

Curriculum Vitae

Christopher Knight, Assistant Computational Scientist

Argonne National Laboratory

Work Address

Leadership Computing Facility
Argonne National Laboratory
Bldg. 240, Rm. 1132
9700 South Cass Avenue
Argonne, IL 60439

office: 630-252-9793
email: knightc@anl.gov

Appointments

<i>Position</i>	<i>Institution (Contact)</i>	<i>Dates</i>
Assistant Computational Scientist	Argonne LCF (Nichols Romero)	9/2013 - present
Joint Appointment	Computation Institute	5/2012 - present

Postdoctoral Appointments

<i>Institution</i>	<i>Supervisor</i>	<i>Dates</i>
Argonne National Laboratory	Gregory A. Voth	4/2010 - 9/2013
University of Chicago	Gregory A. Voth	4/2010 - 10/2010
University of Utah	Gregory A. Voth	4/2009 - 3/2010

Education

<i>Institution (Preceptor)</i>	<i>Degree</i>	<i>Date</i>	<i>Field</i>
The Ohio State University (Sherwin J. Singer)	Ph.D.	2009	Chemistry
Eastern Michigan University	B.S. w/ Honors	2003	Chemistry

Honors

ANL Computational Postdoctoral Fellowship	2011-2012
Ohio State University Presidential Fellowship	2008-2009
Undergraduate Honors Senior Thesis/Project and Symposium	2003
ACS Organic Chemistry Achievement Award	2002
Honors Undergraduate Fellowship	2002
Maurice Decoster Chemistry Scholarship	2002

Professional Societies and Activities

American Chemical Society
The Electrochemical Society
Referee for Journal of Physical Chemistry
Referee for Journal of Physical Chemistry Letters
Referee for Journal of Chemical Theory and Computation
Referee for Molecular Physics
Referee for Journal of Chemical Physics

Teaching Experience

ALCF Catalyst

In ALCF (9/13-present), I assist researchers with accomplishing their scientific goals on both new and mature projects by facilitating solutions to the unique challenges associated with efficient utilization of leadership-scale computational resources. This effort primarily involves porting scientific software to the IBM Blue Gene/Q architecture, teaching and assisting with profiling and debugging codes, discuss strategies and provide general guidance on code parallelization, I/O, load-balancing, workflow design, and data management.

Mentored undergraduate and high school students

Center for Multiscale Theory and Simulation, University of Chicago (6/12-7/12)

Teaching Assistant

Department of Chemistry, The Ohio State University (9/03-9/04, 9/05-9/06)

Research Interests

- Statistical mechanics
- Classical and quantum dynamics and multiscale processes in condensed phases
- Computer simulations via first principle methods
- Accurate and efficient simulation of chemical reactions in condensed phases
- Phase transitions in molecular solids
- High-performance scientific computing
- Chemistry at interfaces: aqueous systems, batteries and supercapacitors (electrolyte/electrode), fuel cells (proton and hydroxide conductance), biofuels (carboxysome), and smart materials (self-healing, piezoelectric, magnetostrictive), nanoporous materials
- Interactions of matter with soft and hard X-ray pulses

Publications: 397 citations; h-index 12; i10-index 13 (May 2015)

