Amir Mahdi Zardoshti

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Curriculum Vitae

Last Updated November 2025

Education

2024 – present

• Master of Science in Physical Chemistry Chemistry Department, Sharif University of Technology, Tehran, Iran.

2020 - 2024

• Bachelor of Science in Applied Chemistry Chemistry Faculty, School of Chemistry, University of Tehran, Tehran, Iran.

Research of Interest

- Research interests centre on the application of machine learning methods and artificial intelligence techniques in computational chemistry and molecular modelling. The initial motivation arose from studies in Quantitative Structure–Activity Relationship (QSAR) analysis, employing methods such as Principal Component Analysis (PCA), Support Vector Machines (SVMs), and voxel-based convolutional neural network. Further work has explored the representation of molecular structures as point clouds, leading to the use of point cloud-based neural network architectures. During the master's programme, attention expanded to Message Passing Neural Networks (MPNNs), which are capable of accurately modelling potential energy surfaces (PESs) and benefit from transfer learning.
- Current research interests also include data preparation and sampling strategies, with particular emphasis on uncertainty estimation through query-by-committee frameworks, and the use of Gaussian Process Regression (GPR) as a probabilistic approach to energy-surface prediction and uncertainty quantification.
- A long-term objective is the development of machine learning interatomic potentials that integrate
 the advantages of MPNNs and GPR to provide accurate, transferable models for quantum-chemical
 systems and molecular dynamics.

Research Experience

2023 - 2024

Industrial Intern. Tarmim Darou Behavar Company (brand name: "Behrokh"). Synthesis of Metal-Organic Frameworks (MOFs): UiO-66(Zr) and derivatives. Explored applications in catalysis and drug delivery. Heterocyclic Compounds Synthesis & Nucleotide Chemistry. Preparation and characterisation of: – Oxazole derivatives as a novel class of VEGFR2 kinase inhibitors. – 2-Arylbenzimidazoles via copper-catalysed condensation of anilines with benzoic acids.

Research Experience (continued)

2022 - 2024

Undergraduate Researcher. Artificial Intelligence in Drug Design, IBB Institute, University of Tehran, Tehran, Iran. Worked under the supervision of Dr. Mohammad Taheri-Ledari and Dr. Kaveh Kavousi; represented molecular structures as point clouds and applied point cloud-based deep learning to predict drug efficacy.

Current Research

 MSc proposal research: Learning Potential Energy Surfaces for Quantum-Accurate Molecular Dynamics and Spectroscopy Under the supervision of Prof. Zahra Jamshidi, Sharif University of Technology.

Research applies **Machine Learning Interatomic Potentials (MLIPs)** with quantum chemistry methods to predict molecular properties accurately and efficiently.

The main applications of this work include:

- Finding minimum-energy structures of molecules and small metal clusters using Neural Networks and Gaussian Process Regression.
- Enhancing Molecular Dynamics simulations for Infrared (IR) spectral prediction.
- I would like to learn and work in the area of Excited-State Non-Adiabatic Molecular Dynamics in the future.
- I would like to learn and work in the area of Tunnelling Splitting Calculations in the future.
- Generative point-cloud models combined with transformer-based SMILES captioning for de novo drug design. Conceived and currently developing this project independently (since February 2025); the core idea is to translate a molecule's point-cloud representation—derived from its electrostatic potential energy surface—into a SMILES format. The work involves employing the FoldingNet architecture as a point cloud auto-encoder, integrated with a Transformer-based molecular captioning model.

Skills

Programming Languages

• **Python** (intermediate), **C**++ (beginner)

Deep Learning

 Experienced with PyTorch and TensorFlow 1.x; worked with point cloud-based deep learning models such as PointNet, PointNet++, and Point Cloud Auto-encoder (FoldingNet), as well as message passing neural networks including SchNet, PhysNet, and PAINN.

Quantum Chemistry

Experienced with **geometry optimization**, **vibrational frequency**, and **TDDFT** calculations using **Gaussian** and **ORCA**; performed **ab initio molecular dynamics (AIMD)** simulations in **CP2K** using both **KS-DFT** and **SCC-DFTB** methods; currently learning **SHARC** and **MCTDH** for **excited-state dynamics**.

Visualisation Tools

Avogadro, VMD, PyMOL.

Languages

• Native in Persian; fluent in English; interested in learning French and German.

Miscellaneous Experience

Awards and Achievements

• 16th Rank Nationwide in Iran's Master's Entrance Exam — Admitted to Sharif University of Technology.

2018, 2019, 2024

• Iranian National Chemistry Olympiad Finalist — Qualified three times (twice in high school, once in university).

• **Top Rank in National University Entrance Exam**, admitted to the University of Tehran for BSc in Applied Chemistry.

• Exceptional Talents School, Studied at a nationally recognized school.

• Member of Iran National Youth Football Team.

Teaching Experience

Fall 2025

 Teaching Assistant — Computational Chemistry, Department of Chemistry Sharif University of Technology Tehran, Iran.

Spring 2025

- Teaching Assistant General Mathematics I, Faculty of Chemistry, School of Chemistry, University of Tehran, Tehran, Iran. View Teaching Certificate
- Teaching Assistant Analytical Chemistry I, Faculty of Chemistry, School of Chemistry, University of Tehran, Tehran, Iran. View Teaching Certificate

Certification

February 2024

 Executive Committee Certificate — Recognized for active participation in organising the 5th Winter School of Computational Chemistry, Sharif University of Technology, under the supervision of Prof Zahra Jamshidi. View Certificate

Publications

Zardoshti, Amir Mahdi; Zarvani, Maral; Taheri-Ledari, Mohammad; Kavousi, Kaveh.
 Enhancing Drug Design for VEGFR2: Integrating Quantum Mechanics-Driven 3D QSAR with Deep Learning to Predict Drug Efficacy. The 12th National Conference and the 3rd
 International Bioinformatics Conference, Behshahr. Proceedings link

References

• Prof. Zahra Jamshidi (M.Sc. Supervisor)
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