

Ramesh Singh

List of Publications by Year in descending order

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Version: 2024-02-01

8
papers

77
citations

1684188
5
h-index

1588992
8
g-index

8
all docs

8
docs citations

8
times ranked

116
citing authors

#	ARTICLE	IF	CITATIONS
1	A Monte Carlo Simulation Study To Predict the Solubility of Carbon Dioxide, Hydrogen, and Their Mixture in the Ionic Liquids 1-Alkyl-3-methylimidazolium bis(trifluoromethanesulfonyl)amide ([C _n mim] ⁺ [Tf ₂ N] ⁻], <i>n</i> = 4, 6). <i>Industrial & Engineering Chemistry Research</i> , 2015, 54, 4385-4395.	3.7	21
2	Effect of CO ₂ and H ₂ O on the behavior of shale gas confined inside calcite [104] slit-like nanopore: a molecular dynamics simulation study. <i>Journal of Molecular Modeling</i> , 2019, 25, 293.	1.8	18
3	Molecular dynamics simulation of shale gas confined inside slit-like calcite [104] nanopore. <i>Molecular Simulation</i> , 2019, 45, 104-110.	2.0	12
4	COSMO-RS Analysis of CO ₂ Solubility in <i>N</i> -Methyldiethanolamine, Sulfolane, and 1-Butyl-3-methyl-imidazolium Acetate Activated by 2-Methylpiperazine for Postcombustion Carbon Capture. <i>ACS Omega</i> , 2021, 6, 747-761.	3.5	10
5	Utilizing Simtronics, a chemical engineering process simulation software, in chemical engineering coursework to reduce the skills gap. <i>Computer Applications in Engineering Education</i> , 2019, 27, 519-525.	3.4	5
6	Thermodynamic analysis using COSMO-RS studies of reversible ionic liquid 3-aminopropyl triethoxysilane blended with amine activators for CO ₂ absorption. <i>Journal of Molecular Liquids</i> , 2021, 324, 114713.	4.9	5
7	Size effect of oscillating columns on mixing: A CFD study. <i>European Journal of Mechanics, B/Fluids</i> , 2019, 77, 230-238.	2.5	4
8	A Monte Carlo simulation study to predict the solubility of H ₂ S in ionic liquids with 1-butyl-3-methylimidazolium ([C ₄ mim] ⁺), tetrafluoroborate ([BF ₄] ⁻), hexafluorophosphate ([PF ₆] ⁻) and bis(trifluoromethanesulfonyl)amide ([Tf ₂ N] ⁻) anions. <i>Molecular Simulation</i> , 2017, 43, 291-297.	2.0	2