

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

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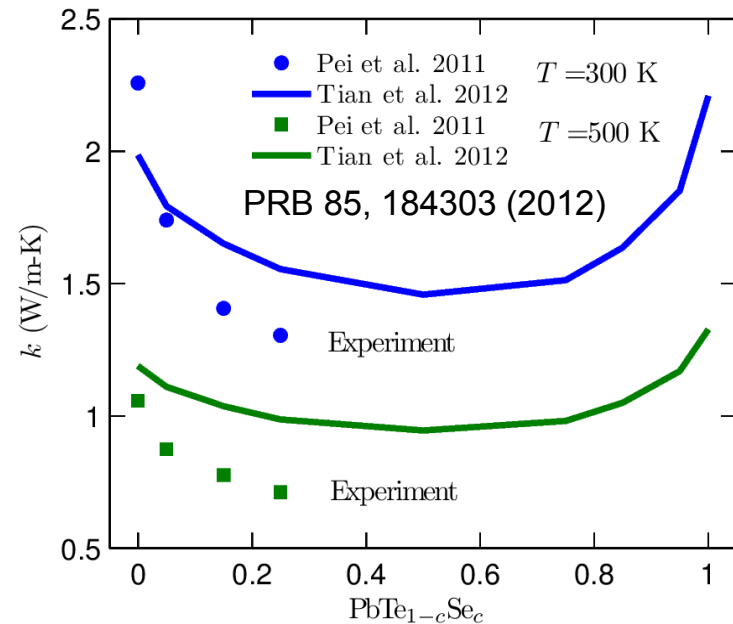
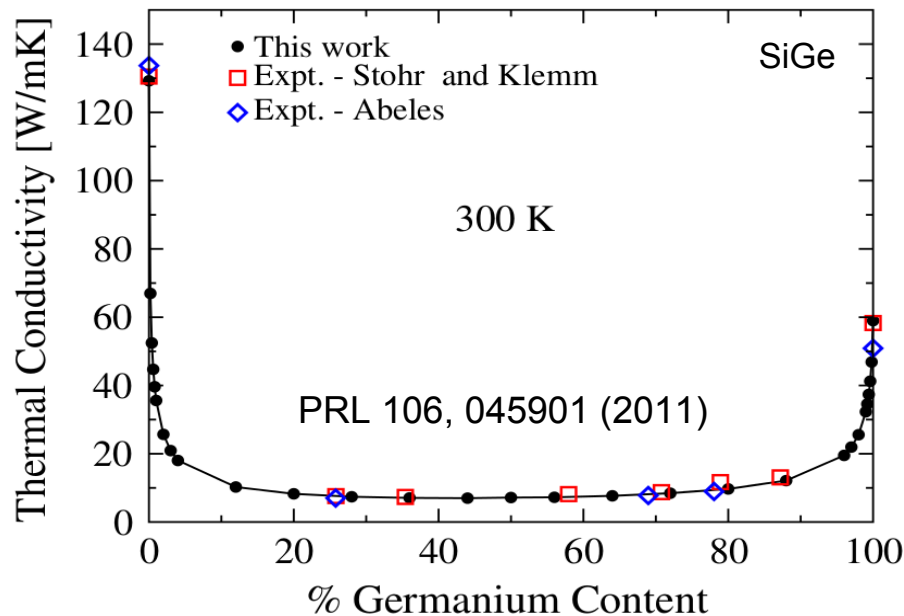
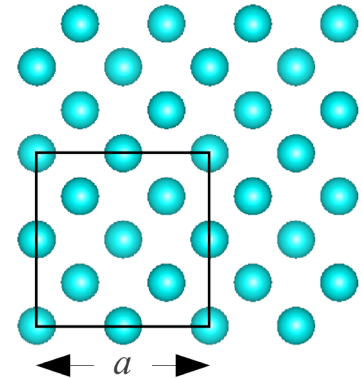
[\*\*http://ntpl.me.cmu.edu/\*\*](http://ntpl.me.cmu.edu/)

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# Motivation: experimental accuracy

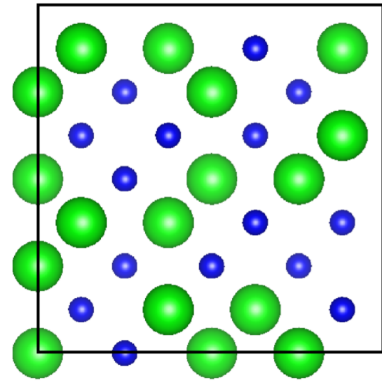
Density Functional Theory (DFT) + **(VC-ALD)**

**(VC-ALD)** = **(VC)** Virtual Crystal approximation  
+  
**(ALD)** Anharmonic Lattice Dynamics

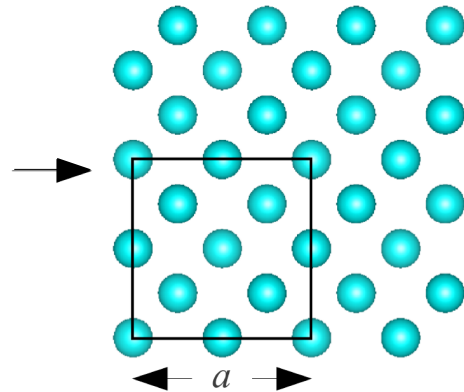


# Virtual Crystal Approximation

Gamma



VC



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

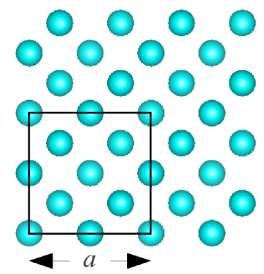
$$k_{ph,\mathbf{n}} = \sum_{\mathbf{\kappa}} \sum_{\nu} \frac{k_B}{V} D_{ph,\mathbf{n}}(\mathbf{\kappa}_{\nu})$$

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

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

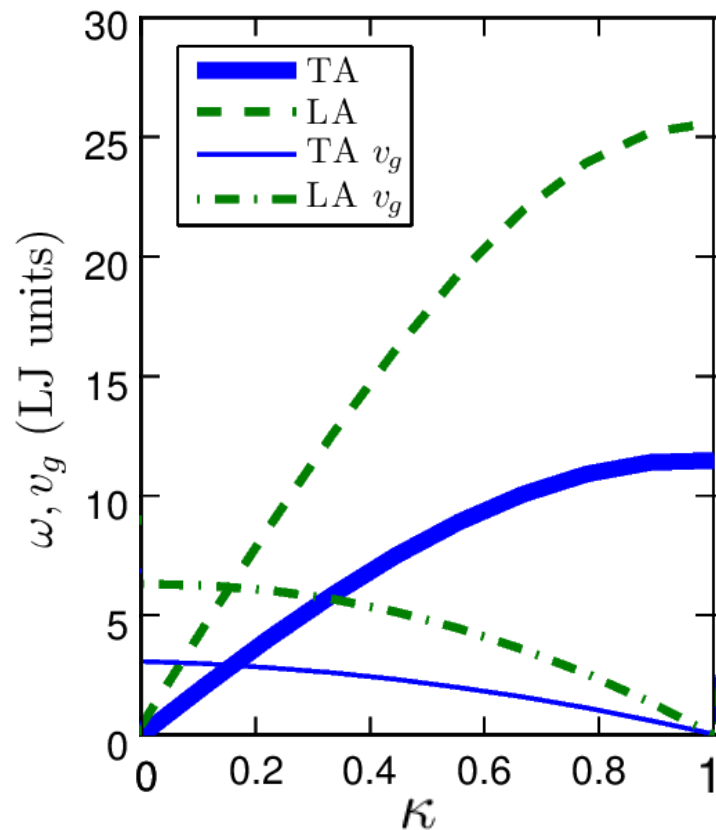
# VC-ALD Diffusivities

VC Unit Cell



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

$$v_{g,n}(\kappa) = \frac{\partial \omega(\kappa)}{\partial \kappa}$$



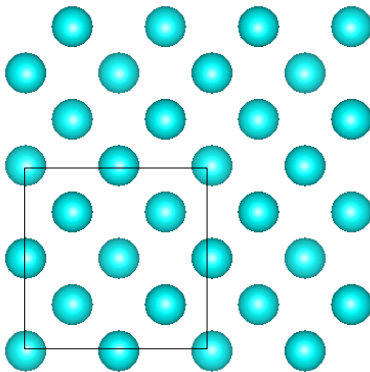
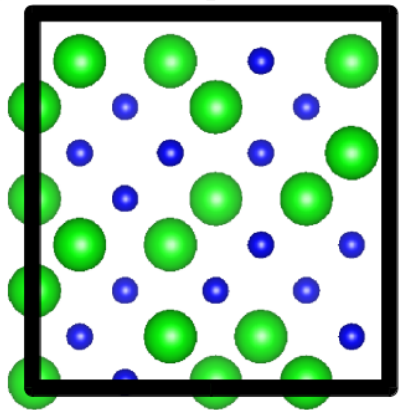
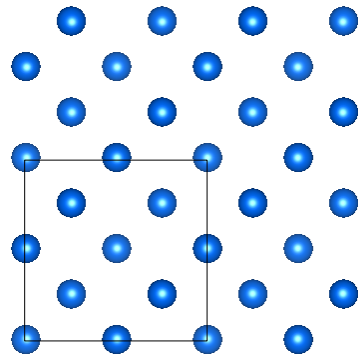
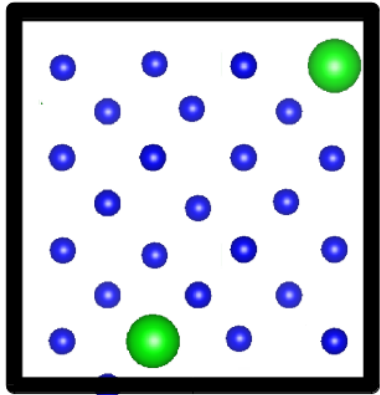
## Matthiessen's Rule

