



Manas Sharma

PHD (PHYSICS)

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Research Interests

My passion lies in **developing codes and methods** for modeling various phenomena and processes in materials science such as **light-matter interactions**, **adsorbate-substrate interactions**, and **thin film deposition**. During my Ph.D., I focused on improving and accelerating **quantum embedding techniques** for molecular and periodic systems. I implemented **density functional theory (DFT)-based embedding** methods and coupled them with **real-time time-dependent DFT (RT-TDDFT)** and **wavefunction theory (WFT)** methods. Notably, I calculated excitation energies of solvated molecules and adsorption energies of molecule-periodic systems using WFT-in-DFT, achieving high accuracy at a fraction of the computational cost. My work also extended to **molecule-in-periodic embedding** and **high harmonic generation** via RT-TDDFT.

In my postdoctoral research, I have diversified into **machine-learned interatomic potentials (MLIPs)**, particularly **graph neural network (GNN)**-based potentials, to study **chemical vapor deposition (CVD)** or **physical vapor deposition (PVD)** of thin films on substrates. To explore deposition processes at the atomic scale, I use DFT calculations and **nudged elastic band (NEB)** methods to study **reaction pathways** and **energy barriers**, as well as **ab initio molecular dynamics (AIMD)** to study dynamic evolution. Using the data from these simulations, I develop MLIPs trained on high-fidelity data and also leverage **foundation models** to extend their applicability.

Beyond simulations, I have a keen interest in **developing computational tools** for researchers. I have built a performant DFT code as well as a neural network library from scratch, optimized for parallelization and GPU acceleration. Additionally, I have experience creating GUIs for materials modeling, visualization, and input file generation.

Outside research, I enjoy creating [YouTube tutorials](#), web & [Android apps](#), and computer software/[libraries](#) for researchers and students.

About me

I am a postdoctoral researcher in the Chemical Engineering department at the Indian Institute of Science (IISc), jointly advised by [Prof. Ananth Govind Rajan](#) and [Prof. Sudeep Punathanam](#).

Previously, I obtained my [PhD](#) (in Physics) from Friedrich Schiller University Jena (FSU) in Germany, with the highest distinction of summa cum laude. I was supervised by [Prof. Dr. Marek Sierka](#) during my PhD.

I possess strong verbal, presentation, and written communication skills as demonstrated by extensive **participation in >17 conferences (7 talks; 10 posters)** as well as publishing **7 scientific articles**.

Prior to joining FSU, I obtained a Master's in Physics from the University of Delhi (India), where I worked with [Dr. Debabrata Mishra](#) and published three papers. I obtained my Bachelor's in Physics (Hons) from the University of Delhi as well.

I am originally from a small, but beautiful town Haldwani, a few minutes from Nainital in India.

Awards and Achievements

2025	Top Viewed Article , Paper investigating high harmonic generation in molecular thin films was one of the most viewed article among the works published in Advanced Optical Materials between January 1 to December 31 2023	–
2024	Summa Cum Laude , Grade awarded for PhD Thesis	Germany
2024	Travel and Accommodation Bursary , Awarded for presenting my work at Turbomole 2024 Conference	UK
2024	3.5 Million Views and 19.9k Subscribers , on YouTube channel , where I share my knowledge on computational materials science, physics and scientific computing	Online
2022	Travel Bursary , Awarded for presenting a talk at ETSF 2022 Conference	Germany
2021	Best Oral Presentation , Awarded for my talk at EMRS Fall 2021 Meeting	Germany
2019	Highly Downloaded Paper , My paper on CrysX was among the top 10 percent of most downloaded papers published in Journal of Applied Crystallography between January 2018 and December 2019	India
2017	Rank #3 , Ranked at the 3rd position in the Physics department of my college during M.Sc. Physics	India
2016	DC Arora Scholarship , Awarded for being a meritorious student during B.Sc. Physics	India
2016	Rank #2 , Ranked at the 2nd position in the Physics department of my college during B.Sc. Physics	India

Education

University of Delhi (Sri Venkateswara College)

New Delhi, India

Master of Science (MSc) Physics

July 2016 - July 2018 (2 years)

- Completed a project on density functional theory and material modeling that resulted in two publications in 2019
- Specialized in Nanoscience
- Ranked 3rd in the Physics department of my college

University of Delhi (Acharya Narendra Dev College)

New Delhi, India

B.Sc (Honours) Physics

July 2013 - July 2016 (3 years)

