

Peng 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.

22

papers

298

citations

10

h-index

17

g-index

22

ext. papers

430

ext. citations

4.3

avg, IF

3.36

L-index

#	Paper	IF	Citations
22	Two-Dimensional π -Conjugated Metal Bis(dithiolene) Complex Nanosheets as Selective Catalysts for Oxygen Reduction Reaction. <i>Journal of Physical Chemistry C</i> , 2015 , 119, 28028-28037	3.8	64
21	From two-dimension to one-dimension: the curvature effect of silicon-doped graphene and carbon nanotubes for oxygen reduction reaction. <i>Physical Chemistry Chemical Physics</i> , 2014 , 16, 17479-86	3.6	42
20	Mechanistic analysis of multiple processes controlling solar-driven HO ₂ synthesis using engineered polymeric carbon nitride. <i>Nature Communications</i> , 2021 , 12, 3701	17.4	35
19	Size effect of oxygen reduction reaction on nitrogen-doped graphene quantum dots. <i>RSC Advances</i> , 2018 , 8, 531-536	3.7	23
18	Oxygen reduction reaction on nitrogen-doped graphene nanoribbons: A density functional theory study. <i>Chemical Physics Letters</i> , 2016 , 663, 123-127	2.5	21
17	Active sites and mechanism of the direct conversion of methane and carbon dioxide to acetic acid over the zinc-modified H-ZSM-5 zeolite. <i>Catalysis Science and Technology</i> , 2019 , 9, 6297-6307	5.5	15
16	First-principles design of bifunctional oxygen reduction and evolution catalysts through bimetallic centers in metal-organic frameworks. <i>Catalysis Science and Technology</i> , 2018 , 8, 3666-3674	5.5	13
15	Oxygen reduction reaction on M-S ₄ embedded graphene: A density functional theory study. <i>Chemical Physics Letters</i> , 2015 , 641, 112-116	2.5	12
14	Adjusting the electronic properties of silicon carbide nanoribbons by introducing edge functionalization. <i>RSC Advances</i> , 2014 , 4, 35042-35047	3.7	12
13	Curvature effect of O ₂ adsorption and dissociation on SiC nanotubes and nanosheet. <i>Chemical Physics Letters</i> , 2015 , 619, 92-96	2.5	11
12	Transition metal-doped borophene as potential oxygen and hydrogen evolution electrocatalyst: A density functional theory study. <i>Catalysis Communications</i> , 2020 , 144, 106090	3.2	10
11	Copper-Carbon: An Efficient Catalyst for Oxygen Reduction. <i>ACS Applied Energy Materials</i> , 2019 , 2, 6295-6301	6.3	8
10	Metal-Organic Framework Derived N/C Supported Austenite Nanoparticles as Efficient Oxygen Reduction Catalysts. <i>ChemNanoMat</i> , 2019 , 5, 525-530	3.5	7
9	Enhanced electrocatalytic activity of nitrogen-doped olympicene/graphene hybrids for the oxygen reduction reaction. <i>Physical Chemistry Chemical Physics</i> , 2016 , 18, 22799-804	3.6	5
8	Oxygen reduction reaction on M ₃ (hexaiminobenzene) ₂ : A density function theory study. <i>Catalysis Communications</i> , 2018 , 115, 17-20	3.2	5
7	Metal-bipyridine complexes as electrocatalysts for the reduction of CO: a density functional theory study. <i>Physical Chemistry Chemical Physics</i> , 2019 , 21, 23742-23748	3.6	5
6	Bimetallic Cu/Zn Co-Doped Porous N/C as Efficient Catalysts for Oxygen Reduction Reaction and Oxidation of 1,2-Propanediol. <i>ChemCatChem</i> , 2020 , 12, 584-592	5.2	4

5	Rapid pollutant degradation by peroxymonosulfate an unusual mediated-electron transfer pathway under spatial-confinement.. <i>RSC Advances</i> , 2022 , 12, 5236-5244	3.7	2
4	Mechanistic studies of oxygen reduction and evolution reactions on Ni ₃ S ₂ surfaces. <i>Applied Catalysis A: General</i> , 2021 , 624, 118324	5.1	2
3	The effects of surface group functionalization and strain on the electronic structures of two-dimensional silicon carbide. <i>Chemical Physics Letters</i> , 2015 , 628, 60-65	2.5	1
2	Transition metal-doped tetra-MoN ₂ monolayers as an electrochemical catalyst for CO ₂ reduction: A density functional theory study. <i>Catalysis Communications</i> , 2021 , 149, 106212	3.2	1
1	Computational screening of transition-metal doped boron nanotubes as efficient electrocatalysts for water splitting.. <i>RSC Advances</i> , 2022 , 12, 6841-6847	3.7	