Rushi Gong

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Education

• The Pennsylvania State University

Ph.D. Materials Science and Engineering

University Park, PA

2020 - 2024

- 4.0 GPA

• Beihang University (Beijing University of Aeronautics & Astronautics)

B.S. Materials Science and Engineering; Minor, Mathematics

Beijing, China

2015 - 2019

- 3.76 GPA

- Merit Student (Top 4%)

Research Experience

• Phases Research Lab, The Pennsylvania State University

University Park, PA

2020 - Present

Graduate Research Assistant (Advisor: Prof. Zi-Kui Liu)

- First to apply thermodynamic modeling in the investigation of site nuclearity on Pd-Zn-based catalysts surfaces
- Developed the Pd-Zn-based alloy thermodynamic databases and quantified uncertainty by leveraging the distribution of model parameters during optimization for accurate nuclearity design
- Built framework to drive selections for stable intermetallic catalysts candidates with Machine Learning and firstprinciples calculations tools
- Developed DFTTK structure builders to automatically generate structures for high-throughput computations
- The Nuclear Science and Engineering Division, Argonne National Laboratory Research Aide Technical PhD (Advisor: Dr. Shayan Shahbazi)

Lemont, IL

07/2022 - 09/2022

- Developed LiF-LnF3 thermodynamic databases to predict relative volatility of Ln
- Quantified uncertainty and sensitivity of thermodynamic modeling of fluoride molten salts
- International Research Institute for Multidisciplinary Science, Beihang University Undergraduate Research Assistant (Advisor: Prof. Qianfan Zhang)

Beijing, China

2017 - 2019

- Built stable substrates with transition metal adsorbed on two-dimensional materials as catalysts
- Designed performance analysis via computational methods to examine the thermodynamics and kinetics of hydrogen evolution reaction on the catalytic substrates
- Department of Materials Science and Engineering, Rensselaer Polytechnic Institute Undergraduate Research Assistant (Advisor: Prof. Yunfeng Shi)

Troy, NY 2018

- Investigated the corrosion process of the glass nanowire by using Molecular Dynamics simulations

- Quantified the corrosion rate and analyzied the relationship between pre-tension and corrosion

Teaching Experience

• Department of Materials Science and Engineering, Penn State University Teaching Assistant University Park, PA

2021

- MatSE 410: Phase Relations in Materials Systems

Technical Skills

Computational Tools and Software: Python, MongoDB, VASP, Thermo-Calc, Matlab, ATAT, PyCalphad, ESPEI **Software Developing:** DFTTK (github.com/phasesresearchlab/dfttk)

Publications

- 3. **R. Gong**, S. L. Shang, H. Sun, M. J. Janik, and Z. K. Liu, Thermodynamic modeling of the Pd-Zn system with uncertainty quantification and its implication to tailor catalysts, **Calphad**, 2022, doi.org/10.1016/j.calphad.2022.102491.
- 2. A. Dasgupta, H. He, **R. Gong**, S. L. Shang, E. K. Zimmerer, R. J. Meyer, Z. K. Liu, M. J. Janik, and R. M. Rioux, Atomic control of active site ensembles in ordered alloys to enhance hydrogenation selectivity, **Nature Chemistry**, 14, 523–529 (2022), doi: 10.1038/s41557-021-00855-3.
- 1. J. P. S. Palma, **R. Gong**, B. J. Bocklund, R. Otis, M. Poschmann, M. Piro, Y. Wang, T. G. Levitskaia, S. Hu, H. Kim, S. L. Shang, and Z. K. Liu, Thermodynamic modeling with uncertainty quantification using the modified quasichemical model in quadruplet approximation: Implementation into PyCalphad and ESPEI, arxiv.org/abs/2204.09111.