

University of Illinois, Urbana-Champaign
 Materials Science and Engineering
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Professional Preparation

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| Xavier University | B.S., Physics and Mathematics | 1996 |
| Ohio State University | Ph.D., Physics | 2003 |
| | Advisors: Prof. John W. Wilkins and Dr. Robert C. Albers | |
| Air Force Research Laboratory | Postdoc, Materials and Manufacturing Directorate | 2004-06 |
| | Advisor: Dr. Christopher Woodward | |

Appointments

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| 2021- | Ivan Racheff Professor of Materials Science and Engineering; Materials Research Laboratory, and Computational Science and Engineering, UIUC |
| 2018-2021 | Professor, Materials Science and Engineering, Materials Research Laboratory, and Computational Science and Engineering, UIUC |
| 2018- | Associate Head, Materials Science and Engineering, UIUC |
| 2012-2018 | Associate Professor, Materials Science and Engineering, Materials Research Laboratory, and Computational Science and Engineering, UIUC |
| 9-12/2017 | Senior Research Fellow, Institute for Pure and Applied Mathematics, UCLA |
| 9-12/2012 | Senior Research Fellow, Institute for Pure and Applied Mathematics, UCLA |
| 2006-2012 | Assistant Professor, Materials Science and Engineering, UIUC |
| 2011-2012 | Assistant Professor, Materials Research Laboratory, UIUC |
| 2009-2012 | Affiliated Assistant Professor, Nuclear, Plasma, and Radiological Eng., UIUC |
| 2005-2008 | Adjunct Assistant Professor of Physics, Ohio State University |
| 2004-2006 | National Research Council Postdoctoral Associate, Air Force Research Laboratory |
| 2001-2003 | Graduate Research Assistant, Ohio State University |
| 2000-2001 | Fowler Fellow, Ohio State University |
| 1998-2000 | Graduate Research Assistant, Los Alamos National Laboratory, Group T-11 |
| 1996-1998 | Fowler Fellow, Ohio State University |

Academic Honors

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|---|--------------------------------------|
| Ivan Racheff Professor of Materials Science and Engineering (2021) | <i>Univ. Illinois</i> |
| Collins Award for Innovative Teaching (2020) | <i>Univ. Illinois</i> |
| Brimacombe Medal (2019) | <i>TMS</i> |
| IAS International Visiting Fellow (2018) | <i>Univ. Warwick</i> |
| NCSA Faculty Fellow (2017) | <i>Univ. Illinois</i> |
| Center for Advanced Study Associate (2017) | <i>Univ. Illinois</i> |
| Willett Faculty Scholar Award (2015) | <i>Univ. Illinois</i> |
| AIME Robert Lansing Hardy Award (2014) | <i>TMS-AIME</i> |
| Xerox Faculty Research Award (2011) | <i>Univ. Illinois</i> |
| Early Career Fellow, Honorable Mention (2011) | <i>TMS</i> |
| Faculty Early Career Development Award (CAREER, 2009) | <i>NSF/CMMI</i> |
| 3M Untenured Faculty Research Award (2008, 2009, 2010) | <i>3M</i> |
| Young Leader International Scholar Award (2008) | <i>TMS</i> |
| Young Leader Professional Development Award, Structural Materials Division (2006) | <i>TMS</i> |
| National Research Council Associate (2004-2006) | <i>Air Force Research Laboratory</i> |

Fowler Fellow (1996–1998, 2000–2001)

Professional Activities

Editorial Review Board for *Integrating Materials and Manufacturing Innovation*, 2012–2017

Co-chair, 2011 Physical Metallurgy Gordon conference

Co-chair, 2025 Physical Metallurgy Gordon conference

TMS technical committees: Computational Materials Science and Engineering (vice-chair 2012–2014; chair 2014–2016); Chemistry and Physics of Materials (vice-chair 2015–2018; chair 2018–2020); Magnesium; Integrated Computational Materials Engineering

TMS service committees: Diversity Committee (2018–), Nominating Committee (2012–2016), Young Leaders (2006–2012)

Organizer, 2017 TMS annual meeting “Computational Materials Discovery and Optimization: From 2D to Bulk Materials” symposium

Organizer, 2016 TMS annual meeting “Computational Methods for Spatio-temporal Scale-bridging: from Atomistics to Mesoscale” symposium

Organizer, 2016 TMS annual meeting “Computational Materials Discovery and Optimization: From 2D to Bulk Materials” symposium

Organizer, 2015 TMS annual meeting “Computational Discovery of Novel Materials” symposium

Organizer, 2014 TMS annual meeting “Computational Discovery of Novel Materials” symposium

Organizer, 2014 TMS annual meeting “Computational Thermodynamics and Kinetics” symposium

Organizer, 2014 TMS annual meeting “Multiscale Perspectives on Plasticity in HCP Metals” symposium

Organizer, 2013 TMS annual meeting “Computational Discovery of Novel Materials” symposium

Organizer, 2012 TMS annual meeting “Stochastic Methods in Materials Research” symposium

Organizer, 2012 TMS annual meeting “Minerals, Metals & Materials under Pressure” symposium

Organizer, 2010 TMS annual meeting “Stochastic Methods in Materials Research” symposium

Organizer, 2010 TMS annual meeting “Computational Thermodynamics and Kinetics” symposium

Organizer, 2008 TMS annual meeting “Minerals, Metals & Materials under Pressure” symposium

Organizer, 2006 TMS annual meeting “Point Defects in Materials” symposium

Organizer, 2008 “Recent Developments in Electronic Structure Methods” workshop, UIUC

Organizer, 2006 “Recent Developments in Electronic Structure Methods” workshop, OSU

NSF TeraGrid Allocations Committee Reviewer (2008–2011)

Proposal reviews: DOE: BES, NSF: MPS, NSF: ENG

Manuscript reviews: Phys. Rev. Lett., Phys. Rev. B, J. Phys. Chem., J. Phys: CM, Acta Mat., Phil. Mag., J. Math. Phys., Comp. Mat. Sci., Model. Sim. Mater. Sci. Eng.

Organizer, 2007–2014 ASM Materials Camp for high school teachers, UIUC

Organizer, 2011–2014 Girls Learning About Materials (GLAM) summer camp for high school seniors, UIUC

2014–2017 Education Innovation Fellow, College of Engineering, UIUC

Current Research Grants

4. *NRT-HDR: Data and Informatics Graduate Intern-traineeship: Materials at the Atomic Scale (DIGI-MAT)*. D. R. Trinkle, NSF/NRT.
3. *Development of High Energy Density Thermomagnetic Processing Technology for Intensification of Industrial Heat-treatment and Increased Material Performance*. M. Manuel, DOE/EERE.
2. *Sandia LDRD: Machine learning the thermodynamics and kinetics of vacancies for acceleration*

edmaterials discovery. D. R. Trinkle, Sandia.

1. *Learning the Universal Free Energy Function*. D. R. Trinkle, NSF/HDR.

