

# Evaluation of the **Virtual Crystal Approximation** for Predicting **Alloy** Vibrational Mode Properties and **Thermal Conductivity**

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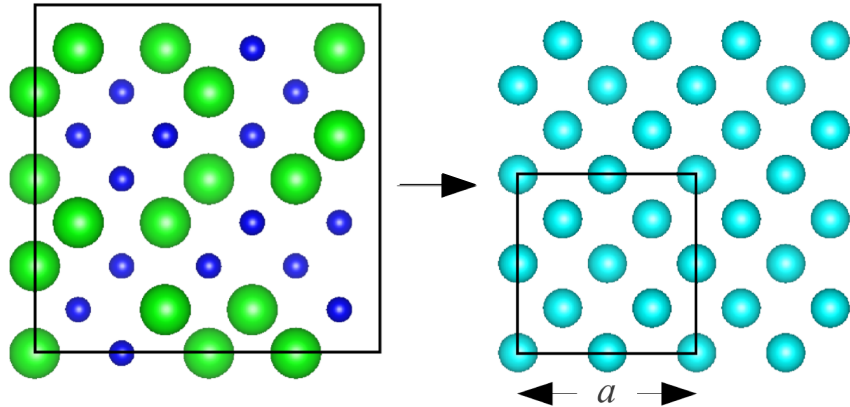
**<http://ntpl.me.cmu.edu/>**

04/04/13

# Virtual Crystal Approximation

Gamma

VC



$$\bar{m}^{\mu} = (1 - c)m^i + cm^j$$

$$k_{ph,\mathbf{n}} = \sum_{\boldsymbol{\kappa}} \sum_{\nu} \frac{k_B}{V} D_{ph,\mathbf{n}}(\boldsymbol{\kappa}_{\nu})$$

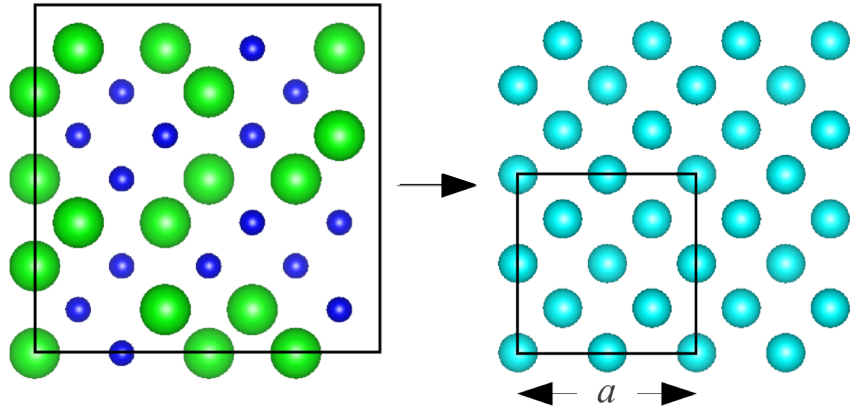
$$D_{ph,\mathbf{n}}(\boldsymbol{\kappa}_{\nu}) = v_{g,\mathbf{n}}^2(\boldsymbol{\kappa}_{\nu}) \tau(\boldsymbol{\kappa}_{\nu})$$

B. Abeles, Phys. Rev. 131, 19061911 (1963)

# VC-ALD Diffusivities: Lifetimes

Gamma

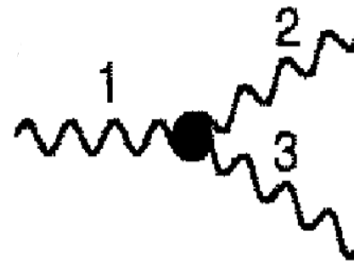
VC



$$D_{ph,n}(\kappa) = v_{g,n}^2(\kappa) \tau(\kappa)$$

Anharmonic Lattice Dynamics (**ALD**)

Tamura Phonon-Defect<sup>1</sup>



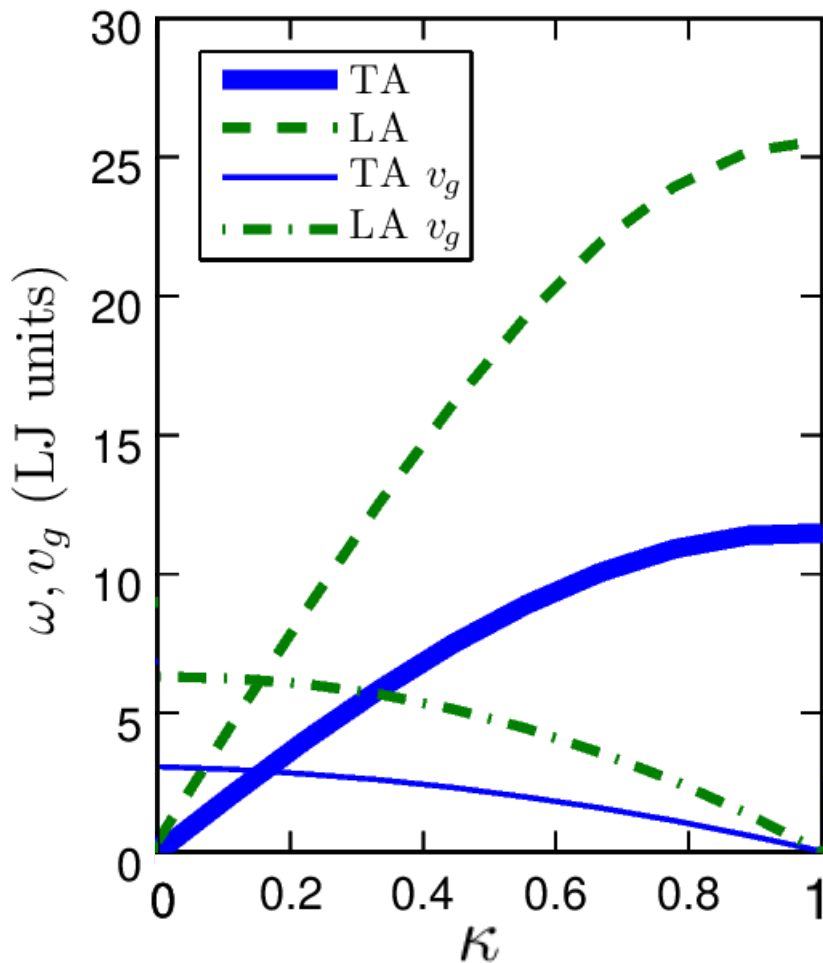
Perturbation theory:

Matthiessen's Rule:

$$\frac{1}{\tau(\kappa)} = \frac{1}{\tau_{p-p}(\kappa)} + \frac{1}{\tau_{p-d}(\kappa)}$$

<sup>1</sup>Phys. Rev. B 27, 858866 (1983)

