

Structures and Properties of Ionic Liquids examined by Molecular Dynamics Simulations

Collection Editor: Junko Habasaki

Description

Some newly-developed ionic salts are liquids even at room temperature, and are now called the ionic liquids (ILs). These systems are designable to exhibit new functions by combinations of complex cations and anions. Molecular Dynamics simulations are a useful tool to examine such systems. The purpose of this collection is to examine the structures and properties of ionic liquids using MD and related methods.

About the Editor

Dr. Junko Habasaki graduated from Tokyo University of Education, Japan, and took her PhD at Tokyo Metropolitan University, Japan. She has published more than 100 papers in international journals and has done more than 40 invited (keynote, plenary) talks at international conferences. She is a co-author of *Dynamics of Glassy, Crystalline and Liquid Ionic Conductors* (Springer, 2017) and is the author of *Molecular Dynamics of Nanostructures and Nanoionics* (Jenny Stanford, 2020).

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