

James J Mccarty

List of Publications by Year in descending order

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23
papers

1,545
citations

394286

19
h-index

642610

23
g-index

26
all docs

26
docs citations

26
times ranked

1410
citing authors

#	ARTICLE	IF	CITATIONS
1	Amyloid Oligomers: A Joint Experimental/Computational Perspective on Alzheimer's Disease, Parkinson's Disease, Type II Diabetes, and Amyotrophic Lateral Sclerosis. <i>Chemical Reviews</i> , 2021, 121, 2545-2647.	23.0	406
2	Insight Into Seeded Tau Fibril Growth From Molecular Dynamics Simulation of the Alzheimer's Disease Protofibril Core. <i>Frontiers in Molecular Biosciences</i> , 2021, 8, 624302.	1.6	17
3	Heightened Cold-Denaturation of Proteins at the Ice-Water Interface. <i>Journal of the American Chemical Society</i> , 2020, 142, 5722-5730.	6.6	59
4	The proline-rich domain promotes Tau liquid-liquid phase separation in cells. <i>Journal of Cell Biology</i> , 2020, 219, .	2.3	58
5	Small ion effects on self-coacervation phenomena in block polyampholytes. <i>Journal of Chemical Physics</i> , 2019, 151, 034904.	1.2	46
6	Complete Phase Diagram for Liquid-Liquid Phase Separation of Intrinsically Disordered Proteins. <i>Journal of Physical Chemistry Letters</i> , 2019, 10, 1644-1652.	2.1	204
7	Molecular design of self-coacervation phenomena in block polyampholytes. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2019, 116, 8224-8232.	3.3	88
8	Narrow equilibrium window for complex coacervation of tau and RNA under cellular conditions. <i>ELife</i> , 2019, 8, .	2.8	111
9	Searching for Entropically Stabilized Phases: The Case of Silver Iodide. <i>Journal of Physical Chemistry C</i> , 2018, 122, 1786-1790.	1.5	18
10	Accuracy, Transferability, and Efficiency of Coarse-Grained Models of Molecular Liquids. <i>Journal of Physical Chemistry B</i> , 2018, 122, 10257-10278.	1.2	26
11	Effect of Surfactants on Surface-Induced Denaturation of Proteins: Evidence of an Orientation-Dependent Mechanism. <i>Journal of Physical Chemistry B</i> , 2018, 122, 11390-11399.	1.2	33
12	Variational Flooding Study of a S_N2 Reaction. <i>Journal of Physical Chemistry Letters</i> , 2017, 8, 580-583.	2.1	23
13	A variational conformational dynamics approach to the selection of collective variables in metadynamics. <i>Journal of Chemical Physics</i> , 2017, 147, 204109.	1.2	101
14	Bespoke Bias for Obtaining Free Energy Differences within Variationally Enhanced Sampling. <i>Journal of Chemical Theory and Computation</i> , 2016, 12, 2162-2169.	2.3	11
15	Variationally Optimized Free-Energy Flooding for Rate Calculation. <i>Physical Review Letters</i> , 2015, 115, 070601.	2.9	35
16	Fast equilibration of coarse-grained polymeric liquids. <i>Journal of Computational Science</i> , 2015, 9, 33-38.	1.5	8
17	An analytical coarse-graining method which preserves the free energy, structural correlations, and thermodynamic state of polymer melts from the atomistic to the mesoscale. <i>Journal of Chemical Physics</i> , 2014, 140, 204913.	1.2	65
18	Effective potentials for representing polymers in melts as chains of interacting soft particles. <i>Journal of Chemical Physics</i> , 2013, 139, 124906.	1.2	36

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19	Thermodynamic Consistency in Variable-Level Coarse Graining of Polymeric Liquids. <i>Physical Review Letters</i> , 2012, 109, 168301.	2.9	55
20	Thermodynamic Consistency between Analytic Integral Equation Theory and Coarse-Grained Molecular Dynamics Simulations of Homopolymer Melts. <i>Macromolecules</i> , 2012, 45, 8482-8493.	2.2	35
21	Analytical rescaling of polymer dynamics from mesoscale simulations. <i>Journal of Chemical Physics</i> , 2010, 132, 224903.	1.2	44
22	Effective Soft-Core Potentials and Mesoscopic Simulations of Binary Polymer Mixtures. <i>Macromolecules</i> , 2010, 43, 3964-3979.	2.2	25
23	Multiscale Modeling of Coarse-Grained Macromolecular Liquids. <i>Journal of Physical Chemistry B</i> , 2009, 113, 11876-11886.	1.2	36