

ZACHARY STREETER

PERSONAL INFORMATION

<i>email</i>	zacharylouis42@gmail.com
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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 in your position. 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>LBNL</i>	<p><i>Summer</i> Lawrence Berkeley National Laboratory <i>2016</i> to <i>Present</i></p> <p>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 parallel batch jobs (e.g. 40+ physical cores with 3000+ processors). Became proficient in parallel programming using PETSC, MPI, CUDA, and OpenMP. Reference: Clyde W. McCurdy +1 (510) 486 4283 · cwmccurdy@lbl.gov</p>
<i>BNL</i>	<p><i>Spring</i> Brookhaven National Laboratory, SULI internship <i>2015</i></p> <p>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</p>
<i>CAMD</i>	<p><i>Summer</i> Center for Advanced Microstructures and Devices <i>2014</i></p> <p>Became a user in order to continue research from SRC. Reference: Cherice EVANS +1 (718) 997 4216 · cherice.evans@qc.cuny.edu</p>
<i>SRC</i>	<p><i>2012–2013</i> Synchrotron Radiation Center</p> <p>Built gas handling systems, ran leak checks for high vacuum line, wrote Igor Pro code for data analysis, and worked on calibrating the monochromator. Also attended lectures in relativistic electrodynamics and worked on electrodynamic problem sets. Reference: Gary FINDLEY +1 (318) 342 1835 · findley@ulm.edu</p>

OPEN SOURCE PROJECTS

Spring 2020 **quantumGrid**

Author quantumGrid is a python package for solving a 1-D Schrödinger equation for an arbitrary potential on any interval. The heart of this package is using a Finite Element Method with a Discrete Variable Representation (FEM-DVR) grid to solve the time-dependent or time-independent Schrödinger equation. This grid provides a compact supported foundation for numerically accurate integration and also allows for a natural application of outgoing scattering boundary conditions by adding a complex tail as the last finite element of the FEM-DVR grid, called exterior complex scaling (ECS). This project was created for a graduate course in time-dependent quantum mechanics at UC Davis. .

EDUCATION

2015-present **The University of California, Davis**

Doctor of Philosophy GPA: 3.9 · School: Chemistry
Description: This degree is a PhD in Theoretical Chemical Physics.
Advisors: Prof. Clyde W. McCURDY, Prof. Robert. LUCCHESI (LBNL)

Fall 2019 **The University of California, Berkeley**

Notable Course CS294 – 73 *Software Engineering for Scientific Computing*
School: Computer Science
Grade: A+
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
Grade: A+
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

Teaching Assistant	Spring 2020	University of California, Davis
	Time-Dependent Quantum Mechanics: The first part of this graduate course covers the basic concepts and techniques for solving the time-dependent Schrödinger equation. The initial portion explores the concepts of quantum superpositions, Gaussian wave packets for free and interacting particles, time propagation, the Schrödinger, 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. .	
Teaching Assistant	Winter 2020	University of California, Davis
	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..	
Teaching Assistant	2015-2016	University of California, Davis
	Taught freshman chemistry for two quarters. My third quarter I taught quantum mechanics for physical chemistry students. This course laid the foundation for quantum mechanics needed later in spectroscopy courses.	
Teaching Assistant	Spring 2015	Queens College
	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 CH ₃ I perturbed by diatomic molecules: electron scattering in H ₂ , HD, D ₂ , O ₂ 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 CH ₃ I Perturbed by Diatomic Molecules: Electron Scattering in H ₂ , D ₂ , O ₂ , and CO," 45 th 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 H ₂ , D ₂ , and O ₂ ," 45 th 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 H ₂ , D ₂ , and O ₂ : 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 H_2O ," 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 H_2O^{++} 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 H_2 , D_2 and O_2 : 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).
- Published* Kirk A. Larsen, Thomas N. Rescigno, Travis Severt, Zachary L. Streeter, Wael Iskandar, Saijoscha Heck, Averell Gatton, Elio G. Champenois, Richard Strom, Bethany Jochim, Dylan Reedy, Dimitri Call, Robert Moshhammer, Reinhard Dörner, Allen L. Landers, Joshua B. Williams, C. William McCurdy, Robert R. Lucchese, Itzik Ben-Itzhak, Daniel S. Slaughter, Thorsten Weber, "Photoelectron and fragmentation dynamics of the $\text{H}^+ + \text{H}^+$ dissociative channel in NH_3 following direct single-photon double ionization," Phys. Rev. Res., **2**, 043056 (2020).
- Published* Kirk A. Larsen, Thomas N. Rescigno, Zachary L. Streeter, Wael Iskandar, Saijoscha Heck, Averell Gatton, Elio G. Champenois, Travis Severt, Richard Strom, Bethany Jochim, Dylan Reedy, Dimitri Call, Robert Moshhammer, Reinhard Dörner, Allen L. Landers, Joshua B. Williams, C. William McCurdy, Robert R. Lucchese, Itzik Ben-Itzhak, Daniel S. Slaughter, Thorsten Weber, "Photoionization and dissociation dynamics of the $\text{NH}_3^{2+} + \text{H}^+$ and $\text{NH}_3^{2+} + \text{H}^+ + \text{H}$ fragmentation channels upon single-photon double ionization of NH_3 at 61.5 eV" Journal of Physics B., **53**, 24 (2020).

January 9, 2021