

Thermal conductivity by non-equilibrium molecular dynamics: Inescapable non-locality and how best to deal with it.

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Non-equilibrium molecular dynamics (NEMD) enables “direct” simulation of thermal conductivity κ . A steady state heat current is driven by heat inputs $H(x)$. A corresponding temperature (T) gradient is extracted. Since the system is finite, non-local relations between $H(x)$ and $T(x)$ (and thus between $\vec{J}(x)$, the heat current, and $\vec{\nabla}T(x)$) are inevitable. Computation can study $\tilde{\kappa}(q)$, the Fourier transform of $\kappa(x - x')$. The $q \rightarrow 0$ limit is the desired macroscopic κ , but the best that can be done is $\tilde{\kappa}(q = 2\pi/L)$ where L is the long axis of the simulation cell. Therefore it makes sense to add heat $H(x) \propto \sin(qx)$, where $q = 2\pi/L$. Analysis of the steady state temperature $T(x) - T_{\text{avg}} \propto \sin(qx)$ allows an accurate estimate of $\tilde{\kappa}(q)$. A recent variation of the popular Müller-Plath algorithm adds and removes chosen amounts of heat with minimal disruption and any designated spatial pattern. The method is tested by computation of $\tilde{\kappa}(q)$ for the Lennard-Jones (LJ) liquid and crystal. Peierls-Boltzmann theory in relaxation time approximation assists analysis of convergence in the crystal. Assuming $1/\tau_Q \propto \omega_Q^2$, an isotropic cell ($N_x = N_y = N_z$) behaves as $\tilde{\kappa}(q) = \kappa - A \times q^{1/2}$. For the more typical anisotropic cell with one length (N_x) large compared to the others, there is an additional term $\propto q^{-1/2}/N_y N_z$. This divergent contribution disappears in the bulk limit. Strategies for extrapolation of simulations are tested.

I. INTRODUCTION

Molecular dynamics (MD) simulation applies to insulating solids at high temperature T . If $k_B T$ is reasonably high on the scale of the lattice energies $\hbar\omega$, classical Newtonian trajectories seem to give correct thermal averages. Going beyond harmonic lattice dynamics is difficult with quantum mechanics, but easy with MD, not requiring perturbation theory. A “direct” simulation of thermal conductivity κ by non-equilibrium molecular dynamics (NEMD) was first done by Payton *et al.*¹ in 1967, and has become common. Good examples are the comparative study by Schelling *et al.*² and the careful study of GaN by Zhou *et al.*³. The biggest difficulty is that small ω_Q vibrational modes have long mean free paths Λ_Q , but only modes with $\Lambda_Q < L$ (where L is the length of the simulation cell) have their contribution to κ correctly treated by NEMD.

This brings up the issue of non-locality of κ . Although non-locality is widely acknowledged, it was seldom^{4–8} regarded as a direct topic of study. Recent interest in boundary effects and in spatially varying heat input $H(\vec{r})$ give new impetus to non-local thinking^{9–15}.

A good illustration of non-local behavior is the simulation by Zhou *et al.*³ reprinted as Fig 1. The temperature excursion $\Delta T(x)$ is generated by a non-uniform profile of heating $H(x)$. Heating is confined to separate slabs under the constraint $\int dx H(x) = 0$ which allows a steady state to emerge. For weak enough heating $|H(x)|$,

the linear relation is $\Delta T(x) = \int dx' S(x - x') H(x')$. In Fourier variables, this is $\tilde{\Delta T}(q) = S(q) \tilde{H}(q)$. Using energy conservation $dJ(x)/dx = H(x)$ and the non-local Fourier law¹⁶ $J(x) = -\int dx' \kappa(x - x') dT(x')/dx'$, we find $\tilde{\kappa}(q) = 1/q^2 S(q)$. The procedure normally used is to compute dT/dx at the midpoint between heat injection and removal, and thus estimate the bulk κ as the ratio of J (the current known from heat input and energy conservation) to $-dT/dx$. We denote κ derived in this way as $\kappa_{\text{eff}}(L)$, where L is the size of the system. In this paper, we show that we can do better than $\kappa_{\text{eff}}(L)$ by thinking non-locally. We show that NEMD calculations are, at best, only able to compute $\tilde{\kappa}(q)$ where $q = 2\pi/L$ and L is the length of the simulation cell. This is as close as such calculations can get to their ultimate aim, the true bulk $\kappa = \tilde{\kappa}(q \rightarrow 0) = \int d\vec{r} \kappa(\vec{r})/V$. An optimum strategy is therefore to aim directly for a computation of $\tilde{\kappa}(2\pi/L)$ before trying to extrapolate to $L \rightarrow \infty$. The reason why $\kappa = \tilde{\kappa}(0)$ is unavailable is because $\tilde{H}(0) = \int d\vec{r} H(\vec{r})/V$ must vanish in order for a steady state to be allowed.

This paper does four things. (1) We Fourier-analyze the common version of NEMD where the system is periodic with length L along the direction of heat flow, and heat input and removal occurs in slabs separated by $L/2$. A rigorous relation between $\kappa_{\text{eff}}(L)$ and $\tilde{\kappa}(q)$ is worked out. (2) A method is given for direct MD simulation of $\tilde{\kappa}(q)$, by applying and extracting heat in a sinusoidal pattern¹⁷. This has numerical advantages over other protocols. (3) Convergence of $\tilde{\kappa}(q)$ towards κ is studied by NEMD simulation for the Lennard-Jones (LJ) liquid and crystal. (4) The Peierls “Boltzmann Transport Equation” (BTE)^{18,19} is solved for $\tilde{\kappa}(q)$ in Debye approximation ($\omega_Q = v|\vec{Q}|$) with $1/\tau_Q = (1/\tau_D)(\omega_Q/\omega_D)^2$. This

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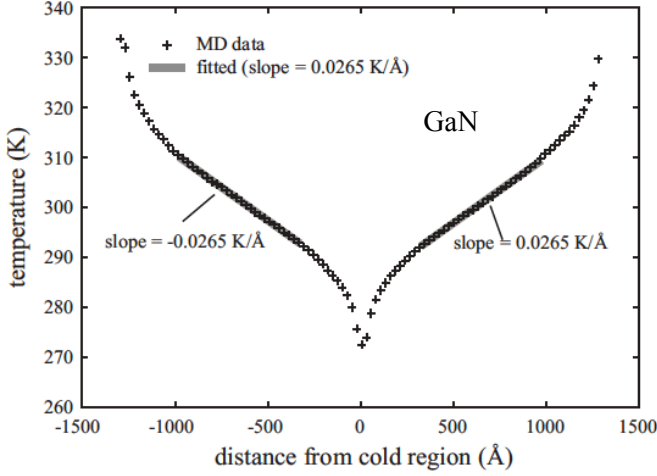


FIG. 1. Simulated temperature profile, computed by Zhou *et al.*³ for wurtzite structure GaN. This is Fig. 6d, reprinted without permission, so far. Heat is removed at a constant rate ($H(0) = -H$) at a slab located at $x = 0$, and inserted at the same rate ($H(\pm L/2) = H$) in equivalent slabs at $x = \pm 250c = \pm 1300\text{\AA}$ along the crystallographic c -axis. There are 30,000 GaN pairs and 1000 GaN double layers in the simulation cell. The temperature is averaged over the final 20 ns of a 24 ns MD run. Since current is independent of x outside the heat-exchange slabs (except changing sign as x changes sign), the curvature seen in the temperature profile is a direct indicator of the non-local relation between current and temperature gradient.

helps understand convergence as the MD simulation-cell size L increases.

II. PRELIMINARIES

This paper assumes a homogeneous single crystal, represented by a simulation cell with periodic boundary conditions. The i^{th} atom at \vec{r}_i and the atom at $\vec{r}_i + \vec{R}$ are equivalent, and have the same trajectory $\vec{r}_i(t)$. For simplicity, the primitive translation vectors \vec{R} of the simulation cell ($\vec{A}_1, \vec{A}_2, \vec{A}_3$), are assumed orthogonal. For example, in the sample calculations presented later, they are taken to be $N_x a \hat{x}, N_y a \hat{y}, N_z a \hat{z}$, where a is the lattice constant, the edge-length of the fcc conventional cube. A typical cell has size $(N_x, N_y, N_z) = (80, 6, 6)$; with 4 atoms in the conventional cube, this means 11,520 atoms. Heating is done in slabs perpendicular to the long vector \vec{A}_1 . Therefore, current flows parallel to \vec{A}_1 , and $L = L_x = |\vec{A}_1| = N_x a$ is chosen as large as computation permits, trying to surpass the distance Λ of non-local thermal memory.

