

Jingheng Wu

List of Publications by Year in descending order

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Version: 2024-02-01

10
papers

382
citations

1040056

9
h-index

1372567

10
g-index

10
all docs

10
docs citations

10
times ranked

648
citing authors

#	ARTICLE	IF	CITATIONS
1	Rational design of NIR-emitting iridium(III) complexes for multimodal phosphorescence imaging of mitochondria under two-photon excitation. <i>Chemical Communications</i> , 2017, 53, 10374-10377.	4.1	48
2	Internal force corrections with machine learning for quantum mechanics/molecular mechanics simulations. <i>Journal of Chemical Physics</i> , 2017, 147, 161732.	3.0	28
3	Topoisomerase II \pm poisoning and DNA double-strand breaking by chiral ruthenium(II) complexes containing 2-furanyl-imidazo[4,5-f][1,10]phenanthroline derivatives. <i>Dalton Transactions</i> , 2016, 45, 10546-10555.	3.3	25
4	Multiscale Quantum Mechanics/Molecular Mechanics Simulations with Neural Networks. <i>Journal of Chemical Theory and Computation</i> , 2016, 12, 4934-4946.	5.3	94
5	Human Ferrochelatase: Insights for the Mechanism of Ferrous Iron Approaching Protoporphyrin IX by QM/MM and QTCP Free Energy Studies. <i>Journal of Chemical Information and Modeling</i> , 2016, 56, 2421-2433.	5.4	14
6	Direct imaging of biological sulfur dioxide derivatives in vivo using a two-photon phosphorescent probe. <i>Biomaterials</i> , 2015, 63, 128-136.	11.4	58
7	Molecular docking and QSAR analysis on maleimide derivatives selective inhibition against human monoglyceride lipase based on various modeling methods and conformations. <i>Chemometrics and Intelligent Laboratory Systems</i> , 2014, 131, 22-30.	3.5	6
8	Investigation by MD simulation of the key residues related to substrate-binding and heme-release in human ferrochelatase. <i>Journal of Molecular Modeling</i> , 2013, 19, 2509-2518.	1.8	9
9	Thiol-specific phosphorescent imaging in living cells with an azobis(2,2'-bipyridine)-bridged dinuclear iridium(III) complex. <i>Chemical Communications</i> , 2013, 49, 2040.	4.1	51
10	A self-adaptive genetic algorithm-artificial neural network algorithm with leave-one-out cross validation for descriptor selection in QSAR study. <i>Journal of Computational Chemistry</i> , 2010, 31, 1956-1968.	3.3	49