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## EDUCATION

<b>University of Pennsylvania</b> Ph.D., Theoretical Chemistry	Philadelphia, PA 2025 - Present
<b>The Cooper Union for the Advancement of Science and Art</b> B.E., Chemical Engineering, <i>Summa Cum Laude</i> Minors in Mathematics and Chemistry	New York, NY 2021 – 2025

## FELLOWSHIPS AND SCHOLARSHIPS

Goldwater Scholar, <i>The Barry Goldwater Scholarship and Excellence in Education Foundation</i>	2024
Full Tuition Scholarship, <i>The Cooper Union</i>	2024

## AWARDS

New York ACS Excellence in Chemistry, <i>New York Section of American Chemical Society</i>	2025
Outstanding Presentation in Computational/Theoretical Chemistry, <i>GCURS Rice University</i>	2024
Norman Perry ChE '49 Internship Fund, <i>The Cooper Union</i>	2024
Daniel E. Kowler ChE '65 Memorial Prize Fund, <i>The Cooper Union</i>	2023

## RESEARCH EXPERIENCE

<b>University of Pennsylvania Department of Chemistry</b> <i>Graduate Researcher</i> , Advisor: Andrew M. Rappe	Philadelphia, PA 2025 - Present
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- Applied pseudopotential planewave DFT to study the relationship between surface chemistry and ferroelectric polarization switching in wurtzite aluminum nitride
- Integrated first-principles simulations with analytical “pencil and paper” theory to study surface charge passivation on AlN slab models relevant for surface engineering for electronic devices

<b>The Cooper Union Department of Chemical Engineering</b> <i>Undergraduate Researcher</i> , Advisor: Abhishek K. Sharma	New York, NY 2024 - 2025
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- Used differentiable molecular dynamics (JAX-MD) to map sequences to network structures analogous to amino acid sequences to three-dimensional protein structures
- Developed original code to simulate and optimize network structures using gradient descent
- Created a model to simulate protein evolution in the presence of mutational stresses

**Stony Brook Institute of Advanced Computational Sciences**

Stony Brook, NY

*Undergraduate Researcher (NSF REU)*, Advisor: Benjamin G. Levine

May – August 2024

- Conducted static electronic structure calculations and non-adiabatic *ab initio* molecular dynamics simulations in an independently led project studying the photochemistry of 2-hydroxyazobenzene
- Mapped out the potential energy surfaces by discovering multiple non-radiative relaxation pathways involving photoisomerization and excited-state intramolecular proton transfer

**The Cooper Union Department of Chemistry**

New York, NY

*Undergraduate Researcher*, Advisor: Robert Q. Topper

2022 - 2024

- Analyzed the decomposition and growth pathways of ammonium nitrate clusters using simulated annealing Monte Carlo and density functional theory in an independently led project
- Explained, for the first time, the patterns for peaks in both the positive and negative ion mass spectra of ammonium nitrate from experimental literature
- Presented the first study on larger ammonium nitrate particles and predicted new morphologies

**Penn State University Department of Material Science**

State College, PA

*Undergraduate Researcher (NSF REU)*, Advisor: Stephanie Law

May – August 2023

- Used Fourier transform infrared (FTIR) spectroscopy and analyzed spectra to assess the potential of Dirac semimetals for infrared photodetection applications
- Compared optical properties of samples of Dirac semimetals fabricated via different methods

**Purdue Energetics Research Center**

West Lafayette, IN

*Undergraduate Researcher (DEVCOM funded)*, Advisor: Stephen Beaudoin

May – August 2022

- Studied particle adhesion in mock polymer-bonded explosives by quantifying Van der Waals forces
- Independently operated an atomic force microscope for 50+ hours to determine Hamaker constants

**TEACHING AND MENTORSHIP EXPERIENCE****University of Pennsylvania Department of Chemistry**

Philadelphia, PA

*Graduate TA*

2025 – Present

- Lectured recitations and graded for General Chemistry covering atomic structure, Bohr model, particle-in-a-box, bonding, gas laws, thermochemistry, chemical kinetics, and electrochemistry

**The Cooper Union Department of Chemistry**

New York, NY

*Physical Chemistry Tutor*

2024 – Present

- Helped students in introductory quantum chemistry with topics such as particle-in-a-box, harmonic oscillator, rigid rotator, variational methods, and perturbation theory

**The Cooper Union Department of Mathematics**

New York, NY

*Math Tutor*

2022 – Present

- Provided tutoring support to undergraduate students in calculus, linear algebra, differential equations, probability, and discrete mathematics
- Delivered clear and effective instruction on mathematical concepts to groups of students in a weekly “Help Room” setting

