## Viacheslav Bolnykh

## List of Publications by Citations

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The third column is the impact factor (IF) of the journal, and the fourth column is the number of citations of the article.

10 87 5 9 g-index

14 156 9.8 2.24 ext. papers ext. citations avg, IF L-index

#	Paper	IF	Citations
10	Extreme Scalability of DFT-Based QM/MM MD Simulations Using MiMiC. <i>Journal of Chemical Theory and Computation</i> , <b>2019</b> , 15, 5601-5613	6.4	20
9	MiMiC: A Novel Framework for Multiscale Modeling in Computational Chemistry. <i>Journal of Chemical Theory and Computation</i> , <b>2019</b> , 15, 3810-3823	6.4	18
8	A universal co-solvent dilution strategy enables facile and cost-effective fabrication of perovskite photovoltaics <i>Nature Communications</i> , <b>2022</b> , 13, 89	17.4	14
7	Accuracy of Molecular Simulation-Based Predictions of Values: A Metadynamics Study. <i>Journal of Physical Chemistry Letters</i> , <b>2020</b> , 11, 6373-6381	6.4	14
6	Molecular Basis of CLC Antiporter Inhibition by Fluoride. <i>Journal of the American Chemical Society</i> , <b>2020</b> , 142, 7254-7258	16.4	10
5	MiMiC: Multiscale Modeling in Computational Chemistry. Frontiers in Molecular Biosciences, <b>2020</b> , 7, 45	5.6	3
4	Recent Advances in First-Principles Based Molecular Dynamics <i>Accounts of Chemical Research</i> , <b>2022</b> ,	24.3	3
3	Expanding the boundaries of ligandEarget modeling by exascale calculations. <i>Wiley Interdisciplinary Reviews: Computational Molecular Science</i> , <b>2021</b> , 11, e1535	7.9	2
2	On the accuracy of molecular simulation-based predictions of koff values: a Metadynamics study		1
1	Biomolecular Simulation: A Perspective from High Performance Computing. <i>Israel Journal of Chemistry</i> , <b>2020</b> , 60, 694-704	3.4	О