Lixue Cheng

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EDUCATION

California Institute of Technology Pasadena, CA

Ph.D. in Chemistry Mar 2022

University of Wisconsin-Madison Madison, WI

Bachelor of Sciences in Honors, GPA: 3.933/4.0 May 2016

Majors: Chemistry (Honor), Mathematics (Honor), Biochemistry, Molecular Biology
Minor (Certificate): Computer Sciences

RESEARCH INTERESTS

Apply computational and mathematical tools, especially machine learning (ML), to improve the current computational chemistry study strategies in

(1) Quantum simulation and electronic structure (MOB-ML)

(2) Molecular computing and nucleic acid design (INSPIRE)

(3) Bayesian optimization for experimental design (ODBO)

RESEARCH AND TEACHING EXPERIENCES

Tencent Quantum Lab July 2022– present

Research scientist

Miller Group, California Institute of Technology July 2017– Mar 2022

PhD student

MOB-ML: Molecular Orbital-Based Machine Learning

• Use the molecular orbital representations to predict the molecular energies via ML tools to give high accuracy as wave function theory (WFT) method but a low computational cost as density functional theory (DFT) method

• Perform electronic structure calculations and quantum simulations for different molecular systems

• Develop different supervised and unsupervised learning methodologies to model different molecular properties

• Develop a ML module in open-source quantum chemistry software (*entos*) and build a MOB-ML database

INSPIRE: Computational Parameterization of Nucleic Acid Secondary Structure Models

• Combine ML tools with molecular dynamics (MD) to derive new thermodynamics and kinetics nucleic acid models with secondary structure information

• Develop and revise ML algorithms that can improve the efficiency of the strategy

• Perform MD simulations for different nucleic acid reaction systems as a source of data for ML model

Record Group, University of Wisconsin-Madison January 2013–July 2017

Undergraduate researcher

Interpreting and predicting Hofmeister salt ion, polyol, PEG and sugar effects on biopolymer processes using solubility assay and solute-partitioning model

• Design and lead projects using solubility experiments to measure different solutes-biomolecule interactions

• Recruit and mentor other undergraduate students to collect solute–nucleobase thermodynamics data

• Develop mathematical models and analyze the data to interpret functional group chemical interactions

• Determine interaction constants between different bases and different solutes and apply them to predict and interpret solute-biopolymer interaction and drug solubility data

Math studies supervised by Qin Li, University of Wisconsin-Madison August 2015–May 2016

Undergraduate researcher

Randomized Singular Value Decomposition algorithm for low rank matrices and its numerical applications

• Reprove and improve the theorems in the literature as a mathematical exercise

• Develop MATLAB codes to implement the algorithm studied in the literature and perform numerical tests for it

• Compare and analyze the numerical test results and explain the mathematical origins in different cases

Grader/ TA for Chem/Biochem 565/665 Biophysical Chemistry

University of Wisconsin-Madison 2015 Fall and 2016 Fall

Grader/ TA for Math 341 Linear Algebra (Honor) and Math 561 Differential Geometry

University of Wisconsin-Madison 2016 Spring

VOLUNTEERING AND WORK EXP ERIENCES

Tencent Quantum Lab, Tencent July 2021– September 2021

Research Intern

• Work on the interface between machine learning and physical problems, including molecular modeling, drug design and material design.

• Develop, implement and test Bayesian learning algorithm in real world problems in the fields of biology, chemistry and material sciences.

Department of Mathematics, University of Wisconsin–Madison

Administrative Office and Front Desk Student Help August 2015–May 2016

Sonoco Products Co., 455 Science Drive, Madison, WI

Research Project Assistant September 2014–January 2015

• Measure geometries and properties of different materials and perform mathematical modeling of the products

• Analyzing data using various software, for example, Excel, Auto CAD, FEA

• Develop test devices and search the literature to support our own research projects