Boundaries create challenging problems. Nanoscale heat transfer is typically dominated by thermal boundary (or Kapitza) resistance^{20,21}. For study of bulk conductivity, the aim is to reduce the influence of boundaries. One can argue²² that periodic boundary conditions are not

the most favorable way to do this. However, in this paper, periodicity offers simplicity for analysis, overruling other considerations.

A further simplification follows computational necessity, and discretizes the temperature $T(\vec{r})$ into slab values $T(\ell)$. The slabs are layered in the \vec{A}_1 -direction, and have width $d = n_S a$ where n_S is a small integer and a factor of N_x . The number of slabs is $N_S = N_x/n_S$. Let the variable x denote distance along the \vec{A}_1 -axis, perpendicular to slabs. The slab indexed by the integer ℓ occupies the interval $\ell d - d/2 < z < \ell d + d/2$. The temperature $T(\ell)$ is found from the average kinetic energy of the atoms in the ℓ^{th} slab. Heat is transferred externally into the ℓ^{th} cell at a volume-average rate $H(\ell)$. In steady state, this is balanced by an outward flux $[J(\ell+1/2) - J(\ell-1/2)]/d$. Both current and temperature gradient are properties of the junction of two adjacent slabs. Their steady-state linear relation is $J(\ell+1/2) = -\sum_{\ell'} \kappa(\ell, \ell') \nabla T(\ell'+1/2)$, where the sum runs over the N_S slabs. This definition is required by linear math. Periodicity requires $\kappa(\ell, \ell') = \kappa(\ell + mN_S, \ell' + nN_S)$ and homogeneity (if the medium is in fact homogeneous) requires that $\kappa(\ell, \ell') = \kappa(\ell - \ell')$. Corresponding (*via* a unitary Fourier transformation) to the N_S distinct slabs, there are N_S distinct wave-vectors $q = 2\pi n_q/N_S d$, indexed by integers n_q and distributed in a one-dimensional Brillouin zone. In the homogeneous case, the relation is $\tilde{J}(q) = -\tilde{\kappa}(q) \tilde{\nabla} T(q)$. The tilde on top of a function indicates that it is in reciprocal space rather than coordinate space. These ideas were introduced in Ref. 17, where further properties are explained.

It is not hard to extend the usual derivation of the Kubo formula (ref. 23, for example) to derive a Kubo formula for $\kappa(x, x')$ or $\kappa(\ell, \ell')$. The classical limit is

$$\kappa(\ell, \ell') = -\frac{1}{k_B T^2} \int_0^\infty dt \langle J(\ell + 1/2, t) J(\ell' + 1/2, 0) \rangle \quad (1)$$

III. DISCRETE SLAB HEATING

A common version of NEMD simulation, illustrated in Fig. 1, removes energy only from slab $\ell = 0$, at a volume-average rate H , and inserts energy at the same rate into slab $\ell = N_S/2$. Zhou *et al.*³ did a careful study of κ for GaN by this method. They discuss, but do not completely resolve, the issue of how the answer for κ scales with system size. Here we analyze this version with periodic boundaries ($\ell = \ell + N_S$), rather than the rigid or free boundaries sometimes used. Heat current $J_x(\ell + 1/2) = \pm(d/2)H \equiv \pm J$ flows across slab boundaries, the plus sign for ℓ to the right of an input and left of an output slab, and the minus sign for opposite cases. Thermal conduction is found using the relation $\kappa_{\text{eff}}(L) \equiv -J/\nabla_x T(\text{mid})$. The rigorous non-local Fourier law is $\tilde{J}(q) = -\tilde{\kappa}(q) \tilde{\nabla} T(q)$ in the one-dimensional discrete Fourier space $q = 2\pi n_q/L$. Analysis given in Ap-

pendix A shows that

$$\frac{1}{\kappa_{\text{eff}}} = \frac{2}{N_S} \sum_q^{n_q=\text{odd}} \frac{(-1)^{(n_q-1)/2}}{\sin(qd/2)} \left[\frac{\cos(qd/2)}{\kappa(q)} \right] \quad (2)$$

This holds if $N_S/4$ is an integer. If it is a half-integer, then $\cos(qd/2)$ is replaced by 1. Liquids have a quite local conductivity, where $\tilde{\kappa}(q) \approx \kappa$ is nearly independent of q . Then the sum in Eq.(2) converges correctly to $\kappa_{\text{eff}} \rightarrow \kappa$ without needing a small $q_{\text{min}} = 2\pi/L$. For crystals like GaN, where non-local behavior is caused by long mean free paths of small- Q phonons, $\tilde{\kappa}(q)$ peaks around $|q| = 0$. Accurate results require a small q_{min} or a large L . When q_{min} is not very small, the first two terms $n_q = \pm 1$ in Eq.(2) dominate. Approximating $\sin(q_{\text{min}}d/2)$ by π/N_S and $\cos(q_{\text{min}}d/2)$ by 1, the result is $\kappa_{\text{eff}} \approx \pi\tilde{\kappa}(q_{\text{min}})/4$, which is smaller than $\pi\kappa/4$ (and still smaller than the true κ .) It would be better to calculate $\tilde{\kappa}(q_{\text{min}})$ directly. A method is given in the next section.

IV. SINUSOIDAL HEATING ALGORITHM

The simulation cell is divided into slabs centered at $x(\ell), \ell = 1, \dots, N_S$. It is shown schematically in Fig. 2. We want to modify the heat input profile. Instead of the two slab version, let the heat input be of the form $H(\ell) = \tilde{H}(q_{\text{min}}) \cos(q_{\text{min}}x(\ell))$. The temperature variation then has the form $T(\ell) = T_0 + \tilde{\Delta T}(q_{\text{min}}) \cos(q_{\text{min}}x(\ell))$. Since slab temperatures $T(\ell)$ at all N_S values of ℓ are used to compute the Fourier amplitude $\tilde{\Delta T}(q_{\text{min}})$, numerical noise averages out faster.

Equation 26 of the previous paper¹⁷ says, for arbitrary heating $H(\ell)$, heating rate, temperature, and conductivity in Fourier space are related by (for $q \neq 0$),

$$\tilde{\Delta T}(q) = \frac{\tilde{H}(q)d^2}{4 \sin^2(qd/2) \tilde{\kappa}(q)} \quad (3)$$

For simple sinusoidal heating, $\tilde{H}(q)$ and $\tilde{\Delta T}(q)$ are zero except at $q = q_{\text{min}}$, where their values are denoted as H and ΔT . Then the conductivity is

$$\tilde{\kappa}(q_{\text{min}}) = \frac{Hd^2/\Delta T}{4 \sin^2(q_{\text{min}}d/2)} \rightarrow \left(\frac{N_S d}{2\pi} \right)^2 \frac{H}{\Delta T} \quad (4)$$

This makes sense: $2\pi\Delta T/N_S d$ is the maximum temperature gradient, and $H N_S d/2\pi$ is the maximum heat current.

Now we need a good numerical algorithm to drive the oscillatory heat input. Furtado *et al.*²⁴ have a nice modification of the popular algorithm by Müller-Plathe²⁵. The usual Müller-Plathe method gives equal heating and cooling to two chosen slabs, normally repeated always on the same two slabs. The hottest atom in the slab chosen for heat removal, and the coldest atom in the slab for heat insertion, have their velocities interchanged, conserving

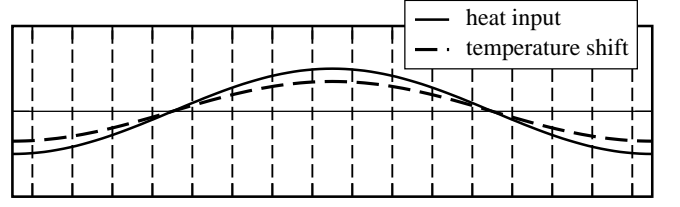


FIG. 2. Schematic picture of an ($N_S = 16$)-slab simulation cell with periodic boundary conditions. Heat is added or subtracted according to position, in the ℓ^{th} slab, as $H \cos(2\pi\ell/N_S)$. The central cell is numbered $\ell = 0$. Temperature $T(\ell)$ is determined by averaging kinetic energy of atoms in each slab. In linear approximation, temperature must also vary as $\Delta T \cos(2\pi\ell/N_S)$.

energy and momentum. The Furtado version adds and subtracts a $\Delta \vec{v}$ from these two atoms in such a way that energy and momentum are still conserved, and the magnitude $|\Delta \vec{v}|$, and the resulting disruption, is minimized. This allows heat input at a predetermined rate which can be spatially varied. Appendix B gives details.

