

Claudia RÃ¶dl

List of Publications by Year in descending order

Source: <https://exaly.com/author-pdf/5387096/publications.pdf>

Version: 2024-02-01

32
papers

2,343
citations

236925

25
h-index

414414

32
g-index

33
all docs

33
docs citations

33
times ranked

3037
citing authors

| # | ARTICLE | IF | CITATIONS |
|----|--|------|-----------|
| 1 | Efficient strain-induced light emission in lonsdaleite germanium. <i>Physical Review Materials</i> , 2021, 5, . | 2.4 | 16 |
| 2 | Layered CuI: a path to 2D p-type transparent conducting materials. <i>Journal of Materials Chemistry C</i> , 2021, 9, 11284-11291. | 5.5 | 7 |
| 3 | From pseudo-direct hexagonal germanium to direct silicon-germanium alloys. <i>Physical Review Materials</i> , 2021, 5, . | 2.4 | 7 |
| 4 | Direct-bandgap emission from hexagonal Ge and SiGe alloys. <i>Nature</i> , 2020, 580, 205-209. | 27.8 | 231 |
| 5 | Stable Ordered Phases of Cuprous Iodide with Complexes of Copper Vacancies. <i>Chemistry of Materials</i> , 2019, 31, 7877-7882. | 6.7 | 17 |
| 6 | Accurate electronic and optical properties of hexagonal germanium for optoelectronic applications. <i>Physical Review Materials</i> , 2019, 3, . | 2.4 | 41 |
| 7 | The ground state of two-dimensional silicon. <i>2D Materials</i> , 2018, 5, 035010. | 4.4 | 25 |
| 8 | Low-energy electronic excitations and band-gap renormalization in CuO. <i>Physical Review B</i> , 2017, 95, . | 3.2 | 7 |
| 9 | Wurtzite silicon as a potential absorber in photovoltaics: Tailoring the optical absorption by applying strain. <i>Physical Review B</i> , 2015, 92, . | 3.2 | 54 |
| 10 | Quasiparticle excitations in the photoemission spectrum of CuO from first principles: $A \rightarrow G \rightarrow W \rightarrow X$. <i>Physical Review B</i> , 2015, 91, . | 3.2 | 48 |
| 11 | Photoemission spectra and effective masses of n- and p-type oxide semiconductors from first principles: ZnO, CdO, SnO ₂ , MnO, and NiO. <i>Physica Status Solidi (A) Applications and Materials Science</i> , 2014, 211, 74-81. | 1.8 | 40 |
| 12 | Band discontinuities at Si-TCO interfaces from quasiparticle calculations: Comparison of two alignment approaches. <i>Physical Review B</i> , 2012, 85, . | 3.2 | 62 |
| 13 | Crystalline and magnetic anisotropy of the 3d-transition metal monoxides MnO, FeO, CoO, and NiO. <i>Physical Review B</i> , 2012, 86, . | 3.2 | 97 |
| 14 | Optical and energy-loss spectra of the antiferromagnetic transition metal oxides MnO, FeO, CoO, and NiO including quasiparticle and excitonic effects. <i>Physical Review B</i> , 2012, 86, . | 3.2 | 59 |
| 15 | First-Principles Optical Spectra for F Centers in MgO. <i>Physical Review Letters</i> , 2012, 108, 126404. | 7.8 | 157 |
| 16 | Tin dioxide from first principles: Quasiparticle electronic states and optical properties. <i>Physical Review B</i> , 2011, 83, . | 3.2 | 145 |
| 17 | Electronic and optical properties of Mg _x Zn _{1-x} O and Cd _x Zn _{1-x} O from <i>ab initio</i> calculations. <i>New Journal of Physics</i> , 2011, 13, 085012. | 2.9 | 60 |
| 18 | Optical Absorption in Degenerately Doped Semiconductors: Mott Transition or Mahan Excitons?. <i>Physical Review Letters</i> , 2011, 107, 236405. | 7.8 | 61 |

| # | ARTICLE | IF | CITATIONS |
|----|--|-----|-----------|
| 19 | Interplay of excitonic effects and van Hove singularities in optical spectra: CaO and AlN polymorphs. Physical Review B, 2011, 84, . | 3.2 | 46 |
| 20 | <i>Ab initio</i> description of heterostructural alloys: Thermodynamic and structural properties of $\text{Mg}_x\text{Zn}_{1-x}\text{O}$. Physical Review B, 2010, 81, . | 3.2 | 49 |
| 21 | Influence of Strong Electron Correlation on Magnetism in Transition-Metal Doped Si Nanocrystals. Journal of Chemical Theory and Computation, 2010, 6, 353-358. | 5.3 | 11 |
| 22 | Band lineup between silicon and transparent conducting oxides. Applied Physics Letters, 2010, 97, . | 3.3 | 48 |
| 23 | Energetic stability and magnetic properties of MnO in the rocksalt, wurtzite, and zinc-blende structures: Influence of exchange and correlation. Physical Review B, 2010, 82, . | 3.2 | 62 |
| 24 | Optical and energy-loss spectra of MgO, ZnO, and CdO from <i>ab initio</i> many-body calculations. Physical Review B, 2009, 80, . | 3.2 | 142 |
| 25 | Band structure and optical transition parameters of wurtzite MgO, ZnO, and CdO from quasiparticle calculations. Physica Status Solidi (B): Basic Research, 2009, 246, 2150-2153. | 1.5 | 68 |
| 26 | Quasiparticle band structures of the antiferromagnetic transition-metal oxides MnO, FeO, CoO, and NiO. Physical Review B, 2009, 79, . | 3.2 | 243 |
| 27 | Branch-point energies and band discontinuities of III-nitrides and III-III-oxides from quasiparticle band-structure calculations. Applied Physics Letters, 2009, 94, . | 3.3 | 177 |
| 28 | LiNbO_3 ground- and excited-state properties from first-principles calculations. Physical Review B, 2008, 77, . | 3.2 | 86 |
| 29 | <i>Ab initio</i> theory of excitons and optical properties for spin-polarized systems: Application to antiferromagnetic MnO. Physical Review B, 2008, 77, . | 3.2 | 79 |
| 30 | Efficient O^2 to solve the Bethe-Salpeter equation for excitonic bound states. Physical Review B, 2008, 78, . | 3.2 | 117 |
| 31 | Ab-Initio Studies of Electronic and Spectroscopic Properties of MgO, ZnO and CdO. Journal of the Korean Physical Society, 2008, 53, 2811-2815. | 0.7 | 26 |
| 32 | Strain influence on valence-band ordering and excitons in ZnO: An <i>ab initio</i> study. Applied Physics Letters, 2007, 91, 241915. | 3.3 | 55 |