

# Alexey Lukoyanov

## List of Publications by Citations

**Source:** <https://exaly.com/author-pdf/9316433/alexey-lukoyanov-publications-by-citations.pdf>

**Version:** 2024-04-25

This document has been generated based on the publications and citations recorded by exaly.com. For the latest version of this publication list, visit the link given above.

The third column is the impact factor (IF) of the journal, and the fourth column is the number of citations of the article.

164  
papers

1,533  
citations

19  
h-index

34  
g-index

174  
ext. papers

1,688  
ext. citations

2.2  
avg, IF

4.58  
L-index

#	Paper	IF	Citations
164	Collapse of magnetic moment drives the Mott transition in MnO. <i>Nature Materials</i> , <b>2008</b> , 7, 198-202	27	145
163	Magnetic state and electronic structure of the $\Gamma$ and $\Phi$ phases of metallic Pu and its compounds. <i>Physical Review B</i> , <b>2005</b> , 72,	3.3	135
162	NiO: correlated band structure of a charge-transfer insulator. <i>Physical Review Letters</i> , <b>2007</b> , 99, 156404	7.4	122
161	Local correlations and hole doping in NiO: A dynamical mean-field study. <i>Physical Review B</i> , <b>2007</b> , 75,	3.3	86
160	Doped Mott insulator as the origin of heavy-fermion behavior in LiV <sub>2</sub> O <sub>4</sub> . <i>Physical Review Letters</i> , <b>2007</b> , 98, 166402	7.4	50
159	The role of transition metal impurities and oxygen vacancies in the formation of ferromagnetism in Co-doped TiO <sub>2</sub> . <i>Journal of Physics Condensed Matter</i> , <b>2006</b> , 18, 1695-1704	1.8	46
158	Metal-insulator transitions and magnetism in correlated band insulators: FeSi and Fe <sub>1-x</sub> Co <sub>x</sub> Si. <i>Physical Review B</i> , <b>2010</b> , 81,	3.3	43
157	Origin of large thermopower in LiRh <sub>2</sub> O <sub>4</sub> : Calculation of the Seebeck coefficient by the combination of local density approximation and dynamical mean-field theory. <i>Physical Review B</i> , <b>2008</b> , 78,	3.3	42
156	The semiconductor-to-ferromagnetic-metal transition in FeSb <sub>2</sub> . <i>European Physical Journal B</i> , <b>2006</b> , 53, 205-207	1.2	42
155	Vacancies in ordered and disordered titanium monoxide: Mechanism of B1 structure stabilization. <i>Journal of Solid State Chemistry</i> , <b>2013</b> , 204, 146-152	3.3	31
154	Ab initio exchange interactions and magnetic properties of the Gd <sub>2</sub> Fe <sub>17</sub> iron sublattice: Rhombohedral versus hexagonal phases. <i>Physical Review B</i> , <b>2009</b> , 80,	3.3	31
153	Evolution of the electronic structure and physical properties of Fe <sub>2</sub> MeAl (Me = Ti, V, Cr) Heusler alloys. <i>Journal of Physics Condensed Matter</i> , <b>2008</b> , 20, 045212	1.8	31
152	Electronic structure of nonstoichiometric compounds in the coherent potential approximation. <i>JETP Letters</i> , <b>2012</b> , 94, 806-810	1.2	29
151	Magnetic and electrical properties of the half-metallic ferromagnets Co <sub>2</sub> CrAl. <i>Physics of the Solid State</i> , <b>2013</b> , 55, 977-985	0.8	26
150	Effect of complex magnetic structure on the magnetocaloric and magneto-transport properties in GdCuSi. <i>Journal of Materials Science</i> , <b>2015</b> , 50, 5723-5728	4.3	25
149	Electronic structure, magnetic, and optical properties of the intermetallic compounds R <sub>2</sub> Fe <sub>17</sub> (R=Pr,Gd). <i>Physical Review B</i> , <b>2006</b> , 73,	3.3	24
148	LDA+DMFT spectral functions and effective electron mass enhancement in the superconductor LaFePO. <i>Physical Review B</i> , <b>2010</b> , 81,	3.3	21

