

ZHENYU WEI

+86 150-0553-2910 \diamond zhenyuwei@seu.edu.cn

Meiyuan M1D # 436 \diamond Southeast Road # 2

Jiangning, Nanjing, Jiangsu, China 211189

EDUCATION

Southeast Univeristy

June 2021

B.S. candidate in School of Mechanical Engineering (GPA: 3.62 / 4.00, Ranking: 18 / 196)

Relevant Courses:

- Theoretical Mechanics (A⁺)
- Materials Mechanics (A⁺)
- Heat Transfer Theory (A)
- Thermal Dynamics (A⁺)
- Electronic Technique (A⁺)
- Programming and Algorithm Language (A⁺)

Stanford Univeristy, Summer session

June 2019 - September 2019

- ME346C, Advanced technology for MD simulation by Prof. Wei Cai. (A⁺)
- ME414, Introduction to Solid State Physics by Prof. Thomas W. Kenny (A⁺)

ACADEMIC AND RESEARCH EXPERIENCE

Independent research of ion hydration and friction in aqueous solution

Directed by Prof. Yunfei Chen, Southeast University

April 2020 - Present

- Obtained deeper understanding of research of aqueous solution, a review under progress.
- Conducted Molecular Dynamic (MD) simulation with LAMMPS.
- Analyzed structure change of ions' hydration shell.
- Analyzed the friction of ions that results from hydrodynamic and dielectirc effects.
- Finding the reason of the abnomral increasing of ion mobility under high external electric fields (0.1V/nm - 3V/nm).

Development of toolboxes for MD simulation

Self-developed and open-sourced

October 2019 - Present

- *Lammps_Toolbox* : A comprehensive toolbox for MD simulation, based on Matlab. Utilized by graduate students of Tsinghua Univeristy, Peking Univeristy and Centralsouth Univeristy.
- *lmp_str* : Structure maker for MD simulation, based on Julia. Utilized by a graduate student in the Baylor University.
- *lmp_data* : Data analyzer for results from MD simulation, based on Julia.

Research of abnormal ions' transportation in nanopore of Si3N4 under external electric field

Directed by Prof. Yunfei Chen, Southeast University

September 2019 - April 2020

- Constructed model of nanopore and aqueous solution with *lmp_str*.
- Conducted MD simulations with LAMMPS.
- Analyzed the data of MD simulation with *Lammps_Toolbox*.

Implementation of Artificial Neural Network (ANN) for hydrogen bonding detecting

Self-developed

January 2020 - February 2020

- Acquired the basic knowledge and techniques for Machine Learning.
- Implemented the Backpropagation (BP) algorithm based on Julia.
- Tested the stability of my neural network package through a familiar topic.
- Thought about the intersection research between MD and ML.

Lecture of Quantum Computation and Communication

Produced by Prof. Jianwei Pan, University of Science and Technology of China

January 2020

- Obtained an overview of the current research of quantum computation.
- Thought about the research of many-body questions (MD, and ab init) with the assistance of a quantum computer.
- Thought about the possible revolution of simulation after the development of a robust quantum computer.

Independent development of MD Programming (basic) based on C++

Research of the melting points of copper

July 2018 - September 2018

- Obtained a basic understanding about MD simulation
- Implemented Embedded-Atom-Method (EAM) potential.

Participating Sakura Exchange Program by MSTC and JST

Start point of my research travel

June 2016 - July 2016

- Conducted experiment with Hideki Shirakawa, co-recipient of the 2000 Nobel Prize in Chemistry.
- Took lecture by Toshihide Maskawa, co-recipient of the 2008 Nobel Prize in Physics.
- Took lecture by Mamoru Mohri, the first astronomer in Japan.

TECHNICAL STRENGTHS

Computer Languages	Julia, C++, Matlab, Python
MD simulation package	LAMMPS, Hoomd-Blue
Visualizing Software	OVITO

PATENT

An energy resolver for passengers' protection

June 2018

Patent ID: ZL201610068802.5