# Tanmoy Sanyal

Scientist, Computational drug design & Structural biology

#### About

Greater Seattle area, WA

(805) 637-0375

tanmoy7989.github.io

twitter.com/hiddenvariable2

**n** tanmoy7989

# Molecular modeling

Rosetta openMM LAMMPS GROMACS IMP Modeller

UCSF ChimeraX

PyMOL VMD

# ML frameworks

PyTorch Tensorflow Tensorflow-probability PyMC3

#### Workflows

Snakemake AWS Batch

#### Languages

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

#### Open Source

LAMMPS Local density potential

nbspike
Integrative epitope prediction

IMP Crosslink guided domain discovery

PDB-Dev deposition Structure of SMC5/6

### Education

# University of California Santa Barbara

Santa Babara, CA, USA

Ph.D Chemical Engineering

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

# Indian Institute of Technology Kharagpur

2008 - 2013

2013 - 2018

Kharagpur, WB, India

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

# Research Experience

#### **Novo Nordisk Research Center Seattle**

2022 - Present

Scientist, Computational drug discovery

- **Biologics Design**: Design of peptides and antibodies for rare endocrine disorders, and SAR (structure activity relationship) analysis based on PLMs (protein language models)
- **Scientific Software development**: Computational workflows for automating molecular model building and MD simulation of modified peptides, such as small molecule acylations or peptide-biopolymer conjugates.

# Department of Bioengineering, University of California San Francisco

2019 - 2022

Postdoctoral scholar, Andrej Sali lab

- **Protein structure modeling using chemical crosslinks**: Combined 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.
- **Computational mapping of nanobody epitopes**: Developed software to integrate chemical crosslinks with escape mutation data for protein-protein docking. Applications include epitope prediction of nanobodies on wild-type SARS-CoV-2 spike and variants.
- 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 digitial twin Served as software lead within the Pancreatic Beta-Cell consortium (collaborating across UCSF, USC, ShanghaiTech and HUJI)

# Chemical Engineering, University of California Santa Barbara

2013 - 2018

PhD candidate, Shell lab

- **Thesis**: Next-Generation Coarse-Grained Models for Molecular Dynamics Simulations of Fluid Phase Equilibria and Protein Biophysics Using the Relative Entropy
- 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: 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.

#### Research Interests

Volunteer

Molecular Sciences

Protein design

2019 - Present Peer-review

Structure, Review of Scientific Instruments, Rapid Reviews Covid-19, Life, Biomolecules, International Journal of

Forcefield development

Amyloidogenic self assembly

Whole cell models

Bayesian networks

Variational inference

Graduate Simulation Seminar Series, University of California Santa Barbara

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, (virtiual) (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 orgal presentation

44th Indian Biophysical Society Meeting

2022

#### **Technopreneurship Promotion Programme**

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

2009

2014 - 2016