Evaluation of the Virtual Crystal Approximation for Predicting Alloy Thermal Conductivity

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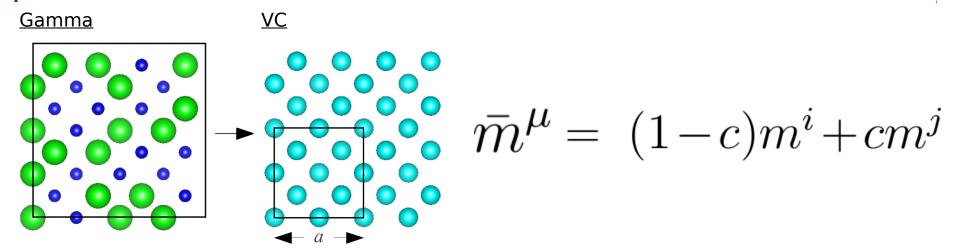
Department of Mechanical Engineering Carnegie Mellon University

http://ntpl.me.cmu.edu/

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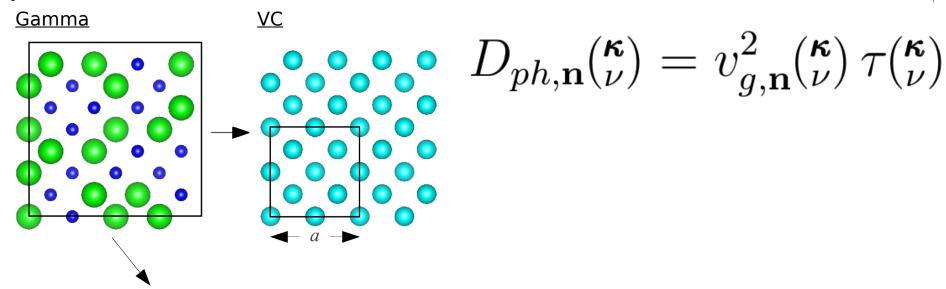
Virtual Crystal Approximation



<u>Lennard-Jones Argon and Alloys</u>

$$k_{ph,\mathbf{n}} = \sum_{\boldsymbol{\kappa}} \sum_{\nu} \frac{\kappa_B}{V} D_{ph,\mathbf{n}} {\kappa \choose \nu}$$
$$D_{ph,\mathbf{n}} {\kappa \choose \nu} = v_{g,\mathbf{n}}^2 {\kappa \choose \nu} \tau {\kappa \choose \nu}$$

Virtual Crystal: Diffusivities

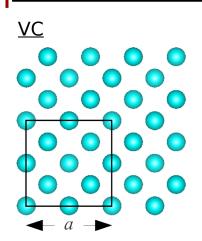


Allen-Feldman (AF) Theory:

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

$$D_{AF,i}(\omega_i) = v_g^2 \tau$$

VC-ALD Diffusivities: Lifetimes



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

Perturbation theory:

<u>Anharmonic Lattice</u> <u>Dynamics (**ALD**)</u>

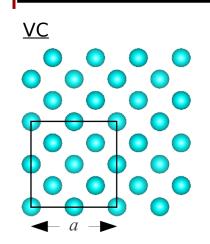


$$\frac{1}{\tau({}^{\kappa}_{\nu})}$$

$$= \frac{1}{\tau_{p-p}\binom{\kappa}{\nu}}$$



VC-ALD Diffusivities: Lifetimes



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

<u>Perturbation theory:</u>

<u>Anharmonic Lattice</u> <u>Dynamics (**ALD**)</u> Phonon-Defect¹



$$\frac{1}{\tau({}^{\kappa}_{
u})}$$

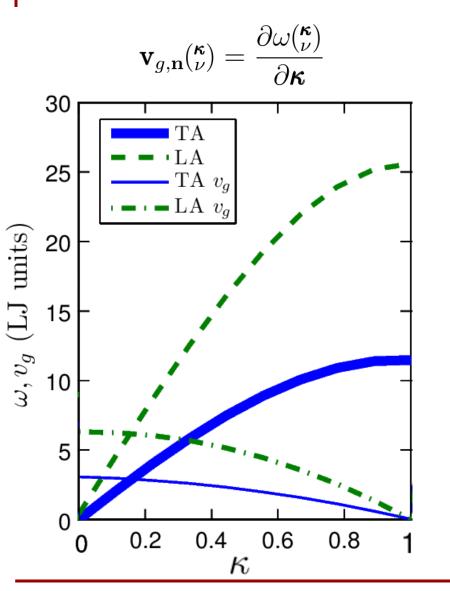
$$= \frac{1}{\tau_{p-p}\binom{\kappa}{\nu}}$$

$$\frac{1}{\tau_{p-d}\binom{\kappa}{\nu}}$$

¹Tamura, PRB 27, 858866 (1983)



