

# Polina A Agzamova

## List of Publications by Year in descending order

Source: <https://exaly.com/author-pdf/1507260/publications.pdf>

Version: 2024-02-01

14  
papers

39  
citations

2682572

2  
h-index

1872680

6  
g-index

14  
all docs

14  
docs citations

14  
times ranked

39  
citing authors

#	ARTICLE	IF	CITATIONS
1	Structure and the electronic and magnetic properties of LaTiO <sub>3</sub> . Physics of the Solid State, 2008, 50, 1795-1798.	0.6	21
2	Elastic properties of rare earth pyrochlores R <sub>2</sub> Ti <sub>2</sub> O <sub>7</sub> (R = Gd, Tb, Dy, Ho, Er, Tm, Yb, Lu): Ab initio calculations. Optical Materials, 2017, 72, 565-570.	3.6	8
3	Magnetic hyperfine interactions at the <sup>51</sup> V nucleus in the pyrochlore Lu <sub>2</sub> V <sub>2</sub> O <sub>7</sub> . Low Temperature Physics, 2015, 41, 34-38.	0.6	2
4	Structural stability of CuAl <sub>2</sub> O <sub>4</sub> under pressure. Journal of Physics Condensed Matter, 2021, 33, 035403.	1.8	2
5	Electronic states in ferromagnetic Cr <sub>2</sub> studied by $\mu$ SR. Physical Review B, 2021, 104, ..	3.2	2
6	Hyperfine Interaction in Charge-Ordered Manganites. Physics of the Solid State, 2005, 47, 1523.	0.6	1
7	Hyperfine interactions at the lanthanum ion nucleus in LaMnO <sub>3</sub> . Low Temperature Physics, 2007, 33, 226-228.	0.6	1
8	Hyperfine interactions in titanates: Study of orbital ordering and local magnetic properties. Journal of Experimental and Theoretical Physics, 2013, 116, 828-833.	0.9	1
9	Phonon Spectrum of Eu <sub>2</sub> Sn <sub>2</sub> O <sub>7</sub> : Ab Initio Calculation. Optics and Spectroscopy (English Translation of) Tj ETQq1 1,0,784314 rgBT /O	1,0,784314	1
10	Specific features of magnetic and electrical hyperfine interactions in titanates according to ab initio calculations. Physics of Metals and Metallography, 2014, 115, 1194-1199.	1.0	0
11	Hyperfine Interactions on Magnetic Ions Nuclei in Titanates. Solid State Phenomena, 0, 215, 109-112.	0.3	0
12	Magnetic Hyperfine Fields in Lu <sub>2</sub> V <sub>2</sub> O <sub>7</sub> : A Model Approach. Journal of Low Temperature Physics, 2016, 185, 544-550.	1.4	0
13	Structure and Lattice Dynamics of MF <sub>2</sub> (M = Ca, Sr, Ba, Pb) in Cubic and Orthorhombic Phases: The ab initio Calculations. Physics of the Solid State, 2019, 61, 11-22.	0.6	0
14	Structure and Lattice Dynamics of R <sub>2</sub> Sn <sub>2</sub> O <sub>7</sub> and R <sub>2</sub> Zr <sub>2</sub> O <sub>7</sub> (R =) Tj ETQq0 0,0 rgBT /O	0,0	0