

Chahra Sifi

List of Publications by Year in descending order

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7
papers

66
citations

1937685

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1872680

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#	ARTICLE	IF	CITATIONS
1	First principle calculations of structural, electronic, thermodynamic and optical properties of $\text{Pb}_{1-x}\text{Ca}_x\text{S}$, $\text{Pb}_{1-x}\text{Ca}_x\text{Se}$ and $\text{Pb}_{1-x}\text{Ca}_x\text{Te}$ ternary alloys. Journal of Physics Condensed Matter, 2009, 21, 195401.	1.8	33
2	Ab initio investigations of calcium chalcogenide alloys. Journal of Alloys and Compounds, 2009, 485, 642-647.	5.5	17
3	Computational investigations of the band structure and thermodynamic properties of calcium-doped BaS using the FP-LAPW approach. Chinese Journal of Physics, 2017, 55, 367-377.	3.9	7
4	First-principles calculations on the origins of the gap bowing in $\text{Pb}_{1-x}\text{Ca}_x\text{S}$, $\text{Pb}_{1-x}\text{Ca}_x\text{Se}$ and $\text{Pb}_{1-x}\text{Ca}_x\text{Te}$ alloys. IOP Conference Series: Materials Science and Engineering, 2012, 28, 012031.	0.6	4
5	Results of optical and thermal study of lead sulfide (PbS), Barium sulfide (BaS) and their ternary ($\text{Pb}_{1-x}\text{Ba}_x\text{S}$) ($0 \leq x \leq 1$) alloys. Computational Condensed Matter, 2019, 21, e00398.	2.1	3
6	Theoretical analysis of the electronic, optical and thermal properties of lead strontium telluride alloys $\text{Pb}_{1-x}\text{Sr}_x\text{Te}$ ($x = 0.0 \sim 1.0$). Philosophical Magazine, 2018, 98, 295-311.	1.6	2
7	The calculations of thermodynamic and opto-electronics properties of $\text{Pb}_{1-x}\text{Ca}_x\text{Se}$ semiconducting ternary Alloys. EPJ Web of Conferences, 2012, 29, 00033.	0.3	0