

David W.H. Swenson

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Education

- 2005 – 2011 **University of California, Berkeley**, *Doctor of Philosophy*
Field: Chemistry, Advisor: William H. Miller
Dissertation: “Quantum Effects from Classical Trajectories: New Methods and Applications for Semiclassical Dynamics.”
- 2003 – 2005 **Université Louis Pasteur**, *Diplôme d'Études Universitaires Générales*
Field: Mathématiques, Informatique, et Applications aux Sciences
Undergraduate degree in “mathematics, computer science, and applications to the sciences.”
- 1999 – 2003 **Colorado College**, *Bachelor of Arts*
Majors: Chemistry, French Literature, and Physics
- 1998 – 1999 **Indiana University-Purdue University, Indianapolis**
Upper-division classes in French taken concurrently with senior year in high school.

Publications

- 2017 Arthur C. Newton, Ramses Kools, David W.H. Swenson, and Peter G. Bolhuis. “The two opposing effects of isotropic and anisotropic attraction on the association kinetics of patchy particles.” *J. Chem. Phys.* **147**, 115101 (2017).
- 2014 David W.H. Swenson and Peter G. Bolhuis. “A replica exchange transition interface sampling method with multiple interface sets for investigating networks of rare events.” *J. Chem. Phys.* **141**, 044101 (2014).
- 2013 Bin Li, Tal J. Levy, David W.H. Swenson, Eran Rabani, and William H. Miller. “A Cartesian Quasi-classical Model to Nonequilibrium Quantum Transport: The Anderson Impurity Model.” *J. Chem. Phys.* **138**, 104110 (2013).
- 2012 David W.H. Swenson, Guy Cohen, and Eran Rabani. “A semiclassical model for the non-equilibrium quantum transport of a many-electron Hamiltonian coupled to phonons.” *Mol. Phys.* **110**, 743 (2012).
- 2011 David W.H. Swenson, Tal Levy, Guy Cohen, Eran Rabani, and William H. Miller. “Application of a semiclassical model for the second-quantized many-electron Hamiltonian to nonequilibrium quantum transport: The resonant level model.” *J. Chem. Phys.* **134**, 164103 (2011).
- 2006 David W.H. Swenson, Heather M. Jaeger, and Clifford E. Dykstra. “Clustering of molecular hydrogen around benzene.” *Chem. Phys.* **326**, 329 (2006).
- 2006 Heather M. Jaeger, David W.H. Swenson, and Clifford E. Dykstra. “Feature Article: Remarkable Features in the Interactions of Quadrupolar Molecules.” *J. Phys. Chem. A* **110**, 6399 (2006).

Publication in preparation

Jan-Hendrik Prinz, David W.H. Swenson, Peter G. Bolhuis, and John D. Chodera. “OpenPathSampling: A flexible, open framework for path sampling simulations. I. Basics.”

David W.H. Swenson, Jan-Hendrik Prinz, John Chodera, and Peter G. Bolhuis. “OpenPathSampling: A flexible, open framework for path sampling simulations. II. Advanced Topics.”

Jocelyne Vreede, Peter G. Bolhuis, and David W.H. Swenson. “Mechanism and rates for the Watson-Crick to Hoogsteen ‘base rolling’ transition in B-DNA.”

David W.H. Swenson, Martin Spruijt, Imke van Rees, and Peter G. Bolhuis. “Studying ‘rarer’ transitions using multiple interface set transition interface sampling.”