147	Pressure-driven metal-insulator transition in BiFeO <sub>3</sub> from dynamical mean-field theory. <i>Physical Review B</i> , <b>2015</b> , 92,	3.3	19
146	Electronic structure and optical spectroscopy studies of HoNi <sub>5</sub> and ErNi <sub>5</sub> compounds doped with Cu. <i>Physica Status Solidi (B): Basic Research</i> , <b>2012</b> , 249, 824-828	1.3	19
145	Magnetism in RRhGe (R=Tb, Dy, Er, Tm): An experimental and theoretical study. <i>Journal of Alloys and Compounds</i> , <b>2015</b> , 640, 56-63	5.7	18
144	The magnetic, electronic and optical properties of HoRhGe. <i>Journal Physics D: Applied Physics</i> , <b>2014</b> , 47, 365002	3	17
143	Inclusion of the correlation short-range order in Ab initio calculations of the energy of the ground state by example of titanium monoxide TiO <sub>1.0</sub> . <i>JETP Letters</i> , <b>2015</b> , 102, 85-90	1.2	17
142	Electronic structure of disordered titanium monoxide TiO <sub>y</sub> depending on stoichiometry. <i>JETP Letters</i> , <b>2012</b> , 95, 647-651	1.2	17
141	LDA+DMFT study of magnetic transition and metallization in CoO under pressure. <i>JETP Letters</i> , <b>2012</b> , 96, 56-60	1.2	17
140	Calculation of temperature dependence of electrical resistivity in the transuranium metals and their alloys. <i>Physical Review B</i> , <b>2007</b> , 76,	3.3	16
139	Optical conductivity of ortho-II YBa <sub>2</sub> Cu <sub>3</sub> O <sub>6.5</sub> . <i>Physical Review B</i> , <b>2005</b> , 71,	3.3	16
138	Impression of magnetic clusters, critical behavior and magnetocaloric effect in FeAl alloys. <i>Physical Chemistry Chemical Physics</i> , <b>2019</b> , 21, 10823-10833	3.6	15
137	Simulation of the short-range order in disordered cubic titanium monoxide TiO <sub>1.0</sub> . <i>JETP Letters</i> , <b>2013</b> , 97, 616-620	1.2	14
136	Specific features of the electrical resistance of half-metallic ferromagnetic alloys Co <sub>2</sub> CrAl and Co <sub>2</sub> CrGa. <i>Physics of the Solid State</i> , <b>2013</b> , 55, 2487-2490	0.8	14
135	Sm <sub>2</sub> Fe <sub>17</sub> and Tm <sub>2</sub> Fe <sub>17</sub> : electronic structure, magnetic and optical properties. <i>Journal of Physics Condensed Matter</i> , <b>2007</b> , 19, 116215	1.8	14
134	Internal energy and parameters of the order-disorder phase transition in titanium monoxide TiO <sub>y</sub> . <i>Journal of Experimental and Theoretical Physics</i> , <b>2013</b> , 116, 945-951	1	12
133	Effect of the long-range order in the vacancy distribution on the electronic structure of titanium monoxide TiO <sub>1.0</sub> . <i>JETP Letters</i> , <b>2012</b> , 96, 507-510	1.2	12
132	Effect of Cu-doping on the electronic structure and optical properties of LaNi <sub>5</sub> . <i>Journal of Alloys and Compounds</i> , <b>2011</b> , 509, 5238-5241	5.7	12
131	Effect of the Structural Disorder and Short-Range Order on the Electronic Structure and Magnetic Properties of the Fe <sub>2</sub> VAl Heusler Alloy. <i>JETP Letters</i> , <b>2018</b> , 107, 126-128	1.2	10
130	Cobalt-related features of spectral and magnetic properties of RNi <sub>4</sub> Co (R=Ho, Er). <i>Journal of Magnetism and Magnetic Materials</i> , <b>2014</b> , 368, 87-90	2.8	10

