

Niels E Christensen

List of Publications by Year in descending order

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218381

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times ranked

2719
citing authors

#	ARTICLE	IF	CITATIONS
1	Rocksalt ZnMgO alloys for ultraviolet applications: Origin of band-gap fluctuations and direct-indirect transitions. <i>Physical Review B</i> , 2020, 101, .	1.1	16
2	ZnO/(Zn)MgO polar and nonpolar superlattices. <i>Journal of Applied Physics</i> , 2019, 125, .	1.1	14
3	Theoretical study of nitride short period superlattices. <i>Journal of Physics Condensed Matter</i> , 2018, 30, 063001.	0.7	31
4	Comparison of wurtzite GaN/AlN and ZnO/MgO short-period superlattices: Calculation of band gaps and built-in electric field. <i>Physica Status Solidi (B): Basic Research</i> , 2017, 254, 1600704.	0.7	15
5	Bandgap behavior of InGaN/GaN short period superlattices grown by metal-organic vapor phase epitaxy. <i>Physica Status Solidi (B): Basic Research</i> , 2017, 254, 1600710.	0.7	10
6	Band gap engineering of In(Ga)N/GaN short period superlattices. <i>Scientific Reports</i> , 2017, 7, 16055.	1.6	22
7	Structural and electronic properties of wurtzite MgZnO and BeMgZnO alloys and their thermodynamic stability. <i>Journal of Applied Physics</i> , 2016, 120, .	1.1	17
8	ZnGeSb ₂ : a promising thermoelectric material with tunable ultra-high conductivity. <i>Physical Chemistry Chemical Physics</i> , 2016, 18, 26275-26283.	1.3	11
9	Band gaps and built-in electric fields in InAlN/GaN short period superlattices: Comparison with (InAlGa)N quaternary alloys. <i>Physical Review B</i> , 2016, 93, .	1.1	10
10	Thermoelectric properties of binary LnN (Ln=La and Lu): First principles study. <i>AIP Conference Proceedings</i> , 2015, , .	0.3	4
11	Influence of internal electric fields on band gaps in short period GaN/GaAlN and InGaN/GaN polar superlattices. <i>Journal of Applied Physics</i> , 2015, 118, .	1.1	18
12	Structural, electronic and hyperfine characterization of pure and Ta-doped ZrSiO ₄ . <i>Physical Review B</i> , 2015, 91, .	1.1	5
13	Band gaps and internal electric fields in semipolar short period InN/GaN superlattices. <i>Applied Physics Letters</i> , 2014, 104, .	1.5	5
14	The discrepancies between theory and experiment in the optical emission of monolayer In(Ga)N quantum wells revisited by transmission electron microscopy. <i>Applied Physics Letters</i> , 2014, 104, .	1.5	48
15	Short period polar and nonpolar In _m /In _n /GaN superlattices. <i>Physica Status Solidi C: Current Topics in Solid State Physics</i> , 2014, 11, 678-681.	0.8	2
16	Theoretical investigation of pressure-induced structural transitions in americium using $GGA+U$ and hybrid density functional theory methods. <i>Physical Review B</i> , 2013, 88, .	1.1	18
17	Electronic correlation strength of Pu. <i>Physical Review B</i> , 2013, 87, .	1.1	7
18	Photoluminescence and pressure effects in short period InN/nGaIn superlattices. <i>Journal of Applied Physics</i> , 2013, 113, 123101.	1.1	26

