

ShinYoung Kang

List of Publications by Year in descending order

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Version: 2024-02-01

24
papers

2,205
citations

623734

14
h-index

610901

24
g-index

24
all docs

24
docs citations

24
times ranked

3548
citing authors

#	ARTICLE	IF	CITATIONS
1	The structural and chemical origin of the oxygen redox activity in layered and cation-disordered Li-excess cathode materials. <i>Nature Chemistry</i> , 2016, 8, 692-697.	13.6	1,022
2	Nanostructured Metal Hydrides for Hydrogen Storage. <i>Chemical Reviews</i> , 2018, 118, 10775-10839.	47.7	461
3	A Facile Mechanism for Recharging Li_2O_2 in Li^+O_2 Batteries. <i>Chemistry of Materials</i> , 2013, 25, 3328-3336.	6.7	179
4	Nanoscale Stabilization of Sodium Oxides: Implications for Na^+O_2 Batteries. <i>Nano Letters</i> , 2014, 14, 1016-1020.	9.1	162
5	Hierarchically Controlled Inside-Out Doping of Mg Nanocomposites for Moderate Temperature Hydrogen Storage. <i>Advanced Functional Materials</i> , 2017, 27, 1704316.	14.9	72
6	Intrinsic stoichiometry and oxygen-induced p -type conductivity of pyrite FeS_2 . <i>Physical Review B</i> , 2011, 84, .	3.2	65
7	A Mechanistic Analysis of Phase Evolution and Hydrogen Storage Behavior in Nanocrystalline $\text{Mg}(\text{BH})_4$ within Reduced Graphene Oxide. <i>ACS Nano</i> , 2020, 14, 1745-1756.	14.6	29
8	Edge-Functionalized Graphene Nanoribbon Encapsulation To Enhance Stability and Control Kinetics of Hydrogen Storage Materials. <i>Chemistry of Materials</i> , 2019, 31, 2960-2970.	6.7	26
9	Elucidating the mechanism of MgB_2 initial hydrogenation via a combined experimental-theoretical study. <i>Physical Chemistry Chemical Physics</i> , 2017, 19, 22646-22658.	2.8	23
10	Colloidal quantum dot based infrared detectors: extending to the mid-infrared and moving from the lab to the field. <i>Journal of Materials Chemistry C</i> , 2022, 10, 790-804.	5.5	21
11	Identifying the Role of Dynamic Surface Hydroxides in the Dehydrogenation of Ti-Doped NaAlH_4 . <i>ACS Applied Materials & Interfaces</i> , 2019, 11, 4930-4941.	8.0	19
12	Heteroatom-Doped Graphenes as Actively Interacting 2D Encapsulation Media for Mg-Based Hydrogen Storage. <i>ACS Applied Materials & Interfaces</i> , 2022, 14, 20823-20834.	8.0	19
13	Fully Exploited Oxygen Redox Reaction by the Interdiffused Cations in Co^0 -Free Li^+ -Rich Materials for High Performance Li^+ Batteries. <i>Advanced Science</i> , 2020, 7, 2001658.	11.2	17
14	Beyond Idealized Models of Nanoscale Metal Hydrides for Hydrogen Storage. <i>Industrial & Engineering Chemistry Research</i> , 2020, 59, 5786-5796.	3.7	15
15	Spontaneous dynamical disordering of borophenes in MgB_2 and related metal borides. <i>Nature Communications</i> , 2021, 12, 6268.	12.8	14
16	Morphology-Dependent Stability of Complex Metal Hydrides and Their Intermediates Using First-Principles Calculations. <i>ChemPhysChem</i> , 2019, 20, 1340-1347.	2.1	11
17	Hydrogen Storage Performance of Preferentially Oriented Mg/rGO Hybrids. <i>Chemistry of Materials</i> , 2022, 34, 2963-2971.	6.7	8
18	Enhancement of effective thermal conductivity of rGO/Mg nanocomposite packed beds. <i>International Journal of Heat and Mass Transfer</i> , 2022, 192, 122891.	4.8	8

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19	Flexible machine-learning interatomic potential for simulating structural disordering behavior of Li ₇ La ₃ Zr ₂ O ₁₂ solid electrolytes. <i>Journal of Chemical Physics</i> , 2022, 156, .	3.0	8
20	Understanding the effects of oxygen defects on the redox reaction pathways in LiVPO ₄ F by combining <i>ab initio</i> calculations with experiments. <i>Journal of Materials Chemistry A</i> , 2019, 7, 13060-13070.	10.3	7
21	Understanding Charge Transfer at Mg/MgH ₂ Interfaces for Hydrogen Storage. <i>ECS Transactions</i> , 2017, 77, 81-90.	0.5	6
22	Chemomechanical effect of reduced graphene oxide encapsulation on hydrogen storage performance of Pd nanoparticles. <i>Journal of Materials Chemistry A</i> , 2021, 9, 11641-11650.	10.3	6
23	Understanding Hydrogenation Chemistry at MgB ₂ Reactive Edges from <i>Ab Initio</i> Molecular Dynamics. <i>ACS Applied Materials & Interfaces</i> , 2022, 14, 20430-20442.	8.0	4
24	An Analytical Bond Order Potential for Mg-H Systems. <i>ChemPhysChem</i> , 2019, 20, 1404-1411.	2.1	3