# VC-ALD Diffusivities: Group Velocity

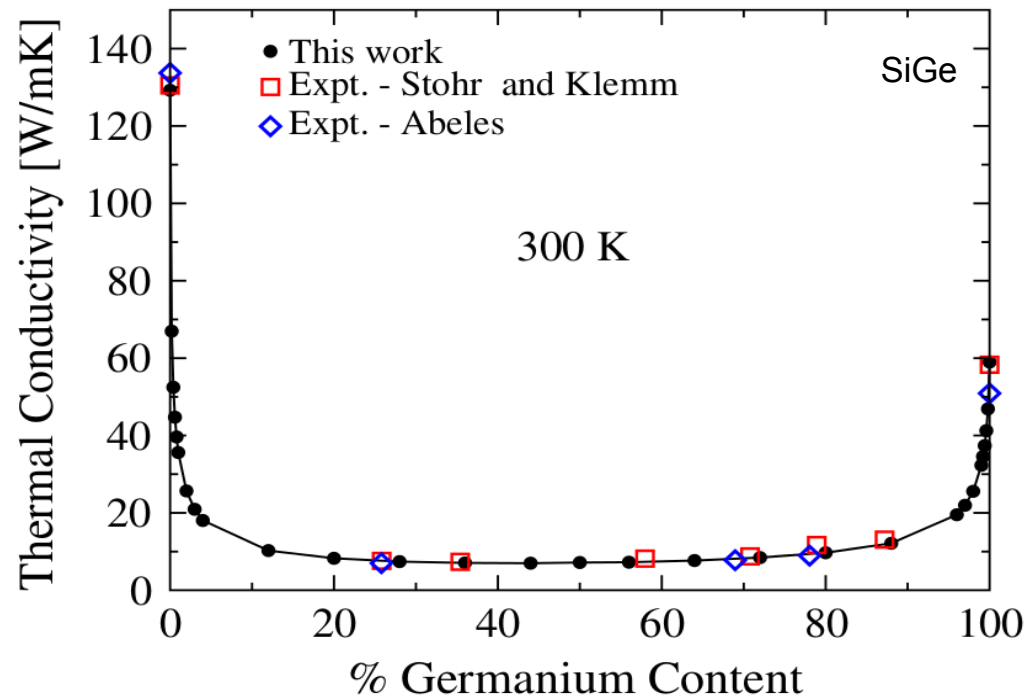


$$D_{ph,n}(\kappa_\nu) = v_{g,n}^2(\kappa_\nu) \tau(\kappa_\nu)$$

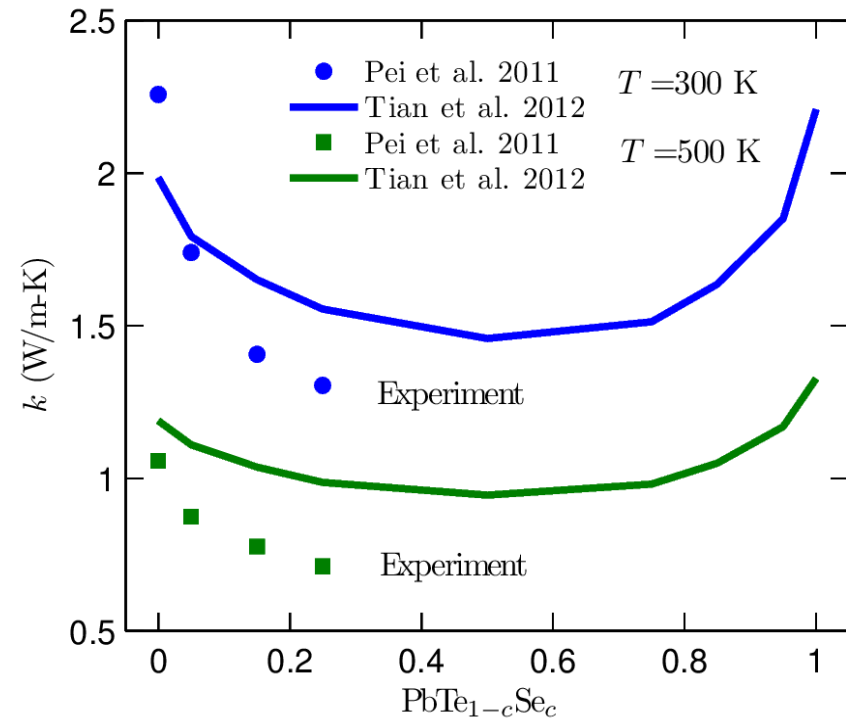
$$\mathbf{v}_{g,n}(\kappa_\nu) = \frac{\partial \omega(\kappa_\nu)}{\partial \kappa}$$

# VC-ALD: experimental accuracy

Density Functional Theory (DFT) + **(VC-ALD)**



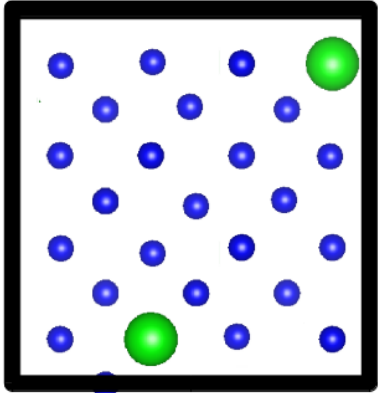
PRL 106, 045901 (2011)



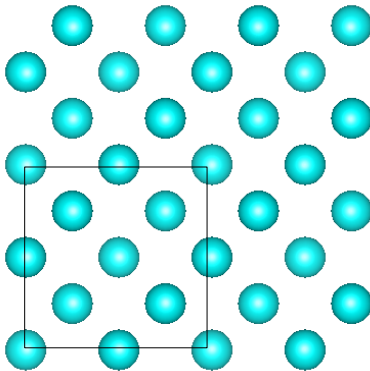
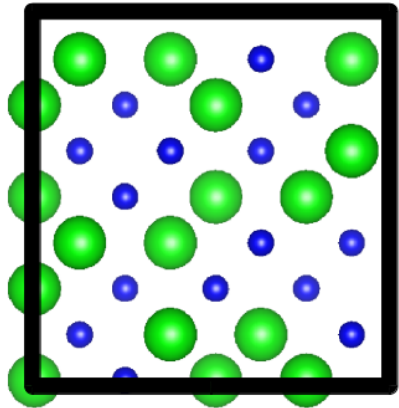
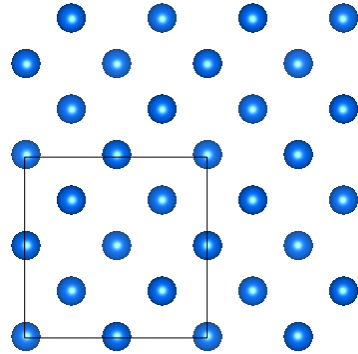
PRB 85, 184303 (2012)

