

Payal Chatterjee

List of Publications by Year in descending order

Source: <https://exaly.com/author-pdf/9394891/publications.pdf>

Version: 2024-02-01

13
papers

255
citations

840776

11
h-index

1125743

13
g-index

13
all docs

13
docs citations

13
times ranked

446
citing authors

#	ARTICLE	IF	CITATIONS
1	Deep Neural Network Model to Predict the Electrostatic Parameters in the Polarizable Classical Drude Oscillator Force Field. <i>Journal of Chemical Theory and Computation</i> , 2022, 18, 1711-1725.	5.3	13
2	Harnessing Deep Learning for Optimization of Lennard-Jones Parameters for the Polarizable Classical Drude Oscillator Force Field. <i>Journal of Chemical Theory and Computation</i> , 2022, 18, 2388-2407.	5.3	17
3	<scp>CHARMMâ€GUI</scp> Drude prepper for molecular dynamics simulation using the classical Drude polarizable force field. <i>Journal of Computational Chemistry</i> , 2022, 43, 359-375.	3.3	24
4	Cancer Cell Metabolism Featuring Nrf2. <i>Current Drug Discovery Technologies</i> , 2020, 17, 263-271.	1.2	2
5	Solvation dynamics: improved reproduction of the time-dependent Stokes shift with polarizable empirical force field chromophore models. <i>Physical Chemistry Chemical Physics</i> , 2019, 21, 17703-17710.	2.8	12
6	Toward Prediction of Electrostatic Parameters for Force Fields That Explicitly Treat Electronic Polarization. <i>Journal of Chemical Theory and Computation</i> , 2019, 15, 2460-2469.	5.3	21
7	Ranking Reversible Covalent Drugs: From Free Energy Perturbation to Fragment Docking. <i>Journal of Chemical Information and Modeling</i> , 2019, 59, 2093-2102.	5.4	35
8	The connexin26 human mutation N14K disrupts cytosolic intersubunit interactions and promotes channel opening. <i>Journal of General Physiology</i> , 2019, 151, 328-341.	1.9	16
9	Molecular Mechanism of Resveratrolâ€™s Lipid Membrane Protection. <i>Scientific Reports</i> , 2018, 8, 1587.	3.3	37
10	Effects of bioactive constituents in the Traditional Chinese Medicinal formula Siâ€™Wuâ€™Tang on Nrf2 signaling and neoplastic cellular transformation. <i>Phytomedicine</i> , 2018, 40, 1-9.	5.3	17
11	Can Relative Binding Free Energy Predict Selectivity of Reversible Covalent Inhibitors?. <i>Journal of the American Chemical Society</i> , 2017, 139, 17945-17952.	13.7	44
12	Polymodal allosteric regulation of Type 1 Serine/Threonine Kinase Receptors via a conserved electrostatic lock. <i>PLoS Computational Biology</i> , 2017, 13, e1005711.	3.2	16
13	Overlapping structure features selection in linear and non-linear QSAR. <i>Journal of Pharmacy Research</i> , 2013, 6, 183-187.	0.4	1