## **Erhong Song**

List of Publications by Year in descending order

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#	Article	IF	CITATIONS
1	Engineering Metallic Heterostructure Based on Ni <sub>3</sub> N and 2Mâ€MoS <sub>2</sub> for Alkaline Water Electrolysis with Industryâ€Compatible Current Density and Stability. Advanced Materials, 2022, 34, e2108505.	11.1	104
2	Dynamic coordination transformation of active sites in single-atom MoS <sub>2</sub> catalysts for boosted oxygen evolution catalysis. Energy and Environmental Science, 2022, 15, 2071-2083.	15.6	33
3	Nano gold coupled black titania composites with enhanced surface plasma properties for efficient photocatalytic alkyne reduction. Applied Catalysis B: Environmental, 2022, 309, 121222.	10.8	11
4	Partial‣ingleâ€Atom, Partialâ€Nanoparticle Composites Enhance Water Dissociation for Hydrogen Evolution. Advanced Science, 2021, 8, 2001881.	5.6	85
5	Designing fluorographene with FeN4 and CoN4 moieties for oxygen electrode reaction: A density functional theory study. Applied Surface Science, 2021, 537, 147846.	3.1	23
6	Identifying Metallic Transition-Metal Dichalcogenides for Hydrogen Evolution through Multilevel High-Throughput Calculations and Machine Learning. Journal of Physical Chemistry Letters, 2021, 12, 2102-2111.	2.1	43
7	Mo decoration on graphene edge for nitrogen fixation: A computational investigation. Applied Surface Science, 2021, 568, 150867.	3.1	11
8	Assembling organic–inorganic building blocks for high-capacity electrode design. Materials Horizons, 2021, 8, 1825-1834.	6.4	1
9	DFT Calculation of NO Adsorption on Cr Doped Graphene. Wuji Cailiao Xuebao/Journal of Inorganic Materials, 2021, 36, 1047.	0.6	2
10	Optimized electron occupancy of solid-solution transition metals for suppressing the oxygen evolution of Li <sub>2</sub> MnO <sub>3</sub> . Journal of Materials Chemistry A, 2021, 9, 9337-9346.	5.2	7
11	DFT Calculationof NOAdsorption on CrDoped Graphene. Wuji Cailiao Xuebao/Journal of Inorganic Materials, 2021, , 78.	0.6	0
12	Manipulation on active electronic states of metastable phase β-NiMoO4 for large current density hydrogen evolution. Nature Communications, 2021, 12, 5960.	5.8	86
13	Bond Electronegativity as Hydrogen Evolution Reaction Catalyst Descriptor for Transition Metal (TM) Tj ETQq1 1	0.784314 3.2	4 rgBT /Over
14	Design of Effective Graphene with the TM/O Moiety for the Oxygen Electrode Reaction. ACS Applied Energy Materials, 2020, 3, 260-267.	2.5	24
15	Stacking faults modulation for scattering optimization in GeTe-based thermoelectric materials. Nano Energy, 2020, 68, 104347.	8.2	77
16	Dualâ€Metal Interbonding as the Chemical Facilitator for Singleâ€Atom Dispersions. Advanced Materials, 2020, 32, e2003484.	11.1	90
17	Theoretical Study of Fast Calculation of Damping Loss Factors for Rubber Polymers. Journal of Physical Chemistry Letters, 2020, 11, 6025-6031.	2.1	7
18	The FeN <sub>3</sub> Doped Fluorographene for N <sub>2</sub> Fixation: A Density Functional Theory Study. ChemistrySelect, 2020, 5, 9370-9376.	0.7	3

