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Advances in Computer-Aided Drug Design and Molecular Dynamics Simulations

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Deadline for manuscript submissions: **28 February 2025**



With the rapid advancement of science and technology, computer-aided drug design (CADD) and molecular dynamics (MD) simulations have become increasingly prominent in drug development.

This Special Issue aims to compile the latest research findings related to CADD and MD simulations, particularly their innovative applications in biomolecular crystal studies. We welcome submissions in the following areas:

Computer-Aided Drug Design: Including structure-based drug design, virtual screening, quantitative structureactivity relationship (QSAR) models, etc.

Molecular Dynamics Simulations: MD simulations, free energy calculations, and large-scale molecular simulations et.al.

Biomolecular Crystal Studies: Involving the analysis of biomacromolecular crystal structures, growth mechanisms, and the interactions between drugs and biomolecular crystals.

Multiscale Simulations and Integrated Methods: Combining quantum mechanics, molecular mechanics, and coarsegrained models et.al.

Case Studies and Applications: Demonstrating the successful applications of CADD and MD simulations in actual drug development, including the design and optimization of anticancer, antiviral, and antibacterial drugs et.al.

Specialsue



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Message from the Editor-in-Chief

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