Past Research Grants

20. *Machine Learning methods for multi-disciplinary multi-scales problems*. O. Pauluis, NSF/HDR (01/01/2020–12/31/2023).
19. *Sandia LDRD: Opening the “Black Box”: An Experimentally-Validated Explainable Machine Learning Framework for Materials Design and Discovery*. D. R. Trinkle, Sandia (10/01/2019–09/30/2022).
18. *USAMP Low-Cost Mg Sheet Component Development and Demonstration Project*. USAMP; DOE/EERE (10/01/2016–06/30/2020).
17. *PIRE: Integrated Computational Materials Engineering for Active Materials and Interfaces in Chemical Fuel Production*. N. R. Aluru; NSF (10/01/2015–10/01/2020).
16. *DMREF/GOALI/Collaborative Research: Computational Design, Rapid Processing and Characterization of Multiple Classes of Materials to Accelerate Materials Innovation*. K. S. Ravi Chandran (Univ. Utah), D. R. Trinkle, and A. P. Sanders (Ortho Development); NSF/CMMI (09/01/2014–08/31/2018).
15. *Surface and Interface Free Energies of Epitaxial Nanocrystals*. J. Zuo and D. R. Trinkle; NSF/DMR (09/01/2014–08/31/2017).
14. *Materials for Extreme Irradiation Environments*. Illinois MRL Cluster; DOE/BES (2012–2018).
13. *GOALI: Experimentally-Validated Computational Approach to Developing and Predicting Kinetics in Anisotropic Systems*. D. R. Trinkle, M. Manuel (Univ. Florida), C. Jiang (ThermoCalc), and K. Wu (ThermoCalc); NSF/CDSE (09/01/2014–08/31/2017).
12. *High-throughput discovery of thermally-functional materials*. D. Cahill, D. Shoemaker, D. R. Trinkle; Toyota (3/1/16–2/28/17).
11. *Integrated Computational Approach to Development of Lightweight 3GAHSS Vehicle Sub-Assembly*. Multi-institutional grant, USAMP/ASP; DOE/EERE (10/01/12–09/30/16).
10. *Enhanced Diffusivity Along Dislocations from Quantum Tunneling to Classical Transport in the PD-H System*. B. Heuser and D. R. Trinkle; NSF/DMR (08/15/12–07/31/15).
9. *First Principles Calculation of Oxygen and Oxygen-Solute Effects on Deformation in Titanium*. D. R. Trinkle; ONR (07/01/12–06/30/15).
8. *Engineered Zircaloy Cladding and Fuel Pellet Modifications for Improved Accident Tolerance of LWR Nuclear Fuel*. Multi-institutional grant; DOE/NEUP (01/01/12–12/31/15).
7. *The Role of Point Defects in the Structure and Properties of Perovskites for Functional Applications*. R. Uvic (Boise State), W. Kriven, D. R. Trinkle; NSF (10/01/2010–09/30/2014).
6. *Surface and Interface Free Energies of Epitaxial Nanocrystals*. J. Zuo and D. R. Trinkle; NSF/DMR (07/01/10–06/30/14).
5. *CAREER: First-Principles Modeling of Titanium-Oxygen-Solute Interaction: Materials Design for Improved Energy Efficiency*. D. R. Trinkle; NSF/CMMI (01/15/09–01/14/14).
4. *Hydrogen-dislocation interactions at low temperature in deformed Pd: Spatial and vibrational characterization using neutron scattering and advanced computational techniques*. B. Heuser and D. R. Trinkle; NSF/DMR (06/15/08–08/14/12).
3. *GOALI: Modeling solute effects in magnesium alloys: First-principles to predictive finite-element*. D. R. Trinkle and L. G. Hector, Jr. (General Motors); NSF/CMMI (08/01/08–07/31/12).
2. *Designing Oxygen-tolerant Titanium Alloys with First-principles*. D. R. Trinkle; Boeing (06/15/09–06/15/11).

1. *Heteroepitaxial metal nanostructure diffusion through collective slip*. J. Weaver and D. R. Trinkle; NSF/DMR (06/15/07–06/14/10).

Refereed Publications

87. *Contributions to diffusion in complex materials quantified with machine learning*. S. Chattopadhyay and D. R. Trinkle. *Phys. Rev. Lett.* **132**, 186301 (2024), doi://10.1103/PhysRevLett.132.186301, (Editor's Suggestion)
86. *Spline-based neural network interatomic potentials: Blending classical and machine learning models*. J. A. Vita and D. R. Trinkle. *Comp. Mater. Sci.* **232**, 112655 (2024), doi://10.1016/j.commatsci.2023.112655
85. *Explainable machine learning for hydrogen diffusion in metals and random binary alloys*. G. M. Lu, M. Witman, S. Agarwal, V. Stavila, and D. R. Trinkle. *Phys. Rev. Mater.* **7**, 105402 (2023), doi://10.1103/PhysRevMaterials.7.105402
84. *Numerical calculation of interstitial dumbbell-mediated transport coefficients in dilute crystalline systems with non-truncated correlations*. S. Chattopadhyay and D. R. Trinkle. *Philos. Mag.* **102**(24), 2459–2505 (2022), doi://10.1080/14786435.2022.2120646
83. *First-principles core energies of isolated basal and prism screw dislocations in magnesium*. Y. Dan and D. R. Trinkle. *Mater. Res. Lett.* **10**, 360–368 (2022), doi://10.1080/21663831.2022.2051763
82. *Solutes that reduce yield strength anisotropies in magnesium from first principles*. M. R. Fellinger, L. G. Hector, Jr., and D. R. Trinkle. *Phys. Rev. Mater.* **6**, 013607 (2022), doi://10.1103/PhysRevMaterials.6.013607
81. *Exploring the necessary complexity of interatomic potentials*. J. A. Vita and D. R. Trinkle. *Comp. Mater. Sci.* **200**, 110752 (2021), doi://10.1016/j.commatsci.2021.110752
80. *Accelerated molecular dynamics simulations of dislocation climb in nickel*. L. T. W. Fey, A. M. Z. Tan, T. D. Swinburne, D. Perez, and D. R. Trinkle. *Phys. Rev. Materials* **5**, 083603 (2021), doi://10.1103/PhysRevMaterials.5.083603
79. *Stress-dependent dislocation core structures leading to non-schmid behavior*. D. Qiu, P. Zhao, D. R. Trinkle, and Y. Wang. *Mater. Res. Lett.* **9**(3), 134–140 (2021), doi://10.1080/21663831.2020.1854359
78. *Split-vacancy defect complexes of oxygen in hcp and fcc cobalt*. S. J. Honrao, Q. Rizzardi, R. Maaß, D. R. Trinkle, and R. G. Hennig. *Phys. Rev. Materials* **4**, 103608 (2020), doi://10.1103/PhysRevMaterials.4.103608
77. *Uncertainty quantification of solute transport coefficients*. R. Agarwal and D. R. Trinkle. In Y. Wang and D. L. McDowell, eds., *Uncertainty Quantification in Multiscale Materials Modeling*, Elsevier Series in Mechanics of Advanced Materials, chapter 4, 93–118 (Woodhead Publishing, 2020), doi://10.1016/B978-0-08-102941-1.00004-3
76. *Designing optimal perovskite structure for high ionic conduction*. R. Gao, A. C. P. Jain, S. Pandya, Y. Dong, Y. Yuan, H. Zhou, L. R. Dedon, V. Thoréton, S. Saremi, R. Xu, A. Luo, T. Chen, V. Gopalan, E. Ertekin, J. Kilner, T. Ishihara, N. H. Perry, D. R. Trinkle, and L. W. Martin. *Adv. Mater.* **32**(1), 1905178 (2020), doi://10.1002/adma.201905178

75. *Impact of solutes on the lattice parameters and elastic stiffness coefficients of hcp Fe from first-principles calculations.*
M. R. Fellingner, L. G. Hector, Jr., and D. R. Trinkle.
Comp. Mater. Sci. **164**, 116–126 (2019), doi://10.1016/j.commatsci.2019.03.056.
Database: hdl.handle.net/11256/982
74. *Dislocation core structures in Ni-based superalloys computed using density functional theory based flexible boundary condition approach.*
A. M. Z. Tan, C. Woodward, and D. R. Trinkle.
Phys. Rev. Mater. **3**, 033609 (2019), doi://10.1103/PhysRevMaterials.3.033609
73. *First principles calculations of solute transport in zirconium: Vacancy-mediated diffusion with metastable states and interstitial diffusion.*
A. C. P. Jain, P. A. Burr, and D. R. Trinkle.
Phys. Rev. Mater. **3**, 033402 (2019), doi://10.1103/PhysRevMaterials.3.033402.
Database: dx.doi.org/10.18126/M2R04K
72. *Variational principle for mass transport.*
D. R. Trinkle.
Phys. Rev. Lett. **121**, 235901 (2018), doi://10.1103/PhysRevLett.121.235901, (Editor's Suggestion)
71. *Geometries of edge and mixed dislocations in bcc Fe from first principles calculations.*
M. R. Fellingner, A. M. Z. Tan, L. G. Hector, Jr., and D. R. Trinkle.
Phys. Rev. Mater. **2**, 113605 (2018), doi://10.1103/PhysRevMaterials.2.113605.
Database: hdl.handle.net/11256/978
70. *Effect of solutes on the lattice parameters and elastic stiffness coefficients of body-centered tetragonal Fe.*
M. R. Fellingner, L. G. Hector, Jr., and D. R. Trinkle.
Comp. Mater. Sci. **152**, 308–323 (2018), doi://10.1016/j.commatsci.2018.05.021.
Database: hdl.handle.net/11256/959
69. *Modeling the long-term evolution of dilute solid solutions in the presence of vacancy fluxes.*
T. Schuler, P. Bellon, D. R. Trinkle, and R. S. Averback.
Phys. Rev. Mater. **2**, 073605 (2018), doi://10.1103/PhysRevMaterials.2.073605
68. *First-principles prediction of oxygen diffusivity near the (10 $\bar{1}$ 2) twin boundary in titanium.*
M. S. Hooshmand, C. Niu, D. R. Trinkle, and M. Ghazisaeidi.
Acta mater. **156**, 11–19 (2018), doi://10.1016/j.actamat.2018.05.076
67. *Ab initio magnesium-solute transport database using exact diffusion theory.*
R. Agarwal and D. R. Trinkle.
Acta mater. **150**, 339–350 (2018), doi://10.1016/j.actamat.2018.03.025.
Database: dx.doi.org/10.18126/M20G83
66. *Solute-induced solid-solution softening and hardening in bcc tungsten.*
Y.-J. Hu, M. R. Fellingner, B. G. Bulter, Y. Wang, K. A. Darling, L. J. Kecskes, D. R. Trinkle, and Z.-K. Liu.
Acta mater. **141**, 304–316 (2017), doi://10.1016/j.actamat.2017.09.019
65. *Automatic numerical evaluation of vacancy-mediated transport for arbitrary crystals: Onsager coefficients in the dilute limit using a Green function approach.*
D. R. Trinkle.
Philos. Mag. **97**, 2514–2563 (2017), doi://10.1080/14786435.2017.1340685
64. *Mechanical properties and phase stability of monoborides using density functional theory calculations.*

