

Oleg N Starovoytov

List of Publications by Year in descending order

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Version: 2024-02-01

10
papers

229
citations

1163117

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1372567

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10
times ranked

374
citing authors

| # | ARTICLE | IF | CITATIONS |
|----|---|-----|-----------|
| 1 | In situ study on the compression deformation of MoNbTaVW high-entropy alloy. Journal of Alloys and Compounds, 2021, 871, 159557. | 5.5 | 7 |
| 2 | Development of a Polarizable Force Field for Molecular Dynamics Simulations of Lithium-Ion Battery Electrolytes: Sulfone-Based Solvents and Lithium Salts. Journal of Physical Chemistry B, 2021, 125, 11242-11255. | 2.6 | 9 |
| 3 | Induced polarization restricts the conformational distribution of a light-harvesting molecular triad in the ground state. Physical Chemistry Chemical Physics, 2017, 19, 22969-22980. | 2.8 | 10 |
| 4 | Development of an AMOEBA water model using GEM distributed multipoles. Theoretical Chemistry Accounts, 2015, 134, 1. | 1.4 | 20 |
| 5 | GEM*: A Molecular Electronic Density-Based Force Field for Molecular Dynamics Simulations. Journal of Chemical Theory and Computation, 2014, 10, 1361-1365. | 5.3 | 64 |
| 6 | Development of AMOEBA Force Field for 1,3-Dimethylimidazolium Based Ionic Liquids. Journal of Physical Chemistry B, 2014, 118, 7156-7166. | 2.6 | 40 |
| 7 | Effects of the Hydroxyl Group on Phenyl Based Ligand/ERR ¹³ Protein Binding. Chemical Research in Toxicology, 2014, 27, 1371-1379. | 3.3 | 9 |
| 8 | Molecular dynamics simulation studies of the influence of imidazolium structure on the properties of imidazolium/azide ionic liquids. Journal of Chemical Physics, 2012, 136, 194506. | 3.0 | 16 |
| 9 | Development of a Polarizable Force Field for Molecular Dynamics Simulations of Poly (Ethylene) Tj ETQq1 1 0.784314 rgBT /Overlock | 5.3 | 39 |
| 10 | Dissolution behavior of silver in ammoniacal solutions using bromine, iodine and hydrogen-peroxide as oxidants. Hydrometallurgy, 2007, 86, 114-119. | 4.3 | 15 |