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19	Thermoelectric properties of chalcopyrite type CuGaTe ₂ and chalcostibite CuSbS ₂ . Journal of Applied Physics, 2013, 114, .	1.1	67
20	Band gaps in InN/GaN superlattices: Nonpolar and polar growth directions. Journal of Applied Physics, 2013, 114, 223102.	1.1	13
21	Electronic topological transition in LaSn ₃ under pressure. Physical Review B, 2012, 85, .	1.1	50
22	Structural, electronic, and hyperfine properties of pure and Ta-doped m-ZrO ₂ . Physical Review B, 2012, 85, .	1.1	16
23	Hydrostatic pressure and strain effects in short period InN/GaN superlattices. Applied Physics Letters, 2012, 101, 092104.	1.5	22
24	Band Structure and Quantum Confined Stark Effect in InN/GaN superlattices. Crystal Growth and Design, 2012, 12, 3521-3525.	1.4	37
25	Band gap bowing in quaternary nitride semiconducting alloys. Applied Physics Letters, 2011, 98, .	1.5	29
26	Band gap bowings and anomalous pressure effects in III-V nitride alloys: Role of In ϵ segregation. Physica Status Solidi (A) Applications and Materials Science, 2011, 208, 1558-1561.	0.8	6
27	Calculations of quasi ϵ particle spectra of semiconductors under pressure. Physica Status Solidi (B): Basic Research, 2011, 248, 1096-1101.	0.7	2
28	Quasiparticle band structures of \hat{I}^2 -HgS, HgSe, and HgTe. Physical Review B, 2011, 84, .	1.1	82
29	Size effects in band gap bowing in nitride semiconducting alloys. Physical Review B, 2011, 83, .	1.1	28
30	Atomic structure and electronic properties of the Si _x Sb _{1-x} alloys. Physica Status Solidi (A) Applications and Materials Science, 2010, 207, 1369-1371.	1.1	7
31	In ϵ clustering effects in InAlN and InGaN revealed by high pressure studies. Physica Status Solidi (A) Applications and Materials Science, 2010, 207, 1369-1371.	0.8	16
32	In-clustering induced anomalous behavior of band gap in InAlN and InGaN. Physica Status Solidi C: Current Topics in Solid State Physics, 2010, 7, 1283-1286.	0.8	7
33	Limitations to band gap tuning in nitride semiconductor alloys. Applied Physics Letters, 2010, 96, .	1.5	38
34	Electronic properties of CuAlO_3 under pressure: Three theoretical approaches. Physical Review B, 2010, 81, .	1.1	54
35	Full-zone analysis of relativistic spin splitting at band anticrossings: The case of zinc-blende semiconductors. Physical Review B, 2010, 81, .	1.1	17
36	Quasiparticle self-consistent GW theory of III-V nitride semiconductors: Bands, gap bowing, and effective masses. Physical Review B, 2010, 82, .	1.1	53

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37	High-pressure structural study of fluoro-perovskite CsCdF_3 to 60 GPa: A combined experimental and theoretical study. <i>Physical Review B</i> , 2010, 81, .	1.1	46
38	Coexistence of different charge states in Ta-doped monoclinic HfO_2 . Theoretical and experimental approaches. <i>Physical Review B</i> , 2010, 82, .	1.1	19
39	Anomalous composition dependence of the band gap pressure coefficients in In-containing nitride semiconductors. <i>Physical Review B</i> , 2010, 81, .	1.1	27
40	Quasiparticle self-consistent GW calculations for PbS, PbSe, and PbTe: Band structure and pressure coefficients. <i>Physical Review B</i> , 2010, 81, .	1.1	134
41	Pressure variation of the valence band width in Ge: A self-consistent GW study. <i>Physical Review B</i> , 2009, 79, .	1.1	9
42	The magnetization of Fe_4N : theory vs. experiment. <i>Physica Status Solidi (B): Basic Research</i> , 2009, 246, 909-928.	0.7	79
43	Electronic and optical properties of III nitrides under pressure. <i>Physica Status Solidi (B): Basic Research</i> , 2009, 246, 570-575.	0.7	12
44	Gap bowing in $\text{In}_x\text{Ga}_{1-x}\text{N}$ and $\text{In}_x\text{Al}_{1-x}\text{N}$ under pressure. <i>Physica Status Solidi C: Current Topics in Solid State Physics</i> , 2009, 6, S368.	0.8	7
45	Influence of indium clustering on the band structure of semiconducting ternary and quaternary nitride alloys. <i>Physical Review B</i> , 2009, 80, .	1.1	132
46	Conduction band filling in In-rich InGaN and InN under hydrostatic pressure. <i>Physica Status Solidi C: Current Topics in Solid State Physics</i> , 2008, 5, 1488-1490.	0.8	1
47	Band structure and effective mass of InN under pressure. <i>Physica Status Solidi (B): Basic Research</i> , 2008, 245, 887-889.	0.7	5
48	Bowing of the band gap pressure coefficient in $\text{In}_x\text{Ga}_{1-x}\text{N}$ alloys. <i>Journal of Applied Physics</i> , 2008, 103, 033514.	1.1	53
49	Electronic structure and effective masses of InN under pressure. <i>Journal of Applied Physics</i> , 2008, 104, 013704.	1.1	27
50	Strain-induced conduction-band spin splitting in GaAs from first-principles calculations. <i>Physical Review B</i> , 2008, 78, .	1.1	16
51	Ab initio study of structural and electronic properties and hyperfine interactions at the Ta nucleus in Ta-doped monoclinic HfO_2 . <i>Physical Review B</i> , 2008, 78, .	1.1	24
52	Role of conduction-band filling in the dependence of InN photoluminescence on hydrostatic pressure. <i>Physical Review B</i> , 2007, 76, .	1.1	27
53	Pressure and composition dependence of the electron effective mass in $\text{GaAs}_{1-x}\text{N}_x$. <i>Physica Status Solidi (B): Basic Research</i> , 2006, 243, 1599-1603.	0.7	2
54	Calculated superconductive properties of Li and Na under pressure. <i>Physical Review B</i> , 2006, 73, .	1.1	39

