

Tanmoy Sanyal

Scientist, Protein Design || Structural biology

About

📍 San Francisco Bay Area, CA
✉ tanmoy.7989@gmail.com
📞 (805) 637-0375
💻 tanmoy7989.github.io
🐦 twitter.com/hiddenvariable2
🔗 tanmoy7989

Protein Design

Rosetta
AlphaFold
RFDiffusion
ProteinMPNN

Molecular modeling

openMM
LAMMPS
GROMACS
IMP
Modeller
UCSF ChimeraX
PyMOL

ML frameworks

PyTorch
Tensorflow-probability
PyMC3

Languages

Python (+Cython) (~20000+ lines)
Fortran-90 (~1000+ lines)
C++ (~4000+ lines)
Bash
TCL

Open Source

nbSPIKE
Integrative epitope prediction

IMP
Crosslink guided domain
discovery

PDB-Dev deposition
Structure of SMC5/6

LAMMPS
Local density potential

Education

University of California Santa Barbara
Santa Barbara, CA, USA

2013 - 2018

Ph.D Chemical Engineering
Graduate emphasis in Computational Science & Engineering (Parallel Computing & Numerical methods for ODEs / PDEs)

Indian Institute of Technology Kharagpur
Kharagpur, WB, India

2008 - 2013

Masters and Bachelors of Technology (Hons.) integrated dual degree Chemical Engineering

Research Experience

Amgen

2023 - Present

Senior protein design scientist

Novo Nordisk Research Center Seattle

2022 - 2023

Protein design scientist

- **Biologics design:** (i) Design of mini-proteins and antibodies for multiple therapy areas using Rosetta and AlphaFold / RFDiffusion / ProteinMPNN pipelines (ii) SAR analysis based on protein language models. (iii) Optimizing non-canonical modifications to peptides for half-life protraction.
- **Scientific software development:** Worked closely with AWS to re-tool AlphaFold and deploy a faster cloud-native in-house version for robust and scalable (~100k sequences) structure prediction; regularly applied now in variant design within the company.
- **Leadership:** (i) Program lead for refining protein interface prediction ML models in collaboration with the Institute for Protein Design at UW, Seattle. (ii) Led and evaluated a proof of concept for onboarding virtual reality based structural rendering software among collaborative teams of computational design scientists and synthetic chemists.

Department of Bioengineering, University of California San Francisco

2019 - 2022

Postdoctoral scholar, Andrej Sali lab

- **Protein structure modeling using chemical crosslinks:** Homology models + chemical-crosslinks + cryo-EM data for protein complex structure modeling; graph sampling methods to extract domain information from crosslinks. Applications include SMC and DNA helicase protein families.
- **Nanobody biophysics:** Developed integrative docking scores for nanobody docking on target receptors (SARS-CoV-2 and variant spike proteins) using chemical crosslinking and escape mutation data. Developed design principles for increasing nanobody thermostability.
- **Whole cell models:** Developed Bayesian Network models to combine continuum scale (pharmacokinetic), particle scale (Brownian dynamics) and network scale (enzyme pathways) models of the pancreatic beta-cell into build a proof-of-concept multimodal digital twin. Served as software lead within the Pancreatic Beta-Cell consortium (collaborating across UCSF, USC, ShanghaiTech and HUJI)

PhD candidate, Shell lab

Research Interests

Protein language models

Integrative structure modeling

Non-canonical amino acids

Nanobodies

Forcefield development

- **Coarse-graining algorithms for bulk liquids:** Introduced the local density potential for structurally and thermodynamically accurate coarse-grained MD simulations of liquid-liquid phase separation, which are relatively rare in the chemical engineering literature
- **Coarse-graining algorithms for protein folding using variational inference techniques:** Developed protein backbone models for (i) template-free folding of 200+ residue coarse-grained protein domains using MD simulations, and (ii) studying self-assembly in amyloidogenic peptides
- **High Performance Computing:** As system administrator, set up 144 core Linux (Rocks 6.2) cluster and wrote utilities for automating (i) network-attached-storage management, and (ii) conflict-free compute resource sharing between lab members.

Volunteer

Peer-review

2019 - Present

Structure, Journal of Physical Chemistry, Proteins: Structure, Function and Bioinformatics, Review of Scientific Instruments, Rapid Reviews Covid-19, Life

Graduate Simulation Seminar Series, University of California Santa Barbara

2014 - 2016

Co-founder

Conception, organization, public outreach and long-term funding support from ChE, Materials Science and Computer Science departments.

Selected Publications

Highly synergistic combinations of nanobodies that target SARS-CoV-2 and are resistant to escape
F.D. Mast, P.C. Fridy, N.E. Karen, J. Wang, E.Y. Jacobs, J.P. Olivier, T. Sanyal, et al, *eLife*, 2021

Bayesian metamodeling of complex biological systems across varying representations
B. Raveh, L. Sun, K.L. White, T. Sanyal, et al, *PNAS*, 2021

Integrative analysis reveals unique structural and functional features of the Smc5/6 complex
F.D. Mast, P.C. Fridy, N.E. Karen, J. Wang, E.Y. Jacobs, J.P. Olivier, T. Sanyal, et al, *PNAS*, 2021

A hybrid, bottom up, structurally-accurate, Go-like coarse-grained protein model
T. Sanyal, J. Mittal & M. Scott Shell, *JCP*, 2019

Transferable coarse-grained models of liquid-liquid equilibrium using local density potentials optimized with the relative entropy
T. Sanyal & M. Scott Shell, *JPC-B*, 2018

Coarse-grained models using local-density potentials optimized with the relative entropy:
Application to implicit solvation
T. Sanyal & M. Scott Shell, *JCP*, 2016

Recent conference presentations

7th Annual Computational Drug Development for Biologics Summit, Boston, MA, USA
(Talk): *Unpacking a workflow for automating protractor placement on peptides for enhanced half life*, Oct-2022

Indian Biophysical Society Meeting, Navi Mumbai, India
(Talk): *Integrative modeling of higher order complexes from binary binding modes*, Apr-2022

Biophysical Society Annual Meeting, San Francisco, CA, USA
(Talk): *Rigid body assignment in integrative determination of protein complex structures*, Feb-2022

PDB50: A special symposium celebrating the 50th anniversary of the Protein Data Bank, (virtual)
(Poster): *Integrative modeling of the SMC5/6 complex*, May-2021

USC Brige Institute - Pancreatic Beta Cell Consortium Retreat, Catalina Island, CA, USA
(Workshop): *Meta-modeling: Probabilistic graphical models and your data*, Nov-2019

Berkeley Mini Stat Mech Meeting, Berkeley, CA, USA
(Poster): *Coarse-grained models for protein folding and self-assembly with the relative entropy*, Jan-2019

Awards

IUPAB best oral presentation

2022

Technopreneurship Promotion Programme

2009

Ministry of Science and Technology, Govt. of India

Seed money for prototype scale-up of bioreactors for algal biodiesel generation

Jagdis Bose National Science Talent Search (JBNSTS) fellowship

2009