V. TEST ON LENNARD-JONES LIQUID

The LJ liquid is a simple case, used by Müller-Plathe²⁵ to test his algorithm. The pair potential is

$$V_{\text{LJ}} = 4\epsilon \left[\left(\frac{\sigma}{r} \right)^{12} - \left(\frac{\sigma}{r} \right)^6 \right] \quad (5)$$

The parameters for argon²⁶ are $\epsilon/k_B = 119.6\text{K}$ and $\sigma = 3.82\text{\AA}$. First, we reproduce Müller-Plathe's results, at the same $(N, V, T) = 2592$ atoms, $\rho = N/V = 0.849/\sigma^3$, and $T = 0.7\epsilon/k_B = 84\text{K}$. The same simulation cell is used, consisting of $18 \times 6 \times 6$ fcc conventional cubes, and periodic boundary conditions. The cut-off distance for the LJ potential is 3.0σ . We get the same answer, $\kappa = 7.1$ in LJ units.

As shown in Fig. 2 of ref. 25, and confirmed by our calculation in Fig. 3, the temperature gradient is essentially constant all the way to, and including, the slabs 0 and $N_S/2 = 10$. This is because thermal conductivity in a liquid is very local. This can be contrasted with Fig. 1 or fig. 4 of ref. 17, for crystals with non-local κ . Gas theory is certainly not correct for a liquid; the concept of a mean-free path is not valid. However, we can get an idea of what happens by unlicensed use of the gas formula $\kappa = C\bar{v}\Lambda/3$. The measured thermal conductivity of liquid argon (0.132 W/mK at temperature near 100K and pressure near 1Mbar²⁷⁻³⁰) then corresponds to a mean free path $\Lambda \approx 0.14\text{\AA}$, more than 30 times smaller than the slab separation $d = 1.51\sigma$. In other words, the non-local conductivity $\kappa(z-z')$ decays to zero by the first neighbor slab, or $\kappa(q)$ is independent of q out to values of q larger than $q_{\text{max}} = \pi/d$.

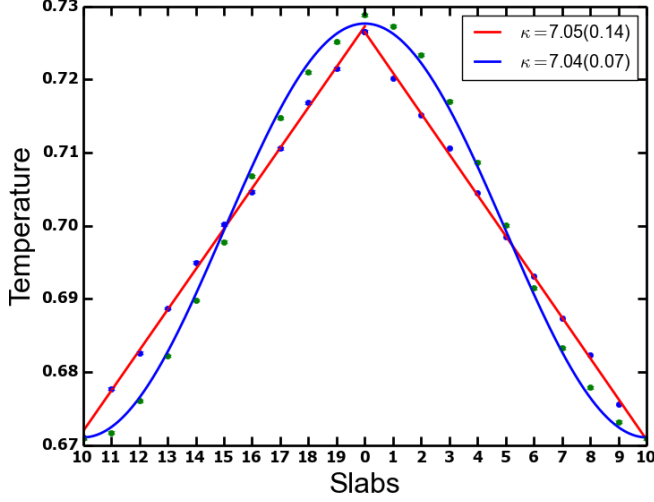


FIG. 3. Temperature profiles from NEMD simulations for the LJ liquid. The points fit by the straight line are a reproduction of the calculation of Müller-Plathe. The points fit by the sine curve use the same simulation cell, with spatially sinusoidal heating and cooling.

Also shown in Fig. 3 is the sinusoidal temperature profile gotten numerically from our sinusoidal heating. The computational system is unaltered. The 2592 atoms are in the same cell, divided into 20 slabs, at the same T . Sinusoidal heat $N_S H$ between 1 and $5 \epsilon / (\Delta t \cdot \text{slab})$ is used. **Yerong – was heating done to each slab simultaneously, or in a sequence of one slab for each Δt ?** The interval between heat injections is $\Delta t = 60\delta t$, where the time step of the “velocity Verlet” Newton’s-law integration algorithm^{31,32} is $\delta t = 0.007t_{LJ}$. The LJ unit of time for argon is $t_{LJ} = \sigma\sqrt{(m/\epsilon)} = 2.16$ ps. Equilibration required $10^4 \delta t$ of constant T simulation, and $T(\ell)$ averaging was done for $2 \times 10^5 \delta t$; good convergence was found in $5 \times 10^4 \delta t$ as shown in Fig. 4.

To estimate errors, consider that there are 130 atoms per slab, each with mean energy $k_B T$ and rms deviation of $k_B T$ from the mean, according to Maxwell-Boltzmann statistics. Thus the mean slab energy per atom, at any particular moment, should be about $k_B T \pm k_B T / \sqrt{130}$. Therefore, if averaged over 100 random and independent thermalized configurations, the temperature error in a slab will be less than 1%. A run of $5 \times 10^4 \delta t$ should be more than sufficient for this purpose. Fig. 3 suggests errors of order $0.001k_B T$ in the slab temperatures.

Compared with the Müller-Plathe²⁵ result, both of the current ones give the same value, $\kappa = 7.1$ in LJ units, which is 0.133 W/mK, very close to the experimental value for argon, 0.132 W/mK^{27–30}. The sinusoidal algorithm gives faster convergence and a slightly more accurate final answer, as shown in Fig 4.

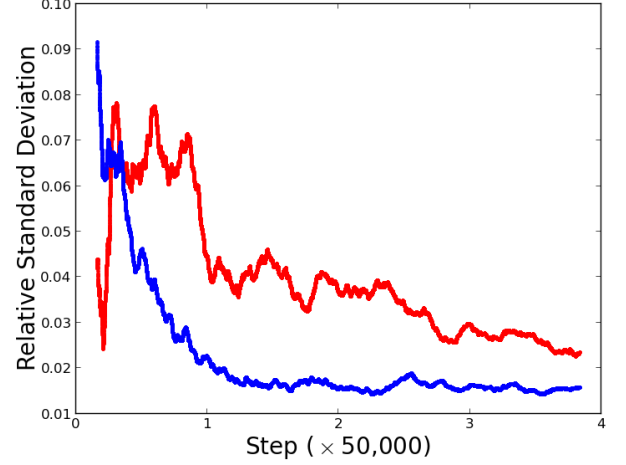


FIG. 4. Time evolution of the error of simulation of κ for the LJ liquid. The upper and noisier curve uses the method of Müller-Plathe. The lower curve is the result of the sinusoidal algorithm of this paper.

VI. EXTRAPOLATION

NEMD answers for κ are computed for finite size L . Therefore extrapolation is required to estimate the bulk ($L \rightarrow \infty$) answer. Sellan *et al.*³³ have analyzed this. It was also analyzed in the previous paper¹⁷, using a Debye model. Here we continue the analysis. Equations (22,23) of ref. 17 are

$$\kappa(q) = \frac{1}{\Omega} \sum_Q \hbar \omega_Q \frac{\partial n_Q}{\partial T} v_{Qz}^2 \tau_Q \cos^2(qd/2) F(q, \Lambda_{Qz}) \quad (6)$$

$$F(q, \Lambda) = \left[1 + 4 \sin^2(qd/2) \left\{ \left(\frac{\Lambda}{d} \right) + \left(\frac{\Lambda}{d} \right)^2 \right\} \right]^{-1}, \quad (7)$$

where v_{Qz} , τ_Q , and $\Lambda = v_{Qz} \tau_Q$ are the group velocity, quasiparticle lifetime, and z -component of mean free path of the phonon mode of frequency ω_Q . These equations solve the BTE in relaxation-time approximation (RTA; also known as the “Single Mode Approximation”) for the case where heat is applied as $\dot{e} \propto \cos(q\ell d)$. In a simulation, the relevant wave vector is $q = 2\pi/L$ where L is the periodic repeat length. This can be visualized as an infinite sample divided into super cells of length L , or as a finite lattice with periodic boundary conditions (equivalent to being wrapped on a torus.) This is the style of simulation used in the calculations of secs. V and VII).

