

Xiaolu Cheng

List of Publications by Year in descending order

Source: <https://exaly.com/author-pdf/3729362/publications.pdf>

Version: 2024-02-01

12
papers

361
citations

1039880

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h-index

1281743

11
g-index

15
all docs

15
docs citations

15
times ranked

426
citing authors

#	ARTICLE	IF	CITATIONS
1	TMEM16 scramblases thin the membrane to enable lipid scrambling. <i>Nature Communications</i> , 2022, 13, 2604.	5.8	22
2	Membrane lipids are both the substrates and a mechanistically responsive environment of TMEM16 scramblase proteins. <i>Journal of Computational Chemistry</i> , 2020, 41, 538-551.	1.5	15
3	Structural Basis of Lipid Scrambling and Ion Conduction by TMEM16 Scramblases. <i>FASEB Journal</i> , 2020, 34, 1-1.	0.2	0
4	Dynamic modulation of the lipid translocation groove generates a conductive ion channel in Ca ²⁺ -bound nhTMEM16. <i>Nature Communications</i> , 2019, 10, 4972.	5.8	23
5	Monitoring Water Clusters "Melt" Through Vibrational Spectroscopy. <i>Journal of the American Chemical Society</i> , 2017, 139, 7082-7088.	6.6	69
6	Vibrational Signatures of Electronic Properties in Oxidized Water: Unraveling the Anomalous Spectrum of the Water Dimer Cation. <i>Journal of the American Chemical Society</i> , 2016, 138, 11936-11945.	6.6	22
7	Tuning vibrational mode localization with frequency windowing. <i>Journal of Chemical Physics</i> , 2016, 145, 124112.	1.2	47
8	Vibrational Signatures of Conformer-Specific Intramolecular Interactions in Protonated Tryptophan. <i>Journal of Physical Chemistry A</i> , 2016, 120, 5598-5608.	1.1	32
9	Accelerating <i>Ab Initio</i> Path Integral Simulations via Imaginary Multiple-Timestepping. <i>Journal of Chemical Theory and Computation</i> , 2016, 12, 1627-1638.	2.3	12
10	Efficient anharmonic vibrational spectroscopy for large molecules using local-mode coordinates. <i>Journal of Chemical Physics</i> , 2014, 141, 104105.	1.2	102
11	Variational mixed quantum/semiclassical simulation of dihalogen guest and rare-gas solid host dynamics. <i>Journal of Chemical Physics</i> , 2014, 141, 034113.	1.2	9
12	Numerical Tests of a Fixed Vibrational Basis/Gaussian Bath Theory for Small Molecule Dynamics in Low-Temperature Media. <i>Journal of Physical Chemistry A</i> , 2011, 115, 3980-3989.	1.1	8