

# Xi Zhang

## List of Publications by Year in descending order

Source: <https://exaly.com/author-pdf/7333738/publications.pdf>

Version: 2024-02-01

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papers

352

citations

1040056

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1372567

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docs citations

10

times ranked

374

citing authors

#	ARTICLE	IF	CITATIONS
1	Ab initio prediction of vacancy energetics in HCP Al-Hf-Sc-Ti-Zr high entropy alloys and the subsystems. <i>Acta Materialia</i> , 2022, 227, 117677.	7.9	22
2	Zr diffusion in BCC refractory high entropy alloys: A case of “non-sluggish” diffusion behavior. <i>Acta Materialia</i> , 2022, 233, 117970.	7.9	33
3	‘Ab initio’ simulations of the surface free energy of TiN(001). <i>Physical Review B</i> , 2021, 103, .	3.2	9
4	Kinetically driven ordering in phase separating alloys. <i>Physical Review Materials</i> , 2019, 3, .	2.4	4
5	Calculating free energies of point defects from ab initio. <i>Computational Materials Science</i> , 2018, 148, 249-259.	3.0	47
6	Temperature dependence of the stacking-fault Gibbs energy for Al, Cu, and Ni. <i>Physical Review B</i> , 2018, 98, . Accurate electronic free energies of the $\text{Al}_{\text{Mn}}\text{Cu}_{\text{Mn}}$ transition metals at high temperatures.	3.2	61
7	$\text{d} \left( \frac{\partial \ln f}{\partial \mu_i} \right) = \frac{\partial \ln f}{\partial \mu_i} + \frac{\partial \ln f}{\partial T} \frac{\partial T}{\partial \mu_i}$ , and $\frac{\partial \ln f}{\partial \mu_i} = \frac{\partial \ln f}{\partial \mu_i} + \frac{\partial \ln f}{\partial T} \frac{\partial T}{\partial \mu_i}$	3.2	70
8	‘Ab initio’ prediction of vacancy properties in concentrated alloys: The case of fcc Cu-Ni. <i>Physical Review B</i> , 2015, 91, .	3.2	29
9	Thermodynamic assessment of Co-Al-W system and solidification of Co-enriched ternary alloys. <i>Journal of Materials Science</i> , 2011, 46, 2611-2621.	3.7	38
10	Thermodynamic assessment of the U-Mo-Al system. <i>Journal of Nuclear Materials</i> , 2010, 402, 15-24.	2.7	39