Recently there has been much progress in numerical solution of the BTE, using input parameters (phonons and anharmonic coupling) from density-functional theory^{34–37}. These methods enable careful comparison of RTA against exact solutions. For many materials, RTA answers are very accurate at room temperature, graphene^{38,39} being a notable exception. Here

we adopt both the RTA and the equally unrealistic Debye model. We do not expect an accurate description of κ , but we do expect that this model should give useful insight about the size-dependence of $\kappa(q_{\min} = 2\pi/L)$, and how to extrapolate to the $L \rightarrow \infty$ limit.

In ref. 17, Eqns. 6, 7 were approximately evaluated using a Debye model, Q being integrated over a Debye sphere. The Debye version is

$$\kappa_D(q) = \kappa_0 \cos^2(qd/2) \frac{1}{N} \sum_Q \left(\frac{Q_z}{Q}\right)^2 \left(\frac{Q_D}{Q}\right)^p F(q, \Lambda_{Qz}), \quad (8)$$

where κ_0 is a convenient scale factor. It is the Debye answer for $\kappa(q = 0)$, exponent $p = 2$, and an isotropic Q -mesh,

$$\kappa_0 = \frac{3N}{\Omega} k_B v^2 \tau_D. \quad (9)$$

The mean free path increases for small Q as Q^{-p}

$$\Lambda_{Qz} = v \tau_D \frac{Q_z}{Q} \left(\frac{Q_D}{Q}\right)^p. \quad (10)$$

The Debye wavevector has its usual value, $(6\pi^2 N/\Omega)^{1/3}$, and the exponent p is chosen to be 2.

Here, instead of integrating over the Debye sphere, we use direct numerics to do the discrete sum of Eq. 8 over the actual discrete Q -mesh in the Brillouin zone of the *fcc* simulation cell that will be used in the next section for the LJ crystal. That is, we use only those \vec{Q} 's in the *fcc* Brillouin zone such that $\exp(i\vec{Q} \cdot \vec{A}_i) = 1$, where \vec{A}_i , for $i = 1, 2, 3$, are the orthogonal translation vectors of the simulation supercell. As an example, the mesh used in Sec. V for the LJ liquid corresponds to $\vec{A}_1 = 6a(1, 0, 0)$, $\vec{A}_2 = 6a(0, 1, 0)$, and $\vec{A}_3 = 18a(0, 0, 1)$, where $a = 1.68\sigma$ is the lattice constant of the *fcc* conventional cubic cell, using the liquid argon density, $0.849/\sigma^3$. Our simulation cells in this and the next section will be very similar, but longer in the \vec{A}_3 direction, and with a readjusted to 1.56σ to give the higher density^{32,40}, $1.053/\sigma^3$, of the low pressure LJ crystal. The corresponding \vec{Q} 's are the vectors $\ell\vec{G}_1 + m\vec{G}_2 + n\vec{G}_3$ of the lattice reciprocal to the \vec{A} 's. This is an anisotropic reciprocal-space mesh, being coarse in the directions \vec{A}_1 and \vec{A}_2 , but finer in the direction \vec{A}_3 , corresponding to the actual distribution of normal modes of the atoms in the simulation cell of the LJ crystal (where $|\vec{G}_{1,2}|$ is larger than $|\vec{G}_3|$.) We are guessing that the Debye model, with frequency $\omega_Q = v|\vec{Q}|$ for all 3 branches, and $1/\tau_Q = (1/\tau_D)(Q/Q_D)^2$, sufficiently captures the physics of the LJ crystal for purposes of learning how to extrapolate to infinite simulation cell size.

Results are shown in Fig. 5 and in appendices C, D, and E. The figure shows two things. First, quite smooth extrapolation to the correct $q = 0$ answer appears when $\kappa(q)$ is plotted *versus* \sqrt{qa} , as anticipated in Refs. 33 and 17, and clarified in Appendix C. Second, an unexpected

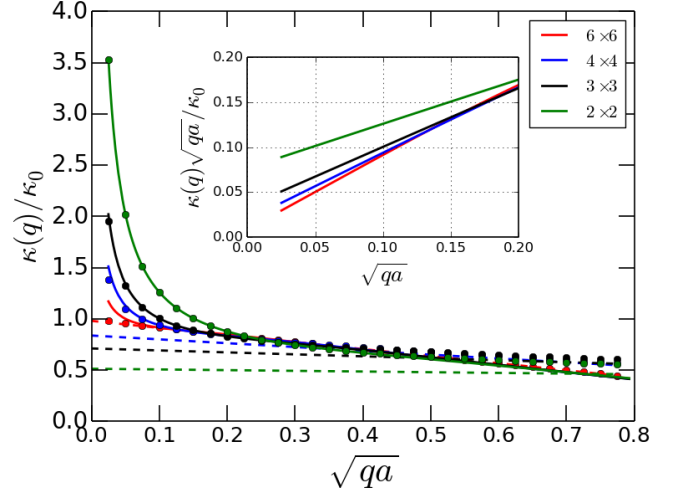


FIG. 5. Convergence of $\kappa(q)$ toward κ_0 found by numerical summation of the Boltzmann-Debye RTA model for an anisotropic cell of size $N_x = N_y$ equalling 2, 3, 4, and 6 (top to bottom) and $N_z = 2\pi/qa$ varying from 10 to 10,000. The Q -sum in Eq.(8) is evaluated not over the Debye sphere, but over the anisotropic Q -mesh corresponding to the normal modes of the *fcc* crystalline LJ lattice; results are shown as solid lines. The inset shows that the small q divergence has a $1/\sqrt{qa}$ form with coefficient diminishing as N_x increases. The simulation results are fitted (for $50 < N_z < 1000$ or $.08 < \sqrt{qa} < .35$) to a 3-term form $\kappa(q) \approx A/\sqrt{qa} + B + C\sqrt{qa}$. The fits are indicated by dots whose colors agree with the lines. The value of B (for $N_x \geq 6$) is a reasonable choice for $\kappa(q = 0)$. The dashed lines show the non-diverging parts $\kappa(q) \approx B + C\sqrt{qa}$ of the 3-term fits. The coefficient B is the slope of the lines in the inset, and the $\sqrt{qa} = 0$ intercept of the dashed lines on the main graph.

divergence (of the form $1/\sqrt{qa}$ begins to appear for cells with small enough q (relative to the transverse size N_x .) Specifically, the onset of the upward turn appears roughly when $\sqrt{qa} < 1/2N_x$, which corresponds to $N_z > 25N_x^2$, a limit not always achieved in NEMD calculations. The origin and significance of this divergence is discussed in Appendix D. The idea for extrapolation is discussed in the caption to Fig. 5 and in Appendix E.

VII. LENNARD-JONES CRYSTAL

Unlike the LJ liquid, for an LJ crystal, phonon gas theory applies well, but only up to half the melting temperature, where higher-order anharmonic terms become important¹⁵. This non-Boltzmann regime is where an MD simulation is worth doing. The resulting shorter phonon mean free paths permit shorter simulation cells. We simulate crystalline LJ argon at $T=80K$, close to the experimental triple point (84K and 0.7 atmospheres) and atmospheric pressure melting temperature (84K).