# Explicit disorder: VC vs Gamma

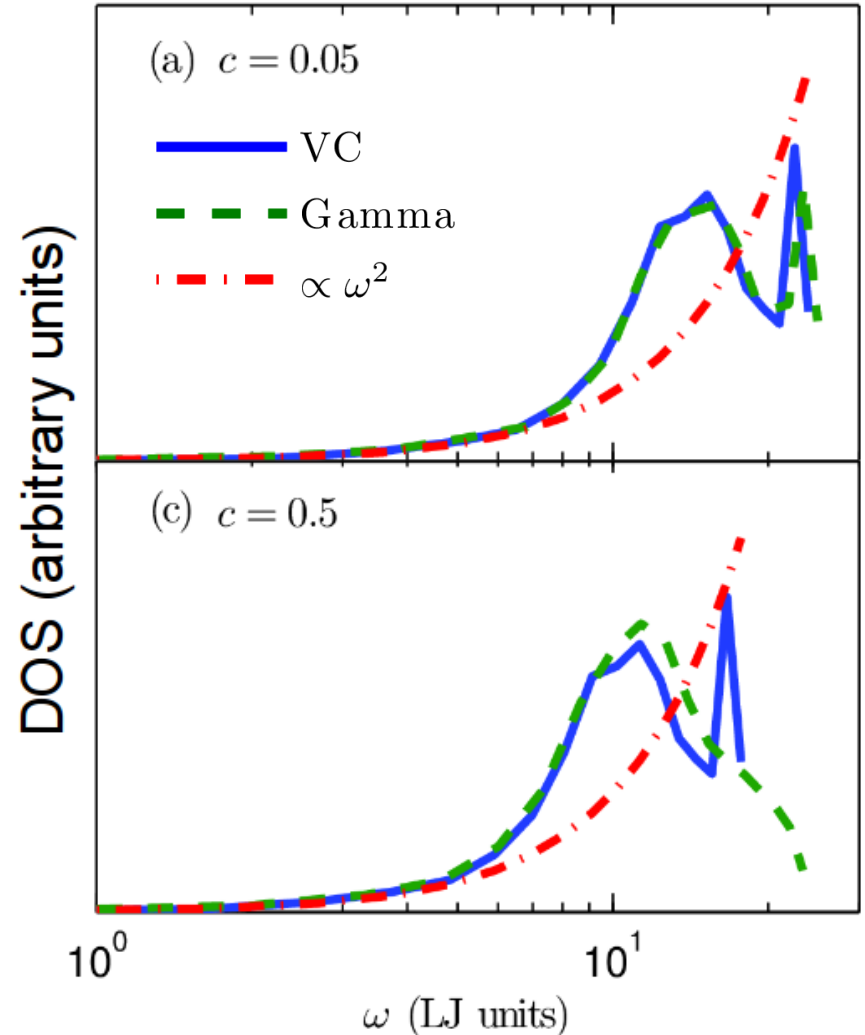
Gamma



VC

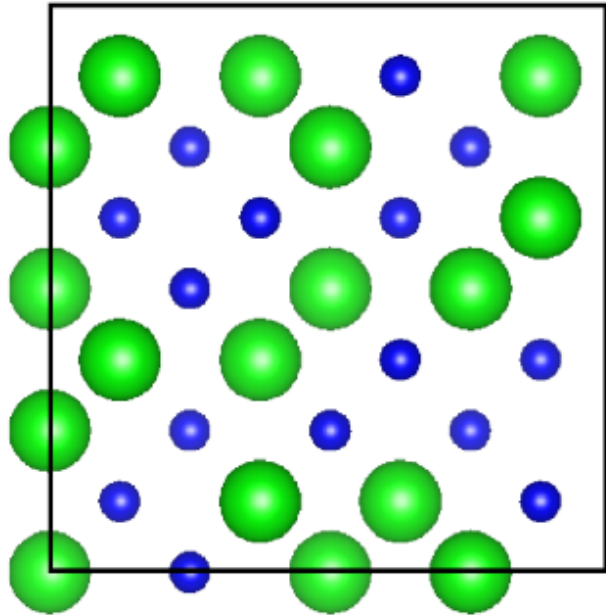


LJ Alloys

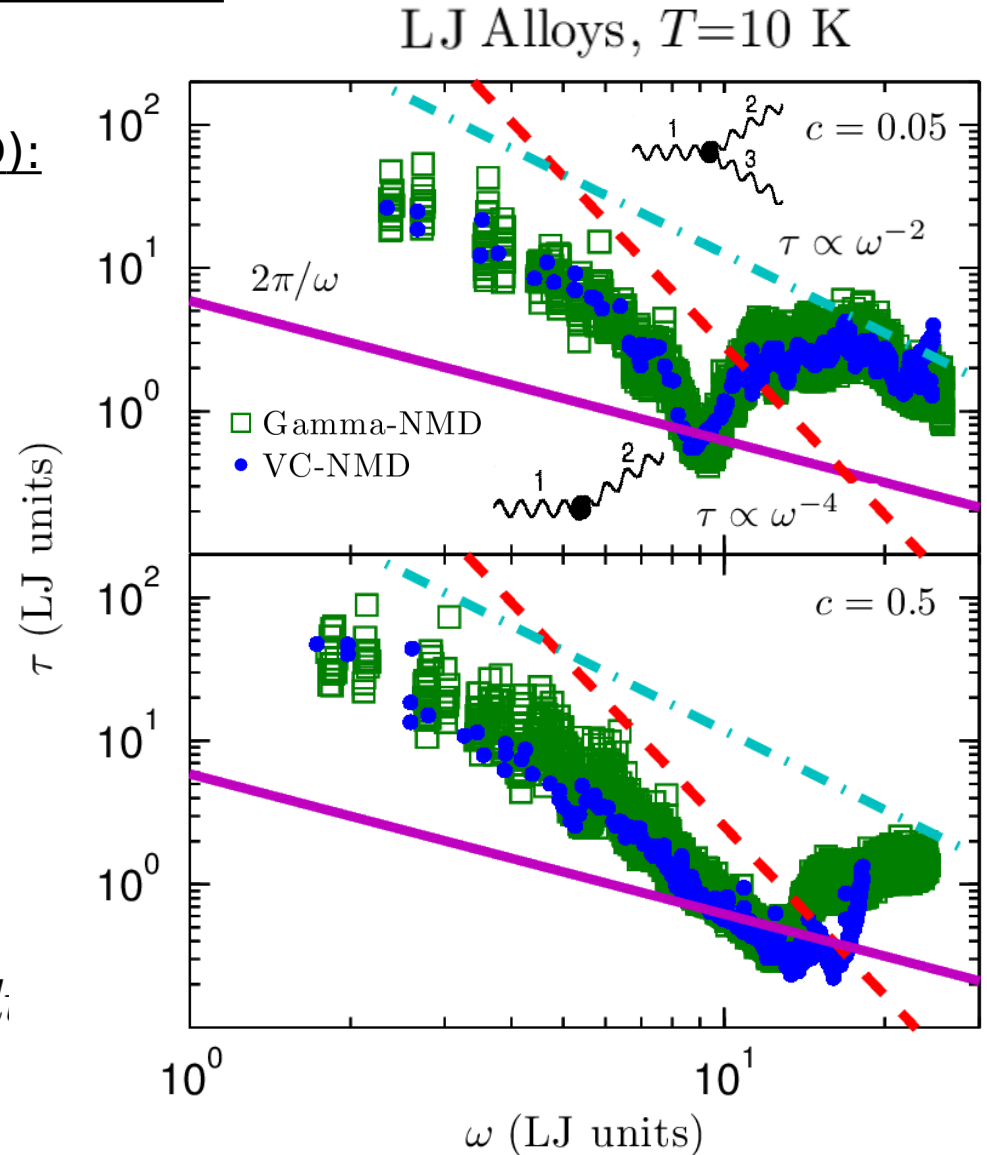


# Explicit disorder: NMD

Normal Mode Decomposition (**NMD**):  
Molecular Dynamics (MD)-based

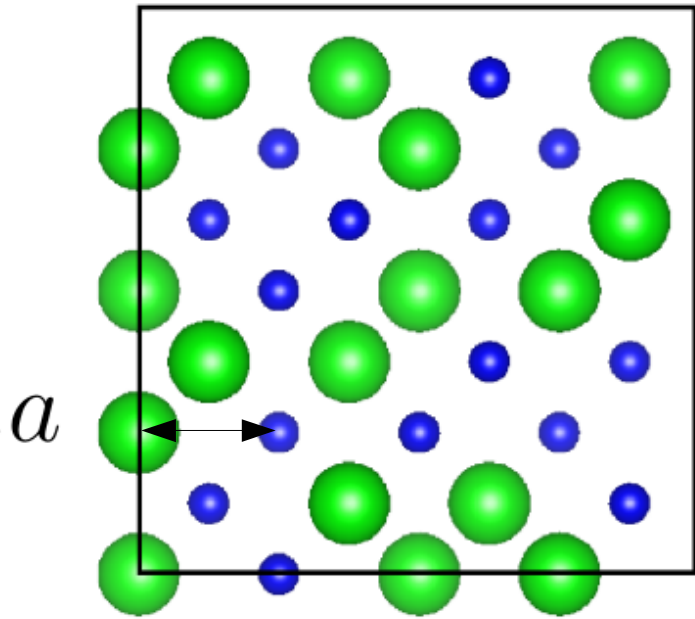


$$\tau(\kappa) = \int_0^{t^*} \frac{\langle E(\kappa; t) E(\kappa; 0) \rangle}{\langle E(\kappa; 0) E(\kappa; 0) \rangle} dt$$



