

Hua Ning

List of Publications by Year in descending order

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Version: 2024-02-01

10
papers

281
citations

1307594

7
h-index

1372567

10
g-index

10
all docs

10
docs citations

10
times ranked

106
citing authors

#	ARTICLE	IF	CITATIONS
1	Roles of in situ-formed NbN and Nb ₂ O ₅ from N-doped Nb ₂ C MXene in regulating the re/hydrogenation and cycling performance of magnesium hydride. <i>Chemical Engineering Journal</i> , 2022, 431, 133985.	12.7	47
2	In situ incorporation of highly dispersed nickel and vanadium trioxide nanoparticles in nanoporous carbon for the hydrogen storage performance enhancement of magnesium hydride. <i>Chemical Engineering Journal</i> , 2022, 446, 137261.	12.7	42
3	Combinations of V ₂ C and Ti ₃ C ₂ MXenes for Boosting the Hydrogen Storage Performances of MgH ₂ . <i>ACS Applied Materials & Interfaces</i> , 2021, 13, 13235-13247.	8.0	111
4	Effect of BaO on Hydrogen Sorption Performance of Mg ₁₇ Al ₁₂ : Experimental and Theoretical Studies. <i>ACS Applied Materials & Interfaces</i> , 2021, 13, 11901-11910.	8.0	2
5	Facile and low-cost synthesis of carbon-supported manganese monoxide nanocomposites and evaluation of their superior catalytic effect toward magnesium hydride. <i>Journal of Alloys and Compounds</i> , 2021, 887, 161380.	5.5	16
6	Synergistic effect of hydrogen absorption on (Ti+Ni), (Ti+V), (Ni+V) doped Mg ₁₇ Al ₁₂ (110) surfaces: A theoretical study. <i>Applied Surface Science</i> , 2020, 514, 145884.	6.1	6
7	Ni catalytic effects for the enhanced hydrogenation properties of Mg ₁₇ Al ₁₂ (110) surface. <i>Applied Surface Science</i> , 2019, 464, 644-650.	6.1	24
8	Hydrogenation properties of Mg ₁₇ Al ₁₂ doped with alkaline-earth metal (Be, Ca, Sr and Ba). <i>Journal of Alloys and Compounds</i> , 2019, 774, 865-872.	5.5	15
9	Effects of Li on hydrogen absorption properties of Mg ₁₇ Al ₁₂ (110) surface: A density functional theory study. <i>International Journal of Hydrogen Energy</i> , 2018, 43, 18330-18338.	7.1	7
10	Hydrogen penetration and diffusion on Mg ₁₇ Al ₁₂ (110) surface: A density functional theory investigation. <i>International Journal of Hydrogen Energy</i> , 2017, 42, 26013-26019.	7.1	11