

# Ruijuan Xiao

## List of Publications by Year in descending order

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30  
papers

3,144  
citations

331670

21  
h-index

501196

28  
g-index

30  
all docs

30  
docs citations

30  
times ranked

4096  
citing authors

#	ARTICLE	IF	CITATIONS
1	Localized domains staging structure and evolution in lithiated graphite. , 2023, 5, .		21
2	Mn-Rich Phosphate Cathodes for Na-Ion Batteries with Superior Rate Performance. ACS Energy Letters, 2022, 7, 97-107.	17.4	91
3	New insights into the mechanism of cation migration induced by cation-anion dynamic coupling in superionic conductors. Journal of Materials Chemistry A, 2022, 10, 3093-3101.	10.3	11
4	Accelerated strategy for fast ion conductor materials screening and optimal doping scheme exploration. Journal of Materiomics, 2022, 8, 1038-1047.	5.7	1
5	Screening LiMn <sub>2</sub> O <sub>4</sub> Surface Modification Schemes under Theoretical Guidance. ACS Applied Materials & Interfaces, 2022, 14, 10353-10362.	8.0	14
6	Ionic Conductivity of LiSiON and the Effect of Amorphization/Heterovalent Doping on Li <sup>+</sup> Diffusion. Inorganics, 2022, 10, 45.	2.7	2
7	Interfacial engineering to achieve an energy density of over 200 Wh kg <sup>-1</sup> in sodium batteries. Nature Energy, 2022, 7, 511-519.	39.5	130
8	The Role of Electron Localization in Covalency and Electrochemical Properties of Lithium-Ion Battery Cathode Materials. Advanced Functional Materials, 2021, 31, 2001633.	14.9	21
9	First-Principles Simulations for the Surface Evolution and Mn Dissolution in the Fully Delithiated Spinel LiMn <sub>2</sub> O <sub>4</sub> . Langmuir, 2021, 37, 5252-5259.	3.5	17
10	Reaction Mechanisms of Ta-Substituted Cubic Li <sub>7</sub> La <sub>3</sub> Zr <sub>2</sub> O <sub>12</sub> with Solvents During Storage. ACS Applied Materials & Interfaces, 2021, 13, 38384-38393.	8.0	14
11	Mn Ion Dissolution Mechanism for Lithium-Ion Battery with LiMn <sub>2</sub> O <sub>4</sub> Cathode: <i>In Situ</i> Ultraviolet-Visible Spectroscopy and <i>Ab Initio</i> Molecular Dynamics Simulations. Journal of Physical Chemistry Letters, 2020, 11, 3051-3057.	4.6	60
12	High-throughput computational discovery of K <sub>2</sub> CdO <sub>2</sub> as an ion conductor for solid-state potassium-ion batteries. Journal of Materials Chemistry A, 2020, 8, 5157-5162.	10.3	23
13	Correlated Migration Invokes Higher Na <sup>+</sup> Ion Conductivity in NaSiCON Type Solid Electrolytes. Advanced Energy Materials, 2019, 9, 1902373.	19.5	162
14	Trace doping of multiple elements enables stable battery cycling of LiCoO <sub>2</sub> at 4.6 V. Nature Energy, 2019, 4, 594-603.	39.5	572
15	Design and preparation of nanoporous Ag-Cu alloys by dealloying Mg-(Ag,Cu)-Y metallic glasses for antibacterial applications. Journal of Materials Chemistry B, 2019, 7, 4169-4176.	5.8	30
16	Another Strategy, Detouring Potential Decay by Fast Completion of Cation Mixing. Advanced Energy Materials, 2018, 8, 1703092.	19.5	30
17	Temperature-Sensitive Structure Evolution of Lithium-Manganese-Rich Layered Oxides for Lithium-Ion Batteries. Journal of the American Chemical Society, 2018, 140, 15279-15289.	13.7	163
18	Three-dimensional atomic-scale observation of structural evolution of cathode material in a working all-solid-state battery. Nature Communications, 2018, 9, 3341.	12.8	60

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19	In Situ Atomic-Scale Observation of Electrochemical Delithiation Induced Structure Evolution of $\text{LiCoO}_2$ Cathode in a Working All-Solid-State Battery. <i>Journal of the American Chemical Society</i> , 2017, 139, 4274-4277.	13.7	142
20	Structural stability and stabilization of $\text{Li}_2\text{MoO}_3$ . <i>Physical Chemistry Chemical Physics</i> , 2017, 19, 17538-17543.	2.8	20
21	Oxysulfide $\text{LiAlSO}$ : A Lithium Superionic Conductor from First Principles. <i>Physical Review Letters</i> , 2017, 118, 195901.	7.8	58
22	Oxygen-driven transition from two-dimensional to three-dimensional transport behaviour in $\text{Li}_3\text{PS}_4$ electrolyte. <i>Physical Chemistry Chemical Physics</i> , 2016, 18, 21269-21277.	2.8	66
23	High-throughput design and optimization of fast lithium ion conductors by the combination of bond-valence method and density functional theory. <i>Scientific Reports</i> , 2015, 5, 14227.	3.3	117
24	$\text{P2-Na}_{0.6}[\text{Cr}_{0.6}\text{Ti}_{0.4}]\text{O}_2$ cation-disordered electrode for high-rate symmetric rechargeable sodium-ion batteries. <i>Nature Communications</i> , 2015, 6, 6954.	12.8	426
25	Candidate structures for inorganic lithium solid-state electrolytes identified by high-throughput bond-valence calculations. <i>Journal of Materiomics</i> , 2015, 1, 325-332.	5.7	50
26	Direct Observation of Ordered Oxygen Defects on the Atomic Scale in $\text{Li}_2\text{O}_2$ for $\text{Li-O}_2$ Batteries. <i>Advanced Energy Materials</i> , 2015, 5, 1400664.	19.5	32
27	Screening possible solid electrolytes by calculating the conduction pathways using Bond Valence method. <i>Science China: Physics, Mechanics and Astronomy</i> , 2014, 57, 1526-1536.	5.1	36
28	A zero-strain layered metal oxide as the negative electrode for long-life sodium-ion batteries. <i>Nature Communications</i> , 2013, 4, 2365.	12.8	515
29	Atomic Structure of $\text{Li}_2\text{MnO}_3$ after Partial Delithiation and Re-lithiation. <i>Advanced Energy Materials</i> , 2013, 3, 1358-1367.	19.5	211
30	Compressibility and hardness of Co-based bulk metallic glass: A combined experimental and density functional theory study. <i>Applied Physics Letters</i> , 2011, 99, .	3.3	49