

# Qi Song

## List of Publications by Year in descending order

Source: <https://exaly.com/author-pdf/9503904/publications.pdf>

Version: 2024-02-01

10  
papers

62  
citations

1937685

4  
h-index

1588992

8  
g-index

10  
all docs

10  
docs citations

10  
times ranked

77  
citing authors

#	ARTICLE	IF	CITATIONS
1	A theoretical study on the role of ammonium ions in the double-layered $V_2O_5$ electrode. <i>Physical Chemistry Chemical Physics</i> , 2021, 23, 4187-4194.	2.8	3
2	Doping Kinetics and Structural Studies for Lithium-rich Mn-based Lithium Ion Cathode Materials. <i>Energy Technology</i> , 2021, 9, 2000894.	3.8	4
3	Origin of the structural diversity of the alkaline metal borohydride $MBH_4$ ( $M = Li, Na, K, Rb$ and $Cs$ ): Insights from first-principles calculations. <i>International Journal of Hydrogen Energy</i> , 2020, 45, 9946-9958.	7.1	1
4	Highly Selective Electrocatalytic $CO_2$ Reduction to Methanol on Iridium Dioxide with $CO^*$ Spectators. <i>ChemElectroChem</i> , 2020, 7, 5036-5043.	3.4	9
5	A first-principles simulation of the metal borohydride ammonia borane complex $(LiBH_4)_2(NH_3BH_3)$ and the decomposition reaction pathway for hydrogen storage. <i>International Journal of Hydrogen Energy</i> , 2019, 44, 20121-20132.	7.1	3
6	First-principles study of the reorientational motion of the $[BH_4]^-$ groups in $CsBH_4$ 's fractionally occupied ground-state phase and the thermal properties. <i>Physica B: Condensed Matter</i> , 2019, 564, 69-79.	2.7	1
7	Modulating the magnetic properties of $MoS_2$ monolayers by group VIII doping and vacancy engineering. <i>RSC Advances</i> , 2018, 8, 18837-18850.	3.6	15
8	A theoretical explanation of $RbBH_4$ 's fractionally occupied ground-state phase and the reorientational motion of the $[BH_4]^-$ group. <i>Journal Physics D: Applied Physics</i> , 2017, 50, 455501.	2.8	3
9	From orientation disordered to ordered-Anab initio simulation on ammonia borane phase transition within van der Waals corrections. <i>Journal of Computational Chemistry</i> , 2015, 36, 22-32.	3.3	4
10	New schemes for internally contracted multi-reference configuration interaction. <i>Journal of Chemical Physics</i> , 2014, 141, 164114.	3.0	19