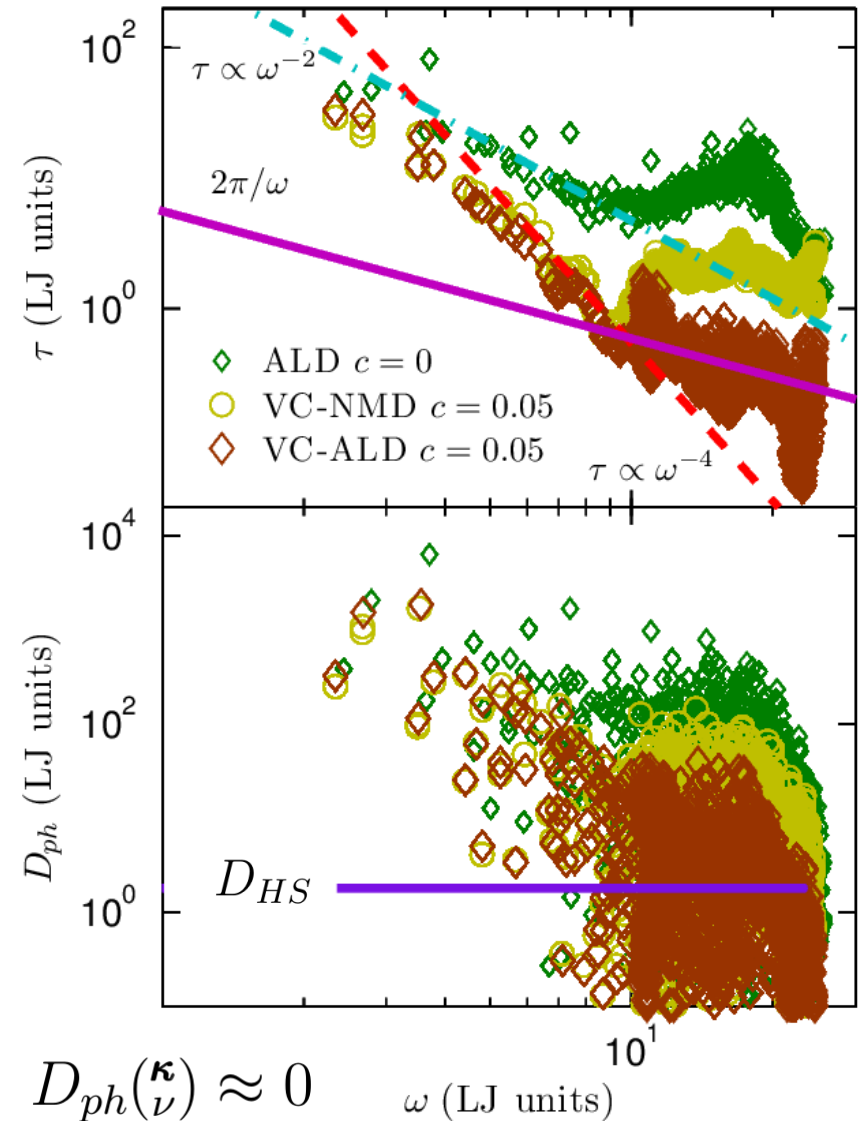
# VC Diffusivities

$$D_{ph,n}(\kappa) = v_{g,n}^2(\kappa) \tau(\kappa)$$



$$D_{HS} = \frac{1}{3} v_s a$$

LJ Argon and Alloys,  $T = 10$  K

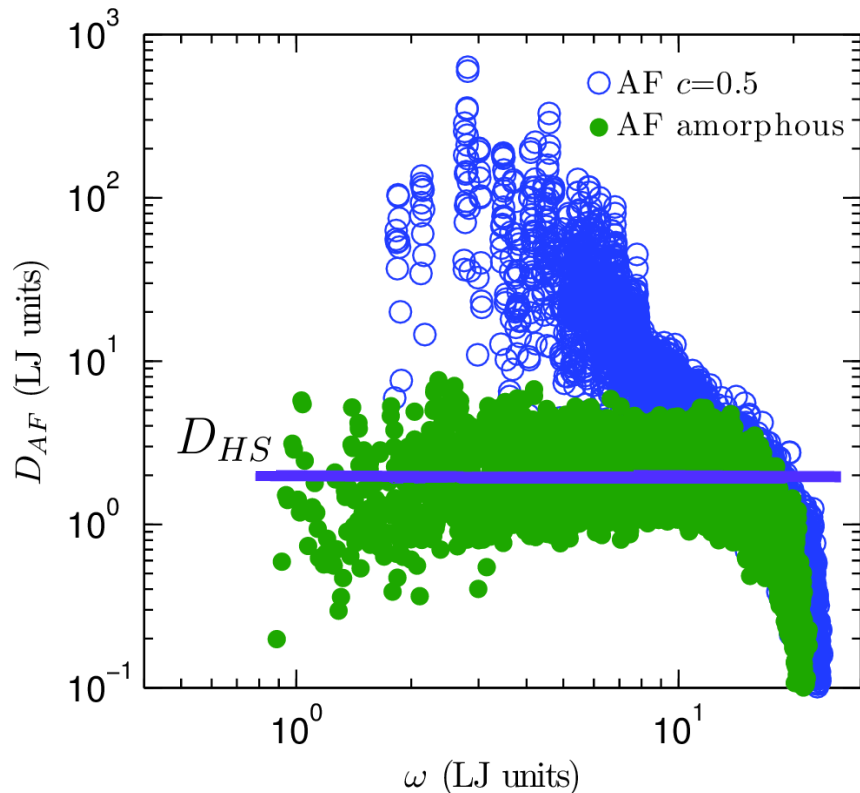




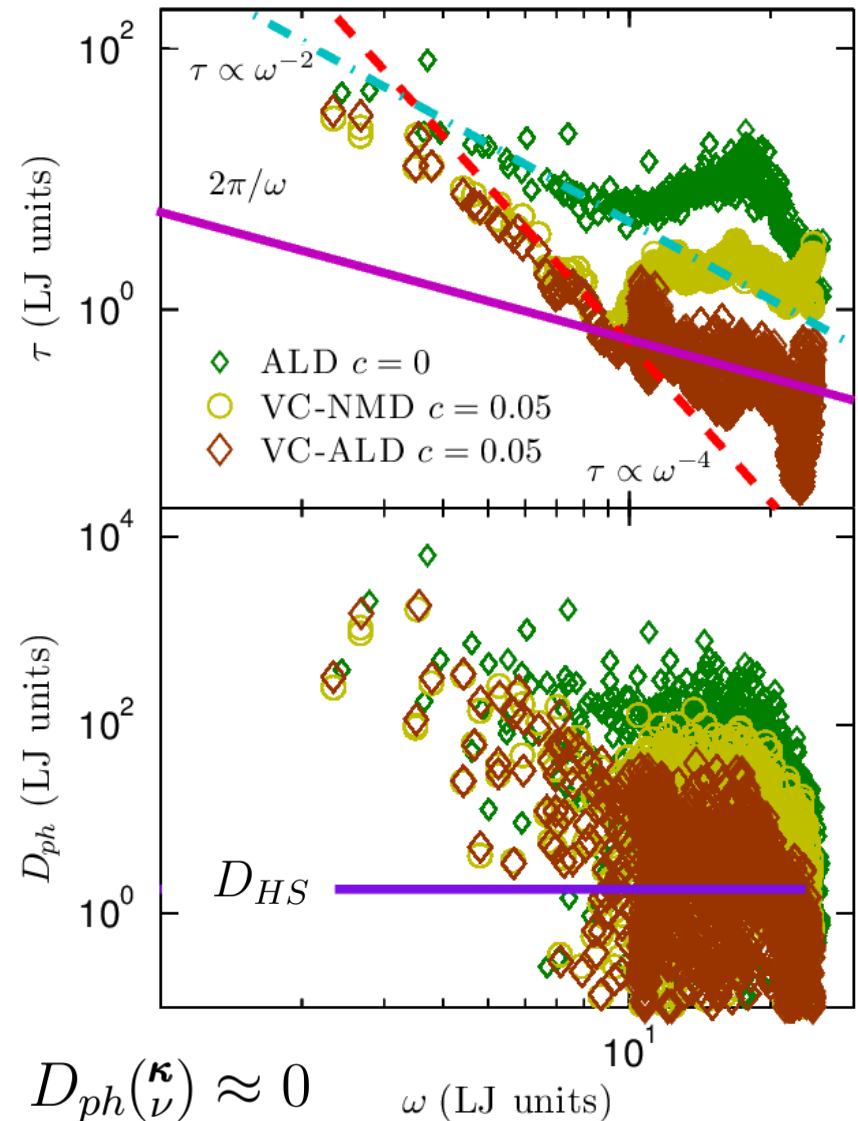
# AF Diffusivities

Allen-Feldman (AF) Theory:

$$k_{AF} = \sum_{\text{diffusons}} \frac{k_B}{V} D_{AF,i}(\omega_i)$$

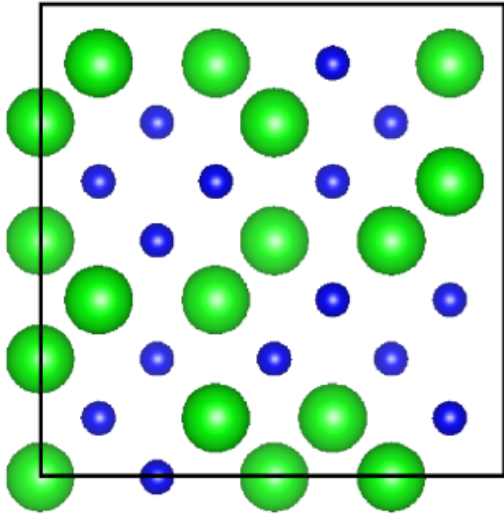


LJ Argon and Alloys,  $T = 10$  K



# Thermal conductivity

MD-based Green-Kubo (GK)

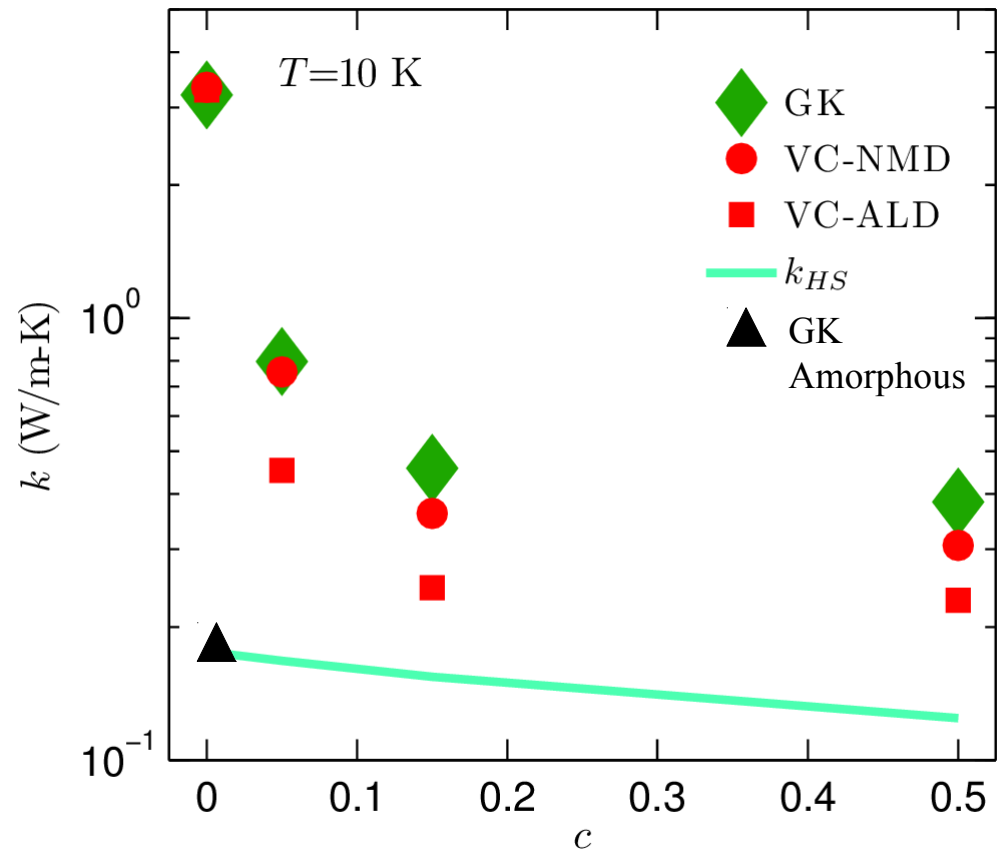


High-scatter adjustment\*:

$$D_{ph}(\kappa) < D_{HS}$$

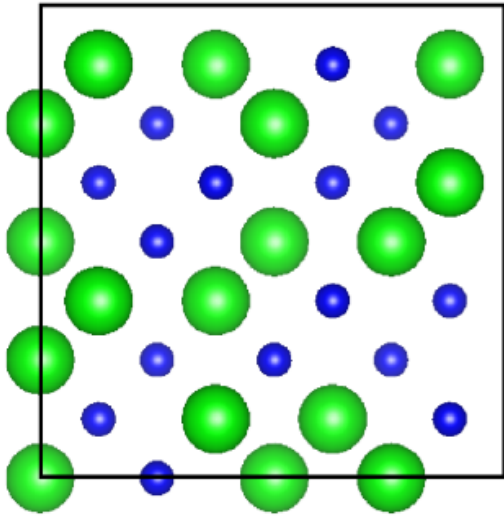
$$D_{ph}(\kappa) = D_{HS}$$

LJ Argon and Alloys,  $T = 10$  K



# Thermal conductivity

MD-based Green-Kubo (GK)