Oral Presentations

- 2017 David W.H. Swenson. “OpenPathSampling: A Python package for investigating biomolecular rare event simulations.” Automation in Biomolecular Simulation and Modeling: Amsterdam, Netherlands. September 12, 2017.
- 2017 David W.H. Swenson. “Scalability of path sampling simulations.” E-CAM Extreme Scale State-of-the-Art Workshop: Barcelona, Spain. July 6-7, 2017.
- 2016 Jocelyne Vreede, Peter G. Bolhuis, David W.H. Swenson. “Simulating the mechanisms and rates of transitions between Watson-Crick and Hoogsteen base pairing.” NWO CHAINS 2016: Veldhoven, Netherlands. December 8, 2016.
- 2016 David W.H. Swenson. “Tutorial: Software testing in scientific programming.” CECAM Extended Software Development Workshop: Traunkirchen, Austria. November 21, 2016.
- 2016 David W.H. Swenson. “Introduction to OpenPathSampling: Overview and Usage.” CECAM Extended Software Development Workshop: Traunkirchen, Austria. November 17, 2016.
- 2016 David W.H. Swenson, Jan-Hendrik Prinz, John D. Chodera, Peter G. Bolhuis. “Generation and Analysis of Arbitrary Path Ensembles using OpenPathSampling.” CECAM Workshop: Reaction Coordinates from Molecular Trajectories: Leiden, Netherlands. September 1, 2016.
- 2015 David W.H. Swenson. “Quantum dynamics from classical trajectory simulations.” Brown University: Providence, Rhode Island. December 8, 2015.
- 2015 David W.H. Swenson. “OpenPathSampling: An open, flexible Python framework for rare event simulations.” Dutch Molecular Dynamics Day 2015: Amsterdam, Netherlands. February 27, 2015.
- 2014 David W.H. Swenson and Peter G. Bolhuis. “Numerical study of correlated rare events using replica exchange multiple state transition interface sampling.” Eleventh International Conference on Monte Carlo and Quasi-Monte Carlo Methods in Scientific Computing.: Leuven, Belgium. April 6-11, 2014.

- 2014 David W.H. Swenson. "Sampling networks of rare events with multiple interface set transition interface sampling." Dutch Molecular Dynamics Day 2014: Enschede, Netherlands. March 7, 2014.
- 2013 David W.H. Swenson and Peter G. Bolhuis. "Numerical study of correlated rare events using replica exchange multiple state transition interface sampling." 246th American Chemical Society National Meeting: Indianapolis, Indiana. September 8, 2013.
- 2013 David W.H. Swenson. "Important initial phase space distributions in semiclassical dynamics and the nature of quantum coherences." 246th American Chemical Society National Meeting: Indianapolis, Indiana. September 8, 2013.
- 2012 David W.H. Swenson. "A Classical Model for Nonequilibrium Quantum Transport: The Dynamics for Classically Mapped Fermions Method." Amsterdam Center for Multiscale Modeling Symposium: Amsterdam, The Netherlands. June 28, 2012.
- 2012 David W.H. Swenson. "Classical Dynamics for Nonequilibrium Quantum Transport: The Dynamics for Classically Mapped Fermions Method." Dutch Molecular Dynamics Day: Groningen, The Netherlands. March 23, 2012.
- 2011 David W.H. Swenson, Tal Levy, Guy Cohen, Eran Rabani, and William H. Miller. "Semiclassical model for fermion dynamics (with applications to molecular electronics)." 242nd American Chemistry Society National Meeting: Denver, Colorado. August 28, 2011.

Poster Presentations

- 2017 David W.H. Swenson. "OpenPathSampling: A flexible Python framework for rare events." Recent Advances in Modeling Rare Events (RARE 2017): Agra, India. December 7-10, 2017.
- 2017 David W.H. Swenson. "Path sampling for unbinding kinetics." 3rd E-CAM General Assembly: Barcelona, Spain. November 29-30, 2017.
- 2014 David W.H. Swenson. "Efficient sampling of large rare event networks with multiple state replica exchange transition interface sampling." CECAM Workshop: Long time dynamics from short time simulations: Lugano, Switzerland. March 12-14, 2014.
- 2013 David W.H. Swenson. "Beyond basic molecular dynamics: New tools for the study of molecules in motion." 246th American Chemical Society National Meeting: Indianapolis, Indiana. September 9, 2013.
- 2011 David W.H. Swenson. "Transition Path Sampling Applied to Semiclassical IVRs." Mini Statistical Mechanics Meeting: Berkeley, California. January 14-16, 2011.
- 2010 David W.H. Swenson and William H. Miller. "Monodromy matrix calculation by the precision finite difference method." 239th American Chemical Society National Meeting: San Francisco, California. March 21-25, 2010.
- 2007 David W.H. Swenson and Cristian Predescu. "A Method for Solving Poisson Problems." Mini Statistical Mechanics Meeting: Berkeley, California. January 12-14, 2007.