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

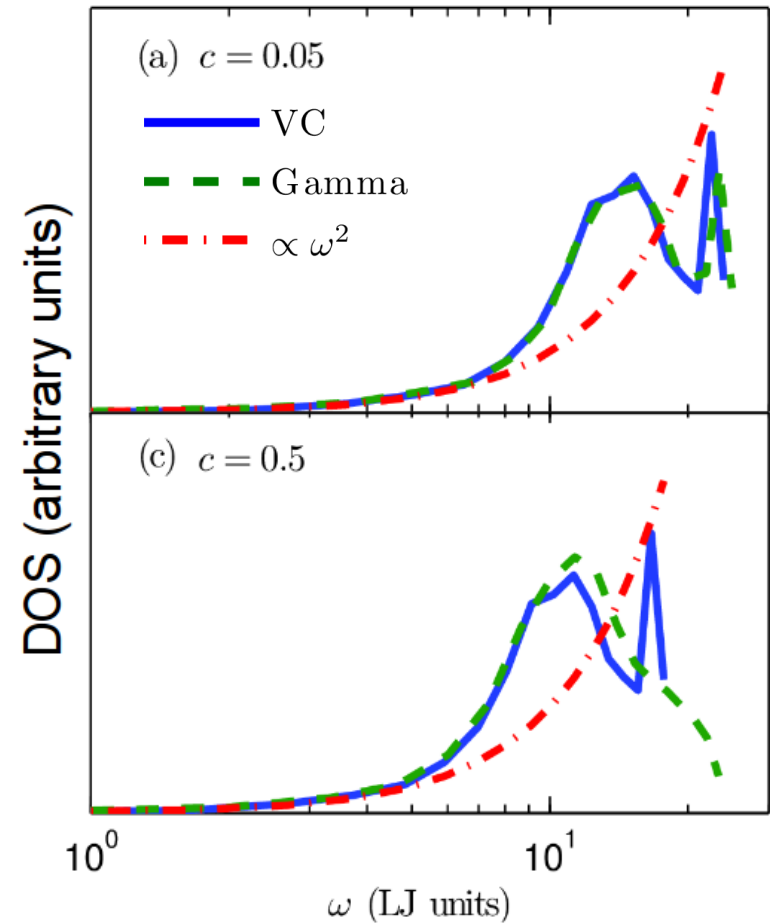


# Explicit disorder: VC vs Gamma

Gamma

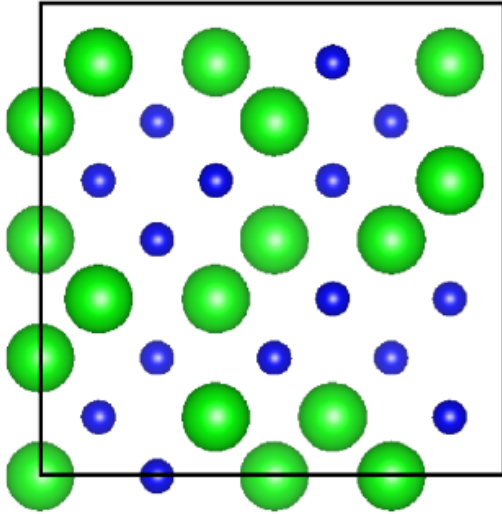


## Lennard-Jones Argon Alloys



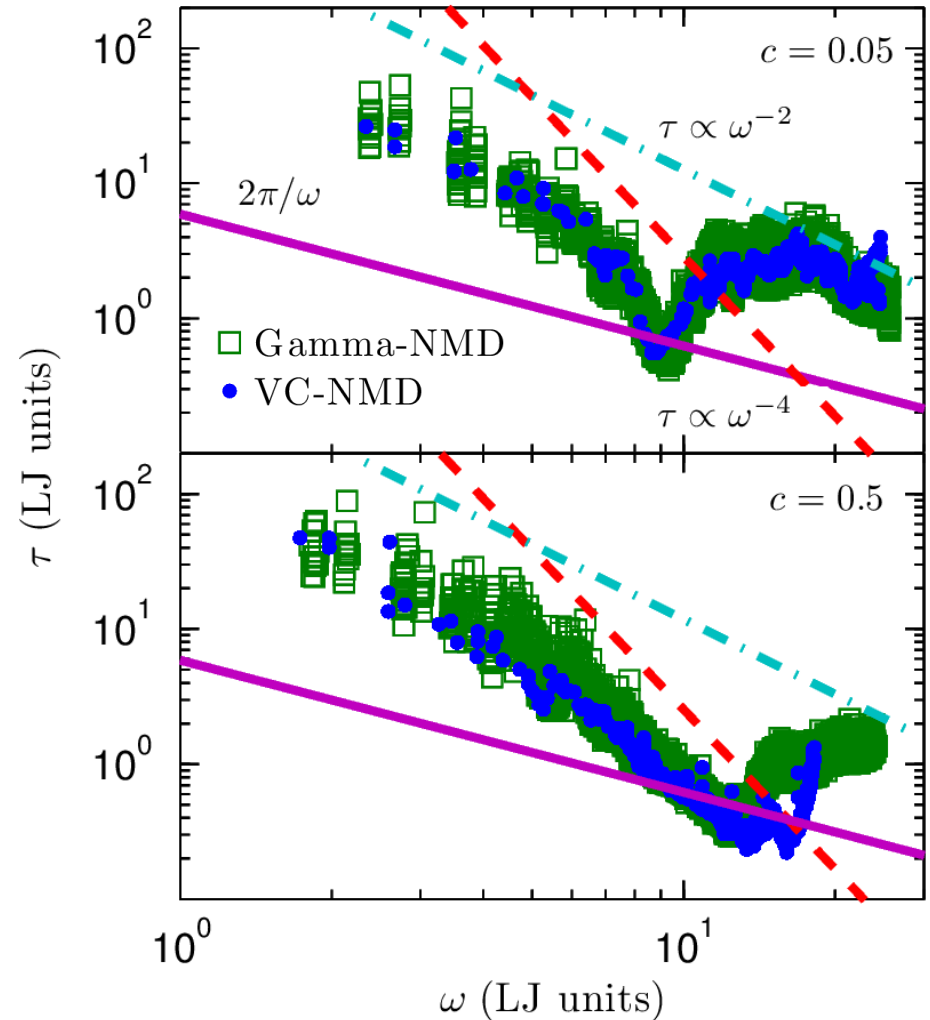
# Normal Mode Decomposition (NMD)

## Molecular Dynamics Gamma



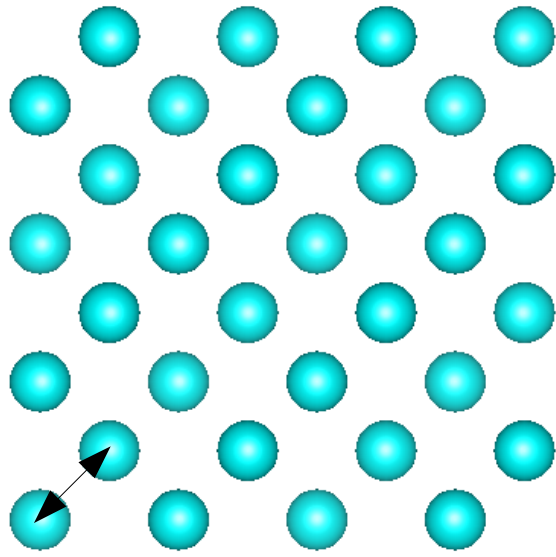
$$\tau(\kappa) = \int_0^{t^*} \frac{\langle E(\kappa; t) E(\kappa; 0) \rangle}{\langle E(\kappa; 0) E(\kappa; 0) \rangle} dt$$

