

ZACHARY STREETER

PERSONAL INFORMATION

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<i>Address</i>	2314 Bonar St. Berkeley, CA 94702
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<i>Familial Status</i>	Single, no children

BRIEF INTRODUCTION

A passionate, creative, and curious individual, with all qualities situated firmly on a foundation of work ethic. My passion leads me to problems that hopefully will create a better universal experience. My creativity leads me to problems that force me to mold or drop notions that I previously held. My curiosity allows me to listen and work with others that are not in my field. My work ethic allows my résumé to be read by people like you. Thank you for taking your time to read about me!

If you would like, please follow the red links above to email me, link up on LinkedIn, and/or check out my github page!

TECHNICAL SKILLS

Software

COMPILED · C, C++, Fortran, Cython.

PARALLEL API · MPI, OpenMP, Cuda, PETSC.

SCRIPTING · Posix, Bash, Python.

BUILD PROCESS · CMake, Make.

MARKUP · L^AT_EX, Markdown, ReStructuredText.

DEBUGGER · gdb, lldb, TotalView.

PROFILER · Nsight/visual profiler, VTune.

Workflow

EDITOR · Vim/Neovim.

MULTIPLEXER · Tmux.

VERSION-CONTROL · Git.

RESEARCH INTERESTS

Computer Science High Performance Computing

- Numerical algorithms/methods.
- Computational geometry.
- Hybrid CPU/GPU Architectures.

- HPC in coordination with Quantum Computing.

Theoretical Physics and Chemistry

- Quantum Information and Computation.
- Quantum Computers applied to quantum physics/chemistry.
- Nonlinear chemical reaction kinetics.
- Scattering Theory.
- Symplectic Mechanics.
- Underlying Symmetries throughout Physics.

INTERNSHIPS AND RESEARCH POSITIONS

<i>Summer</i> 2016	<i>to</i> <i>Present</i>	LBNL	Lawrence Berkeley National Laboratory	Created fully dimensional potential energy surfaces for H_2O^{++} using MOLPRO and Columbus Quantum Chemistry packages. These hypersurfaces were then used in a MPI parallelized classical trajectory simulation of H_2O^{++} breakup following double ionization. This work was essential to deduce the body-frame of the water molecule at the momentum of photo-absorption and resulted in two immediate papers while also providing a benchmark for intense field experimentalist that will be in print shortly. Created a novel suite of high-performance codes that calculate double-ionization cross section for water and can be easily modified to other polyatomics. In general, honed programming skills in C, C++, Fortran, and Python, while becoming a learned software developer devoted to best practices and good documentation. Used NERSC supercomputers EDISON and CORI, and also a cluster called Lawrencium, for running large (e.g. 40+ physical cores with 3000+ processors) parallel batch jobs. Became proficient in parallel programming using PETSC, MPI, CUDA, and OpenMP. Reference: Clyde W. McCurdy +1 (510) 486 4283 · cwmccurdy@lbl.gov
<i>Spring</i> 2015		BNL	Brookhaven National Laboratory	Performed experiments with soft X-rays utilizing the Linear Electron Accelerator Facility (LEAF) and the van de Graaff. Prepared samples in glove box and worked on purifying Xenon and CO. This work was essential in studying electron mobility through CO. Once this work was completed, we calculated the quasi-free electron energy resulting in a publication. Understanding the free-electron energy in various liquids is critical in order for those liquids to be used in scattering experiments. Reference: Richard Holroyd +1 (631) 344 4329 · holroydr@optonline.net
<i>Summer</i> 2014		CAMD	Center for Advanced Microstructures and Devices	Became a user in order to continue research from SRC. Reference: Cherice EVANS +1 (718) 997 4216 · cherice.evans@qc.cuny.edu
2012–2013		SRC	Synchrotron Radiation Center	Built gas handling systems, ran leak checks for high vacuum line, wrote Igor Pro code for data analysis, and worked on calibrating the monochrometer. Also attended lectures in relativistic electrodynamics and worked on electrodynamic problem sets. Reference: Gary FINDLEY +1 (318) 342 1835 · findley@ulm.edu

