

# Alexander G Kvashnin

## List of Publications by Year in descending order

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69  
papers

4,845  
citations

201385

27  
h-index

95083

68  
g-index

69  
all docs

69  
docs citations

69  
times ranked

6536  
citing authors

#	ARTICLE	IF	CITATIONS
1	Diamane quasicrystals. <i>Applied Surface Science</i> , 2022, 572, 151362.	3.1	16
2	Nanohardness from First Principles with Active Learning on Atomic Environments. <i>Journal of Chemical Theory and Computation</i> , 2022, 18, 1109-1121.	2.3	10
3	Computational Design of Gas Sensors Based on V3S4 Monolayer. <i>Nanomaterials</i> , 2022, 12, 774.	1.9	7
4	Sr <sup>2+</sup> -Doped Superionic Hydrogen Glass: Synthesis and Properties of SrH <sub>22</sub> . <i>Advanced Materials</i> , 2022, 34, e2200924.	11.1	10
5	Efficient Synthesis of WB <sub>5</sub> and WB <sub>2</sub> Powders with Selectivity for WB <sub>5</sub> Content. <i>Inorganic Chemistry</i> , 2022, 61, 6773-6784.	1.9	3
6	GIPAW Pseudopotentials of d Elements for Solid-State NMR. <i>Materials</i> , 2022, 15, 3347.	1.3	4
7	Map of Two-Dimensional Tungsten Chalcogenide Compounds (W <sub>2</sub> S, W <sub>2</sub> Se, W <sub>2</sub> Te) Based on USPEX Evolutionary Search. <i>JETP Letters</i> , 2022, 115, 292-296.	0.4	3
8	Large-Scale Synthesis and Applications of Hafnium-Tantalum Carbides. <i>Advanced Functional Materials</i> , 2022, 32, .	7.8	8
9	Crystal Structure Evolution of Fluorine under High Pressure. <i>Journal of Physical Chemistry C</i> , 2022, 126, 11358-11364.	1.5	5
10	Novel Strongly Correlated Europium Superhydrides. <i>Journal of Physical Chemistry Letters</i> , 2021, 12, 32-40.	2.1	33
11	Synthesis of molecular metallic barium superhydride: pseudocubic BaH <sub>12</sub> . <i>Nature Communications</i> , 2021, 12, 273.	5.8	66
12	Anomalous High-Temperature Superconductivity in YH <sub>6</sub> . <i>Advanced Materials</i> , 2021, 33, e2006832.	11.1	196
13	Computational Modeling of 2D Materials under High Pressure and Their Chemical Bonding: Silicene as Possible Field-Effect Transistor. <i>ACS Nano</i> , 2021, 15, 6861-6871.	7.3	18
14	Superconductivity at 253 K in lanthanum-yttrium ternary hydrides. <i>Materials Today</i> , 2021, 48, 18-28.	8.3	119
15	Novel two-dimensional boron oxynitride predicted using the USPEX evolutionary algorithm. <i>Physical Chemistry Chemical Physics</i> , 2021, 23, 26178-26184.	1.3	4
16	Superconductivity at 161 K in thorium hydride ThH <sub>10</sub> : Synthesis and properties. <i>Materials Today</i> , 2020, 33, 36-44.	8.3	187
17	Computational Search for New W-Mo-B Compounds. <i>Chemistry of Materials</i> , 2020, 32, 7028-7035.	3.2	22
18	Environmentally Friendly Method of Silicon Recycling: Synthesis of Silica Nanoparticles in an Aqueous Solution. <i>ACS Sustainable Chemistry and Engineering</i> , 2020, 8, 14006-14012.	3.2	9