129	Magnetism, electronic structure and optical properties of TbNiGe <sub>2</sub> . <i>Journal of Alloys and Compounds</i> , <b>2016</b> , 664, 120-124	5.7	9
128	Optical properties and electronic structure of alloys Co <sub>2</sub> Cr <sub>1-x</sub> Fe <sub>x</sub> Al (x = 0, 0.4, 0.6, 1.0). <i>Physics of the Solid State</i> , <b>2016</b> , 58, 164-169	0.8	9
127	Orbital density functional as a means to restore the discontinuities in the total-energy derivative and the exchange-correlation potential. <i>Journal of Physics Condensed Matter</i> , <b>2007</b> , 19, 106206	1.8	9
126	Features of electronic properties of band ferromagnets Co <sub>2</sub> MeAl and Fe <sub>2</sub> MeAl (Me = Ti, V, Cr, Mn, Fe, Ni). <i>Materials Research Express</i> , <b>2017</b> , 4, 116102	1.7	8
125	Optical properties and the electronic structure of Co <sub>2</sub> TiGe and Co <sub>2</sub> TiSn Heusler alloys. <i>Physics of Metals and Metallography</i> , <b>2017</b> , 118, 965-969	1.2	8
124	Structural stability and magnetic properties of Mn <sub>2</sub> FeAl alloy with a $\sqrt{5}$ Mn structure. <i>Journal of Magnetism and Magnetic Materials</i> , <b>2020</b> , 513, 167205	2.8	8
123	Magnetic properties and exchange interactions in TbNi <sub>5-x</sub> M <sub>x</sub> (M=Co and Fe) compounds: Ab initio calculations. <i>Journal of Applied Physics</i> , <b>2011</b> , 109, 07E152	2.5	8
122	Specific features of the behavior of the optical properties of TbNi <sub>5-x</sub> Cu <sub>x</sub> intermetallic compounds. <i>Optics and Spectroscopy (English Translation of Optika i Spektroskopiya)</i> , <b>2008</b> , 104, 360-365	0.7	8
121	Ordering Sequence in Strongly Nonstoichiometric Niobium Carbide with the Formation of Nb <sub>6</sub> C <sub>5</sub> -Type Superstructures. <i>Journal of Experimental and Theoretical Physics</i> , <b>2019</b> , 129, 863-876	1	8
120	Electron Structure and Optical Properties of the Mn <sub>1.8</sub> Co <sub>1.2</sub> Al Alloy and Spin Gapless Semiconductor State. <i>Physics of Metals and Metallography</i> , <b>2018</b> , 119, 1068-1072	1.2	8
119	A comparative study of the optical properties of TbRhGe and DyRhGe. <i>Solid State Sciences</i> , <b>2015</b> , 44, 22-26	3.4	7
118	Understanding the magnetic, electronic and optical properties of ternary rare earth intermetallic compound HoNiSi. <i>Journal of Alloys and Compounds</i> , <b>2015</b> , 650, 542-546	5.7	7
117	Two successive spin transitions in a wide range of pressure and coexistence of high- and low-spin states in clinoferrosilite FeSiO <sub>3</sub> . <i>Physical Review B</i> , <b>2016</b> , 93,	3.3	7
116	Electronic structure and magnetic state of transuranium metals under pressure. <i>Journal of Physics Condensed Matter</i> , <b>2010</b> , 22, 495501	1.8	7
115	Transition of iron ions from high-spin to low-spin state and pressure-induced insulator-metal transition in hematite Fe <sub>2</sub> O <sub>3</sub> . <i>Journal of Experimental and Theoretical Physics</i> , <b>2007</b> , 105, 1035-1042	1	7
114	Anomalous concentration dependence of residual electrical resistivity in Fe-Cr alloys. <i>Physical Review B</i> , <b>2005</b> , 72,	3.3	7
113	Correlational short-range order in superstructures. <i>Mendeleev Communications</i> , <b>2017</b> , 27, 147-149	1.9	6
112	Influence of aluminum impurity on the electronic structure and optical properties of the TbNi <sub>5</sub> intermetallic compound. <i>Physics of the Solid State</i> , <b>2013</b> , 55, 385-388	0.8	6