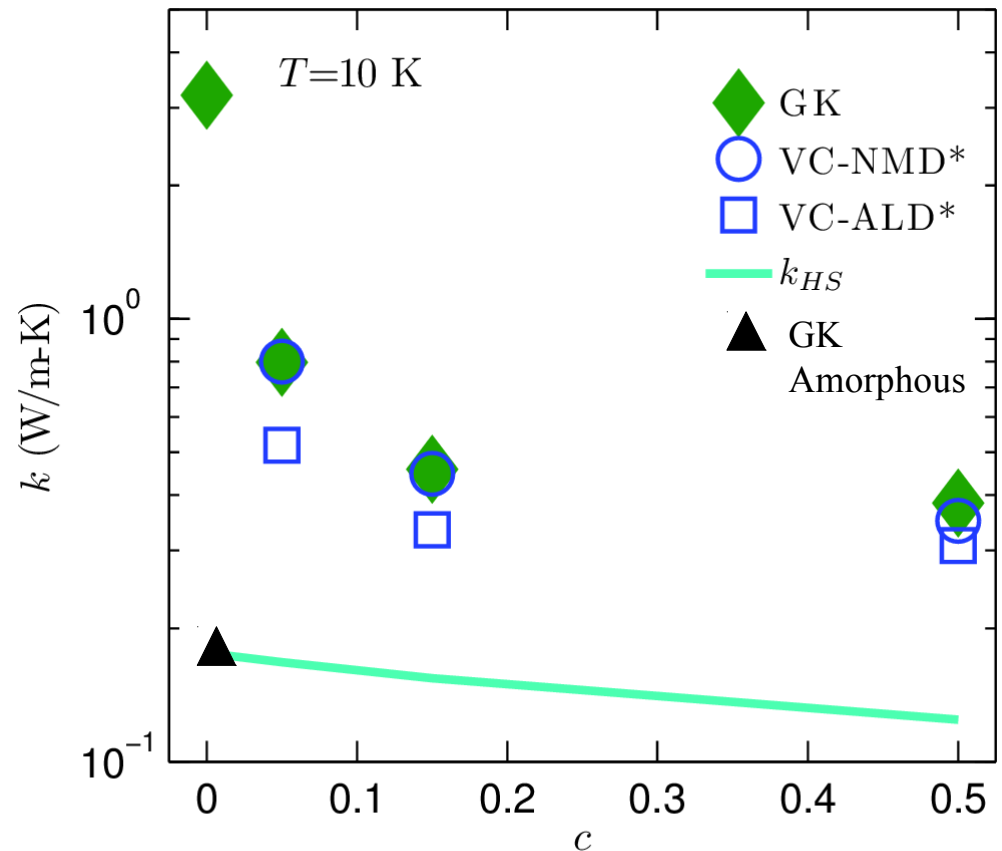


High-scatter adjustment\*:

$$D_{ph}(\boldsymbol{\kappa}) < D_{HS}$$

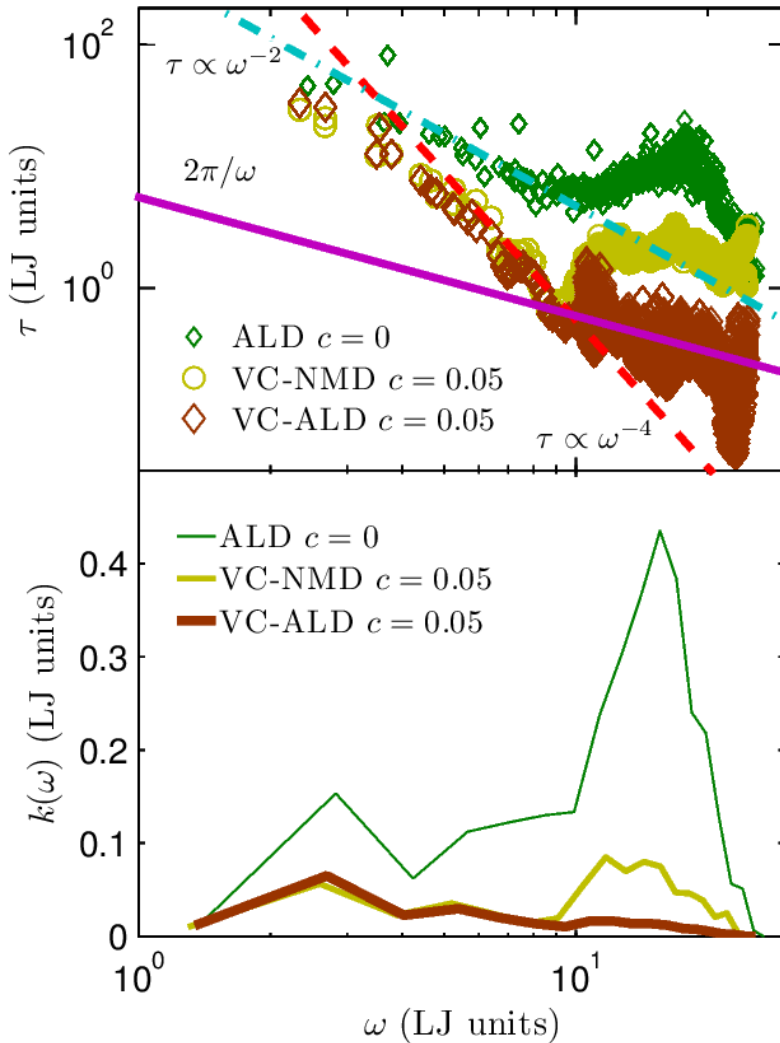
$$D_{ph}(\boldsymbol{\kappa}) = D_{HS}$$

