

## David Williams–Young

### Academic Address

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### OBJECTIVE

To obtain a research position relating to high–performance scientific computing and numerical linear algebra.

### RESEARCH INTERESTS

- Development of high–performance and reduced scaling relativistic electronic structure methods.
- Algorithmic development for high–performance quaternionic linear algebra software.
- Application of abstract mathematical paradigms, such as algebraic topology, to develop elegant and novel solutions to the problems that arise in ab initio electronic structure theory.

### EDUCATION

*Doctor of Philosophy (Pursuant)*, Chemistry  
University of Washington, Seattle, WA  
Adviser: Dr. Xiaosong Li

(Projected) May 2018

*Bachelor of Science (Magna Cum Laude)*, Chemistry, Mathematics  
Indiana University of Pennsylvania, Indiana, PA  
Adviser: Dr. Jaeju Ko

May 2013

### PROFESSIONAL EXPERIENCE

*Graduate Research Assistant*  
University of Washington, Seattle, WA

July 2013 – Present

*Graduate Teaching Assistant*  
University of Washington, Seattle, WA

September 2013 – Present

- Provided supplementary lecture instruction for the undergraduate General Chemistry course series as well as full instruction for the associated lab coursework. This job entailed full responsibility for two sections of thirty students during a supplemental lecture series referred to as a quiz section. I was responsible for the evaluation of students' performance both in lecture coursework (exams, homework assignments) as well as the lab (lab reports).
- Provided supplementary lecture instruction for the final lecture course (in a series of three) of the undergraduate Organic Chemistry course series. Primarily, I was responsible for evaluating students' performance on Exams. In addition, I was responsible for making myself available for course related questions through structured office hours.
- Aided in the development of new course material for the second lecture course (in a series of two), of the graduate Quantum Chemistry course series. The developed course material focused on the practical development of basic quantum chemical methods, such as Hartree-Fock and Density Functional Theory, using the MATLAB development environment. The students developed, from scratch, a working implementation of these methods with the aid of provided course material developed by my self and

the primary course instructor. For this course, I also had the primary responsibility for evaluating students' progress through the grading of exams and assigned course work.

*Undergraduate Research Assistant*

September 2011 – May 2013

Indiana University of Pennsylvania, Indiana, PA

*Chemistry Tutor*

August 2010 – May 2011

Indiana University of Pennsylvania Disability Services, Indiana, PA

- Provided supplemental course instruction for the General and Organic Chemistry series for students with disabilities. This primarily entailed meeting with students individually on a weekly basis to aid in their understanding of course material and to ensure that they were not falling behind in coursework.

*Information Technology Technician*

August 2009 – January 2010

Central Michigan University Information Technology, Mt. Pleasant, MI

- Primarily responsible for solving network connectivity issues for incoming students at Central Michigan University.

## COMPUTATIONAL PROFICIENCY

I am very proficient in the following computational areas:

- *Programming Languages*: C/C++/C#, FORTRAN 77/95/03, Java
- *Scripting Languages*: Python, Julia, R, MATLAB, Octave, C/Bash Shell
- *Libraries / Paradigms*: OpenMP, TBB, MPI, OpenGL, CUDA, OpenACC, PThreads
- *Software*: Gaussian, Mathematica, MATLAB

I have contributed to the development of the following software packages:

- Chronus Quantum Chemistry (ChronusQ) Software Package (*Principle Developer*)
- Gaussian (*Contributor*)
  - Analytical Hessians of time-dependent density functional theory
  - Chebyshev expansion of the quantum propagator in real-time density functional theory