VC-ALD Diffusivities: Group Velocity



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

$$D_{ph}(^{\kappa}_{\nu}) \approx 0$$

High-Scatter limit:1

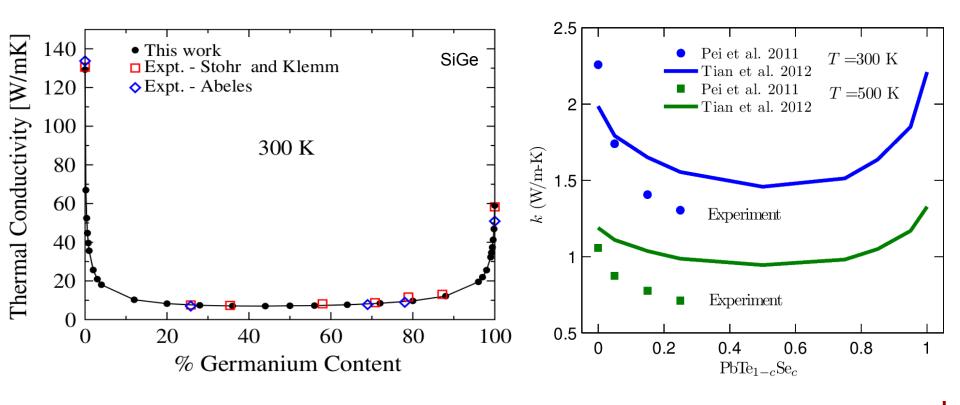
$$D_{HS} = \frac{1}{3}v_s a$$

Mechanical Engineering

VC-ALD: experimental accuracy

Density Functional Theory (DFT)

+ (VC-ALD)



PRL 106, 045901 (2011)

PRB 85, 184303 (2012)

