

Jun Zhang

List of Publications by Citations

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The third column is the impact factor (IF) of the journal, and the fourth column is the number of citations of the article.

68

papers

1,223

citations

20

h-index

32

g-index

69

ext. papers

1,566

ext. citations

6.2

avg, IF

4.78

L-index

#	Paper	IF	Citations
68	Improved hydrogen storage in the modified metal-organic frameworks by hydrogen spillover effect. <i>International Journal of Hydrogen Energy</i> , 2007 , 32, 4005-4010	6.7	146
67	Nanoparticles mimicking viral surface topography for enhanced cellular delivery. <i>Advanced Materials</i> , 2013 , 25, 6233-7	24	129
66	Adsorption mechanism of oil components on water-wet mineral surface: A molecular dynamics simulation study. <i>Energy</i> , 2013 , 59, 295-300	7.9	80
65	Study on the transformation from linear to branched wormlike micelles: An insight from molecular dynamics simulation. <i>Journal of Colloid and Interface Science</i> , 2017 , 494, 47-53	9.3	42
64	Molecular dynamics simulation on volume swelling of CO ₂ /alkane system. <i>Fuel</i> , 2015 , 143, 194-201	7.1	40
63	Highly efficient water desalination in carbon nanocones. <i>Carbon</i> , 2018 , 129, 374-379	10.4	40
62	Enantioselective Molecular Transport in Multilayer Graphene Nanopores. <i>Nano Letters</i> , 2017 , 17, 6742-6746	14.5	37
61	Competitive adsorption and diffusion of CH ₄ /CO ₂ binary mixture within shale organic nanochannels. <i>Journal of Natural Gas Science and Engineering</i> , 2018 , 53, 329-336	4.6	36
60	Displacement Mechanism of Oil in Shale Inorganic Nanopores by Supercritical Carbon Dioxide from Molecular Dynamics Simulations. <i>Energy & Fuels</i> , 2017 , 31, 738-746	4.1	34
59	Enhanced oil recovery mechanism of CO ₂ water-alternating-gas injection in silica nanochannel. <i>Fuel</i> , 2017 , 190, 253-259	7.1	33
58	A graphene-like membrane with an ultrahigh water flux for desalination. <i>Nanoscale</i> , 2017 , 9, 18951-18957	7.7	30
57	Oil detachment mechanism in CO ₂ flooding from silica surface: Molecular dynamics simulation. <i>Chemical Engineering Science</i> , 2017 , 164, 17-22	4.4	27
56	Coarse-grained molecular dynamics study on the self-assembly of Gemini surfactants: the effect of spacer length. <i>Physical Chemistry Chemical Physics</i> , 2017 , 19, 4462-4468	3.6	24
55	Electrical field facilitates selective transport of CO ₂ through a laminated MoS ₂ supported ionic liquid membrane. <i>Journal of Materials Chemistry A</i> , 2019 , 7, 10041-10046	13	24
54	CO ₂ -philic WS ₂ laminated membranes with a nanoconfined ionic liquid. <i>Journal of Materials Chemistry A</i> , 2018 , 6, 16566-16573	13	24
53	Unusual, photo-induced self-assembly of azobenzene-containing amphiphiles. <i>Soft Matter</i> , 2014 , 10, 8758-64	3.6	24
52	Nanoconfined deep eutectic solvent in laminated MXene for efficient CO ₂ separation. <i>Chemical Engineering Journal</i> , 2021 , 405, 126961	14.7	24

