

Research Scientist: Molecular Dynamics

(Actual position will be determined based on qualifications)

UST Corporation, a biotech company, headquartered in Greater Boston with a branch in San Diego, is looking for an outstanding computational biophysicist to join our R&D team in Boston. He/she will lead the company's efforts at molecular modeling and simulation of biomolecular systems. A successful candidate should have at least four years hands-on experience on computational structural dynamics of proteins and their complexes using different molecular simulation techniques. This position is a unique opportunity to work on the development of a very advanced DNA sequencing system.

Specific Requirements

Desired Skills:

- Strong background in biophysics, biochemistry and computational simulation as well as statistics, thermodynamics, quantum mechanics (QM) and hybrid QM/molecular mechanics (MM)
- Extensive knowledge of DNA and protein structures, including X-ray and NMR techniques
- Hands-on experience on molecular dynamics simulation with expertise in commonly used MD force fields, and coarse-grained elastic network models for sampling molecular intrinsic motions
- Hands-on experience on one or more common software package, such as MAPS, Abalone, Ascalaph Designer, LAMMPS, HyperChem, Discovery Studio, Culgi, COSMOS.
- Hands-on programming and scripting experience with C/C++/C#, Java, Python, etc.
- Strong analytical and quantitative skills, able to analyze experimental data using statistical and machine learning methods and compare with numerical simulation data
- Cloud Computing experience with AWS, GCP and Azure a plus

Duties and Responsibilities:

- Perform numerical modeling and simulation of biopolymers using a commercially available MD simulation software package, interacting with external experts and internal software group for the design, development and improvement of simulation code
- Work with experimentalists and product manager for the optimization of product and process designs, providing the simulation data for the development of products
- Work with software group for data acquisition and analysis software development
- Help scientists and engineers troubleshooting and resolving complex problems using numerical methods, improving their experiment design and output
- Other work/project assignments as required

Qualifications:

- Advanced degree in biophysics or computational physics/chemistry or related scientific discipline; PhD desired.
- Minimum 4 years of substantial experience in molecular dynamics (MD) and Monte Carlo (MC) simulation of biomolecules, preferably proteins and

their complexes, familiar with common molecular dynamics simulation software packages

- **Excellent oral and written communication skills and a productive team player.**

Interested candidates please submit your updated resume with a cover letter as well as a list of at least three references to info@ustcorporation.com