### Jiahui Chen

CONTACT Information Department of Mathematics

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ACADEMIC PREPARATION

Michigan State University, East Lansing, Michigan

Postdoc, 2019-present (Mentor: Guo-Wei Wei) Southern Methodist University, Dallas, Texas

PhD, Computational and Applied Mathematics, 2014-2019 (Advisor: Weihua Geng)

Xi'an Jiaotong-Liverpool University, Suzhou, China

BSc Applied Mathematics, May 2014

WORKING EXPERIENCE Michigan State University, East Lansing, MI

Visiting Assistant Professor

August 2019 - Present

Michigan State University, East Lansing, MI

Visiting scholar

May - July, 2018

Pacific Northwest National Laboratory, Richland, WA

PhD internship

May - August, 2016

RESEARCH INTERESTS

Mathematical Biology/Biophysics; Topological Data Analysis; Differential Geometry; Numerical PDB; Boundary Integral Equations; Boundary Element Method; Treecode and Fast Multipole Method; Machine Learning/Deep Learning; and Parallel/GPU Computing

**PUBLICATIONS** 

(\* co-first author)
Published/Accepted

- 19. **J. Chen**, K. Gao, R. Wang, and G.W. Wei, Revealing the threat of emerging SARS-CoV-2 mutations to antibody therapies, *Chemical Science*, in publication (2021)
- 18. R. Wang, **J. Chen**, K. Gao, G.W. Wei, Vaccine-escape and fast-growing mutations in the United Kingdom, the United States, Singapore, Spain, India, and other COVID-19-devastated countries, *Genomics*, 113 (4), 2158-2170, 2021
- 17. **J. Chen\***, K. Gao\*, R. Wang\*, and G.W. Wei, Prediction and mitigation of mutation threats to COVID-19 vaccines and antibody therapies, *Chemical Science*, 12 (20), 6929-6948, 2021
- 16. R. Wang, R. Zhao, E. Ribando-Gros, **J. Chen**, Y. Tong, and G.W. Wei, HERMES: Persistent spectral graph software, *Foundations of Data Science*, in publication (2021)
- J. Strubbe-Rivera, J. Chen, B. A. West, K. N. Parent, G.W. Wei, and J. N. Bazil, Modeling the Effects of Calcium Overload on Mitochondrial Ultrastructural Remodeling, Applied Sciences, DOI:10.3390/app11052071 (2021)
- 14. R. Wang, **J. Chen**, K. Gao, Y. Hozumi, C. Yin, and G.W. Wei, Analysis of SARS-CoV-2 mutations in the United States suggests presence of four substrains and novel variants, *Communications Biology*, 4 (1), 1-14 (2021)
- 13. **J. Chen**, R. Wang, and G.W. Wei, SARS-CoV-2 becoming more infectious as revealed by algebraic topology and deep learning, *Communications in Information and Systems*, 21 (1), 31-36, (2021)
- 12. **J. Chen**, W. Geng, and D. Reynolds, Cyclically parallelized treecode for fast computations of electrostatic interactions on molecular surfaces, *Computer Physics Communication*, 260, 107742 (2021)

- J. Chen, J. Hu, Y. Xu, R. Krasny, and W. Geng, Computing protein pKas using the TABI Poisson-Boltzmann Solver, Journal of Computational Biophysics and Chemistry, 2042006, (2020)
- R. Wang, J. Chen, Y. Hozumi, C. Yin, G.W. Wei, Decoding asymptomatic COVID-19 infection and transmission, The Journal of Physical Chemistry Letters, 11 (23), 10007-10015, (2020)
- 9. **J. Chen**, R. Wang, M. Wang, and G.W. Wei, Mutations strengthened SARS-CoV-2 infectivity, *Joiunnal of molecular biology*, 432 (19), 5212-5226, (2020)
- 8. R. Zhao, M. Wang, **J. Chen**, Y. Tong, and G.W. Wei, The de Rham-Hodge Analysis, and Modeling of Biomolecules, *Bulletin of Mathematical Biology*, 82 (8), 1-38, (2020)
- 7. **J. Chen\***, K. Gao\*, R. Wang, D. Nguyen, and G.W. Wei, Review of COVID-19 antibody therapies, Annual Review of Biophysical, 50, (2020)
- 6. K. Gao, D. Nguyen, **J. Chen**, R. Wang, and G.W. Wei, Repositioning of 8565 existing drugs for COVID-19, *The Journal of Physical Chemistry Letters*, 11, (13), 5373-5382, (2020)
- D. Nguyen, K. Gao, J. Chen, R. Wang, and G.W. Wei, Unveiling the molecular mechanism of SARS-CoV-2 main protease inhibition from 92 crystal structures, *Chemical Science*, 11, 12046-12046, (2020)
- 4. **J. Chen**, R. Zhao, Y. Tong, and G.W. Wei, Evolutionary de Rham-Hodge method, *Discrete & continuous Dynamical Systems B*, doi: 10.3934/dcdsb.2020257 (2020)
- 3. **J. Chen** and W. Geng, On Preconditioning the Treecode-Accelerated Boundary Integral (TABI) Poisson-Boltzmann Solver, *J. Comput. Phys.*, 373, 750-762 (2018)
- E. Jurrus, D. Engel, K. Star, K. Monson, J. Brandi, L. E. Felberg, D. H. Brookes, L. Wilson, J. Chen, K. Liles, M. Chun, P. Li, D. W. Gohara, T. Dolinsky, R. Konecny, D. R. Koes, J. E. Nielsen, T. Head-Gordon, W. Geng, R. Krasny, G.-W. Wei, M. J. Holst, J. A. McCammon, and N. A. Baker, Improvements to the APBS Biomolecular Solvation Software Suite, Protein Science, 27(1), 112-128 (2018)
- J. Chen, W. Geng, Parallel Computing of the Adaptive N-body Treecode Algorithm for Solving Boundary Integral Poisson-Boltzmann Equation, Lecture Notes in Computer Science, 9576, 1-9, Springer (2016)