51	A solution-processable and ultra-permeable conjugated microporous thermoset for selective hydrogen separation. <i>Nature Communications</i> , 2020 , 11, 1633	17.4	23
50	Correlated Rectification Transport in Ultranarrow Charged Nanocones. <i>Journal of Physical Chemistry Letters</i> , 2017 , 8, 435-439	6.4	22
49	Static and dynamic behavior of CO ₂ enhanced oil recovery in nanoslits: Effects of mineral type and oil components. <i>International Journal of Heat and Mass Transfer</i> , 2020 , 153, 119583	4.9	20
48	Bottom-up self-assembly of heterotrimeric nanoparticles and their secondary Janus generations. <i>Chemical Science</i> , 2019 , 10, 10388-10394	9.4	20
47	The molecular mechanism of the inhibition effects of PVCaps on the growth of sl hydrate: an unstable adsorption mechanism. <i>Physical Chemistry Chemical Physics</i> , 2018 , 20, 8326-8332	3.6	19
46	Molecular insight into the miscible mechanism of CO ₂ /C ₁₀ in bulk phase and nanoslits. <i>International Journal of Heat and Mass Transfer</i> , 2019 , 141, 643-650	4.9	18
45	CO activating hydrocarbon transport across nanopore throat: insights from molecular dynamics simulation. <i>Physical Chemistry Chemical Physics</i> , 2017 , 19, 30439-30444	3.6	17
44	Vesicle formation of cationic mixtures of CTAC/SDS induced by ratio: a coarse-grained molecular dynamic simulation study. <i>RSC Advances</i> , 2016 , 6, 13442-13449	3.7	17
43	How the oil recovery in deep oil reservoirs is affected by injected gas types: A molecular dynamics simulation study. <i>Chemical Engineering Science</i> , 2021 , 231, 116286	4.4	17
42	Oil extraction mechanism in CO ₂ flooding from rough surface: Molecular dynamics simulation. <i>Applied Surface Science</i> , 2019 , 494, 80-86	6.7	16
41	Oil detachment by modified nanoparticles: A molecular dynamics simulation study. <i>Computational Materials Science</i> , 2019 , 170, 109177	3.2	15
40	Mechanism of asphaltene aggregation induced by supercritical CO ₂ : insights from molecular dynamics simulation. <i>RSC Advances</i> , 2017 , 7, 50786-50793	3.7	15
39	The self-assembly structure and the CO-philicity of a hybrid surfactant in supercritical CO: effects of hydrocarbon chain length. <i>Soft Matter</i> , 2016 , 12, 8177-8185	3.6	13
38	Photothermal-Responsive Microporous Nanosheets Confined Ionic Liquid for Efficient CO Separation. <i>Small</i> , 2020 , 16, e2002699	11	13
37	Molecular dynamics study of di-CF ₄ based reverse micelles in supercritical CO. <i>Physical Chemistry Chemical Physics</i> , 2016 , 18, 29156-29163	3.6	13
36	One-pot production of porous assemblies by PISA of star architecture copolymers: a simulation study. <i>Physical Chemistry Chemical Physics</i> , 2018 , 20, 10069-10076	3.6	12
35	Water desalination of a new three-dimensional covalent organic framework: a molecular dynamics simulation study. <i>Physical Chemistry Chemical Physics</i> , 2020 , 22, 16978-16984	3.6	10
34	Molecular Insights into the Effect of a Solid Surface on the Stability of a Hydrate Nucleus. <i>Journal of Physical Chemistry C</i> , 2020 , 124, 2664-2671	3.8	10

33	Cooperative assembly of Janus particles and amphiphilic oligomers: the role of Janus balance. <i>Nanoscale</i> , 2019 , 11, 7221-7228	7.7	9
32	Why synthetic virus-like nanoparticles can achieve higher cellular uptake efficiency?. <i>Nanoscale</i> , 2020 , 12, 14911-14918	7.7	9
31	Effect of organic salt on the self-assembly of ammonium gemini surfactant: An experiment and simulation study. <i>Colloids and Surfaces A: Physicochemical and Engineering Aspects</i> , 2018 , 548, 198-205	5.1	8
30	Ionic liquid gated 2D-CAP membrane for highly efficient CO ₂ /N ₂ and CO ₂ /CH ₄ separation. <i>Applied Surface Science</i> , 2019 , 494, 477-483	6.7	8
29	Migration of oil/methane mixture in shale inorganic nano-pore throat: A molecular dynamics simulation study. <i>Journal of Petroleum Science and Engineering</i> , 2020 , 187, 106784	4.4	8
28	Molecular-Scale Design of Hydrocarbon Surfactant Self-Assembly in Supercritical CO ₂ . <i>Langmuir</i> , 2017 , 33, 5291-5297	4	7
27	Composite Nanotube Ring Structures Formed by Two-Step Self-Assembly for Drug Loading/Release. <i>Langmuir</i> , 2019 , 35, 3108-3115	4	7
26	Prediction of efficient promoter molecules of sH hydrogen hydrate: An ab initio study. <i>Chemical Physics</i> , 2019 , 516, 15-21	2.3	7
25	Molecular insight into the fluidity of cement pastes: Nano-boundary lubrication of cementitious materials. <i>Construction and Building Materials</i> , 2022 , 316, 125800	6.7	7
24	Molecular insights into the separation mechanism of imidazole-based ionic liquid supported membranes. <i>Journal of Molecular Liquids</i> , 2021 , 340, 117173	6	7
23	Optimal aggregation number of reverse micelles in supercritical carbon dioxide: a theoretical perspective. <i>Soft Matter</i> , 2019 , 15, 3323-3329	3.6	6
22	Molecular insight into the aggregation and dispersion behavior of modified nanoparticles. <i>Journal of Petroleum Science and Engineering</i> , 2020 , 191, 107193	4.4	6
21	Molecular insight into the oil charging mechanism in tight reservoirs. <i>Chemical Engineering Science</i> , 2020 , 211, 115297	4.4	6
20	Voltage-gated multilayer graphene nanochannel for K ⁺ /Na ⁺ separation: A molecular dynamics study. <i>Journal of Molecular Liquids</i> , 2020 , 317, 114025	6	6
19	Two-dimensional hydrogen hydrates: structure and stability. <i>Physical Chemistry Chemical Physics</i> , 2020 , 22, 5774-5784	3.6	5
18	Tuning the self-assembly of surfactants by the confinement of carbon nanotube arrays: a cornucopia of lamellar phase variants. <i>Nanoscale</i> , 2015 , 7, 6069-74	7.7	5
17	Insight into the pressure-induced displacement mechanism for selecting efficient nanofluids in various capillaries. <i>Environmental Science: Nano</i> , 2020 , 7, 2785-2794	7.1	5
16	Manipulating Hybrid Nanostructures by the Cooperative Assembly of Amphiphilic Oligomers and Triblock Janus Nanoparticles. <i>Journal of Physical Chemistry Letters</i> , 2020 , 11, 3369-3375	6.4	4