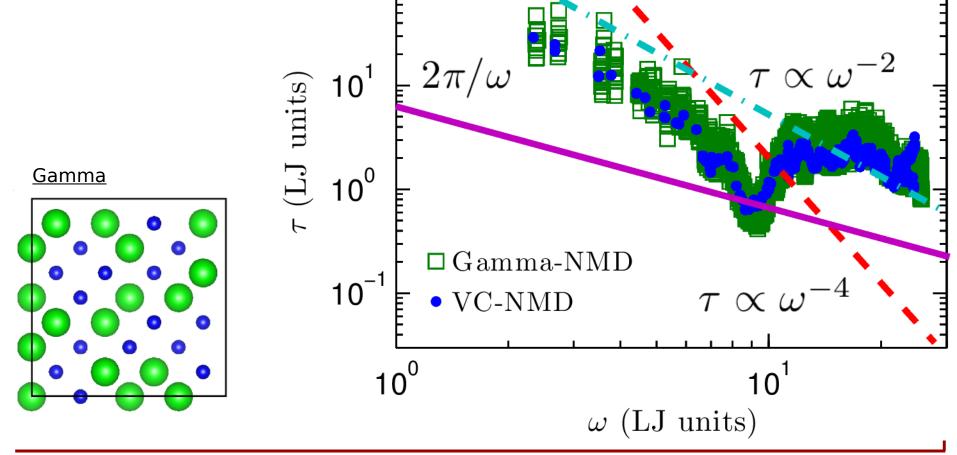


Explicit disorder: NMD

10²

Normal Mode Decomposition (NMD): Molecular Dynamics

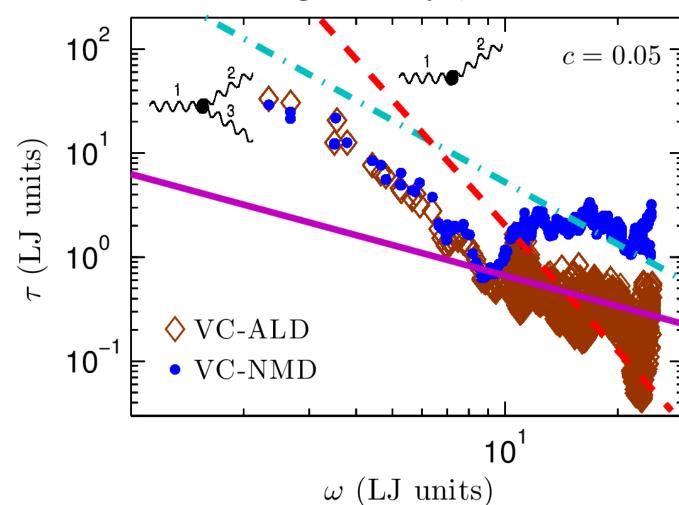
LJ Argon Alloys, T = 10 K

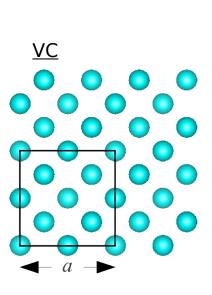


c = 0.05

VC-NMD vs VC-ALD

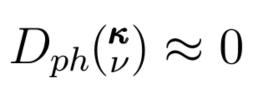




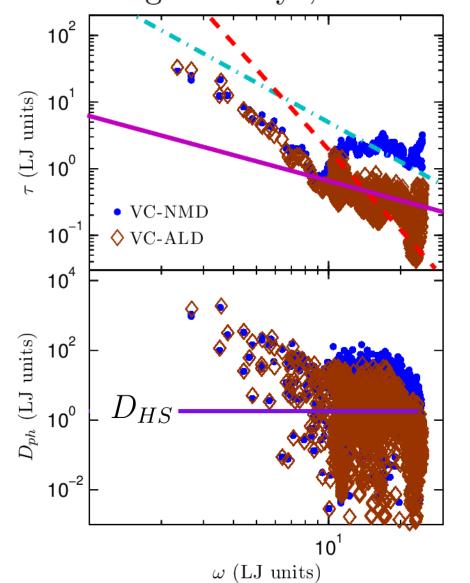


VC Diffusivities

LJ Argon Alloys, T = 10 K



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

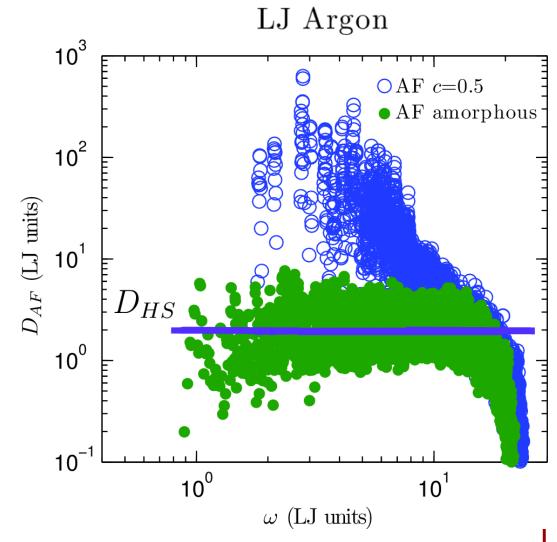




AF Diffusivities

Allen-Feldman (AF) Theory:

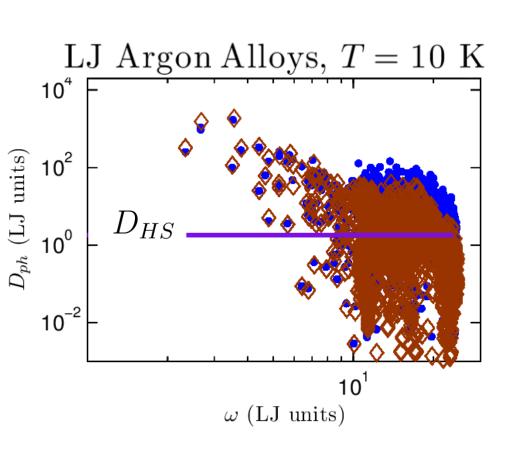
$$k_{AF} = \sum_{diffusions} \frac{k_B}{V} D_{AF,i}(\omega_i)$$