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19	Liquidâ€Phase Assisted Engineering of Highly Strong SiC Composite Reinforced by Multiwalled Carbon Nanotubes. Advanced Science, 2020, 7, 2002225.	5.6	11
20	Enhanced performance of in-plane transition metal dichalcogenides monolayers by configuring local atomic structures. Nature Communications, 2020, 11, 2253.	5.8	112
21	Transition metal-doped α-borophene as potential oxygen and hydrogen evolution electrocatalyst: A density functional theory study. Catalysis Communications, 2020, 144, 106090.	1.6	20
22	Reducing the charge overpotential of Li–O <sub>2</sub> batteries through band-alignment cathode design. Energy and Environmental Science, 2020, 13, 2540-2548.	15.6	30
23	Predicting Transitionâ€Metal/Benzenehexathiol Systems as Effective Cathodes of Liâ€ <del>S</del> batteries with Conjugate Conductivity and Synergetic Discharge. ChemistrySelect, 2020, 5, 7783-7788.	0.7	5
24	The VN3 embedded graphane with the improved selectivity for nitrogen fixation. Applied Surface Science, 2020, 513, 145855.	3.1	23
25	Cooperative Effect of Multiple Active Sites and Hierarchical Chemical Bonds in Metal–Organic Compounds for Improving Cathode Performance. ACS Energy Letters, 2020, 5, 477-485.	8.8	10
26	Feasibility of N <sub>2</sub> Reduction on the V Anchored 1Tâ^'MoS <sub>2</sub> Monolayer: A Density Functional Theory Study. ChemPhysChem, 2020, 21, 1235-1242.	1.0	14
27	Recent progress in theoretical and computational investigations of structural stability and activity of single-atom electrocatalysts. Progress in Natural Science: Materials International, 2019, 29, 256-264.	1.8	27
28	High strain rate compressive response of the Cf/SiC composite. Ceramics International, 2019, 45, 6812-6818.	2.3	12
29	In-situ liquid cell transmission electron microscopy investigation on oriented attachment of gold nanoparticles. Nature Communications, 2018, 9, 421.	5.8	171
30	Adsorption-energy-based activity descriptors for electrocatalysts in energy storage applications. National Science Review, 2018, 5, 327-341.	4.6	129
31	Auto-optimizing Hydrogen Evolution Catalytic Activity of ReS <sub>2</sub> through Intrinsic Charge Engineering. ACS Nano, 2018, 12, 4486-4493.	7.3	111
32	Ultrafine WC nanoparticles anchored on co-encased, N-doped carbon nanotubes for efficient hydrogen evolution. Energy Storage Materials, 2017, 6, 104-111.	9.5	48
33	Surface synergism of Pd/H 2 Ti 3 O 7 composite nanowires for catalytic and photocatalytic hydrogen production from ammonia borane. International Journal of Hydrogen Energy, 2016, 41, 3428-3435.	3.8	29
34	Revealing the Active Intermediates in the Oxidation of Formic Acid on Au and Pt(111). Chemistry - A European Journal, 2014, 20, 11005-11012.	1.7	28
35	Tuning electronic and magnetic properties of partially hydrogenated graphene by biaxial tensile strain: a computational study. Nanoscale Research Letters, 2014, 9, 491.	3.1	2
36	External electric field induced hydrogen storage/release on calcium-decorated single-layer and bilayer silicene. Physical Chemistry Chemical Physics, 2014, 16, 23985-23992.	1.3	47

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37	Density functional theory calculations of adsorption of hydrogen fluoride on titanium embedded graphene. Thin Solid Films, 2013, 546, 124-127.	0.8	12
38	Titanium Decorated Graphene as CO Detection Sensor. Nanoscience and Nanotechnology Letters, 2013, 5, 198-203.	0.4	10
39	External Electric Field Catalyzed N <sub>2</sub> O Decomposition on Mn-Embedded Graphene. Journal of Physical Chemistry C, 2012, 116, 20342-20348.	1.5	44
40	CO Catalytic Oxidation on Copper-Embedded Graphene. Journal of Physical Chemistry C, 2011, 115, 3678-3683.	1.5	337
41	Critical Role of Interfacial Charge Transfer in Reducing Charge Potential of Li–O2 Battery. Journal of Physical Chemistry C, 0, , .	1.5	1
42	Effective Descriptor for Nitrogen Reduction on Atomic Catalysts. Catalysis Letters, 0, , 1.	1.4	1