## PUBLICATIONS

**9 in press publications, 2 as first author**

\* Indicates equal contribution to published work

10. Kasper, J.; **Williams–Young, D. B.**; Vecharynski, E.; Yang, C.; Li, X.; “A Well-Tempered Hybrid Method of Davidson and GPLHR for Solving Challenging TDDFT Systems”; *J. Chem. Theor. Comp.*, **2018**, Submitted.
9. Lestrangé, P.; **Williams–Young, D. B.**; Jimenez–Hoyos, C.; Li, X.; “An Efficient Implementation of Variation After Projection Generalized Hartree–Fock” *J. Chem. Theor. Comp.*, **2018**, Just Accepted.
8. Barclay, M. S.; Quincy, T. J.; **Williams–Young, D. B.**; Caricato, M.; Elles, C. G.; “Accurate Assignments of Excited-State Resonance Raman Spectra: A Benchmark Study Combining Experiment and Theory” *J. Phys. Chem. A*, **2017**, 121(41), 7937–7946.
7. Van Beeuman, R.; **Williams–Young, D. B.**; Kasper, J.; Yang, C.; Ng, E. G.; Li, X.; “A Model Order Reduction Algorithm for Estimating the Absorption Spectrum” *J. Chem. Theor. Comp.*, **2017**, 13(10), 4950–4961.
6. Egidi, F.\*; **Williams–Young, D. B.\***; Baiardi, A.\*; Bloino, J.; Scalmani, G.; Frisch, M.; Li, X.; Barone, V.; “Effective Inclusion of Mechanical and Electrical Anharmonicity in Excited Electronic States: the VPT2-TDDFT Route” *J. Chem. Theor. Comp.*, **2017**, 13(6), 2789–2803.

5. Petrone, A.\*; **Williams–Young, D. B.\***; Lingerfelt, D. B.; Li, X.; “*Ab Initio Transient Raman Analysis*” *J. Phys. Chem. A.*, **2017**, 121(20), 3958–3965.
4. Petrone, A.; Lingerfelt, D. B.; **Williams–Young, D. B.**; Li, X.; “*Ab Initio Transient Vibrational Spectral Analysis*” *J. Phys. Chem. Lett.*, **2016**, 7, 4501–4508.
3. **Williams–Young, D. B.**; Goings, J.; Li, X.; “*Accelerating Real-Time Time-Dependent Density Functional Theory with a Non-Recursive Chebyshev Expansion of the Quantum Propagator*” *J. Chem. Theor. Comp.*, **2016**, 12(11) 5333–5338.
2. **Williams–Young, D. B.**; Egidi, F.; Li, X.; “*Relativistic Two-Component Particle-Particle Tamm–Dancoff Approximation*” *J. Chem. Theor. Comp.*, **2016**, 12(11), 5379–5384.
1. Lingerfelt, D. B.; **Williams–Young, D. B.**; Petrone, A.; Li, X.; “*Direct ab Initio (Meta-)Surface-Hopping Dynamics*”, *J. Chem. Theor. Comp.*, **2016**, 12(3), 935–945.

### CURRENT SOFTWARE CITATIONS

3. **Williams–Young, D. B.**; *HAXX: Hamilton’s Quaternion Algebra for CXX*, <http://github.com/wavefunction91/HAXX>, **2017**.
2. Li, X.; Valeev, E. F.; **Williams–Young, D. B.**; Ding, F.; Liu, H.; Goings, J. J.; Petrone, A.; Lestranger, P.; *Chronus Quantum, Beta Version*, <http://www.chronusquantum.org>, **2016**.
1. Frisch, M. J.; Trucks, G. W.; Schlegel, H. B.; Scuseria, G. E.; Robb, M. A.; Cheeseman, J. R.; Scalmani, G.; Barone, V.; Petersson, G. A.; Nakatsuji, H.; Li, X.; Caricato, M.; Marenich, A.; Bloino, J.; Janesko, B. G.; Gomperts, R.; Mennucci, B.; Hratchian, H. P.; Izmaylov, A. F.; Sonnenberg, J. L.; **Williams–Young, D. B.**; Ding, F.; Lipparini, F.; Egidi, F.; Goings, J.; Peng, B.; Petrone, A.; Ortiz, J. V.; Zakrzewski, V. G.; Gao, J.; Rega, N.; Zheng, G.; Liang, W.; Hada, M.; Ehara, M.; Toyota, K.; Fukuda, R.; Hasegawa, J.; Ishida, M.; Nakajima, T.; Honda, Y.; Kitao, O.; Nakai, H.; Vreven, T.; Throssell, K.; Montgomery Jr., J. A.; Peralta, J. E.; Ogliaro, F.; Bearpark, M.; Heyd, J. J.; Brothers, E.; Kudin, K. N.; Staroverov, V. N.; Keith, T.; Kobayashi, R.; Normand, J.; Raghavachari, K.; Rendell, A.; Burant, J. C.; Iyengar, S. S.; Tomasi, J.; Cossi, M.; Millam, J. M.; Klene, M.; Adamo, C.; Cammi, R.; Ochterski, J. W.; Martin, R. L.; Morokuma, K.; Farkas, O.; Foresman, J. B.; and Fox, D. J.; *Gaussian 16, A.03*, Gaussian, Inc., Wallingford CT, **2016**.

