

# Srimukh Prasad Veccham Krishna Prasad

Research Interests: Development of theoretical models and computational techniques and applying them to chemical systems

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## EDUCATION

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- 2015 – 2020 **University of California Berkeley**, Berkeley, CA  
(expected) Candidate for Ph.D in Physical Chemistry | GPA: 3.95/4  
Advisor: Prof. Martin Head-Gordon
- 2010 – 2015 **Indian Institute of Technology Bombay**, Mumbai, India  
Masters of Science in Chemistry | GPA: 9.29/10  
Minor in Electrical Engineering  
Ranked #1 in Department of Chemistry

## CURRENT RESEARCH PROJECTS

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### Polarized Many-Body Expansion

*Central computation engine implemented in C++ with end-to-end Python wrappers for creation, execution, and collection of computation processes on high performance computing architectures*

- Formulated a computationally inexpensive technique to compute molecular energies and properties using Density Functional Theory (DFT)
- Formal complexity reduced from  $\mathcal{O}(N^3)$  to  $\mathcal{O}(N^2)$ , providing speedups of 10 to 100 times
- Algorithm is “embarrassingly parallelizable,” practically achieving perfect scaling upto 64 processors
- Demonstrated the success of this technique in over 50 molecular systems with errors
- Currently working on extensions to correlated wavefunction methods

### Performance of Density Functionals for Hydrogen Storage

*Implemented Python wrappers for spawning jobs and collecting data using Q-Chem, MRCC, and Psi4. Used Pandas for data analysis and interpretation*

- Compiled a dataset of 195 data points to accurately represent the hydrogen storage problem in porous materials
- Assessed the performance of over 55 density functionals against the reference binding energies
- Identified density functionals which can predict hydrogen binding energies with about 3% error
- Reduced computational cost by upto 1000 times in predicting binding energies

### Variational Decomposition of Pair-wise Charge Transfer

*Implemented numerical quadrature algorithms and iterative algorithms for non-linear objective function optimization in C++ using Armadillo*

- Implemented a variational alternative to the perturbative pair-wise charge decomposition in Energy Decomposition Analysis scheme
- Analyzed and carefully characterized the physical interpretation of each of the charge transfer terms such that charge transfer is comprehensively categorized

## PUBLICATIONS

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- Srimukh Prasad Veccham, Joonho Lee, Martin Head-Gordon “Making Many-Body Interactions Nearly Pairwise Additive: The Polarized Many-Body Expansion Approach” *J. Chem. Phys.* 2019, 151, 194101
- Srimukh Prasad Veccham, Martin Head-Gordon “Density Functionals for Hydrogen Storage: Defining the H2Bind195 Test Set with Ab Initio Benchmarks and Assessment of 55 Functionals” *Submitted to Journal of Chemical Theory and Computation*
- Srimukh Prasad Veccham, Joonho Lee, Martin Head-Gordon “Variational Charge-Transfer Analysis in ALMO-based Energy Decomposition Analysis” *Manuscript in preparation*
- Mark D. Allendorf, Zeric Hulvey, Thomas Gennett, ... Srimukh Prasad Veccham, Brandon C. Wood “An assessment of strategies for the development of solid-state adsorbents for vehicular hydrogen storage” *Energy & Environmental Science* 2018, 11, 2784

- E. Tsvion, **Srimukh Prasad Veccham**, Martin Head-Gordon "High-Temperature Hydrogen Storage of Multiple Molecules: Theoretical Insights from Metalated Catechols" *Chem. Phys. Chem.* 2017, 18, 184
- Sajesh P. Thomas, **Srimukh Prasad Veccham**, Louis J. Farrugia, Tayur N. Guru Row " 'Conformational Simulation' of Sulfamethizole by Molecular Complexation and Insights from Charge Density Analysis: Role of Intramolecular S...O Chalcogen Bonding" *Crystal Growth & Design* 2015, 15, 2110

## CONFERENCE PRESENTATIONS

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|---|-------------|
| <b>Polarized Many-body Expansion: A Perfect Marriage between Embedded Mean-field Theory and Variational Many-body Expansion</b> | April, 2019 |
| Oral presentation at the Spring National Meeting of the American Chemical Society, Orlando, FL                                  |             |
| <b>Assessment of the Performance of Density Functionals for Hydrogen Storage in Sorbents</b>                                    | April, 2019 |
| Oral presentation at the Spring National Meeting of the American Chemical Society, Orlando, FL                                  |             |
| <b>Polarized Many-Body Expansion</b>  | May, 2019   |
| Poster presentation at the Northern California Theoretical Chemistry Meeting, Berkeley, CA                                      |             |

## AWARDS AND HONORS

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| <b>Institute Silver Medal</b> for highest GPA in Department of Chemistry, IIT Bombay       | August, 2015           |
| <b>Young researcher</b> at 65 <sup>th</sup> Lindau Nobel Laureate Meeting, Lindau, Germany | June, 2015             |
| <b>Institute Academic Prize</b> for highest GPA in each semester for 6 semesters           | July, 2011 – May, 2014 |
| <b>Summer Undergraduate Research Fellowship</b> awarded Duke University                    | May, 2013              |
| <b>Summer Research Fellowship</b> awarded by Indian Academy of Sciences                    | May, 2012              |
| <b>INSPIRE fellowship</b> by Department of Science & Technology, Government of India       | July, 2010 – May, 2015 |

## SKILLS

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| SCIENTIFIC PROGRAMMING: | Extensive production-level programming experience in Q-Chem, a comprehensive electronic structure package. Contributed to embedding and energy decomposition analysis libraries. |
| SOFTWARE AND PACKAGES:  | Psi4, MRCC, Gaussian, NAMD, GROMACS, VMD, NBO, ChemOffice, $\text{\LaTeX}$   |
| PROGRAMMING LANGUAGES:  | C++ (Armadillo), Python (NumPy, SciPy, Pandas, Matplotlib)   |

## RELEVANT COURSEWORK

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- **Computer science:** Machine Learning, Machine Learning in Chemistry and Biology, Computer Programming and Utilization
- **Mathematics/Statistics:** Vector Calculus, Linear Algebra, Differential Equations, Data Analysis and Interpretation, Probability and Random Processes
- **Numerical methods:** Advanced Simulation Techniques in Physics, Numerical Solutions to Differential Equations (audit), Advanced Matrix Computations (audit)
- **Chemistry:** Advanced Quantum Mechanics I, Advanced Quantum Mechanics II, Thermodynamics and Statistical Mechanics, Chemical Kinetics

## TEACHING EXPERIENCE

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| <b>Upper Division Physical Chemistry</b>   | Spring 2017, UC Berkeley    |
| <ul style="list-style-type: none"> <li>• Lead discussion sections and formulated and graded assignments and exams</li> </ul>   |                             |
| <b>Freshman General Chemistry</b>  | Fall 2015, 2016 UC Berkeley |
| <ul style="list-style-type: none"> <li>• Lead a laboratory section, solved tutorial problems and lead discussion sections</li> <li>• Graded laboratory reports, assignments, quizzes, and exams with feedback</li> </ul> |                             |
| <b>Freshman Physical Chemistry</b>   | Spring 2014, IIT Bombay     |
| <ul style="list-style-type: none"> <li>• Lead discussion section, solved tutorial problems, and handled student questions</li> </ul>   |                             |