

Yong Zhang

List of Publications by Year in descending order

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25
papers

1,967
citations

471061

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580395

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27
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docs citations

27
times ranked

1416
citing authors

#	ARTICLE	IF	CITATIONS
1	Deep Eutectic Solvents: A Review of Fundamentals and Applications. <i>Chemical Reviews</i> , 2021, 121, 1232-1285.	23.0	1,334
2	Liquid Structure and Transport Properties of the Deep Eutectic Solvent Ethaline. <i>Journal of Physical Chemistry B</i> , 2020, 124, 5251-5264.	1.2	84
3	Solvation Structure and Dynamics of Li ⁺ in Ternary Ionic Liquid–Lithium Salt Electrolytes. <i>Journal of Physical Chemistry B</i> , 2019, 123, 516-527.	1.2	62
4	Water-in-Salt LiTFSI Aqueous Electrolytes. 1. Liquid Structure from Combined Molecular Dynamics Simulation and Experimental Studies. <i>Journal of Physical Chemistry B</i> , 2021, 125, 4501-4513.	1.2	52
5	PyLAT: Python LAMMPS Analysis Tools. <i>Journal of Chemical Information and Modeling</i> , 2019, 59, 1301-1305.	2.5	49
6	Water or Anion? Uncovering the Zn ²⁺ Solvation Environment in Mixed Zn(TFSI) ₂ and LiTFSI Water-in-Salt Electrolytes. <i>ACS Energy Letters</i> , 2021, 6, 3458-3463.	8.8	45
7	Signatures of Ion Pairing and Aggregation in the Vibrational Spectroscopy of Super-Concentrated Aqueous Lithium Bistriflimide Solutions. <i>Journal of Physical Chemistry C</i> , 2020, 124, 3470-3481.	1.5	44
8	Solvation Dynamics of Wet Ethaline: Water is the Magic Component. <i>Journal of Physical Chemistry B</i> , 2021, 125, 8888-8901.	1.2	32
9	Computational Design of New Magnesium Electrolytes with Improved Properties. <i>Journal of Physical Chemistry C</i> , 2017, 121, 16126-16136.	1.5	26
10	Water-In-Salt LiTFSI Aqueous Electrolytes (2): Transport Properties and Li ⁺ Dynamics Based on Molecular Dynamics Simulations. <i>Journal of Physical Chemistry B</i> , 2021, 125, 13246-13254.	1.2	26
11	Evaluation and Refinement of the General AMBER Force Field for Nineteen Pure Organic Electrolyte Solvents. <i>Journal of Chemical & Engineering Data</i> , 2018, 63, 3488-3502.	1.0	23
12	Effect of alkyl-group flexibility on the melting point of imidazolium-based ionic liquids. <i>Journal of Chemical Physics</i> , 2020, 153, 044504.	1.2	20
13	Sigma profiles in deep learning: towards a universal molecular descriptor. <i>Chemical Communications</i> , 2022, 58, 5630-5633.	2.2	20
14	Temperature Dependence of Volumetric and Dynamic Properties of Imidazolium-Based Ionic Liquids. <i>Journal of Physical Chemistry B</i> , 2018, 122, 2414-2424.	1.2	19
15	Anion Enhancement at the Liquid–Vacuum Interface of an Ionic Liquid Mixture. <i>Journal of Physical Chemistry C</i> , 2018, 122, 27392-27401.	1.5	19
16	Comparison of fixed charge and polarizable models for predicting the structural, thermodynamic, and transport properties of molten alkali chlorides. <i>Journal of Chemical Physics</i> , 2020, 153, 214502.	1.2	19
17	Impact of anion shape on Li ⁺ solvation and on transport properties for lithium–air batteries: a molecular dynamics study. <i>Physical Chemistry Chemical Physics</i> , 2020, 22, 15842-15852.	1.3	19
18	Refined Classical Force Field for Choline Chloride and Ethylene Glycol Mixtures over Wide Composition Range. <i>Journal of Chemical & Engineering Data</i> , 2022, 67, 1864-1871.	1.0	19

#	ARTICLE	IF	CITATIONS
19	Exchange-Mediated Transport in Battery Electrolytes: Ultrafast or Ultraslow?. Journal of the American Chemical Society, 2022, 144, 8591-8604.	6.6	18
20	Melting points of alkali chlorides evaluated for a polarizable and non-polarizable model. Journal of Chemical Physics, 2020, 153, 011101.	1.2	15
21	From Networked to Isolated: Observing Water Hydrogen Bonds in Concentrated Electrolytes with Two-Dimensional Infrared Spectroscopy. Journal of Physical Chemistry B, 2022, 126, 5305-5319.	1.2	9
22	Structure of water-in-salt and water-in-bisalt electrolytes. Physical Chemistry Chemical Physics, 2022, 24, 10727-10736.	1.3	5
23	Layer-based thermal migration of an ionic liquid nano-droplet on a graphene surface: a molecular dynamics study. Molecular Simulation, 2020, 46, 829-836.	0.9	3
24	Functionalized Phosphonium Cations Enable Zn Metal Reversibility in Aqueous Electrolytes. ECS Meeting Abstracts, 2021, MA2021-02, 14-14.	0.0	0
25	Liquid Structure and Transport Properties of the Deep Eutectic Solvent Ethaline. ECS Meeting Abstracts, 2020, MA2020-02, 2910-2910.	0.0	0