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19	The Volumetric Source Function: Looking Inside van der Waals Interactions. <i>Scientific Reports</i> , 2020, 10, 7816.	1.6	10
20	Phase Transitions in Tungsten Monoborides. <i>JETP Letters</i> , 2020, 111, 343-349.	0.4	3
21	Structure, Stability, and Mechanical Properties of Boron-Rich Mo $\alpha$ B Phases: A Computational Study. <i>Journal of Physical Chemistry Letters</i> , 2020, 11, 2393-2401.	2.1	30
22	WB 5 $\alpha$ x : Synthesis, Properties, and Crystal Structure $\alpha$ New Insights into the Long $\alpha$ Debated Compound. <i>Advanced Science</i> , 2020, 7, 2000775.	5.6	17
23	Young $\alpha$ TM's Modulus and Tensile Strength of Ti <sub>3</sub> C <sub>2</sub> MXene Nanosheets As Revealed by <i>In Situ</i> TEM Probing, AFM Nanomechanical Mapping, and Theoretical Calculations. <i>Nano Letters</i> , 2020, 20, 5900-5908.	4.5	88
24	Superconductivity of $\text{LaH}_{10}$ and $\text{LaH}_{16}$ polyhydrides. <i>Physical Review B</i> , 2020, 101, .	1.1	62
25	High-Pressure Synthesis of Magnetic Neodymium Polyhydrides. <i>Journal of the American Chemical Society</i> , 2020, 142, 2803-2811.	6.6	59
26	Exotic Two-Dimensional Structure: The First Case of Hexagonal NaCl. <i>Journal of Physical Chemistry Letters</i> , 2020, 11, 3821-3827.	2.1	38
27	On Distribution of Superconductivity in Metal Hydrides. <i>Current Opinion in Solid State and Materials Science</i> , 2020, 24, 100808.	5.6	104
28	Mechanical Engineering Effect in Electronic and Optical Properties of Graphene Nanomeshes. <i>ACS Applied Materials &amp; Interfaces</i> , 2020, 12, 55189-55194.	4.0	9
29	Computational discovery of hard and superhard materials. <i>Journal of Applied Physics</i> , 2019, 126, .	1.1	46
30	Nonstoichiometric Phases of Two-Dimensional Transition-Metal Dichalcogenides: From Chalcogen Vacancies to Pure Metal Membranes. <i>Journal of Physical Chemistry Letters</i> , 2019, 10, 6492-6498.	2.1	15
31	Novel Unexpected Reconstructions of (100) and (111) Surfaces of NaCl: Theoretical Prediction. <i>Scientific Reports</i> , 2019, 9, 14267.	1.6	21
32	Stability and magnetism of FeN high-pressure phases. <i>Physical Chemistry Chemical Physics</i> , 2019, 21, 5262-5273.	1.3	12
33	Stable and hard hafnium borides: A first-principles study. <i>Journal of Applied Physics</i> , 2019, 125, .	1.1	13
34	Two-Dimensional CuO Inside the Supportive Bilayer Graphene Matrix. <i>Journal of Physical Chemistry C</i> , 2019, 123, 17459-17465.	1.5	12
35	Iron Superhydrides FeH <sub>5</sub> and FeH <sub>6</sub> : Stability, Electronic Properties, and Superconductivity. <i>Journal of Physical Chemistry C</i> , 2018, 122, 4731-4736.	1.5	48
36	Actinium Hydrides AcH <sub>10</sub> , AcH <sub>12</sub> , and AcH <sub>16</sub> as High-Temperature Conventional Superconductors. <i>Journal of Physical Chemistry Letters</i> , 2018, 9, 1920-1926.	2.1	100