## PUBLICATIONS

- [2] **Hassan, U. S.**; Mehmood, A.; and Levine, B. G. Static Quantum Mechanical Calculations and Non-Adiabatic Dynamics of 2-Hydroxyazobenzene. (*In preparation*)
- [2] Vazquez, J. M.; Ellis, Lucas G.; **Hassan, U. S.**; Sapp, J. R.; Jensen, R. J.; Vaz, S.; Corti, D. S.; and Beaudoin, S. P. The Effect of Particulate Solids and Plasticizer Content of Polymer Formulation on the Measured Forces Via AFM. *The Journal of Adhesion Science and Technolog.* (Submitted 2025)
- [1] **Hassan, U. S.**; Amat, M. A.; and Topper, R. Q. Decomposition and Growth Pathways of Ammonium Nitrate Clusters and Nanoparticles. *Journal of Physical Chemistry A*, 128, 9184-9194 (2024). DOI: [10.1021/acs.jpca.4c04630](https://doi.org/10.1021/acs.jpca.4c04630)

## TALKS

- [5] **Hassan, U. S.**; Behrendt, D.; and Rappe, A. M. *Polar AlN Surfaces and Defect Migration*. The Center for 3D Ferroelectric Microelectronics Manufacturing (3DFeM<sup>2</sup>) Thrust 1, Virtual, 2025.
- [5] **Hassan, U. S.**; Mehmood, A.; and Levine, B. G. *Computational Insights into the Photodynamics of 2-Hydroxyazobenzene*. Gulf Coast Undergraduate Research Symposium, Rice University, Houston, TX, 2024.
- [5] **Hassan, U. S.**; Mehmood, A.; and Levine, B. G. *Computational Insights into the Photodynamics of 2-Hydroxyazobenzene*. Gulf Coast Undergraduate Research Symposium, Rice University, Houston, TX, 2024.
- [3] **Hassan, U. S.**; Trice, R.; and Law, S. *The Potential of Dirac Semimetals for Infrared Photodetection*. Penn State Materials Research Institute 2D Materials REU Talks, State College, PA, 2023.
- [2] **Hassan, U. S.** and Topper, R. Q. *Computational Analysis of Mass Spectra and Growth Patterns of Ammonium Nitrate Nanoparticles*. New York Chapter of ACS Undergraduate Symposium, LaGuardia Community College, NY, 2023.
- [1] **Hassan, U. S.**; Vazquez, J. M.; and Beaudoin, S. *Adhesion of Mock Polymer-Bonded Explosives*. Purdue Energetics Research Symposium, West Lafayette, IN, 2022.

## POSTER PRESENTATIONS

- [6] **Hassan, U. S.**; Mehmood, A.; and Levine, B. G. *Photochemistry of 2-Hydroxyazobenzene: A Computational Study*. Stony Brook Summer Research Symposium, Stony Brook, NY, 2024.
- [5] **Hassan, U. S.**; Amat, M. A.; and Topper, R. Q. *Decomposition and Growth Pathways of Aerosolized Ammonium Nitrate Particles*. American Conference on Theoretical Chemistry, Chapel Hill, NC, 2024.
- [4] **Hassan, U. S.** and Topper, R. Q. *Patterns in Growth of Ammonium Nitrate Clusters*. Virtual Winter School on Computational Chemistry, 2024.
- [3] **Hassan, U. S.**; Amat, M. A.; and Topper, R. Q. *Growth and Decomposition Pathways for Ammonium Nitrate Clusters*. AIChE Annual Conference, Orlando, FL, 2023.
- [2] **Hassan, U. S.**; Trice, R.; and Law, S. *Dirac Semimetals Potential in Infrared Photodetection*. Penn State REU Symposium, State College, PA, 2023.
- [1] **Hassan, U. S.**; Vazquez, J. M.; and Beaudoin, S. *Quantifying Van der Waals Adhesion of Energetic Particles*. Purdue Energetics Research Symposium, West Lafayette, IN, 2022.

## TECHNICAL SKILLS

**Computational:** Python, Bash, Slurm, High-Performance Computing, JAX-MD, Excel

**Quantum Chemistry Codes:** Quantum Espresso, ORCA, TeraChem, Spartan, OpenMolcas, Psi4

**Lab:** AFM, FTIR, UV-VIS, MS, XRD

## RELEVANT COURSEWORK

### Graduate level

Quantum Chemistry I, Statistical Mechanics I, Computational Chemistry, Mathematics for Chemistry

### Undergraduate

**Chemistry:** Statistical Mechanics, Physical Chemistry I & II, Organic Chemistry I & II, Organometallic Chemistry, Biochemistry

**Mathematics:** Linear Algebra, Abstract Algebra, Real Analysis I & II, Ordinary and Partial Differential Equations, Discrete Math, Probability, Vector Calculus

**Physics:** Quantum Mechanics, Electricity and Magnetism, Optics and Modern Physics

**Engineering:** Material Science, Chemical Reaction Engineering, Heat and Mass Transfer, Fluid Dynamics, Separation Principles, Chemical Process Evaluation and Design I & II, Chemical Process Simulation, Chemical Dynamics & Control