

# Zhaoming Fu

## List of Publications by Year in descending order

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31  
papers

977  
citations

394421

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434195

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times ranked

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citing authors

#	ARTICLE	IF	CITATIONS
1	Large-Scale Domain Engineering in Two-Dimensional Ferroelectric $\text{CuInP}_{2}\text{S}_{6}$ via Giant Flexoelectric Effect. <i>Nano Letters</i> , 2022, 22, 3275-3282.	9.1	43
2	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
3	Interface-induced ferroelectric domains and charged domain walls in $\text{BiFeO}_{3}$ superlattices. <i>Physical Review B</i> , 2021, 103, .	7.8	53
4	Understanding the Activity of Single-Atom Catalysis from Frontier Orbitals. <i>Physical Review Letters</i> , 2020, 125, 156001.	7.8	53
5	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
6	Tailoring the Electronic Structure of Transition Metals by the $\text{V}_{2}\text{C}$ MXene Support: Excellent Oxygen Reduction Performance Triggered by Metal-Support Interactions. <i>ACS Applied Materials &amp; Interfaces</i> , 2020, 12, 28206-28216.	8.0	39
7	MXene ( $\text{Ti}_{3}\text{C}_{2}\text{T}_{x}$ ) and Carbon Nanotube Hybrid-Supported Platinum Catalysts for the High-Performance Oxygen Reduction Reaction in PEMFC. <i>ACS Applied Materials &amp; Interfaces</i> , 2020, 12, 19539-19546.	8.0	67
8	Understanding the correlation between the electronic structure and catalytic behavior of $\text{TiC}(001)$ and $\text{TiN}(001)$ surfaces: DFT study. <i>Applied Surface Science</i> , 2019, 494, 57-62.	6.1	8
9	Substrate effects on the in-plane ferroelectric polarization of two-dimensional $\text{SnTe}$ . <i>Physical Review B</i> , 2019, 99, .	3.2	17
10	Comparative Study of $\text{C}_{3}\text{N}$ - and Graphene-Supported Single-Atom Pt. <i>Journal of Physical Chemistry C</i> , 2019, 123, 5731-5735.	3.1	12
11	An electronic perturbation in $\text{TiC}$ 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
12	Repairing single and double atomic vacancies in a $\text{C}_{3}\text{N}$ monolayer with CO or NO molecules: a first-principles study. <i>Physical Chemistry Chemical Physics</i> , 2018, 20, 13517-13527.	2.8	41
13	Synthesis, Structure, and Properties of the Layered Oxyselenide $\text{Ba}_{2}\text{CuO}_{2}\text{Cu}_{2}\text{Se}_{2}$ . <i>Inorganic Chemistry</i> , 2018, 57, 5108-5113.	4.0	5
14	Strong metal-support interactions impart activity in the oxygen reduction reaction: Au monolayer on $\text{Mo}_{2}\text{C}$ (MXene). <i>Journal of Physics Condensed Matter</i> , 2018, 30, 475201.	1.8	20
15	Gap Symmetry of the Heavy Fermion Superconductor $\text{CeCu}_{2}\text{Si}_{2}$ at Ambient Pressure. <i>Physical Review Letters</i> , 2018, 120, 217001.	7.8	32
16	Dopant segregation and CO adsorption on doped $\text{Fe}_{3}\text{O}_{4}$ ( $1\text{e}^{-}1\text{e}^{-}1$ ) surfaces: A first-principle study. <i>Journal of Catalysis</i> , 2018, 364, 291-296.	6.2	22
17	Effects of Sm doping content on the ionic conduction of $\text{CeO}_{2}$ in SOFCs from first principles. <i>Applied Physics Letters</i> , 2017, 111, .	3.3	20
18	Tuning the Physical and Chemical Properties of 2D $\text{InSe}$ with Interstitial Boron Doping: A First-Principles Study. <i>Journal of Physical Chemistry C</i> , 2017, 121, 28312-28316.	3.1	11

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19	Heavy fermion behavior in the quasi-one-dimensional Kondo lattice CeCo <sub>2</sub> Ga <sub>8</sub> . Npj Quantum Materials, 2017, 2, .	5.2	27
20	The magnetism and spin-dependent electronic transport properties of boron nitride atomic chains. Journal of Chemical Physics, 2016, 145, 044301.	3.0	9
21	Electronic transport properties of the first all-boron fullerene B <sub>40</sub> and its metallofullerene Sr@B <sub>40</sub> . Physical Chemistry Chemical Physics, 2016, 18, 12024-12028.	2.8	35
22	Rectifications in organic single-molecule diodes alkanethiolate-terminated heterocyclics. Physics Letters, Section A: General, Atomic and Solid State Physics, 2016, 380, 923-926.	2.1	11
23	Effects of a TiC substrate on the catalytic activity of Pt for NO reduction. Physical Chemistry Chemical Physics, 2016, 18, 13304-13309.	2.8	17
24	The growth modes of graphene in the initial stage of a chemical vapor-deposition process. RSC Advances, 2016, 6, 91157-91162.	3.6	4
25	Strong enhancement of spin ordering by A-site magnetic ions in the ferrimagnet Ca <sub>3</sub> Co <sub>2</sub> Si <sub>2</sub> O <sub>12</sub> . Physical Chemistry Chemical Physics, 2016, 18, 27976-27980.	2.8	36
26	The rectifying and negative differential resistance effects in graphene/h-BN nanoribbon heterojunctions. Physical Chemistry Chemical Physics, 2016, 18, 27976-27980.	2.8	36
27	The electronic transport properties of transition-metal dichalcogenide lateral heterojunctions. Journal of Materials Chemistry C, 2016, 4, 10962-10966.	5.5	59
28	Spin-dependent electronic transport properties of zigzag silicon carbon nanoribbon. RSC Advances, 2015, 5, 107136-107141.	3.6	22
29	Influence of electric fields on the structure and structure transition of water confined in a carbon nanotube. Journal of Chemical Physics, 2014, 140, 154508.	3.0	26
30	Intrinsic negative differential resistance characteristics in zigzag boron nitride nanoribbons. RSC Advances, 2014, 4, 46934-46939.	3.6	27
31	Importance of oxygen spillover for fuel oxidation on Ni/YSZ anodes in solid oxide fuel cells. Physical Chemistry Chemical Physics, 2014, 16, 8536.	2.8	16