

Zhaoming Fu

List of Publications by Year in descending order

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#	ARTICLE	IF	CITATIONS
1	Ordered Macroporous Superstructure of Nitrogen-Doped Nanoporous Carbon Implanted with Ultrafine Ru Nanoclusters for Efficient pH-Universal Hydrogen Evolution Reaction. <i>Advanced Materials</i> , 2021, 33, e2006965.	21.0	213
2	MXene (Ti ₃ C ₂ T _x) and Carbon Nanotube Hybrid-Supported Platinum Catalysts for the High-Performance Oxygen Reduction Reaction in PEMFC. <i>ACS Applied Materials & Interfaces</i> , 2020, 12, 19539-19546.	8.0	67
3	The electronic transport properties of transition-metal dichalcogenide lateral heterojunctions. <i>Journal of Materials Chemistry C</i> , 2016, 4, 10962-10966.	5.5	59
4	Understanding the Activity of Single-Atom Catalysis from Frontier Orbitals. <i>Physical Review Letters</i> , 2020, 125, 156001.	7.8	53
5	Strong enhancement of spin ordering by site magnetic ions in the ferrimagnet CaC ₂ F ₃ . <i>Physical Review Letters</i> , 2020, 125, 156001.	7.8	44
6	Large-Scale Domain Engineering in Two-Dimensional Ferroelectric CuInP ₂ S ₆ via Giant Flexoelectric Effect. <i>Nano Letters</i> , 2022, 22, 3275-3282.	9.1	43
7	Repairing single and double atomic vacancies in a C ₃ N monolayer with CO or NO molecules: a first-principles study. <i>Physical Chemistry Chemical Physics</i> , 2018, 20, 13517-13527.	2.8	41
8	Tailoring the Electronic Structure of Transition Metals by the V ₂ C MXene Support: Excellent Oxygen Reduction Performance Triggered by Metal-Support Interactions. <i>ACS Applied Materials & Interfaces</i> , 2020, 12, 28206-28216.	8.0	39
9	The rectifying and negative differential resistance effects in graphene/h-BN nanoribbon heterojunctions. <i>Physical Chemistry Chemical Physics</i> , 2016, 18, 27976-27980.	2.8	36
10	Electronic transport properties of the first all-boron fullerene B ₄₀ and its metallofullerene Sr@B ₄₀ . <i>Physical Chemistry Chemical Physics</i> , 2016, 18, 12024-12028.	2.8	35
11	Gap Symmetry of the Heavy Fermion Superconductor CeCu ₂ Si ₂ at Ambient Pressure. <i>Physical Review Letters</i> , 2018, 120, 217001.	7.8	32
12	Intrinsic negative differential resistance characteristics in zigzag boron nitride nanoribbons. <i>RSC Advances</i> , 2014, 4, 46934-46939.	3.6	27
13	Heavy fermion behavior in the quasi-one-dimensional Kondo lattice CeCo ₂ Ga ₈ . <i>Npj Quantum Materials</i> , 2017, 2, .	5.2	27
14	Influence of electric fields on the structure and structure transition of water confined in a carbon nanotube. <i>Journal of Chemical Physics</i> , 2014, 140, 154508.	3.0	26
15	Tailoring tantalum doping into a perovskite ferrite to obtain a highly active and stable anode for solid oxide fuel cells. <i>Journal of Materials Chemistry A</i> , 2020, 8, 18778-18791.	10.3	24
16	Spin-dependent electronic transport properties of zigzag silicon carbon nanoribbon. <i>RSC Advances</i> , 2015, 5, 107136-107141.	3.6	22
17	Dopant segregation and CO adsorption on doped Fe ₃ O ₄ (111) surfaces: A first-principle study. <i>Journal of Catalysis</i> , 2018, 364, 291-296.	6.2	22
18	Effects of Sm doping content on the ionic conduction of CeO ₂ in SOFCs from first principles. <i>Applied Physics Letters</i> , 2017, 111, .	3.3	20

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19	Strong metal–support interactions impart activity in the oxygen reduction reaction: Au monolayer on Mo ₂ C (MXene). <i>Journal of Physics Condensed Matter</i> , 2018, 30, 475201.	1.8	20
20	Effects of a TiC substrate on the catalytic activity of Pt for NO reduction. <i>Physical Chemistry Chemical Physics</i> , 2016, 18, 13304-13309.	2.8	17
21	Substrate effects on the in-plane ferroelectric polarization of two-dimensional SnTe. <i>Physical Review B</i> , 2019, 99, .	3.2	17
22	Importance of oxygen spillover for fuel oxidation on Ni/YSZ anodes in solid oxide fuel cells. <i>Physical Chemistry Chemical Physics</i> , 2014, 16, 8536.	2.8	16
23	Comparative Study of C ₃ N- and Graphene-Supported Single-Atom Pt. <i>Journal of Physical Chemistry C</i> , 2019, 123, 5731-5735.	3.1	12
24	Rectifications in organic single-molecule diodes alkanethiolate-terminated heterocyclics. <i>Physics Letters, Section A: General, Atomic and Solid State Physics</i> , 2016, 380, 923-926.	2.1	11
25	Tuning the Physical and Chemical Properties of 2D InSe with Interstitial Boron Doping: A First-Principles Study. <i>Journal of Physical Chemistry C</i> , 2017, 121, 28312-28316.	3.1	11
26	The magnetism and spin-dependent electronic transport properties of boron nitride atomic chains. <i>Journal of Chemical Physics</i> , 2016, 145, 044301.	3.0	9
27	Interface-induced ferroelectric domains and charged domain walls in Bi ₂ FeO ₃ superlattices. <i>Physical Review B</i> , 2021, 103, .	6.1	9
28	Understanding the correlation between the electronic structure and catalytic behavior of TiC(001) and TiN(001) surfaces: DFT study. <i>Applied Surface Science</i> , 2019, 494, 57-62.	6.1	8
29	An electronic perturbation in TiC supported platinum monolayer catalyst for enhancing water–gas shift performance: DFT study. <i>Journal of Physics Condensed Matter</i> , 2019, 31, 305201.	1.8	8
30	Synthesis, Structure, and Properties of the Layered Oxyselenide Ba ₂ CuO ₂ Cu ₂ Se ₂ . <i>Inorganic Chemistry</i> , 2018, 57, 5108-5113.	4.0	5
31	The growth modes of graphene in the initial stage of a chemical vapor-deposition process. <i>RSC Advances</i> , 2016, 6, 91157-91162.	3.6	4