



## **Advances in Computational Biology: Molecular Dynamics Simulation of Biomolecules**

Guest Editor:

**Dr. Paulino Gómez-Puertas**

Molecular Modeling Group,  
Centro de Biología Molecular  
Severo Ochoa (CBM, CSIC-UAM),  
CL Nicolas Cabrera, 1. Campus  
UAM, 28049 Madrid, Spain

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### **Message from the Guest Editor**

Dear Colleagues,

The simulation of important processes in biomedicine and biotechnology using computational methods has become a well-established field in recent years. We can simulate motions associated with the function of key macromolecules in cellular processes. Using molecular dynamics systems, we can study the effect of a mutation that causes a rare disease or a carcinogenic process on the structure and/or function of a macromolecule. We can also design molecules with the ability to act as antimicrobial, antiviral or anticancer agents. All these processes, as well as the methodology that makes them possible, are welcome in this Special Issue of *IJMS*.

Dr. Paulino Gómez-Puertas  
*Guest Editor*





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### Prof. Dr. Maurizio Battino

Department of  
Odontostomatologic and  
Specialized Clinical Sciences,  
Sez-Biochimica, Faculty of  
Medicine, Università Politecnica  
delle Marche, Via Ranieri 65,  
60100 Ancona, Italy

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*International Journal of Molecular  
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MDPI, Grosspeteranlage 5  
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