

Weiqing Jiang

List of Publications by Year in descending order

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#	ARTICLE	IF	CITATIONS
1	Effect of substituting Al for Co on the hydrogen-storage performance of $\text{La}_{0.7}\text{Mg}_{0.3}\text{Ni}_{2.6}\text{Al}_x\text{Co}_{0.5-x}$ ($x=0.0\text{--}0.3$) alloys. <i>International Journal of Hydrogen Energy</i> , 2009, 34, 2986-2991.	7.1	28
2	Rare earth-Mg-Ni-based alloys with superlattice structure for electrochemical hydrogen storage. <i>Journal of Alloys and Compounds</i> , 2021, 887, 161381.	5.5	25
3	Annealing effect on hydrogen storage property of Co-free $\text{La}_{1.8}\text{Ti}_{0.2}\text{MgNi}_{8.7}\text{Al}_{0.3}$ alloy. <i>Journal of Alloys and Compounds</i> , 2013, 565, 37-43.	5.5	22
4	A study on the hydrogen-storage properties of $\text{La}_{2-x}\text{Ti}_x\text{MgNi}_9$ ($x=0.1, 0.2, 0.3, 0.4$) alloys. <i>International Journal of Hydrogen Energy</i> , 2009, 34, 4827-4832.	7.1	20
5	Influence of annealing treatment on the hydrogen storage properties of $\text{La}_{2-x}\text{Ti}_x\text{MgNi}_9$ ($x=0.2, 0.3$) alloys. <i>International Journal of Hydrogen Energy</i> , 2010, 35, 11016-11024.	7.1	20
6	Effect of annealing on the structure and electrochemical properties of $\text{La}_{1.8}\text{Ti}_{0.2}\text{MgNi}_{8.9}\text{Al}_{0.1}$ hydrogen storage alloy. <i>Journal of Power Sources</i> , 2013, 221, 84-89.	7.8	19
7	Dehydrogenation properties of LiBH_4 modified by Mg from first-principles calculations. <i>Journal of Alloys and Compounds</i> , 2018, 735, 668-676.	5.5	17
8	First-principles study on the dehydrogenation characteristics of LiBH_4 modified by Ti. <i>Results in Physics</i> , 2017, 7, 3236-3242.	4.1	13
9	A comparative study on dehydrogenation of Mg-doped LiBH_4 and $\text{Li}_2\text{B}_{12}\text{H}_{12}$ from first-principles calculations. <i>Computational Materials Science</i> , 2018, 154, 187-193.	3.0	8
10	First-principles study on the dehydrogenation of $\text{Li}_4\text{BN}_3\text{H}_{10}$ modified by Co. <i>International Journal of Hydrogen Energy</i> , 2021, 46, 11815-11823.	7.1	7
11	The first-principles investigation on the electronic structure and mechanism of $\text{LiH} + \text{NH}_3 \rightarrow \text{LiNH}_2 + \text{H}_2$ reaction. <i>International Journal of Hydrogen Energy</i> , 2012, 37, 18937-18943.	7.1	5
12	First-principles investigation of dehydrogenation of Cu-doped LiBH_4 . <i>Solid State Communications</i> , 2021, 326, 114184.	1.9	5
13	First-principles investigation of LaMg_2Ni and its hydrides. <i>Scientific Reports</i> , 2020, 10, 12167.	3.3	2
14	Hydrogen storage properties of $\text{La}_{1.8}\text{Ti}_{0.2}\text{MgNi}_{9-x}\text{Co}_x$ ($x = 0, 0.1, 0.3, 0.5$) alloys. <i>Russian Journal of Electrochemistry</i> , 2016, 52, 435-440.	0.9	1
15	The Effects of Sn Element on Hydrogen Storage Characteristics of $\text{Mg}_{2-x}\text{Sn}_x\text{Ni}$ ($x=0, 0.05, 0.1, 0.15, 0.2$) Alloys. <i>Materials Research Society Symposia Proceedings</i> , 2006, 971, 1.	0.1	0