- Aggregate percentage: 87.08
- Ranked second in the Physics department of my college (Acharya Narendra Dev College)
- Earned the D.C. Arora meritorious Scholarship, Acharya Narendra Dev College (Apr 2016)
- Was an active member of the Physics society there and held various posts

Work Experience

Indian Institute of Science

Bengaluru, India

Postdoctoral Researcher

1st August 2024 - present

- Development of graph neural network-based machine learning interatomic potentials (MLIPs) for investigating thin film deposition and for materials design.
- Extensive first-principles (DFT, AIMD and NEB) calculations for reaction pathways and transition state barriers.
- Molecular dynamics simulations powered by foundation MLIPs for various properties such as radial distribution functions, density, and so on.
- I also served as the teaching assistant for a course on Quantum-Mechanical Modeling of Nanomaterials, where I graded assignments and student exams.
- Serving as a teaching assistant for an NPTEL course on Machine Learning for Core Engineering Disciplines where I help create question papers, assignments, and reviewing course materials and lectures.

Friedrich Schiller University Jena

Jena, Germany

Wissenschaftlicher Mitarbeiter (Scientific Employee)

3rd June 2019 - 4th April 2024 (4 yrs 10 mos)

- I developed computationally efficient ab-initio methods and codes to study the properties of hybrid systems. I mainly programmed in Fortran and Python for this role
- Furthermore, I served as a teaching assistant for a number of courses:
 - Modelling and Simulation (Curriculum included basics of quantum mechanics, overview of standard algorithms utilized in quantum chemistry codes such as geometry optimization, and so on.)
 - Basics of Stochastics and Experimental Design
 - Algorithms for Scientific Computing (Curriculum included: Python programming, introduction to machine learning, theory and applications of classical ML models such as SVM, decision trees, and so on.)

As a teaching assistant, I created exercise sheets and examination papers, conducted weekly tutorials, and graded answer sheets. Additionally, I also developed course materials for the course on machine learning and Python.

Phys Whiz (Self-Employed)

Content Creator and Founder

January 2016 - present

- Founded Phys Whiz to create content like tutorials and lectures on physics, computational material science, and scientific computing
- Currently at ~ 18k subscribers on [YouTube](#) and ~ 22k followers on [Instagram](#)
- Acquired various practical skills through this endeavour like image and video editing, graphic designing, animations, basic social media marketing and management skills
- The YouTube channel has garnered 3.2 Million views (approximately)

BragitOff.com (Self-Employed)

Blogger and Founder

June 2014 - present

- Owner and author of [bragitoff.com](#), an educational blog, especially for physics students with a focus on computational material science, numerical techniques, and scientific computing
- The blog gets decent traffic with > 30k monthly views
- Searching '[bragitoff.com](#)' on [Google Scholar](#) gives 28 unique results that have cited various blog posts of mine

Publications

Here is a link to my [Google Scholar](#) 

- *Optical Gaps of Ionic Materials from GW/BSE-in-DFT and CC2-in-DFT*
M. Sharma, and M. Sierka
J. Chem. Theo. Comput. 20, 21, 9592-9605 (2024)
- *TURBOMOLE: Today and Tomorrow*
Y. J. Franzke, C. Holzer, J. H. Andersen, T. Begušić, F. Bruder, S. Coriani, F. Della Sala, E. Fabiano, D. A. Fedotov, S. Fürst, S. Gillhuber, R. Grotjahn, M. Kaupp, M. Kehry, M. Krstić, F. Mack, S. Majumdar, B. D. Nguyen, S. M. Parker, F. Pauly, A. Pausch, E. Perlt, G. S. Phun, A. Rajabi, D. Rappoport, B. Samal, T. Schrader, **M. Sharma**, E. Tapavicza, R. S. Treß, V. Voora, A. Wodyński, J. M. Yu, B. Zerulla, F. Furche, C. Hättig, M. Sierka, D. P. Tew, and F. Weigend
J. Chem. Theo. Comput. 19, 20, 6859-6890 (2023)
Note: The author list for the above publication was determined using alphabetical order, with the exception of the first two authors and the last five corresponding authors.
- *Resonance Effect in Brunel Harmonic Generation in Thin Film Organic Semiconductors*
W. Li, A. Saleh, **M. Sharma**, C. Huenecke, M. Sierka, M. Neuhaus, L. Hedewig, B. Bergues, M. Alharbi, H. ALQahtani, A. M. Azeer, S. Graefe, M. F. Kling, A. F. Alharbi, and Z. Wang
Adv. Optical Mater. 2203070 (2023) [ON COVER]
- *Efficient Implementation of Density Functional Theory based Embedding for Molecular and Periodic Systems using Gaussian Basis Functions*
M. Sharma, and M. Sierka
J. Chem. Theo. Comput. 18, 11, 6892-6904 (2022) [ON SUPPLEMENTARY COVER]
- *Real-time time-dependent density functional theory using density fitting and the continuous fast multipole method*
C. Mueller, **M. Sharma**, and M. Sierka
J. Comput. Chem. 41, 2573–2582 (2020)
- *CrysX: crystallographic tools for the Android platform*
M. Sharma, and D. Mishra
J. Appl. Cryst. 52, 1449-1454 (2019) [ON COVER PAGE]
- *First-principles study of the structural and electronic properties of bulk ZnS and small ZnSn nanoclusters in the framework of the DFT+U method*
M. Sharma, D. Mishra, and J. Kumar
Phys. Rev. B 100, 045151 (2019)
- *DFT+U study of small ZnO nanoclusters*
M. Sharma, and D. Mishra
AIP Conference Proceedings 2142, 110025 (2019)