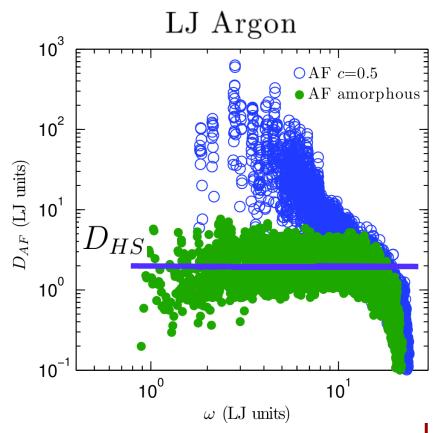


AF and VC Diffusivities

Phonons

Diffusons

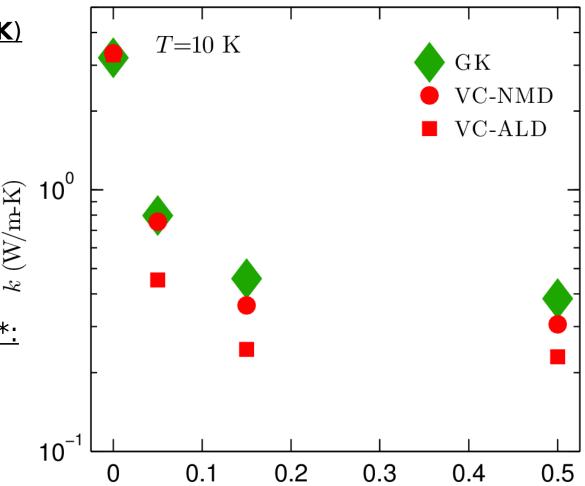




Thermal conductivity

LJ Argon and Alloys, T = 10 K

MD-based Green-Kubo (GK)



