

Dmitry V Ivlev

List of Publications by Year in descending order

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14
papers

180
citations

1163117

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h-index

1058476

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all docs

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docs citations

14
times ranked

150
citing authors

| # | ARTICLE | IF | CITATIONS |
|----|--|-----|-----------|
| 1 | Computation of drug solvation free energy in supercritical CO ₂ : Alternatives to all-atom computer simulations. <i>Fluid Phase Equilibria</i> , 2021, 544-545, 113096. | 2.5 | 1 |
| 2 | Computation of the Solubility of 1,2,4-Thiadiazole Derivatives in Supercritical Carbon Dioxide. <i>Russian Journal of Physical Chemistry B</i> , 2021, 15, 1166-1170. | 1.3 | 2 |
| 3 | Carbamazepine solubility in supercritical CO ₂ : A comprehensive study. <i>Journal of Molecular Liquids</i> , 2020, 311, 113104. | 4.9 | 20 |
| 4 | Conformational equilibria of pharmaceuticals in supercritical CO ₂ , IR spectroscopy and quantum chemical calculations. <i>Spectrochimica Acta - Part A: Molecular and Biomolecular Spectroscopy</i> , 2020, 230, 118072. | 3.9 | 18 |
| 5 | High-pressure NMR spectroscopy in studies of the conformational composition of small molecules in supercritical carbon dioxide. <i>Journal of Molecular Liquids</i> , 2020, 309, 113113. | 4.9 | 23 |
| 6 | Possibility of pressure crossover prediction by classical DFT for sparingly dissolved compounds in scCO ₂ . <i>Journal of Molecular Liquids</i> , 2019, 276, 801-805. | 4.9 | 12 |
| 7 | Screening of conformational polymorphism of ibuprofen in supercritical CO ₂ . <i>Journal of Molecular Liquids</i> , 2017, 239, 49-60. | 4.9 | 30 |
| 8 | Conformational lability of ibuprofen in supercritical carbon dioxide. <i>Russian Journal of Physical Chemistry B</i> , 2016, 10, 1153-1162. | 1.3 | 5 |
| 9 | Helium diffusion in aqueous sodium chloride solution at high pressures. <i>Russian Journal of Physical Chemistry A</i> , 2012, 86, 974-978. | 0.6 | 1 |
| 10 | Topology of hydrogen-bonded clusters in sub- and supercritical n-butanol. <i>Molecular dynamics simulation. Russian Journal of Physical Chemistry A</i> , 2010, 84, 2077-2081. | 0.6 | 6 |
| 11 | The volume characteristics and molecular dynamics simulation of nonaqueous solutions of aliphatic alcohols. <i>Russian Journal of Physical Chemistry A</i> , 2009, 83, 209-213. | 0.6 | 8 |
| 12 | Concentration dependence of the viscosity of tert-butanol-water mixtures: Physical experiment and computer simulation. <i>Russian Journal of General Chemistry</i> , 2004, 74, 1156-1162. | 0.8 | 10 |
| 13 | The study of hydrophobicity in water-methanol and water-tert-butanol mixtures. <i>Journal of Molecular Liquids</i> , 2004, 110, 193-199. | 4.9 | 29 |
| 14 | Preferential solvation and elasticity of the hydrogen bonds network in tertiary butyl alcohol-water mixture. <i>Chemical Physics Letters</i> , 2003, 379, 581-587. | 2.6 | 15 |