

Special Issue

Modelling Materials and Devices at Atomistic Level

Message from the Guest Editor

In the past few years, atomistic simulations have become essential tools for exploring new materials and investigating the working principles of novel electronic devices. The success of ab initio methods based on the density functional theory (DFT) ensures valuable guidance for the synthesis of proper candidates. Complementary, molecular dynamics simulations are able to cover mechanical properties at significantly larger timescales and system sizes. In addition, in recent years, machine learning (ML) techniques combined with atomistic descriptions have boosted the search for new materials and the optimization of device properties. This Special Issue shall cover topics concerning the modeling of materials and devices at atomistic level. These shall be focused but not limited to materials and interfaces for opto-electronic applications, spin and charge transport in atomic-scale devices, and novel ML approaches coupled to ab initio approaches.

- density functional theory
- electronic structure calculations
- molecular dynamics
- machine learning
- spin and charge transport
- atomic-scale devices
- ab initio modeling of interfaces
- molecular dynamics simulations

Guest Editor

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About the Journal

Message from the Editor-in-Chief

Materials (ISSN 1996-1944) was launched in 2008. The journal covers twenty-five comprehensive topics: biomaterials, energy materials, advanced composites, advanced materials characterization, porous materials, manufacturing processes and systems, advanced nanomaterials and nanotechnology, smart materials, thin films and interfaces, catalytic materials, carbon materials, materials chemistry, materials physics, optics and photonics, corrosion, construction and building materials, materials simulation and design, electronic materials, advanced and functional ceramics and glasses, metals and alloys, soft matter, polymeric materials, quantum materials, mechanics of materials, green materials, general. *Materials* provides a unique opportunity to contribute high quality articles and to take advantage of its large readership.

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