<u>High-scatter adjustment*:</u>

$$D_{ph}({}^{\kappa}_{\nu}) < D_{HS}$$

$$D_{ph}({}^{\kappa}_{\nu}) = D_{HS}$$

Thermal conductivity

LJ Argon and Alloys, T = 10 K

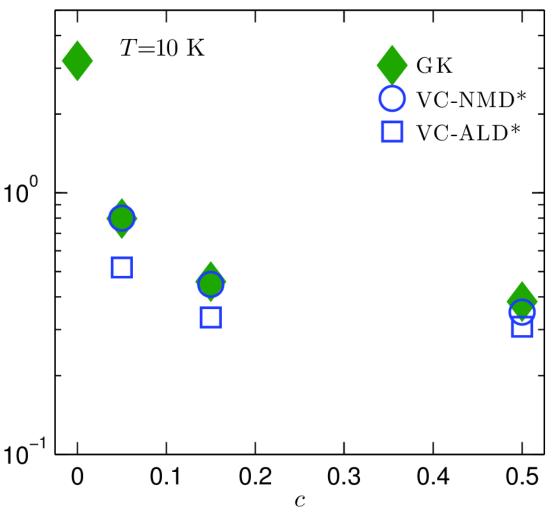




<u>High-scatter adjustment*:</u>

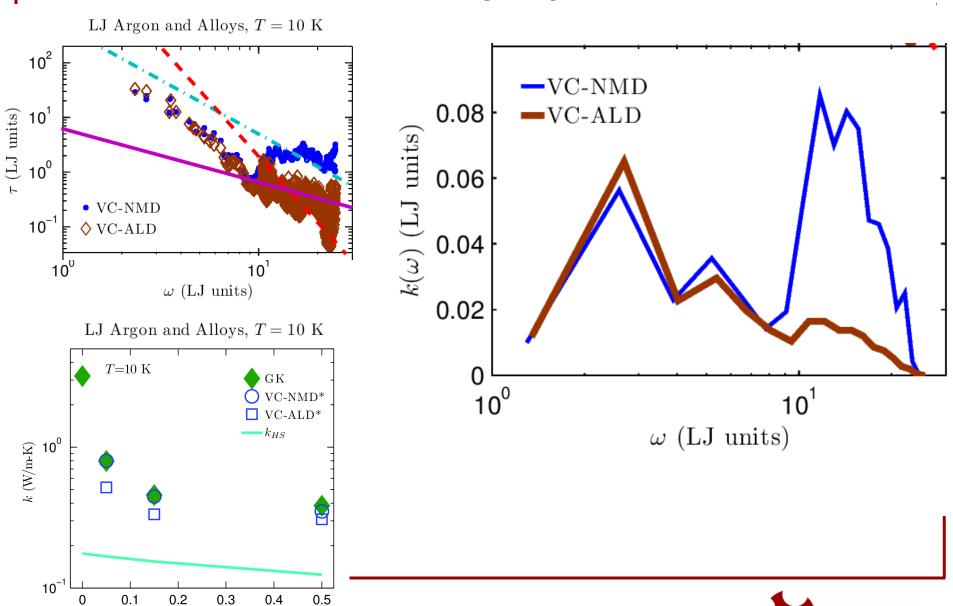
$$D_{ph}({}^{\kappa}_{\nu}) < D_{HS}$$

$$D_{ph}({}^{\kappa}_{\nu}) = D_{HS}$$

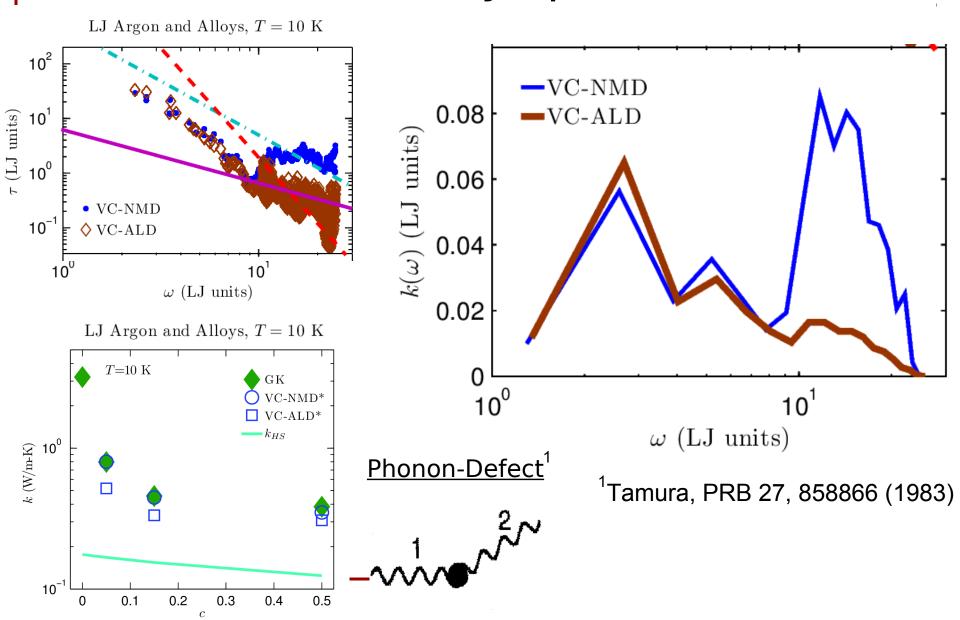


Thermal conductivity spectrum

