## Yuxiang Zhang

Undergraduate of Shanghai Jiao Tong University E-mail: zhangyuxiangSJTU@outlook.com

## **EDUCATION**

Shanghai Jiao Tong University, School of Medicine, Shanghai, China Bachelor of Biomedical Science (major)

GPA: 3.75/4.3 Sep. 2020-Present Expected in June 2024

Shanghai Jiao Tong University, SEIEE, Shanghai, China Bachelor of Computer Science and Technology (minor) Sep. 2021-Present Expected in June 2024

## RESEARCH EXPERIENCES

Molecular Design Laboratory, Shanghai Jiao Tong University, School of Medicine

Aug. 2022-Present

Molecular Dynamics Simulation in Biomolecules; supervisor: Jian Zhang and Shaoyong Lu

Yan Lab, Shanghai Jiao Tong University, School of Life Sciences and Biotechnology

Mar. 2022-Aug.2022

➤ Protein Function Analysis and Drug Design; supervisor: Wupeng Yan

Xiong Lab, Shanghai Jiao Tong University, School of Life Sciences and Biotechnology

Apr. 2021-Mar. 2022

Machine Learning in Drug-drug Interaction; supervisor: Yi Xiong

# **SKILLS**

- ➤ Conducting MD simulations of complex protein system
- > Trajectory analysis of MD simulations
- ➤ Enhance sampling methods (NEB and GaMD)

- > Virtual scanning of small molecules
- > C/C++, Python, R and bash shell
- ➤ Machine Learning Framework

# RELATED PROJECTS&COURSES

## **Projects**

#### **Molecular Dynamics Simulation in Biomolecules**

- Decoupling the dynamic mechanism revealed by FGFR2 mutation-induced population shift, J. BIO. STR. DYN., 2023
  - Utilized GaMD simulation in 6 different systems, revealing the long range allosteric pathway in FGFR2 kinase
  - ➤ Co-first author | Received 22 Feb 2023, Accepted 10 Apr 2023.
- Large-scale computational analysis of Peptide-GPCR-G protein system (Unpublished Data)

#### **Protein Function Analysis and Drug design**

- > Drug design of small molecule inhibitor targeting ras family protein Rab43 phosphokinase
  - > Applied virtual scanning on LRRK2 kinase domain, obtaining 10 potential compounds

#### **Machine Learning in Drug-drug Interaction**

- > Drug-drug Interaction Prediction Method Based on Deep Learning
  - > Using deep neural network and convolutional neural network to predict the subtype of drug-drug interaction

#### **Courses**

JCCX0001 Machine Learning for Biomedical Signal Processing Grade: A

➤ Basic biostatistics analysis skills using R; Implementation classic machine learning and DL algorithm using Python

PHY1202H Introduction to Physics (HONOR) Grade: A+ CS3322 Data Structure Grade: A

MATH1607H Mathematical Analysis (HONOR) Grade: A

MATH1205H Linear Algebra (HONOR) Grade: A

BIO2355 Biochemistry Grade: A+ BIO3350 Cell Biology Grade: A

BIO2357 Applied Data Science Grade: A CS2501 Discrete Mathematics Grade: A

#### **HONORS&AWARDS**

2022 Zhiyuan Honor Scholarship	Dec. 2022
2022 Xiaomi Scholarship	Dec. 2022
2022-2023 Academic Year SJTU Excellent Undergraduate Scholarship (B)	Nov. 2022
The 6th Biomedical Science Forum (School of Medicine, SJTU) Personal Report First Prize	Sep. 2022
Shanghai Jiao Tong University 40th PRP Excellent Project	Mar. 2022
2021 Zhiyuan Honor Scholarship	Dec. 2021
2021-2022 Academic Year SJTU Excellent Undergraduate Scholarship (B)	Dec. 2021
2020 Zhivuan Honor Scholarship	Dec. 2020