

# Hao-Qing Ji

## List of Publications by Year in descending order

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29  
papers

1,842  
citations

304743

22  
h-index

477307

29  
g-index

29  
all docs

29  
docs citations

29  
times ranked

1965  
citing authors

#	ARTICLE	IF	CITATIONS
1	Over 56.55% Faradaic efficiency of ambient ammonia synthesis enabled by positively shifting the reaction potential. <i>Nature Communications</i> , 2019, 10, 341.	12.8	412
2	Proton-filtering covalent organic frameworks with superior nitrogen penetration flux promote ambient ammonia synthesis. <i>Nature Catalysis</i> , 2021, 4, 322-331.	34.4	216
3	Facilitating nitrogen accessibility to boron-rich covalent organic frameworks via electrochemical excitation for efficient nitrogen fixation. <i>Nature Communications</i> , 2019, 10, 3898.	12.8	191
4	Salting-out effect promoting highly efficient ambient ammonia synthesis. <i>Nature Communications</i> , 2021, 12, 3198.	12.8	105
5	A new high ionic conductive gel polymer electrolyte enables highly stable quasi-solid-state lithium sulfur battery. <i>Energy Storage Materials</i> , 2019, 22, 256-264.	18.0	89
6	Single-Atom Iron as Lithiophilic Site To Minimize Lithium Nucleation Overpotential for Stable Lithium Metal Full Battery. <i>ACS Applied Materials &amp; Interfaces</i> , 2019, 11, 32008-32014.	8.0	64
7	Altering the rate-determining step over cobalt single clusters leading to highly efficient ammonia synthesis. <i>National Science Review</i> , 2021, 8, nwaal136.	9.5	64
8	Boosting the Optimization of Lithium Metal Batteries by Molecular Dynamics Simulations: A Perspective. <i>Advanced Energy Materials</i> , 2020, 10, 2002373.	19.5	56
9	All-Liquid-Phase Reaction Mechanism Enabling Cryogenic Li <sup>+</sup> S Batteries. <i>ACS Nano</i> , 2021, 15, 13847-13856.	14.6	55
10	Molecular Dynamics Simulation of Methane Hydrate Formation and Dissociation in the Clay Pores with Fatty Acids. <i>Journal of Physical Chemistry C</i> , 2018, 122, 1318-1325.	3.1	49
11	Unveiling the Essential Nature of Lewis Basicity in Thermodynamically and Dynamically Promoted Nitrogen Fixation. <i>Advanced Functional Materials</i> , 2020, 30, 2001244.	14.9	49
12	Pyridinic and graphitic nitrogen-enriched carbon paper as a highly active bifunctional catalyst for Zn-air batteries. <i>Electrochimica Acta</i> , 2020, 334, 135562.	5.2	45
13	Single-atom scale metal vacancy engineering in heteroatom-doped carbon for rechargeable zinc-air battery with reduced overpotential. <i>Chemical Engineering Journal</i> , 2020, 393, 124702.	12.7	43
14	In-situ observation as activity descriptor enables rational design of oxygen reduction catalyst for zinc-air battery. <i>Energy Storage Materials</i> , 2020, 27, 226-231.	18.0	42
15	Effects of Asphaltenes on the Formation and Decomposition of Methane Hydrate: A Molecular Dynamics Study. <i>Energy &amp; Fuels</i> , 2016, 30, 5643-5650.	5.1	41
16	Interfacial Microextraction Boosting Nitrogen Feed for Efficient Ambient Ammonia Synthesis in Aqueous Electrolyte. <i>Advanced Functional Materials</i> , 2022, 32, .	14.9	41
17	Microsecond Molecular Dynamics Simulation of Methane Hydrate Formation in Humic-Acid-Amended Sodium Montmorillonite. <i>Energy &amp; Fuels</i> , 2016, 30, 7206-7213.	5.1	35
18	Super lithiophilic SEI derived from quinones electrolyte to guide Li uniform deposition. <i>Energy Storage Materials</i> , 2020, 24, 426-431.	18.0	34

#	ARTICLE	IF	CITATIONS
19	Atomic Metal Vacancy Modulation of Single-Atom Dispersed Co/N/C for Highly Efficient and Stable Air Cathode. ACS Applied Materials & Interfaces, 2020, 12, 15298-15304.	8.0	33
20	Effects of Salt Ions on the Methane Hydrate Formation and Dissociation in the Clay Pore Water and Bulk Water. Energy & Fuels, 2018, 32, 12486-12494.	5.1	32
21	CO <sub>2</sub> and CH <sub>4</sub> Hydrates: Replacement or Cogrowth?. Journal of Physical Chemistry C, 2019, 123, 13401-13409.	3.1	27
22	Rapid leakage responsive and self-healing Li-metal batteries. Chemical Engineering Journal, 2021, 404, 126470.	12.7	26
23	Molecular Mechanisms for Cyclodextrin-Promoted Methane Hydrate Formation in Water. Journal of Physical Chemistry C, 2017, 121, 20967-20975.	3.1	25
24	Molecular Simulations Guided Polymer Electrolyte towards Superior Low-Temperature Solid Lithium-Metal Batteries. ACS Applied Materials & Interfaces, 2021, 13, 48810-48817.	8.0	16
25	Surpassing the Redox Potential Limit of Organic Cathode Materials via Extended $\pi$ -Conjugation of Dioxin. Nano Letters, 2022, 22, 3473-3479.	9.1	14
26	New Type of Dynamically "Solid" Liquid Interconvertible Electrolyte for High-Rate Zn Metal Battery. Nano Letters, 2022, 22, 2898-2906.	9.1	13
27	Unravelling critical role of metal cation engineering in boosting hydrogen evolution reaction activity of molybdenum diselenide. Rare Metals, 2022, 41, 1851-1858.	7.1	10
28	Identifying the Lewis Base Chemistry in Preventing the Deposition of Metal Oxides on Ketone-Enriched Carbon Cathodes for Highly Durable Metal-Air Batteries. ACS Applied Materials & Interfaces, 2020, 12, 3603-3609.	8.0	9
29	Enhanced utilization of active sites of Fe/N/C catalysts by pore-in-pore structures for ultrahigh mass activity. Nanotechnology, 2020, 31, 315401.	2.6	6