribution $P(\sigma) = \frac{A}{\sigma^3}$, where $\sigma \in [0.73, 1.63]$ and A is a normalization factor. The cross-diameter σ_{ij} is determined according to a non-additive mixing rule, $\sigma_{ij} = \frac{\sigma_i + \sigma_j}{2}(1 - \epsilon|\sigma_i - \sigma_j|)$ with $\epsilon = 0.2$. The interaction between two particles i and j is given by the inverse power law potential, $V(r_{ij}) = (\sigma_{ij}/r_{ij})^{12} + V_{cut}(r_{ij})$, when the separation r_{ij} is smaller than the potential cutoff $r_{ij}^c = 1.25\sigma_{ij}$, and zero otherwise. Here, $V_{cut}(r_{ij}) = c_0 + c_2 (r_{ij}/\sigma_{ij})^2 + c_4 (r_{ij}/\sigma_{ij})^4$, and the coefficients c_0 , c_2 and c_4 are chosen to guarantee the continuity of $V(r_{ij})$ at r_{ij}^c up to the second derivative.

The number of particles N varied between 48000 and 192000. The largest systems had to be analyzed to determine sound attenuation at the lowest wavevectors re-

ported.

Appendix C: Plane-wave approximation

We assume that for small wavevectors we can approximate eigenvectors of the Hessian matrix by plane waves. We note that strictly speaking, for our amorphous solids the normalization factor is configuration-dependent. We checked that this dependence is weak and for this reason we use the following approximation,

$$\mathcal{E}_j^p \approx N^{-1/2} \hat{\mathbf{e}}_{\mathbf{q}} e^{-i\mathbf{q} \cdot \mathbf{R}_j}. \quad (\text{C.1})$$

Approximate plane-wave eigenvectors are labeled by their wavevector \mathbf{q} and their polarization $\hat{\mathbf{e}}_{\mathbf{q}}$. For each wavevector \mathbf{q} we have one longitudinal and two transverse modes. We assume that the associated eigenvalues are given by $(v_L q)^2$ and $(v_T q)^2$ for the longitudinal and transverse modes, respectively.

Here we will present the derivation of approximate formula for the contribution to the transverse sound damping coefficient originating from Θ , Eq. (23) of the main text. The contribution originating from Δ , Eq. (24) of the main text and the approximate expression for the longitudinal sound damping can be derived in a similar way.

First, we need to calculate

$$-iN^{-1/2} \sum_j \Xi_{j,\gamma\delta} \hat{\mathbf{e}}_{\gamma} k_{\delta} \cdot \mathcal{E}_j^p \approx -iN^{-1} \sum_j \Xi_{j,\gamma\delta} \hat{\mathbf{e}}_{\gamma} k_{\delta} \cdot \hat{\mathbf{e}}_{\mathbf{q}} e^{-i\mathbf{q} \cdot \mathbf{R}_j} = iN^{-1} \sum_j \sum_{l \neq j} \frac{\partial^2 V(R_{jl})}{\partial R_{j,\gamma} \partial \mathbf{R}_j} R_{jl,\delta} \hat{\mathbf{e}}_{\gamma} k_{\delta} \cdot \hat{\mathbf{e}}_{\mathbf{q}} e^{-i\mathbf{q} \cdot \mathbf{R}_j} \quad (\text{C.2})$$