LJ Argon and Alloys,  $T=10$  K



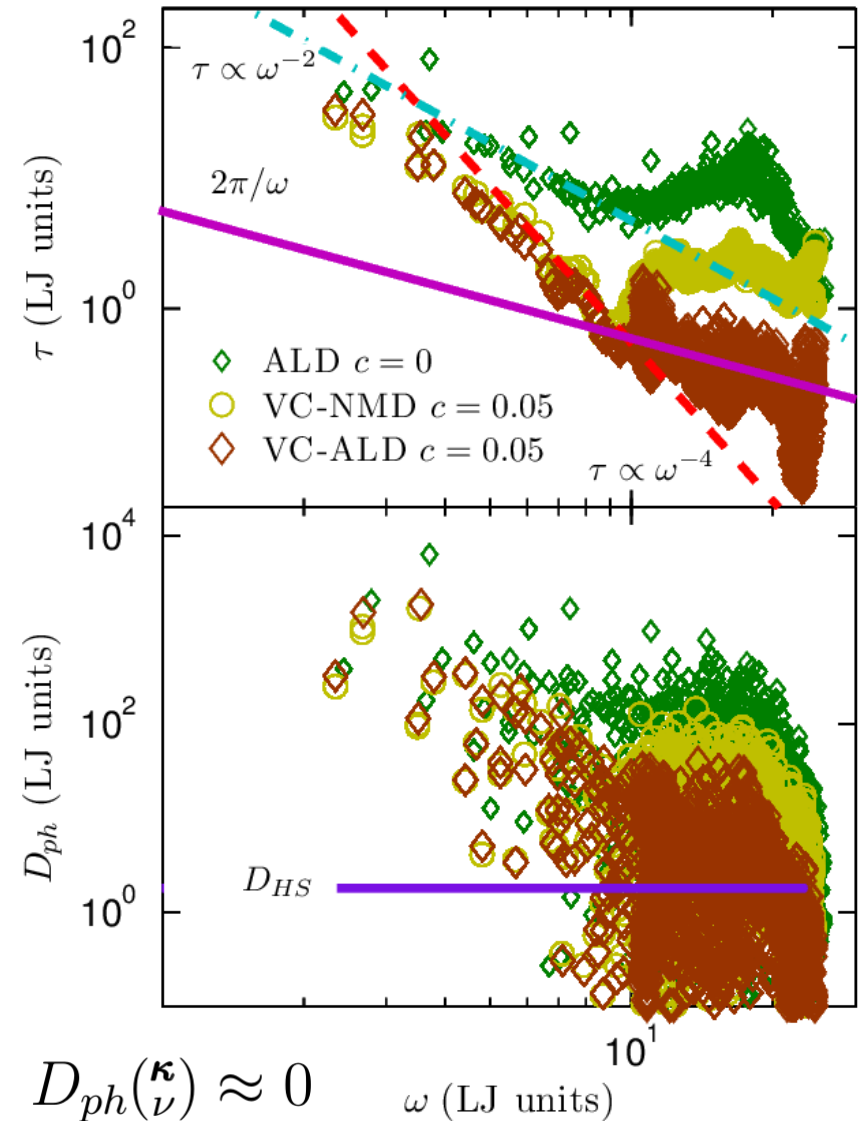
# VC Diffusivities

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



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

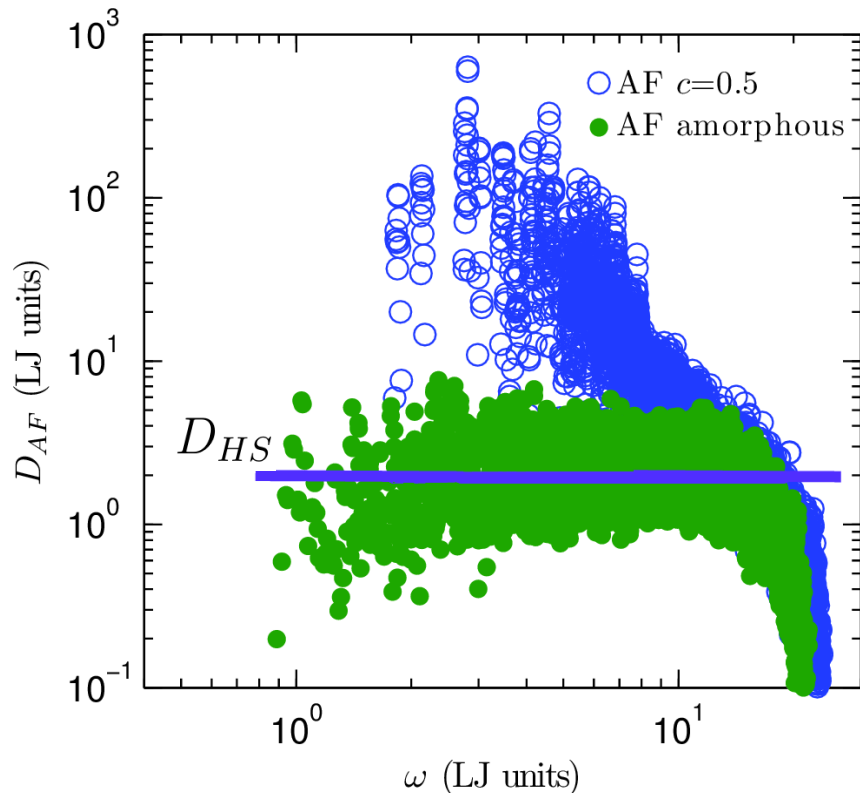
LJ Argon and Alloys,  $T = 10$  K



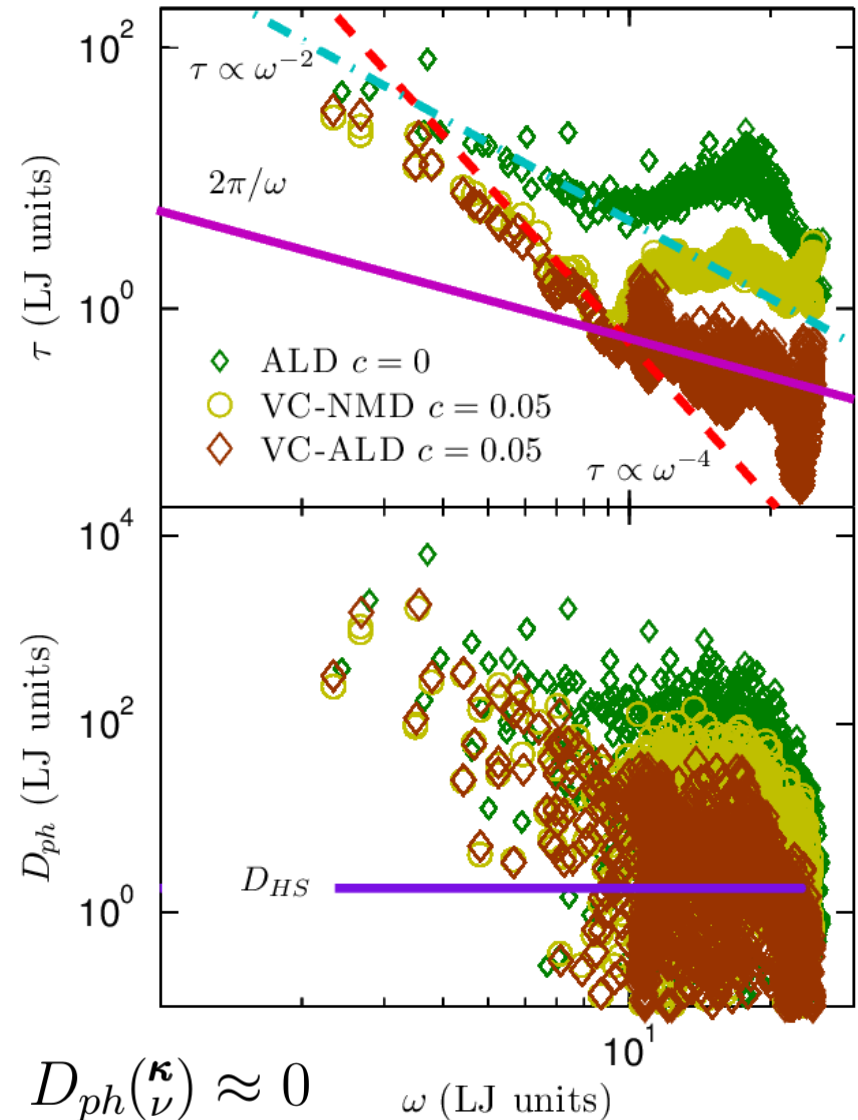
# AF Diffusivities

Allen-Feldman (AF) Theory:

