David R Cole

List of Publications by Year in descending order

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361413 434195 1,109 31 20 31 citations h-index g-index papers 31 31 31 1229 docs citations times ranked citing authors all docs

#	Article	IF	Citations
1	Molecular Structure of Adsorbed Water Phases in Silica Nanopores. Journal of Physical Chemistry C, 2022, 126, 2885-2895.	3.1	8
2	Effect of Pore Connectivity on the Behavior of Fluids Confined in Sub-Nanometer Pores: Ethane and CO2 Confined in ZSM-22. Membranes, 2021, 11, 113.	3.0	7
3	Comparative geochemistry of flowback chemistry from the Utica/Point Pleasant and Marcellus formations. Chemical Geology, 2021, 564, 120041.	3.3	11
4	CO2 Adsorption in Metal-Organic Framework Mg-MOF-74: Effects of Inter-Crystalline Space. Nanomaterials, 2020, 10, 2274.	4.1	18
5	Effects of inter-crystalline space on the adsorption of ethane and CO ₂ in silicalite: implications for enhanced adsorption. Physical Chemistry Chemical Physics, 2020, 22, 13951-13957.	2.8	14
6	Structure and dynamics of ethane confined in silica nanopores in the presence of CO2. Journal of Chemical Physics, 2020, 152, 084707.	3.0	14
7	In situ transformation of ethoxylate and glycol surfactants by shale-colonizing microorganisms during hydraulic fracturing. ISME Journal, 2019, 13, 2690-2700.	9.8	18
8	Genome-Resolved Metagenomics Extends the Environmental Distribution of the <i>Verrucomicrobia</i> Phylum to the Deep Terrestrial Subsurface. MSphere, 2019, 4, .	2.9	38
9	Effects of water on the stochastic motions of propane confined in MCM-41-S pores. Physical Chemistry Chemical Physics, 2019, 21, 25035-25046.	2.8	16
10	Sorption, Structure and Dynamics of CO2 and Ethane in Silicalite at High Pressure: A Combined Monte Carlo and Molecular Dynamics Simulation Study. Molecules, 2019, 24, 99.	3.8	18
11	Structure and dynamics of water on the forsterite surface. Physical Chemistry Chemical Physics, 2018, 20, 27822-27829.	2.8	10
12	Comparative genomics and physiology of the genus <i>Methanohalophilus</i> , a prevalent methanogen in hydraulically fractured shale. Environmental Microbiology, 2018, 20, 4596-4611.	3.8	28
13	Members of Marinobacter and Arcobacter Influence System Biogeochemistry During Early Production of Hydraulically Fractured Natural Gas Wells in the Appalachian Basin. Frontiers in Microbiology, 2018, 9, 2646.	3.5	33
14	Aqueous Hydrogen Sulfide in Slit-Shaped Silica Nanopores: Confinement Effects on Solubility, Structural, and Dynamical Properties. Journal of Physical Chemistry C, 2018, 122, 14744-14755.	3.1	20
15	Effects of Confinement and Pressure on the Vibrational Behavior of Nano-Confined Propane. Journal of Physical Chemistry A, 2018, 122, 6736-6745.	2.5	20
16	Coupled laboratory and field investigations resolve microbial interactions that underpin persistence in hydraulically fractured shales. Proceedings of the National Academy of Sciences of the United States of America, 2018, 115, E6585-E6594.	7.1	69
17	Propane–Water Mixtures Confined within Cylindrical Silica Nanopores: Structural and Dynamical Properties Probed by Molecular Dynamics. Langmuir, 2017, 33, 11310-11320.	3.5	30
18	Molecular dynamics simulations of propane in slit shaped silica nano-pores: direct comparison with quasielastic neutron scattering experiments. Physical Chemistry Chemical Physics, 2017, 19, 32320-32332.	2.8	22

#	Article	IF	CITATION
19	Sulfide Generation by Dominant <i>Halanaerobium</i> Microorganisms in Hydraulically Fractured Shales. MSphere, 2017, 2, .	2.9	62
20	Confined Water Determines Transport Properties of Guest Molecules in Narrow Pores. ACS Nano, 2016, 10, 7646-7656.	14.6	66
21	Microbial metabolisms in a 2.5-km-deep ecosystem created by hydraulic fracturing in shales. Nature Microbiology, 2016, 1, 16146.	13.3	207
22	Role of Confinement on Adsorption and Dynamics of Ethane and an Ethane–CO ₂ Mixture in Mesoporous CPG Silica. Journal of Physical Chemistry C, 2016, 120, 4843-4853.	3.1	28
23	Factors governing the behaviour of aqueous methane in narrow pores. Philosophical Transactions Series A, Mathematical, Physical, and Engineering Sciences, 2016, 374, 20150019.	3.4	20
24	N-octane diffusivity enhancement via carbon dioxide in silica slit-shaped nanopores – a molecular dynamics simulation. Molecular Simulation, 2016, 42, 745-752.	2.0	24
25	CO ₂ –C ₄ H ₁₀ Mixtures Simulated in Silica Slit Pores: Relation between Structure and Dynamics. Journal of Physical Chemistry C, 2015, 119, 15274-15284.	3.1	86
26	Dynamics of Propane in Nanoporous Silica Aerogel: A Quasielastic Neutron Scattering Study. Journal of Physical Chemistry C, 2015, 119, 18188-18195.	3.1	29
27	Propane simulated in silica pores: Adsorption isotherms, molecular structure, and mobility. Chemical Engineering Science, 2015, 121, 292-299.	3.8	43
28	Sorption Phase of Supercritical CO ₂ in Silica Aerogel: Experiments and Mesoscale Computer Simulations. Journal of Physical Chemistry C, 2014, 118, 15525-15533.	3.1	24
29	Direct Measurements of Pore Fluid Density by Vibrating Tube Densimetry. Langmuir, 2012, 28, 5070-5078.	3.5	29
30	Pore Size Effects on the Sorption of Supercritical CO ₂ in Mesoporous CPG-10 Silica. Journal of Physical Chemistry C, 2012, 116, 917-922.	3.1	50
31	Microstructural Characterization of Adsorption and Depletion Regimes of Supercritical Fluids in	3.1	47