

# Xingfeng He

## List of Publications by Year in descending order

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Version: 2024-02-01

19  
papers

6,492  
citations

516561

16  
h-index

794469

19  
g-index

20  
all docs

20  
docs citations

20  
times ranked

5930  
citing authors

#	ARTICLE	IF	CITATIONS
1	Negating interfacial impedance in garnet-based solid-state Li metal batteries. <i>Nature Materials</i> , 2017, 16, 572-579.	13.3	1,583
2	Origin of Outstanding Stability in the Lithium Solid Electrolyte Materials: Insights from Thermodynamic Analyses Based on First-Principles Calculations. <i>ACS Applied Materials &amp; Interfaces</i> , 2015, 7, 23685-23693.	4.0	1,314
3	Electrochemical Stability of $\text{Li}_{10}\text{GeP}_2\text{S}_{12}$ and $\text{Li}_7\text{La}_3\text{Zr}_2\text{O}_{12}$ Solid Electrolytes. <i>Advanced Energy Materials</i> , 2016, 6, 1501590.	10.2	781
4	First principles study on electrochemical and chemical stability of solid electrolyte-electrode interfaces in all-solid-state Li-ion batteries. <i>Journal of Materials Chemistry A</i> , 2016, 4, 3253-3266.	5.2	748
5	Origin of fast ion diffusion in super-ionic conductors. <i>Nature Communications</i> , 2017, 8, 15893.	5.8	570
6	Computation-Accelerated Design of Materials and Interfaces for All-Solid-State Lithium-Ion Batteries. <i>Joule</i> , 2018, 2, 2016-2046.	11.7	266
7	Super-Aligned Carbon Nanotube Films as Current Collectors for Lightweight and Flexible Lithium Ion Batteries. <i>Advanced Functional Materials</i> , 2013, 23, 846-853.	7.8	258
8	Statistical variances of diffusional properties from ab initio molecular dynamics simulations. <i>Npj Computational Materials</i> , 2018, 4, .	3.5	240
9	Strategies Based on Nitride Materials Chemistry to Stabilize Li Metal Anode. <i>Advanced Science</i> , 2017, 4, 1600517.	5.6	185
10	Unsupervised discovery of solid-state lithium ion conductors. <i>Nature Communications</i> , 2019, 10, 5260.	5.8	150
11	Accelerated materials design of $\text{Na}_{0.5}\text{Bi}_{0.5}\text{TiO}_3$ oxygen ionic conductors based on first principles calculations. <i>Physical Chemistry Chemical Physics</i> , 2015, 17, 18035-18044.	1.3	104
12	Crystal Structural Framework of Lithium Super-Ionic Conductors. <i>Advanced Energy Materials</i> , 2019, 9, 1902078.	10.2	93
13	Hybrid super-aligned carbon nanotube/carbon black conductive networks: A strategy to improve both electrical conductivity and capacity for lithium ion batteries. <i>Journal of Power Sources</i> , 2013, 233, 209-215.	4.0	66
14	Enhanced rate capabilities of $\text{Co}_3\text{O}_4$ /carbon nanotube anodes for lithium ion battery applications. <i>Journal of Materials Chemistry A</i> , 2013, 1, 11121.	5.2	50
15	Computation-Guided Design of $\text{LiTaSiO}_5$ , a New Lithium Ionic Conductor with Sphene Structure. <i>Advanced Energy Materials</i> , 2019, 9, 1803821.	10.2	35
16	First-Principles Study of Oxyhydride H Ion Conductors: Toward Facile Anion Conduction in Oxide-Based Materials. <i>ACS Applied Energy Materials</i> , 2018, 1, 1626-1634.	2.5	26
17	First principles hybrid functional study of small polarons in doped $\text{SrCeO}_3$ perovskite: towards computation design of materials with tailored polaron. <i>Ionics</i> , 2018, 24, 1139-1151.	1.2	12
18	$\text{Li}_{15}\text{P}_4\text{S}_{16}\text{Cl}_3$ , a Lithium Chlorothiophosphate as a Solid-State Ionic Conductor. <i>Inorganic Chemistry</i> , 2020, 59, 226-234.	1.9	9

#	ARTICLE	IF	CITATIONS
19	Lithium Superionic Conductors: Crystal Structural Framework of Lithium Superionic Conductors (Adv. Energy Mater. 43/2019). Advanced Energy Materials, 2019, 9, 1970169.	10.2	2