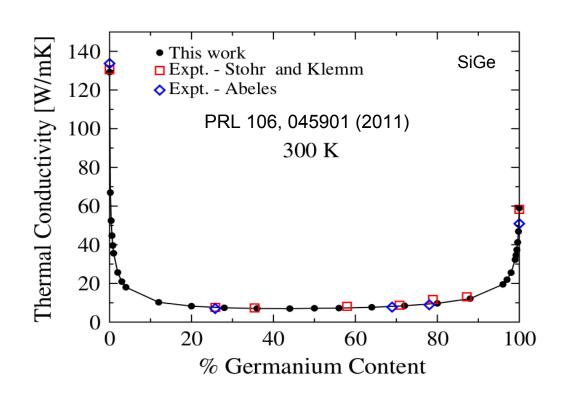
c



Thermal conductivity spectrum

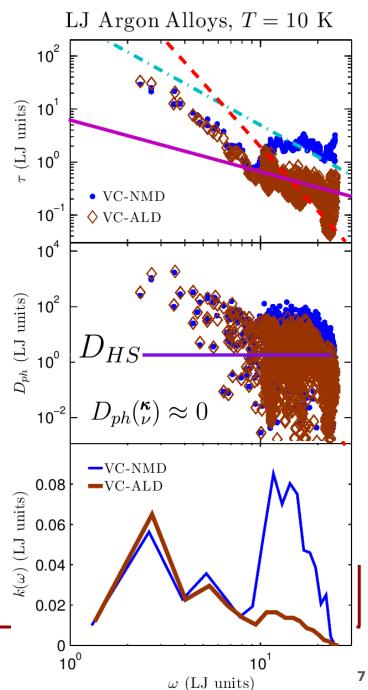


<u>Summary</u>

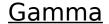


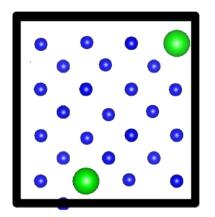
This work was supported by AFOSR award FA95501010098 and by a grant of computer time from the DOD High Performance Computing Modernization Program at the US Army Engineer Research and Development Center.