Skills

Programming languages	C, C++, C#, Python, FORTRAN, Java, shell scripting
DFT methods development	PySCF, TURBOMOLE
Quantum Chemistry packages	Quantum ESPRESSO, PySCF, TURBOMOLE, NWChem, Serenity, ORCA, VASP
Molecular Dynamics packages	LAMMPS, ASE
Modeling and visualization	VESTA, Avogadro, Jmol, VMD, CrysX-3D Viewer, BURAI, Py3Dmol, ASE
Machine/Deep Learning	scikit-learn, TensorFlow, PyTorch, Nequip, CrysX-NN
App development	Android, Windows
Web development	Wordpress, HTML
Data processing/wrangling	pandas, numpy
Data visualization	matplotlib, seaborn
Outreach	YouTube videos, Blog posts, Instagram posts
Debugging and profiling	ARM forge for FORTRAN, VS code for python
Miscellaneous	Video editing, Molecular animations, Graphic designing, Photo editing, Git, MS Excel
Documentation	Markdown, LaTeX/Overleaf, Jupyter notebooks, MS Word, MS Powerpoint
Soft Skills	Time Management, Teamwork, Problem-solving, Documentation, Engaging Presentation

Social Media

LinkedIn <https://www.linkedin.com/in/manassharma07>
Blog <https://bragitoff.com/>
YouTube <https://www.youtube.com/@PhysWhiz>
Twitter <https://twitter.com/manassharma07>
GitHub <https://github.com/manassharma07/>

Instagram










- <https://www.instagram.com/phys.whiz>
- <https://www.instagram.com/ducktape07/>
- https://www.instagram.com/crysx_3d/

Facebook

- <https://www.facebook.com/ducktape07>
- <https://www.facebook.com/bragitoff/>
- <https://www.facebook.com/physwhizforum/>


Google Scholar <https://scholar.google.com/citations?user=WYOEL94AAAAJ&hl=en>
ORCID <https://orcid.org/0000-0002-5346-6280>
Researchgate <https://www.researchgate.net/profile/Manas-Sharma-5>