- H. Kim and D. R. Trinkle.
Phys. Rev. Mater. **1**, 013601 (2017), doi://10.1103/PhysRevMaterials.1.013601.
Database: dx.doi.org/10.18126/M24S3J
63. *Design principles for radiation-resistant solid solutions.*
T. Schuler, D. R. Trinkle, P. Bellon, and R. Averback.
Phys. Rev. B **95**, 174102 (2017), doi://10.1103/PhysRevB.95.174102
62. *Mesoscale modeling of vacancy-mediated Si segregation near an edge dislocation in Ni under irradiation.*
Z. Li and D. R. Trinkle.
Phys. Rev. B **95**, 144107 (2017), doi://10.1103/PhysRevB.95.144107
61. *Exact model of vacancy-mediated solute transport in magnesium.*
R. Agarwal and D. R. Trinkle.
Phys. Rev. Lett. **118**, 105901 (2017), doi://10.1103/PhysRevLett.118.105901
60. *Ab initio calculations of the lattice parameter and elastic stiffness coefficients of bcc Fe with solutes.*
M. R. Fellingner, L. G. Hector, Jr., and D. R. Trinkle.
Comp. Mater. Sci. **126**, 503–513 (2017), doi://10.1016/j.commatsci.2016.09.040.
Database: hdl.handle.net/11256/671
59. *First principles calculations of beryllium stability in zirconium surfaces.*
A. C. P. Jain and D. R. Trinkle.
Acta mater. **122**, 359–368 (2017), doi://10.1016/j.actamat.2016.10.003
58. *Reforming an undergraduate materials science curriculum with computational modules.*
R. Mansbach, A. Ferguson, K. Kilian, J. Krogstad, C. Leal, A. Schleife, D. R. Trinkle, M. West, and G. L. Herman.
J. Mater. Edu. **38**, 161–174 (2016)
57. *Ab initio modeling of quasielastic neutron scattering of hydrogen pipe diffusion in palladium.*
E. J. Schiavone and D. R. Trinkle.
Phys. Rev. B **94**, 054114 (2016), doi://10.1103/PhysRevB.94.054114
56. *Light element diffusion in Mg using first principles calculations: Anisotropy and elastodiffusion.*
R. Agarwal and D. R. Trinkle.
Phys. Rev. B **94**, 054106 (2016), doi://10.1103/PhysRevB.94.054106.
Database: hdl.handle.net/11256/694
55. *Computation of the lattice Green function for a dislocation.*
A. M. Z. Tan and D. R. Trinkle.
Phys. Rev. E **94**, 023308 (2016), doi://10.1103/PhysRevE.94.023308
54. *A modified embedded atom method potential for interstitial oxygen in titanium.*
P. Zhang and D. R. Trinkle.
Comp. Mater. Sci. **124**, 204–210 (2016), doi://10.1016/j.commatsci.2016.07.039.
Database: hdl.handle.net/11256/782
53. *Diffusivity and derivatives for interstitial solutes: Activation energy, volume, and elastodiffusion tensors.*
D. R. Trinkle.
Philos. Mag. **96**, 2714–2735 (2016), doi://10.1080/14786435.2016.1212175
52. *Oxygen diffusion in HCP metals.*
H. H. Wu, P. Wisesa, and D. R. Trinkle.
Phys. Rev. B **94**, 014307 (2016), doi://10.1103/PhysRevB.94.014307

51. *Kinetic monte carlo investigation of tetragonal strain on Onsager matrices.*
Z. Li and D. R. Trinkle.
Phys. Rev. E **93**, 053305 (2016), doi://10.1103/PhysRevE.93.053305
50. *Size misfits of point defects in aliovalently doped SrTiO₃ using density functional theory.*
H. Kim and D. R. Trinkle.
Comp. Mater. Sci. **119**, 41–45 (2016), doi://10.1016/j.commatsci.2016.02.040
49. *A study of stress relaxation in AZ31 using high-energy X-ray diffraction.*
W. Tang, K. L. Halm, D. R. Trinkle, M. Koker, U. Lienert, P. Kenesei, and A. J. Beaudoin.
Acta mater. **101**, 71–79 (2015), doi://10.1016/j.actamat.2015.08.072
48. *Energetics of rutile TiO₂ vicinal surfaces with <001> steps from the energy density method.*
B. Lee and D. R. Trinkle.
J. Phys. Chem. C **119**, 18203–18209 (2015), doi://10.1021/acs.jpcc.5b03623
47. *Database optimization for empirical interatomic potential models.*
P. Zhang and D. R. Trinkle.
Model. Simul. Mater. Sci. Eng. **23**, 065011 (2015), doi://10.1088/0965-0393/23/6/065011
46. *Importance of coordination number and bond length in titanium revealed by electronic structure investigations.*
L.-F. Huang, B. Grabowski, E. McEniry, D. R. Trinkle, and J. Neugebauer.
Phys. Stat. Solid. B **252**, 1907–1924 (2015), doi://10.1002/pssb.201552280
45. *Thermal transport across high-pressure semiconductor-metal transition in Si and Si_{0.991}Ge_{0.009}.*
G. T. Hohensee, M. R. Fellinger, D. R. Trinkle, and D. G. Cahill.
Phys. Rev. B **91**, 205104 (2015), doi://10.1103/PhysRevB.91.205104
44. *Calculation of strain effects on vacancy-mediated diffusion of impurities in FCC structures: General approach and application to Ni_{1-x}Si.*
T. Garnier, Z. Li, M. Nastar, P. Bellon, and D. R. Trinkle.
Phys. Rev. B **90**, 184301 (2014), doi://10.1103/PhysRevB.90.184301
43. *Diffusion of Si impurities in Ni under stress: A first-principles study.*
T. Garnier, V. R. Manga, P. Bellon, and D. R. Trinkle.
Phys. Rev. B **90**, 024306 (2014), doi://10.1103/PhysRevB.90.024306.
Database: hdl.handle.net/11115/239
42. *Direct measurement of hydrogen dislocation pipe diffusion in deformed polycrystalline Pd using quasielastic neutron scattering.*
B. J. Heuser, D. R. Trinkle, N. Jalarvo, J. Serio, E. J. Schiavone, E. Mamontov, and M. Tyagi.
Phys. Rev. Lett. **113**, 025504 (2014), doi://10.1103/PhysRevLett.113.025504
41. *Interaction of oxygen interstitials with lattice faults in Ti.*
M. Ghazisaeidi and D. R. Trinkle.
Acta mater. **76**, 82–86 (2014), doi://10.1016/j.actamat.2014.05.025
40. *First-principles study of interfacial boundaries in Ni–Ni₃Al.*
C. Woodward, A. van de Walle, M. Asta, and D. R. Trinkle.
Acta mater. **75**, 60–70 (2014), doi://10.1016/j.actamat.2014.04.056
39. *Quantitative modeling of solute drag by vacancies in face-centered-cubic alloys.*
T. Garnier, D. R. Trinkle, M. Nastar, and P. Bellon.
Phys. Rev. B **89**, 144202 (2014), doi://10.1103/PhysRevB.89.144202
38. *Stress-induced anisotropic diffusion in alloys: Complex Si solute flow near a dislocation core in Ni.*
T. Garnier, V. R. Manga, D. R. Trinkle, M. Nastar, and P. Bellon.
Phys. Rev. B **88**, 134108 (2013), doi://10.1103/PhysRevB.88.134108

