# SADMAN SADEED OMEE

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# **EXPERIENCE**

# **University of South Carolina**

January 2022 - Current

Graduate Research Assistant | Machine Learning and Evolution Laboratory

Columbia, SC, United States

- Research topic: We use deep learning techniques to solve materials informatics problems (e.g., crystal structure prediction, materials property prediction) in our lab. My work on crystal structure prediction is similar to the protein structure prediction problem solved by AlphaFold2. I have also developed a state-of-the art global attention-based Graph Neural Network architecture for the materials property prediction problem using differentiable group normalization and skip-connections.
- Concept / skills usage: I mostly use Graph Neural Networks (GNNs) for my research. I have also used Convolutional Neural Networks, Transformers and Generative Models for my other researches. I am fluent in using PyTorch and PyTorch Geometric.

## **University of South Carolina**

**August 2021 - December 2021** 

Graduate Teaching Assistant | Course: CSCE102 (General Applications Programming)

Columbia, SC, United States

• Taught HTML, CSS, and JavaScript to a total of 72 students of three sections students, and two lab group of total 50 students.

#### **Bangladesh University of Engineering and Technology**

May 2019 - July 2021

Research Student | Graph Drawing and Information Visualization Laboratory

Dhaka, Bangladesh

• Worked on unsupervised learning, mostly developing hierarchical clustering algorithms.

#### **PUBLICATIONS**

#### **Journal Articles**

- 1. **Omee, S. S.**, Louis, S. Y., Fu, N., Wei, L., Dey, S., Dong, R., Li, Q., & Hu, J. (2022). Scalable deeper graph neural networks for high-performance materials property prediction. *Patterns*.
- 2. Varivoda, D., Dong, R., **Omee, S. S.**, & Hu, J. (2023). Materials property prediction with uncertainty quantification: A benchmark study. *Applied Physics Reviews*.
- 3. Dong, R., Zhao, Y., Song, Y., Fu, N., **Omee, S. S.**, Dey, S., Li, Q., Wei, L., & Hu, J. (2022). DeepXRD: A deep learning model for predicting XRD spectrum from material composition. *ACS Applied Materials & Interfaces*.
- 4. Hu, J., Stefanov, S., Song, Y., **Omee, S. S.**, Louis, S. Y., Siriwardane, E. M., Zhao, Y., & Wei, L. (2022). MaterialsAtlas.org: A materials informatics web app platform for materials discovery and survey of state-of-the-art. *npj Computational Materials*.
- 5. Fu, N., Wei, L., Song, Y., Li, Q., Xin, R., **Omee, S. S.**, Dong. R., & Hu, J. (2023). Material transformers: Deep learning language models for generative materials design. *Machine Learning: Science and Technology*.
- 6. Louis, S. Y., Siriwardane, E. M. D., Joshi, R. P., **Omee, S. S.**, Kumar, N., & Hu, J. (2022). Accurate prediction of voltage of battery electrode materials using attention-based graph neural networks. *ACS Applied Materials & Interfaces*.
- 7. Li, Q., Dong, R., Fu, N., **Omee, S. S.**, Wei, L., & Hu, J. (2023). Global mapping of structures and properties of crystal materials. *Journal of Chemical Information and Modeling*.
- 8. Wei, L., Fu, N., Siriwardane, E. M., Yang, W., Omee, S. S., Dong, R., Xin, R., & Hu, J. (2022). TCSP: a template-based crystal structure prediction algorithm for materials discovery. *Inorganic Chemistry*.

# RELEVANT SKILLS

**Programming languages**: Python, C, C++, Java, JavaScript.

ML / DL frameworks: PyTorch, Tensorflow, PyTorch Lightning, Keras, Scikit-learn.

ML / DL libraries: PyTorch Geometric (PyG), Deep Graph Library (DGL), Transformers (Hugging Face), PyTorch-Ignite, Pandas, NumPy, SciPy, Pymatgen, ASE, Matminer.

Other skills: LATEX, Git, Matplotlib, Seaborn, HTML, CSS, MySQL.

# **EDUCATION**

## **University of South Carolina**

August 2021 - Present

Columbia, SC, United States

*Ph.D. in Computer Science* Current GPA: 3.95/4.00

**Bangladesh University of Engineering and Technology** 

February 2015 – April 2019

B.S. in Computer Science

Dhaka, Bangladesh