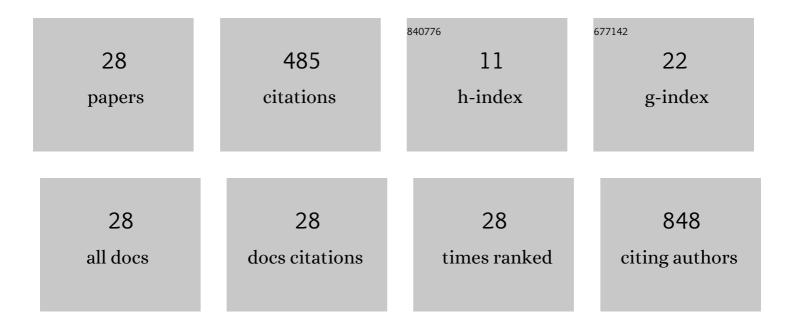
Dawei Zhou

List of Publications by Year in descending order

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ΠΛΜΕΙ ΖΗΟΙΙ

#	Article	IF	CITATIONS
1	Formation Mechanism of Impure Phases and Crystallinity Investigation of YAG Powders Synthesized via the Co-precipitation Method. Journal Wuhan University of Technology, Materials Science Edition, 2021, 36, 517-525.	1.0	1
2	Pressure-induced the formation of Mg(CH3)2 and Ca(CH3)2 studied by the first principles. Solid State Communications, 2020, 320, 114027.	1.9	0
3	Exploring the real ground-state structures of W3Si silicides from first-principles calculations. Computational Materials Science, 2020, 180, 109719.	3.0	16
4	Hydrogenated PtP ₂ monolayer: theoretical predictions on the structure and charge carrier mobility. Journal of Materials Chemistry C, 2019, 7, 12231-12239.	5.5	21
5	Crystal structure prediction of ReN at high pressure: a new incompressible phase. Phase Transitions, 2019, 92, 595-602.	1.3	1
6	Unexpected Xe Cations and Superconductivity in Y–Xe Intermediate Compounds under Pressure. Journal of Physical Chemistry C, 2019, 123, 9323-9330.	3.1	6
7	Prediction of stable Cu-Li binary intermetallics from first-principles calculations: Stoichiometries, crystal structures, and physical properties. Journal of Alloys and Compounds, 2018, 766, 640-648.	5.5	12
8	Two-Dimensional C ₄ N Global Minima: Unique Structural Topologies and Nanoelectronic Properties. Journal of Physical Chemistry C, 2017, 121, 2669-2674.	3.1	49
9	High-pressure synthesis and thermoelectric performance of tellurium doped with bismuth. Journal of Materials Science, 2017, 52, 10526-10532.	3.7	28
10	Pressure-induced metallization and superconducting phase in ReS 2. Npj Quantum Materials, 2017, 2, .	5.2	53
11	Origin of superconductivity in the Weyl semimetal <mml:math xmlns:mml="http://www.w3.org/1998/Math/MathML"><mml:mrow><mml:mi>WT</mml:mi><mml:msub><mml:m mathvariant="normal">e<mml:mn>2</mml:mn></mml:m </mml:msub></mml:mrow> under pressure. Physical Review B, 2016, 94, .</mml:math 	^{າi} 3.2	91
12	Crossover from metal to insulator in dense lithium-rich compound CLi ₄ . Proceedings of the United States of America, 2016, 113, 2366-2369.	7.1	21
13	Pressure-induced phase transition of BiOF: novel two-dimensional layered structures. Physical Chemistry Chemical Physics, 2015, 17, 4434-4440.	2.8	15
14	Phase transition and thermodynamic properties of YAg alloy from first-principles calculations. Computational Materials Science, 2015, 102, 21-26.	3.0	5
15	Transport and Magnetic Properties of K0.8Fe2â^'x Cu x Se2(0 ⩽ x ⩽ 2) System. Journal of Superconductiv and Novel Magnetism, 2015, 28, 219-222.	ity 1.8	3
16	Ab-initio study of phase stability, elastic and thermodynamic properties of AlY alloy under pressure. Journal of Alloys and Compounds, 2015, 648, 67-74.	5.5	7
17	Elastic and thermodynamic properties of vanadium nitride under pressure and the effect of metallic bonding on its hardness. Chinese Physics B, 2014, 23, 026201.	1.4	9
18	First-principles investigation of elastic and thermodynamic properties of SiCN under pressure. Computational Materials Science, 2014, 95, 228-234.	3.0	5

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#	Article	IF	CITATIONS
19	Superconducting properties of barium in three phases under high pressure from first principles. Chinese Physics B, 2013, 22, 087403.	1.4	3
20	Theoretical investigation on structural and thermodynamic properties of the intermetallic compound in Mg–Zn–Ag alloy under high pressure and high temperature. Journal of Alloys and Compounds, 2013, 550, 406-411.	5.5	18
21	The Predicted fcc Superconducting Phase for Compressed Se and Te. Chinese Physics Letters, 2013, 30, 027401.	3.3	5
22	<i>Ab initio</i> study revealing a layered structure in hydrogen-rich KH <mml:math xmlns:mml="http://www.w3.org/1998/Math/MathML" display="inline"><mml:msub><mml:mrow /><mml:mn>6</mml:mn></mml:mrow </mml:msub>under high pressure. Physical Review B, 2012, 86, .</mml:math 	3.2	79
23	Melting curve of the cl16 sodium at high pressure from <i>ab initio</i> calculations. Physica Status Solidi (B): Basic Research, 2011, 248, 1143-1148.	1.5	2
24	A New High-Pressure Polar Phase of Crystalline Bromoform: A First-Principles Study. Journal of Physical Chemistry B, 2010, 114, 13933-13939.	2.6	7
25	Peierls transition in sodium under high pressure: a first-principles study. Journal of Physics Condensed Matter, 2009, 21, 025508.	1.8	9
26	Near-edge X-ray absorption fine structure of solid oxygen under high pressure: A density functional theory study. Solid State Communications, 2008, 147, 126-129.	1.9	5
27	Annealing temperature dependance of magnetic properties and magneto-impedance effect in CoZrB alloys. Journal of Alloys and Compounds, 2008, 464, 28-32.	5.5	5
28	Influence of Fe on magnetoimpedance effect of Co72â^'xFexZr8B20alloys. Journal Physics D: Applied Physics, 2006, 39, 4299-4303.	2.8	9