

Gulou Shen

List of Publications by Year in descending order

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times ranked

209
citing authors

#	ARTICLE	IF	CITATIONS
1	Effect of surface roughness on partition of ionic liquids in nanopores by a perturbed-chain SAFT density functional theory. <i>Journal of Chemical Physics</i> , 2022, 157, .	3.0	4
2	Modeling interfacial properties of ionic liquids with ePC-SAFT combined with density gradient theory. <i>Fluid Phase Equilibria</i> , 2021, 536, 112984.	2.5	5
3	Accelerate the ePC-SAFT-DFT Calculation with the Chebyshev Pseudospectral Collocation Method. <i>Industrial & Engineering Chemistry Research</i> , 2021, 60, 9269-9285.	3.7	7
4	Partition and selectivity of electrolytes in cylindrical nanopores with heterogeneous surface charge. <i>Journal of Molecular Liquids</i> , 2021, 340, 116839.	4.9	5
5	Accelerate the Electrolyte Perturbed-Chain Statistical Associating Fluid Theory's Density Functional Theory Calculation With the Chebyshev Pseudo-Spectral Collocation Method. Part II. Spherical Geometry and Anderson Mixing. <i>Frontiers in Chemistry</i> , 2021, 9, 801551.	3.6	4
6	Interfacial structure and differential capacitance of ionic liquid/graphite interface: A perturbed-chain SAFT density functional theory study. <i>Journal of Molecular Liquids</i> , 2020, 310, 113199.	4.9	14
7	Modeling Thermodynamic Derivative Properties and Gas Solubility of Ionic Liquids with ePC-SAFT. <i>Industrial & Engineering Chemistry Research</i> , 2019, 58, 8401-8417.	3.7	33
8	Modeling Viscosity of Ionic Liquids with Electrolyte Perturbed-Chain Statistical Associating Fluid Theory and Free Volume Theory. <i>Industrial & Engineering Chemistry Research</i> , 2018, 57, 8784-8801.	3.7	28
9	Developing Electrolyte Perturbed-Chain Statistical Associating Fluid Theory Density Functional Theory for CO ₂ Separation by Confined Ionic Liquids. <i>Journal of Physical Chemistry C</i> , 2018, 122, 15464-15473.	3.1	12
10	Modelling interfacial properties of ionic liquids with ePC-SAFT combined with density gradient theory. <i>Molecular Physics</i> , 2016, 114, 2492-2499.	1.7	14
11	Modeling thermodynamic derivative properties of ionic liquids with ePC-SAFT. <i>Fluid Phase Equilibria</i> , 2015, 405, 73-82.	2.5	43
12	Modeling the Viscosity of Ionic Liquids with the Electrolyte Perturbed-Chain Statistical Association Fluid Theory. <i>Industrial & Engineering Chemistry Research</i> , 2014, 53, 20258-20268.	3.7	32
13	A hybrid perturbed-chain SAFT density functional theory for representing fluid behavior in nanopores. <i>Journal of Chemical Physics</i> , 2013, 138, 224706.	3.0	31
14	A hybrid perturbed-chain SAFT density functional theory for representing fluid behavior in nanopores: Mixtures. <i>Journal of Chemical Physics</i> , 2013, 139, 194705.	3.0	30