Joshua J. Goings, Ph.D.

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Education
Ph.D.  University of Washington, Seattle, WA  Theoretical Chemistry  2017
B.S.  Seattle Pacific University, Seattle, WA  Chemistry & Biochemistry  2012

Summary
Experienced researcher at the interface of chemistry, quantum computing, and machine learning. Recent highlights:
- advancing chemistry simulations on fault-tolerant quantum computers (preprints #33 & #34)
- reinforcement learning for accurate wave function compression (publication #31)
- interpretable machine learning for molecular dynamics (publication #28)

Experience
Senior Applications Engineer  02/2022 to present
IonQ
Visiting Research Scientist  07/2021 to 02/2022
Google Quantum AI
Data Science Postdoctoral Research Fellow  03/2021 to 07/2021
eScience Institute, University of Washington
Postdoctoral Research Associate  07/2020 to 07/2021
University of Washington
Advisor: Prof. Xiaosong Li
Postdoctoral Research Associate  08/2017 to 06/2020
Yale University (lab moved Jan 2018)
University of Illinois at Urbana-Champaign
Advisor: Prof. Sharon Hammes-Schiffer
Graduate Research Assistant  07/2012 to 06/2017
NSF Graduate Research Fellow
University of Washington
Advisor: Prof. Xiaosong Li

Publications & Preprints
34  Reliably assessing the electronic structure of cytochrome P450 on today’s classical computers and tomorrow’s quantum computers
J. J. Goings, A. White, J. Lee, C. S. Tautermann, M. Degroote, C. Gidney, T. Shiozaki, R. Babbush, N. C. Rubin
arXiv preprint, arXiv:2202.01244.

33  Efficient quantum computation of molecular forces and other energy gradients
T. E. O’Brien, M. Streif, N. C. Rubin, R. Santagati, Y. Su, W. J. Huggins, J. J. Goings, N. Moll, E. Kyoseva, M. Degroote, C. S. Tautermann, J. Lee, D. W. Berry, N. Wiebe, R Babbush
arXiv preprint, arXiv:2111.12437.

32  Determination of the SmO+ bond energy by threshold photodissociation of the cryogenically cooled ion*
A. Lachowicz, E. H. Perez, N. S. Shuman, S. Ard, A. A. Viggiano, P. B. Armentrout, J. J. Goings, P. Sharma, X. Li, and M. A. Johnson
J. Chem. Phys., 2021, 155, 174303.
* 2021 JCP Editors’ Pick
Reinforcement Learning Configuration Interaction
J. J. Goings, H. Hu, C. Yang, and X. Li
*J. Chem. Theory Comput.*, 2021, 17, 5482–5491.

Role of Intact Hydrogen-Bond Networks in Multiproton-Coupled Electron Transfer
W. D. Guerra, E. Odella, M. Secor, J. J. Goings, M. N. Urrutia, B. L. Wadsworth, M. A. Gervaldo, L. E. Sereno, T. A. Moore, G. F. Moore, S. Hammes-Schiffer, A. L. Moore
*J. Amer. Chem. Soc.* 2020, 142, 21842–21851.

Formation of an Unusual Glutamine Tautomer in a Blue Light Using Flavin Photocycle Characterizes the Light-adapted State
J. J. Goings, P. Li, Q. Zhu, S. Hammes-Schiffer
*Proc. Natl. Acad. Sci.* 2020, 117, 26626–26632.

Nonequilibrium Dynamics of Proton-Coupled Electron Transfer in Proton Wires: Concerted but Asynchronous Mechanisms
J. J. Goings, S. Hammes-Schiffer
*ACS Cent. Sci.* 2020, 6, 1594–1601.

The Chronus Quantum Software Package
D. B. Williams-Young, A. Petrone, S. Sun, T. F. Stetina, P. Lestrange, C. E. Hoyer, D. R. Nascimento, L. Koulias, A. Wildman, J. Kasper, J. J. Goings, F. Ding, A. E. DePrince III, E. F. Valeev, X. Li
*WIREs Comput. Mol. Sci.* 2020, e1436.

Proton-Coupled Electron Transfer Across Benzimidazole Bridges in Bioinspired Proton Wires
E. Odella†, S. J. Mora†, B. L. Wadsworth†, J. J. Goings†, M. A. Gervaldo, L. E. Sereno, T. L. Groy, D. Gust, T. A. Moore, G. F. Moore, S. Hammes-Schiffer, A. L. Moore
*Chem. Sci.* 2020, 11, 3820–3828.
†Equal contribution.

Early Photocycle of Slr1694 Blue-Light Using Flavin Photoreceptor Unraveled through Adiabatic Excited-State Quantum Mechanical/Molecular Mechanical Dynamics
J. J. Goings, S. Hammes-Schiffer
*J. Amer. Chem. Soc.* 2019, 141, 20470–20479.