### HONORS

CCG Excellence Award for Graduate Students	The Chemical Computing Group (2017)
MolSSI Software Fellow	Molecular Sciences Software Institute (2017-2018)
Lloyd E. and Florence M. West Fellowship in Chemistry	Lloyd E. and Florence M. West (2016)
Early Bird Research Assistantship (EBRA)	University of Washington (2013)
Excellence in Chemistry Graduate Fellowship Award (ECGFA)	University of Washington (2013)
Provost Scholar	Indiana University of Pennsylvania (2013)

### PRESENTATIONS

\* Indicates an invited presentation

10. **Williams–Young, D. B.**; van Beeuman, R.; Yang, C.; Li, X.; “*A Novel Model Reduction Algorithm for the Efficient Evaluation of Molecular Response Properties*”; 254th American Chemical Society National Meeting & Exposition, Washington, D.C. **2017**. Poster.

9. \* **Williams–Young, D. B.**; “*Studying Semi-Classical Molecular Light–Matter Interaction through Time-Dependent Density Function Theory*”; High Performance Computing Seminar, University of Washington. **2017**. Oral Presentation.
8. **Williams–Young, D. B.**; Goings, J. J.; Li, X.; “*Accelerating Real-Time Time-Dependent Density Functional Theory with a Chebyshev Expansion of the Quantum Propagator*”; Theory and Applications of Computational Chemistry (TACC) 2016, Seattle, WA. **2016**. Poster.
7. Egidi F.; **Williams–Young, D. B.**; Li, X.; “*Electronic Structure Methods for Relativistic Effects in Excited States*”; Low Scaling and Unconventional Electronic Structure Theory (LUEST) 2016, Telluride, CO. **2016**. Poster.
6. **Williams–Young, D. B.**; Yang, W.; Li, X.; “*Moving past the particle-hole description of excited states: Affordable methodologies*”; Chemical Congress of Pacific Basin Societies (PacifiChem) 2015, Honolulu, HI. **2015**. PHYS: Recent Progress in Molecular Theory for Excited-state Electronic Structure and Dynamics. Oral Presentation.
5. **Williams–Young, D. B.**; Ko, J.; Ondrechen, M. J.; “*Computational approach to the prediction of enzyme specificities*”; 245th American Chemical Society National Meeting & Exposition, New Orleans, LA. **2013**. Sci-Mix Poster Session. Poster.
4. **Williams–Young, D. B.**; Ko, J.; Ondrechen, M. J.; “*Computational approach to the prediction of enzyme specificities*”; 245th American Chemical Society National Meeting & Exposition, New Orleans, LA. **2013**. Division of Computers in Chemistry Poster Session. Poster.
3. **Williams–Young, D. B.**; Ko, J.; “*Prediction of Relative Activities of Enzymes using Computed Chemical Properties*”; American Chemical Society Student Member Symposium, Duquesne University. Duquesne, PA. **2012**. Poster.
2. **Williams–Young, D. B.**; Ko, J.; “*Prediction of Relative Activities of Enzymes using Computed Chemical Properties*”; Undergraduate Research Forum, Indiana University of Pennsylvania. Indiana, PA. **2012**. Poster.
1. Ford, J.; Ko, J.; Mintmier, B.; Machovia, T.; Kang, M.; Owens, A.; **Williams–Young, D. B.**; “*Undergraduate Research in Biomass Utilization Symposium*”, Pennsylvania Association of the Council of Trustees Conference, Indiana University of Pennsylvania. Indiana, PA. **2011**. Oral Presentation.

## MEMBERSHIPS

Alpha Chi Sigma (AXΣ), GT Chapter, Professional Chemistry Fraternity  
American Chemical Society (ACS), Computers in Chemistry Division

## INTERESTS

World percussion, kayaking, hiking, climbing, Linux kernel development