111	Electronic structure and magnetic properties of PuMgAs compounds within the LDA + U + SO method. <i>JETP Letters</i> , <b>2012</b> , 96, 452-455	1.2	6
110	Nature of the electronic states involved in the chemical bonding and superconductivity at high pressure in SnO. <i>JETP Letters</i> , <b>2011</b> , 94, 142-146	1.2	6
109	Optical absorption and structure of energy bands of GdNi <sub>5</sub> & Cu <sub>x</sub> intermetallic compounds. <i>Physics of Metals and Metallography</i> , <b>2009</b> , 107, 173-178	1.2	6
108	Theoretical and experimental investigations on the magnetic and related properties of RAgSn <sub>2</sub> (R=Ho, Er) compounds. <i>Journal of Materials Science</i> , <b>2016</b> , 51, 6341-6347	4.3	6
107	Vacancy ordered structures in a nonstoichiometric niobium carbide NbC <sub>0.83</sub> . <i>Mendeleev Communications</i> , <b>2019</b> , 29, 707-709	1.9	6
106	Experimental and theoretical investigations on magnetic and related properties of ErRuSi. <i>Materials Research Express</i> , <b>2015</b> , 2, 046101	1.7	5
105	Crystal and electronic structure of high temperature superconducting compound Y <sub>1-x</sub> Ca <sub>x</sub> Ba <sub>2</sub> Cu <sub>3</sub> O <sub>y</sub> in the temperature interval 80-300K. <i>Journal of Alloys and Compounds</i> , <b>2016</b> , 658, 891-897	5.7	5
104	Electronic structure and optical properties of Nd <sub>5</sub> Ge <sub>3</sub> compound. <i>Journal of Alloys and Compounds</i> , <b>2014</b> , 588, 725-727	5.7	5
103	Electronic structure and stability of nonstoichiometric titanium monoxide TiO <sub>y</sub> with structural vacancies in one of the sublattices. <i>Physics of the Solid State</i> , <b>2013</b> , 55, 2108-2115	0.8	5
102	Investigation of real materials with strong electronic correlations by the LDA+DMFT method. <i>Acta Crystallographica Section C, Structural Chemistry</i> , <b>2014</b> , 70, 137-59	0.8	5
101	Correlation effects in Ni 3d states of LaNiPO. <i>Physical Review B</i> , <b>2010</b> , 81,	3.3	5
100	Theoretical investigation of the residual electrical resistivity concentration dependence of transuranium metal alloys. <i>Physical Review B</i> , <b>2009</b> , 80,	3.3	5
99	Electronic structure of the intermetallic compounds Ce <sub>2</sub> Fe <sub>17</sub> and Ce <sub>2</sub> Fe <sub>15.3</sub> M <sub>1.7</sub> (M = Al, Si): Experiment and theory. <i>Physics of the Solid State</i> , <b>2007</b> , 49, 99-106	0.8	5
98	Remarkable increase of Curie temperature in doped GdFeSi compound. <i>Intermetallics</i> , <b>2021</b> , 133, 107183,5	3.5	5
97	Electronic Structure and Electronic Properties of PtSn <sub>4</sub> Single Crystal. <i>Journal of Experimental and Theoretical Physics</i> , <b>2019</b> , 128, 939-945	1	4
96	Structural, Electronic, Optical, and Magnetic Properties of Fe <sub>3</sub> Al Alloys. <i>Journal of Superconductivity and Novel Magnetism</i> , <b>2019</b> , 32, 2995-3000	1.5	4
95	Short-range order in disordered transition metal oxides, carbides, and nitrides with the B1 structure. <i>Physics of the Solid State</i> , <b>2015</b> , 57, 637-651	0.8	4
94	Itinerant metamagnetic transition in the ferromagnet LuCo <sub>3</sub> induced by high field: Instability of the 3d-electron subsystem. <i>Physical Review B</i> , <b>2020</b> , 101,	3.3	4

93	Pressure-induced modification of the electron structure of metallic thorium. <i>Journal of Experimental and Theoretical Physics</i> , <b>2014</b> , 118, 148-152	1	4
92	Influence of copper impurities on the evolution of the electronic structure and optical spectra of the LuNi5 compound. <i>Physics of the Solid State</i> , <b>2015</b> , 57, 866-870	0.8	4
91	Coherent potential approximation simulation of the evolution of the electronic structure of titanium monoxide with the degree of vacancy ordering. <i>Journal of Experimental and Theoretical Physics</i> , <b>2014</b> , 119, 761-765	1	4
90	Disorder-order and order-order phase transformations in TaC phases predicted using the evolutionary algorithm and symmetry analysis. <i>Physical Chemistry Chemical Physics</i> , <b>2020</b> , 22, 24116-24132 <sup>3,6</sup>	3.6	4
89	Specific features of the electronic structure and spectral characteristics of the Gd5Si3 compound. <i>Physics of the Solid State</i> , <b>2017</b> , 59, 429-433	0.8	3
88	Evolution of the electronic structure and optical spectra of intermetallides DyNi5 1-x Cu x under changes of concentration. <i>Optics and Spectroscopy (English Translation of Optika i Spektroskopiya)</i> , <b>2015</b> , 118, 357-363	0.7	3
87	Electronic and Spectral Properties of RRhSn (R = Gd, Tb) Intermetallic Compounds. <i>Physics of the Solid State</i> , <b>2018</b> , 60, 225-229	0.8	3
86	Electronic structure and spectral properties of RCuSi (R=Nd,Gd) compounds. <i>Physica B: Condensed Matter</i> , <b>2016</b> , 487, 85-89	2.8	3
85	Optical spectroscopy and electronic structure of the GdCu x compounds (x = 1, 2, 5). <i>Physics of the Solid State</i> , <b>2013</b> , 55, 140-144	0.8	3
84	Calculation of the electronic structure of the intermetallic compounds ErNi5 1-x Al x (x = 0, 1, 2). <i>Physics of the Solid State</i> , <b>2015</b> , 57, 1-4	0.8	3
83	Low-temperature heat capacity upon the transition from paramagnetic to ferromagnetic Heusler alloys Fe2 MeAl (Me = Ti, V, Cr, Mn, Fe, Co, Ni). <i>Physics of the Solid State</i> , <b>2016</b> , 58, 1500-1504	0.8	3
82	Electronic Structure of Intermetallic Antiferromagnet GdNiGe. <i>Symmetry</i> , <b>2019</b> , 11, 737	2.7	3
81	Electronic structure of DyRhSn and HoRhSn compounds: band calculations and optical study. <i>European Physical Journal B</i> , <b>2019</b> , 92, 1	1.2	2
80	Characterization of d and f Electronic States in RSn1.1Ge0.9 (R = Gd, Tb) Compounds by Optical Spectroscopy and Electronic-Structure Calculations. <i>Physica Status Solidi (B): Basic Research</i> , <b>2018</b> , 255, 1700579	1.3	2
79	Raman scattering by electron and phonon excitations in FeSi. <i>JETP Letters</i> , <b>2016</b> , 103, 316-320	1.2	2
78	Electronic structure of nitrides PuN and UN. <i>Journal of Experimental and Theoretical Physics</i> , <b>2016</b> , 123, 864-868	1	2
77	Magnetically driven phase transitions with a large volume collapse in MnSe under pressure: A DFT+DMFT study. <i>Physical Review B</i> , <b>2018</b> , 98,	3.3	2
76	Magnetization, resistivity, specific heat and ab initio calculations of GdSb. <i>Journal of Physics Condensed Matter</i> , <b>2018</b> , 30, 295802	1.8	2