Proton-Coupled Electron Transfer Drives Long-Range Proton Translocation in Bioinspired Systems
E. Odella†, B. L. Wadsworth†, S. J. Mora†, J. J. Goings†, M. T. Huynh, D. Gust, T. A. Moore, G. F. Moore, S. Hammes-Schiffer, A. L. Moore
*J. Amer. Chem. Soc.* 2019, 141, 14057–14061.
†Equal contribution.

Electron-Coupled Double Proton Transfer in the Slr1694 BLUF Photoreceptor: A Multireference Electronic Structure Study
E. R. Sayfutyarova, J. J. Goings, S. Hammes-Schiffer
*J. Phys. Chem. B.* 2019, 123, 439–447.

Propensity for Proton Relay and Electrostatic Impact of Protein Reorganization in Slr1694 BLUF Photoreceptor
J. J. Goings, C. R. Reinhardt, S. Hammes-Schiffer
*J. Amer. Chem. Soc.* 2018, 140, 15241–15251.

Controlling Proton-Coupled Electron Transfer in Bio-Inspired Artificial Photosynthetic Relays
E. Odella, S. J. Mora, B. L. Wadsworth, M. T. Huynh, J. J. Goings, P. A. Liddell, T. L. Groy, M. Gervaldo, L. E. Sereno, D. Gust, T. A. Moore, G. F. Moore, S. Hammes-Schiffer, A. L. Moore
*J. Amer. Chem. Soc.* 2018, 140, 15450–15460.
19 Current Development of Noncollinear Electronic Structure Theory  
   J. J. Goings, F. Egidi, X. Li  
   *Int. J. Quantum Chem. 2018, 118, e25398.*

18 Real-Time Time-Dependent Electronic Structure Theory  
   J. J. Goings, P. J. Lestrange, X. Li  
   WIREs Comput. Mol. Sci. 2017, e1341.

17 Two-Component Non-Collinear Time-Dependent Spin Density Functional Theory for Excited State Calculations  
   F. Egidi, S. Sun, J. J. Goings, G. Scalmani, M. J. Frisch, X. Li  
   J. Chem. Theory Comput. 2017, 13, 2591–2603.

16 Can Quantized Vibrational Effects Be Obtained from Ehrenfest Mixed Quantum-Classical Dynamics?  
   J. J. Goings, D. B. Lingerfelt, X. Li  
   J. Phys. Chem. Lett. 2016, 7, 5193–5197.

15 Accelerating Real-Time Time-Dependent Density Functional Theory with a Non-Recursive Chebyshev Expansion of the Quantum Propagator  
   D. Williams-Young, J. J. Goings, X. Li  
   J. Chem. Theory Comput. 2016, 12, 5333–5338.

14 Quantum Confinement Effects on Optical Transitions in Nanodiamonds Containing Nitrogen Vacancies  
   A. Petrone, J. J. Goings, X. Li  
   Phys. Rev. B. 2016, 94, 165402.

13 Real Time Propagation of the Exact Two Component Time-Dependent Density Functional Theory  
   J. J. Goings, J. M. Kasper, F. Egidi, S. Sun, X. Li  
   J. Chem. Phys. 2016, 145, 104107.

12 A Direct Atomic-Orbital Based Relativistic Two-Component Linear Response Method for Calculating Excited State Fine Structures  
   F. Egidi, J. J. Goings, M. J. Frisch, X. Li  
   J. Chem. Theory Comput. 2016, 12, 3711–3718.

11 An Atomic Orbital Based Real-Time Time-Dependent Density Functional Theory for Computing Electronic Circular Dichroism Band Spectra  
   J. J. Goings, X. Li  
   J. Chem. Phys. 2016, 144, 234102.

10 Imaging Excited Orbitals of Quantum Dots: Experiment and Electronic Structure Theory  
   L. Nienhaus†, J. J. Goings†, D. Nguyen, S. Wieghold, J. Lyding, X. Li, M. Gruebele  
   J. Amer. Chem. Soc. 2015, 137, 14743–14750.  
   †Equal contribution.

9 Approximate Singly Excited States from a Two-Component Hartree-Fock Reference  
   J. J. Goings, F. Ding, E. R. Davidson, X. Li  
   J. Chem. Phys. 2015, 143, 144106.
8 Ab Initio Two-Component Ehrenfest Dynamics
F. Ding, J. J. Goings, H. Liu, D. Lingerfelt, X. Li
*J. Chem. Phys.* 2015, 143, 114105.

7 Energy-Specific Equation-of-Motion Coupled-Cluster Methods for High-Energy Excited States: Application to K-Edge X-ray Absorption Spectroscopy
B. Peng, P. J. Lestrange, J. J. Goings, M. Caricato, X. Li
*J. Chem. Theory Comput.* 2015, 11, 4146–4153.

