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Simulation of Crystallization in Complex Atomic and Particulate Systems

Guest Editor:

Message from the Guest Editor

Dear Colleagues,

Deadline for manuscript submissions:

closed (20 August 2020)

The purpose of the present Special Issue is to host the most recent advances in the computer simulation of crystal nucleation and growth with the intention of highlighting the similarities and differences of the phenomenon as it spans highly diversified atomic and particulate systems. Accordingly, this Special Issue is open to simulation works crossing interdisciplinary fields and coming from different research backgrounds.

Given the focus on the simulation aspects of crystallization, this Special Issue also accepts manuscripts detailing developments in related algorithms and methodologies.

Prof. Dr. Nikos Ch. Karayiannis Guest Editor







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Editor-in-Chief

Prof. Dr. Alessandra Toncelli Department of Physics, University of Pisa, 56126 Pisa, Pl, Italy

Message from the Editor-in-Chief

Welcome to *Crystals*, the journal dedicated to the fascinating world of crystallographic research! Crystals are more than mere decorative elements; they hold the key to understanding the fundamental structure of matter. Our mission is to explore the crucial significance of this research across various fields. From medicine to technology, chemistry to geology, crystals play a vital role. Their structure provides insights into new advanced materials, innovative drugs, and groundbreaking technologies. Through *Crystals*, we delve into the microscopic world to discover solutions that will shape the future. Join us on a journey through the *Crystals*, where science merges with beauty and innovation.

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