

KAI ZHENG, Ph.D.

Lecturer & Master Supervisor | Computational Biology & AI

CONTACT

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KEY METRICS

21
h-index
1600+
Total Citations
NeurIPS
Premier AI Venue

EXPERTISE

- Geometric Deep Learning
- Graph Neural Networks
- AI for Genomics
- Non-coding RNA
- Drug-Target Prediction
- Bioinformatics Tools

HONORS

🏆 Best Paper Award, ICIC 2019.

FUNDING

Principal Investigator:
Fundamental Research Funds
for the Central Universities (No.
25CX06033A).

EXECUTIVE SUMMARY

Researcher bridging **Artificial Intelligence** and **Computational Biology**. Specializes in developing **Geometric Deep Learning** frameworks to solve complex problems in genomics and therapeutics. Proven track record of publishing in top-tier methodology conferences (**NeurIPS**) and leading application journals (*PLOS Computational Biology*).

EXPERIENCE & EDUCATION

Lecturer / Master Supervisor	2024.07 – Present
China University of Petroleum (East China)	
Ph.D. in Computer Science	2020 – 2024
Central South University	
M.Sc. in Computer Science	2017 – 2020
China University of Mining and Technology	

EDITORIAL & ACADEMIC SERVICE

- PC Member: KDD 2025/2026, BIBM 2025/2026.
- Review Editor: *Frontiers in Bioinformatics*.
- Session Chair: ISAICS 2025.
- Reviewer: *Briefings in Bioinformatics*, *IEEE/ACM TCBB*.

FULL PUBLICATION LIST

* Categorized by research topic. All listed papers are indexed by SCI.

📁 I. AI Methodology & Graph Learning

1. Graph-Theoretic Insights into Bayesian Personalized Ranking for Recommendation.
Zheng K, Wang JX, Xu JH. *Advances in Neural Information Processing Systems (NeurIPS)*, 2025. [**Premier AI Venue**]
- 📁 II. Genomics & Non-coding RNA Analysis
1. Line graph attention networks for predicting disease associated Piwi-interacting RNAs.
Zheng K, Zhang XL, Wang L, et al. *Briefings in Bioinformatics*, 2022 [Q1, IF: 9.5].
2. SPRDA: a link prediction approach based on the structural perturbation to infer disease-associated Piwi-interacting RNAs.
Zheng K, Zhang XL, Wang L, et al. *Briefings in Bioinformatics*, 2023 [Q1, IF: 9.5].
3. iCDA-CGR: Identification of CircRNA-Disease Associations based on Chaos Game Representation.
Zheng K, You ZH, Wang L, et al. *PLOS Computational Biology*, 2020 [Q1].
4. DBMDA: A unified embedding for sequence-based miRNA similarity measure.
Zheng K, You ZH, Wang L, et al. *Molecular Therapy - Nucleic Acids*, 2020 [Q1, IF: 8.8].

5. **MISSIM: an incremental learning-based model for miRNA-disease association.**
Zheng K, You ZH, Wang L, et al. *IEEE/ACM Transactions on Computational Biology and Bioinformatics*, 2020 [Q1].
6. **iMDA-BN: Identification of miRNA-disease associations based on the biological network.**
Zheng K, You ZH, Wang L, et al. *Computational and Structural Biotechnology Journal*, 2020 [Q1].
7. **MLMDA: a machine learning approach to predict miRNA-disease associations.**
Zheng K, You ZH, Wang L, et al. *Journal of Translational Medicine*, 2019 [Q1, IF: 7.4].

III. Computational Drug Discovery

1. **LRTM: Left-Right Transition Matrices for Molecular Association Prediction.**
Zheng K, Duan GH, Yang MY, et al. *IEEE/ACM Transactions on Computational Biology and Bioinformatics*, 2025 [Q1].
2. **DLP: duplex link prediction via subspace segmentation for predicting drug-miRNA associations.**
Zheng K, Zhao QC, Liang X, et al. *IEEE/ACM Transactions on Computational Biology and Bioinformatics*, 2024 [Q1].
3. **NASMDR: a framework for miRNA-drug resistance prediction using efficient neural architecture search.**
Zheng K, Zhao H, Zhao Q, et al. *Briefings in Bioinformatics*, 2022 [Q1, IF: 9.5].