1. Hasan Metin Aktulga, Paul Coffman, Wei Jiang, Chris Knight, and Tzu-Ray Shan, "Implementation and Performance Optimization of the LAMMPS/ReaxC Package for Multi-core Architectures" (*in preparation*)
2. Revati Kumar, Chris Knight, Collin D. Wick, and Bin Chen, "Bringing Reactivity to the Aggregation-Volume-Bias Monte Carlo Based Simulation Framework: Water Nucleation Induced by a Reactive Proton", *J. Phys. Chem. B (ASAP)*
3. Peng Bai, Mi Yeong Jeon, Limin Ren, Chris Knight, Michael W. Deem, Michael Tsapatsis, and J. Ilja Siepmann, "Finding optimal zeolites for challenging separations and chemical conversions", *Nature Commun.*, **6**:5912 (2015)
4. Ying-Lung Steve Tse, Chris Knight, and Gregory A. Voth, "An Analysis of Hydrated Proton Diffusion in Ab Initio Molecular Dynamics", *J. Chem. Phys.*, **142**:014104 (2015)
5. Alistar Ottochian, Guilhem Dezanneau, Clement Gilles, Paolo Raiteri, Chris Knight, and Julian D. Gale, "Influence of isotropic and biaxial strain on proton conduction in Y-doped BaZrO₃: a reactive molecular dynamics study", *J. Mat. Chem. A*, **2**(9):3127-3133 (2014).
6. Revati Kumar, Chris Knight, and Gregory A. Voth, "Exploring the Behavior of the Hydrated Excess Proton at Hydrophobic Interfaces", *Farad. Discuss.*, **167**:263-278 (2013)
7. Takefumi Yamashita, Yuxing Peng, Chris Knight, and Gregory A. Voth, "Computationally Efficient Multiconfigurational Reactive Molecular Dynamics", *J. Chem. Theory Comput.*, **8**(12):4863-4875 (2012) (Berny Schlegel Festschrift)
8. Chris Knight, Gerrick E. Lindberg, and Gregory A. Voth, "Multiscale Reactive Molecular Dynamics", *J. Chem. Phys.*, **137**(22):22A525 (2012) (Nonadiabatic Dynamics special issue)
9. Chris Knight and Gregory A. Voth, "Coarse-graining Away Electronic Structure: A Rigorous Route to Accurate Condensed Phase Interaction Potentials", *Mol. Phys.*, **110**(9-10):935-944 (2012) (Bill Miller special issue).
10. Chris Knight and Gregory A. Voth, "The Curious Case of the Hydrated Proton", *Acc. Chem. Res.*, **45**(1):101-109 (2012) (Water in Chemistry special issue).
11. Sherwin J. Singer and Chris Knight, "Hydrogen bond topology and proton ordering in ice and water clusters", *Adv. Chem. Phys.*, **147**:1 (2012).
12. Hui Zhang, Ali A. Hassanali, Yun Kyung Shin, Chris Knight, and Sherwin J. Singer, "The water-amorphous silica interface: Analysis of the Stern layer and surface conduction", *J. Chem. Phys.*, **134**(2):024705 (2011).
13. Ali Hassanali, Hui Zhang, Chris Knight, Yun Kyung Shin, and Sherwin J. Singer, "The dissociated amorphous silica surface: Model development and evaluation", *J. Chem. Theory Comput.*, **6**(11):3456 (2010).
14. Chris Knight, C. Mark Maupin, Sergei Izvekov, and Gregory A. Voth, "Defining condensed phase reactive force fields from *ab initio* molecular dynamics simulations: The case of the hydrated excess proton", *J. Chem. Theory Comput.*, **6**(10):3223 (2010).

15. Chris Knight and Sherwin J. Singer, "Site disorder in ice VII arising from hydrogen bond fluctuations", *J. Phys. Chem. A*, **113**(45):12433 (2009) (Russell Pitzer Festschrift).
16. Chris Knight and Sherwin J. Singer, "Hydrogen bond ordering in ice V and the transition to ice XIII", *J. Chem. Phys.*, **129**(16):164513 (2008).
17. Chris Knight and Sherwin J. Singer, "A re-examination of the ice III/IX hydrogen bond ordering phase transition", *J. Chem. Phys.*, **125**(6):064506 (2006).
18. Chris Knight, Sherwin J. Singer, Jer-Lai Kuo, Tomas K. Hirsch, Lars Ojamäe and Michael L. Klein, "Hydrogen bond topology and the ice VII/VIII and Ih/XI proton ordering phase transitions", *Phys. Rev. E*, **73**(5):056113 (2006).
19. Chris Knight and Sherwin J. Singer, "Prediction of a phase transition to a hydrogen bond ordered form of ice VI", *J. Phys. Chem. B*, **109**(44):21040 (2005).
20. Sherwin J. Singer, Jer-Lai Kuo, Tomas K. Hirsch, Chris Knight, Lars Ojamäe and Michael L. Klein, "Hydrogen bond topology and the ice VII/VIII and Ih/XI proton ordering phase transitions", *Phys. Rev. Lett.*, **94**(13):135701 (2005).
21. Christopher Knight and M. C. Milletti, "Theoretical characterization of two reaction pathways for the intramolecular cyclization of 2-(3-benzylaminopropanoyl-amino)benzamide", *J. Mol. Struct.: THEOCHEM*, **724**:143 (2005).

