

# Xue Wenjuan

## List of Publications by Year in descending order

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23  
papers

1,184  
citations

516710

16  
h-index

642732

23  
g-index

23  
all docs

23  
docs citations

23  
times ranked

940  
citing authors

#	ARTICLE	IF	CITATIONS
1	Self-adaptive dual-metal-site pairs in metal-organic frameworks for selective CO <sub>2</sub> photoreduction to CH <sub>4</sub> . <i>Nature Catalysis</i> , 2021, 4, 719-729.	34.4	406
2	Morphology controlled synthesis of $\gamma$ -Fe <sub>2</sub> O <sub>3-x</sub> with benzimidazole-modified Fe-MOFs for enhanced photo-Fenton-like catalysis. <i>Applied Catalysis B: Environmental</i> , 2021, 291, 120129.	20.2	105
3	IL-induced formation of dynamic complex iodide anions in IL@MOF composites for efficient iodine capture. <i>Journal of Materials Chemistry A</i> , 2019, 7, 18324-18329.	10.3	91
4	Ionic Liquid/Metal-Organic Framework Composites for H <sub>2</sub> S Removal from Natural Gas: A Computational Exploration. <i>Journal of Physical Chemistry C</i> , 2015, 119, 3674-3683.	3.1	86
5	Rigidifying induced fluorescence enhancement in 2D porous covalent triazine framework nanosheets for the simultaneously luminous detection and adsorption removal of antibiotics. <i>Chemical Engineering Journal</i> , 2020, 384, 123382.	12.7	83
6	Effects of ionic liquid dispersion in metal-organic frameworks and covalent organic frameworks on CO <sub>2</sub> capture: A computational study. <i>Chemical Engineering Science</i> , 2016, 140, 1-9.	3.8	53
7	Synergistic effect of the metal-support interaction and interfacial oxygen vacancy for CO <sub>2</sub> hydrogenation to methanol over Ni/In <sub>2</sub> O <sub>3</sub> catalyst: A theoretical study. <i>Journal of Energy Chemistry</i> , 2022, 65, 623-629.	12.9	51
8	Unveiling Secondary-Ion-Promoted Catalytic Properties of Cu-SSZ-13 Zeolites for Selective Catalytic Reduction of NO <sub>x</sub> . <i>Journal of the American Chemical Society</i> , 2022, 144, 12816-12824.	13.7	51
9	Polymer-supported ultra-thin ZIF-67 membrane through in situ interface self-repair. <i>Journal of Membrane Science</i> , 2021, 625, 119139.	8.2	45
10	Synergy Effect of Pore Structure and Amount of Carboxyl Site for Effective Removal of Pb <sup>2+</sup> in Metal-Organic Frameworks. <i>Journal of Chemical &amp; Engineering Data</i> , 2019, 64, 2728-2735.	1.9	36
11	Theoretical Insights into the Initial Hydrolytic Breakdown of HKUST-1. <i>Journal of Physical Chemistry C</i> , 2020, 124, 1991-2001.	3.1	30
12	Monodentate AlEgen Anchored on Metal-Organic Framework for Fast Fluorescence Sensing of Phosphate. <i>Chinese Journal of Chemistry</i> , 2021, 39, 99-105.	4.9	21
13	Postsynthetic Oxidation of the Coordination Site in a Heterometallic Metal-Organic Framework: Tuning Catalytic Behaviors. <i>Chemistry of Materials</i> , 2020, 32, 5192-5199.	6.7	20
14	Water: A promoter of ammonia selective catalytic reduction over copper-exchanged LTA zeolites. <i>Applied Catalysis B: Environmental</i> , 2021, 294, 120244.	20.2	20
15	Theoretical Insights into CO Oxidation over MOF-808-Encapsulated Single-Atom Metal Catalysts. <i>Journal of Physical Chemistry C</i> , 2021, 125, 17097-17108.	3.1	19
16	Efficient separation of vitamins mixture in aqueous solution using a stable zirconium-based metal-organic framework. <i>Journal of Colloid and Interface Science</i> , 2019, 555, 714-721.	9.4	18
17	Steam Etched Construction of Hierarchically Porous Metal-Organic Frameworks. <i>Chinese Journal of Chemistry</i> , 2021, 39, 1538-1544.	4.9	13
18	Structural and Hydrolytic Stability of Coordinatively Unsaturated Metal-Organic Frameworks M <sub>3</sub> (BTC) <sub>2</sub> (M = Cu, Co, Mn, Ni, and Zn): A Combined DFT and Experimental Study. <i>Journal of Physical Chemistry C</i> , 2021, 125, 5832-5847.	3.1	11

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19	Rational Design of Synergistic Active Sites for Catalytic Ethene/2-Butene Cross-Metathesis in a Rhenium-Doped $\gamma$ Zeolite Catalyst. <i>ACS Catalysis</i> , 2021, 11, 3530-3540.	11.2	9
20	Ultramicroporous Metal-Organic Framework with Polar Groups for Efficiently Recovering Propylene from Polypropylene Off-Gas. <i>Industrial &amp; Engineering Chemistry Research</i> , 2019, 58, 14333-14339.	3.7	6
21	Density Functional Theory Study on the Morphology Evolution of Hydroxylated $\beta$ -Cristobalite Silica and Desilication in the Presence of Methanol. <i>Journal of Physical Chemistry C</i> , 2021, 125, 7868-7879.	3.1	6
22	Effects of hydroxylation on the acidic and basic strengths of anatase $\text{TiO}_2$ surfaces. <i>Molecular Simulation</i> , 2022, 48, 829-843.	2.0	2
23	CO Oxidation over HKUST-1 Catalysts: The Role of Defective Sites. <i>Journal of Physical Chemistry C</i> , 2022, 126, 9652-9664.	3.1	2