Conferences, Workshops and Seminars













Attended	IISc-Fujitsu Workshop on Quantum Computing, IISc Bengaluru India (link)	23-24 Jan 2025
Attended	AI/ML Workshop 2025, IISc Bengaluru India (link)	07-08 Jan 2025
Talk	TURBOMOLE Users Meet Developers Conference, Oxford, UK (link)	19 Sep 2024
Lecture+Tutorials	Mastering DFT: A Hands-on Training Workshop, Janardan Singh Foundation (ONLINE)	27 Jul 2024
Poster 	17th International Congress of Quantum Chemistry 2023 in Bratislava, Slovakia (link)	25 Jun-01 Jul 2023
Talk	VISTA Seminars 2022 (VIRTUAL) (link)	30 Nov 2022
Poster 	BIOVIA Conference 2022 (VIRTUAL) (link)	11-13 Oct 2022
Poster 	ICQNN 2022 Conference in Jena, Germany (link)	05-09 Sep 2022
Poster 	Psi-K 2022 Conference at EPFL in Lasuane, Switzerland (link)	22-25 Aug 2022
Talk 	25th ETSF Workshop on Electronic Excitations 2022 in Leuven, Belgium (link)	13-17 Jun 2022
Talk 	Turbomole Developers Seminar Series 2021-2022	21 Feb 2022
Talk	NOA Spring Meeting 2022 (ONLINE) (link)	21-23 Feb 2022
Talk	Computational Methods in Materials Science (CMMS 2021) (ONLINE) (link)	24-25 Sep 2021
Poster	57th Symposium on Theoretical Chemistry (STC 2021) (ONLINE) (link)	20-24 Sep 2021
Talk 	EMRS Fall 2021 meeting (Online) 🏆BEST ORAL PRESENTATION AWARD🏆 (link)	20-23 Sep 2021
Talk 	The Materials and Molecular Modelling Hub++ Annual Conference 2021 (ONLINE) (link)	14-15 Sep 2021
Talk 	DokDok Lite 2021 in Jena, Germany (link)	01-03 Sep 2021
Poster	International Workshop on Recent Developments in Electronic Structure (ES21) (ONLINE) (link)	12-15 Jul 2021
Poster	eSSSENCE-eMMC eMeeting MMMM 2021 (Online) (link)	07-08 Jun 2021
Attended	International Workshop on Computer-Aided Materials Discovery (Weekly Online ZOOM Webinars) (link)	27 May-24 Jun 2021
Talk	NOA CRC 1375 Spring Meeting (ONLINE) (link)	15-17 Feb 2021
Attended	Intel® Software Development Tools for HPC (Webinar) (link)	11-20 Nov 2020
Attended	Intel® Software Development Tools for Artificial Intelligence (Webinar) (link)	17 Nov 2020
Attended	NOA Fall PhD School (ONLINE) (link)	26-28 Oct 2020
Poster+Presentation	NOA Spring School and PI Workshop in Jena, Germany (link)	02-06 Mar 2020
Attended	NOA seminar by Dr. Heiko Appel (MPI Hamburg) in Jena, Germany (link)	21 Jan 2020
Attended	24th ETSF Workshop on Electronic Excitations in Jena, Germany (link)	16-20 Sep 2019
Poster	ICABS 2019 in Bhiwani, India (link)	07-09 Feb 2019
Poster	6th ISIF 2017 in New Delhi, India (link)	13 Dec 2017

Code Development

Contributions

TURBOMOLE | DFT based embedding coupled with WFT and RT-TDDFT methods within the RIPER module of the popular TURBOMOLE package |  www.turbomole.org/

Independently Developed Libraries and Android/PC/Web Applications

CrysX-NN	An efficient neural network library from scratch that supports parallelization and GPUs	 GitHub link
CrysX-3D Viewer	A molecule and crystal viewer that renders high quality visualizations using complex shaders developed using Unity gaming engine. Available on Android, Windows, Mac and Linux	 Project Home page
CrysX-AR	An Android app for augmented reality visualization of molecules and crystals	 Google Play Store link
CrysX-Crystallographic Tools	A set of crystallographic tools (XRD pattern simulation, CIF creator/parser, equation of state fitting, etc.) available as an Android app	 Google Play Store link
CrysX-Demo DFT based Embedding	Online demo of frozen density embedding and projection based embedding	 Web App link
RIPER-Tools	An online set of tools that allows to create input files for the RIPER module of TURBOMOLE from MaterialsProject or PubChem database	 Web App link
VASP-GUI	An online GUI for VASP input preparation and parsing/visualizing output files	 Web App link
CrysX-CompChem File Converter	Web app that allows inter-conversion between various chemical file formats in current use	 Web App link
Basis Set Converter	Web app to inter-convert between various basis set formats, powered by Basis Set Exchange	 Web App link
The Math App	A suite of mathematical tools that has the potential to act as a substitute for Computer softwares like Matlab/Scilab on Android devices	 Google Play Store link
ML: Microstructure Classification Demo	A web demo app of a neural network (crysx_nn) model trained to classify microstructures	 Web App link
ML: MNIST_Plus Digit Classification Demo	A web app that classifies user given handwritten digits using a convolutional network model (PyTorch) trained on a modified MNIST dataset	 Web App link

Languages

English	Bilingual proficiency
Hindi	Native proficiency
German	Elementary proficiency

References available upon request.