



Simulation of Abrasive Processes for Crystalline Materials by Means of Molecular Dynamics and Multiscale Methods

Guest Editors:

Dr. Nikolaos Karkalos

School of Mechanical Engineering, National Technical University of Athens, 15780 Zografou, Greece

Prof. Dr. Umberto Prisco

Department of Chemical, Materials and Production Engineering, University of Napoli Federico II, piazzale Tecchio, 80, 80125 Napoli, Italy

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Message from the Guest Editors

Abrasive processes are essential in manufacturing in order to render high-quality surfaces of mechanical parts. Given that experiments at the micro and nanoscale are very expensive, simulation methods can be employed in order to evaluate the capabilities of abrasive processes in these scales, and optimize them with respect to various materials. At the nanoscale, the molecular dynamics method is the most frequently used method to achieve reliable simulation of material removal.

This Special Issue plans to present an overview of the most recent advances in the field of simulations of abrasive processes at the nanoscale by means of molecular dynamics and multiscale methods, including:

- Simulation of abrasive processes using molecular dynamics;
- Meshless methods for the simulation of abrasive processes;
- Development of multiscale frameworks for abrasive process simulations;
- Prediction of microstructure during abrasive processes;
- Simulation of hybrid abrasive processes;
- Simulation of abrasive processes for nanostructured materials;
- Simulation of abrasive processes for high-entropy alloys;
- Simulation of abrasive processes for polycrystalline materials.





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Editor-in-Chief

Prof. Dr. Alessandra Toncelli

Department of Physics, University
of Pisa, 56126 Pisa, Italy

Message from the Editor-in-Chief

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Crystals Editorial Office
MDPI, Grosspeteranlage 5
4052 Basel, Switzerland

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