

# Yuwei Li

## List of Publications by Year in descending order

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24  
papers

756  
citations

687363

13  
h-index

580821

25  
g-index

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all docs

25  
docs citations

25  
times ranked

1362  
citing authors

#	ARTICLE	IF	CITATIONS
1	Copper(I)-Based Flexible Organic-Inorganic Coordination Polymer and Analogues: High-Power Factor Thermoelectrics. ACS Applied Materials & Interfaces, 2020, 12, 53841-53851.	8.0	14
2	Prediction of ternary alkaline-earth metal Sn(II) and Pb(II) chalcogenide semiconductors. Physical Review Materials, 2020, 4, .	2.4	1
3	First-principles investigation of structural and electronic properties of oxygen adsorbing phosphorene. Progress in Natural Science: Materials International, 2019, 29, 316-321.	4.4	12
4	Discovery of TaFeSb-based half-Heuslers with high thermoelectric performance. Nature Communications, 2019, 10, 270.	12.8	227
5	First principles based screen for identification of transparent conductors. Journal of Materials Chemistry C, 2019, 7, 2436-2442.	5.5	5
6	Thermoelectric properties of layered NaSbSe <sub>2</sub> . Journal of Physics Condensed Matter, 2018, 30, 225501.	1.8	10
7	Bismuth and antimony-based oxyhalides and chalcogenides as potential optoelectronic materials. Npj Computational Materials, 2018, 4, .	8.7	86
8	Frustrated Structural Instability in Superconducting Quasi-One-Dimensional $KxMn_2O_3$ . Physical Review Letters, 2018, 121, 187002.	7.8	16
9	Thermoelectric properties of p-type cubic and rhombohedral GeTe. Journal of Applied Physics, 2018, 123, .	2.5	40
10	Optimal Bandgap in a 2D Ruddlesden-Popper Perovskite Chalcogenide for Single-Junction Solar Cells. Chemistry of Materials, 2018, 30, 4882-4886.	6.7	49
11	Tunability of electronic and optical properties of the BaZrS system via dimensional reduction. European Physical Journal B, 2018, 91, 1.	1.5	16
12	Rational design of new phases of tin monosulfide by first-principles structure searches. Science China: Physics, Mechanics and Astronomy, 2018, 61, 1.	5.1	15
13	Optical and electronic properties of doped p-type CuI: Explanation of transparent conductivity from first principles. Physical Review Materials, 2018, 2, .	2.4	7
14	Sn <sub>2</sub> Se <sub>3</sub> : A conducting crystalline mixed valent phase change memory compound. Journal of Applied Physics, 2017, 121, .	2.5	9
15	Infrared absorption and visible transparency in heavily doped p-type BaSnO <sub>3</sub> . Applied Physics Letters, 2017, 110, 051904.	3.3	11
16	Sn(II)-Containing Phosphates as Optoelectronic Materials. Chemistry of Materials, 2017, 29, 2459-2465.	6.7	17
17	Prediction of nontrivial band topology and superconductivity in $MxMn_2O_3$ . Physical Review Materials, 2017, 1, .	2.4	8
18	New stable ternary alkaline-earth metal Pb(II) oxides: $CaPbO_3$ and $Pb_3O_3$ . Physical Review Materials, 2017, 1, .	2.4	10

#	ARTICLE	IF	CITATIONS
19	Electronic fitness function for screening semiconductors as thermoelectric materials. Physical Review Materials, 2017, 1, .	2.4	98
20	Properties of the ferroelectric visible light absorbing semiconductors: $P\text{Sn}_2$ and $S\text{Sn}_6$ and $\text{Sn}_2$ $\text{Sn}_6$ Determination of the high pressure phases of $\text{CaWO}_4$ by CALYPSO and X-ray diffraction studies. Physica Status Solidi (B): Basic Research, 2016, 253, 1947-1951.	2.4	23
21	Determination of the high pressure phases of $\text{CaWO}_4$ by CALYPSO and X-ray diffraction studies. Physica Status Solidi (B): Basic Research, 2016, 253, 1947-1951.	1.5	8
22	Identification and properties of the non-cubic phases of $\text{Mg}_2\text{Pb}$ . AIP Advances, 2016, 6, 125108.	1.3	3
23	Design of ternary alkaline-earth metal $\text{Sn}(\text{II})$ oxides with potential good p-type conductivity. Journal of Materials Chemistry C, 2016, 4, 4592-4599.	5.5	29
24	Tuning optical properties of transparent conducting barium stannate by dimensional reduction. APL Materials, 2015, 3, .	5.1	29