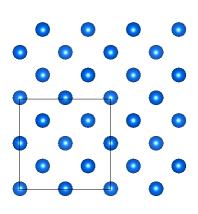


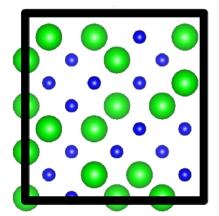
Explicit disorder: VC vs Gamma

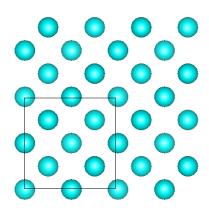




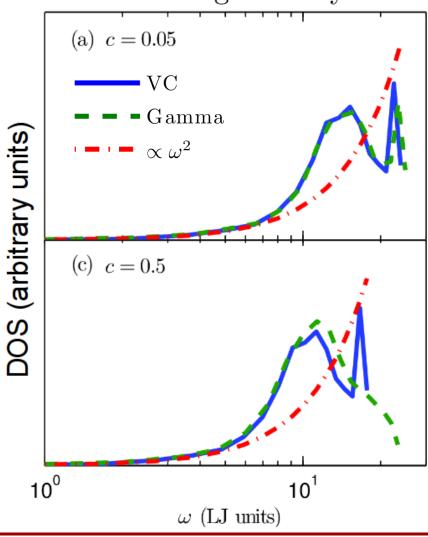




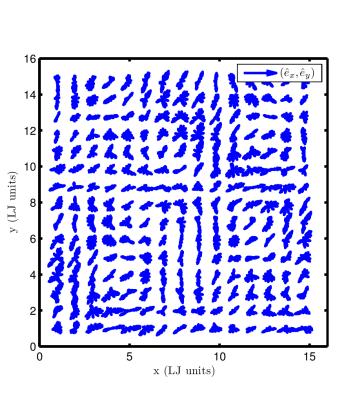


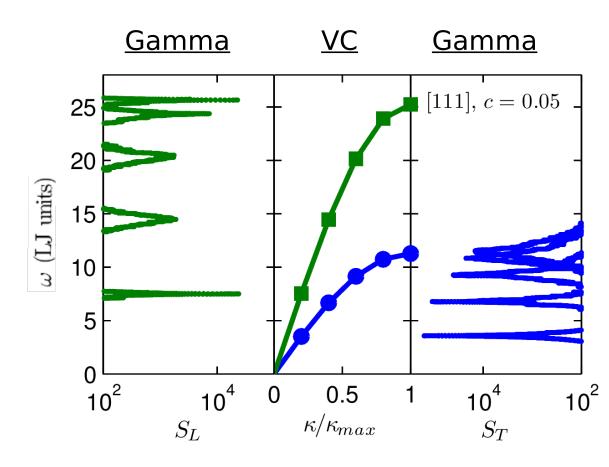


LJ Argon Alloys



Explicit disorder: Structure Factor

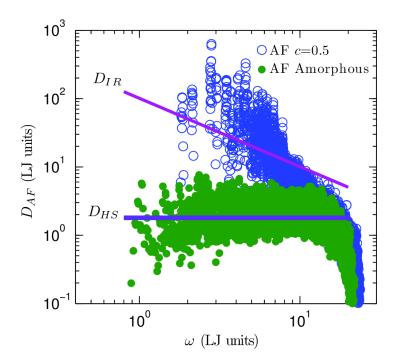




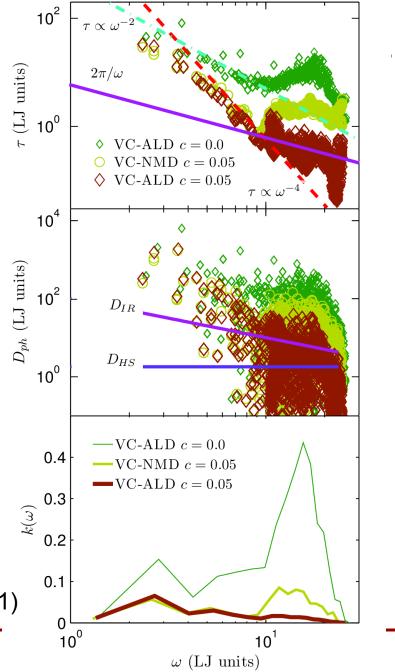


HS/IR Limit

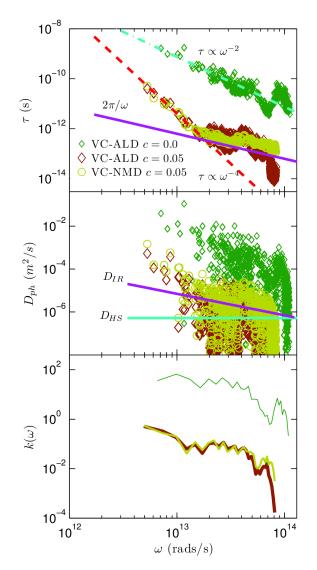
$$D_{IR} = \frac{2\pi}{3} \frac{v_s^2}{\omega}.$$

