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education

**Ph.D. Chemical Engineering,** Brigham Young University, Provo, UT 2017

* Dissertation: How a Systematic Approach to Uncertainty Quantification Renders Molecular

Simulation a Quantitative Tool in Predicting the Critical Constants for Large *n*-Alkanes

* Expertise: Force Field Development, Computational Chemistry, Configuration Reweighting,

Uncertainties in Molecular Simulation, Thermodynamic Data Analysis

* Elective Courses: Quantum Chemistry, Statistical Mechanics, Nonlinear Statistical Analysis,

Polymer Science and Engineering, Advanced Organic Chemistry,

Classical Mechanics, Instrumental Analysis Lecture/Lab GPA: 4.0

**B.S. Chemical Engineering,** Brigham Young University, Provo, UT 2012

* Elective Courses: Molecular Modeling, Introduction to Partial Differential Equations
* Excelled in: Thermodynamics, Physical Chemistry, Reaction Engineering,

Separations, Process Control, Statistics

* Minors: Spanish, French Overall GPA: 3.78

Work experience

**NRC Postdoc Associate,** Thermodynamics Research Center, Boulder, CO Feb. 2017- 2019

* Implemented alchemical free energy methods to accelerate optimizing force field parameters
* Collaborated with researchers at the National Institute of Standards and Technology, University of Colorado, University of Akron, and Wayne State University
* Mentored undergraduate student during three-month project for the 10th Industrial Fluid Properties Simulation Challenge

**Research Assistant,** Design Institute for Physical Properties, Provo, UT Jan. 2012-Feb. 2014

* Performed experimental work that involved: preparation, execution, cleaning, and processing data
* Evaluated literature experimental data and property prediction models for two biofuels

**Teaching Assistant,** BYU Chemical Engineering, Provo, UT

* Courses: Chemical Process Principles, Dr. Thomas H. Fletcher Jan.-Apr. 2012

Plant Design & Synthesis, Dr. W. Vincent Wilding Jan.-Apr. 2013/2014

Molecular Modeling, Dr. Thomas A. Knotts IV Jan.-Apr. 2015

* Conducted exam reviews, held office hours, and graded homework assignments

Volunteer work

**Church Representative,** The Church of Jesus Christ of Latter-day Saint, Guatemala Nov. 2006-2008

* Led a regional group of 12 representatives

**Boy Scout Leader,** Boy Scouts of America 1999-2006

* Inspired younger scouts to achieve their Eagle while organizing campouts and teaching activities

SKILLS/AWARDS

* **Trained in:** C++, VBA, MATLAB, Python, Towhee, Cassandra, Gromacs, GOMC
* **Spanish, French, Portuguese -** proficient reading, writing, and speaking
* **Dean’s List Student -** achieved a 4.0 semester GPA as undergraduate Apr. 2009 & Jun. 2010
* **Eagle Scout Award -** erected a flag pole in front of a religious center Sept. 11th, 2002

Presentations

**Uncertainty Quantification of Non-bonded Potentials for Prediction of Thermophysical Properties with Molecular Simulation,** *20th Symposium on Thermophysical Properties*, Boulder Colorado, 2018.

**Improved Prediction of High Pressure Viscosities with Molecular Simulation,** *Thermodynamics Research Center Consortium,* Boulder, Colorado, 2018.

**Accelerating Force Field Parameterization to Improve the Quantitative Predictability of Thermophysical Properties,** *American Institute of Chemical Engineers Annual Meeting,* Minneapolis, Minnesota, 2017.

**Pushing the Frontier of Data-Driven Force Field Development,** *Thermodynamics Research Center Consortium,* Boulder, Colorado, 2017.

**A Novel Force Field Development Algorithm to Improve the Quantitative Predictability of Thermophysical Properties with Molecular Simulation**, *European Symposium on Applied Thermodynamics 2017,* Bucharest, Romania, 2017.

**How Uncertainty Quantification Renders Molecular Simulation a Quantitative Tool for Thermophysical Property Evaluation,** *American Institute of Chemical Engineers Annual Meeting*, San Francisco CA, 2016.

**Uncertainty Quantification of Intermolecular Parameters with a Transferable Potential Model for**

***n*-Alkanes,** *American Institute of Chemical Engineers Annual Meeting*, San Francisco CA, 2016.

**Uncertainty Quantification and Propagation Associated with a Transferable Intermolecular Potential Model for *n*-Alkanes,** *International Conference on Properties and Phase Equilibria for Product and Process Design*, Porto Portugal, 2016.

**A Statistical Analysis of the Propagation of Error Associated with the Intermolecular Potential Model When Simulating Large Compounds,** *American Institute of Chemical Engineers Annual Meeting*, Salt Lake City Utah, 2015.

**A Statistical Approach to Reducing Finite-Size Effects from Gibbs Ensemble Monte Carlo Simulations for Predicting the Critical Point,** *American Institute of Chemical Engineers Annual Meeting*, Salt Lake City Utah, 2015.

**Group Contribution Model for Predicting Critical Volume with the Flory-Huggins Theory Asymptotic Behavior,** *19th Symposium on Thermophysical Properties*, Boulder Colorado, 2015.

**Molecular Simulations of the Critical Point for Molecules that Decompose Experimentally,** *American Institute of Chemical Engineers Annual Meeting*, Atlanta Georgia, 2014.