37. *Solute drag by vacancies in body-centered cubic alloys.*
T. Garnier, M. Nastar, P. Bellon, and D. R. Trinkle.
Phys. Rev. B **88**, 134201 (2013), doi://10.1103/PhysRevB.88.134201
36. *Solute effect on oxygen diffusion in α -titanium.*
H. H. Wu and D. R. Trinkle.
J. Appl. Phys. **113**, 223504 (2013), doi://10.1063/1.4808283.
Database: hdl.handle.net/11115/125
35. *Hydrogen trapping at dislocation cores at room temperature in deformed Pd.*
B. J. Heuser, D. R. Trinkle, T.-N. Yang, and L. He.
J. Alloy. Compd. **577**, 189–191 (2013), doi://10.1016/j.jallcom.2013.04.082
34. *Ab initio based empirical potential used to study the mechanical properties of molybdenum.*
H. Park, M. R. Fellingner, T. J. Lenosky, W. W. Tipton, D. R. Trinkle, S. P. Rudin, C. Woodward, J. W. Wilkins, and R. G. Hennig.
Phys. Rev. B **85**, 214121 (2012), doi://10.1103/PhysRevB.85.214121
33. *Direct calculation of lattice green function with arbitrary interactions for general crystals.*
J. A. Yasi and D. R. Trinkle.
Phys. Rev. E **85**, 066706 (2012), doi://10.1103/PhysRevE.85.066706
32. *Prediction of thermal cross-slip stress in magnesium alloys from a geometric interaction model.*
J. A. Yasi, L. G. Hector, Jr., and D. R. Trinkle.
Acta mater. **60**, 2350–2358 (2012), doi://10.1016/j.actamat.2012.01.004
31. *Core structure of a screw dislocation in Ti from density functional theory and classical potentials.*
M. Ghazisaeidi and D. R. Trinkle.
Acta mater. **60**, 1287–1292 (2012), doi://10.1016/j.actamat.2011.11.024
30. *Au/TiO₂(110) interfacial reconstruction stability from ab initio.*
M. Yu and D. R. Trinkle.
J. Phys. Chem. C **115**, 17799–17805 (2011), doi://10.1021/jp2017133
29. *Direct diffusion through interpenetrating networks: Oxygen in titanium.*
H. H. Wu and D. R. Trinkle.
Phys. Rev. Lett. **107**, 045504 (2011), doi://10.1103/PhysRevLett.107.045504
28. *Prediction of thermal cross-slip stress in magnesium alloys from direct first principles data.*
J. A. Yasi, L. G. Hector, Jr., and D. R. Trinkle.
Acta mater. **59**, 5652–5660 (2011), doi://10.1016/j.actamat.2011.05.040
27. *Nanoscale-hydride formation at dislocations in palladium: Ab initio theory and inelastic neutron scattering measurements.*
D. R. Trinkle, H. Ju, B. J. Heuser, and T. J. Udovic.
Phys. Rev. B **83**, 174116 (2011), doi://10.1103/PhysRevB.83.174116
26. *Thermal conductivity of compressed H₂O to 22 GPa: A test of the Leibfried-Schlömann equation.*
B. Chen, W.-P. Hsieh, D. G. Cahill, D. R. Trinkle, and J. Li.
Phys. Rev. B **83**, 132301 (2011), doi://10.1103/PhysRevB.83.132301
25. *Energy density in density functional theory: Application to crystalline defects and surfaces.*
M. Yu, D. R. Trinkle, and R. M. Martin.
Phys. Rev. B **83**, 115113 (2011), doi://10.1103/PhysRevB.83.115113

24. *Accurate and efficient algorithm for Bader charge integration.*
M. Yu and D. R. Trinkle.
J. Chem. Phys. **134**, 064111 (2011), doi://10.1063/1.3553716
23. *First-principles calculation of H vibrational excitations at a dislocation core of Pd.*
H. M. Lawler and D. R. Trinkle.
Phys. Rev. B **82**, 172101 (2010), doi://10.1103/PhysRevB.82.172101
22. *Misfit-dislocation-mediated island diffusion: Cu on Ag(111).*
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14. *Convergence rate for numerical computation of the lattice Green's function.*
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 1. *Bound states in waveguides and bent quantum wires.*
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Submitted Papers

1. *Discovering melting temperature prediction models of inorganic solids by combining supervised and unsupervised learning.*
V. Gharakhanyan, L. J. Wirth, J. A. G. Torres, E. Eisenberg, T. Wang, D. R. Trinkle, S. Chatterjee, and A. Urban, (under review), [arXiv:2403.03092](https://arxiv.org/abs/2403.03092) (2024)

Open-source Software and Databases

16. Software: ONSAGER, D. R. Trinkle.
<http://dallastrinkle.github.io/Onsager>, v1.2.2 (2017), [doi://10.5281/zenodo.494987](https://doi.org/10.5281/zenodo.494987)

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14. Database: *Data citation: Kinson data for fcc random and high entropy alloys*, D. R. Trinkle and S. Chattopadhyay (2024), [doi://10.5281/zenodo.10214334](https://doi.org/10.5281/zenodo.10214334).
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Database: dx.doi.org/10.25584/1811899
12. Database: *Data citation: Impact of solutes on the lattice parameters and elastic stiffness coefficients of hcp Fe from first-principles calculations*, M. R. Fellingner, L. G. Hector, Jr., and D. R. Trinkle (2019).
Database: hdl.handle.net/11256/982
11. Database: *Data citation: Vacancy mediated and interstitial solute transport in Zr from density functional theory calculations*, A. C. P. Jain, P. A. Burr, and D. R. Trinkle (2018).
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Database: hdl.handle.net/11256/671
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6. Database: *Data citation: Fitting database entries for a modified embedded atom method potential for interstitial oxygen in titanium*, D. R. Trinkle and P. Zhang (2016).
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5. Database: *Data citation: Light element diffusion in Mg using first principles calculations: Anisotropy and elastodiffusion*, R. Agarwal and D. R. Trinkle (2016).
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Data files for ab initio calculations of the lattice parameter and elastic stiffness coefficients of BCC Fe with solutes.
M. R. Fellingner, L. G. Hector, Jr., and D. R. Trinkle.
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3. Database: *Data citation: Diffusion of Si impurities in Ni under stress: A first-principles study*, T. Garnier, V. R. Manga, P. Bellon, and D. R. Trinkle (2014).
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1. Database: *Dallas R. Trinkle Research Group Magnesium Database*, J. A. Yasi and D. R. Trinkle (2013).
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Conference Proceedings and Other Publications

7. *Applying for computational time on NSF's TERAgrid—the world's largest cyberinfrastructure supporting open research.*
K. Thornton, D. R. Trinkle, and R. G. Hennig.
JOM **62**, 17–18 (2010), [doi://10.1007/s11837-010-0042-6](https://doi.org/10.1007/s11837-010-0042-6)
6. *Hydrogen trapping in dislocations in palladium at low temperature: Results from inelastic neutron scattering and advanced computations.*
B. J. Heuser, H. Ju, D. R. Trinkle, and T. J. Udovic.
In *Effects of Hydrogen on Materials*, 2008 International Hydrogen Conference, 464–468 (2009)
5. *Ab-initio molecular dynamics simulations of molten Ni-based superalloys.*
M. Asta, D. R. Trinkle, and C. Woodward.
In *2007 High Performance Computing User's Group Conference*, 147–152 (IEEE Computing Society, Los Alamitos, CA, 2007)
4. *Ab-initio molecular dynamics simulations of molten Ni-based superalloys.*
M. Asta, D. R. Trinkle, and C. Woodward.
In *2006 High Performance Computing User's Group Conference*, 177–181 (IEEE Computing Society, Los Alamitos, CA, 2007)
3. *Alpha-omega transition in Ti: Equation of state and kinetics.*
C. W. Greeff, D. R. Trinkle, and R. C. Albers.
AIP Conference Proceedings **620**(1), 225–228 (2002)
2. *Actinide electronic structure and atomic forces.*
R. C. Albers, S. P. Rudin, D. R. Trinkle, and M. D. Jones.
AIP Conference Proceedings **532**, 412–413 (2000)
1. *Amazing magnetic moments.*
R. Singleton, D. Trinkle, and T. Toepker.
Phys. Teach. **35**(2), 122–125 (1997), [doi://10.1119/1.2344615](https://doi.org/10.1119/1.2344615)

