

AAMOD ATRE

In Pursuit | M. Sc. Quantum Science and Technology

About Me

1. **INTERESTED** in understanding the interplay of quantum dynamical processes across various timescales
2. **ENJOY** delving into the theoretical underpinnings of a problem and exploring systems computationally
3. **INTEND** to pursue the development of theoretical and numerical tools in quantum optics and many-body physics

Education History

Post-Graduate GPA 1.5/5.0 Max. 1.0	Technical University of Munich M. Sc. Quantum Science and Technology	Oct 2022 - Present
Undergraduate GPA 8.65/10.0	Birla Institute of Technology and Science, Pilani B.E. Chemical Engg. Minor : Physics	Aug 2017 - Aug 2021
Higher Secondary Percentage : 96.5 %	FIITJEE Junior College, Kukatpally, Hyderabad Grades XI & XII	Apr 2015 - Aug 2017
Senior Secondary GPA : 10.0/10.0	Delhi Public School, Nacharam, Hyderabad Grades IX & X	Mar 2013 - Mar 2015

Academic Research

Master's Thesis **Apr. 2024 - Present**
Prof. Frank Pollmann | Technical University of Munich

Numerically studying measurement-induced entanglement transitions, in random quantum circuits of varying local Hilbert-space dimensions, to understand the transition universality and distill a classical analogue of the transition.

Undergraduate Thesis **Jan. 2021 - June 2021**
Prof. Jeremy Richardson | ETH Zürich

Gauged the efficacy of a spin-mapping-based semiclassical dynamics technique by a Python-based numerical study of the population and coherence dynamics of exciton relaxation in a 1D polymer chain. Formulated a spin-mapping dynamics algorithm for tight-binding polymer chains which conform to $SU(2)$ symmetry.

Work Experience

Student Assistant | HiWi **May. 2024 - Present**
Quantum Computing and Technologies Dept. | Leibniz Supercomputing Centre (LRZ), Munich
Benchmarking C++ and Python-based quantum simulation programs using the QAOA algorithm to solve the Max-Cut problem. Facilitating the integration of the quantum simulation programs into LRZ's HPC infrastructure.

Student Assistant | HiWi **Apr. 2023 - Sept. 2023**
Prof. Peter Rabl | Walther Meißner Institute, Munich
Compared various numerical integration approaches in stochastic Master equation simulations of a dissipative cascading quantum network, to facilitate simulations of larger system sizes.

Remote Research Assistant **Oct. 2021 - Oct. 2022**
Dr. Aaron Kelly | Max Planck Institute for the Structure and Dynamics of Matter, Hamburg
Benchmarked a semiclassical mapping-based dynamics method (spin-PLDM) by modelling the interactions between a two-level atomic subsystem and a cavity-modified field.

Remote Research Internship **July 2021 - Sept. 2021**
Prof. Pengfei Huo | University of Rochester, New York
Numerically compared the performance of standard and spin-based partially linearized density matrix (PLDM) algorithms in calculating linear absorption spectra of a bi-exciton coupled dimer model.

Research Internship **May 2019 - July 2019**
Prof. Bibek Dash | CSIR Institute of Minerals and Materials Technolgy, India
DFT-based computational designing of triazole-based molecular precursors for selective CO_2 capture. Statistically determined the optimum functional-basis combination to model the CO_2 -triazole interactions with DFT. Studied CO_2 interactions with the aromatic building blocks to propose a new triazole moiety design.

Technical Skills		Languages
Programming	Software Packages	English (Native/Bilingual)
Python	MATLAB	German (B1)
C++	Maple	French (Elementary)
Julia/C (Elementary)	Mathematica	Marathi Hindi (Native)
	LAMMPS	
	Quantum Espresso	
	Gaussian 09	
Standardized Tests		
GRE General Test : 335/340 Quant : 168/170 Verbal : 167/170 Analytical Writing : 5/6		Aug. 2021
TOEFL : 115/120 Reading : 30/30 Listening : 30/30 Speaking : 25/30 Writing : 29/30		Sept. 2021
Undergraduate Teaching Experience		
Teaching Assistant <i>Dr. S. D. Manjare Process Design Principles - I</i>		Aug. 2020 - Dec. 2020
Teaching Assistant <i>Dr. Radhika Vathsan Quantum Mechanics - II</i>		Aug. 2020 - Dec. 2020
Featured Undergraduate Projects		
Designing Lithium-based metal organic frameworks for hydrogen production <i>Design Project Dr. Paramita Halder</i>		Jan. 2020 - June 2021
<i>Ab initio</i> computational study of graphene-based Li-MOFs, employing density functional theory implementation in Quantum Espresso for electronic structure calculations & reactive molecular dynamics within LAMMPS framework to study hydrogen evolution amongst proposed models.		
Study of cavity QED formalism and modern quantum control techniques <i>Study Project Dr. Raghunath Ratabole</i>		Aug. 2020 - Dec. 2020
Literature survey of modern qubit implementations, entangled state preparation and quantum gate implementations in molecule-coupled cavity systems. Study of field quantization, cavity QED formalism and the applications of Jaynes-Cummings model.		
Modelling kinetics of photo-catalytic reactions involved in waste-water treatment <i>Study Project Dr. Sharad Sontakke</i>		Aug. 2019 - Dec. 2019
Modelled TiO_2 -based photocatalytic degradation of phenol and extraction of Cr and Cu ions with MATLAB. Optimized TiO_2 catalyst concentration and reaction rates for varying contamination levels.		
Study of metal-organic frameworks as tools for adsorptive CO_2 capture <i>Study Project Dr. Richa Singhal</i>		Jan. 2019 - May. 2019
Literature review of thermodynamic and electronic properties of MOFs. Studied the methodologies and developments in the field of CO_2 capture, focusing on MOFs.		
Extracurricular Activities		
Quantum Computing	· Cleared Quantum Open Source Foundation (QOSF) cohort 4 assesment task	Sept. 2021
	· Completed IBM Quantum's Global Summer School	June 2020, July 2022
Social Groups	· Member - Offerings Dept. : <i>PushQuantum</i>	Nov. 2023 - Present
	· Core Member : <i>Kala - Fine Arts Club</i>	Mar. 2018 - May 2020
Hobbies	· Long distance running, and hiking in the Bavarian Alps.	
	· Avid reader of fiction and non fiction literature	

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