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55	Linear and nonlinear optical properties of Li under pressure. <i>Physical Review B</i> , 2006, 73, .	1.1	9
56	Electronic structure of rare-earth impurities in GaAs and GaN. <i>Physical Review B</i> , 2006, 74, .	1.1	65
57	First Principles Determination of Hyperfine Parameters on fcc-Fe ₈ X (X = C, N) Arrangements. <i>Hyperfine Interactions</i> , 2005, 161, 197-202.	0.2	4
58	Ab-initio study of excitonic effects in conventional and organic semiconductors. <i>Physica Status Solidi (B): Basic Research</i> , 2005, 242, 1754-1758.	0.7	14
59	Pressure strengthening and its application to the analysis of hydrogen sample-radius behaviour in a tungsten gasket as a function of the initial sample-radius-to-tip-radius ratio. <i>Journal of Physics Condensed Matter</i> , 2002, 14, 10907-10910.	0.7	0
60	Predicted Superconductive Properties of Lithium under Pressure. <i>Physical Review Letters</i> , 2001, 86, 1861-1864.	2.9	80
61	Electronic structure of materials under pressure. <i>International Journal of Quantum Chemistry</i> , 2000, 77, 880-894.	1.0	11
62	High-pressure Cmca and hcp phases of germanium. <i>Physical Review B</i> , 2000, 62, R10603-R10606.	1.1	64
63	First-principles investigation of SrBi ₂ Ta ₂ O ₉ . <i>Ferroelectrics</i> , 2000, 237, 49-56.	0.3	5
64	Computer simulations of defects in perovskite KNbO ₃ crystals. <i>Ferroelectrics</i> , 1999, 229, 69-75.	0.3	6
65	Solids under Pressure. Ab Initio Theory. <i>Physica Status Solidi (B): Basic Research</i> , 1999, 211, 5-16.	0.7	33
66	Structural and vibronic properties of perovskites studied by using the Perdewâ€“Burkeâ€“Ernzerhof GGA. , 1998, , .		0
67	Native defects and carbon impurity in cubic BN. <i>MRS Internet Journal of Nitride Semiconductor Research</i> , 1998, 3, 1.	1.0	5
68	Cubic Inclusions as the Cause for the Unusually Weak Pressure Shift of the Luminescence in InGaN. <i>Materials Research Society Symposia Proceedings</i> , 1997, 482, 720.	0.1	0
69	Theoretical study of point defects in GaN and AlN; lattice relaxations and pressure effects. <i>MRS Internet Journal of Nitride Semiconductor Research</i> , 1997, 2, 1.	1.0	12
70	Electronic Structures of Semiconductors under Pressure. <i>Physica Status Solidi (B): Basic Research</i> , 1996, 198, 23-34.	0.7	34
71	First-Principles Simulations of Interstitial Atoms in Ionic Solids. <i>Materials Research Society Symposia Proceedings</i> , 1995, 408, 509.	0.1	1
72	Optical and structural properties of III-V nitrides under pressure. <i>Physical Review B</i> , 1994, 50, 4397-4415.	1.1	640

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73	Semiconductor heterojunctions and superlattices: Band offsets and electronic structures. The Philosophical Magazine: Physics of Condensed Matter B, Statistical Mechanics, Electronic, Optical and Magnetic Properties, 1994, 70, 567-581.	0.6	4
74	Bonding and ionicity in semiconductors. Physical Review B, 1987, 36, 1032-1050.	1.1	274
75	Electronic structure of GaAs under strain. Physical Review B, 1984, 30, 5753-5765.	1.1	299