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



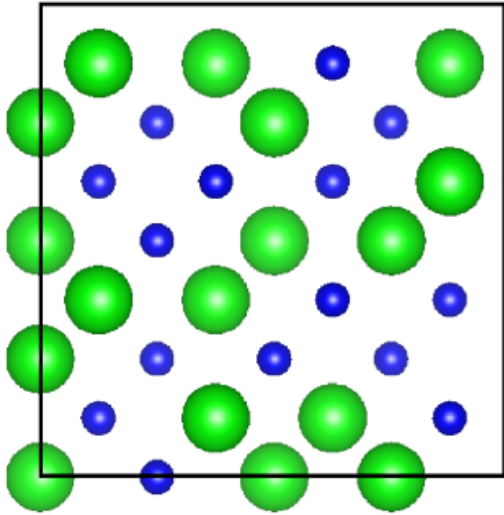
LJ Argon and Alloys,  $T = 10$  K





# Thermal conductivity

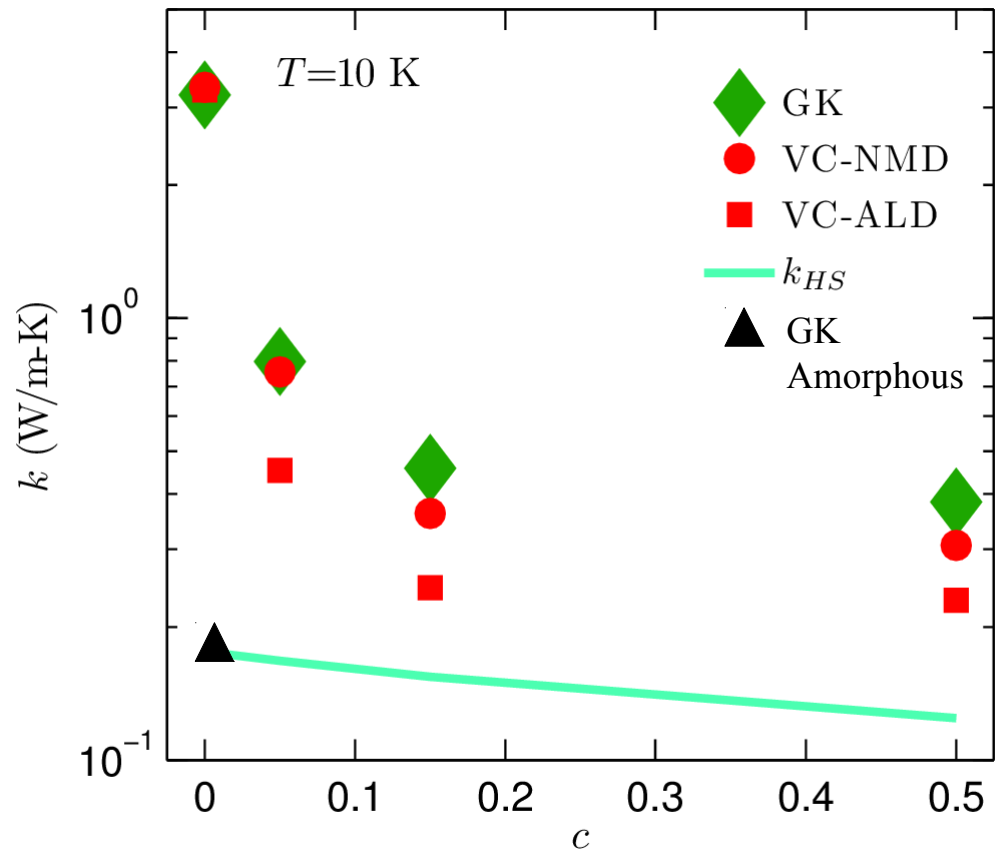
MD-based Green-Kubo (GK)



High-scatter adjustment\*:

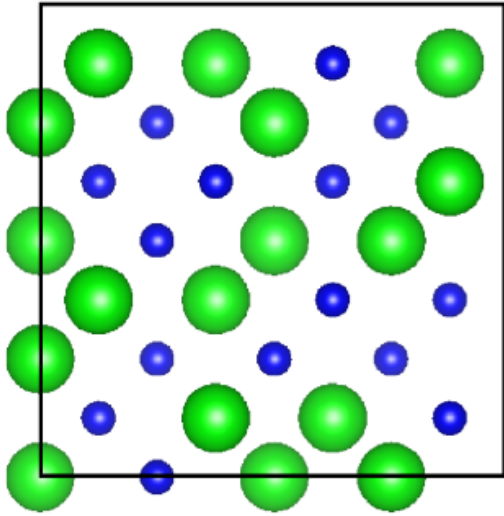
$$D_{ph}(\boldsymbol{\kappa}) < D_{HS}$$

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



# Thermal conductivity

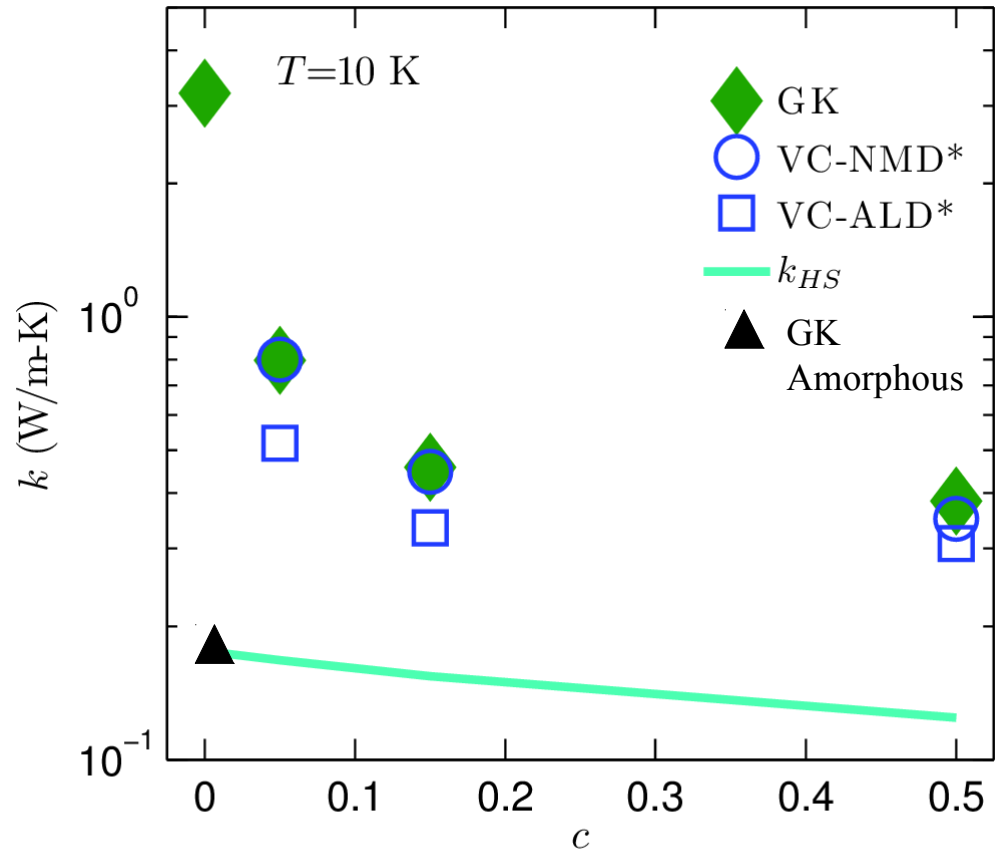
## Gamma MD-based Green-Kubo



High-scatter adjustment\*:

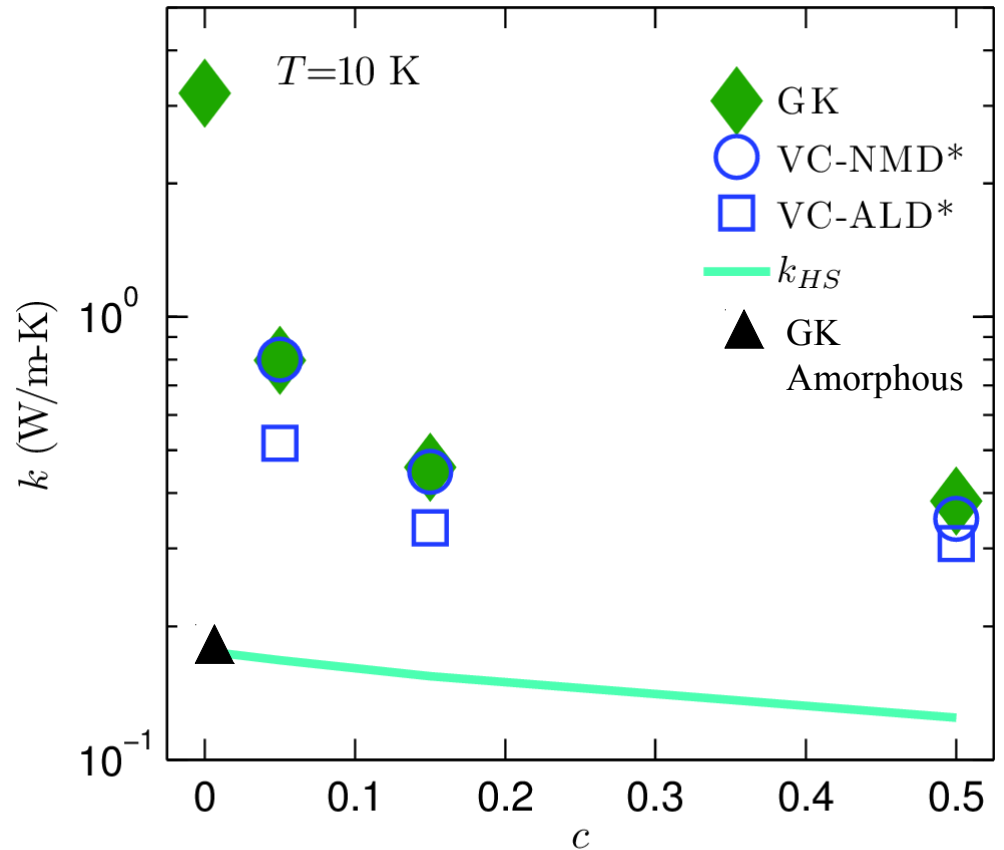
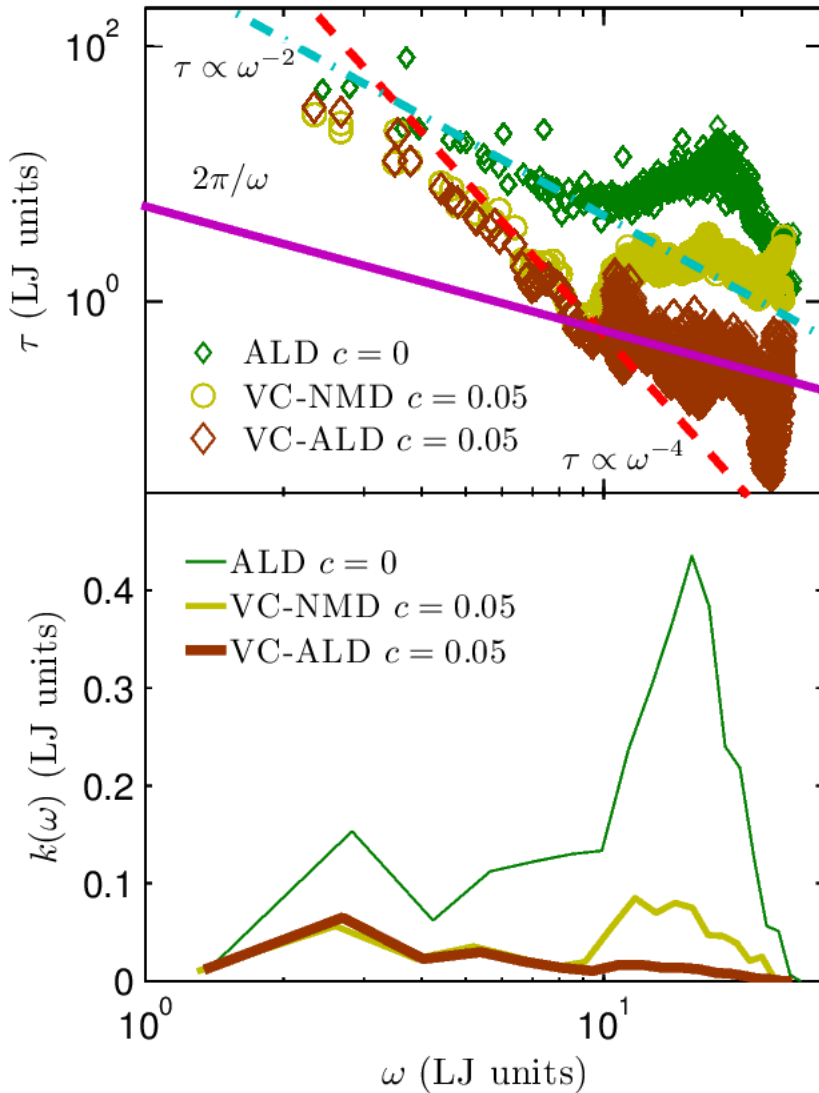
$$D_{ph}(\boldsymbol{\kappa}) < D_{HS}$$

