Topical Collection

Computational Studies of Biomolecules

Message from the Collection Editors

The current Topical Collection aims to attract highquality contributions of modeling biomolecular structures, dynamics, functions, and interactions with the potential of interpretation of experimental data and applications in drug design and protein design. Topics of interest:

- Development and validation of new computational modeling methods;
- Computational studies of proteins' structure–function relationships;
- Computational investigations of nucleic acids' structure-function relationships;
- Modeling of protein and nucleic acid dynamics;
- Protein docking;
- Protein-ligand interactions;
- Nucleic acid-ligand interactions;
- Protein design:
- Computational enzymology-enzymatic reaction mechanisms;
- Protein homology modeling.

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Message from the Editor-in-Chief

The International Journal of Molecular Sciences (*IJMS*, ISSN 1422-0067) is an open access journal, which was established in 2000. The journal aims to provide a forum for scholarly research on a range of topics, including biochemistry, molecular and cell biology, molecular biophysics, molecular medicine, and all aspects of molecular research in chemistry. *IJMS* publishes both original research and review articles, and regularly publishes special issues to highlight advances at the cutting edge of research. We invite you to read recent articles published in *IJMS* and consider publishing your next paper with us.

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