

Shuzhe Wang

List of Publications by Year in descending order

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Version: 2024-02-01

11
papers

241
citations

1307594

7
h-index

1372567

10
g-index

11
all docs

11
docs citations

11
times ranked

290
citing authors

#	ARTICLE	IF	CITATIONS
1	Incorporating NOE-Derived Distances in Conformer Generation of Cyclic Peptides with Distance Geometry. <i>Journal of Chemical Information and Modeling</i> , 2022, , .	5.4	7
2	Evaluation of nine condensed-phase force fields of the GROMOS, CHARMM, OPLS, AMBER, and OpenFF families against experimental cross-solvation free energies. <i>Physical Chemistry Chemical Physics</i> , 2021, 23, 13055-13074.	2.8	9
3	Volume-scaled common nearest neighbor clustering algorithm with free-energy hierarchy. <i>Journal of Chemical Physics</i> , 2021, 154, 084106.	3.0	8
4	Passing the Barrier – How Computer Simulations Can Help to Understand and Improve the Passive Membrane Permeability of Cyclic Peptides. <i>Chimia</i> , 2021, 75, 518-521.	0.6	4
5	Effect of Flexibility, Lipophilicity, and the Location of Polar Residues on the Passive Membrane Permeability of a Series of Cyclic Decapeptides. <i>Journal of Medicinal Chemistry</i> , 2021, 64, 12761-12773.	6.4	22
6	Use of molecular dynamics fingerprints (MDFPs) in SAMPL6 octanol-water log P blind challenge. <i>Journal of Computer-Aided Molecular Design</i> , 2020, 34, 393-403.	2.9	19
7	Combining Machine Learning and Molecular Dynamics to Predict P-Glycoprotein Substrates. <i>Journal of Chemical Information and Modeling</i> , 2020, 60, 4730-4749.	5.4	30
8	Improving Conformer Generation for Small Rings and Macrocycles Based on Distance Geometry and Experimental Torsional-Angle Preferences. <i>Journal of Chemical Information and Modeling</i> , 2020, 60, 2044-2058.	5.4	82
9	Machine Learning with and for Molecular Dynamics Simulations. <i>Chimia</i> , 2019, 73, 1024.	0.6	5
10	Rationalization of the Membrane Permeability Differences in a Series of Analogue Cyclic Decapeptides. <i>Journal of Chemical Information and Modeling</i> , 2019, 59, 294-308.	5.4	55
11	Simulation Studies on the Lipid Interaction and Conformation of Novel Drug-Delivery Pseudopeptidic Polymers. <i>Journal of Physical Chemistry B</i> , 2017, 121, 9113-9125.	2.6	0