Using the $i \leftrightarrow j$ symmetry we get

$$\begin{aligned} iN^{-1} \sum_j \sum_{l \neq j} \frac{\partial^2 V(R_{jl})}{\partial R_{j,\gamma} \partial \mathbf{R}_j} R_{jl,\delta} \hat{\mathbf{e}}_{\gamma} k_{\delta} \cdot \hat{\mathbf{e}}_{\mathbf{q}} e^{-i\mathbf{q} \cdot \mathbf{R}_j} &= \frac{i}{2N} \sum_j \sum_{l \neq j} \frac{\partial^2 V(R_{jl})}{\partial R_{j,\alpha} \partial R_{j,\gamma}} R_{jl,\delta} \hat{\mathbf{e}}_{\alpha} \hat{\mathbf{e}}_{\gamma} k_{\delta} \left[1 - e^{i\mathbf{q} \cdot (\mathbf{R}_j - \mathbf{R}_l)} \right] e^{-i\mathbf{q} \cdot \mathbf{R}_j} \\ &= \frac{1}{2N} \sum_j \sum_{l \neq j} \frac{\partial^2 V(R_{jl})}{\partial R_{j,\alpha} \partial R_{j,\gamma}} R_{jl,\delta} R_{jl,\beta} \hat{\mathbf{e}}_{\alpha} \hat{\mathbf{e}}_{\beta} \hat{\mathbf{e}}_{\gamma} k_{\delta} e^{-i\mathbf{q} \cdot \mathbf{R}_j} + o(q) = (\rho N)^{-1} \sum_j A_{j,\alpha\beta\gamma\delta}^{\text{Born}} \hat{\mathbf{e}}_{\alpha} \hat{\mathbf{e}}_{\beta} \hat{\mathbf{e}}_{\gamma} k_{\delta} e^{-i\mathbf{q} \cdot \mathbf{R}_j} + o(q). \quad (\text{C.3}) \end{aligned}$$

Next, we need to take the square of the absolute value of expression (C.3) for a given wavevector \mathbf{q} and polarization $\hat{\mathbf{e}}_{\mathbf{q}}$ and then integrate over spherical shell with frequency $qv_L = kv_T$ for longitudinal and $qv_T = kv_T$ for transverse modes. We shall note that since the spherical shell is specified in the frequency space, there will be additional factors, $1/v_L$ for longitudinal and $1/v_T$ for transverse modes. Finally, we need to multiply the result by $\pi/(2v_T^2 k^2)$ to get the contribution to the transverse sound damping coefficient.

To perform these calculations we assume that wavevector \mathbf{k} is parallel to the y axis and the sound polarization $\hat{\mathbf{e}}$ is along the x axis. Furthermore, we specify the polarization vector for the approximate plane-wave eigenvectors as $\hat{\mathbf{e}}_{\mathbf{q}}^L = \hat{\mathbf{q}} \equiv (\sin \theta \cos \phi, \sin \theta \sin \phi, \cos \theta)$ for the longi-

tudinal modes and $\hat{\mathbf{e}}_{\mathbf{q}}^{\text{T1}} = (\cos \theta \cos \phi, \cos \theta \sin \phi, -\sin \theta)$ and $\hat{\mathbf{e}}_{\mathbf{q}}^{\text{T2}} = (-\sin \phi, \cos \phi, 0)$ for the two transverse modes.

The contribution of the longitudinal modes reads

$$\begin{aligned} &\frac{\pi}{2(v_T k)^2} \frac{v_T^4}{v_L^5} \frac{V k^6}{(2\pi)^3} \quad (\text{C.4}) \\ &\times \int d\hat{\mathbf{q}} \left| (\rho N)^{-1} \sum_j A_{j,\alpha\beta\gamma\delta}^{\text{Born}} \hat{c}_{\mathbf{q}\alpha}^L \hat{q}_{\beta} e^{-i\hat{\mathbf{q}}(kv_T/v_L) \cdot \mathbf{R}_j} \right|^2. \end{aligned}$$

Guided by our numerical calculations we assume that the following small wavevector limit is finite and does not

depend on the direction

$$\lim_{q \rightarrow 0} N^{-1} \sum_j A_{j,\alpha\beta xy}^{\text{Born}} e^{i\mathbf{q} \cdot \mathbf{R}_j} \sum_l A_{l,\gamma\delta xy}^{\text{Born}} e^{-i\mathbf{q} \cdot \mathbf{R}_l} \\ \equiv \langle A_{\alpha\beta xy}^{\text{Born}} A_{\gamma\delta xy}^{\text{Born}} \rangle. \quad (\text{C.5})$$

