

Houda Ben Abdallah

List of Publications by Year in descending order

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Version: 2024-02-01

11
papers

197
citations

1307594

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h-index

1281871

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12
all docs

12
docs citations

12
times ranked

276
citing authors

| # | ARTICLE | IF | CITATIONS |
|----|--|-----|-----------|
| 1 | Electronic structure and optical properties of Sb ₂ S ₃ crystal. <i>Physica B: Condensed Matter</i> , 2011, 406, 287-292. | 2.7 | 87 |
| 2 | First principles calculations of electronic and optical properties of Ag ₂ S. <i>Solid State Sciences</i> , 2013, 26, 65-71. | 3.2 | 26 |
| 3 | First-principles study of the electronic and the optical properties of In ₆ Se ₇ compound. <i>Physica B: Condensed Matter</i> , 2010, 405, 3427-3432. | 2.7 | 20 |
| 4 | Electronic structure of the hexaindium heptasulfide In ₆ S ₇ . <i>Physica B: Condensed Matter</i> , 2006, 382, 181-188. | 2.7 | 12 |
| 5 | Spin-orbit coupling effect on electronic, linear and nonlinear optical properties of Bi ₂ S ₃ and the ternary bismuth sulfide Bi ₂ S _{2.75} Se _{0.25} : Ab-initio calculations. <i>Optical and Quantum Electronics</i> , 2022, 54, 1. | 3.3 | 10 |
| 6 | First-principles calculations on magnetism and exchange interactions in GaMnAs and GaMnAsP. <i>Physica Status Solidi (B): Basic Research</i> , 2017, 254, 1700115. | 1.5 | 9 |
| 7 | First-principles calculations of the electronic and optical properties of In ₆ S ₇ compound. <i>Physica B: Condensed Matter</i> , 2009, 404, 194-198. | 2.7 | 8 |
| 8 | Hybrid functional calculations of electro-optical properties of novel Ga _{1-x} In _x Te ternary chalcogenides. <i>Applied Physics A: Materials Science and Processing</i> , 2020, 126, 1. | 2.3 | 8 |
| 9 | DFT calculations on ZnO _{1-x} compounds for optoelectronic applications. <i>Journal of Computational Electronics</i> , 2021, 20, 467-479. | 2.5 | 8 |
| 10 | Ab Initio Study of Structural, Electronic, and Magnetic Properties of A _{1-x} B _x III. <i>Journal of Superconductivity and Novel Magnetism</i> , 2018, 31, 2089-2097. | 1.8 | 6 |
| 11 | Selenium alloying of indium sulfide: Ab-initio study of structural, electronic and optical features. <i>Materials Science in Semiconductor Processing</i> , 2015, 31, 56-67. | 4.0 | 3 |