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



# Thermal conductivity spectrum

LJ Argon and Alloys,  $T = 10$  K



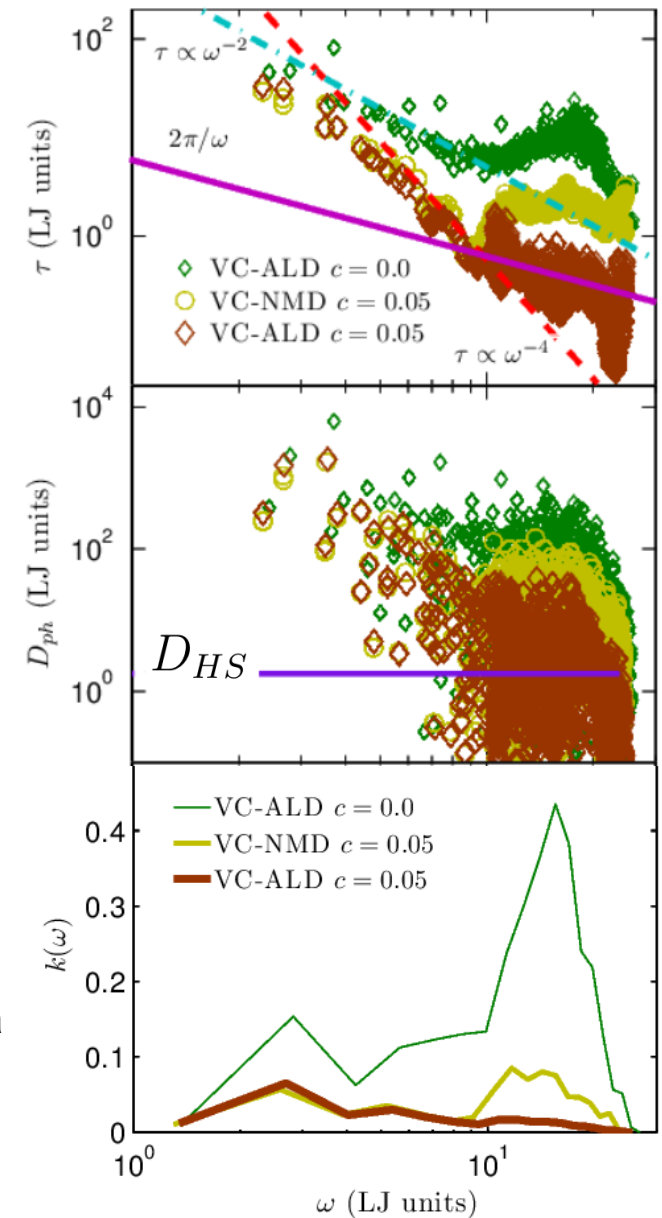
# Summary

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

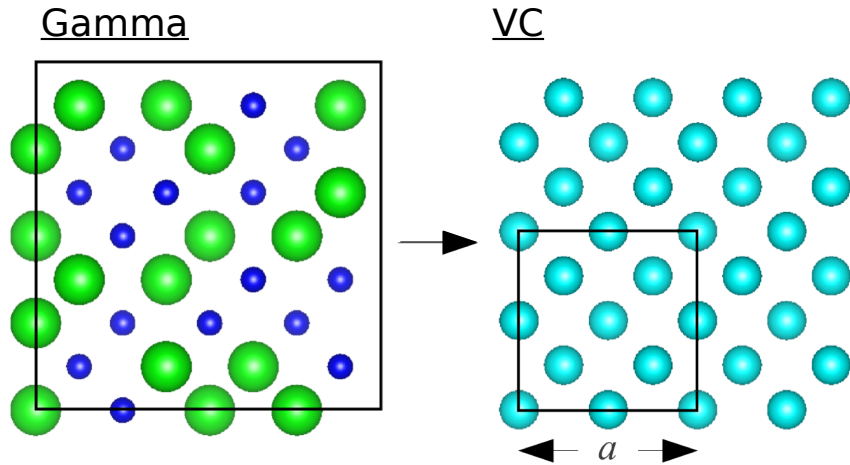
**VC-ALD underpredicts lifetimes** at high-frequency.

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

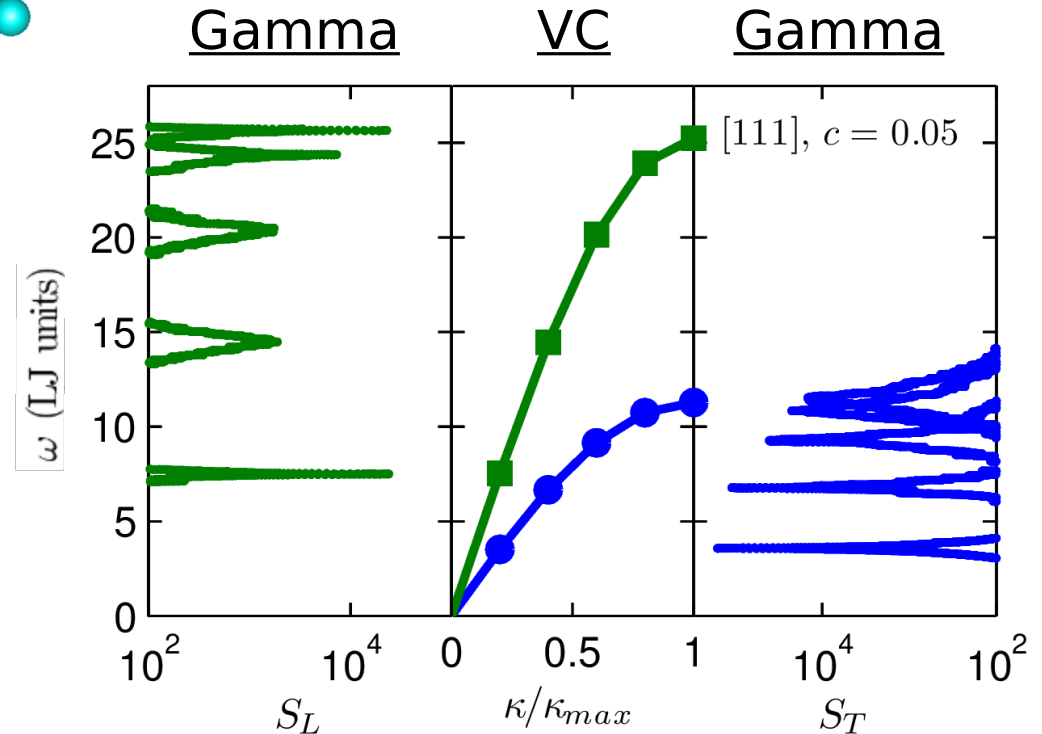
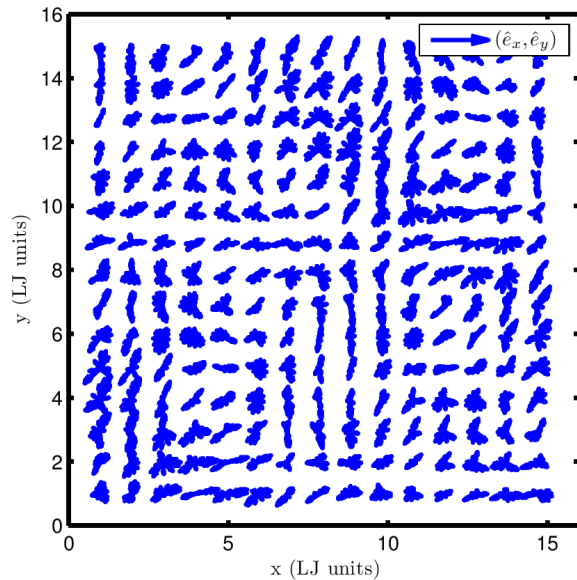
This work was supported by AFOSR award FA95501010098 and by a grant of computer time from the DOD High Performance Computing Modernization Program at the US Army Engineer Research and Development Center. We thank Davide Donadio, Jivtesh Garg, Asad Hasan, Craig Maloney, and Zhiting Tian for helpful discussions.