EDUCATION

2015-
present The University of California, Davis

Doctor of
Philosophy

GPA: 3.7 · School: Chemistry
Description: This degree is a PhD in Theoretical Chemical Physics.
Advisor: Prof. Clyde W. McCURDY

Fall 2019 The University of California, Berkeley

Notable Course

CS294 – 73 *Software Engineering for Scientific Computing*
School: Computer Science
Description: This graduate course focused on the seven motifs in scientific computing: dense and sparse linear algebra, structured and unstructured grid methods, particle methods, fast Fourier transforms (FFT), and Monte Carlo.
Professor: Phillip COLELLA · colella@eecs.berkeley.edu

Spring
2020 The University of California, Berkeley

Notable Course

CS267 *Applications of Parallel Computers*
School: Computer Science
Description: Graduate course focused on models for parallel programming. Overview of parallelism on scientific applications and study of parallel algorithms for linear algebra, particles, meshes, sorting, FFT, graphs, machine learning, etc. Survey of parallel machines and machine structures. Programming shared- and distributed-memory parallel computers, GPUs, and cloud platforms. Parallel programming languages, compilers, libraries and toolboxes. Data partitioning techniques. Techniques for synchronization and load balancing. Detailed study and algorithm/program development of medium sized applications.
Professor: Katherine A. YELICK · yelick@cs.berkeley.edu
Professor: James DEMMEL · demmel@cs.berkeley.edu
Professor: Aydin BULUÇ · aydin@eecs.berkeley.edu

2007-2009,
2011-2015 The University of Louisiana, Monroe

Bachelor of Science

GPA: 3.46 · School: School of Sciences
Major (Concentration): Biology (Chemical Biology)
Personal Courses: Attended formal lectures in Statistical Mechanics, Quantum Mechanics, Electricity and Magnetism, and Relativistic Electrodynamics.
Advisor: Prof. Gary FINDLEY & Prof. Ann FINDLEY

TEACHING

Spring
2020 University of California, Davis

Teaching Assistant

Time-Dependent Quantum Mechanics: The first part of this graduate course covers the basic concepts and techniques for solving the time-dependent Schrodinger equation. The initial portion explores the concepts of quantum superpositions, Gaussian wave packets for free and interacting particles, time propagation, the Schrodinger, interaction and Heisenberg representations, time-dependent density matrices, the Wigner phase space distribution, Ehrenfest's theorem, the connection between quantum and classical mechanics in the context of molecular dynamics, the semiclassical wave packet approximation, and time-dependent perturbation theory. The second part of the course turned to applications. Those included absorption and emission of electromagnetic radiation, correlation functions and spectra, molecular dynamics, potential energy surfaces, conical intersections, nonadiabatic transitions and variational transition state theory. .

Winter
2020 University of California, Davis

Teaching Assistant

Quantum Chemistry: a graduate level discussion of the principles of quantum mechanics and its application to (primarily) stationary state problems in atoms and

molecules, including Hartree-Fock calculations of their electronic structure. Using the Psi4 quantum chemistry codes and the Python programming language we performed calculations on small molecules using restricted Hartree-Fock, unrestricted Hartree-Fock, Møller-Plesset perturbation theory (MP2), and configuration interaction (CI) and coupled cluster (CCSD) methods..

2015-2016 University of California, Davis

Teaching Assistant

Taught freshman chemistry for two quarters. Also taught quantum mechanics for physical chemistry students.

Spring Queens College
2015

Teaching Assistant

Taught second semester of freshman chemistry and the corresponding lab. Created lab and recitation quizzes and was the sole arbiter as to how the courses were conducted .

