

# SÃ¼leyman Ãabuk

## List of Publications by Year in descending order

Source: <https://exaly.com/author-pdf/11613361/publications.pdf>

Version: 2024-02-01

19

papers

196

citations

1163117

8

h-index

1058476

14

g-index

19

all docs

19

docs citations

19

times ranked

260

citing authors

#	ARTICLE	IF	CITATIONS
1	Urbach rule and optical properties of the LiNbO <sub>3</sub> and LiTaO <sub>3</sub> . Journal of Optics, 1999, 1, 424-427.	1.5	28
2	Structural, electronic, elastic and vibrational properties of BiAlO <sub>3</sub> : A first principles study. Journal of Alloys and Compounds, 2013, 574, 520-525.	5.5	24
3	First-principles study of the structural, elastic and electronic properties of SbXI (X=S, Se, Te) crystals. Journal of Molecular Modeling, 2018, 24, 66.	1.8	23
4	<i>Ab initio</i>study of the structural, electronic and optical properties of NaTaO <sub>3</sub> . Philosophical Magazine, 2010, 90, 2965-2976.	1.6	18
5	Ab initio volume-dependent elastic and lattice dynamics properties of KTaO <sub>3</sub> . Physica Status Solidi (B): Basic Research, 2010, 247, 93-97.	1.5	17
6	First-principles study of structural, elastic, electronic and vibrational properties of BiCoO <sub>3</sub> . Solid State Sciences, 2014, 34, 1-7.	3.2	16
7	The nonlinear optical susceptibility and electro-optic tensor of ferroelectrics: first-principle study. Open Physics, 2012, 10, .	1.7	15
8	First-principles studies of the electronic structure and optical properties of AgBO <sub>3</sub> (B=Nb,Ta) in the paraelectric phase. Open Physics, 2008, 6, .	1.7	12
9	Mechanical, electronic and optical properties of SeZnO <sub>3</sub> : a GGA+U study. Philosophical Magazine, 2020, 100, 601-618.	1.6	8
10	Ab initio study of the lattice dynamical and thermodynamic properties of SbXI (X= S, Se, Te) compounds. Computational Condensed Matter, 2018, 16, e00320.	2.1	7
11	Investigation of structural and mechanical properties of rutile SnO <sub>2</sub> . Materials Research Express, 2019, 6, 085069.	1.6	7
12	Ab initio study of the linear and nonlinear optical responses in BiAlO <sub>3</sub> . Philosophical Magazine, 2016, 96, 190-207.	1.6	6
13	DFT - based study of electronic structures and mechanical properties of LiTaO <sub>3</sub> : ferroelectric and paraelectric phases. Philosophical Magazine, 2017, 97, 2469-2483.	1.6	6
14	FIRST-PRINCIPLES STUDY OF THE ELECTRONIC, LINEAR, AND NONLINEAR OPTICAL PROPERTIES OF <font>Li(Nb, Ta)O</font><sub>3</sub>. International Journal of Modern Physics B, 2010, 24, 6277-6290.	2.0	5
15	Argand Diagram and Oscillation Description of Electron State in Ferroelastic Crystals. Ferroelectrics, 2004, 307, 19-23.	0.6	1
16	First-principles study of electronic structure and optical properties of Sr(Ti,Zr)O <sub>3</sub> . Open Physics, 2013, 11, .	1.7	1
17	Magnetic, electronic and mechanical properties of SeXO <sub>3</sub> (X=Mn, Ni) with the LSDA+U framework. Journal of Alloys and Compounds, 2021, 850, 156674.	5.5	1
18	Structure and electronic properties of LnScO <sub>3</sub> compounds: A GGA+U calculation. Computational Materials Science, 2022, 208, 111350.	3.0	1

# ARTICLE

IF CITATIONS

19	Electret States in Some Electrooptical Materials. Ferroelectrics, 2003, 296, 117-125.	0.6	0
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