Higher energy phonons have mean free paths a bit longer than the unit cell $a = 5.32\text{\AA}$, which we choose

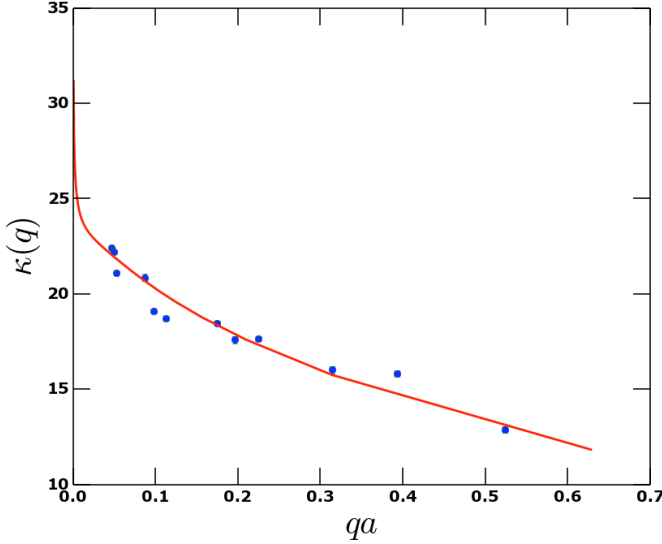


FIG. 6. Results of NEMD for the LJ crystal. The wave-vector $q = 2\pi/L_z$ is the smallest compatible with the simulation cell, whose size is $6a \times 6a \times (L_z = N_z a)$. Values of N_z vary from 12 to 120. The red line is a numerical summation of the Debye-RTA model, Eq. 8, using a Q -mesh with $N_x = N_y = 6$ compatible with the discrete normal mode quantum numbers of the simulation cell. The $N_z = 2\pi/qa$ values vary from 10 at the right, up to 10^4 at the left. The numerical answers are scaled for comparison with the NEMD simulation. An anomalous small- q behavior is better seen in Fig. 7, and is shown in Appendix D to be an artifact of the coarse N_{xy} mesh.

to be the slab thickness d . Lower energy phonons have mean free paths Λ_Q increasing, roughly as $1/\omega_Q^2$. The values of Λ_Q are not as long as in GaN, modeled by Zhou *et al.*³. Nevertheless, doing a converged calculation by MD methods is challenging. We use this to test whether our algorithm is helpful. We use a time step of the “velocity Verlet” Newton’s-law integration algorithm^{31,32} $\delta t = 0.007t_{LJ}$ for smaller-size samples, and $0.014t_{LJ}$ for larger ones.

Figs. 6 and 7 show results for $\kappa(q = 2\pi/L)$ where $L = N_z d$ is the length of the simulation cell, and $d = a$. These calculations used a heat input per volume $\dot{\epsilon}\tau$ of 1.265ϵ , the interval τ being $60\delta t$. The value of κ in argon at $T = 80$ K is measured⁴¹ to be in the range 0.4 to 0.6 W/mK. Christen and Pollack⁴² found $\kappa(80K) \approx 0.30$ W/mK. Fig. 6 looks as if it might extrapolate linearly in the region $qa < 0.1$ to a value around 23 in LJ units, whereas Fig. 7 seems more convincingly linear in \sqrt{qa} , extrapolating to a value $\kappa(q = 0) \approx 26$ in LJ units. The LJ unit of thermal conductivity is $(k_B/\sigma^2)\sqrt{(\epsilon/M)} = 0.0188$ W/mK. Our LJ crystal simulation thus gives $\kappa \approx 0.49$ W/mK. This compares with the value 0.16 – 0.17 W/mK found by Turney *et al.*¹⁵, 0.236 W/mK found by Omini and Sparavigna³⁴, 0.25 W/mK found (at 69.2K) by Kaburaki *et al.*⁴³, 0.33 – 0.49 W/mK (depending on density) found by Christen and

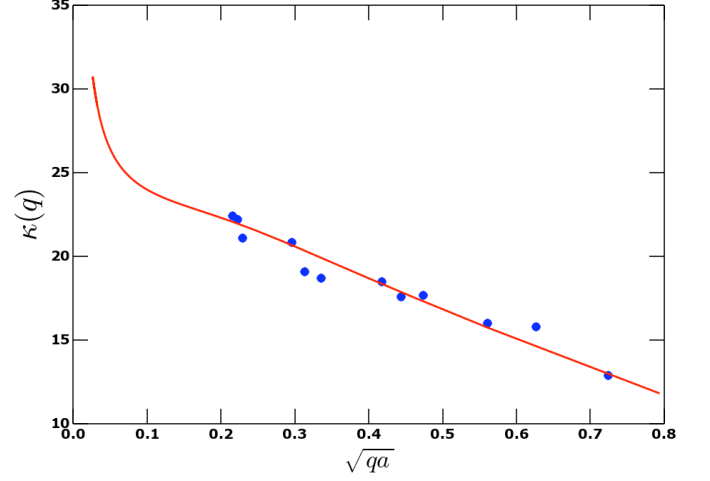


FIG. 7. Points are numerical NEMD, identical to those in Fig. 6. Values of κ are shown in LJ units ($\kappa_{LJ} = (k_B/\sigma^2)\sqrt{(\epsilon/M)} = 0.0188$ W/mK). The red line is also as in Fig. 6, a scaled RTA-Debye model with $p = 2$ (Eq. 8) summed using the same coarse transverse Q -mesh as the simulation cell, and varying N_z and thus $q = q_{\min} = 2\pi/L = 2\pi/N_z a$. Use of \sqrt{qa} for the horizontal coordinate gives better linearity and a more reliable extrapolation to $q = 0$. The anomalous low- q behavior of the red line is discussed in Appendix D.

Pollack⁴², and 0.19 W/mK found by Chernatynskiy and Phillpot⁴⁴.

VIII. CONCLUSIONS

The algorithm of Sec. IV works smoothly and converges more rapidly than the common two-slab heating. It generates a reliable value of $\kappa(q = 2\pi/L)$. The macroscopic conductivity, $\kappa = \lim_{q \rightarrow 0} \kappa(q)$, found by extrapolating the long sample dimension to $L \rightarrow \infty$, is problematic, although less so than for the discrete-slab heating algorithm. Even if Boltzmann transport theory fails because phonon mean free paths are so short that quasiparticles are not well-defined, nevertheless, Boltzmann theory should correctly model the long-wavelength phonon contribution to $\kappa(q)$, which is the problematic part.

Our analysis using the BTE reveals two effects responsible for slower than expected convergence of $\kappa(q)$ to κ . These are an inevitable correction which scales as \sqrt{qa} , and the anisotropic mesh artifact that scales as $1/[N_x N_y \sqrt{qa}] \propto \sqrt{(N_z)/N_x N_y}$. These are found by numerical summation of the Boltzmann-Debye RTA version, but should faithfully model the effects seen in NEMD models. Simple graphical extrapolation to $q = 0$ by assuming linear behavior in \sqrt{qa} is less justifiable than had been hoped, because the contamination by the $1/\sqrt{qa}$ term alters the appearance of the $\kappa(q)$ versus \sqrt{qa} graph. The cure is to be sure that the ratio $\sqrt{(N_z)/N_x N_y}$ does not get too small. Appendix E dis-

cusses this further. There are reasons for mild insecurity on the issue of to what extent extrapolation is justified.

Appendix A: Discrete slab heating

Adding periodic boundary conditions permits a simple Fourier representation of J_x ,

$$\begin{aligned}\tilde{J}_x(q) &= \frac{1}{N_S} \sum_{\ell=0}^{N_S-1} e^{-iqx(\ell+1/2)} J_x(\ell+1/2) \\ &= \frac{J}{N_S} \sum_{\ell=0}^{N_S/2-1} \left[e^{-iqx(\ell+1/2)} - \text{c.c.} \right] \\ &= \frac{J}{N_S} \left(\frac{1 - e^{-iqL/2}}{1 - e^{iqd}} \right) e^{-iqd/2} - \text{c.c.} \\ &= \frac{J}{iN_S} \left[\frac{1 - (-1)^{n_q}}{\sin(qd/2)} \right]\end{aligned}\quad (\text{A1})$$

At $q = 0$, the expression $[]$ in the last line of Eq.(A1) needs definition; the correct value is 0, as is also true for all q 's with even n_q . Now let us analyze the approximate thermal conductivity,

$$\kappa_{\text{eff}}(L) \equiv -J/\nabla_z T(\text{mid}). \quad (\text{A2})$$

That is, the thermal conductivity is approximated by the ratio of the actual current J , controlled by the heating rate \dot{e} , to the temperature gradient $-\nabla_z T(\text{mid})$ found midway between the heat input slab ($\ell = 0$) and output slab ($\ell = N_S/2$). This temperature gradient has the Fourier representation

$$\begin{aligned}\nabla_z T(\text{mid}) &= \sum_q e^{iqN_S d/4} \nabla_z T(q) \quad [N_S/4 = \text{half integer}] \\ &= \sum_q e^{iqN_S d/4} \cos(qd/2) \nabla_z T(q) \quad [N_S/4 = \text{integer}]\end{aligned}\quad (\text{A3})$$

