

Zhiye Tang

List of Publications by Year in descending order

Source: <https://exaly.com/author-pdf/4833575/publications.pdf>

Version: 2024-02-01

13
papers

296
citations

1040056

9
h-index

1125743

13
g-index

20
all docs

20
docs citations

20
times ranked

426
citing authors

#	ARTICLE	IF	CITATIONS
1	All-atomistic molecular dynamics study of the glass transition of amorphous polymers. <i>Polymer</i> , 2022, 254, 125044.	3.8	2
2	A comparison of the brittle PMMA with the ductile PC on the elasticity and yielding from a molecular dynamics perspective. <i>Polymer</i> , 2021, 226, 123809.	3.8	5
3	All-atom molecular dynamics study of impact fracture of glassy polymers. II: Microscopic origins of stresses in elasticity, yielding, and strain hardening. <i>Polymer</i> , 2020, 207, 122908.	3.8	19
4	Transient States and Barriers from Molecular Simulations and the Milestoning Theory: Kinetics in Ligand-Protein Recognition and Compound Design. <i>Journal of Chemical Theory and Computation</i> , 2020, 16, 1882-1895.	5.3	19
5	All-atom molecular dynamics study of impact fracture of glassy polymers. I: Molecular mechanism of brittleness of PMMA and ductility of PC. <i>Polymer</i> , 2019, 178, 121570.	3.8	34
6	Potential Mean Force from Umbrella Sampling Simulations: What Can We Learn and What Is Missed?. <i>Journal of Chemical Theory and Computation</i> , 2019, 15, 2433-2443.	5.3	96
7	Binding Thermodynamics and Kinetics Calculations Using Chemical Host and Guest: A Comprehensive Picture of Molecular Recognition. <i>Journal of Chemical Theory and Computation</i> , 2018, 14, 303-318.	5.3	42
8	An MM and QM Study of Biomimetic Catalysis of Diels-Alder Reactions Using Cyclodextrins. <i>Catalysts</i> , 2018, 8, 51.	3.5	5
9	A molecular dynamics investigation of CDK8/CycC and ligand binding: conformational flexibility and implication in drug discovery. <i>Journal of Computer-Aided Molecular Design</i> , 2018, 32, 671-685.	2.9	12
10	Adaptation of a Genetic Screen Reveals an Inhibitor for Mitochondrial Protein Import Component Tim44. <i>Journal of Biological Chemistry</i> , 2017, 292, 5429-5442.	3.4	18
11	Systematic Dissociation Pathway Searches Guided by Principal Component Modes. <i>Journal of Chemical Theory and Computation</i> , 2017, 13, 2230-2244.	5.3	7
12	Understanding ligand-receptor non-covalent binding kinetics using molecular modeling. <i>Frontiers in Bioscience - Landmark</i> , 2017, 22, 960-981.	3.0	18
13	Facet Selectivity of Ligands on Silver Nanoplates: Molecular Mechanics Study. <i>Journal of Physical Chemistry C</i> , 2014, 118, 21589-21598.	3.1	14