75	Electronic structure and magnetic susceptibility of monoclinic plutonium. <i>JETP Letters</i> , <b>2014</b> , 99, 656-660	2
74	Electronic structure and optical properties of Er <sub>5</sub> Si <sub>3</sub> . <i>Physica B: Condensed Matter</i> , <b>2014</b> , 442, 12-15	2.8 2
73	Optical spectroscopy and electronic structure of compounds HoNi <sub>5</sub> [x Al x (x = 0, 1, 2). <i>Optics and Spectroscopy (English Translation of Optika I Spektroskopiya)</i> , <b>2013</b> , 115, 690-695	0.7 2
72	Effect of manganese doping on the electronic structure and optical properties of Ce <sub>2</sub> Fe <sub>17-x</sub> Mn <sub>x</sub> (x = 0, 1, 2). <i>European Physical Journal B</i> , <b>2017</b> , 90, 1	1.2 2
71	Electronic structure of RSn <sub>1.1</sub> Ge <sub>0.9</sub> (R = Dy, Ho) ternary compounds: Band calculation and optical properties. <i>Physica B: Condensed Matter</i> , <b>2017</b> , 521, 98-101	2.8 2
70	Copper-doping effects in electronic structure and spectral properties of SmNi <sub>5</sub> . <i>Low Temperature Physics</i> , <b>2015</b> , 41, 1024-1028	0.7 2
69	Optical spectroscopy and electronic structure of TmRhGe compound. <i>Physics of the Solid State</i> , <b>2015</b> , 57, 2357-2360	0.8 2
68	Role of Fe and Co in optical conductivity and electronic structure of TbNi <sub>4</sub> Fe and TbNi <sub>4</sub> Co. <i>Optics and Spectroscopy (English Translation of Optika I Spektroskopiya)</i> , <b>2014</b> , 117, 414-418	0.7 2
67	Electrical resistivity of pure transuranium metals under pressure. <i>Journal of Nuclear Materials</i> , <b>2011</b> , 413, 41-46	3.3 2
66	Optical Properties and Electronic Structure of LaNi <sub>5</sub> -XCu <sub>x</sub> (x=0-1.2) Intermetallic System. <i>Solid State Phenomena</i> , <b>2010</b> , 168-169, 529-532	0.4 2
65	Temperature and concentration dependences of the electrical resistivity for alloys of plutonium with americium under normal conditions. <i>Journal of Experimental and Theoretical Physics</i> , <b>2010</b> , 111, 1019-1027	2
64	Neutron and Raman studies of lattice distortions in Zn <sub>1-x</sub> Ni <sub>x</sub> Se induced by nickel. <i>Physica Status Solidi C: Current Topics in Solid State Physics</i> , <b>2004</b> , 1, 3150-3153	2
63	Optical properties and electronic structure of the CeFeSi-type GdTIGe and GdTisi compounds. <i>Journal of Alloys and Compounds</i> , <b>2004</b> , 384, 57-61	5.7 2
62	Electronic Structure and Optical Properties of the Mn <sub>2</sub> CrAl Heusler Alloy. <i>Physics of Metals and Metallography</i> , <b>2020</b> , 121, 532-536	1.2 2
61	Electronic properties of WTe <sub>2</sub> and MoTe <sub>2</sub> single crystals. <i>Journal of Physics: Conference Series</i> , <b>2019</b> , 1389, 012149	0.3 2
60	Electronic Structure and Magnetic Properties of Strongly Correlated Transition Metal Compounds. <i>Physics of Metals and Metallography</i> , <b>2018</b> , 119, 1254-1258	1.2 2
59	A wide energy range ab initio modeling of the electronic structure of valence states in Cu(In,Ga)Se <sub>2</sub> : Comparison with photoelectron spectra. <i>Journal of Alloys and Compounds</i> , <b>2019</b> , 802, 19-24	5.7 1
58	Electronic structure of the TbMn <sub>0.33</sub> Ge <sub>2</sub> compound: Band calculation and optical experiment. <i>Physics of the Solid State</i> , <b>2016</b> , 58, 2373-2378	0.8 1