Invited Talks

—2024

116. *Understanding diffusion in complex materials using machine learning and a variational approach.*
Data Sciences for Mesoscale and Macroscale Materials Models workshop. Institute for Mathematical and Statistical Innovation, Univ. Chicago, Chicago, IL, May 13–17, 2024.
115. *Kinosons: Quantifying the contributions to diffusion in complex materials.*
Materials Science and Engineering Colloquium. Carnegie Mellon University, Pittsburgh, PA, April 4, 2024.
114. *Understanding diffusion processes in a multicomponent alloy using a variational approach.*
Annual Meeting of the Minerals, Metals & Materials Society (TMS), "Algorithm Development in Materials Science and Engineering" symposium. Orlando, FL, March 4–March 7, 2024.

113. *Isolated dislocation core energy from first principles energy density method.*
Annual Meeting of the Minerals, Metals & Materials Society (TMS), “Mechanical Behavior at the Nanoscale VII” symposium. Orlando, FL, March 4–March 7, 2024.
- 2023
112. *Diffusivity in a multicomponent alloy using machine learning and variational approaches.*
New Horizons in materials design workshop. Max-Planck-Institut für Eisenforschung, Düsseldorf, Germany, August 18, 2023.
111. *Designing magnesium alloys from density-functional theory and atomistic models.*
American Physical Society March Meeting. Las Vegas, NV, March 5, 2023.
- 2022
110. *Efficient mass transport calculations using a variational principle.*
Multiscale Modeling of Materials Meeting, Stochastic methods in materials simulation symposium. Baltimore, MD, October 10, 2022.
109. *Isolated Dislocation Core Energy from First Principles Energy Density Method.*
Multiscale Modeling of Materials Meeting, Multiscale Materials Modeling using Ab-initio Accuracy Methods symposium. Baltimore, MD, October 10, 2022.
108. *Solute transport in zirconium with nonequilibrium vacancy concentration.*
University of Wisconsin, 4th Computing in Engineering Forum. Virtual, September 20, 2022.
107. *Machine learning in diffusivity calculations using a variational principle.*
Los Alamos National Laboratory, Center for Nonlinear Science. Los Alamos, NM, July 25, 2022.
106. *Machine learning in diffusivity calculations using a variational principle.*
Annual Meeting of the Minerals, Metals & Materials Society, “Hume-Rothery Symposium” symposium. Anaheim, CA, February 27–March 2, 2022.
- 2021
105. *Dislocation environment and hydrogen: Nanoscale hydrides and pipe diffusion in palladium*
SFB-Workshop on “Defect Phases”. RWTH Aachen (virtual), September 9, 2021.
104. *Machine learning in diffusivity calculations using a variational principle.*
“Complex High-Dimensional Energy Landscapes” reunion workshop. Institute for Pure and Applied Mathematics, UCLA, June 7, 2021.
- 2020
103. *Variational Principle for Mass Transport.*
CECAM meeting, Ab Initio Description of Iron and Steel. Ringsberg, Germany, November 2–5, 2020.
102. *Achieving machine-learning-potential accuracy with traditional classical potential forms*
NIST “Atomistic Simulations for Industrial Needs” workshop. Virtual, August 5–7, 2020.
101. *Isolated dislocation core energy from first principles energy density method*
Annual Meeting of the Minerals, Metals & Materials Society, “Algorithm Development in Materials Science and Engineering” symposium. San Diego, CA, February 24–27, 2020.
- 2019
100. *Efficient mass transport calculations using a variational principle*
Fall Meeting of the Materials Research Society. Boston, Massachusetts, December 3, 2019.
99. *Efficient mass transport calculations using a variational principle*
Materials Science and Engineering colloquium, University of British Columbia, Vancouver, BC, November 25, 2019.

98. *Efficient mass transport calculations using a variational principle*
Materials Science and Engineering colloquium, Ohio State University, Columbus, OH, November 15, 2019.
97. *A variational principle for mass transport calculations*
PRISMS workshop. Ann Arbor, MI, August 8, 2019.
96. *A variational principle for mass transport calculations II*
“Complex High-Dimensional Energy Landscapes” reunion workshop. Institute for Pure and Applied Mathematics, UCLA, June 13, 2019.
95. *A variational principle for mass transport calculations*
“Complex High-Dimensional Energy Landscapes” reunion workshop. Institute for Pure and Applied Mathematics, UCLA, June 12, 2019.
94. *A variational principle for mass transport calculations*
Annual Meeting of the Minerals, Metals & Materials Society, “Algorithm Development in Materials Science and Engineering” symposium. San Antonio, TX, March 10–14, 2019.
93. *Uncertainty quantification for solute transport modeling*
Annual Meeting of the Minerals, Metals & Materials Society, “Hume-Rothery Symposium” symposium. San Antonio, TX, March 10–14, 2019.
92. *A variational principle for mass transport calculations*
Annual Meeting of the Minerals, Metals & Materials Society, “Algorithm Development in Materials Science and Engineering” symposium. San Antonio, TX, March 10–14, 2019.
91. *A variational principle for mass transport calculations.*
Center for Advanced Study, University of Illinois, Urbana-Champaign, February 5, 2019

—2018

90. *Computing mass transport in solids.*
University of Warwick. Warwick, England, August 20, 2018.
89. *Lattice Green function calculations in the presence of a dislocation.*
University of Warwick. Warwick, England, August 15, 2018.
88. *A variational principle for mass transport calculations.*
CNLS 38th Annual Conference: Rate Theory and Full Timescale Simulations. Center for Nonlinear Science, Santa Fe, New Mexico, May 29, 2018.
87. *Exact theory of solute transport in zirconium.*
Joint ICTP-IAEA School and Workshop on Fundamental Methods for Atomic, Molecular and Materials Properties in Plasma Environments. International Centre for Theoretical Physics, Trieste, Italy, April 16, 2018.
86. *Lattice Green function calculations in the presence of a dislocation.*
Applied Math Seminar. University of California, Berkeley, April 4, 2018.

—2017

85. *Uncertainty quantification for solute transport modeling.*
“Complex High-Dimensional Energy Landscapes” workshop. Institute for Pure and Applied Mathematics, UCLA, December 14, 2017.
84. *Uncertainty quantification for solute transport modeling.*
Fall Meeting of the Materials Research Society. Boston, Massachusetts, November 27, 2017.
83. *Exact theory of vacancy-mediated solute transport in magnesium.*
EuroMat 2017: European Congress and Exhibition on Advanced Materials and Processes. Thessaloniki, Greece, September 18, 2017.

82. *Dislocation cores from first principles: Ni and Ni₃Al for superalloys.*
US National Congress on Computational Mechanics (USNCCM14). Montreal, Canada, July 17, 2017.
81. *Dislocation cores and defect interactions from first principles: Current state of the art and new challenges.*
“Density Functional Theory and Beyond: Analysis and Computation” workshop, Mathematics Institute of the University of Warwick. Warwick, England, July 3, 2017.
80. *Automated Green function calculation of dilute solute diffusivity.*
DIMAT 2017: International Conference on Diffusion in Materials. Haifa, Israel, May 7, 2017.
79. *Exact theory of vacancy-mediated solute transport in magnesium.*
Magnesium Workshop: Alloys and Lightweight Structural Systems, Johns Hopkins University. Baltimore, Maryland, April 3, 2017.
78. *Dislocation core structures in Ni-based superalloys computed using density functional theory based flexible boundary condition approach.*
Modeling and Simulation of Superalloys International Workshop, Ruhr-Universität. Bochum, Germany, March 29, 2017.
77. *Computing mass transport in crystals: Theory, computation, and applications.*
Max-Planck-Institut für Eisenforschung, Düsseldorf, Germany, March 27, 2017.
76. *Computing mass transport in crystals: Theory, computation, and applications.*
Materials Science and Engineering Seminar, Univ. California, Berkeley. February 24, 2017.

