Xue Wenjuan

List of Publications by Year in descending order

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#	Article	IF	CITATIONS
1	Synergistic effect of the metal-support interaction and interfacial oxygen vacancy for CO2 hydrogenation to methanol over Ni/In2O3 catalyst: A theoretical study. Journal of Energy Chemistry, 2022, 65, 623-629.	12.9	51
2	Effects of hydroxylation on the acidic and basic strengths of anatase TiO ₂ surfaces. Molecular Simulation, 2022, 48, 829-843.	2.0	2
3	CO Oxidation over HKUST-1 Catalysts: The Role of Defective Sites. Journal of Physical Chemistry C, 2022, 126, 9652-9664.	3.1	2
4	Unveiling Secondary-Ion-Promoted Catalytic Properties of Cu-SSZ-13 Zeolites for Selective Catalytic Reduction of NO <i>_x</i> . Journal of the American Chemical Society, 2022, 144, 12816-12824.	13.7	51
5	Monodentate AlEgen Anchored on Metalâ€Organic Framework for Fast Fluorescence Sensing of Phosphate. Chinese Journal of Chemistry, 2021, 39, 99-105.	4.9	21
6	Density Functional Theory Study on the Morphology Evolution of Hydroxylated β-Cristobalite Silica and Desilication in the Presence of Methanol. Journal of Physical Chemistry C, 2021, 125, 7868-7879.	3.1	6
7	Structural and Hydrolytic Stability of Coordinatively Unsaturated Metal–Organic Frameworks M ₃ (BTC) ₂ (M = Cu, Co, Mn, Ni, and Zn): A Combined DFT and Experimental Study. Journal of Physical Chemistry C, 2021, 125, 5832-5847.	3.1	11
8	Rational Design of Synergistic Active Sites for Catalytic Ethene/2-Butene Cross-Metathesis in a Rhenium-Doped Y Zeolite Catalyst. ACS Catalysis, 2021, 11, 3530-3540.	11.2	9
9	Polymer-supported ultra-thin ZIF-67 membrane through in situ interface self-repair. Journal of Membrane Science, 2021, 625, 119139.	8.2	45
10	<scp>Airâ€Steam</scp> Etched Construction of Hierarchically Porous <scp>Metalâ€Organic</scp> Frameworks. Chinese Journal of Chemistry, 2021, 39, 1538-1544.	4.9	13
11	Theoretical Insights into CO Oxidation over MOF-808-Encapsulated Single-Atom Metal Catalysts. Journal of Physical Chemistry C, 2021, 125, 17097-17108.	3.1	19
12	Self-adaptive dual-metal-site pairs in metal-organic frameworks for selective CO2 photoreduction to CH4. Nature Catalysis, 2021, 4, 719-729.	34.4	406
13	Morphology controlled synthesis of α-Fe2O3-x with benzimidazole-modified Fe-MOFs for enhanced photo-Fenton-like catalysis. Applied Catalysis B: Environmental, 2021, 291, 120129.	20.2	105
14	Water: A promoter of ammonia selective catalytic reduction over copper-exchanged LTA zeolites. Applied Catalysis B: Environmental, 2021, 294, 120244.	20.2	20
15	Rigidifying induced fluorescence enhancement in 2D porous covalent triazine framework nanosheets for the simultaneously luminous detection and adsorption removal of antibiotics. Chemical Engineering Journal, 2020, 384, 123382.	12.7	83
16	Theoretical Insights into the Initial Hydrolytic Breakdown of HKUST-1. Journal of Physical Chemistry C, 2020, 124, 1991-2001.	3.1	30
17	Postsynthetic Oxidation of the Coordination Site in a Heterometallic Metal–Organic Framework: Tuning Catalytic Behaviors. Chemistry of Materials, 2020, 32, 5192-5199.	6.7	20
18	Efficient separation of vitamins mixture in aqueous solution using a stable zirconium-based metal-organic framework. Journal of Colloid and Interface Science, 2019, 555, 714-721.	9.4	18

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19	Ultramicroporous Metal–Organic Framework with Polar Groups for Efficiently Recovering Propylene from Polypropylene Off-Gas. Industrial & Engineering Chemistry Research, 2019, 58, 14333-14339.	3.7	6
20	IL-induced formation of dynamic complex iodide anions in IL@MOF composites for efficient iodine capture. Journal of Materials Chemistry A, 2019, 7, 18324-18329.	10.3	91
21	Synergy Effect of Pore Structure and Amount of Carboxyl Site for Effective Removal of Pb ²⁺ in Metal–Organic Frameworks. Journal of Chemical & Engineering Data, 2019, 64, 2728-2735.	1.9	36
22	Effects of ionic liquid dispersion in metal-organic frameworks and covalent organic frameworks on CO2 capture: A computational study. Chemical Engineering Science, 2016, 140, 1-9.	3.8	53
23	lonic Liquid/Metal–Organic Framework Composites for H ₂ S Removal from Natural Gas: A Computational Exploration. Journal of Physical Chemistry C, 2015, 119, 3674-3683.	3.1	86