

Konstantin E Glukhov

List of Publications by Year in descending order

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40
papers

270
citations

1040056

9
h-index

1058476

14
g-index

45
all docs

45
docs citations

45
times ranked

274
citing authors

#	ARTICLE	IF	CITATIONS
1	The first-principle study of substitutional impuritiesâ€™ effect on elastic properties of TlInS ₂ layered crystal. <i>Low Temperature Physics</i> , 2022, 48, 57-63.	0.6	0
2	Phonon spectra and phase transitions in van der Waals ferroics MMâ€™P ₂ X ₆ . <i>Molecular Crystals and Liquid Crystals</i> , 2022, 747, 14-22.	0.9	4
3	Theoretical and Experimental Studies of Electronic and Optical Properties of Layered TlIn(S _{0.75} Se _{0.25}) ₂ Ferroelectric Crystal. <i>Integrated Ferroelectrics</i> , 2021, 220, 18-29.	0.7	4
4	Layered GeP ₂ S ₆ , GeP ₂ Se ₆ , GeP ₂ Te ₆ , SnP ₂ S ₆ , SnP ₂ Se ₆ , and SnP ₂ Te ₆ Polar Crystals with Semiconductorâ€™Metal Transitions Induced by Pressure or Chemical Composition. <i>Integrated Ferroelectrics</i> , 2021, 220, 90-99.	0.7	5
5	Nature of thermoelectric properties occurring in defected Sn ₂ P ₂ S ₆ chalcogenide crystals. <i>CrystEngComm</i> , 2020, 22, 2336-2349.	2.6	7
6	Electronic and vibrational properties of pure MnPS ₃ crystal: Theoretical and experimental investigation. <i>Computational Materials Science</i> , 2020, 177, 109592.	3.0	9
7	Raman study of a magnetic phase transition in the MnPS ₃ single crystal. <i>Low Temperature Physics</i> , 2019, 45, 1082-1091.	0.6	9
8	Layered ferrielectric crystals CuInP ₂ S(Se) ₆ : a study from the first principles. <i>Phase Transitions</i> , 2019, 92, 440-450.	1.3	14
9	Band structures and optical properties related to substitutional impurities in TlGaSe ₂ layered crystals: first-principles study. <i>Phase Transitions</i> , 2019, 92, 451-460.	1.3	6
10	Thermal diffusivity and thermal conductivity in layered ferrielectric materials M ¹⁺ M ³⁺ P ₂ (S,Se) ₆ (M ¹⁺ =Cu, Ag; Tj ETQ 0 0 0 BT /Overl		
11	Cation role in the thermal properties of layered materials <mml:math xmlns:mml="http://www.w3.org/1998/Math/MathML"><mml:mrow><mml:msup><mml:mrow><mml:mi>M</mml:mi></mml:mrow></mml:math>		

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19	Electron-Deformational Phase Transitions in a TlGaSe_2 Layered Crystal. <i>Acta Physica Polonica A</i> , 2016, 129, A-123-A-125.	0.5	3
20	Construction of the Adiabatic Potential of a Symmetric Molecule in the Vicinity of Charged Semiconductor Surface. <i>Acta Physica Polonica A</i> , 2016, 129, A-120-A-122.	0.5	0
21	Vibronic interaction in crystals with the Jahn-Teller centers in the elementary energy bands concept. <i>Condensed Matter Physics</i> , 2015, 18, 33705.	0.7	4
22	Peculiarities of Chemical Bonding in Crystals of the In-Se System. <i>Acta Physica Polonica A</i> , 2014, 126, 1146-1148.	0.5	3
23	Electron-Phonon Interaction as a Mechanism of Phase Transition in the CuInP_2S_6 Crystal. <i>Acta Physica Polonica A</i> , 2014, 126, 1143-1145.	0.5	9
24	Chemical Bonding and Polarons in $\text{Sn}_2\text{P}_2\text{S}_6(\text{Se})_6$ Ferroelectrics. <i>Ferroelectrics</i> , 2014, 462, 117-128.	0.6	6
25	Exact ground state for the four-electron problem in a 2D finite honeycomb lattice. <i>Philosophical Magazine</i> , 2014, 94, 2195-2223.	1.6	0
26	Electronic structure of low-pressure and high-pressure phases of silicon disulfide. <i>Applied Physics A: Materials Science and Processing</i> , 2014, 117, 1499-1514.	2.3	6
27	Optical properties and band structure of a layered Tl_2S crystal. <i>Physics of the Solid State</i> , 2013, 55, 2317-2323.	0.6	4
28	Electron structure of the equilibrium and metastable phases in superionic Li_2SiS_3 . <i>Semiconductor Physics, Quantum Electronics and Optoelectronics</i> , 2013, 16, 48-54.	1.0	2
29	Ferroelectricity and Polarons in $\text{Sn}_2\text{P}_2\text{S}_6$ Crystals. <i>Ferroelectrics</i> , 2012, 440, 31-41.	0.6	12
30	Electronic Structure and Phase Transition in Ferroelectric $\text{Sn}_2\text{P}_2\text{S}_6$ Crystal. <i>International Journal of Molecular Sciences</i> , 2012, 13, 14356-14384.	4.1	41
31	Parameters of an Unique Condensation State in the Structure of the In_4Se_3 Crystal. <i>Acta Physica Polonica A</i> , 2012, 122, 1115-1117.	0.5	2
32	Charge Transfer and Anharmonicity in $\text{Sn}_2\text{P}_2\text{S}_6$ Ferroelectrics. <i>Ferroelectrics</i> , 2011, 414, 30-40.	0.6	16
33	Ferroelectric and Semiconducting Properties of $\text{Sn}_2\text{P}_2\text{S}_6$ Crystals with Intrinsic Vacancies. <i>Ferroelectrics</i> , 2011, 418, 124-133.	0.6	21
34	XPS of Impurities Influence on Electronic Structure of $\text{Sn}_2\text{P}_2\text{S}_6$ Ferroelectrics. <i>Ferroelectrics</i> , 2011, 418, 134-142.	0.6	5
35	Non-standard anisotropy of the energy spectrum of a layered TlGaSe_2 crystal. <i>Physica Status Solidi (B): Basic Research</i> , 2011, 248, 1446-1452.	1.5	2
36	Gain Spectrum for the In_4Se_3 Crystal with a Non-Standard Dispersion Law of Charge Carriers. <i>Acta Physica Polonica A</i> , 2011, 119, 720-722.	0.5	2

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37	Elementary energy bands concept, band structure, and peculiarities of bonding in $\hat{1}^2$ -InSe crystal. <i>Physica Status Solidi (B): Basic Research</i> , 2010, 247, 318-324.	1.5	7
38	Elementary energy bands in the band structure of AIV, AIIIBV crystals and superlattices built upon them. <i>Physica Status Solidi (B): Basic Research</i> , 2007, 244, 1318-1336.	1.5	7
39	Energy states in short-period symmetrical and asymmetrical (GaAs)N/(AlAs)M superlattices: The effect of the boundary conditions. <i>Semiconductors</i> , 2004, 38, 410-418.	0.5	2
40	Covalent bridges - an induced modification of the conduction band in layered crystals. <i>Physica Status Solidi (B): Basic Research</i> , 2004, 241, 845-855.	1.5	0