

# Oleg V Bovgyra

## List of Publications by Year in descending order

Source: <https://exaly.com/author-pdf/2320015/publications.pdf>

Version: 2024-02-01

26  
papers

117  
citations

1478505

6  
h-index

1372567

10  
g-index

27  
all docs

27  
docs citations

27  
times ranked

81  
citing authors

#	ARTICLE	IF	CITATIONS
1	First principle study of electronic properties of ZnO nanoclusters with native point defects during gas adsorption. Applied Nanoscience (Switzerland), 2022, 12, 983-993.	3.1	1
2	Manifestation of the ferroelastoelectric phase transition in the absorption spectra of (NH <sub>4</sub> ) <sub>2</sub> CuCl <sub>4</sub> ·2D <sub>2</sub> O crystals. Physica B: Condensed Matter, 2022, , 413929.	2.7	1
3	Growth, crystal structure and optical properties of Al-doped ZnO thin films. Molecular Crystals and Liquid Crystals, 2021, 717, 72-79.	0.9	2
4	Structural, Electronic and Optical Properties of CsPbBr <sub>3</sub> and CH <sub>3</sub> NH <sub>3</sub> PbBr <sub>3</sub> : First-Principles Modeling. , 2021, , .		0
5	Structure and electronic properties of CsPbBr <sub>3</sub> perovskite: first principle calculations. Journal of Physical Studies, 2021, 25, .	0.5	1
6	A DFT study for adsorption of CO and H <sub>2</sub> on Pt-doped ZnO nanocluster. SN Applied Sciences, 2020, 2, 1.	2.9	13
7	Electronic Properties of Al-, Ga-, and In-Doped Armchair ZnO Nanoribbons. , 2019, , .		3
8	First principle study of native point defects in (ZnO) <sub>n</sub> nanoclusters (n=34, 60). Applied Nanoscience (Switzerland), 2019, 9, 1067-1074.	3.1	6
9	Specific Features of Content Dependences for Energy Gap in In <sub>x</sub> Tl <sub>1-x</sub> Solid State Crystalline Alloys. Acta Physica Polonica A, 2018, 133, 68-75.	0.5	16
10	Lattice vibration spectra of A <sub>4</sub> BX <sub>6</sub> group crystals. Journal of Physical Studies, 2018, 22, .	0.5	4
11	Ab Initio Study of Structural and Electronic Properties of (ZnO) <sub>n</sub> "Magical" Nanoclusters (n=34, 60). Nanoscale Research Letters, 2017, 12, 76.	5.7	19
12	Temperature behavior of thermal expansion and birefringence of In <sub>x</sub> Tl <sub>1-x</sub> substitution solid solutions. Optics and Spectroscopy (English Translation of Optika i Spektroskopiya), 2017, 123, 177-180.	0.6	3
13	Birefringence of In <sub>x</sub> Tl <sub>1-x</sub> solid state solution. Functional Materials, 2017, 23, 026-030.	0.1	4
14	Elastic Properties of Substitutional Solid Solutions In <sub>x</sub> Tl <sub>1-x</sub> and Sound Wave Velocities in Them. Ukrainian Journal of Physics, 2017, 62, 679-684.	0.2	7
15	Nonlocal Model Pseudopotential Calculations of the Electronic Structure of AlIBVI (CdS, CdSe) Bulk Crystals and Nanocrystals. Journal of Nano- and Electronic Physics, 2017, 9, 02030-1-02030-7.	0.5	2
16	Anisotropy of inter-band transitions and band structure of Cs <sub>3</sub> Zn <sub>6</sub> B <sub>9</sub> O <sub>21</sub> nonlinear optical crystals. Optical Materials, 2016, 56, 129-133.	3.6	2
17	Electronic Structure, Optical and Sensor Properties of ZnO Nanowires. Journal of Nano- and Electronic Physics, 2016, 8, 02031-1-02031-5.	0.5	3
18	Piezo-optical properties of incommensurately modulated Rb <sub>2</sub> ZnCl <sub>4</sub> crystals. Optics and Spectroscopy (English Translation of Optika i Spektroskopiya), 2015, 118, 547-551.	0.6	0

#	ARTICLE	IF	CITATIONS
19	Band Structure and Birefringence of RbKSO <sub>4</sub> Crystals. Journal of Applied Spectroscopy, 2015, 82, 755-759.	0.7	1
20	DFT study of electronic and magnetic properties of bare and substitutionally doped ZnO nanoribbons. , 2015, , .		6
21	First-principles simulations of the electronic density of states for superionic Ag <sub>2</sub> CdI <sub>4</sub> crystals. Solid State Ionics, 2011, 188, 31-35.	2.7	3
22	Fast intrinsic emission in single crystal. Radiation Measurements, 2007, 42, 869-873.	1.4	1
23	The band energy structure of RbKSO <sub>4</sub> crystals. Condensed Matter Physics, 2007, 10, 39.	0.7	2
24	Energy band structure and refractive properties of LiRbSO <sub>4</sub> crystals. Physics of the Solid State, 2006, 48, 1268-1272.	0.6	12
25	Optical constants of indium bromide. Low Temperature Physics, 2001, 27, 153-157.	0.6	1
26	Title is missing!. Ukrainian Journal of Physical Optics, 2000, 1, 49-52.	13.0	0