6 Stability of the Complex Generalized Hartree-Fock Equations
J. J. Goings, F. Ding, M. J. Frisch, X. Li
*J. Chem. Phys.* 2015, 142, 154109.

5 Ab Initio Non-Relativistic Spin Dynamics
F. Ding, J. J. Goings, M. J. Frisch, X. Li
*J. Chem. Phys.* 2014, 141, 214111.

4 Theoretical Characterization of Conduction-Band Electrons in Photodoped and Aluminum-Doped Zinc Oxide (AZO) Quantum Dots
J. J. Goings, A. M. Schimpf, J. W. May, R. W. Johns, D. R. Gamelin, X. Li
*J. Phys. Chem. C.* 2014, 118, 26584–26590.

3 Assessment of Low-Scaling Approximations to the Equation of Motion Coupled-Cluster Singles and Doubles Equations
J. J. Goings, M. Caricato, M. Frisch, X. Li
*J. Chem. Phys.* 2014, 141, 164116.

2 Sorption of H\textsubscript{2} to Open Metal Sites in a Metal Organic Framework: A Symmetry Adapted Perturbation Analysis
J. J. Goings, S. Ohlsen, K. Blaisdell, D. Schofield
*J. Phys. Chem. A.* 2014, 118, 7411–7417.

1 Self-Consistent-Field Using Direct Inversion in Iterative Subspace Method and Quasi-Newton Vectors
J. J. Goings, F. Ding, X. Li
*Adv. Quantum Chem.* 2014, 68, 77–86.

### Awards

- University of Washington Data Science Postdoctoral Fellow 2021
- Chemical Computing Group (CCG) Excellence Award 2016
- National Science Foundation (NSF) Graduate Research Fellow 2014
- Bernard and Claudine Nist Endowed Fellowship in Chemistry 2014
- Peter Salamon Award, Telluride Science Research Center 2014
- Natt-Lingafelter Graduate Fellowship Award 2012
- David T. & Christina L. Wong Scholarship in Chemistry 2011

### Oral Presentations

- *Impact of protein conformational changes and electrostatics on a BLUF photoreceptor.*
  258\textsuperscript{th} ACS National Meeting, San Diego, CA (2019)

- *Engineering a redox-active proton wire with proton-coupled electron transfer.*
  258\textsuperscript{th} ACS National Meeting, San Diego, CA (2019)

- *Can quantized vibrational effects be obtained from Ehrenfest mixed quantum-classical dynamics?*
  253\textsuperscript{rd} ACS National Meeting, San Francisco, CA (2017)
• Atomic-orbital based real-time TDDFT for circular dichroism spectroscopy.
  251st ACS National Meeting, San Diego, CA (2016)

• Theoretical investigation of magnetic exchange interactions in dilute magnetic semiconductor quantum dots induced by defects.
  Pacificchem 2015, Honolulu, HI (2015)

• Linear Response Complex Generalized Hartree-Fock.
  249th ACS National Meeting, Denver, CO (2015)

• Equation of motion formalism of second order many-body perturbation theory (EOM-MBPT2) and second-order approximate coupled-cluster (CC2).
  248th ACS National Meeting, San Francisco, CA (2014)

• Low-Scaling Approximations to the Equation of Motion Coupled-Cluster Singles and Doubles Equations.
  Excited States and Time-Dependent Electronic Structure Theory, Telluride, CO (2014)

Poster Presentations
• Propensity for Proton Relay and Impact of Protein Reorganization in Slr1694 BLUF Photoreceptor.
  Gordon Research Conference on Computational Chemistry, West Dover, VT (2018)

• Real Time Propagation of the Exact Relativistic Two-Component Equations.
  Theory and Applications of Computational Chemistry (TACC), Seattle, WA (2016)

• Recent advances in real-time TDDFT for the description of optical activity.
  Low-scaling and Unconventional Electronic Structure Techniques Conference (LUEST), Telluride, CO (2016)

• Towards a Real-Time Description of Magnetic Systems with Applications to Magnetic Circular Dichroism Spectroscopy.
  251st ACS National Meeting, San Diego, CA (2016)

Teaching
Lecturer, Chemistry, University of Washington
CHEM 465/565: Computational Chemistry
• Lead instructor; re-designed course for asynchronous online learning.
• Managed team of TAs to facilitate hands-on student experience with modern computational chemistry software and UW’s high-performance computing (HPC) resources.
• Lectures highlighted fundamental and recent developments in computational chemistry.
• Guided students in novel research projects utilizing state-of-the-art methods and computing resources directly applicable to individual research interests.

Teaching Assistant, Chemistry, Yale University
CHEM 426/526: Computational Chemistry and Biochemistry
• Designed a computational chemistry course for graduate and undergraduate students.
• Gave students hands-on guidance using computational chemistry software.
• Facilitated student access to Yale’s high-performance computing center.
• Performed computational “demos” at the end of each lecture; guest lectured as necessary.