LJ Argon and Alloys,  $T = 10$  K

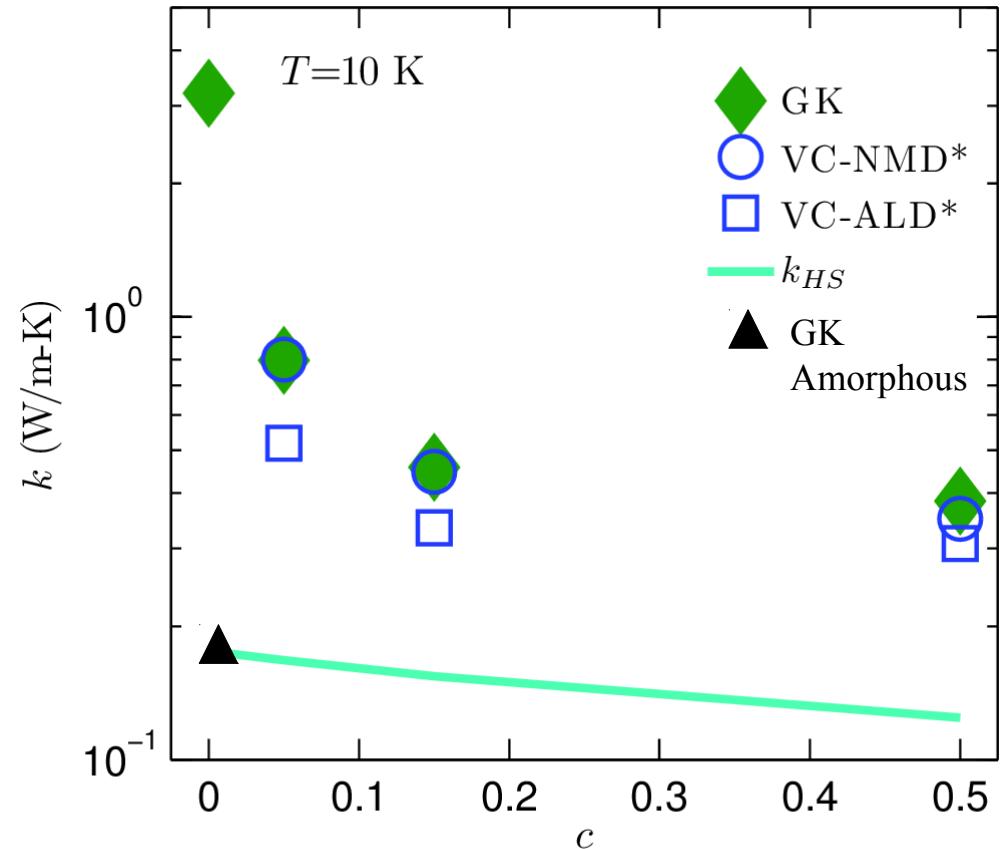


# Thermal conductivity spectrum

LJ Argon and Alloys,  $T = 10$  K



LJ Argon and Alloys,  $T = 10$  K



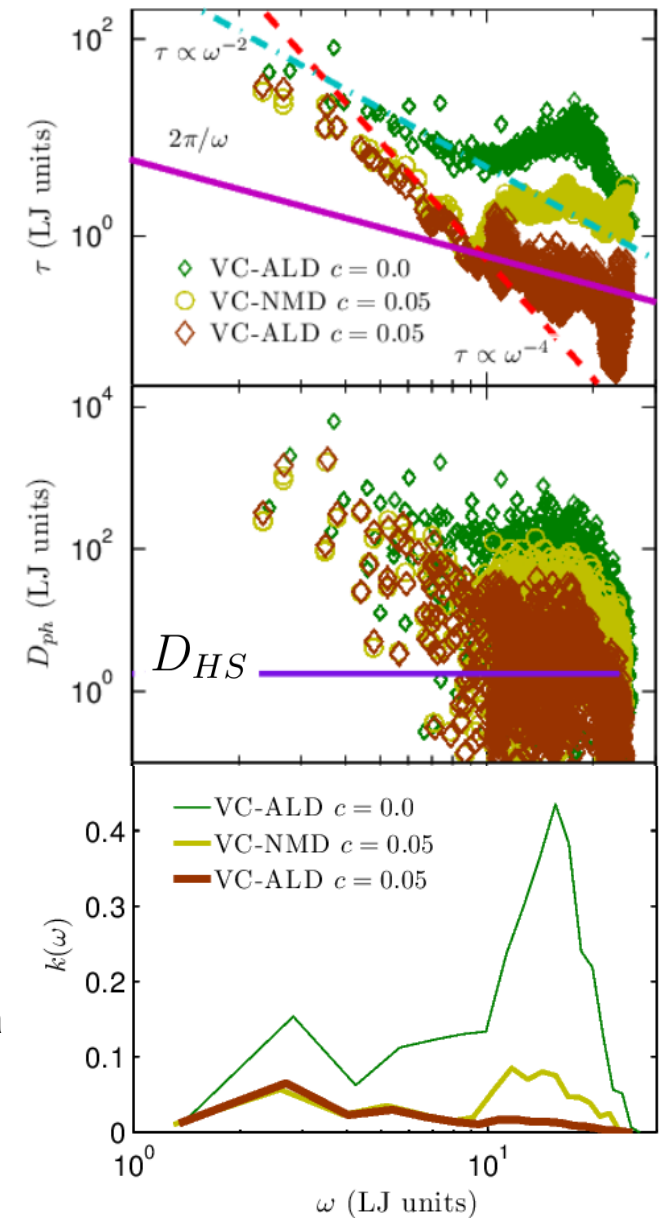
# Summary

**VC approximation underpredicts mode group velocities** at high frequency.

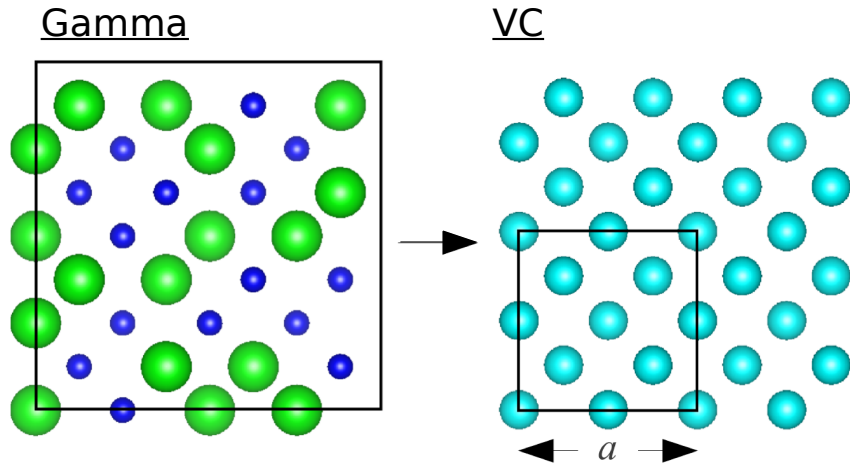
**VC-ALD underpredicts lifetimes** at high-frequency.

**Breakdown of VC-ALD method** is likely for materials near HS limit.

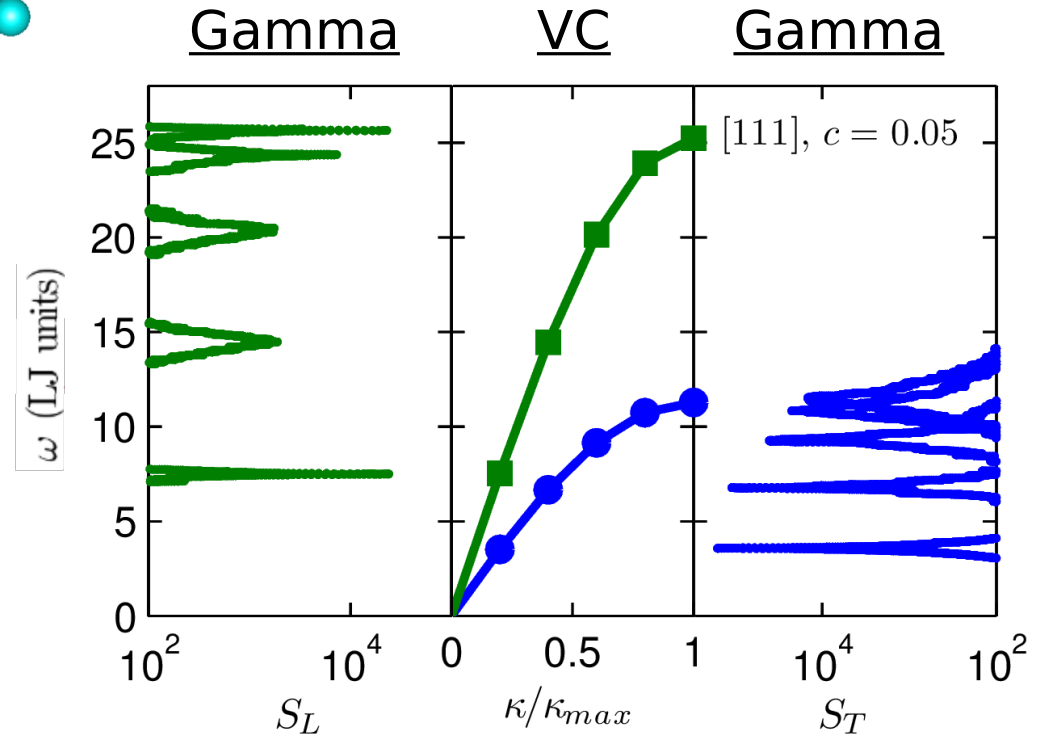
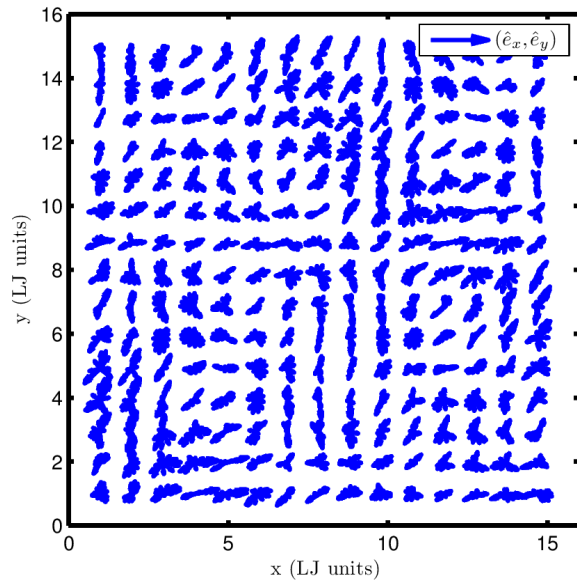
This work was supported by AFOSR award FA95501010098 and by a grant of computer time from the DOD High Performance Computing Modernization Program at the US Army Engineer Research and Development Center. We thank Davide Donadio, Jivtesh Garg, Asad Hasan, Craig Maloney, and Zhiting Tian for helpful discussions.



