



# Sartaaj Khan

## Contact Information

Email: [sartaaj.khan@mail.utoronto.ca](mailto:sartaaj.khan@mail.utoronto.ca)

## Education

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### University of Toronto

Sept. 2023 – Present

**PhD** Chemical Engineering and Applied Chemistry  
Principal Investigator: Dr. Seyed Mohamad Moosavi  
cGPA: 4.00/4.00

### University of Toronto

Sept. 2018 – May 2023

**BASc** Chemical Engineering and Applied Chemistry  
Minor in Bioengineering  
cGPA: 3.73/4.00

## Research Experience

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### Doctoral Researcher

Sept. 2023 – Present  
Toronto, ON

AI4ChemS Laboratory, University of Toronto

- Developed a deep learning model that utilizes powder x-ray diffraction (PXRD) patterns and SMILES precursors for metal-organic framework property predictions.
- Enhanced model accuracy through self-supervised and representation learning techniques, and fine-tuned for properties such as band gap, gas uptake at varying pressures, pore volume and Henry coefficient.
- Built a recommendation system that can give a suggestion on the optimal application of the MOF given its PXRD and precursors.
- Aided in the construction of a physics-informed machine learning model to predict dynamic viscosity of thermal fluids from small datasets and conducted uncertainty assessments of predictions.
- Leading a project focusing on constructing a model that can find the mathematical representations of different MOF topologies and validating against mechanical properties.
- Supervised undergraduate students in pretraining/finetuning of physics-informed machine learning models and in the construction of mathematical representations of topologies.

### Undergraduate Researcher

July 2021 – Sept. 2022  
Toronto, ON

Advanced Membranes Laboratory, University of Toronto

- Developed a model showing the concentration profiles of common ions found in brackish water in a reverse osmosis process.
- Created a model involving different process modules in series such as a reverse osmosis membrane and a custom membrane specified.
- Used elemental-based modelling to compute the concentration profiles, flux of water and the flux of solute with respect to position and recovery ratio while including more complexities.
- Created and published a Python solution chemistry package called pySolution for the purpose of calculating parameters that are involved in non-ideal solutions from the Pitzer model.
  - GitHub repository is available [here](#).

### Undergraduate Researcher

June 2021 – Sept. 2021  
Los Angeles, CA

Ping Lab, University of California, Los Angeles



- Developed a bioinformatics tool for generating a more reliable protein sequence database for sus scrofa cardiac proteomic studies.
  - GitHub repository is available [here](#).
- Researched and developed features relevant in proteomics such as sequence length, PE score and number of confirmed isoforms.
- Created a classifier trained against homo sapien protein data from UniProt and tested against mus musculus and sus scrofa protein data to determine “reliable” and “unreliable” proteins in those respective organisms.
- Constructed cardiovascular knowledge graphs and created link prediction algorithms to find unknown connections between proteins in different organisms – specifically in cardiac conduction and muscle contraction.
  - GitHub repository is available [here](#).

## Undergraduate Researcher

May 2020 – Sept. 2020

Laboratory of Metabolic Systems Engineering, University of Toronto

Toronto, ON

- Created algorithms that genetically modify E. coli models and performed constrained optimization to optimize the flux of around 70 different industrially important chemicals through flux balance analysis.
- Modified mcPECASO simulator to obtain results with two-stage bioprocesses to analyze the impact of phenotypic switches on produced flux in iJO1366 and iML1515 E. coli models.
  - GitHub repository is available [here](#).

## Publications

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- **Khan ST**, Moosavi SM. Connecting metal-organic framework synthesis to applications with a self-supervised multimodal model. ChemRxiv. 2024; doi:10.26434/chemrxiv-2024-mq9b4 (under review)
- **Khan, ST**, Moosavi, SM. XRayPro: A self-supervised multimodal model for MOF application recommendations using PXRD and precursors. In *AI for Accelerated Materials Design-NeurIPS 2024*.
- **Khan, S. et al.** (2024). Reflections from the 2024 Large Language Model (LLM) Hackathon for Applications in Materials Science and Chemistry. arXiv preprint arXiv:2411.15221.

## Invited Presentations

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- **Khan, S.T.**, Moosavi, S.M. "XRayPro: A Self-Supervised Multimodal Model for MOF Application Recommendations Using PXRD and Precursors." AI for Accelerated Materials Design - NeurIPS, 2024, [\[Paper\]](#).
- **Khan, S.T.** (2024, October 8). Unveiling Property Insights and Recommending MOF Applications from X-Ray Diffraction Patterns using Self-Supervised Transformer Models. [\[Presentation\]](#). Canadian Society for Chemical Engineering (CSCHE) 2024, Toronto, Canada
- **Khan, S.T.**, Moosavi, S.M. (2024, August 7). Using a self-supervised model for property predictions and MOF application recommendations with only powder x-ray diffraction patterns. [Poster]. Accelerate Conference 2024, Vancouver, Canada

## Work Experience

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### Artificial Intelligence Engineer

Sept. 2021 – Sept. 2022

MGRID Energy Inc.

Whitehorse, YK

- Created and tested a variety of object detection models (primarily YOLO) using convolutional neural networks for the purpose of vehicle classification.
- Designed the architecture of a hybrid nonlinear regression-convolutional neural network model for the automatic detection of wildfires.



- Researched and extracted different features to include in the wildfire models through literature done on previous machine learning models in the area.
- Generated a synthetic dataset for the purpose of training a natural language processing (NLP) model to send alerts on new projects for bidding.

## Research Mentorship Experience

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1. **Hari Om Chadha** (Summer 2024): University of Toronto undergraduate student in Engineering Science; Pretraining graph convolutional networks using Barlow-Twin for predicting fluid properties.
2. **Josh Goldman** (Summer 2024): McGill University undergraduate student in Statistics and Computer Science; Topological deep learning in metal-organic frameworks.

## Awards

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### LLMs in Chemistry and Materials Hackathon

May 2024

Awarded 3<sup>rd</sup> place (\$250 prize funded by Radical AI and Anthropic) in the LLM Hackathon for Applications in Materials and Chemistry for the development of **PoreVoyant** – a chemistry-informed AI agent that can generate new linkers to decrease the band gap in metal-organic frameworks while leveraging MOF literature. This work was featured in a [Medium article](#). The GitHub repository can be found [here](#).

### Bayesian Optimization for Chemistry and Materials Hackathon

March 2024

Awarded 2<sup>nd</sup> place (\$500 CAD prize funded by Acceleration Consortium) in the [Bayesian Optimization for Chemistry and Materials Hackathon](#) hosted by Acceleration Consortium at the University of Toronto. The project involved the application of Bayesian Optimization to accelerate the discovery of fluids with the highest heat transfer coefficients. The repository can be found [here](#).

### Dean's List

Sept. 2018 – May 2023

Award issued for achieving an average of at least an A- per semester with a full course load. This was achieved for all semesters of undergraduate.

### NSERC USRA

May 2020

Award of \$6000 CAD issued for completion of 16 weeks of research work for the summer 2020 term and high academic achievements.

## Skills

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**Programming:** Python, MATLAB and Simulink, Cypher Query Language, Neo4j

**Software:** ChemDraw, AutoCAD, Aspen Plus, Aspen Dynamics

**Applications:** Microsoft Word, Microsoft PowerPoint, Microsoft Excel, Microsoft VISIO