

Kazuaki Z Takahashi

List of Publications by Year in descending order

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papers

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759233

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32
all docs

32
docs citations

32
times ranked

253
citing authors

#	ARTICLE	IF	CITATIONS
1	Cutoff radius effect of the isotropic periodic sum and Wolf method in liquid–vapor interfaces of water. <i>Journal of Chemical Physics</i> , 2011, 134, 174112.	3.0	39
2	Cutoff radius effect of the isotropic periodic sum method in homogeneous system. II. Water. <i>Journal of Chemical Physics</i> , 2010, 133, 014109.	3.0	32
3	Cutoff radius effect of isotropic periodic sum method for transport coefficients of Lennard-Jones liquid. <i>Journal of Chemical Physics</i> , 2007, 127, 114511.	3.0	28
4	Machine learning-aided analysis for complex local structure of liquid crystal polymers. <i>Scientific Reports</i> , 2019, 9, 16370.	3.3	27
5	Molecular Dynamics Simulations for Resolving Scaling Laws of Polyethylene Melts. <i>Polymers</i> , 2017, 9, 24.	4.5	25
6	Multistep nucleation of anisotropic molecules. <i>Nature Communications</i> , 2021, 12, 5278.	12.8	23
7	A combination of the tree-code and IPS method to simulate large scale systems by molecular dynamics. <i>Journal of Chemical Physics</i> , 2011, 135, 174108.	3.0	19
8	An Improved Isotropic Periodic Sum Method That Uses Linear Combinations of Basis Potentials. <i>Journal of Chemical Theory and Computation</i> , 2012, 8, 4503-4516.	5.3	16
9	Onset of static and dynamic universality among molecular models of polymers. <i>Scientific Reports</i> , 2017, 7, 12379.	3.3	15
10	Truncation Effects of Shift Function Methods in Bulk Water Systems. <i>Entropy</i> , 2013, 15, 3249-3264.	2.2	14
11	Mining of effective local order parameters for classifying crystal structures: A machine learning study. <i>Journal of Chemical Physics</i> , 2020, 152, 214501.	3.0	14
12	Searching local order parameters to classify water structures of ice Ih, Ic, and liquid. <i>Journal of Chemical Physics</i> , 2021, 154, 164505.	3.0	13
13	Cut-off radius effect of the isotropic periodic sum method for polar molecules in a bulk water system. <i>Molecular Simulation</i> , 2012, 38, 397-403.	2.0	12
14	Design of a reaction field using a linear–combination–based isotropic periodic sum method. <i>Journal of Computational Chemistry</i> , 2014, 35, 865-875.	3.3	10
15	Development of Coarse-Grained Liquid-Crystal Polymer Model with Efficient Electrostatic Interaction: Toward Molecular Dynamics Simulations of Electroactive Materials. <i>Materials</i> , 2018, 11, 83.	2.9	10
16	Application of isotropic periodic sum method for 4-pentyl-4'-cyanobiphenyl liquid crystal. <i>Molecular Simulation</i> , 2015, 41, 927-935.	2.0	8
17	Critical test of bead–spring model to resolve the scaling laws of polymer melts: a molecular dynamics study. <i>Molecular Simulation</i> , 2017, 43, 1196-1201.	2.0	8
18	Performance of Coarse Graining in Estimating Polymer Properties: Comparison with the Atomistic Model. <i>Polymers</i> , 2020, 12, 382.	4.5	7

#	ARTICLE	IF	CITATIONS
19	Molecular architecture dependence of mesogen rotation during uniaxial elongation of liquid crystal elastomers. <i>Polymer</i> , 2021, 229, 123970.	3.8	7
20	Comparison of the accuracy of periodic reaction field methods in molecular dynamics simulations of a model liquid crystal system. <i>Journal of Computational Chemistry</i> , 2015, 36, 2406-2411.	3.3	6
21	Critical test of isotropic periodic sum techniques with group-based cut-off schemes. <i>Scientific Reports</i> , 2018, 8, 4185.	3.3	6
22	Searching for local order parameters to classify water structures at triple points. <i>Journal of Computational Chemistry</i> , 2021, 42, 1720-1727.	3.3	6
23	Mining of Effective Local Order Parameters to Classify Ice Polymorphs. <i>Journal of Physical Chemistry A</i> , 2021, 125, 9518-9526.	2.5	6
24	A determination of liquid-vapour interfacial properties for methanol using a linear-combination-based isotropic periodic sum. <i>Molecular Simulation</i> , 2015, 41, 795-800.	2.0	4
25	An improvement of truncation method by a novel reaction field: Accurate computation for estimating methanol liquid-vapor interfacial systems. <i>Computational Materials Science</i> , 2015, 100, 191-194.	3.0	3
26	Trade-off effect between the stress and strain range in the soft elasticity of liquid crystalline elastomers. <i>Polymer Journal</i> , 2022, 54, 1017-1027.	2.7	3
27	A fast and accurate computational method for the linear-combination-based isotropic periodic sum. <i>Scientific Reports</i> , 2018, 8, 11880.	3.3	2
28	Screening toward the Development of Fingerprints of Atomic Environments Using Bond-Orientational Order Parameters. <i>ACS Omega</i> , 2022, 7, 4606-4613.	3.5	2
29	Combined use of periodic reaction field and coarse-grained molecular dynamics simulations. I. phospholipid monolayer systems. <i>Molecular Simulation</i> , 2017, 43, 971-976.	2.0	1
30	Molecular dynamics study on structural change in the surface membrane of an insonified coated microbubble. , 2012, , .		0
31	Cutoff Radius Effect of Water Configuration Using the Wolf Method. , 2011, , .		0