- 2002 David W.H. Swenson, Laurent Bonnet, and Jean-Claude Rayez. “The State Distribution of the Associative Desorption of H₂ from a Pt(1 1 1) Surface.” 2002 Undergraduate Research Poster Session: Gainesville, Florida. October 25-27, 2002.
- 2002 David W.H. Swenson and Jeffrey A. Cina. “Exploring the Ground State Potential of I₂-Ar.” Inter-REU Workshop: 1st US/France Chemistry and Communication Meeting: Strasbourg, France. June 13-15, 2002.

Awards

- 1999 – 2003 **Otis A. and Margaret T. Barnes Scholarship**, *Colorado College*
A four-year, full-tuition scholarship for chemistry majors.
- 1999 – 2003 **Central Newspapers Foundation Scholarship**, *Central Newspapers Foundation*
A four-year, \$10,000 scholarship for children of employees of Central Newspapers, Inc.

Research Employment Experience

- 2012 – Present **Universiteit van Amsterdam**, *Post-Doctoral Fellow*
Theoretical Chemistry (Rare Events), Advisor: Peter G. Bolhuis
- Developed multiple interface set transition interface sampling to study networks of rare events.
 - Co-principal developer of OpenPathSampling, a software package to study rare events.
 - Developed a mathematical framework (path ensemble theory) underlying OpenPathSampling.
 - Planned and guided undergraduate research projects. Assisted graduate students with research.
 - Formed collaborations with other researchers to study systems of biological importance.
- Sep 2015 – **Memorial Sloan Kettering Cancer Center**, *Visiting Post-Doctoral Fellow*
Dec 2015 Computational Biology, Advisor: John D. Chodera
- Continued development of OpenPathSampling.
- Aug 2011 – **Tel Aviv University**, *Visiting Post-Doctoral Fellow*
Dec 2011 Theoretical Chemistry (Molecular Electronics), Advisor: Eran Rabani
- Extended my “dynamics for classically mapped fermions” approach to more complicated systems.
- 2005 – 2011 **University of California, Berkeley**, *Graduate Student Researcher*
Theoretical Chemistry (Semiclassical Dynamics), Advisor: William H. Miller
- Developed new methods for calculating the semiclassical prefactor.
 - Developed methods to reduce the number of trajectories for semiclassical calculations.
 - Created “dynamics for classically mapped fermions,” a semiclassical method for fermion dynamics.
 - Applied my “dynamics for classical mapped fermions” method to models of molecular conduction.
- Sept 2008 – **D. E. Shaw Research, LLC**, *Research Intern*
Jan 2009 Theoretical Chemistry/Computational Biology, Manager: John Klepeis
- Beta-tested a new parallel framework for trajectory analysis, based on the map-reduce approach used by Google.
 - Verified selfdiffusion, viscosity, and orientational correlation of a novel water model.

