# Joseph James Radler

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## **EDUCATION**

#### **UNIV. OF WASHINGTON**

MASTER OF SCIENCE (MS)
June 2018 | Seattle, WA
Conc. Computational & Theoretical
Quantum Chemistry

#### **PURDUE UNIVERSITY**

BACHELOR OF SCIENCE (BS)
May 2012 | West Lafayette, IN
Major Chemistry
Minor Japanese Language

#### IVY TECH COMM. COLL.

ASSOCIATE OF APPLIED SCIENCE (AAS)

Dec. 2010 | Lafayette, IN Major Chemical Technology

## SKILLS

#### SCIENTIFIC COMPUTING

Numerical Modeling • Simulation Design

- Open Source Development
- Data Visualization Machine Learning
- Discrete Time-Frequency Analysis
- High-Performance Computing
- Data Management Automation

#### PROGRAMMING EXPERIENCE

5+ Years: Python • C • MATLAB • LETEX
3-5 Years: Bash Script • C++ • Markdown
1-3 Years: Haskell • SQL • HTML • Scala

#### **QUANTITATIVE**

- Discrete Mathematics Group Theory
- Differential Equations Linear Algebra
- Complex Analysis Numerical Analysis
- •Vector Calculus Statistics

#### **SOFTWARE TOOLS**

Workload Schedulers: PBS • Slurm Version Control Systems: Git • Subversion

#### **OPERATING SYSTEMS**

Proficient: Windows • MacOS Familiar: Linux (Fedora, Debian & Ubuntu)

# RELEVANT WORK & RESEARCH EXPERIENCE

## INDUCED POLARIZATION ASSOCIATES, LLC | SEATTLE, WA

MARINE INDUCED POLARIZATION DEVELOPMENT SR. RESEARCH SCIENTIST

May 2019 - Present

- Develop detection hardware for electrical waveforms in a noisy environment.
- Build and implement automated data processing and analysis tools in Python.
- Design applications for controlling and detecting data from field devices.
- Create and implement documentation workflows for testing.
- Implement a variety of digital signal processing software algorithms.
- Write reports for collaborators, investors, and funding bodies apprising them of our progress.
- Plan and construct data management tools such as code libraries, SQL database schemas, and version control workflows.

## UNIVERSITY OF WASHINGTON | SEATTLE, WA

RESEARCH ASSISTANT | XIAOSONG LI GROUP DEPARTMENT OF CHEMISTRY

Jan 2016 - Sept. 2018

- Wrote and published *three* (3) *original research articles* on simulating coupled quantum dynamics in peer-reviewed journals (pleaser see Publications section below).
- Collaborated with researchers from various experimental disciplines.
- Constructed a novel computational approach to treating fast intersystem crossings with Time-Dependent Density Functional Theory (TD-DFT).
- Developed high-performance computing pipelines, data analysis, and simulation software tools.
- Worked on a development team for an Open Source simulation application.
- Created visualizations for time-resolved frequency analysis.
- Automated parsing and analysis of large data sets from simulations.
- Presented my original work in both small-group settings and public seminars.
- Designed, typeset, and presented award-winning posters for conferences.

## **PUBLICATIONS**

- Valentine, A. J. S.; **Radler, J. J.**; Mills, A.; Kim, P.; Castellano, F. N.; Chen, L. X.; and Li, X.; *Resolving the ultrafast intersystem crossing in a bimetallic platinum complex.* J. Chem. Phys., **2019**. DOI: 10.1063/1.5115169
- Radler, J. J.; Lingerfelt, D. B.; Castellano, F. N.; Chen, L. X.; Li, X.; Role of Vibrational Dynamics on Excited-State Electronic Coherence in a Binuclear Platinum Complex (Featured Cover Article) J. Phys. Chem. A, 2018. DOI: 10.1021/acs.jpca.8b01352
- Lingerfelt, D. B.; Lestrange, P. J.; Radler, J. J.; Brown-Xu, S. E.; Kim, P.; Castellano, F. N.; Chen, L. X.; Li, X.; Can Excited State Electronic Coherence Be Tuned via Molecular Structural Modification? A First-Principles Quantum Electronic Dynamics Study of Pyrazolate-Bridged Pt(II) Dimers. J. Phys. Chem. A., 2017. DOI: 10.1021/acs.jpca.6b12099

