

Hari Sharma

307 Sunderland Dr. Mount Juliet, TN, 37122, USA

[\(+1\) \(860\)-975-7070](tel:+18609757070) | harisharma55@gmail.com | [linkedin](https://www.linkedin.com/in/harisharma55/) | [Google Scholar](https://scholar.google.com/citations?user=...)

Summary

- Organized and detail-oriented computational scientist skilled at working on multiple projects and collaborating to drive scientific efforts to successful outcomes.
- Expertise in quantitative and statistical analysis on complex datasets.
- Excellent communication skills and proven ability to thrive in a fast-paced environment.

Skills

Technical	Machine Learning (supervised & unsupervised) Deep Learning Natural Language Processing (NLP) Exploratory Data Analysis (EDA) Cloud Computing with Amazon Web Service (AWS) Big Data with PySpark
Programming	Python (Experienced) SQL BASH, C++ (Working knowledge) TCL/TK (Familiar) \LaTeX
Frameworks & Libraries	Jupyter Matplotlib Numpy Pandas Scikit-learn TensorFlow NLTK GIT Linux
Languages	Nepali(Native), English (Fluent), Hindi(Understand)

Experience

StemSynergy Therapeutics INC.

Franklin, TN, USA

SENIOR SCIENTIST

Oct. 2021 - Present

- Lead multiple projects involving the computational aspects of drug discovery using state-of-the art technologies.
- Collaborated with experimental chemist and biologist and contributed towards the design of future experiments.
- Analyzed the data (Gigabytes) obtained from large scale molecular dynamics (MD) simulations using self written python scripts.
- Wrote BASH shell scripts for performing fully automated molecular dynamics simulation using GROMACS software suite.
- Administered linux computer cluster nodes with GPU's and CPU's (4 nodes), installed open source molecular modeling simulation softwares relevant for drug discovery and lead computational part of the research.

Vanderbilt University

Nashville, TN, USA

POSTDOC SCHOLAR

Sept. 2019 - Sept. 2021

- Contributed towards the ligand docking algorithm in ROSETTA (C++ codebase) macromolecular modeling suite to carry out docking of ensembles of congeneric series of ligand simultaneously into protein and their mutant variants.
- Communicated results to collaborators and provided recommendations for future experiments.
- Mentored rotating graduate student.

University of Connecticut

Storrs, CT, USA

GRADUATE RESEARCH ASSISTANT

Aug 2012 - Aug 2019

- Assembled and administered linux computer cluster nodes with installation of open source molecular modeling software suites.
- Wrote python scripts from scratch to analyze the gigabytes of data obtained from MD simulations.
- Implemented an algorithm using Graph theory and matrix algebra that is able to compute the end points of a cloud of data points oriented in a flexible contour where data points are dynamic.

Education

University of Connecticut

Storrs, CT, USA

PHD IN PHYSICS

Jan 2015 - Aug. 2019

- Dissertation Title: Computer Simulations of Amphiphile Based Hybrid Nanostructures.

University of Connecticut

Storrs, CT, USA

M. S IN PHYSICS

Aug 2012 - Dec. 2014

Tribhuvan University

Kirtipur, Nepal

M. S IN PHYSICS

May 2008 - May 2010