- Summers 2003 – 2005 **Indiana University-Purdue University Indianapolis**, *Summer Researcher*
Theoretical Chemistry, Advisor: Clifford E. Dykstra
- Developed a potential energy surface to describe molecular hydrogen clustering around benzene.
 - Applied my model to study clusters of up to 20 hydrogen molecules around a single benzene.
- Summer 2002 **Université Bordeaux I**, *REU Exchange Student*
Theoretical Chemistry, Advisors: Jean-Claude Rayez and Laurent Bonnet
- Began work on a quasiclassical trajectory simulation to determine product state distributions from the associative desorption of hydrogen from a Pt(1 1 1) surface.
- Summer 2001 **University of Oregon**, *REU Student*
Theoretical Chemistry, Advisor: Jeffrey A. Cina
- Studied the potential energy surface of the I₂-Ar system in preparation for theoretical studies of the pump-probe spectroscopy of I₂ in an argon matrix.
 - Explored the viability of using a computer algebra system as a tool in theoretical chemistry.
- Summer 2000 **Colorado College**, *Summer Researcher*
Natural Product Synthesis, Advisors: Nick Drapela and Ted Lindeman
- Steps toward the synthesis of candidate anti-tumor agent sequoiatone A.

Service and Volunteer Activities

- 2013–2015 **Group Outing Organization Committee**, *Molecular Simulation Group*
Helped plan activities for a group outing that had scientific value and was also fun for the participants. Designed several games based on useful scientific skills such as reading and presenting scientific literature, and explaining scientific concepts. Organized some logistical aspects of the outing.
- 2002–2003 **Student Representative to Department Meetings**, *Colorado College*
Elected by fellow students to represent their interests at Chemistry Department meetings. Organized student involvement in the hiring of a new professor.
- 2000–2003 **Kids' Science Day**, *Colorado College*
Program to introduce elementary school students to fun parts of science. Helped organize the first annual event, and served as counselor or chemistry/physics demonstrator every year.
- 1999–2003 **Student Affiliates of the American Chemical Society**, *Colorado College*
Served in several roles, including editor of the department newsletter, outreach co-ordinator, co-coordinator of National Chemistry Week activities, and executive vice president.

Teaching Experience

- May 2016 – Present **Master's Thesis Project**, *Universiteit van Amsterdam (Supervisor)*
Daily supervisor for the thesis project of a master's student.
- Apr–Jun 2014 **Bachelor's Thesis Project**, *Universiteit van Amsterdam (Supervisor)*
Proposed, supervised, and helped design a research project to serve as the thesis project for a Bachelor's student.

- Jun 2013 **Second-Year Student Project**, *Universiteit van Amsterdam (Supervisor)*
Designed, planned, and supervised a one-month research project for second-year undergraduates studying chemistry. Results from this project were presented in one of my talks at the 246th ACS National Meeting.
- Every January 2013 – 2015, **Molecular Simulation Tutorial**, *Universiteit van Amsterdam (Teaching Assistant)*
2017 Course for graduate students and post-docs in the field of computational chemistry. Guided several computer lab sessions.
- Fall 2010 **General Chemistry**, *University of California, Berkeley (Teaching Assistant)*
Instructors: John Arnold, Angela Stacy, Marcin Majda, and Michelle Douskey
- Fall 2007 **Advanced Quantum Mechanics**, *University of California, Berkeley (Teaching Assistant)*
Instructor: Daniel M. Neumark
Sole teaching assistant for a graduate-level course taken by most physical chemistry graduate students.
- Fall 2006 **Physical Chemistry**, *University of California, Berkeley (Teaching Assistant)*
Instructors: William H. Miller and Haw Yang
Undergraduate course to introduce chemistry majors to quantum mechanics.
- Fall 2005 **General Chemistry**, *University of California, Berkeley (Teaching Assistant)*
Instructors: Richard Saykally, Mark Kubinec, and Michelle Douskey
Introduction to chemical principles for non-majors.
- 2003-2004 **English Assistant**, *Lycée Le Corbusier (Illkirch, France)*
Taught English (as a foreign language) in a French high school focused on construction-related employment. My students ranged from housepainter apprentices to those who planned to become architects and designers.
- 1999-2003 **Chemistry and Physics Tutor**, *Colorado College*
Helped students in lower division chemistry and physics classes with homework and lab reports.