# Duy-Khoi Dang

San Jose, CA | khoidang@umich.edu | khoidang.com github.com/duykhoidang | ORCID: 0000-0001-7530-0540

## **Education**

University of Michigan, Ann Arbor, MI Ph.D. in Chemistry and Scientific Computing	Sep 2018 - Apr 2023 GPA: 3.9
Northeastern University, Boston, MA Bachelor of Science in Chemistry and Mathematics	Sep 2014 - May 2018 GPA: 4.0
Awards and Honors	
University of Michigan, Rackham One Term Fellowship	Sep 2022 - Dec 2022
NSF XSEDE, Startup Allocation on SDSC Expanse	May 2022 - May 2023
Michigan Institute for Computation Discovery and Engineering, Fellow	Sep 2018 - Apr 2023
National Science Foundation, Graduate Research Fellowship	Sep 2018 - Apr 2023
Northeastern University, President's Award	Mar 2018

#### Skills

**Areas of Expertise:** quantum chemistry, parallel computing, high performance computing **Proficiencies:** Python, C++, MPI+OpenMP, OpenACC, CUDA, CMake, bash, Linux, git, SLURM, PyTorch,

Proficiencies: Python, C++, MPI+OpenMP, OpenACC, CUDA, CMake, bash, Linux, git, SLURM, PyTorch, TensorFlow (python), dask, pandas, NumPy, SymPy, Rust, Q-Chem, PySCF, Molpro, OpenMolcas, Mathematica

# **Experience**

#### **TSMC**

Principal Engineer, San Jose, CA

Northeastern University, Dean's Scholarship

Jul 2023 - Present

Aug 2017

Sep 2014 - May 2018

- Contributed to and maintained scientific software library for lithography simulations using CUDA and C++
- Developed in-house machine-learning inferencing engines in CUDA and C++ to obtain speedup over TensorFlow
- Developed high-performance polygon algorithms and engines for CPU and GPU

Division of Inorganic Chemistry of the ACS, Undergraduate Award

#### **Zimmerman Group**

Graduate Student Research Assistant, Ann Arbor, MI

May 2019 - Apr 2023

- Developed scalable heat-bath CI library in C++ with MPI+OpenMP in Q-Chem software package
- Devloped incremental CASSCF methods for recovering strong correlation in Q-Chem software package
- Developed GPU accelerated library for computing Slater integrals with C++ and OpenACC
- Collaborated with experimental groups to study high-spin organic organic complexes with significant radicaloid character
- Collaborated with experimental groups to study photoisomerization processes using DFT and CASSCF

## University of Michigan

Graduate Student Instructor, Ann Arbor, MI

Sep 2019 - Apr 2020

• Led discussions for CHEM 260, an introductory course in Physical Chemistry

#### Geva Group

Rotation Student, Ann Arbor, MI

Jan 2019 - May 2019

• Studied applications of Redfield equation in quantum dynamics simulations and spectroscopy

#### **Zimmerman Group**

Rotation Student, Ann Arbor, MI

Sep 2018 - Dec 2018

• Developed reaction path-finding methods for systematically searching for spin crossings

## Lopez Lab

Undergraduate Research Assistant, Boston, MA

Sep 2017 - May 2018

• Studied photochemical systems, specifically spiropyrans and diazirines, using DFT and CASSCF in the Gaussian software package

# **Northeastern University**

Grader, Boston, MA

Sep 2017 - May 2018

• Graded exams and problem sets for Physical Chemistry

### Pfizer Inc.

Co-op Student, Cambridge, MA

Jul 2016 - Dec 2016

• Studied allosteric networks in G-protein coupled receptors with molecular dynamics, docking, and statistical methods using using the Schödinger, MOE, and Amber software packages

#### Kirss Lab

Undergraduate Research Assistant, Boston, MA

Dec 2015 - May 2018

• Studied phosphine exchange and ligand isomerization in Cp(PPh)<sub>3</sub>RuX complexes using DFT in the Gaussian software package

# **Northeastern University**

Peer Tutor, Boston, MA

Sep 2015 - May 2018

• Tutored students in Organic and Physical chemistry course series

## **Boeing Co.**

Intern - Student Engineer, Huntsville, AL

May 2015 - Aug 2015

• Investigated next-generation anodizing and electroplating solutions to replace incumbent methods utilizing toxic substances

# **Boeing Co.**

Intern - Student Engineer, Huntington Beach, CA

Jun 2014 - Aug 2014

• Studied CO<sub>2</sub> adsorption properties of zeolites

# **Community Involvement**

# American Chemical Society Graduate and Postdoctoral Scholars Office

Graduate Student Symposium Planning Committee Coordinator

Apr 2022 - Sep 2024

- Acted as a liaison between the American Chemical Society's Graduate Student and Postdoctoral Scholar Advisory Board and active Graduate Student Symposium Planning Committees
- Provided logistical support for Graduate Student Symposium Planning Committees at American Chemical Chemical Society National Meetings

### University of Michigan Graduate Student Symposium Planning Committee

American Chemical Society Correspondent, Ann Arbor, MI

Nov 2018 - Nov 2020

• Planned the Graduate Student Symposium for the American Chemical Society National Meeting in Philadelphia 2020

## Northeastern University Student Affiliates of the ACS

President, Boston, MA

May 2017 - May 2018

- Acted as a bridge between undergraduate chemistry students and the chemistry department
- Hosted weekly speakers to discuss current research and future career opportunities for chemistry graduates