Conference Publications

1. Preeti Malakar, Venkatram Vishwanath, Todd Munson, Christopher Knight, Mark Hereld, Sven Leyffer, Michael Papka, "Optimal Scheduling of Simulation-time Analysis for Large-scale Scientific Simulations", SC15 (submitted).
2. Nayong Kim, Richard Platania, Tom Keyes, Wei Huang, Chris Knight, Seun-jong Park, and Joohyun Kim, "Enabling Large-scale Biomolecular Conformation Search with Replica Exchange Statistical Temperature Molecular Dynamics (RESTMD) over HPC and Cloud Computing Resources" in *IEEE 29th International Conference on Advanced Information Networking and Applications Workshops (WAINA)*, 61-66 (2015).
3. Andrew M. Herring, Melissa A. Vandiver, Asley M. Maes, Himashu Sarode, E. Bryan Coughlin, Daniel M. Knauss, Yushan Yan, Gerrick E. Lindberg, Christopher Knight, Gregory A. Voth, Daniel Herbst, Thomas A. Witten, and Matthew W. Liberatore, "Fundamental Studies of Alkaline Exchange Membranes Towards Optimization in a Fuel Cell Environment", *ECS Trans.*, **50**(2): 2059-2066 (2012).
4. Gerrick E. Lindberg, Chris Knight, Lisa E. Feldberg, and Gregory A. Voth, "Molecular Dynamics Simulations of Hydroxide Solvation and Transport in Anionic Exchange Membranes", *ECS Trans.*, **50**(2): 2053-2058 (2012).
5. Yuxing Peng, Chris Knight, Philip Blood, Lonnie Crosby, Gregory A. Voth, "Extending Parallel Scalability of LAMMPS and Multiscale Reactive Molecular Simulations", XSEDE12, **37**:1-7 (ACM, New York, 2012).
6. Gerrick E. Lindberg, Chris Knight, Ryan Jorn, James F. Dama, and Gregory A. Voth, "Multiscale Simulation of Hydroxide Solvation and Transport in Anion Exchange Membranes", *ECS Trans.*, **41**(1):1785 (2011).

7. Himanshu Sarode, Melissa A. Vandiver, Ashley M. Maes, Benjamin Caire, James L. Horan, Yating Tan, Yifan Li, Gerrick E. Lindberg, James F. Dama, Chris Knight, Ryan Jorn, Martin E. Lenz, Robert Kasper, Shuang Gu, Bingzi Zhang, Sönke Seifert, Tsung-han Tsai, Wen X. Zhang, E. Bryan Coughlin, Daniel M. Knauss, Yushan Yan, Gregory A. Voth, Thomas A. Witten, Matthew W. Liberatore, and Andrew M. Herring, “Designing Alkaline Exchange Membranes from Scratch”, *ECS Trans.*, **41**(1):1761 (2011).
8. Chris Knight and Sherwin J. Singer, “Theoretical study of a hydroxide ion within the ice-Ih lattice”, *Physics and Chemistry of Ice* (Proceedings of the 11th International Conference on the Physics and Chemistry of Ice), ed., Werner F. Kuhs (Royal Soc. of Chemistry, 2007), p.339.
9. Chris Knight and Sherwin J. Singer, “Tackling the problem of hydrogen bond order and disorder”, *Physics and Chemistry of Ice* (Proceedings of the 11th International Conference on the Physics and Chemistry of Ice), ed., Werner F. Kuhs (Royal Soc. of Chemistry, 2007), p.329.

Technical Reports

1. Adrian W. Lange, Gard Nelson, Christopher Knight, and Gregory A. Voth, “Multiscale Molecular Simulations at the Petascale (Parallelization of Reactive Force Field Model for Blue Gene/Q)”, (*ALCF-2 Early Science Program Technical Report*), Argonne National Laboratory, 2013.

Presentations

1. Algorithms for computationally efficient molecular simulations: chemical reactivity, predictive modeling of zeolites, and x-ray damage mechanisms,” LANS Informal Seminar presented at Argonne National Laboratory, October 1, 2014.
2. Algorithms for computationally efficient molecular simulations: chemical reactivity, predictive modeling of zeolites, and x-ray damage mechanisms,” Seminar presented at the Argonne Leadership Computing Facility, Argonne National Laboratory, September 3, 2014.
3. ”Computationally efficient reactive models derived from condensed phase ab initio simulations,” Invited talk presented as the Technologies for Extreme Scale Computing seminar in the Center for Computation & Technology at Louisiana State University, July 17, 2014.
4. ”Computationally efficient reactive models derived from condensed phase ab initio simulations”, Seminar presented in the Chemistry Department at the University of Kansas, December 12, 2013.
5. ”Rigorous route to interaction potentials for reactive models derived from condensed phase ab initio simulations”, Talk presented at the 5th Annual Postdoctoral Research Symposium, Argonne National Laboratory, September 20, 2012.
6. ”Rigorous route to interaction potentials for reactive models derived from condensed phase ab initio simulations”, Invited talk presented at the 244th American Chemical Society National Meeting in symposium titled ”Potential energy surface exploration in quantum chemistry” honoring Berny Schlegel, Philadelphia, PA, August 19-23, 2012.
7. ”Extending Parallel Scalability of LAMMPS and Multiscale Reactive Simulations”, Talk presented at XSEDE12, Chicago, IL, July 16-20, 2012.
8. “Defining reactive forcefields from condensed phase ab initio simulations”, Poster presented at the Gordon Research Conference on Batteries, Ventura, CA, March 4-9, 2012.