## In revision/Revision submitted

1. K. Gao\*, R. Wang\*, **J. Chen**, L. Cheng, J. Frishcosy, Y. Huzumi, Y. Qiu, T. Schluckbler, and G.W. Wei, Methodology-centered review of molecular modeling, simulation, and prediction of SARS-CoV-2, *Chemical Review*, in revision, 2021

#### Honors and Awards

Dedman College, SMU, Graduate Student Travel Grant Award, \$500, 2019

Dedman College, SMU, Graduate School Dissertation Completion Fellowship, \$20,000, 2018

Math Dept., Southern Methodist University, Teaching Assistant Award, Cash award, 2018

#### Papers in Preparation

- With L. Wilson, W. Geng, R. Krasny, "Improvements to the TABIPB C-Version"
- With Q. Sun, D. Chan and W. Geng, "Regularization for Boundary Integral Poisson-Boltzmann Equation and Treecode Acceleration"
- With W. Geng, J. Tausch, "Adaptive Cartesian FMM Accelerated Poisson-Boltzmann Solver"
- With W. Geng, T. Schlick, G. Bascom, "Multipole Method Accelerated Electrostatics Computation for Mesoscale Chromatin Simulation"
- With H-X Zhou, W. Geng, J. Tausch, "Cartesian FMM for Modeling Protein Folding and Binding under All-Atom Crowders"

# Conferences and Presentations

- 10. Workshop on Modeling and Analysis in Molecular Biology and Electrophysiology, Duke Kunshan University, 2021 (presentation)
- 9. Computational Biology Forum, MSU, January 2021 (presentation)
- 8. Computational topology/geometry/graph seminar, webinar, December 2020 (presentation)
- 7. Colorado State University, webinar, November 2020 (presentation)
- 6. NSF-CBMS Conference: Mathematical Molecular Bioscience and Biophysics, the University of Alabama, Tuscaloosa, Alabama, May 2019 (poster)
- 5. IMSM workshop, SAMSI, North Carolina State University, July 2018 (presentation)
- 4. Workshop on the Mathematics of Drug Design/Discovery, the Fields Institute, Toronto, Canada, June 2018 (attendance)
- 3. Multiple Faces of Biomolecular Electrostatics, MBI, Ohio State University, October 2016 (poster)
- 2. Mathematics of Biological Charge Transport: Molecules and Beyond, IMA, University of Minnesota, Twin Cities, July 2016 (attendance)
- 1. 2015 SMU PhD interdisciplinary project of data science and analysis (presentation)

#### TEACHING EXPERIENCE

## MTH496, Introduction of Machine Learning

Fall 2020 / Spring 2021 Michigan State University

Math 3316, Introduction to High Performance Scientific Computing Fall 2015
Laboratory TA Southern Methodist University

Graduate research mentoring (With Professor Guo-Wei Wei at MSU)

Ms. Rui Wang
Mr. Yuta Hozumi
Mr. Xiaoqi Wei
2019-present
2020-present
2020-present

#### Undergraduate research mentoring (With Professor Guo-Wei Wei at MSU)

• Ms. Yongjia Xu Summer, 2021

# Professional Services

#### Journal reviewer

Instructor

International Journal for Numerical Methods in Biomedical Engineering, Computational and Mathematical Biophysics, Journal of Chemical Information and Modeling, Frontiers in Pharmacology **Proposal reviewer** 

The French National Research Agency

## Software

- Wei-Lab (MSU) servers maintenance
- C-version Treecode (http://github.com/jiahuic/treecode)
- Parallel C-version Treecode (http://github.com/jiahuic/treecode\_parallel)
- C-version TABI Poisson-Boltzmann Solver (http://github.com/jiahuic/TABI\_solver)

#### ACTIVATES

• Math Outreach, Glencairn Elementary School, East Lansing, January 9. 2020

#### Professional Membership

- American Mathematical Society (2014-)
- Society for Industrial and Applied Mathematics (2014-)

### Computer Skills

- Mathematical and statistical modeling and programming
- Parallel programming: MPI, Open MP, GPU/CUDA
- $\bullet \ \ Programming \ languages: \ FORTRAN, C/C++, HTML, Python, Java, Cython, HTML, JavaScript$
- Mathematical and statistical softwares: Matlab, Mathematica, Excel
- $\bullet$  Bimolecular simulation and dynamics packages: AMBER, CHARMM, APBS, VMD/NAMD, PyMol