Department of Chemistry and Mathematics, University of Wisconsin–Madison

Private Tutor August 2014–May 2017

College Library, University of Wisconsin–Madison

Circulation Student Assistant August 2013–May 2015

Greater University Tutoring Service (GUTS), University of Wisconsin–Madison

Academic Tutor (Chemistry and Mathematics) September 2012–August 2013

PATENTS, PUBLICATIONS AND THESES

• Miller III, T.F.; Welborn, M.; **Cheng, L**.; Husch, T.; Song, J.; Kovachiki, N.; Burov, D.; Teh, Y.S.; Anandkumar, A.; Ding, F.; Lee, S.J.R.; Qiao, Z.; Lale, A.S. Systems and methods for determining molecular structures with molecular-orbital-based features. U.S. Patent 16817489, 2020 [Link]
• **Cheng, L**.; Sun, J.; Deustua, J. E.; Bhethanabotla,V.C; Miller III, T.F. Molecular-orbital-based machine learning for general ground state electronic structure theories with kernel addition Gaussian process regression: Closed-shell and open-shell systems. *In preparation* (2022)
• Sun, J.; **Cheng, L**.; Miller III, T. F. Molecular dipole moment learning via rotationally equivariant Gaussian process regression with derivatives in molecular-orbital-based machine learning. arXiv:2205.15510 (2022). [Link]
• **Cheng, L**.; Sun, J.; Miller III, T.F. Accurate molecular-orbital-based machine learning energies via unsupervised clustering of chemical space. arXiv: 2204.09831 (2022). [Link]
• Sun, J.; **Cheng, L**.; Miller III, T.F. Molecular energy learning using alternative blackbox matrix-matrix multiplication algorithm for exact Gaussian process. arXiv: 2109.09817 (2022). [Link].
• **Cheng, L**.; Yang, Z.; Liao, B.; Hsieh, C.; Zhang, S. ODBO: Bayesian Optimization with prescreening for directed protein evolution. arXiv:2205.09548 (2022). [Link]
• Lu, F.*; **Cheng, L**.*; DiRisio, R. J.*; Finney, J. M.; Boyer, M. A.; Sun, J.; Lee, S. J. R.; Deustua, J. E.; Miller III, T. F.; McCoy, A. B. Fase near *ab initio* potential energy surfaces using machine learning. *J. Phys. Chem. A,* **2022.** (*co-first author) [Link]
• Husch, T.; Sun, J.; **Cheng, L.**; Lee, S. J. R.; Miller III, T.F. Improved accuracy and transferability of molecular-orbital-based machine learning: Organics, transition-metal complexes, non-covalent interactions, and transition states. *J. Chem. Phys.*, **2021.** [Link]
• **Cheng, L.**; Kovachki, N; Welborn, M.; Miller III, T. F. Regression clustering for improved accuracy and training costs with molecular-orbital-based machine learning. *J. Chem. Theory Comput.*, **2019**. [Link]
• **Cheng, L.**; Welborn, M.; Miller III, T. F. A universal density matrix functional from molecular orbital-based machine learning: Transferability across organic molecules. *J. Chem. Phys.*, **2019.** [Link]
• Welborn, M.; **Cheng, L.**; Miller III, T. F. Transferability in machine learning for electronic structure via the molecular orbital basis. *J. Chem. Theory Comput.* **2018**.[Link] (Highlighted with commentary in C&EN and Caltech News)
• Knowles, D. B.; Shkel, I. A.; Phan, N. M.; Sternke, M.; Lingeman, E.; Cheng, X.; **Cheng, L.**; O'Connor, K.; Record, M. T. Chemical interactions of polyethylene glycols (PEGs) and glycerol with protein functional groups: Applications to effects of PEG and glycerol on protein processes. *Biochemistry* **2015**, *54* (22), 3528–3542. [Link]
• Undergraduate Senior Honor Thesis for Chemistry with Prof. M. Thomas Record: *Interactions of polyols with aromatics and nucleobases: Large opposing contributions from interactions of nonpolar and polar groups*
• Undergraduate Senior Honor Thesis for Mathematics with Prof. Qin Li: *Randomized singular value decomposition (RSVD) algorithm for low rank matrices and its numerical applications*

HONORS AND AWARDS

- *2017* Caltech General Fellowship, Division of Chemistry and Chemical Engineering, Caltech
- *2016* Mary Ellen Rudin Foundation Scholarship Award, Department of Mathematics, UW-Madison
- *2015* David H. Durra Scholarship, College of Letters & Sciences, UW-Madison
- *2015* Eugene and Patricia Kreger Herscher Undergraduate Scholarship for the 2015–16 Academic Year, Department of Chemistry, UW-Madison
- *2015* Walter W. & Young-Ja C. Toy Summer Research Fellowship, Department of Chemistry, UW-Madison
- *2014* Biochemistry Undergraduate Summer Research Scholarship, Department of Biochemistry, UW-Madison
- *2014* Gerald W. and Tui G. Hedstrom Scholarship, College of Letters & Science, UW-Madison
- *2014* Martha Gunhild Weeks Undergraduate Scholarship for the 2014-15 Academic Year, Department of Chemistry, UW-Madison
- *2014* The Honor Society of Phi Kappa Phi Member, University of Wisconsin-Madison Chapter
- *2014* Trewartha Senior Thesis Grant, Letters & Science Honors Program, UW-Madison
- *2014* Wisconsin Hilldale Undergraduate/Faculty Research Fellowship, UW-Madison
- *2013* John & Elizabeth Moore Awards for Excellence, Department of Chemistry, UW-Madison: Excellence for Chem109 Advanced General Chemistry, 1/35