

Dmitry V Ivlev

List of Publications by Year in descending order

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14

papers

180

citations

1163117

8

h-index

1058476

14

g-index

14

all docs

14

docs citations

14

times ranked

150

citing authors

#	ARTICLE	IF	CITATIONS
1	Screening of conformational polymorphism of ibuprofen in supercritical CO ₂ . <i>Journal of Molecular Liquids</i> , 2017, 239, 49-60.	4.9	30
2	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
3	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
4	Carbamazepine solubility in supercritical CO ₂ : A comprehensive study. <i>Journal of Molecular Liquids</i> , 2020, 311, 113104.	4.9	20
5	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
6	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
7	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
8	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
9	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
10	Topology of hydrogen-bonded clusters in sub- and supercritical n-buthanol. Molecular dynamics simulation. <i>Russian Journal of Physical Chemistry A</i> , 2010, 84, 2077-2081.	0.6	6
11	Conformational lability of ibuprofen in supercritical carbon dioxide. <i>Russian Journal of Physical Chemistry B</i> , 2016, 10, 1153-1162.	1.3	5
12	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
13	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
14	Computation of drug solvation free energy in supercritical CO ₂ . xmlns:mml="http://www.w3.org/1998/Math/MathML" altimg="si3.svg"> $\text{Alternatives to all-atom computer simulations.}$ <i>Fluid Phase Equilibria</i> , 2021, 544-545, 113096.	2.5	1