—2016

75. *Computing mass transport in crystals: Theory, computation, and applications.*
William Maxwell Reid Seminar, Univ. Kentucky. Lexington, KY, October 28, 2016.
74. *Solute interactions with dislocations in BCC Fe.*
CECAM meeting, Ab Initio Description of Iron and Steel. Ringsberg, Germany, October 3–7, 2016.
73. *Dislocation cores and defect interactions from first principles: Current state of the art and new challenges.*
Society for Engineering Science, 53rd Annual Technical Meeting. College Park, MD, October 2–5, 2016.
72. *Bridging timescales: Automating diffusivity calculations for interstitial and solute diffusion from first-principles.*
Recent Advances in Computational Methods for Nanoscale Phenomena workshop. Ann Arbor, MI, August 29–31, 2016.
71. *Database optimization for empirical interatomic potentials.*
Annual Meeting of the Minerals, Metals & Materials Society, “Computational Methods for Uncertainty Quantification, Model Validation, and Stochastic Predictions” symposium. Nashville, TN, February 14–18, 2016.
70. *Hydrides and Hydrogen Pipe Diffusion in Palladium: First Principles, Kinetic Monte Carlo, and Neutron Scattering.*
Annual Meeting of the Minerals, Metals & Materials Society, “Hume-Rothery” symposium. Nashville, TN, February 14–18, 2016.

—2015

69. *Automating Computation of Solute Diffusion.*
Materials Science and Technology Meeting, “Data and Tools for Materials Discovery and Design” symposium. Columbus, OH, October 4–8, 2015.

68. *First-principles computation of a stepped surface: TiO₂*
Materials Science and Technology Meeting, "Deformation and Transitions at Grain Boundaries IV" symposium. Columbus, OH, October 4–8, 2015.
 67. *Hydrides and Hydrogen Pipe Diffusion in Palladium: First Principles, Kinetic Monte Carlo, and Neutron Scattering.*
Service de Recherches de Metallurgie Physique seminar, CEA/Saclay, September 14, 2015.
 66. *Automating diffusivity calculations for interstitial and solute diffusion from first-principles.*
2015 Psi-k Conference, San Sebastián / Donostia, Spain, September 7–10, 2015.
 65. *Multiscale evolution of defects near a dislocation: Solutes and vacancies in nickel.*
"Defects in Materials," Lake Arrowhead reunion program. Institute for Pure and Applied Mathematics, UCLA, June 8, 2015.
 64. *Light interstitial diffusion in magnesium.*
NIST Diffusion workshop, Gaithersburg, MD, May 14–15, 2015.
 63. *Evolution of defects near a dislocation: Solutes and vacancies in nickel.*
Annual Meeting of the Minerals, Metals & Materials Society, "Computational Thermodynamics and Kinetics" symposium. Orlando, FL, March 15–19, 2015.
- 2014
62. *Forming nanoprecipitates at dislocations: Hydrides in Pd and Silicides in Ni.*
Max-Planck-Institut für Eisenforschung, Düsseldorf, Germany, November 5, 2014.
 61. *Stress and irradiation effects on solute diffusion to dislocations.*
Multiscale Modeling of Materials Meeting, Berkeley, CA, October 6, 2014.
 60. *Lattice Green function calculations in the presence of a dislocation*
"Multiscale Models of Crystal Defects" program, Banff International Research Station for Mathematical Innovation and Discovery, September 21, 2014.
 59. *Hydrogen Pipe Diffusion in Palladium: First Principles, Kinetic Monte Carlo, and Experiments*
14th International Symposium on Metal-Hydrogen Systems, Univ. Salford, Manchester, UK, July 25, 2014.
 58. *Lattice Green function calculations in the presence of a dislocation*
"Defects in Materials," Lake Arrowhead reunion program. Institute for Pure and Applied Mathematics, UCLA, June 9, 2014.
 57. *Metal solute-oxygen interaction in titanium for changes in oxygen diffusivity*
NIST Diffusion workshop, Gaithersburg, MD, April 28–29, 2014.
 56. *Forming nanoprecipitates at dislocations: Hydrides in Pd and Silicides in Ni.*
Spring Meeting of the Materials Research Society, San Francisco, CA, April 21–25, 2014.
 55. *Stress and Irradiation Effects on Solute Diffusion to Dislocations.*
Annual Meeting of the American Chemical Society, "Radiation Hardened Materials for Accelerators, Reactors and Spacecraft" symposium. Dallas, TX, March 16–20, 2014.
 54. *Deformation of Magnesium from First Principles.*
2nd ESISM Workshop in Kyoto, "Fundamental Issues of Structural Materials." Kyoto, Japan, March 11–12, 2014.
 53. *Density-functional theory methods for interfaces: Lattice Greens Function and Energy Density Methods.*
Annual Meeting of the Minerals, Metals & Materials Society, "Mechanical Behavior Related to Interface Physics II" symposium. San Diego, CA, February 16–20, 2014.

—2013

52. *Interfaces in HCP Ti: twinning, transformation, oxygen.*
Materials Science and Technology Meeting, "Deformation and Transitions at Grain Boundaries III" symposium. Montreal, Canada, October 28–31, 2013.
51. *Dislocations with first principles: From defects to material behavior in magnesium.*
Computational Science and Engineering, Fall Seminar, Univ. Illinois, Urbana-Champaign, October 23, 2013.
50. *First principles modeling of titanium: oxygen, dislocations, and boundaries.*
Keynote lecture; Pacific Rim International Congress on Advanced Materials and Processing (PRICM-8), Light Metals and Alloys symposium, Waikoloa, Hawaii, August 4–9, 2013.
49. *Energy Density Method and Bader Integration for Defect Energies.*
International Symposium on Atomistic Modeling for Mechanics and Multiphysics of Materials (ISAM⁴ 2013), Tokyo, Japan, July 22, 2013.
48. *Deformation in magnesium from first-principles.*
Max-Planck-Institut für Eisenforschung, Düsseldorf, Germany, July 10, 2013.
47. *Beyond binaries: Solid solution strengthening and softening in magnesium.*
Magnesium Workshop Madrid 2013, Madrid, Spain, May 21–24, 2013.
46. *Stress-induced anisotropic diffusion in alloys: Complex Si solute flow near a dislocation core in Ni.*
NIST Diffusion workshop, Gaithersburg, MD, May 9–10, 2013.
45. *Predicting strength and cross-slip of magnesium alloys: First-principles, solute distribution, and deformation.*
Annual Meeting of the Minerals, Metals & Materials Society, "Modeling and Experimental Validation of Multiscale Mechanical Behavior from Atomic Scale to Macro Scale" symposium. San Antonio, TX, March 4–March 7, 2013.

—2012

44. *Solute diffusion under stress around dislocations: Si in Ni.*
Workshop on "Simulation of Complex Microstructure Pathways for Alloy Design." McMaster University, Hamilton, CA, November 29–30, 2012.
43. *Deformation in magnesium from first-principles.*
 - (a) Physics Colloquium, California State Univ., Northridge, November 9, 2012.
 - (b) Materials Science and Engineering Colloquium, Univ. British-Columbia, November 6, 2012.
 - (c) Physics Colloquium, Arizona State Univ., November 5, 2012.
 - (d) Materials Science and Engineering Colloquium, Ohio State Univ., October 5, 2012.
 - (e) Materials Science and Engineering Colloquium, Univ. Florida, September 19, 2012.
42. *Dislocation cores and defect interactions from first principles: Current state of the art and new challenges.*
"Atomistic and Mesoscale Modeling of Materials Defects" workshop. Institute for Pure and Applied Mathematics, UCLA, October 22, 2012.
41. *Solute diffusion to dislocations under irradiation: Si in Ni.*
Materials Science and Technology Meeting, "Materials Development for Nuclear Applications and Extreme Environments" symposium. Pittsburgh, PA, October 7–11, 2012.
40. *From dislocation/solute interactions and solute distribution to deformation in magnesium.*
9th International Conference on Magnesium Alloys and their Applications. Vancouver, BC, Canada, July 8–12, 2012.
39. *Deformation in magnesium from first-principles.*