Expression (C.4) becomes

$$\frac{\pi}{2} \frac{v_T^4}{(v_T k)^2} \frac{V k^6}{v_L^5 (2\pi)^3} \frac{1}{\rho^2 N} \langle A_{\alpha\beta xy}^{\text{Born}} A_{\gamma\delta xy}^{\text{Born}} \rangle \int d\hat{\mathbf{q}} \hat{e}_{\mathbf{q}\alpha}^L \hat{q}_\beta \hat{e}_{\mathbf{q}\gamma}^L \hat{q}_\delta \\ = \frac{\pi}{2} \frac{v_T^4}{(v_T k)^2} \frac{V k^6}{v_L^5 (2\pi)^3} \frac{1}{\rho^2 N} \langle A_{\alpha\beta xy}^{\text{Born}} A_{\gamma\delta xy}^{\text{Born}} \rangle \\ \times \frac{4\pi}{15} (\delta_{\alpha\beta} \delta_{\gamma\delta} + \delta_{\alpha\gamma} \delta_{\beta\delta} + \delta_{\alpha\delta} \delta_{\beta\gamma}) \\ = \frac{1}{60\pi} \frac{k^4}{\rho^3} \frac{v_T^2}{v_L^5} [\langle A_{\alpha\alpha xy}^{\text{Born}} A_{\beta\beta xy}^{\text{Born}} \rangle + \langle A_{\alpha\beta xy}^{\text{Born}} A_{\alpha\beta xy}^{\text{Born}} \rangle \\ + \langle A_{\alpha\beta xy}^{\text{Born}} A_{\beta\alpha xy}^{\text{Born}} \rangle]. \quad (\text{C.6})$$

Assuming again that the small wavevector limit of the correlation functions of local wave propagation coefficients is finite and does not depend on the direction, the contribution of the two transverse modes reads

$$\frac{\pi}{2} \frac{1}{(v_T k)^2} \frac{V k^6}{v_T (2\pi)^3} \frac{1}{\rho^2 N} \langle A_{\alpha\beta xy}^{\text{Born}} A_{\gamma\delta xy}^{\text{Born}} \rangle \\ \times \int d\hat{\mathbf{q}} \left(\hat{e}_{\mathbf{q}\alpha}^{\text{T1}} \hat{q}_\beta \hat{e}_{\mathbf{q}\gamma}^{\text{T1}} \hat{q}_\delta + \hat{e}_{\mathbf{q}\alpha}^{\text{T2}} \hat{q}_\beta \hat{e}_{\mathbf{q}\gamma}^{\text{T2}} \hat{q}_\delta \right) \\ = \frac{\pi}{2} \frac{1}{(v_T k)^2} \frac{V k^6}{v_T (2\pi)^3} \frac{1}{\rho^2 N} \langle A_{\alpha\beta xy}^{\text{Born}} A_{\gamma\delta xy}^{\text{Born}} \rangle \\ \times \frac{4\pi}{15} (4\delta_{\alpha\gamma} \delta_{\beta\delta} - \delta_{\alpha\beta} \delta_{\gamma\delta} - \delta_{\alpha\delta} \delta_{\beta\gamma}) \\ = \frac{1}{60\pi} \frac{k^4}{\rho^3} \frac{1}{v_T^3} [4 \langle A_{\alpha\beta xy}^{\text{Born}} A_{\alpha\beta xy}^{\text{Born}} \rangle - \langle A_{\alpha\alpha xy}^{\text{Born}} A_{\beta\beta xy}^{\text{Born}} \rangle \\ - \langle A_{\alpha\beta xy}^{\text{Born}} A_{\beta\alpha xy}^{\text{Born}} \rangle] \quad (\text{C.7})$$

Adding expressions (C.6) and (C.7) we get Eq. (23) of the main text.

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