# Explicit disorder: Structure Factor

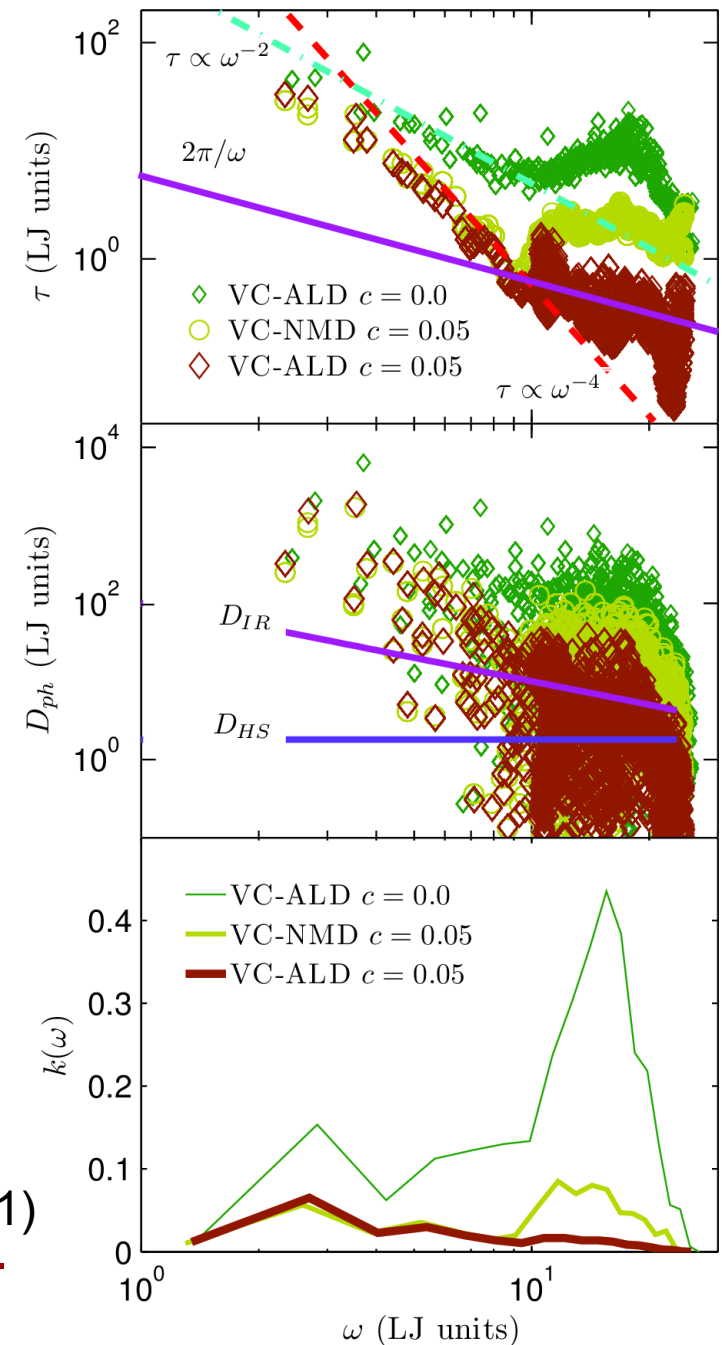
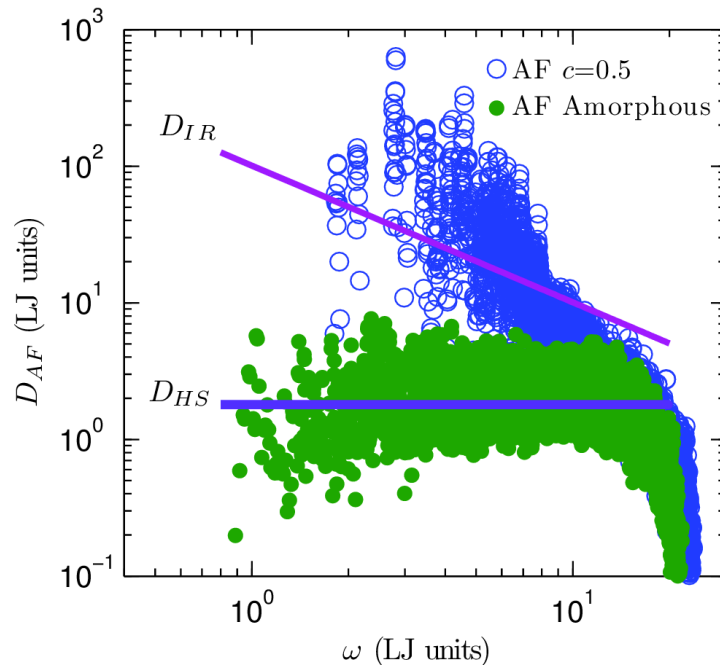


$$D_{ph,n}(\kappa) = v_{g,n}^2(\kappa) \tau(\kappa)$$



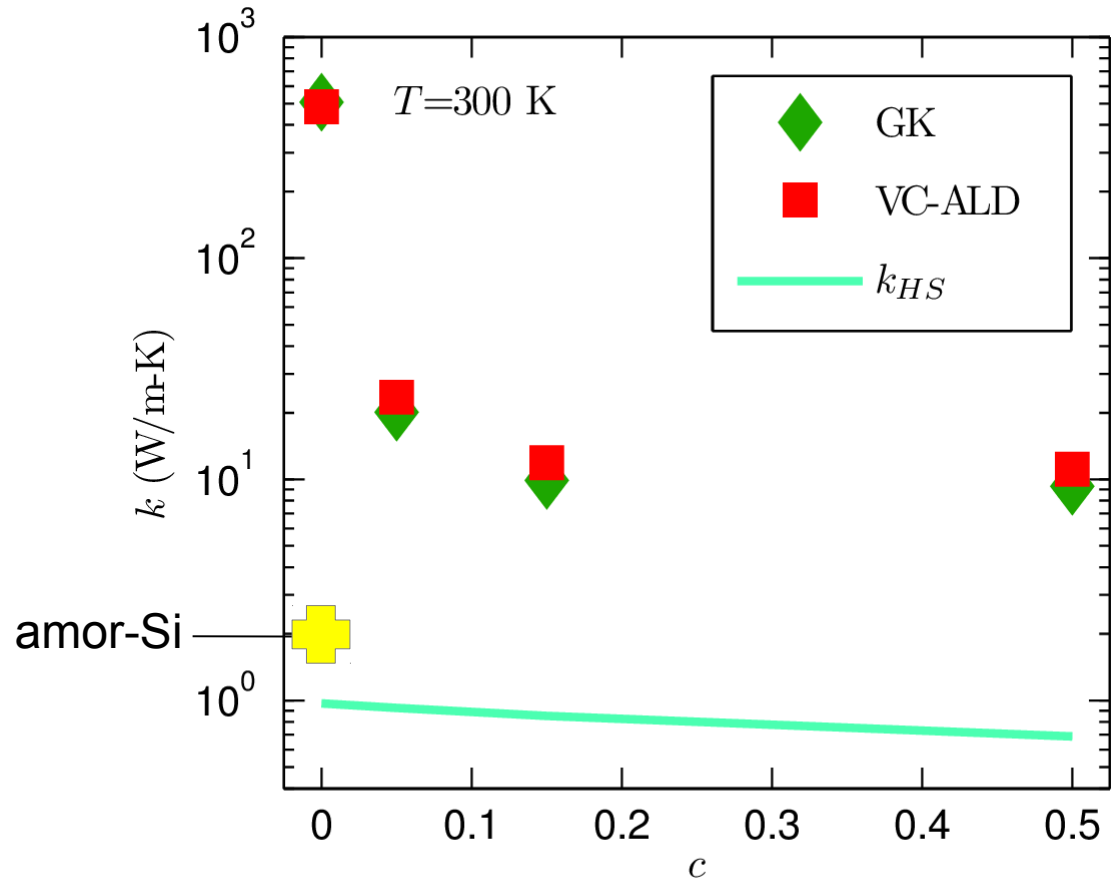
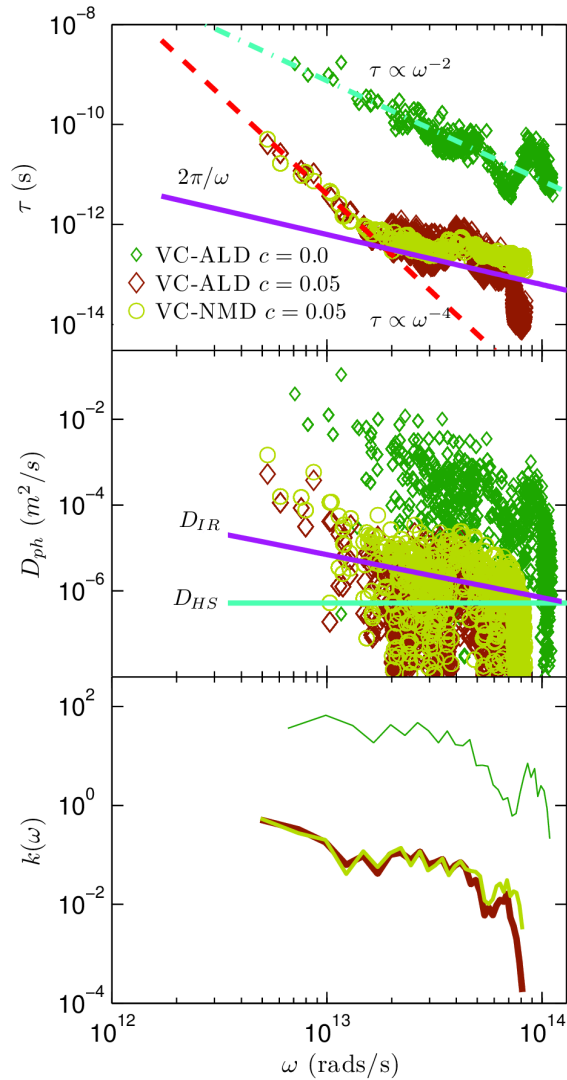
# HS/IR Limit