9. "Defining reactive forcefields from condensed phase ab initio simulations", Poster presented at the American Conference on Theoretical Chemistry, Telluride, CO, July 17-22, 2011.
10. "Hydrogen bond topology, new interpretations of order/disorder transitions in ice, and the behavior of defects in a disordered ice lattice", Poster presented at 11th International Conference on the Physics and Chemistry of Ice, Bremerhaven, Germany, July 23-28, 2006.
11. "Hydrogen bond topology, new interpretations of order/disorder transitions in ice, and the behavior of defects in a disordered ice lattice", Talk presented at 38th Midwest Theoretical Chemistry Conference, Columbus, OH, June 15-17, 2006.
12. "Hydrogen bond topology, new interpretations of order/disorder transitions in ice, and the behavior of defects in a disordered ice lattice", Talk presented at Midwest Thermodynamics and Statistical Mechanics Conference, Akron, OH, May 25-26, 2006.
13. "Theoretical characterization of the potential energy surfaces of 2-(3-benzylamino-propanoylamino) benzamide", Poster presented at 225th American Chemical Society National Meeting, New Orleans, LA, March 23-27, 2003.
14. "Microwave synthesis and investigation of medium-sized ring systems containing nitrogen atoms", Poster presented at 225th American Chemical Society National Meeting, New Orleans, LA, March 23-27, 2003.

Workshops

1. Participation in the Nanoporous Materials Genome Center All-Hands meeting hosted by the University of Minnesota, October 18-19, 2014.
2. Participation as an ALCF Catalyst at the "Mira Performance Boot Camp" organized by the Argonne Leadership Computing Facility (ALCF), May 20-22, 2014, Argonne National Laboratory.
3. Invitation to participate in the "7th Bishop's Lodge Workshop: Materials for Energy Conversion" workshop organized by Plamen Atanassov (The University of New Mexico), November 3-5, 2013, Santa Fe, NM. Talk titled "Progress towards a fundamental understanding of charge transport mechanisms in fuel cell membranes".
4. Invitation to participate in the "ONETEP Master Class 2013" workshop organized by the ONETEP Developers group, August 27-30, 2013, Cambridge, UK.
5. Participation in the "Mira Performance Boot Camp" organized by the Argonne Leadership Computing Facility (ALCF), May 21-24, 2013, Argonne National Laboratory.
6. Talk titled "Multiscale Molecular Simulations at the Petascale" at the "Early Science Program Principle Investigators Meeting" organized by the Argonne Leadership Computing Facility (ALCF), May 15-16, 2013, Argonne National Laboratory.
7. Participation in the "Mira Community Conference" organized by the Argonne Leadership Computing Facility (ALCF), March 4-8, 2013, Argonne National Laboratory.
8. Invitation to participate in the "ONETEP Master Class 2012" workshop organized by the ONETEP Developers group, August 28-31, 2012, Cambridge, UK.

9. Invitation to participate in the "Leadership Computing Platforms, Extreme-scale Applications, and Performance Strategies" workshop organized by the Center for Scalable Application Development Software (CScADS), July 23-26, 2012, Snowbird, UT.
10. Participation in the "Early Science April Workshop - Code for Q" workshop organized by the Argonne Leadership Computing Facility (ALCF), April 30 - May 3, 2012, Argonne National Laboratory.
11. Participation in the "ALCF 2012 Winter Workshop" organized by the Argonne Leadership Computing Facility (ALCF), January 23-26, 2012, Argonne National Laboratory.
12. Talk titled "Multiscale Molecular Simulations at the Petascale" at the "ALCF Early Science Program Kick-Off Workshop" organized by the Argonne Leadership Computing Facility (ALCF), October 18-19, 2010, Argonne National Laboratory.