

# Haozhi Wang

## List of Publications by Year in descending order

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Version: 2024-02-01

20  
papers

1,681  
citations

567281

15  
h-index

752698

20  
g-index

20  
all docs

20  
docs citations

20  
times ranked

2005  
citing authors

#	ARTICLE	IF	CITATIONS
1	Phase Transfer of Mo <sub>2</sub> C Induced by Boron Doping to Boost Nitrogen Reduction Reaction Catalytic Activity. <i>Advanced Functional Materials</i> , 2022, 32, .	14.9	51
2	Boosting Energy Efficiency and Stability of Li <sup>+</sup> /CO <sub>2</sub> Batteries via Synergy between Ru Atom Clusters and Single <sup>+</sup> Atom Ru <sup>+</sup> N <sub>4</sub> sites in the Electrocatalyst Cathode. <i>Advanced Materials</i> , 2022, 34, e2200559.	21.0	83
3	Ni-Doped Mo <sub>2</sub> C Anchored on Graphitized Porous Carbon for Boosting Electrocatalytic N <sub>2</sub> Reduction. <i>ACS Applied Materials &amp; Interfaces</i> , 2022, 14, 17273-17281.	8.0	12
4	Designing Nanoporous Coral <sup>+</sup> -Like Pt Nanowires Architecture for Methanol and Ammonia Oxidation Reactions. <i>Advanced Functional Materials</i> , 2022, 32, .	14.9	27
5	Cobalt sulfides constructed heterogeneous interfaces decorated on N,S-codoped carbon nanosheets as a highly efficient bifunctional oxygen electrocatalyst. <i>Journal of Materials Chemistry A</i> , 2021, 9, 13926-13935.	10.3	27
6	Accelerated exciton dissociation and electron extraction across the metallic sulfide <sup>+</sup> -carbon nitride ohmic interface for efficient photocatalytic hydrogen production. <i>Journal of Materials Chemistry A</i> , 2021, 9, 16522-16531.	10.3	24
7	Strain <sup>+</sup> -Dependent Band Structures and Electronic Properties in Sb/Bi Lateral Heterostructures Calculated by First Principles. <i>Physica Status Solidi - Rapid Research Letters</i> , 2021, 15, 2100148.	2.4	9
8	Band and optical properties of arsenene and antimonene lateral heterostructure by first-principles calculations. <i>Physica E: Low-Dimensional Systems and Nanostructures</i> , 2021, 134, 114933.	2.7	17
9	Mechanistic understanding of ethane dehydrogenation and aromatization over Zn/ZSM-5: effects of Zn modification and CO <sub>2</sub> co-reactant. <i>Catalysis Science and Technology</i> , 2020, 10, 8359-8373.	4.1	17
10	Selective electroreduction of CO <sub>2</sub> to acetone by single copper atoms anchored on N-doped porous carbon. <i>Nature Communications</i> , 2020, 11, 2455.	12.8	265
11	Enhanced heterogeneous activation of peroxydisulfate by Co and N codoped porous carbon for degradation of organic pollutants: the synergism between Co and N. <i>Environmental Science: Nano</i> , 2019, 6, 399-410.	4.3	129
12	Comparative computational study of CO <sub>2</sub> dissociation and hydrogenation over Fe-M (M = Pd, Ni, Co) bimetallic catalysts: The effect of surface metal content. <i>Journal of CO<sub>2</sub> Utilization</i> , 2019, 29, 179-195.	6.8	17
13	Origin of Pd-Cu bimetallic effect for synergetic promotion of methanol formation from CO <sub>2</sub> hydrogenation. <i>Journal of Catalysis</i> , 2019, 369, 21-32.	6.2	80
14	A short review of recent advances in CO <sub>2</sub> hydrogenation to hydrocarbons over heterogeneous catalysts. <i>RSC Advances</i> , 2018, 8, 7651-7669.	3.6	499
15	Mechanistic Understanding of Alloy Effect and Water Promotion for Pd-Cu Bimetallic Catalysts in CO <sub>2</sub> Hydrogenation to Methanol. <i>ACS Catalysis</i> , 2018, 8, 4873-4892.	11.2	171
16	Facet effect on CO <sub>2</sub> adsorption, dissociation and hydrogenation over Fe catalysts: Insight from DFT. <i>Journal of CO<sub>2</sub> Utilization</i> , 2018, 26, 160-170.	6.8	35
17	DFT insight into the effect of potassium on the adsorption, activation and dissociation of CO <sub>2</sub> over Fe-based catalysts. <i>Physical Chemistry Chemical Physics</i> , 2018, 20, 14694-14707.	2.8	40
18	Mechanistic Insight into C <sup>+</sup> -C Coupling over Fe <sup>+</sup> -Cu Bimetallic Catalysts in CO <sub>2</sub> Hydrogenation. <i>Journal of Physical Chemistry C</i> , 2017, 121, 13164-13174.	3.1	91

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19	Computational Investigation of Fe-Cu Bimetallic Catalysts for CO <sub>2</sub> Hydrogenation. Journal of Physical Chemistry C, 2016, 120, 9364-9373.	3.1	49
20	A computational study of adsorption and activation of CO <sub>2</sub> and H <sub>2</sub> over Fe(1 0 0) surface. Journal of CO <sub>2</sub> Utilization, 2016, 15, 107-114.	6.8	38