In the case $N_S/4 = \text{integer}$, the slab $\ell = N_S/4$ lies midway between heat input and output, so the temperature gradient (needed at the slab mid-point) is taken as the average of the left and right slab boundaries. This introduces the factor $\cos(qd/2)$ in the second version of Eq.(A3). Finally, we substitute $\nabla_z T(q) = -J_z(q)/\kappa(q)$ in Eq.(A3), and use Eq.(A1) for $J_z(q)$. Then Eq.(A2) becomes

$$\begin{aligned}\frac{1}{\kappa_{\text{eff}}} &= \frac{2}{N_S} \sum_q^{n_q=\text{odd}} \frac{\sin(qL/4)}{\sin(qd/2)} \left[\frac{1 \text{ or } \cos(qd/2)}{\kappa(q)} \right] \\ &= \left[\frac{4 \text{ or } 4 \cos(\pi/N_S)}{N_S \sin(\pi/N_S)} \right] \frac{1}{\kappa(q_{\text{min}})} + \{ |n_q| \geq 3 \text{ terms} \}.\end{aligned}\quad (\text{A4})$$

This is a surprisingly complicated relation between the size-dependent ‘‘computational’’ value of κ_{eff} and the Fourier representation $\kappa(q)$. From Eq.(A4), the leading term (at small $q_{\text{min}}d/2 = \pi d/L = \pi/N_S$)

is $1/\kappa_{\text{eff}} = 4/\pi\kappa(q_{\text{min}})$, with oscillatory corrections $\sum_{n=1} (4/\pi)(-1)^n / [(2n+1)\kappa((2n+1)q_{\text{min}})]$. In the local limit, $\kappa(q) = \kappa$, Eq.(A4) converges exactly to $\kappa_{\text{eff}} = \kappa$.

Appendix B: Heat exchange algorithm

The Müller-Plathe recipe is: find the hottest atom in the cold ($\ell = N_S/2$) slab, and the coldest atom in the hot ($\ell = 0$) slab. Interchange their velocities. Energy and momentum are both conserved, and the system is driven from equilibrium in a way that must be monitored. Cao and Li⁴⁵ among others, have suggested modified algorithms of this type.

In our method, driving is done by spatially periodic injection and simultaneous removal of heat. Examine a slab ℓ (with $0 \leq \ell < N_S/4$), and its conjugate slab $N_S/2 - \ell$. The former is ‘‘hotter’’ than the latter because $\dot{e}(\ell) > 0 > \dot{e}(N_S/2 - \ell) = -\dot{e}(\ell)$. Find the coldest atom (meaning least kinetic energy) of all atoms of mass m_i in the hotter slab, denoting its velocity as \vec{v}_C . Find the hottest atom of the same mass in the colder slab, denoting its velocity \vec{v}_H . Given the large fluctuations of the Maxwell-Boltzmann ensemble, it is certain that $v_H^2 > v_C^2$. Choose an appropriate velocity \vec{w} and add it to the velocity \vec{v}_C and subtract it from \vec{v}_H :

$$\begin{aligned}\vec{v}_H &\rightarrow \vec{v}_H - \vec{w} \\ \vec{v}_C &\rightarrow \vec{v}_C + \vec{w}\end{aligned}\quad (\text{B1})$$

The same operation should be done for the pair of slabs $-\ell$ and $\ell - N_S/2$. All slabs can be done simultaneously, or different random times can be used for different slabs.

There are three criteria for an appropriate \vec{w} , which uniquely fix the desired choice. (i) The cold atom's kinetic energy should increase by Δ , an energy that can be specified in advance as $(\dot{e}\Omega_S\tau)\cos(2\pi\ell/N_S)$, where τ is the average time interval between random interventions, and Ω_S is the volume of a slab. (ii) The hot atom's kinetic energy should decrease by Δ . Then both momentum and energy are conserved. Heating has the desired form $\dot{e}\cos(2\pi\ell/N_S)$ with \dot{e} chosen not too different from $k_B T/\Omega_S\tau$. Trial calculations should test for the best choices of \dot{e} and τ . (iii) There is still a one-dimensional family of vectors \vec{w} ; from these, choose the smallest $|\vec{w}|$, which gives the least impulse to the affected atoms.

The energy shift criteria (i) and (ii) give equations $-2\vec{v}_H \cdot \vec{w} + w^2 = -2\Delta/m \equiv -\delta$, and $2\vec{v}_C \cdot \vec{w} + w^2 = +\delta$. Adding and subtracting these equations give

$$\begin{aligned}\delta &= (\vec{v}_H + \vec{v}_C) \cdot \vec{w} \\ w^2 &= (\vec{v}_H - \vec{v}_C) \cdot \vec{w},\end{aligned}\quad (\text{B2})$$

a linear and a quadratic equation for \vec{w} . These have a simple geometric interpretation shown in Fig. 8. The first equation restricts the projection of \vec{w} along the vector $\vec{v}_H + \vec{v}_C$. Geometrically, this means that \vec{w} lies on the plane (shown by C, C' in Fig. 8) perpendicular to the vector $\vec{b} \equiv \delta(\vec{v}_C + \vec{v}_H)/|\vec{v}_C + \vec{v}_H|^2$, where the origin of

and in Debye approximation, $v_{Qz} = v \cos \theta$. We are interested in the minimum wavevector $q = 2\pi/L$, so $\cos(qd/2) \rightarrow 1$ and $2 \sin(qd/2)\Lambda/d \rightarrow 2\pi\Lambda/L$. Then Eq. C1 becomes (for the special case $p = 2$)

$$\kappa(q_{\min}) = \frac{\kappa_0}{3N} \sum_Q \frac{\cos^2 \theta (Q_D/Q)^2}{1 + \cos^2 \theta \lambda^2 (Q_D/Q)^4} \quad (\text{C2})$$

where $\lambda \equiv q\ell_{\min} = 2\pi v\tau_D/L$ and $\kappa_0 = 3k_B v^2 \tau_D N/\Omega$. This is the Debye-RTA result for any choice of cell dimensions (N_x, N_y, N_z) and corresponding Q -mesh. There are two parameters, ℓ_{\min} which determines the scale of q through the relation $\lambda = q\ell_{\min}$, and κ_0 which determines the scale of κ . In the numerical results shown in Sec. VI, Fig. 5, and later in Appendix E, Figs. 9,10,11, the choices were made $\ell_{\min} = 2.45a$ and $\kappa_0 = 26.44$ (in LJ units). These choices were made simply to scale the numerical summations to fit the numerical simulations of Sec. VII.

For a uniform and dense mesh (*i.e.* a large crystal), the sum becomes an integral,

$$\frac{1}{N} \sum_Q \rightarrow \frac{9}{2} \int_0^{Q_D} \frac{dQ Q^2}{Q_D^3} \int_{-1}^1 d\cos \theta, \quad (\text{C3})$$

where an extra factor of 3 is for the three modes. After performing the angular integral, C2 becomes

$$\begin{aligned} \kappa(q_{\min}) &= \frac{3\kappa_0}{\lambda^2} \int_0^{Q_D} \frac{dQ Q^4}{Q_D^5} \\ &\times \left[1 - \frac{1}{\lambda} \left(\frac{Q}{Q_D} \right)^2 \tan^{-1} \lambda \left(\frac{Q_D}{Q} \right)^2 \right] \\ &= \frac{3\kappa_0}{2\lambda^2} \int_0^1 du u^{3/2} \left[1 - \frac{u}{\lambda} \cot^{-1} \frac{u}{\lambda} \right]. \end{aligned} \quad (\text{C4})$$

Here the replacements have been made $u = (Q/Q_D)^2$ and $\tan^{-1}(1/x) = \cot^{-1} x$.

