

Viacheslav Bolnykh

List of Publications by Year in descending order

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Version: 2024-02-01

10
papers

248
citations

1307543

7
h-index

1372553

10
g-index

14
all docs

14
docs citations

14
times ranked

201
citing authors

#	ARTICLE	IF	CITATIONS
1	A universal co-solvent dilution strategy enables facile and cost-effective fabrication of perovskite photovoltaics. <i>Nature Communications</i> , 2022, 13, 89.	12.8	77
2	Accuracy of Molecular Simulation-Based Predictions of k_{off} Values: A Metadynamics Study. <i>Journal of Physical Chemistry Letters</i> , 2020, 11, 6373-6381.	4.6	41
3	Extreme Scalability of DFT-Based QM/MM MD Simulations Using MiMiC. <i>Journal of Chemical Theory and Computation</i> , 2019, 15, 5601-5613.	5.3	32
4	MiMiC: A Novel Framework for Multiscale Modeling in Computational Chemistry. <i>Journal of Chemical Theory and Computation</i> , 2019, 15, 3810-3823.	5.3	31
5	Recent Advances in First-Principles Based Molecular Dynamics. <i>Accounts of Chemical Research</i> , 2022, 55, 221-230.	15.6	22
6	Molecular Basis of CLC Antiporter Inhibition by Fluoride. <i>Journal of the American Chemical Society</i> , 2020, 142, 7254-7258.	13.7	20
7	Expanding the boundaries of ligand-target modeling by exascale calculations. <i>Wiley Interdisciplinary Reviews: Computational Molecular Science</i> , 2021, 11, e1535.	14.6	13
8	MiMiC: Multiscale Modeling in Computational Chemistry. <i>Frontiers in Molecular Biosciences</i> , 2020, 7, 45.	3.5	5
9	Biomolecular Simulation: A Perspective from High Performance Computing. <i>Israel Journal of Chemistry</i> , 2020, 60, 694-704.	2.3	2
10	Wavefunction-Based Electrostatic-Embedding QM/MM Using CFOUR through MiMiC. <i>Journal of Chemical Theory and Computation</i> , 2022, 18, 13-24.	5.3	2