$$D_{IR} = \frac{2\pi}{3} \frac{v_s^2}{\omega}$$



P. Sheng and M. Zhou, Science 253, 539542 (1991)

# Thermal conductivity: SW silicon alloy

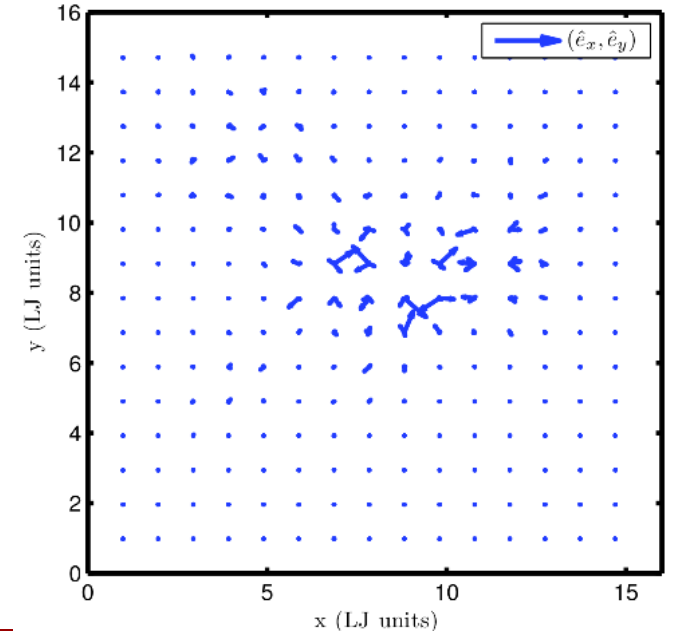
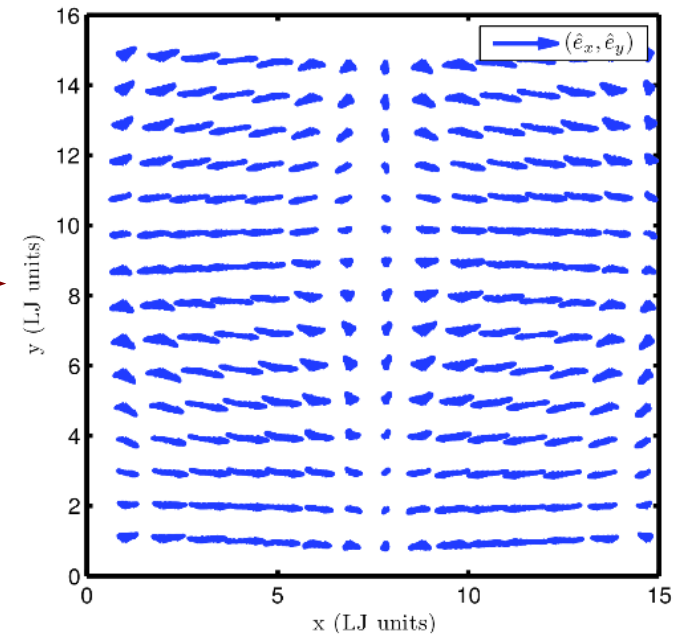
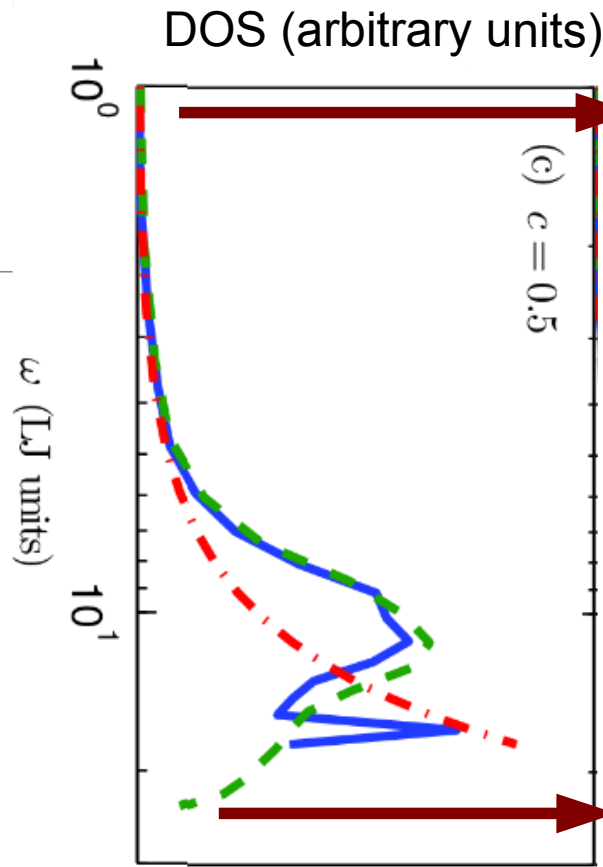
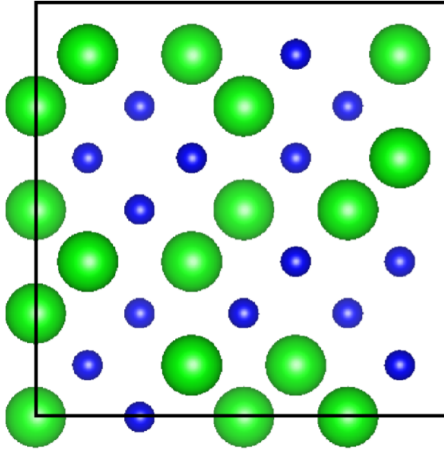




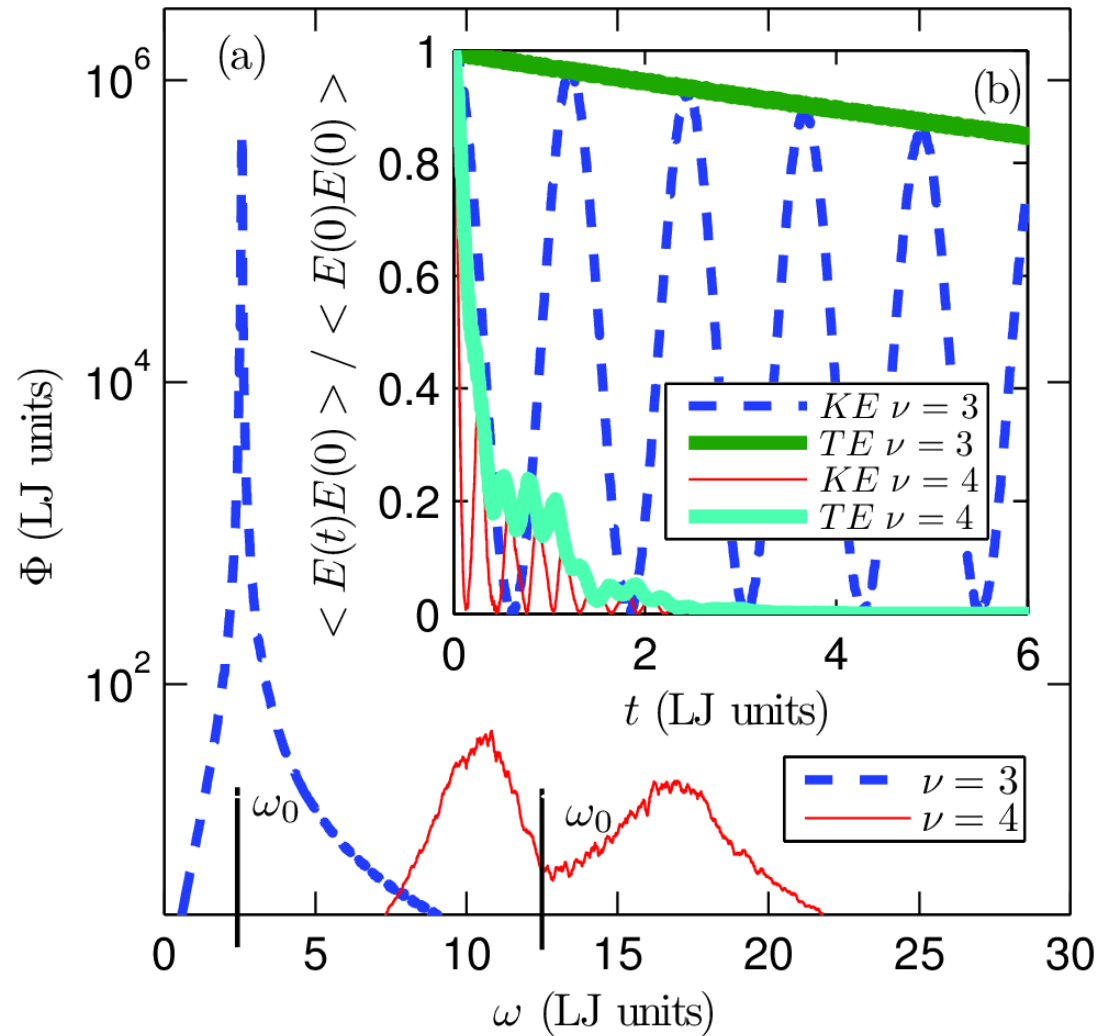
# Gamma modes

$$e\left(\begin{smallmatrix} \kappa=0 & b \\ \nu & \alpha \end{smallmatrix}\right)$$

(a) disordered supercell



# NMD using VC modes



# Phonon Spectrum: LJ Ar vs SW Si

