

# S Labidi

## List of Publications by Year in descending order

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Version: 2024-02-01

19

papers

251

citations

933447

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docs citations

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times ranked

191

citing authors

#	ARTICLE	IF	CITATIONS
1	First principles calculations of structural, electronic and thermodynamic properties of SrS, SrSe, SrTe compounds and $\text{SrS}_{1-x}\text{Sex}$ alloy. Physics Procedia, 2009, 2, 1205-1212.	1.2	37
2	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
3	(FP-LPAW) Investigation of Ternary Alloys: $\text{CdS}_{1-x}\text{TeX}$ . Physics Procedia, 2009, 2, 881-887.	1.2	25
4	Structural, electronic, thermodynamic and optical properties of alkaline earth oxides $\text{MgO}$ , $\text{SrO}$ and their alloys. Physica Scripta, 2010, 82, 045605.	2.5	24
5	FP-LAPW investigations of $\text{SrS}_{1-x}\text{Se}$ , $\text{SrS}_{1-x}\text{Te}$ and $\text{SrSe}_{1-x}\text{Te}$ ternary alloys. Journal of Physics Condensed Matter, 2008, 20, 445213.	1.8	23
6	First principles calculations of structural, electronic and optical properties of $\text{BAs}_{1-x}\text{Px}$ alloy. Physics Procedia, 2009, 2, 933-940.	1.2	17
7	Ab initio investigations of calcium chalcogenide alloys. Journal of Alloys and Compounds, 2009, 485, 642-647.	5.5	17
8	First principles calculations of structural, electronic, optical and thermodynamic properties of $\text{PbS}$ , $\text{SrS}$ and their ternary alloys $\text{Pb}_{1-x}\text{Sr}_x\text{S}$ . Computational Materials Science, 2011, 50, 1077-1082.	3.0	14
9	First principles calculations of structural, electronic, thermodynamic and optical properties of $\text{BAs}_{1-x}\text{P}_{x}$ alloy. Physica Scripta, 2009, 79, 045002.	2.5	12
10	Structural, electronic, thermodynamic and optical properties of $\text{SrS}_{1-x}\text{Ox}$ mixed crystals. Physica B: Condensed Matter, 2009, 404, 4100-4105.	2.7	10
11	STRUCTURAL, ELECTRONIC, OPTICAL AND THERMODYNAMIC PROPERTIES OF $\text{PbS}$ , $\text{PbSe}$ AND THEIR TERNARY ALLOY $\text{PbS}_{1-x}\text{Se}_x$ . Modern Physics Letters B, 2011, 25, 473-486.	1.9	9
12	Density functional calculations of $\text{Pb}_{1-x}\text{Ca}_x\text{SySe}_{1-y}$ alloys lattice matched to different substrates. Journal of Physics and Chemistry of Solids, 2012, 73, 608-613.	4.0	9
13	Investigation of total and partial magnetic moments of $\text{Mn}_2\text{NiAl}$ with pressure at a several temperatures. Phase Transitions, 2019, 92, 699-706.	1.3	6
14	Structural, electronic and thermodynamic properties of $\text{SrxCd}_{1-x}\text{O}$ : A first-principles study. Materials Science in Semiconductor Processing, 2013, 16, 1853-1858.	4.0	5
15	Ab initio calculations of fundamental properties of $\text{SrTe}_{1-x}\text{O}_x$ alloys. Bulletin of Materials Science, 2016, 39, 827-835.	1.7	5
16	Temperature and pressure dependence on structural, electronic and thermal properties of $\text{ZnO}$ wurtzite phase - first principle investigation. Phase Transitions, 2020, 93, 654-665.	1.3	3
17	Theoretical Investigations of the Physical Properties of $\text{Cd}_{1-x}\text{Ca}_x\text{O}$ Ternary Alloys. Chinese Physics Letters, 2014, 31, 046104.	3.3	2
18	Structural, electronic and thermodynamic properties of $\text{Sr}_{x}\text{Ca}_{1-x}\text{S}$ : A first-principles study. Physica Status Solidi C: Current Topics in Solid State Physics, 2016, 13, 629-634.	0.8	0

#	ARTICLE	IF	CITATIONS
19	Structural, electronic, magnetic and thermodynamic properties of Ni <sub>1-x</sub> Ti <sub>x</sub> O alloys an ab initio calculation and Monte Carlo study. Phase Transitions, 2018, 91, 600-609.	1.3	0