- (a) Materials Science and Engineering Colloquium, Los Alamos National Laboratory, June 19, 2012.
- (b) NIST Metallurgy Colloquium, July 6, 2012.
- 38. *The Challenge of Multicomponent Systems: Atomistics, Thermodynamics and Kinetics*. NSF Workshop on Emerging Research Areas in the Science of Metals and Metallic Nanostructures. Univ. California, Santa Barbara, June 13–14, 2012.
- 37. *Atomic-scale modeling and data*. NIST “Materials Genome Initiative Workshop on Building the Materials Innovation Infrastructure: Data and Standards.” U.S. Department of Commerce, Washington, D.C., May 14–15, 2012.
- 36. *Deformation in magnesium from first-principles*. Annual Meeting of the Minerals, Metals & Materials Society (TMS), “Phase Transformations and Deformation in Magnesium Alloys” symposium. Orlando, FL, March 11–March 15, 2012.
- 35. *First Principles Modeling of Dislocation/Twin Boundary Interactions in Titanium*. International Conference on Plasticity. Puerto Rico, January 3–8, 2012.
- 2011
- 34. *Deformation in magnesium from first-principles*. U.S. Army Research Laboratory speaker series on Integrated Computational Materials Engineering, Aberdeen Proving Ground, MD, November 16, 2011.
- 33. *First-principles predictions of dislocation/solute interactions: Rare earths in magnesium*. MagNET Research Network Workshop, “Rare Earth Elements in Magnesium Alloys.” CANMET Facility, Hamilton, Ontario, November 7–8, 2011.
- 32. *Interactions of dislocations and twin boundaries in titanium with first-principles*. Materials Science and Technology Meeting, “Deformation and Transitions at Grain Boundaries” symposium. Columbus, OH, October 16–20, 2011.
- 31. *Dislocations with first principles: From defects to material behavior in magnesium*. Materials Science and Engineering Colloquium, Massachusetts Institute of Technology, October 13, 2011.
- 30. *First principles modeling of dislocations for magnesium alloys*. Keynote lecture; NSF Workshop on Magnesium Science and Technology, Arlington, VA, May 19–20, 2011.
- 29. *Dislocations with first principles: From defects to material behavior*.
 - (a) Materials Science and Engineering Colloquium, Univ. Michigan, April 15, 2011.
 - (b) Materials Science and Engineering Colloquium, Cornell University, April 21, 2011.
- 28. *Formation of hydrogen Cottrell atmosphere in palladium: Theory and measurement from inelastic neutron scattering*. Annual Meeting of the Minerals, Metals & Materials Society (TMS), “Hydrogen Storage in Materials: Theory and Experiment” symposium. San Diego, CA, February 27–March 3, 2011.
- 27. *Predicting Mg strength from first-principles: Solid-solution strengthening, softening, and cross-slip*. Keynote lecture; Annual Meeting of the Minerals, Metals & Materials Society (TMS), “Magnesium Technology 2011” symposium. San Diego, CA, February 27–March 3, 2011.
- 26. *Dislocations with first principles: From defects to material behavior*.
 - (a) Materials Science and Engineering Colloquium, Univ. California, Berkeley, February 24, 2011.

(b) Lawrence Livermore National Laboratory, February 25, 2011.

—2010

25. *Dislocations from first principles: Connecting defects to macroscale behavior in real materials.*

(a) Mechanical, Aerospace, and Nuclear Engineering Colloquium, Rensselaer Polytechnic Institute, November 10, 2010.

(b) Materials Science and Engineering Colloquium, California Institute of Technology, December 8.

(c) Physics and Theory Colloquium, Los Alamos National Laboratory, December 9, 2010.

—2009

24. *Point, line, and planar defects and interactions using electronic structure methods.*

CAMM meeting, “Possibilities and Limitations of Characterization and Modeling Solid/Solid Interfaces.” Bernkastel-Kues, Germany, May 17–20, 2009.

23. *Palladium-hydrogen interaction in dislocations: Trapping and diffusion.*

Annual Meeting of the Minerals, Metals & Materials Society (TMS), “Computational Thermodynamics and Kinetics” symposium. San Francisco, CA, February 15–20, 2009.

22. *Interface mobility for Ti alpha to omega transformation.*

Annual Meeting of the Minerals, Metals & Materials Society (TMS), “Transformations Under Extreme Conditions” symposium. San Francisco, CA, February 15–20, 2009.

21. *Chemical short-range ordering in liquid-phase Ni alloys.*

Annual Meeting of the Minerals, Metals & Materials Society (TMS), “Advanced Characterization and Modeling of Phase Transformations in Metals” symposium. San Francisco, CA, February 15–20, 2009.

20. *Electronic structure in dislocation modeling: Dislocation/solute interactions in Mo, Al, and Mg.*

Joint Materials/Solid Mechanics Seminar Series, Brown Univ., January 30, 2009.

19. *Dislocation/solute interactions via first-principles for modeling strength of materials: Mo, Al, Mg.*

International Conference on Plasticity. St. Thomas, January 3–8, 2009.

—2008

18. *First-principles calculation of dislocation/solute interaction: Solid-solution softening, dynamic strain-aging, and hydrogen trapping.*

General Motors Research and Development, April 24, 2008.

17. *Electronic structure for dislocation/defect interactions.*

Japan Institute of Metals Annual Meeting, Tokyo, Japan, March 28, 2008.

—2007

16. *First-principles calculation of dislocation/solute interaction: Solid-solution softening, dynamic strain-aging, and hydrogen trapping.*

Materials Science and Engineering, Univ. Wisconsin, December 14, 2007.

15. *Chemistry of deformation: Solid-solution softening from first-principles to dislocation mobility.*

Service de Recherches de Metallurgie Physique seminar, CEA/Saclay, June 26, 2007.

14. *Implementing flexible-boundary conditions and computing the lattice Green function.*

Service de Recherches de Metallurgie Physique, Dislocation simulation workshop, CEA/Saclay, June 21, 2007.

13. *Density-functional theory, flexible-boundary conditions, and predicting dislocation-core structures.*
Service de Recherches de Metallurgie Physique, Dislocation simulation workshop, CEA/Saclay, June 21, 2007.
 12. *Material strength from first principles: Electronic structure and the solute/dislocation interaction.*
Electronic Structure Workshop, North Carolina State University, June 12–15, 2007.
 11. *Chemistry of deformation: From atomic bonding to dislocations to strength.*
Joint Metallurgical Engineering and Material Science & Engineering graduate seminar, University of Utah, March 28, 2007.
 10. *Chemistry of deformation: Solid-solution softening from first-principles to dislocation mobility.*
Annual Meeting of the Minerals, Metals & Materials Society (TMS), “Plasticity from the Atomic Scale to Constitutive Laws” symposium. Orlando, FL, February 25–March 1, 2007.
 9. *Martensitic alpha to omega in titanium: Atomic pathway and impurity effects.*
Annual Meeting of the Minerals, Metals & Materials Society (TMS), “Fundamentals of Shape Memory and Related Transitions” symposium. Orlando, FL, February 25–March 1, 2007.
- 2006
8. *The simulation of material deformation.*
Physics in Careers seminar, Xavier University, October 20, 2006.
 7. *The chemistry of deformation.*
 - (a) Materials Science and Engineering colloquium, Univ. Illinois, Urbana-Champaign, January 23, 2006.
 - (b) Mechanical Engineering seminar, Johns Hopkins University, February 7, 2006.
 - (c) Materials Science and Engineering seminar, Northwestern University, March 1, 2006.
- 2005
6. *Alpha to omega in titanium alloys: Martensitic phase transitions at the atomic length scale.*
Materials Science and Technology, “Modeling and Simulation of Titanium Technology: Theory and Practices” symposium. Pittsburgh, PA, September 25–28, 2005.
 5. *Martensitic phase transitions at the atomic length scale: Titanium alpha to omega.*
March Meeting of the American Physical Society, Los Angeles, March 21–25, 2005.
- 2003
4. *Atomistic mechanisms of the martensitic titanium hcp to omega phase transition.*
Air Force Research Laboratory, Wright-Patterson AFB, August 20, 2003.
 3. *Atomistic mechanisms of martensitic phase transformations: Titanium hcp to omega.*
Electronic Structure Workshop, University of Minnesota, May 17–19, 2003.
- 2001
2. *Investigating transition pathways with symmetry analysis and the nudged-elastic band method.*
Condensed Matter seminar, Brigham Young University, October 18, 2001.
- 1999
1. *Total energy and elastic properties of titanium via tight-binding.*
Condensed Matter seminar, University of Cincinnati, July 14, 1999.