# Explicit disorder: Structure Factor

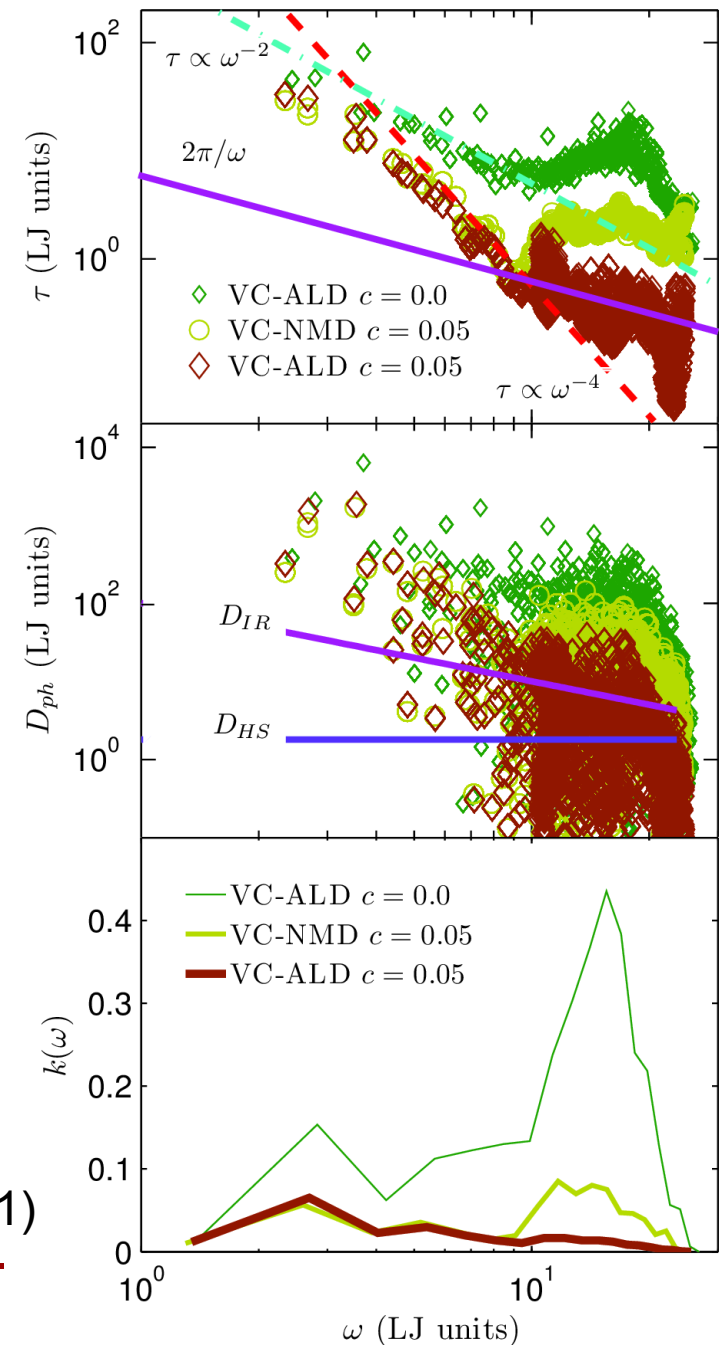
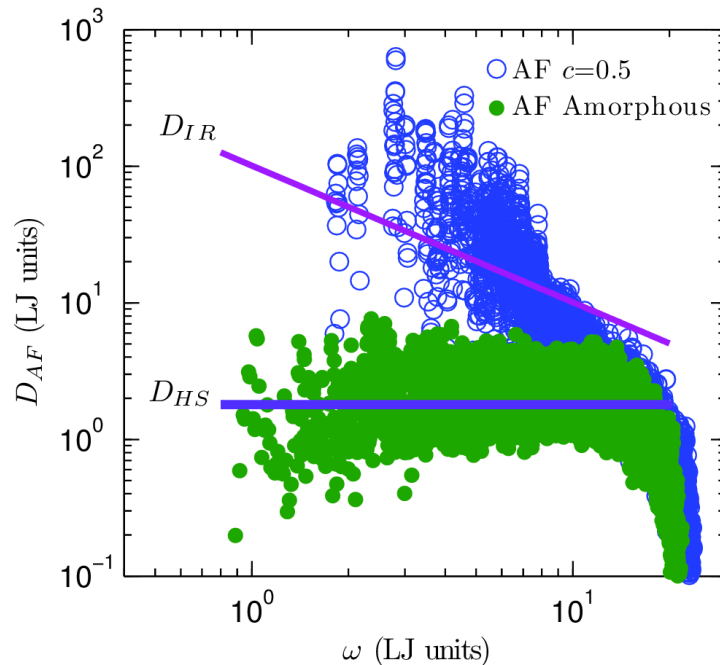


$$D_{ph,n}(\kappa) = v_{g,n}^2(\kappa) \tau(\kappa)$$



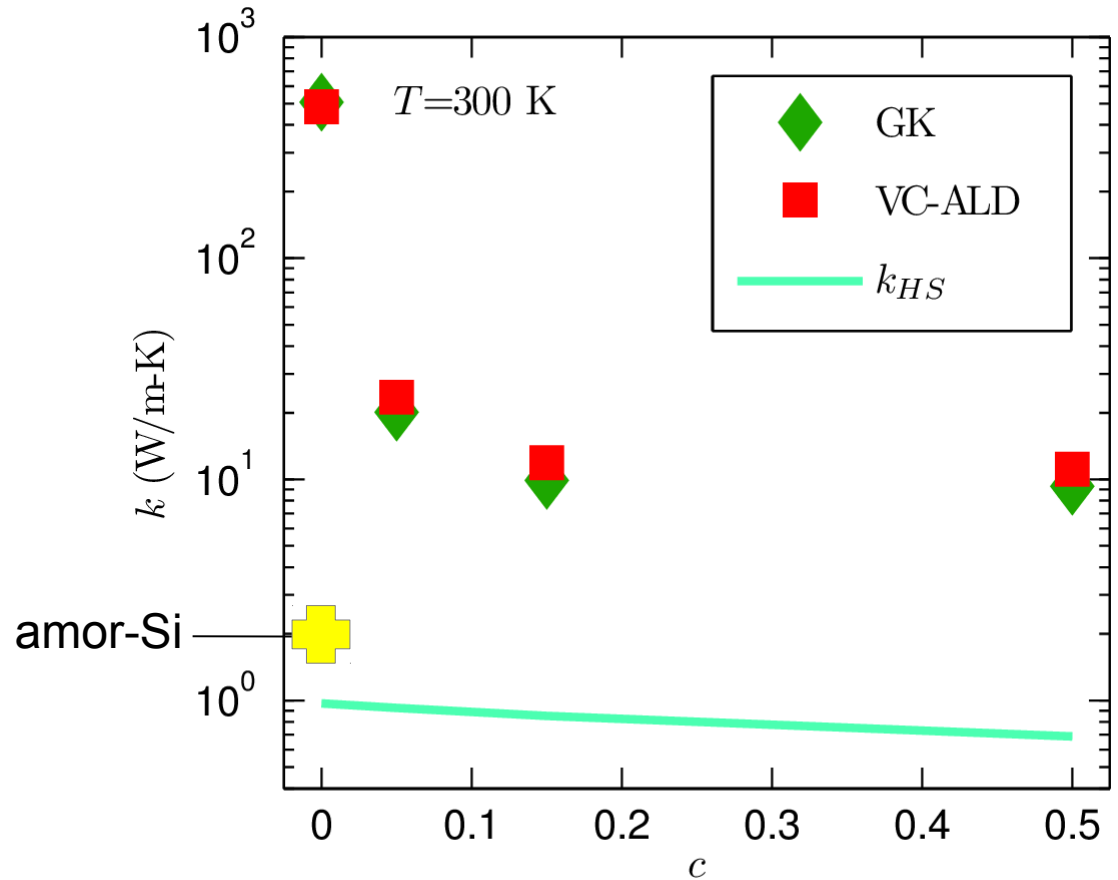
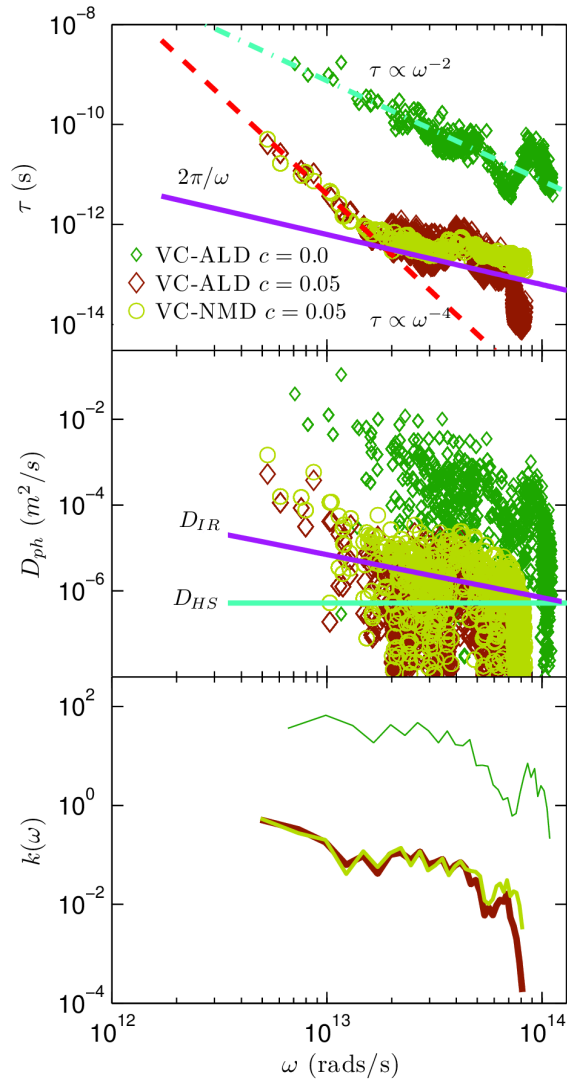
# HS/IR Limit