## Northeastern University College of Science Advisory Council

Chemistry Representative, Boston, MA

Sep 2015 - May 2018

• Advised the deans of the College of Science on important issues of the undergraduate student body

### Northeastern University Student Affiliates of the ACS

Secretary, Boston, MA

May 2016 - May 2017

- Wrote a semi-weekly newsletter about upcoming events and opportunities to the chemistry undergraduates
- Invited professionals in chemistry to discuss relevant research and career opportunities to participate in the club's weekly speaker series

## **Beyond Benign**

Green Chemistry Outreach Fellow, Wilmington, MA

Oct 2015 - May 2018

• Participated in outreach events throughout the Boston Area to promote the principles of green chemistry and engage young students in science

# **Publications**

- 10. **D.-K. Dang**, J.D. Einkauf, X. Ma, R. Custelcean, Y.-Z. Ma, P.M. Zimmerman, V.S. Bryantsev, Photoisomerization mechanism of iminoguanidinium receptors from spectroscopic methods and quantum chemical calculations, *Phys. Chem. Phys.*, 2024, **26**, 24008-24020.
- 9. B. Prajapati, M. Ambhore, **D.-K. Dang**, P. Chmielewski, T. Lis, C. Gómez-García, P.M. Zimmerman, M. Stępień, Tetrafluorenofulvalene: A Sterically Frustrated Open-Shell Alkene, *Nat. Chem.*, 2023, **15**, 1541.
- 8. **D.-K. Dang**, D. Cehreli, B.S. Rich, T.E. Haas, F.R.Fronczek, and R.U. Kirss, Phosphine substitution and linkage isomerization in cyclopentadienylruthenium bis(triphenylphosphine)thiocyanide and selenocyanide, CpRu(PPh<sub>3</sub>)<sub>2</sub>NCS and CpRu(PPh<sub>3</sub>)<sub>2</sub>SeCN, *Dalton Trans.*, 2023, **52**, 13258.
- 7. S. Tribedi, **D.-K. Dang**, B. Kanungo, V. Gavini, P.M. Zimmerman, Exchange Correlation Potentials from Accurate FCI Densities Constructed from Slater Basis Functions, *J. Chem. Phys.* 2023, **159**, 054106.
- 6. **D.-K. Dang**, J. A. Kammeraad, P. M. Zimmerman, Advances in Parallel Heat Bath Configuration Interaction, *J. Phys. Chem. A*, 2023, **127**(1), 400.
- 5. **D.-K. Dang**, L. W. Wilson, P. M. Zimmerman, The numerical evaluation of Slater integrals on graphics processing units, *J. Comput. Chem.*, 2022, **43**(25), 1680.
- 4. B. Prajapati, **D.-K. Dang**, P. J. Chmielewski, M. A. Majewski, T. Lis, C. J. Gómez-García, P. M. Zimmerman, M. Stępień, An Open-Shell Coronoid with Hybrid Chichibabin–Schlenk Conjugation, *Angew. Chem. Int. Ed.* 2021, **60**, 22496.
- 3. D.-K. Dang and P. M. Zimmerman, Fully Variational Incremental CASSCF, J. Chem. Phys., 2021, 154, 014105
- 2. V. M. Breslin, N. A. Barbour, **D.-K. Dang**, S. A. Lopez, and M. A. Garcia-Garibay, Nanosecond laser flash photolysis of a 6-nitroindolinospiropyran in solution and in nanocrystalline suspension under single excitation conditions, *Photochem. Photobiol. Sci.*, 2018, **17**, 741
- 1. D. Hill, C. Delaney, M. Clark, M. Eaton, B. Hassan, O. Hendricks, **D.-K. Dang**, and R.U. Kirss, Kinetics of phosphine substitution in  $CpRu(PPh_3)_2X$  (X = Cl, Br, I, N<sub>3</sub>, and NCO), RSC Adv., 2017, 7, 34425

## **Presentations**

- 6. **D.-K. Dang**, P. M. Zimmerman, Advances in Parallel Heat-Bath Configuration Interaction, Michigan Institute for Computational Discovery and Engineering Symposium, March 24, (2023)
- 5. **D.-K. Dang**, L. W. Wilson, P. M. Zimmerman, GPU-accelerated numerical integration of electron repulsion integrals in a Slater basis, ACS Spring National Meeting & Events, San Diego, CA, United States, March 20-24, (2022). COMP
- 4. **D.-K. Dang**, P. M. Zimmerman, Treating Large CASSCF Active Spaces with the Method of Increments, 33rd Annual Workshop on Recent Developments in Electronic Structure Methods (ES21), (2021)
- 3. **D.-K. Dang**, P. M. Zimmerman, Variational procedure for the many-body expansion of the CASSCF energy, 259th ACS National Meeting & Exposition, Philadelphia, PA, United States, March 22-26, (2020), COMP-0591
- 2. M. C. Enright, A. Giarross, J. Romeo, **D.-K. Dang**, J. Davis, C. Olbrich. Empowering undergraduates to be green chemistry ambassadors in their community through outreach, 21st Annual Green Chemistry & Engineering Conference, Reston, VA, United States, June 13-15 (2017), GC+E-239
- 1. **D.-K. Dang**, and R. U. Kirss. Linkage isomerization in phosphine substitution reactions of CpRu(PPh<sub>3</sub>)<sub>2</sub>NCS, 253rd ACS National Meeting & Exposition, San Francisco, CA, United States, April 2-6, (2017), COMP-323.