57	Electronic structure and optical properties of the HoCoSi and ErNiSi compounds. <i>Journal of Experimental and Theoretical Physics</i> , <b>2016</b> , 123, 638-642	1	1
56	Anisotropy of the Complex Permittivity of the Kagome-Staircase Compounds Co <sub>3</sub> V <sub>2</sub> O <sub>8</sub> and Ni <sub>3</sub> V <sub>2</sub> O <sub>8</sub> : Experiment and Ab Initio Calculations. <i>Journal of Experimental and Theoretical Physics</i> , <b>2018</b> , 126, 779-783	1	1
55	Magnetism and electronic structure of Gd <sub>5</sub> Ge <sub>2</sub> Sb: Experiment and theory. <i>Journal of Alloys and Compounds</i> , <b>2019</b> , 806, 575-579	5.7	1
54	Optical spectroscopy and electronic structure of the Er <sub>5</sub> Ge <sub>3</sub> compound. <i>Physics of the Solid State</i> , <b>2014</b> , 56, 1737-1741	0.8	1
53	Role of structural vacancies in the stabilization of the basic B1 structure in nonstoichiometric titanium monoxide TiO <sub>y</sub> . <i>Bulletin of the Russian Academy of Sciences: Physics</i> , <b>2013</b> , 77, 309-312	0.4	1
52	Specific features of the electronic structure and spectral properties of NdNi <sub>5-x</sub> Cu <sub>x</sub> compounds. <i>Physics of the Solid State</i> , <b>2013</b> , 55, 2191-2195	0.8	1
51	Electronic structure and optical properties of the Pr <sub>5</sub> Ge <sub>3</sub> compound. <i>Physics of the Solid State</i> , <b>2015</b> , 57, 1705-1709	0.8	1
50	Electronic structure of the PuCoIn <sub>5</sub> compound. <i>JETP Letters</i> , <b>2015</b> , 101, 402-406	1.2	1
49	Optical properties and electronic structure of YNi <sub>5-x</sub> Cu <sub>x</sub> intermetallic compounds. <i>Optics and Spectroscopy (English Translation of Optika i Spektroskopiya)</i> , <b>2011</b> , 111, 808-813	0.7	1
48	Studying charge ordering and parameters of exchange interaction in Na <sub>x</sub> CoO <sub>2</sub> . <i>Physics of Metals and Metallography</i> , <b>2006</b> , 101, 255-260	1.2	1
47	Charge state of a transition metal impurity in III-V semiconductors. <i>Physics of the Solid State</i> , <b>2005</b> , 47, 1560	0.8	1
46	Evidence for canonical spin glass behaviour in polycrystalline Mn <sub>1.5</sub> Fe <sub>1.5</sub> Al Heusler alloy. <i>Journal of Magnetism and Magnetic Materials</i> , <b>2021</b> , 168752	2.8	1
45	Electronic Structure and Optical Properties of Heusler Alloy Mn <sub>1.5</sub> Fe <sub>1.5</sub> Al. <i>Journal of Experimental and Theoretical Physics</i> , <b>2021</b> , 133, 471-476	1	1
44	Electronic States and Optical Spectra of ErSn <sub>1.1</sub> Ge <sub>0.9</sub> and TmSn <sub>1.1</sub> Ge <sub>0.9</sub> Compounds. <i>Physics of Metals and Metallography</i> , <b>2020</b> , 121, 537-542	1.2	1
43	Effect of Electronic Correlations on the Electronic Structure, Magnetic and Optical Properties of the Ternary RCuGe Compounds with R = Tb, Dy, Ho, Er. <i>Materials</i> , <b>2020</b> , 13,	3.5	1
42	Electronic structure of the NpMT <sub>5</sub> (M = Fe, Co, Ni; T = Ga, In) series of neptunium compounds. <i>Physics of the Solid State</i> , <b>2016</b> , 58, 438-443	0.8	1
41	Revelation of spin glass behavior in Ru doped MnNiGe: experiment and theory. <i>Journal of Physics Condensed Matter</i> , <b>2019</b> , 31, 125803	1.8	1
40	Composition-Induced Magnetic Transition in GdMn <sub>1-x</sub> Ti <sub>x</sub> Si Intermetallic Compounds for x = 0. <i>Metals</i> , <b>2021</b> , 11, 1296	2.3	1