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



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

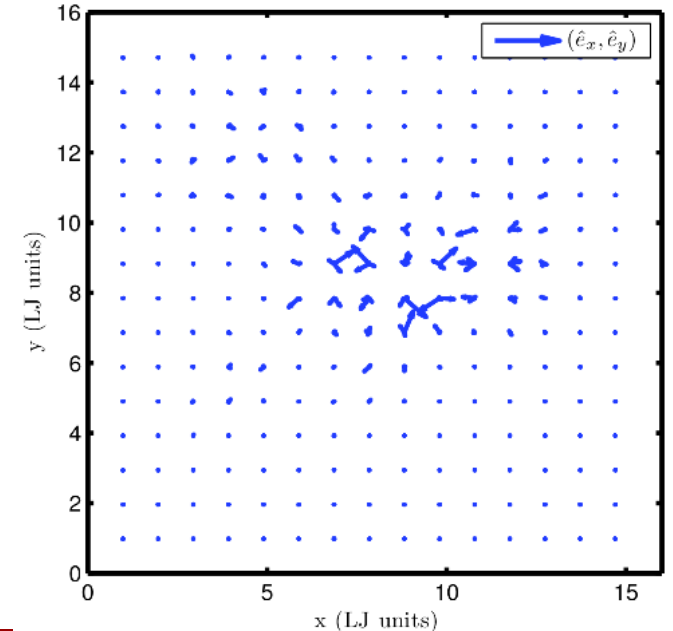
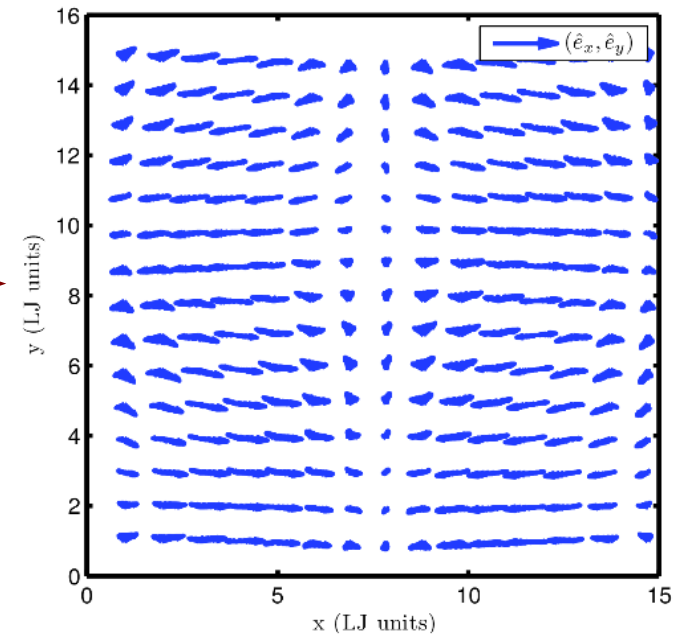
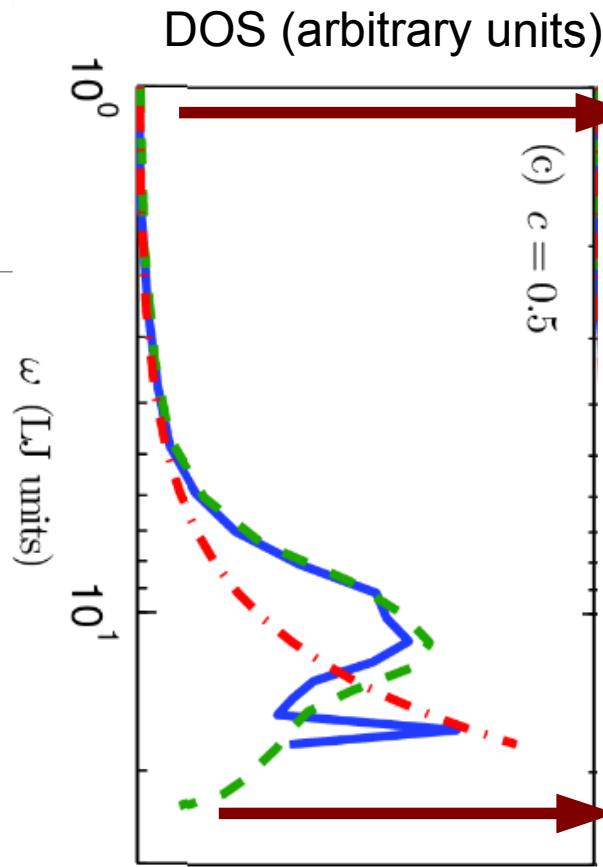
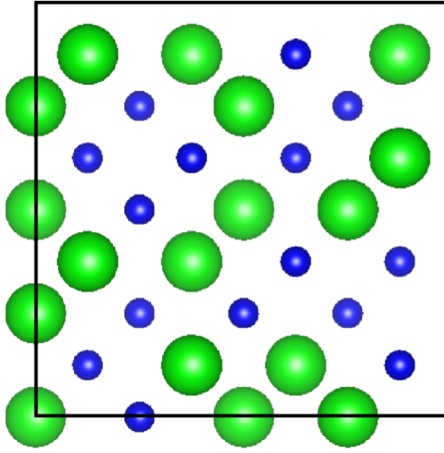
# Thermal conductivity: SW silicon alloy



# Gamma modes

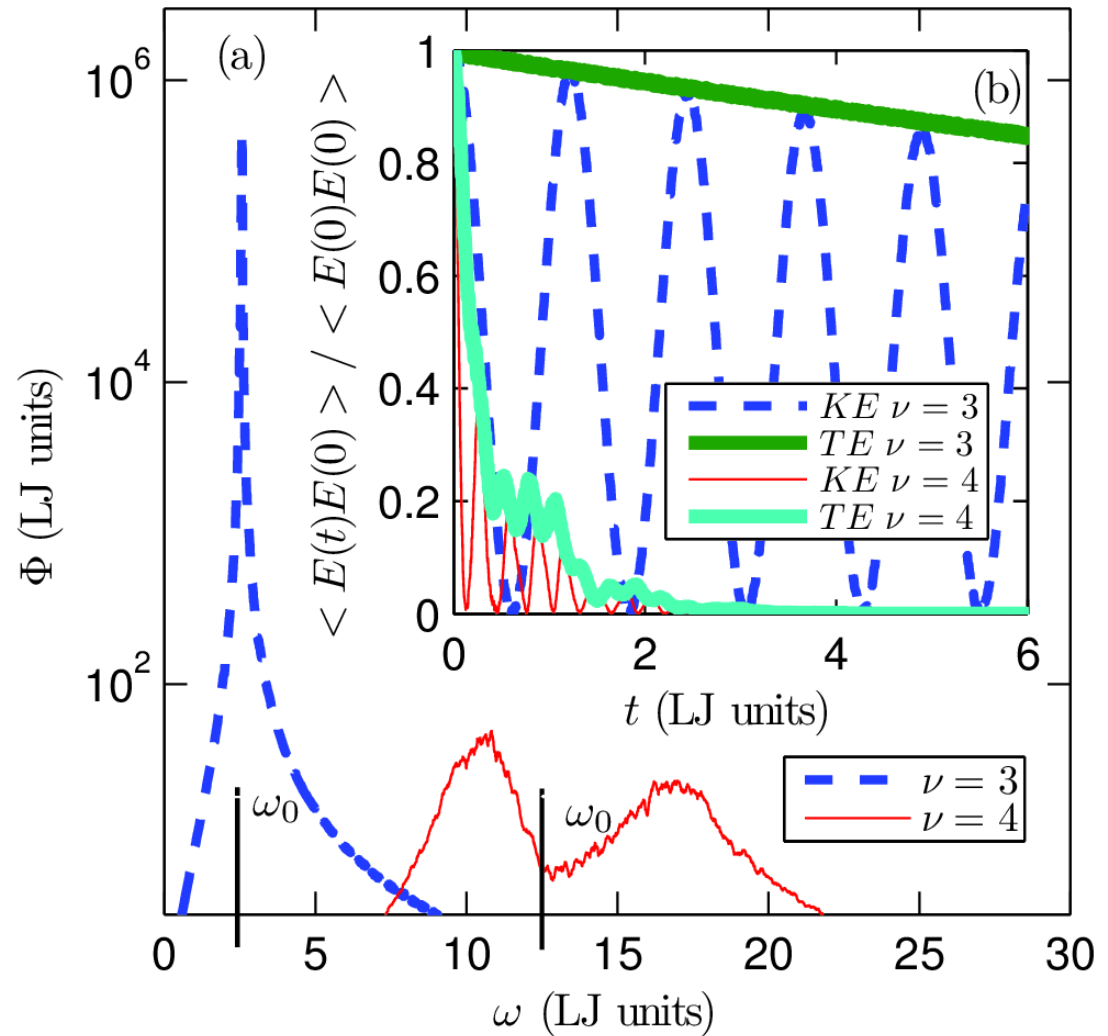
$$e\left(\begin{smallmatrix} \kappa=0 & b \\ \nu & \alpha \end{smallmatrix}\right)$$

