## **JUSTIN Z. TAM**

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#### **SUMMARY**

Ph.D. student in Computer Science at Lehigh University specializing in bioinformatics, with experience as a Bioinformatics research co-op at Moderna. Skilled in machine learning, 3D modeling, full-stack development, and software engineering tools. Developed novel LLM and graph-based deep learning models to enhance Moderna's computational mRNA pipeline, driving innovation in the field. Expert in applying graph-based machine learning methods to structural protein and mRNA research, with a strong focus on biochemical-driven solutions.

#### **EDUCATION**

## Ph.D. Candidate in Computer Science/Bioinformatics

Aug 2020 - Dec 2025

Lehigh University, P.C. Rossin College of Engineering, Bethlehem, PA

Advisor: Brian Y. Chen

## B.A. in Mathematics, Minor in Computer Science and Chemistry

Aug 2015 - May 2019

Skidmore College, Saratoga Springs, NY

#### **TECHNICAL SKILLS**

**Machine Learning:** (proficient) Tensorflow, PyTorch, transformer and deep learning models, GPT Prompt Engineering, (familiar) Scala

Programming: (proficient) Python, C, C++, JavaScript (familiar) Java

Full Stack Web Development: NodeJS, ReactJS, MongoDB

Other Technical Skills: Git, Matlab, Unix, Docker, Google Cloud, AWS, 3D visualization

Other Relevant Skills: Presentation, Project Planning, Independent Research Development and Proposal, Team

Management, Wet Lab Protein Research, Biochemistry

#### **EXPERIENCE**

### Moderna Therapeutics, Cambridge, MA: Bioinformatics Research Co-op

Jan 2024 - June 2024

- Developed a PyTorch-based transformer model and graph neural network model for mRNA analysis
- Discovered a novel method for altering normally inflexible mRNA structural properties using graph features.
- Engineered several production-level features and improvements to the current Moderna computational pipeline.
- Presented comprehensive reviews and project insights to executive leadership at Moderna, including the VP and directors of computational science.

### New York City College of Technology, Brooklyn, NY: Adjunct Professor in Mathematics Aug 2019 - June 2020

- Aided in the development of a web-based interactive homework assignment and grading system, used in all math classes throughout the college.
- Taught several college level mathematics courses including pre-calculus, calculus 1, and calculus 2.
- Utilized a mix of lecture-style and activity-style teaching strategies to maximize participation and understanding of difficult topics.

### **PUBLICATIONS**

# Tam, J.Z. et al. A Containerization Framework for Bioinformatics Software to Advance Scalability, Portability, and Maintainability

Proceedings of the 14th ACM International Conference on Bioinformatics, Computational Biology and Health Informatics, Sept 3–6, 2023

- Independent full-stack development project to build a container framework with reusable templates, streamlining the dissemination and deployment of Bioinformatics software.
- Awarded best paper finalist of CSBW at ACM-BCB 2023

Liu, Y., Armstrong, G., Tam, J.Z., & Chen, B. Y. MechPPI: Binding Mechanism-based Machine-Learning tool for Predicting Protein-Protein Binding Affinity Changes Upon Mutations bioRxiv, Oct 31, 2023

• Presented a state-of-the-art ML method leveraging gradient boosted trees and shapley values to curated mechanistic features for analyzing protein-protein binding affinity changes.

## Tam, J. Z. et al. HBcompare: Classifying Ligand Binding Preferences with Hydrogen Bond Topology Biomolecules, MDPI, Oct 28, 2022

- · Presented a graph neural network pipeline for analyzing hydrogen bond topology in proteins
- Performed a benchmark study using various machine learning methods like graph neighborhood aggregation strategies, graph kernels, principal component analysis, and convolutional neural networks.

## Tam, J. Z., Palumbo, T., Miwa, J. M., & Chen, B. Y. Analysis of Protein-Protein Interactions for Intermolecular Bond Prediction

Molecules, MDPI, Sep 21, 2022

• Designed a novel graph network representation of protein-protein interactions, accurately representing intermolecular bonds with efficient representations using networkx.

## Tam, J. Z., Palumbo, T., Miwa, J. M., & Chen, B. Y. DiffBond: A Method for Predicting Intermolecular Bond Formation

2021 IEEE International Conference on Bioinformatics and Biomedicine (BIBM), IEEE, Dec 9, 2021

### **LANGUAGES**

English (Native)
Mandarin (Fluent)
Cantonese (Limited Working)