Next we assert the following identity, best proved by direct differentiation:

$$\begin{aligned} h(u) &= \frac{7}{2} \int du \left[u^{3/2} - \frac{u^{5/2}}{\lambda} \cot^{-1} \frac{u}{\lambda} \right] \\ &= u^{5/2} + 2\lambda^2 u^{1/2} - \frac{u^{7/2}}{\lambda} \cot^{-1} \frac{u}{\lambda} \\ &- \frac{\lambda^{5/2}}{\sqrt{2}} \left[\frac{1}{2} \log \left(\frac{u + \sqrt{2u\lambda} + \lambda}{u - \sqrt{2u\lambda} + \lambda} \right) + \tan^{-1} \left(\frac{\sqrt{2u\lambda}}{\lambda - u} \right) \right] \end{aligned} \quad (\text{C5})$$

Now we have an answer,

$$\kappa(q_{\min}) = \frac{3}{7} \frac{\kappa_0}{\lambda^2} [h(1) - h(0)] \quad (\text{C6})$$

It is necessary to be careful in evaluating $h(1) - h(0)$. As we integrate u from 0 towards 1 (as in Eq. C4), all the terms in the integral, Eq. C5 have smooth behavior except the last term, $\tan^{-1}[\sqrt{(2u\lambda)/(\lambda - u)}]$. Note that

$0 < \lambda < 1$ since L is large. Therefore, in this term, the argument of the arctangent increases from zero to $+\infty$, where the arctangent equals $\pi/2$, then switches to $-\infty$, with the arctangent still equal to $\pi/2$, and finally increases toward 0 from the negative side, with the arctangent approaching π from the negative side. Therefore, the value of $\tan^{-1}[\sqrt{(2u\lambda)/(\lambda - u)}]$ at $u = 1$ is $\pi - \tan^{-1}[\sqrt{(2\lambda)/(1 - \lambda)}]$. The answer is thus

$$\begin{aligned} \frac{7\lambda^2}{3\kappa_0} \kappa(q_{\min}) &= 1 + 2\lambda^2 - \frac{\tan^{-1} \lambda}{\lambda} \\ &- \frac{\lambda^{5/2}}{\sqrt{2}} \left[\frac{1}{2} \log \left(\frac{1 + \sqrt{2\lambda} + \lambda}{1 - \sqrt{2\lambda} + \lambda} \right) + \pi - \tan^{-1} \left(\frac{\sqrt{2\lambda}}{1 - \lambda} \right) \right] \end{aligned} \quad (\text{C7})$$

Finally, at large sample size L , the smallest mean free path $\ell_{\min} = v\tau_D$ is small compared with L . The parameter λ is $2\pi\ell_{\min}/L$ and the leading size-dependent correction is

$$\frac{\kappa(q_{\min})}{\kappa_0} = 1 - \frac{3\pi}{7} \sqrt{\frac{\pi\ell_{\min}}{L}} = 1 - \frac{3\pi}{7} \sqrt{\frac{\ell_{\min}}{2a}} \sqrt{(qa)}. \quad (\text{C8})$$

This result confirms the conjecture of Refs. 33 and 17 that numerical result might be best extrapolated by plotting κ versus $\sqrt{(1/L)}$.

Appendix D: Inhomogeneous mesh

Fig. 7 indicates that the analytic answer of Eq. C8 gives a good match to the size-dependence when L_z is not too big (and q is not too small.) At smaller q there is an up-turn in the numerical Debye-RTA sum, that is not derived in the analytic integration Eq. C7. This up-turn is strongly enhanced at small transverse cell size $N_x N_y$. The problem is that the ratio L_z/L_x becomes very large at small $q = 2\pi/L_z$. It is necessary to reconsider how Eq. C1 behaves in a finite-size crystal or simulation cell, when one dimension ($L_z \equiv L$) gets large but the other two ($L_x = L_y \equiv L_{xy}$) remain small. The answer is, a new non-analytic piece occurs, and rather than converging to the $L \rightarrow \infty$ limit as $1/\sqrt{L}$ (found above for the homogeneous mesh), it diverges as \sqrt{L} . These results are specific to the power law relaxation $\tau_Q = \tau_D (\omega_D/\omega)^p$ with $p = 2$. The divergence is a property of a one-dimensional wire, indicating that ballistic transport has a dominant effect in such a system. As the area $L_x L_y$ of the wire increases, the divergent term in $\kappa(q_{\min})$ decreases as $a^2/L_x L_y$, restoring the better-behaved homogeneous answer.

The specific system under consideration is an fcc crystal of volume $Na^3/4$, where a is the conventional primitive cube size, $N = 4N_x N_y N_z$ is the total number of atoms, N_x and N_y being small integers held fixed, and N_z being a large integer. We seek the behavior as N_z gets very large. The conductivity is given by Eq. C2,

rewritten as

$$\frac{\kappa(q_{\min})}{\kappa_0} = \frac{1}{N} \sum_{Q_z} \sum_{Q_x, Q_y} \frac{(Q_z/Q)^2 (Q_D/Q)^2}{1 + (Q_z/Q)^2 \lambda^2 (Q_D/Q)^4}, \quad (\text{D1})$$

where $\lambda = 2\pi\ell_{\min}/L$, ℓ_{\min} being the mean free path of the highest frequency phonons. The Q -vectors are $\vec{Q} = (2\pi/a)(n_x/N_x, n_y/N_y, n_z/N_z)$. This choice is required to make vibrational normal modes satisfy periodicity in the supercell. There are N Q -vectors in the Brillouin zone. The Debye model simplifies summation over the Brillouin zone by using the Debye sphere, with a volume equal to n times the volume of the primitive Brillouin zone, n being the number of atoms in the primitive cell, 4 for fcc. The Q -points are dense along the z direction and sparse in the others. Only the Q_z -sum can be turned into an integral. It is consistent with the philosophy of the Debye model, to not use a sphere in this case, but a cube-shaped “pseudo-Brillouin zone”, of volume 4 times $(2\pi/a)^3$, containing the correct number of states. The Q_z sum is then an integral, going from $Q_z = 0$ to the boundary of the “pseudo-Brillouin zone,” $4^{1/3}\pi/a$, and multiplied by 2 to cover both negative and positive Q_z .

$$\frac{\kappa(q_{\min})}{\kappa_0} = \frac{1}{4N_x N_y} \sum_{Q_x, Q_y} \frac{a}{\pi} \int_0^{4^{1/3}\pi/a} dQ \frac{(Q_z/Q)^2 (Q_D/Q)^2}{1 + (Q_z/Q)^2 \lambda^2 (Q_D/Q)^4} \quad (\text{D2})$$

The number of terms in the Q_x, Q_y sum is $\approx n^{2/3}N_x N_y$, typically 50 for an MD simulation, or a few thousand for a small nanowire. Of these terms, the one which requires special attention is the $Q_x = Q_y = 0$ term. We denote this term by $\kappa_{00}(q_{\min})$,

$$\begin{aligned} \frac{\kappa_{00}(q_{\min})}{\kappa_0} &= \frac{1}{4N_x N_y} \frac{a}{\pi} \int_0^{4^{1/3}\pi/a} dQ \frac{Q^2 Q_D^2}{Q^4 + \lambda^2 Q_D^4} \\ &= \frac{(3/\pi)^{1/3}}{2N_x N_y} \int_0^\zeta du \frac{u^2}{u^4 + \lambda^2}, \end{aligned} \quad (\text{D3})$$

where the variable u is Q/Q_D , and the upper limit is $\zeta = (\pi/6)^{1/3}$. Using similar algebra to that of Appendix A, the answer is

$$\frac{\kappa_{00}(q_{\min})}{\kappa_0} = \frac{(3/\pi)^{1/3}}{2N_x N_y} [g(\zeta) - g(0)], \quad (\text{D4})$$

where the indefinite integral $g(u)$ is

$$\begin{aligned} g(u) &= \frac{1}{2\sqrt{(2\lambda)}} \left[-\frac{1}{2} \log \left(\frac{u^2 + \sqrt{2\lambda}u + \lambda}{u^2 + \sqrt{2\lambda}u + \lambda} \right) \right. \\ &\quad \left. + \tan^{-1} \left(\frac{\sqrt{2\lambda}u}{\lambda - u^2} \right) \right]. \end{aligned} \quad (\text{D5})$$

Because N_z is large, $\lambda = 2\pi\ell_{\min}/N_z a$ is small, and to first order, the definite integral $g(\zeta) - g(0)$ is determined