Assisted Professor: Prof. Cherice EVANS

TALKS AND POSTERS PRESENTED AT CONFERENCES

- 2013 SRC Users Meeting Zachary Streeter, Kamil Krynski, C. M. Evans, and G. L. Findley, "Quasi-Free electron in near critical point hydrogen and deuterium," 2013 SRC Users Meeting, University of Wisconsin Synchrotron Radiation Center, Stoughton, WI, September 27 – 28, 2013.
- 2013 SRC Users Meeting Kamil Krynski, Zachary Streeter, C. M. Evans, and G. L. Findley, "Field ionization and photoionization of CH₃I perturbed by diatomic molecules: electron scattering in H₂, HD, D₂, O₂ and CO," 2013 SRC Users Meeting, University of Wisconsin Synchrotron Radiation Center, Stoughton, WI, September 27 – 28, 2013.
- 2014 DAMOP Cherice Evans, Kamil Krynski, Zachary Streeter, and G. L. Findley, "Field Ionization and Photoionization of CH₃I Perturbed by Diatomic Molecules: Electron Scattering in H₂, D₂, O₂, and CO," 45th Annual Meeting of the APS Division of Atomic, Molecular, and Optical Physics, Madison, WI, June 2 – 6, 2014.
- 2014 DAMOP Zachary Streeter, Kamil Krynski, C. M. Evans, and G. L. Findley, "The energy of the quasi-free electron in near critical point H₂, D₂, and O₂," 45th Annual Meeting of the APS Division of Atomic, Molecular, and Optical Physics, Madison, WI, June 2 – 6, 2014.
- 2016 APS Kamil Krynski, Zachary Streeter, C. M. Evans, and G. L. Findley, "Energy of the Quasi-Free Electron in H₂, D₂, and O₂: Probing Intermolecular Potentials within the Local Wigner-Seitz Model," American Physical Society March Meeting, Baltimore, MD, March 14 – 18, 2016.
- 2017 DAMOP Zachary Streeter, Frank Yip, Dylan P. Reedy, Allen Landers, C. William McCurdy, "Classical trajectory studies on the dynamics of one-photon double photionization of H₂O," 48th Annual Meeting of the APS Division of Atomic, Molecular, and Optical Physics, Sacramento, CA, June 5 – 9, 2017.
- 2018 ACS Cherice M. Evans, Jennifer Hare, Baxter Flor, Kamil Krynski, Zachary Streeter, and G. L. Findley, "Energy of the Quasi-Free Electron in CO and HD: Extension of the Local Wigner-Seitz Model to Polar Fluids," 225th ACS National Meeting and Exposition, New Orleans, LA, March 18 – 22, 2018.
- 2019 DAMOP Z. L. Streeter, and C. W. McCurdy, "Sequential dissociation of H₂O⁺⁺ following double photoionization" 50th Annual Meeting of the APS Division of Atomic, Molecular, and Optical Physics, Milwaukee, WI, May 27 – 31, 2019.

PUBLICATIONS

- Published C. M. Evans, Kamil Krynski, Zachary Streeter, and G. L. Findley, "Energy of the Quasi-free Electron in H₂, D₂ and O₂: Probing Intermolecular Potentials within the Local Wigner-Seitz Model," J. Chem. Phys. **143**, 224303 (2015)"

- Published* C. M. Evans, Baxter Flor, Kamil Krynski, Zachary Streeter, and G. L. Findley, "*Energy of the Quasi-Free Electron in CO and HD: Probing Intermolecular Potentials within the Local Wigner-Seitz model*," J. Chem. Phys. **149**, 064307 (2018).
- Published* Zachary L. Streeter, Frank L. Yip, Robert R. Lucchese, Benoit Gervais, and C. William McCurdy, "*Dissociation dynamics of the water dication following one-photon double ionization I: Theory*," Phys. Rev. A, **98**, 053429 (2018).
- Published* D. Reedy, J. B. Williams, B. Gaire, A. Gatton, M. Weller, A. Menssen, T. Bauer, K. Henrichs, Ph. Burzynski, B. Berry, Z. L. Streeter, J. Sartor, I. Ben-Itzhak, T. Jahnke, R. Dörner, Th. Weber, and A. L. Landers, "*Dissociation dynamics of the water dication following one-photon double ionization I: Experiment*," Phys. Rev. A, **98**, 053430 (2018).

September 14, 2020