Weiwei Sun

List of Publications by Year in descending order

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394421 315739 1,585 39 19 38 citations h-index g-index papers 39 39 39 2597 docs citations times ranked citing authors all docs

#	Article	IF	CITATIONS
1	Direct Synthesis of Largeâ€Area 2D Mo ₂ C on In Situ Grown Graphene. Advanced Materials, 2017, 29, 1700072.	21.0	305
2	In situ atomistic insight into the growth mechanisms of single layer 2D transition metal carbides. Nature Communications, 2018, 9, 2266.	12.8	125
3	Computational Discovery and Design of MXenes for Energy Applications: Status, Successes, and Opportunities. ACS Applied Materials & Energy Applications: 91, 24885-24905.	8.0	105
4	Tracking ion intercalation into layered Ti ₃ C ₂ MXene films across length scales. Energy and Environmental Science, 2020, 13, 2549-2558.	30.8	100
5	Double transition metal MXenes with wide band gaps and novel magnetic properties. Nanoscale, 2018, 10, 11962-11968.	5.6	88
6	Effects of Surface Terminations of 2D Bi ₂ WO ₆ on Photocatalytic Hydrogen Evolution from Water Splitting. ACS Applied Materials & Samp; Interfaces, 2020, 12, 20067-20074.	8.0	78
7	Low-Temperature Synthesis of Single Palladium Atoms Supported on Defective Hexagonal Boron Nitride Nanosheet for Chemoselective Hydrogenation of Cinnamaldehyde. ACS Nano, 2021, 15, 10175-10184.	14.6	77
8	Computational Screening of MXene Electrodes for Pseudocapacitive Energy Storage. Journal of Physical Chemistry C, 2019, 123, 315-321.	3.1	69
9	Edge Segregated Polymorphism in 2D Molybdenum Carbide. Advanced Materials, 2019, 31, e1808343.	21.0	56
10	Surface Reorganization Leads to Enhanced Photocatalytic Activity in Defective BiOCl. Chemistry of Materials, 2018, 30, 5128-5136.	6.7	55
11	Interfacial and electronic properties of heterostructures of MXene and graphene. Physical Review B, 2019, 99, .	3.2	53
12	Investigation of Charge-Transfer Interaction in Mixed Stack Donor–Acceptor Cocrystals Toward Tunable Solid-State Emission Characteristics. Crystal Growth and Design, 2018, 18, 6001-6008.	3.0	51
13	A new 2D monolayer BiXene, M ₂ C (M = Mo, Tc, Os). Nanoscale, 2016, 8, 15753-15762.	5.6	46
14	Synergistic effects between polyvinylpyrrolidone and oxygen vacancies on improving the oxidase-mimetic activity of flower-like CeO ₂ nanozymes. Nanoscale, 2020, 12, 19104-19111.	5.6	37
15	Understanding the electrochemical properties of A $<$ sub $>$ 2 $<$ /sub $>$ MSiO $<$ sub $>$ 4 $<$ /sub $>$ (A = Li and Na; M =) Tj ETQq calculations. Journal of Materials Chemistry A, 2016, 4, 17455-17463.	q1 1 0.784 10.3	1314 rgBT (O 35
16	Multiscale and Multimodal Characterization of 2D Titanium Carbonitride MXene. Advanced Materials Interfaces, 2020, 7, 1902207.	3.7	35
17	Structure and energy of point defects in TiC: An <i>ab initio</i> study. Physical Review B, 2015, 91, .	3.2	23
18	High photon energy spectroscopy of NiO: Experiment and theory. Physical Review B, 2016, 93, .	3.2	22

#	Article	IF	CITATIONS
19	Dynamic stability of the single-layer transition metal dichalcogenides. Computational Materials Science, 2014, 92, 206-212.	3.0	19
20	Self-diffusion of Ti interstitial based point defects and complexes in TiC. Acta Materialia, 2019, 165, 381-387.	7.9	18
21	The role of group III, IV elements in Nb ₄ AC ₃ MAX phases (A = Al, Si, Ga, Ge) and the unusual anisotropic behavior of the electronic and optical properties. Physical Chemistry Chemical Physics, 2017, 19, 15471-15483.	2.8	17
22	Role of correlation and relativistic effects in MAX phases. Journal of Materials Science, 2012, 47, 7615-7620.	3.7	16
23	Investigations on electronic, Fermi surface, Curie temperature and optical properties of Zr 2 CoAl. Journal of Solid State Chemistry, 2017, 247, 97-104.	2.9	15
24	Stabilizing a hexagonal Ru2C via Lifshitz transition under pressure. Applied Physics Letters, 2013, 103, .	3.3	14
25	Anisotropic distortion and Lifshitz transition in <mml:math xmlns:mml="http://www.w3.org/1998/Math/MathML"><mml:mi>α</mml:mi></mml:math> -Hf under pressure. Physical Review B, 2017, 95, .	3.2	14
26	Solid emission color tuning of organic charge transfer cocrystals based on planar π-conjugated donors and TCNB. Journal of Solid State Chemistry, 2019, 272, 96-101.	2.9	14
27	Strain and doping effects on the energetics of hydrogen desorption from the MgH ₂ (001) surface. Europhysics Letters, 2013, 101, 27006.	2.0	13
28	Pushing the limit of thermal conductivity of MAX borides and MABs. Journal of Materials Science and Technology, 2022, 97, 79-88.	10.7	12
29	Combining configurational energies and forces for molecular force field optimization. Journal of Chemical Physics, 2017, 147, 161713.	3.0	11
30	A combined machine learning and density functional theory study of binary Ti-Nb and Ti-Zr alloys: Stability and Young's modulus. Computational Materials Science, 2020, 184, 109830.	3.0	10
31	Improvement in the desorption of H2 from the MgH2 (110) surface by means of doping and mechanical strain. Computational Materials Science, 2014, 86, 165-169.	3.0	9
32	Gluing together metallic and covalent layers to form Ru ₂ C under ambient conditions. Physical Chemistry Chemical Physics, 2015, 17, 9730-9736.	2.8	9
33	Photoinduction of palladium single atoms supported on defect-containing Î ³ -AlOOH nanoleaf for efficient trans-stilbene epoxidation. Chemical Engineering Journal, 2022, 429, 132149.	12.7	8
34	Stability of a new cubic monoxide of Thorium under pressure. Scientific Reports, 2015, 5, 13740.	3.3	7
35	The correlation between N deficiency and the mechanical properties of the Ti2AlNy MAX phase. Journal of the European Ceramic Society, 2020, 40, 2279-2286.	5.7	7
36	Li ₂ FePO ₄ F and its metal-doping for Li-ion batteries: an ab initio study. RSC Advances, 2014, 4, 50195-50201.	3.6	6

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#	Article	IF	CITATIONS
37	Behavior of intrinsic defects in BaF2 under uniaxial compressions: An ab initio investigation. Materials Today Communications, 2021, 28, 102730.	1.9	5
38	Observation of Square-Planar Distortion in Lanthanide-Doped Skutterudite Crystals. Journal of Physical Chemistry C, 2019, 123, 14632-14638.	3.1	1
39	Exchange of Re and Mo atoms in MoS2 driven by Scanning Transmission Electron Microscopy. Microscopy and Microanalysis, 2017, 23, 1702-1703.	0.4	0