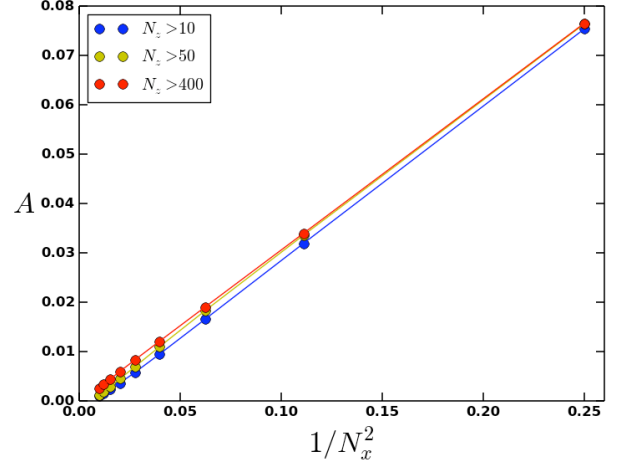


FIG. 9. The coefficient of the singular part $\propto 1/\sqrt{(qa)}$ scales nicely with the reciprocal of the transverse dimension $1/N_x^2$ of the simulation cell, especially when the semi-empirical fit is restricted to smaller values of q .

by the phase of the arctangent, and equals $\pi/2\sqrt{(2\lambda)}$, insensitive to the details of the “pseudo-Brillouin zone” boundary location. Thus to leading order, the piece of $\kappa(q)$ coming from the $Q_x = Q_y = 0$ part of the Q -mesh, is

$$\frac{\kappa_{00}(q_{\min})}{\kappa_0} = \frac{(3/\pi)^{1/3}}{2N_x N_y} \frac{\pi}{4} \sqrt{\frac{L}{\pi\ell_{\min}}} \quad (\text{D6})$$

The Debye model is reliable as a guide for the low frequency behavior, provided the relaxation-time approximation and the associated power law p are correct. The conclusion is a bit surprising. It indicates that if anisotropic simulation cells are used for NEMD simulation of κ , then the extrapolation to very long cells suffers from an unintended 1D singularity. The product $N_x N_y$ should increase at least as rapidly as $N_z^{1/2}$ to prevent this term distorting the extrapolated answer. In actual simulations, this is probably more a sobering thought than a serious warning. But the effect is real, and shows up in the Debye-model numerics shown in Figs. 6 and 7.

Appendix E: Semi-empirical Fitting

Figures 6 and 7 show (as red lines) Debye-RTA discrete Q -sums adjusted to fit NEMD results. These are intended to guide extrapolation, but reveal possible ambiguity about how to correct for the $1/\sqrt{(qa)}$ behavior. This is an artifact of too large a ratio N_z/N_x , not achieved in the NEMD simulations, but easily achieved in the Debye-RTA numerical sums on discrete Q -meshes. Here we show some results of a 3-term fit to the numerical Debye-RTA sums. From Appendices C and D, we are led to expect behavior of the type

$$\kappa(q)/\kappa_0 \approx A/\sqrt{qa} + B + C\sqrt{qa}. \quad (\text{E1})$$

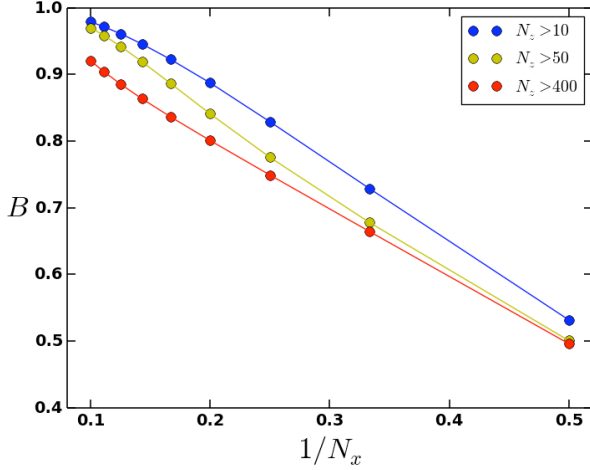


FIG. 10. The expected $q \rightarrow 0$ limit of Debye-RTA numerical sums corresponds to coefficient $B \rightarrow 1$, which happens nicely as the transverse cell size is increased to 10×10 . The fact that it happens most quickly when even short cells $N_z = 10$ are included in the fit, suggests fortuitous error cancellation.

The coefficient A should diminish as $N_x N_y$, since it arises from only the $Q_x = 0$ and $Q_y = 0$ part of the normal mode spectrum, one part out of the total of $N_x N_y$ Q_x , Q_y -values. This is tested in Fig. 9, and found correct. Numerical summation over the Q -mesh was done for 7 choices of $N_x = N_y$, namely 2, 3, 4, 5, 6, 8, and 10, and for a mesh of N_z ranging from the coarse value of 10 to the dense value of 10,000. The 7 resulting $\kappa_{\text{Debye}}(q)$ curves were least-squares fitted to Eq. E1. The fitting was done for the smallest $q = 2\pi/N_z a$, up to a cutoff (all N_z greater than a minimum value, chosen as 10, 50, or 200.) The fit is extremely accurate for the smallest q -cut ($N_z = 200$) and least accurate for the largest ($N_z = 10$.) The scaling with $1/N_x^2$ behaves just as expected. The singular up-turn at very small q is indeed an artifact of an anisotropic simulation cell, and should be avoided by not letting the ratio $N_z/(N_x N_y)$ become too large.

The coefficient B is supposed to represent the true converged $q \rightarrow 0$ limit of $\kappa(q)$. Figure 10 shows how this coefficient behaves as the transverse dimension increases. For $N_x = N_y = 10$, the Debye-RTA calculation is converged to 96% or better, no matter what range of q is used for least-squares fitting. Even down to $N_x = N_y = 5$, the value of B is 95% of the bulk value 1, provided all cell sizes down to the smallest ($N_z = 10$) are included in the least-squares fit, whereas, if only small q 's ($N_z > 200$) are used, convergence is a lot slower. This odd result indicates that the improved convergence found by fitting the less relevant large q 's is likely to be partially a result of a fortuitous cancellation of errors.

Finally, the leading finite size effect for bulk $\kappa(q)$ is contained in the term $C\sqrt{qa}$. This term comes long-wavelength phonons whose mean-free path exceeds the cell size unless the cell is large in all three directions. This negative contribution to κ is apparently suppressed

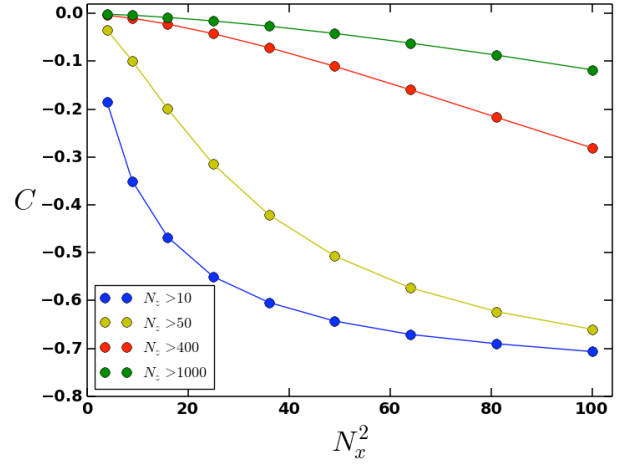


FIG. 11. The coefficient C of the leading finite size correction $\kappa(q)/\kappa_0 \approx 1 + C\sqrt{qa}$ is surprisingly sensitive to transverse simulation cell size and to the choice of q -range to use in the fit.

when the cell becomes more anisotropic. Simultaneously the diverging term is getting larger. These seem to accidentally compensate, making the NEMD answers better than the true convergence estimates lead one to expect. This is nicely illustrated in Fig. 5, which shows that for N_z not too big ($0.22 < \sqrt{qa} < 0.5$, corresponding to $25 < N_z < 130$), the curve of $\kappa(q)$ versus \sqrt{qa} is remarkably independent of N_x and appears to point smoothly to the correct extrapolated value of κ_0 . The more careful fitting to the 3-term expression gives extrapolated values shown by the dotted lines, which converge well for meshes $N_x \geq 6$, but not as rapidly as the less careful extrapolation. The prescription for extrapolation of NEMD with realistic simulation cells seems to be, don't use meshes with $N_x = N_y$ much smaller than 6, and extrapolate linearly if numerical results appear linear when plotted against $\sqrt{qa} \propto 1/\sqrt{N_z}$.

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