(a) disordered supercell

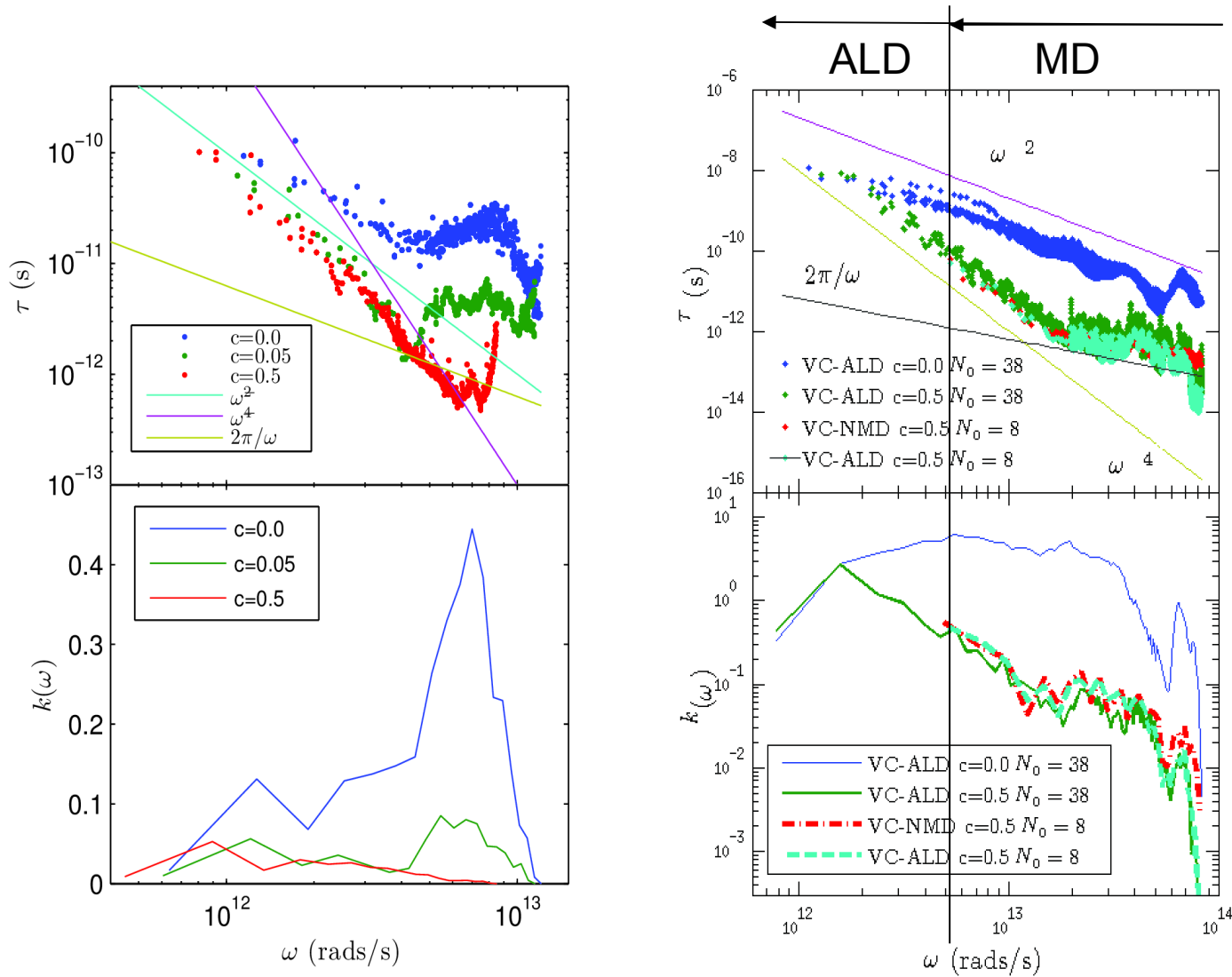




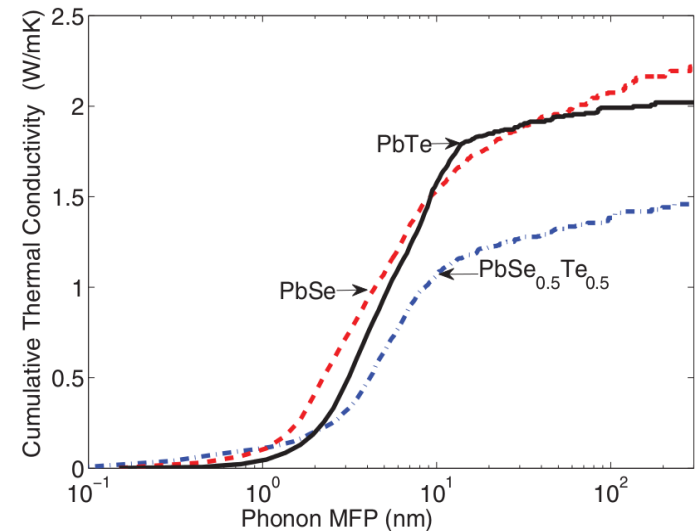
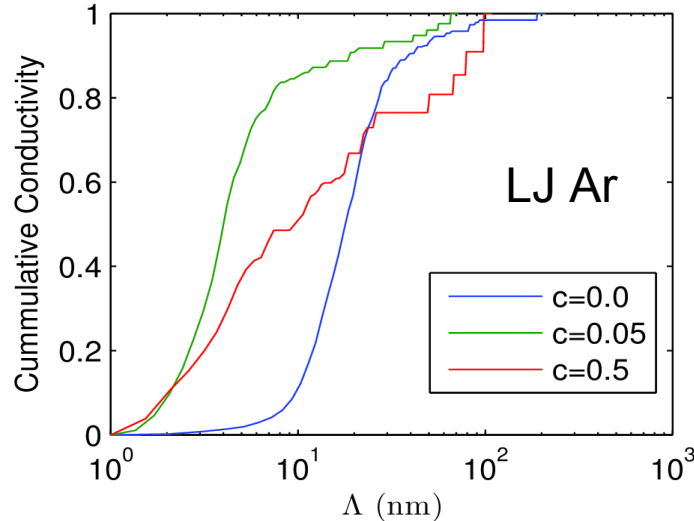
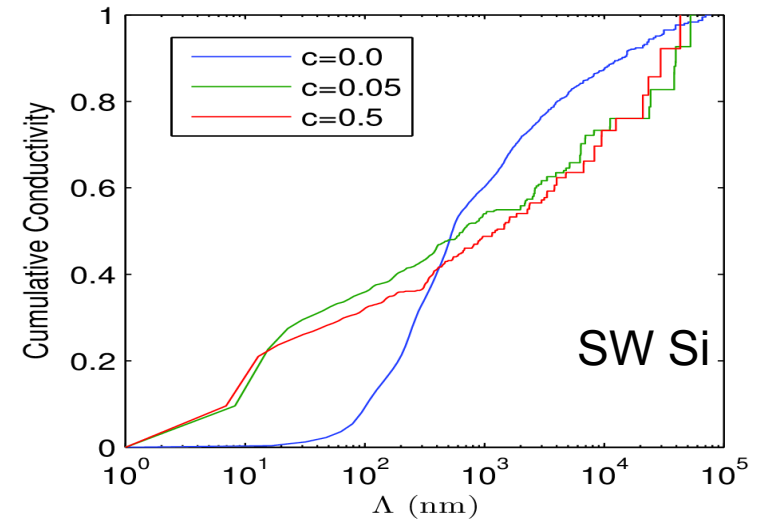
# NMD using VC modes



# Phonon Spectrum: LJ Ar vs SW Si



# Conductivity Accumulation



PHYSICAL REVIEW B 85, 184303 (2012)