39	Field induced metamagnetism and large magnetic entropy change in RRhSi (R = Tb, Dy, Ho) rare earth intermetallics. <i>Journal of Alloys and Compounds</i> , <b>2021</b> , 888, 161493	5.7	1
38	Electronic Structure and Optical Properties of the Co <sub>2</sub> NiAl Heusler Alloy. <i>Physics of Metals and Metallography</i> , <b>2019</b> , 120, 729-732	1.2	0
37	The Structure of Electronic States in FeSb <sub>2</sub> According to Optical Spectroscopy and Band Calculations. <i>Physics of the Solid State</i> , <b>2019</b> , 61, 969-972	0.8	0
36	Electronic Structure and Spectral Characteristics of the Mn <sub>3</sub> Al Compound. <i>Physics of Metals and Metallography</i> , <b>2021</b> , 122, 954-959	1.2	0
35	Site-selective spin transition in LuCo <sub>3</sub> . <i>Journal of Physics and Chemistry of Solids</i> , <b>2022</b> , 163, 110552	3.9	0
34	Effect of Electronic Correlations on the Electronic Structures of the FeAlO <sub>3</sub> and FeSiO <sub>3</sub> Compounds. <i>Journal of Experimental and Theoretical Physics</i> , <b>2021</b> , 132, 548-555	1	0
33	Electronic properties and electronic structure of Co <sub>2</sub> YSi (Y = Ti, V, Cr, Mn, Fe) Heusler alloys. <i>IEEE Transactions on Magnetics</i> , <b>2021</b> , 1-1	2	0
32	Non-collinear antiferromagnetism to compensated ferrimagnetism in Ti(FeCo) (= 0, 0.5 and 1) alloys: experiment and theory. <i>Physical Chemistry Chemical Physics</i> , <b>2021</b> , 23, 5607-5614	3.6	0
31	Crystal and Electronic Structure of High-Temperature Superconductive Layered Cuprates in Temperature Interval 100-300 K. <i>Journal of Superconductivity and Novel Magnetism</i> , <b>2018</b> , 31, 1999-2002	1.5	0
30	Vacancy ordered phases of nonstoichiometric hafnium carbide from evolutionary crystal structure predictions. <i>Journal of Alloys and Compounds</i> , <b>2022</b> , 891, 162063	5.7	0
29	Electronic Structure and Optical Spectra of GdFeAl and GdFeSi Compounds. <i>Physics of the Solid State</i> , <b>2021</b> , 63, 866-871	0.8	0
28	Electronic structure and optical spectroscopy of the GdRhGe compound. <i>Optics and Spectroscopy (English Translation of Optika i Spektroskopiya)</i> , <b>2017</b> , 122, 574-579	0.7	
27	Spontaneous and induced magnetic phase transitions in Tb <sub>0.9</sub> Er <sub>0.1</sub> Ni <sub>5</sub> . <i>Journal of Magnetism and Magnetic Materials</i> , <b>2019</b> , 475, 593-601	2.8	
26	Magnetic Moments, Electronic Structure, and Optical Spectroscopy of Cobalt-Based Intermetallic Compounds YCo <sub>3</sub> , Y <sub>2</sub> Co <sub>7</sub> , and LaCo <sub>5</sub> . <i>Journal of Experimental and Theoretical Physics</i> , <b>2020</b> , 131, 600-606 <sup>1</sup>		
25	Electronic Structure and Optical Properties of the FeAl <sub>2</sub> Compound. <i>Physics of the Solid State</i> , <b>2020</b> , 62, 106-109	0.8	
24	Electronic Structure of the DyFe <sub>2</sub> Si <sub>2</sub> Compound: Energy Band Calculations and Optical Studies. <i>Physics of the Solid State</i> , <b>2020</b> , 62, 414-418	0.8	
23	Features of Electronic Structure of Intermetallic Compounds CeNi <sub>4</sub> M (M = Fe, Co, Ni, Cu). <i>Physics of the Solid State</i> , <b>2018</b> , 60, 466-469	0.8	
22	Electronic structure of Gd-doped MgO. <i>Journal of Experimental and Theoretical Physics</i> , <b>2016</b> , 122, 338-340		