15	Alternating electric field-induced ion current rectification and electroosmotic pump in ultranarrow charged carbon nanocones. <i>Physical Chemistry Chemical Physics</i> , 2018 , 20, 27910-27916	3.6	4
14	Shape transition of water-in-CO reverse micelles controlled by the surfactant midpiece. <i>Physical Chemistry Chemical Physics</i> , 2018 , 20, 15535-15542	3.6	4
13	Dissipative Particle Dynamics Simulation on Vesicles Self-Assembly Controlled by Terminal Groups. <i>Journal of Physical Chemistry B</i> , 2018 , 122, 10607-10614	3.4	3
12	Molecular mechanism of formation of the face-sharing double cages in structure-I methane hydrate. <i>Chemical Physics Letters</i> , 2018 , 691, 155-162	2.5	2
11	The biphasic effect of ABA triblock copolymers on the self-assembly of surfactants: insight from dissipative particle dynamics. <i>Molecular Systems Design and Engineering</i> , 2019 , 4, 921-928	4.6	1
10	Combined DFT and kinetic Monte Carlo study of a bridging-spillover mechanism on fluorine-decorated graphene. <i>Physical Chemistry Chemical Physics</i> , 2021 , 23, 2384-2391	3.6	1
9	Polymerization-Induced Self-Assembly of Comb-like Amphiphilic Copolymers into Onion-like Vesicles. <i>Macromolecules</i> , 2021 , 54, 7448-7459	5.5	1
8	Atypical adsorption of polycarboxylate superplasticizers on calcium silicate hydrate surface: Converting interaction by solvent effects. <i>Construction and Building Materials</i> , 2022 , 330, 127160	6.7	1
7	Quantum Chemical Study of the Carbon Dioxide-Philicity of Surfactants: Effects of Tail Functionalization. <i>Langmuir</i> , 2020 , 36, 15352-15361	4	0
6	Novel joint catalytic properties of Fe and N co-doped graphene for CO oxidation. <i>Physical Chemistry Chemical Physics</i> , 2020 , 22, 28376-28382	3.6	0
5	Manipulating the interactions between the lipid bilayer and triblock Janus nanoparticles: insight from dissipative particle dynamics. <i>Molecular Systems Design and Engineering</i> , 2021 , 6, 156-162	4.6	0
4	Molecular insight into the oil displacement mechanism of gas flooding in deep oil reservoir. <i>Chemical Physics Letters</i> , 2021 , 783, 139044	2.5	0
3	An ion sieving conjugated microporous thermoset ultrathin membrane for high-performance Li-S battery. <i>Energy Storage Materials</i> , 2022 , 49, 1-10	19.4	0
2	A bottom-up strategy to surface assembly: Second growth from metal-rich embryos. <i>Materials Chemistry and Physics</i> , 2018 , 204, 228-235	4.4	
1	Stability of CH ₄ , CO ₂ , and H ₂ S in two-dimensional clathrate hydrates. <i>Computational Materials Science</i> , 2021 , 188, 110174	3.2	