

Yongmao Cai

List of Publications by Year in descending order

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#	ARTICLE	IF	CITATIONS
1	Two-Dimensional V_2N MXene Monolayer as a High-Capacity Anode Material for Lithium-Ion Batteries and Beyond: First-Principles Calculations. ACS Omega, 2022, 7, 17756-17764.	1.6	18
2	Understanding the Hydrogen-Bonded Clusters of Ammonia (NH_3) $_n$ ($n = 3-6$): Insights from the Electronic Structure Theory. ACS Omega, 2020, 5, 31724-31729.	1.6	10
3	The Anchoring Effect of 2D Graphdiyne Materials for Lithium-Sulfur Batteries. ACS Omega, 2020, 5, 13424-13429.	1.6	10
4	A low-cost and energy-saving preparation method for silicon derived from rice husks and lithium ion battery applications. Materials Research Express, 2019, 6, 045505.	0.8	2
5	Structural phase transition and bonding properties of high-pressure polymeric CaN_3 . RSC Advances, 2018, 8, 4314-4320.	1.7	14
6	Crystal structures of transition metal pernitrides predicted from first principles. RSC Advances, 2018, 8, 36412-36421.	1.7	15
7	NASICON-Type $Mg_{0.5}Ti_2(PO_4)_3$ Negative Electrode Material Exhibits Different Electrochemical Energy Storage Mechanisms in Na-Ion and Li-Ion Batteries. ACS Applied Materials & Interfaces, 2017, 9, 4709-4718.	4.0	47
8	Stability and electronic properties of two-dimensional silicene and germanene on graphene. Physical Review B, 2013, 88, .	1.1	173
9	First-Principles Calculations on the $LiMSO_4F/MSO_4F$ ($M = Fe, Co, \text{ and } Ni$) Systems. Journal of Physical Chemistry C, 2011, 115, 7032-7037.	1.5	32
10	Pressure-induced phase transformation and magnetism transition in $BaRuO_3$: A first-principles study. Solid State Sciences, 2011, 13, 350-355.	1.5	9
11	Prediction of the phase transition from ferromagnetic perovskite to non-magnetic post-perovskite in $SrRuO_3$: A first-principles study. Solid State Communications, 2011, 151, 798-801.	0.9	5
12	Electrochemical Kinetics of the $Li[Li_{0.23}Co_{0.3}Mn_{0.47}]O_2$ Cathode Material Studied by GITT and EIS. Journal of Physical Chemistry C, 2010, 114, 22751-22757.	1.5	285
13	Charge disproportionation in AlV_2O_4 : A first-principles study. Journal of Alloys and Compounds, 2010, 505, L23-L26.	2.8	8