- 21 The Influence of Copper Impurity on the Electronic Structure and Optical Properties of TmNi<sub>5</sub> Compound. *Optics and Spectroscopy (English Translation of Optika I Spektroskopiya)*, **2018**, 124, 784-788 0.7
- 20 Electronic Structure of GdCuGe Intermetallic Compound. *Physics of the Solid State*, **2018**, 60, 631-633 0.8
- 19 Electronic structure and optical properties of GdNi<sub>2</sub>Mnx compounds. *Low Temperature Physics*, **2018**, 44, 157-161 0.7
- 18 Electronic Structure, Optical, and Magnetic Properties of Mn<sub>100-x</sub>Gex (x = 20, 25, and 30) Alloys Near Tetragonal-Orthorhombic Structural Phase Transition. *Physica Status Solidi (B): Basic Research*, **2019**, 256, 1900155 1.3
- 17 Ab initio study on the rare-earth iron-pnictides RFeAsO (R = Pr, Nd, Sm, Gd) in the low-temperature Cmma phase. *Journal of Physics Condensed Matter*, **2014**, 26, 045501 1.8
- 16 Ab initio simulation of the electron structure and optical spectroscopy of ErRhGe compound. *Physics of the Solid State*, **2017**, 59, 1275-1278 0.8
- 15 Effect of copper and cobalt impurities on the electronic structure and optical spectra of the intermetallic compound PrNi<sub>5</sub>. *Physics of the Solid State*, **2014**, 56, 1933-1938 0.8
- 14 Role of Electronic Band Structure and Lattice Parameters in Magnetism of the R<sub>2</sub>(Fe,M)<sub>17</sub>Si<sub>3</sub> Al Compounds. *Solid State Phenomena*, **2009**, 152-153, 41-44 0.4
- 13 The origin of an elastic line in the L<sub>3</sub> x-ray emission spectrum of metallic manganese. *Physics of the Solid State*, **2006**, 48, 420-426 0.8
- 12 Specific features of the behavior of the optical properties of TbNi<sub>5</sub> & Cu<sub>x</sub> intermetallic compounds **2010**, 104, 360
- 11 Electronic and Optical Properties of RCuGe Compounds (R = Dy, Ho). *Bulletin of the Russian Academy of Sciences: Physics*, **2020**, 84, 1152-1155 0.4
- 10 Evolution of Electronic Structure of GdT<sub>100-x</sub>MnxFe<sub>0.95-x</sub>Si<sub>1</sub> Compounds According to Band Calculations and Optical Investigations. *Physics of Metals and Metallography*, **2021**, 122, 472-477 1.2
- 9 Magnetism of 3d and 4d doped MnT<sub>2</sub>Ge (T = Fe, Co, Ru and Rh): bulk magnetization and ab initio calculations. *Journal of Physics Condensed Matter*, **2019**, 31, 495804 1.8
- 8 Ab initio computational study of the electronic and magnetic properties of the HoNi<sub>2</sub>Z compounds accounting for electronic correlations. *Journal of Physics: Conference Series*, **2021**, 1740, 012032 0.3
- 7 A Role of the 3d Electron Subsystem in the Evolution of Band Structure and Magnetic and Optical Properties of ErNi<sub>5</sub> & Cox Compounds (x = 0-4). *Physics of the Solid State*, **2018**, 60, 2363-2369 0.8
- 6 The Structure of Electronic States and Optical Properties of Cr<sub>80</sub>Al<sub>20</sub> Compound. *Optics and Spectroscopy (English Translation of Optika I Spektroskopiya)*, **2018**, 125, 195-198 0.7
- 5 Electronic Structure and Exchange Interactions in RNi<sub>4</sub>Co (R = Eu, Yb) Compounds. *Physics of the Solid State*, **2018**, 60, 1682-1685 0.8
- 4 Optical Properties of Heusler Alloy Mn<sub>2</sub>FeAl with the  $\bar{1}11$ Mn Structure. *Physics of Metals and Metallography*, **2021**, 122, 737-741 1.2

- 3 Magnetic properties and electronic structure of Mn-Al alloys in the  $\bar{\Gamma}$ Mn structure. *Journal of Magnetism and Magnetic Materials*, **2021**, 168600 2.8
- 2 Induced by the pressure and the spin fluctuations the phase transitions in chiral itinerant ferromagnetics (for example MnSi). *Journal of Magnetism and Magnetic Materials*, **2021**, 539, 168282 2.8
- 1 Spectral characteristics and electronic structure of semimetallic ScSb and YSb. *Optical Materials*, **2022**, 129, 112466 3.3