

# Mehmet Topsakal

## List of Publications by Year in descending order

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Version: 2024-02-01

38

papers

9,096

citations

218677

26

h-index

361022

35

g-index

38

all docs

38

docs citations

38

times ranked

8808

citing authors

#	ARTICLE	IF	CITATIONS
1	Two- and One-Dimensional Honeycomb Structures of Silicon and Germanium. <i>Physical Review Letters</i> , 2009, 102, 236804.	7.8	2,837
2	Monolayer honeycomb structures of group-IV elements and III-V binary compounds: First-principles calculations. <i>Physical Review B</i> , 2009, 80, .	3.2	1,769
3	First-principles study of two- and one-dimensional honeycomb structures of boron nitride. <i>Physical Review B</i> , 2009, 79, .	3.2	580
4	Electronic and magnetic properties of $\text{Mn}_{3\text{Mn}}$ atom adsorbed graphene and graphene nanoribbons. <i>Physical Review B</i> , 2008, 77, .	3.2	452
5	First-principles study of defects and adatoms in silicon carbide honeycomb structures. <i>Physical Review B</i> , 2010, 81, .	3.2	344
6	The response of mechanical and electronic properties of graphane to the elastic strain. <i>Applied Physics Letters</i> , 2010, 96, .	3.3	344
7	First-principles study of zinc oxide honeycomb structures. <i>Physical Review B</i> , 2009, 80, .	3.2	298
8	A Comparative Study of Lattice Dynamics of Three- and Two-Dimensional MoS <sub>2</sub> . <i>Journal of Physical Chemistry C</i> , 2011, 115, 16354-16361.	3.1	298
9	Structures of fluorinated graphene and their signatures. <i>Physical Review B</i> , 2011, 83, .	3.2	254
10	Elastic and plastic deformation of graphene, silicene, and boron nitride honeycomb nanoribbons under uniaxial tension: A first-principles density-functional theory study. <i>Physical Review B</i> , 2010, 81, .	3.2	219
11	Kinetic pathways of ionic transport in fast-charging lithium titanate. <i>Science</i> , 2020, 367, 1030-1034.	12.6	197
12	Graphene coatings: An efficient protection from oxidation. <i>Physical Review B</i> , 2012, 85, .	3.2	178
13	Current-voltage $\text{I}_{\text{V}}$ of armchair graphene nanoribbons under uniaxial strain. <i>Physical Review B</i> , 2010, 81, .	3.2	170
14	Superlattice structures of graphene-based armchair nanoribbons. <i>Physical Review B</i> , 2008, 78, .	3.2	148
15	Armchair nanoribbons of silicon and germanium honeycomb structures. <i>Physical Review B</i> , 2010, 81, .	3.2	137
16	First-principles approach to monitoring the band gap and magnetic state of a graphene nanoribbon via its vacancies. <i>Physical Review B</i> , 2008, 78, .	3.2	120
17	Frictional Figures of Merit for Single Layered Nanostructures. <i>Physical Review Letters</i> , 2012, 108, 126103.	7.8	110
18	Long-range interactions in carbon atomic chains. <i>Physical Review B</i> , 2010, 82, .	3.2	86

