

**Barbalinardo *et al.* Reply:** Comment [1] shows that the thermal conductivity ( $\kappa$ ) of a (10,0) carbon nanotube (CNT), obtained by inversion of the linearized Boltzmann transport equation (BTE), may not converge when the third-order interatomic force constants (IFC) are computed analytically. Bruns *et al.* [1,2] showed that the lifetimes  $\tau_{\text{RTA}}(q)$  of the acoustic branches, computed in the relaxation time approximation (RTA), obey precise power laws for  $q \rightarrow 0$ , which are violated by the numerical IFCs in our Letter [3].  $\kappa$  in Ref. [1] is much larger than that in Ref. [3], and does not converge even when 400  $q$ -points are used to integrate the first Brillouin zone. These new results are correct and challenge the agreement between BTE (with classical statistics) and molecular dynamics (MD) proposed in Ref. [3], and the statement that  $\kappa$  is finite in the bulk limit.

Here we address the accuracy of finite-differences (FD) calculations, the convergence of  $\kappa$  for infinitely long CNTs, and  $\kappa(L)$  for CNTs of length  $L$ . Bruns *et al.* shows that the long-wavelength trends of  $\tau_{\text{RTA}}$  are disrupted by numerical errors introduced in the FD calculation of the IFCs resulting in nonsymmetrized force constant matrices. We computed IFCs by FD with displacements  $dx$  ranging from  $10^{-6}$  to  $10^{-3}$  Å, and the correct trends for  $\tau_{\text{RTA}}$  of the flexural (FA), longitudinal (LA) and torsional (TW) acoustic modes at low  $q$  are recovered for  $dx > 4 \times 10^{-5}$ . Figure 1 shows that  $\tau_{\text{RTA}}(q)$  from numerical IFCs with  $dx = 10^{-4}$  Å are nearly indistinguishable from those obtained from analytical IFCs (in Ref. [3] we used  $dx > 4 \times 10^{-6}$  Å). This result is useful for future works for those cases in which it is not possible to compute IFCs analytically.

$\tau_{\text{RTA}}$  provides a compelling test of the IFC's accuracy but cannot be used to calculate  $\kappa$  of CNTs due to the predominance of hydrodynamic effects in low-dimensional materials [4]. We computed the lifetimes of the acoustic modes by directly inverting the scattering matrix ( $\tau_{\text{inv}}$ ), using both analytical and numerical IFCs [Fig. 1]. Because of the hydrodynamic effects, the values and low- $q$  trends of  $\tau_{\text{inv}}$  and  $\tau_{\text{RTA}}$  are different. Fitting  $\tau_{\text{inv}}$  for  $q \rightarrow 0$  one may infer whether  $\kappa$  converges or diverges in the infinite-size limit. The modes that may lead to divergence are FA and TW, as their  $\tau_{\text{inv}}$  diverge for  $q \rightarrow 0$ . The group velocity of the FA modes is proportional to  $q$ , and that of TW is constant; thus the onset of  $\kappa$  divergence is either  $\tau_{\text{FA}} \propto 1/q^3$  or  $\tau_{\text{TW}} \propto 1/q$ .  $\tau_{\text{TW}}$  has a similar low- $q$  behavior as in RTA calculations,  $\tau_{\text{inv,TW}} \propto 1/q^{0.5}$ .  $\tau_{\text{inv,FA}}$  from analytical IFCs diverges as  $1/q^3$  or steeper, suggesting divergent  $\kappa$ . Conversely, using numerical IFCs we get a slower than  $1/q^3$  trend, which would lead to finite  $\kappa$ . This calculation may be more representative of systems at finite temperature, possibly reconciling BTE with convergent  $\kappa$  in MD simulations. Calculations with finite temperature IFC renormalization and thicker  $q$ -point sampling will be necessary to prove the convergence of  $\kappa$  in CNTs.

Finally, we computed finite-length  $\kappa(L)$  with analytical IFCs. We find that  $\kappa(L)$  remains in agreement with our

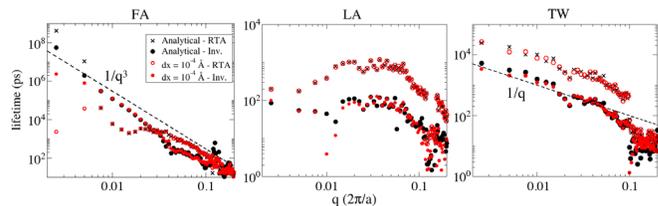


FIG. 1. Lifetimes of the FA, LA, and TW acoustic modes computed with quantum statistics. All the calculations are done by integrating the BTE over a uniform grid of 401  $q$  points.

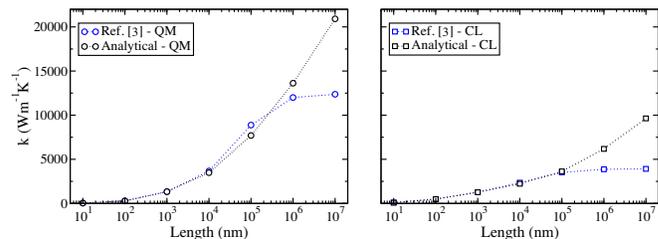


FIG. 2. Thermal conductivity as a function of length computed with analytical and the numerical IFCs from Ref. [3].

former calculations [3] up to  $L = 1$  mm while it gets significantly larger for larger lengths (Fig. 2).

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- [1] D. Bruns, A. Nojeh, A. S. Phani, and J. Rottler, preceding comment, Ultrahigh Convergent Thermal Conductivity of Carbon Nanotubes from Comprehensive Atomistic Modeling, *Phys. Rev. Lett.* **128**, 259601 (2022).
- [2] D. Bruns, A. Nojeh, A. S. Phani, and J. Rottler, Nanotube heat conductors under tensile strain: Reducing the three-phonon scattering strength of acoustic phonons, *Phys. Rev. B* **104**, 075440 (2021).
- [3] G. Barbalinardo, Z. Chen, H. Dong, Z. Fan, and D. Donadio, Ultrahigh Convergent Thermal Conductivity of Carbon Nanotubes from Comprehensive Atomistic Modeling, *Phys. Rev. Lett.* **127**, 025902 (2021).
- [4] A. Cepellotti, G. Fugallo, L. Paulatto, M. Lazzeri, F. Mauri, and N. Marzari, Phonon hydrodynamics in two-dimensional materials, *Nat. Commun.* **6**, 6400 (2015).