Conference Contributions

29. *Ab initio kinetic Monte Carlo study of hydrogen pipe diffusion.*
2013 Gordon Research Conference on Hydrogen-Metal Systems, July 14.
28. *High-throughput solute interaction for deformation in magnesium from first-principles.*
2012 Fall Meeting of the Materials Research Society, Materials Informatics Symposium, Boston, November 25.
27. *Predicting single crystal flow stress for magnesium alloys from first principles dislocation/solute interactions and solute distribution.*
2012 Multiscale Modeling of Materials Meeting, Singapore, October 15, 2012.
26. *First principles dislocation/solute interactions and deformation in magnesium alloys.*
2012 Dislocations, Budapest, Hungary, August 27, 2012.
25. *Interfacial reconstruction of Au/TiO₂ from ab initio.*
2012 Annual Meeting of the Minerals, Metals & Materials Society (TMS), "Solid-State Interfaces II: Toward an Atomistic-Scale Understanding of Structure, Properties, and Behavior through Theory and Experiment" symposium. Orlando, FL, March 11–March 15.
24. *Au/TiO₂ interfacial reconstruction stability from ab initio.*
2011 Materials Science and Technology Meeting, "Interfaces, Grain Boundaries and Surfaces from Atomistic and Macroscopic Approaches—Fundamental and Engineering Issues" symposium. Columbus, OH, October 16–20.
23. *Fundamentals of hydrogen and hydride formation at defects in palladium: First-principles prediction and inelastic neutron scattering measurements.*
2011 Spring Meeting of the Materials Research Society, San Francisco, April 23–27.
22. *First-principles calculation of dislocation/solute interactions and modeling of strength in Mg.*
Multiscale Materials Modeling 2010, Freiburg, Germany, October 4–8.
21. *Ab initio modeling to improve oxygen tolerance of Ti alloys.*
Multiscale Materials Modeling 2010, Freiburg, Germany, October 4–8.
20. *H trapping and diffusion in Pd dislocation cores.*
2008 Fall Meeting of the Materials Research Society, Boston, December 1–5.
19. *Ab initio calculation of H in Pd dislocation core.*
2007 Fall Meeting of the Materials Research Society, Boston, November 26–30.
18. *New quantitative analysis explains softening of pure metals by solutes.*
2006 Annual Meeting of the Minerals, Metals & Materials Society (TMS), San Antonio, March 12–16.
17. *Quantitative modeling of solid-solution softening in bcc Mo at the atomistic scale.*
2005 Fall Meeting of the Materials Research Society, Boston, November 28–December 2.
16. *Intrinsic solid-solution softening in bcc Mo from dislocation-solute interactions.*
2005 March Meeting of the American Physical Society, Los Angeles, March 21–25.
15. *Solid-solution softening trends in bcc Mo by first principles.*
2005 Annual Meeting of the Minerals, Metals & Materials Society (TMS), San Francisco, February 14–17.
14. *Solid-solution softening in bcc molybdenum from first principles.*
2004 MS&T Annual Meeting, New Orleans, LA, September 26–30.
13. *The lattice Green function for phonons: Decoupling short and long range contributions.*
2004 Electronic Structure Workshop, New Brunswick, NJ, May 27–30. (poster)

12. *Impurities block the alpha to omega martensitic transformation in titanium.*
2004 Annual Meeting of the Minerals, Metals & Materials Society (TMS), Charlotte, March 15–18.
11. *Inhomogeneous mechanism for the hcp to omega martensitic transformation in pure titanium.*
2003 March Meeting of the American Physical Society, Austin, March 3–7.
10. *A new mechanism for hcp-omega transformation in pure titanium.*
2003 Annual Meeting of the Minerals, Metals & Materials Society (TMS), San Diego, March 2–6.
9. *The mechanism of the titanium hcp to omega transformation.*
2002 Electronic Structure Workshop, Berkeley, June 5–8. (poster)
8. *Titanium hcp to omega transformation with oxygen impurities.*
2002 March Meeting of the American Physical Society, Indianapolis, March 18–22.
7. *Complete Titanium HCP to Omega Transformation Mechanism.*
2002 March Meeting of the American Physical Society, Indianapolis, March 18–22.
6. *Modeling the hcp to omega phase transition in titanium.*
2000 Psi-k Conference, Schwäbisch Gmünd, Germany, August 22–26. (poster)
5. *An improved tight-binding model for titanium.*
2000 March Meeting of the American Physical Society, Minneapolis, March 20–24.
4. *Modeling the hcp to omega phase transition in titanium.*
1999 Psi-k Workshop on Calculation of Material Properties, Trieste, Italy, August 9–18. (poster)
3. *Total energy and elastic properties of titanium via tight-binding.*
1999 Electronic Structure Workshop, Urbana-Champaign, May 21–24. (poster)
2. *Interstitials on Si {113} surface.*
1998 Electronic Structure Workshop, Philadelphia, May 29–31. (poster)
1. *Interstitials on Si {113} surface.*
1998 March Meeting of the American Physical Society, Los Angeles, March 23–27.

Current Advisees

Yang Dan (Ph.D., 2016–present)

Grace Lu (Ph.D., 2018–present)

Charith DeSilva (Ph.D., 2021–present)

Past Advisees

Min Yu (Ph.D., 2010; Coadvised with Prof. R. Martin, Physics) “Energy density method and its applications to defect energies.”

Maryam Ghazisaeidi (Ph.D., 2011) “A first principles study of defects in Titanium: Interaction of twin boundaries with dislocations and oxygen interstitials.”

Joseph A. Yasi (Ph.D., 2012) “Strength and ductility of Mg alloys from first-principles.”

Henry W. Wu (Ph.D., 2013) “Oxygen diffusion through titanium and other HCP metals.”

Pinchao Zhang (Ph.D., 2016) “Database optimization algorithm for empirical potentials.”

Zebo Li (Ph.D., 2017) “Multi-scale investigation of vacancy-mediated diffusion of Si in Ni near an edge dislocation.”

Emily Schiavone (Ph.D., 2017) “Ab initio modeling of hydrogen pipe diffusion in palladium.”

Ravi Agarwal (Ph.D., 2018) "Solute transport in magnesium and their uncertainty quantification using density functional theory and green function approach."

Anne Marie Tan (Ph.D., 2018) "Computation and application of the lattice Green function to dislocations in metals, intermetallics, and semiconductors."

Hyojung Kim (Ph.D., 2019) "Investigation of the mechanical properties and phase stability of Ti-transition metal alloys using first-principle calculations"

Abhinav Jain (Ph.D., 2019) "Atomic scale diffusion in complex systems from first principles."

Josh Vita (Ph.D., 2023) "Designing efficient, interpretable, and generalizable machine learning interatomic potentials"

Soham Chattopadhyay (Ph.D., 2023) "Methods for mass transport coefficient calculation in crystalline alloys used in irradiated environments"

Dr. Hadley Lawler (postdoc, 2009–2010)

Dr. Venkateswara Rao Manga (postdoc, 2010–2013)

Dr. Thomas Garnier (postdoc, 2013–2014)

Dr. Bora Lee (postdoc, 2013–2015)

Dr. Thomas Schuler (postdoc, 2015–2016)

Dr. Michael Fellingner (postdoc, 2013–2020)

Dr. Lucas Wirth (postdoc, 2021–present)

Chanda Lowrance (B.S., 2009–2010)

Pandu Wisesa (B.S., 2011–2012)

Arvind Srikanth (B.S., 2011–2013)

Ah-Young Song (M.S., 2011–2013)

Lauren Smith (B.S., 2016–2018)

Delma Nieves Rivera (REU, Univ. Puerto Rico, Summer 2014)

Ellango Jothimurugesan (REU, Duke, Summer 2014)

Cal Williams (REU, Univ. Roanoke, Summer 2015)

Randall Pittman (REU, Univ. Roanoke, Summer 2015)

Nitin Srirang (Intern, B.Tech. Eng. Phys. IIT Bombay, Summer 2018)