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37	Layered heterostructures based on graphene, hexagonal zinc oxide and molybdenum disulfide: Modeling of geometry and electronic properties. <i>Computational Materials Science</i> , 2018, 142, 32-37.	1.4	7
38	High-Temperature Superconductivity in a $\text{Th}\epsilon\text{-H}$ System under Pressure Conditions. <i>ACS Applied Materials &amp; Interfaces</i> , 2018, 10, 43809-43816.	4.0	95
39	Uranium polyhydrides at moderate pressures: Prediction, synthesis, and expected superconductivity. <i>Science Advances</i> , 2018, 4, eaat9776.	4.7	82
40	New Tungsten Borides, Their Stability and Outstanding Mechanical Properties. <i>Journal of Physical Chemistry Letters</i> , 2018, 9, 3470-3477.	2.1	61
41	Computational Materials Discovery: Dream or Reality?. , 2018, , 1-14.		3
42	Computational Search for Novel Hard Chromium-Based Materials. <i>Journal of Physical Chemistry Letters</i> , 2017, 8, 755-764.	2.1	62
43	Fullerite-based nanocomposites with ultrahigh stiffness. Theoretical investigation. <i>Carbon</i> , 2017, 115, 546-549.	5.4	15
44	Novel hybrid C/BN two-dimensional heterostructures. <i>Nanotechnology</i> , 2017, 28, 085205.	1.3	8
45	Estimation of graphene surface stability against the adsorption of environmental and technological chemical agents. <i>Physica Status Solidi (B): Basic Research</i> , 2017, 254, 1600702.	0.7	5
46	New allotropic forms of carbon based on $\text{C}_{60}$ and $\text{C}_{70}$ fullerenes with specific mechanical characteristics. <i>JETP Letters</i> , 2017, 105, 419-425.	0.4	3
47	Stable reconstruction of the (110) surface and its role in pseudocapacitance of rutile-like $\text{RuO}_2$ . <i>Scientific Reports</i> , 2017, 7, 10357.	1.6	30
48	Features of Electronic, Mechanical, and Electromechanical Properties of Fluorinated Diamond Films of Nanometer Thickness. <i>Journal of Physical Chemistry C</i> , 2017, 121, 28484-28489.	1.5	29
49	The possible formation of a magnetic $\text{FeS}_2$ phase in the two-dimensional $\text{MoS}_2$ matrix. <i>Physical Chemistry Chemical Physics</i> , 2016, 18, 26956-26959.	1.3	1
50	Heterostructures based on graphene and $\text{MoS}_2$ layers decorated by $\text{C}_{60}$ fullerenes. <i>Nanotechnology</i> , 2016, 27, 365201.	1.3	11
51	Ionic Graphitization of Ultrathin Films of Ionic Compounds. <i>Journal of Physical Chemistry Letters</i> , 2016, 7, 2659-2663.	2.1	9
52	Toward the Ultra-incompressible Carbon Materials. Computational Simulation and Experimental Observation. <i>Journal of Physical Chemistry Letters</i> , 2015, 6, 2147-2152.	2.1	16
53	Flexoelectricity in Carbon Nanostructures: Nanotubes, Fullerenes, and Nanocones. <i>Journal of Physical Chemistry Letters</i> , 2015, 6, 2740-2744.	2.1	68
54	Hydrogen adsorption study. Formation of quantum dots on graphene nanoribbons within tight-binding approach. <i>Nanotechnology</i> , 2015, 26, 175704.	1.3	1

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55	Transport investigation of branched graphene nanoflakes. <i>Nanotechnology</i> , 2015, 26, 385705.	1.3	7
56	Radiation-Induced Nucleation of Diamond from Amorphous Carbon: Effect of Hydrogen. <i>Journal of Physical Chemistry Letters</i> , 2014, 5, 1924-1928.	2.1	26
57	Phase Diagram of Quasi-Two-Dimensional Carbon, From Graphene to Diamond. <i>Nano Letters</i> , 2014, 14, 676-681.	4.5	154
58	Lonsdaleite Films with Nanometer Thickness. <i>Journal of Physical Chemistry Letters</i> , 2014, 5, 541-548.	2.1	56
59	Spontaneous Graphitization of Ultrathin Cubic Structures: A Computational Study. <i>Nano Letters</i> , 2014, 14, 7126-7130.	4.5	31
60	Graphitic Phase of NaCl. Bulk Properties and Nanoscale Stability. <i>Journal of Physical Chemistry Letters</i> , 2014, 5, 4014-4019.	2.1	16
61	Investigation of new superhard carbon allotropes with promising electronic properties. <i>Journal of Applied Physics</i> , 2013, 114, 183708.	1.1	10
62	Strong Influence of Graphane Island Configurations on the Electronic Properties of a Mixed Graphene/Graphane Superlattice. <i>Journal of Physical Chemistry C</i> , 2012, 116, 20035-20039.	1.5	13
63	Influence of Size Effect on the Electronic and Elastic Properties of Diamond Films with Nanometer Thickness. <i>Journal of Physical Chemistry C</i> , 2011, 115, 132-136.	1.5	82
64	Theoretical Study of Elastic Properties of SiC Nanowires of Different Shapes. <i>Journal of Nanoscience and Nanotechnology</i> , 2010, 10, 4992-4997.	0.9	3
65	Large Scale Growth and Characterization of Atomic Hexagonal Boron Nitride Layers. <i>Nano Letters</i> , 2010, 10, 3209-3215.	4.5	2,317
66	Theoretical Study of Atomic Structure and Elastic Properties of Branched Silicon Nanowires. <i>ACS Nano</i> , 2010, 4, 2784-2790.	7.3	4
67	The Theoretical Study of Mechanical Properties of Graphene Membranes. <i>Fullerenes Nanotubes and Carbon Nanostructures</i> , 2010, 18, 497-500.	1.0	17
68	Diamond-like C <sub>2</sub> H nanolayer, diamane: Simulation of the structure and properties. <i>JETP Letters</i> , 2009, 90, 134-138.	0.4	169
69	Density functional study of $\gamma$ thin silicon nanowires. <i>Physical Review B</i> , 2008, 77, .		