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## RELEVANT COURSES

#### **CHEMISTRY & PHYSICS**

- Computational Molecular Dynamics
- Quantum Computational Chemistry
- Group Theory for Physical Chemistry
- Advanced Inorganic Chemistry
- Computational Chemistry
- Quantum Chemistry I
- Quantum Chemistry II

### **MATHEMATICS**

- Numerical Linear Algebra
- Computational Data Analysis
- Ordinary Differential Equations
- Multivariable & Vector Calculus

#### **COMPUTER PROGRAMMING**

- High-Performance Computing
- C Programming
- Intro Engineering with MATLAB
- Developing Quantum Models

#### WRITING AND LANGUAGE

- Grant and Proposal Writing
- •Technical Writing and Editing
- English Composition
- Japanese Language & Literature

## OTHER SKILLS

### **NATURAL LANGUAGES**

Native or Fluent:

English • Japanese

Conversational:

French • Portuguese • Russian • Spanish Basic:

German • Mandarin

#### COMMUNICATION

Technical Writing • Editing
Visualizations • Developing Tutorials
Publishing • Poster & Oral Presentation

# LINKS

Github://jjradler LinkedIn://jjradler Webpage://jjradler.github.io

# SELECTED POSTER PRESENTATIONS

- Radler, J. J.; Lingerfelt, D. B.; Li, X.; Cascading into Coherence Theoretical Investigations of Long-Lived Excited State Coherences in Bimetallic Pt(II) Complexes.
  Conference on Excited State Processes (ESP 2018); 06/2018.
- Radler, J. J.; Lingerfelt, D. B.; Li, X.; Exploring the Role of Nuclear Motion on Excited State Coherences in Binuclear Pyrazolate-Bridged Platinum Complexes. 1st Northwest Theoretical and Computational Chemistry Conference; 10/2017.
- Houferak, C.; Kasper, J.; Radler, J. J.; Sun, S.; Applications of Compressive Sensing to Simulated Chemical Spectra. University of Washington Engineering in Data Science Symposium; 03/2016
- Radler, J. J.; Jackson, G. S.; Koopman, H.; Westgate, A.; Determination of <sup>14</sup>C Pelagic Ocean Values through Atomic Bomb Radiocarbon Dating of Dolphin Teeth. 13<sup>th</sup> Accelerator Mass Spectrometry Conference; **08**/2014.
- Radler, J. J.; Schauer, D. J.; Probing the Nature of Metal-Phosphonate Interactions by FTIR Spectroscopy, Central Regional Meeting of the ACS; 06/2010.

## OTHER WORK EXPERIENCE

## UNIVERSITY OF WASHINGTON | SEATTLE, WA

GRADUATE TEACHING ASSISTANT / ASSOCIATE | GENERAL CHEMISTRY DEPARTMENT OF CHEMISTRY

Sept. 2015 - Mar. 2019

- Taught students the theoretical and technical aspects of chemistry.
- Conveyed abstract concepts from physics in introductory-level chemistry courses.
- Instructed First Year Graduate students at a Teaching Assistant Orientation Seminar.

## PURDUE RARE ISOTOPE MEASUREMENT (PRIME) LABORATORY

Chemist | Geological Sample Processing Chemistry Operations May 2012 - Aug 2015 | West Lafayette, IN

- Processed soil, rock, and water samples for <sup>36</sup>Cl analysis by Accelerator Mass Spectrometry.
- Optimized sample processing, tripling sample throughput.
- Wrote and revised sample processing documentation and procedures.
- Collaborated remotely with researchers to advise them with their own processing facilities.

## **AWARDS**

- 2017 J. Phys. Chem. A Award for Best Original Research Poster (1st NWTCC)
- 2012 Dean's List, Purdue University

# **SOCIETIES**

- 2016 Current Society of Industrial and Applied Mathematicians (SIAM)
- 2015 Current American Chemical Society (ACS)
- 2016 UW High-Performance Computing Club (UW-HPCC)
- 2006 Alpha Chi Sigma Beta Nu Chapter