

# Tanmoy Pal

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## EDUCATION

### Boston University

PhD Candidate in Computational Chemistry

Boston, MA

Spring 2018 –

### University of Wisconsin-Madison

Master of Science in Chemistry

Madison, WI

Fall 2016 – Fall 2017

### Indian Institute of Science

Bachelor of Science (Research)

Bengaluru, India

Fall 2012 – Spring 2016

*Major: Chemistry | Special focus: Computational Chemistry*

*Thesis: Ab-initio Molecular Dynamics Studies of Carborane Clusters*

## PUBLICATIONS

1. Qiang Cui, Tanmoy Pal, and Luke Xie. “Biomolecular QM/MM Simulations: What Are Some of the “Burning Issues”?” In: *The Journal of Physical Chemistry B* 125.3 (2021). PMID: 33401903, pp. 689–702. DOI: [10.1021/acs.jpcb.0c09898](https://doi.org/10.1021/acs.jpcb.0c09898)
2. [Interfacial Polarization and Ionic Structure at the Ionic Liquid-Metal Interface Studied by Vibrational Spectroscopy and Molecular Dynamics Simulations](#), *Matthew J. Voegtle, Tanmoy Pal, Anuj K. Pennathur, Sevan Menachekanian, Joel G. Patrow, Sohini Sarkar, Qiang Cui, and Jahan M. Dawlaty*, *J. Phys. Chem. B*

## RESEARCH

- Boston University | University of Wisconsin-Madison Spring 2017 –  
*Advisor: Prof. Qiang Cui*
  - QM/MM Simulations of Mechanism of Proton Transfer in Cytochrome c Oxidase
  - Study of Two-body and Three-body Energies of Water using Semi-empirical Models, Calculation of Second Virial Coefficients of Small Molecules
  - Multidimensional DFTB3 Free Energy Calculations using Reinforcement Learning
  - Interfacial Polarization and Ionic Structure at the Ionic Liquid-Metal Interface Studied by Vibrational Spectroscopy and Molecular Dynamics Simulations
  - Machine Learning Assisted High-accuracy Multidimensional Free Energy Surfaces
- Indian Institute of Science, Bengaluru
  - Ab-initio Molecular Dynamics (CPMD) and Ab-initio Computational Modelling of Carboranes  
*Advisor: Prof. E D Jemmis* Summer 2015 - Spring 2016
  - Generating Formula for Condensed Polyhedral Boranes and its Computer Implementation  
*Advisor: Prof. E D Jemmis* Summer 2014
  - Generation of Promoter:Reporter Constructs to Study Cytokinin Distribution Profile in Rice (*Oryza sativa*)  
*Advisor: Prof. Usha Vijayraghavan* Summer 2013
- Indian Institute of Science Education and Research, Kolkata
  - Quantum Harmonic Oscillator: A Special Case of Quasi-exactly Solvable Equations  
*Prof. Prasanta K Panigrahi* Summer 2012

## TEACHING

- [BU] Mathematical Methods for Molecular Sciences CH 225 Fall 2020
- [BU] Mathematical Methods for Molecular Sciences CH 225 (Grader) Spring 2020
- [BU] General Chemistry for Engineering Sciences CH 131 Fall 2019
- [BU] General Chemistry CH 102 Spring 2019
- [UWMadison] General Chemistry Chem 104 Spring 2017
- [UWMadison] General Chemistry Chem 103 Fall 2016

## TECHNICAL EXPERIENCE

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- Programming Languages
  - Python, Linux shell, L<sup>A</sup>T<sub>E</sub>X, C/C++, Javascript, OpenMP, MPI, OpenACC, CUDA, Wolfram Language, MySQL
  - Python Packages: TensorFlow/Keras, Bokeh, Matplotlib, NumPy, SciPy, Pandas, cuDF, Requests
- [HackerRank](#)
- Computational Chemistry Software
  - CHARMM, Gaussian, CPMD, CP2K, Quantum Espresso, GROMACS, ASE, SchNetPack, DeePMD
- Art and Design
  - Adobe Photoshop, Illustrator, InDesign, Muse, Edge Animate, Premiere Pro
  - Procreate [Procreate Portfolio](#)

## RELEVANT COURSEWORKS AND WORKSHOPS

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| • Telluride School on Theoretical Chemistry at Telluride Science Research Center, Telluride, CO  | Summer 2017 |
| Project: <i>Programming Davidson's Diagonalization Method</i>  |             |
| • [BU] Machine Learning by Prof. Kate Saenko   | Fall 2018   |
| Project: <i>Classification of Bacteriophage Genome with Harvard Innovation Lab's XGENOMES</i>  |             |
| • [UWMadison] High Performance Computing for Engineering Applications by Prof. Dan Negrut  | Fall 2017   |
| Project: <i>Parallel Implementation of Davidson's Diagonalization Method: Comparison of Performance Between Eigen C++, OpenMP, NumPy, and cuSolver</i> |             |
| • [UWMadison] Molecular Dynamics and Monte Carlo Simulations in Materials Science  | Spring 2017 |
| • [BU] Advanced Molecular Quantum Mechanics by Prof. David Coker   | Spring 2018 |
| • [UWMadison] Special Topics in Physical Chemistry by Prof. Qiang Cui  | Spring 2017 |
| • [IISc] Computational Modeling of Materials   | Spring 2014 |
| • [IISc] Organic Electronics   | Spring 2015 |
| • [IISc] Introduction to Electronics and Electrical Engineering  | Spring 2013 |
| • [IISc] Algorithms and Programming  | Fall 2012   |

## CERTIFICATIONS

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- DeepLearning.AI TensorFlow Developer Certification
  - [Introduction to TensorFlow for Artificial Intelligence, Machine Learning, and Deep Learning](#)
- IBM Data Science Professional Certificate
  - [Machine Learning with Python](#)
  - [Data Science Methodology](#)
  - [Data Science Orientation](#)
  - [Data Visualization with Python](#)
  - [Open Source Tools for Data Science](#)
  - [Python for Applied Data Science](#)

## HONORS AND AWARDS

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- Delegate at 22<sup>nd</sup> Session of Youth Assembly (formerly Youth Assembly at the United Nations)
- Hirschfelder Prize Graduate Fellowship at Theoretical Chemistry Institute, Department of Chemistry, University of Wisconsin-Madison (2016)
- Nominated by Ministry of Human Resource Development, India for Commonwealth Scholarship for PhD in the United Kingdom (2015)
- National Initiative on Undergraduate Science (NIUS) Chemistry fellow (batch IX 2012-14)
- KVPY National Fellowship (India) (2011)