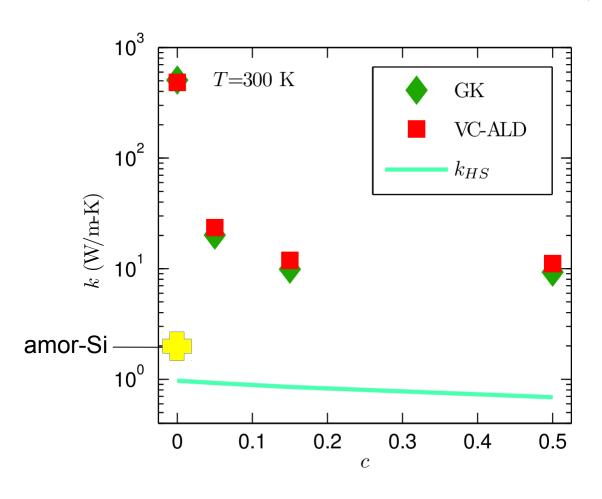


P. Sheng and M. Zhou, Science 253, 539542 (1991)



Thermal conductivity: SW silicon alloy

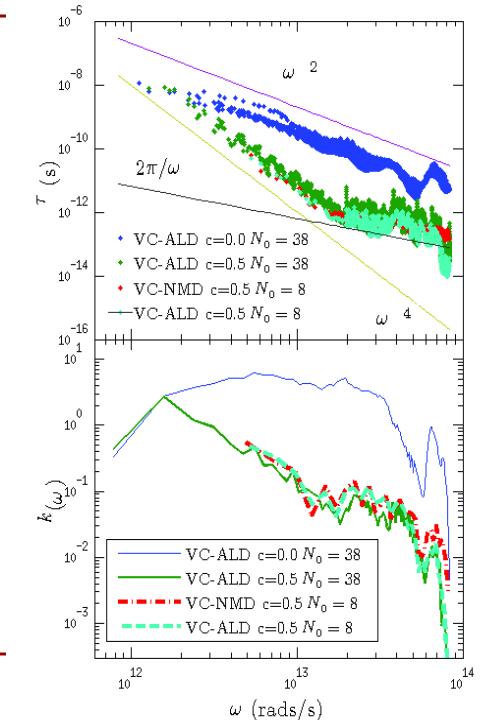


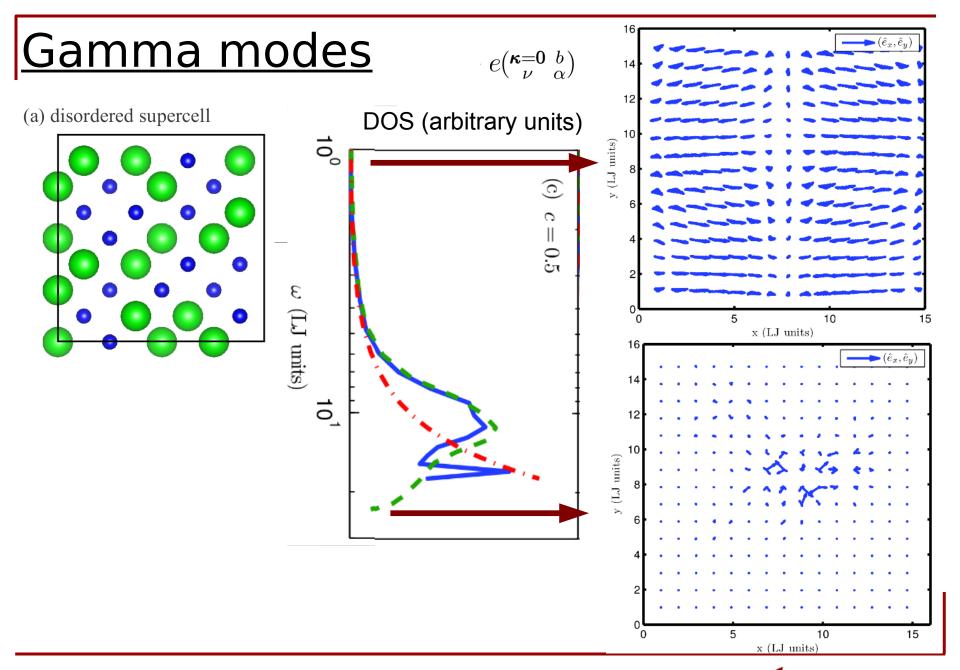




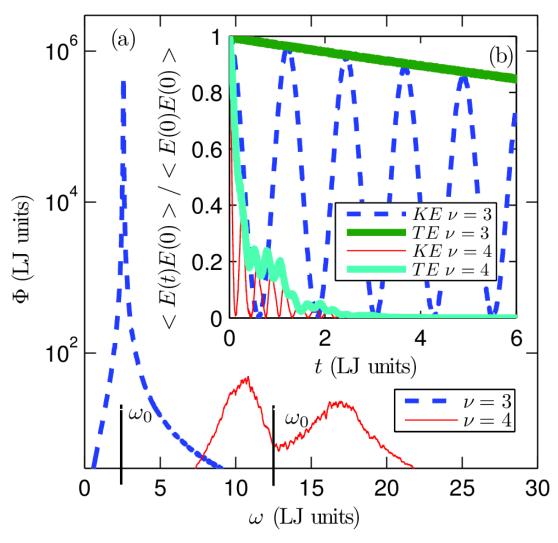


Phonon Spectrum: SW Si

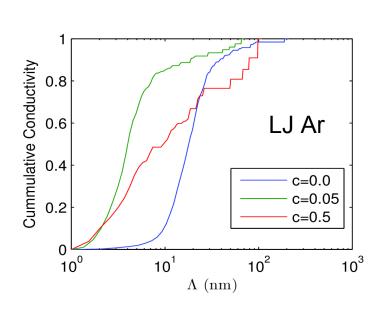


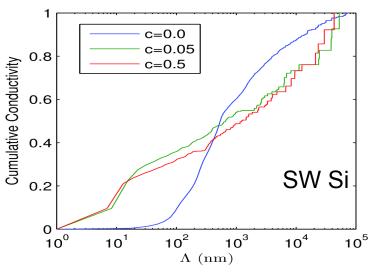


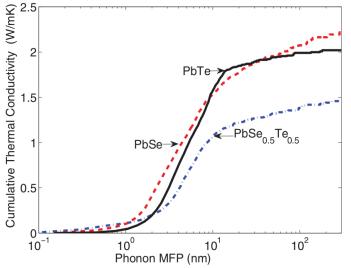
NMD using VC modes



Conductivity Accumulation







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