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19	Spin confinement in the superlattices of graphene ribbons. <i>Applied Physics Letters</i> , 2008, 92, .	3.3	79
20	Atomic and electronic structure of exfoliated black phosphorus. <i>Journal of Vacuum Science and Technology A: Vacuum, Surfaces and Films</i> , 2015, 33, .	2.1	73
21	Classification of local chemical environments from x-ray absorption spectra using supervised machine learning. <i>Physical Review Materials</i> , 2019, 3, .	2.4	66
22	Bandgap Restructuring of the Layered Semiconductor Gallium Telluride in Air. <i>Advanced Materials</i> , 2016, 28, 6465-6470.	21.0	58
23	Multi-Stage Structural Transformations in Zero-Strain Lithium Titanate Unveiled by <i>&lt; i&gt;in Situ&lt;/i&gt;</i> X-ray Absorption Fingerprints. <i>Journal of the American Chemical Society</i> , 2017, 139, 16591-16603.	13.7	57
24	Effects of Charging and Electric Field on Graphene Oxide. <i>Journal of Physical Chemistry C</i> , 2013, 117, 5943-5952.	3.1	47
25	Domain formation on oxidized graphene. <i>Physical Review B</i> , 2012, 86, .	3.2	40
26	Effects of static charging and exfoliation of layered crystals. <i>Physical Review B</i> , 2012, 85, .	3.2	35
27	Static charging of graphene and graphite slabs. <i>Applied Physics Letters</i> , 2011, 98, .	3.3	23
28	<i>&lt; i&gt;Operando&lt;/i&gt;</i> structural and chemical evolutions of TiS <sub>2</sub> in Na-ion batteries. <i>Journal of Materials Chemistry A</i> , 2020, 8, 12339-12350.	10.3	23
29	A New Line Defect in NdTiO <sub>3</sub> Perovskite. <i>Nano Letters</i> , 2016, 16, 6816-6822.	9.1	18
30	Identification of dopant site and its effect on electrochemical activity in Mn-doped lithium titanate. <i>Physical Review Materials</i> , 2018, 2, .	2.4	17
31	Nonresonant valence-to-core x-ray emission spectroscopy of niobium. <i>Physical Review B</i> , 2018, 97, .	3.2	11
32	Electronic structure of BaSnO <sub>3</sub> investigated by high-energy-resolution electron energy-loss spectroscopy and <i>&lt; i&gt;ab initio&lt;/i&gt;</i> calculations. <i>Journal of Vacuum Science and Technology A: Vacuum, Surfaces and Films</i> , 2018, 36, .	2.1	11
33	Metallic line defect in wide-bandgap transparent perovskite BaSnO <sub>3</sub> . <i>Science Advances</i> , 2021, 7, .	10.3	11
34	Probing ferroelectricity by x-ray absorption spectroscopy in molecular crystals. <i>Physical Review Materials</i> , 2020, 4, .	2.4	4
35	Decomposition of <i>&lt; i&gt;and&lt;/i&gt;</i> <i>xml�:ml="http://www.w3.org/1998/Math/MathML"&gt;&lt;mml:mrow&gt;&lt;mml:mi&gt;L&lt;/mml:mi&gt;&lt;mml:msub&gt;&lt;mml:mi&gt;a&lt;/mml:mi&gt;&lt;mml:mrow&gt;&lt;mml:mi&gt;x&lt;/mml:mi&gt;&lt;mml:mo&gt;^&lt;/mml:mo&gt;&lt;mml:mi&gt;2&lt;/mml:mi&gt;&lt;mml:mo&gt;*&lt;/mml:mo&gt;&lt;mml:mi&gt;b&lt;/mml:mi&gt;&lt;/mml:mrow&gt;&lt;/mml:msub&gt;&lt;mml:mi&gt;Cu&lt;/mml:mi&gt;&lt;mml:msub&gt;&lt;mml:mi&gt;x&lt;/mml:mi&gt;&lt;/mml:msub&gt;&lt;mml:mi&gt;2&lt;/mml:mi&gt;&lt;mml:mo&gt;*&lt;/mml:mo&gt;&lt;mml:mi&gt;y&lt;/mml:mi&gt;&lt;/mml:mrow&gt;&lt;/i&gt;</i>	2.4	2
36	Machine-Learning Assisted Structure Determination of Metallic Nanoparticles: A Benchmark. , 2020, , 127-140.	1	

#	ARTICLE	IF	CITATIONS
37	Electronic Structure of New Line Defect in Strained NdTiCb on SrTiO <sub>3</sub> . Microscopy and Microanalysis, 2015, 21, 2073-2074.	0.4	0
38	Probing the Electronic Structure of BaSnO <sub>3</sub> by EELS Analysis and ab initio Calculations. Microscopy and Microanalysis, 2017, 23, 1602-1603.	0.4	0