

# Martin Dahlqvist

## List of Publications by Year in descending order

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66  
papers

3,876  
citations

156536

32  
h-index

139680

61  
g-index

67  
all docs

67  
docs citations

67  
times ranked

2571  
citing authors

#	ARTICLE	IF	CITATIONS
1	Correlation strength, orbital-selective incoherence, and local moments formation in the magnetic MAX-phase Physical Review B, 2022, 105, .	1.1	3
2	Chemical order or disorder – a theoretical stability expose for expanding the compositional space of quaternary metal borides. Materials Advances, 2022, 3, 2908-2917.	2.6	9
3	Investigation of out-of-plane ordered Ti <sub>4</sub> MoSiB <sub>2</sub> from first principles. Journal of Physics Condensed Matter, 2022, , .	0.7	2
4	Theoretical predictions of phase stability for orthorhombic and hexagonal ternary MAB phases. Physical Chemistry Chemical Physics, 2022, 24, 11249-11258.	1.3	18
5	First-principles study on the superconductivity of doped zirconium diborides. Physical Review Materials, 2022, 6, .	0.9	2
6	On the nature of planar defects in transition metal diboride line compounds. Materialia, 2022, 24, 101478.	1.3	4
7	Synthesis, characterization and first principle modelling of the MAB phase solid solutions: (Mn <sub>1-x</sub> Cr <sub>x</sub> ) <sub>2</sub> AlB <sub>2</sub> and (Mn <sub>1-x</sub> Cr <sub>x</sub> ) <sub>3</sub> AlB <sub>4</sub> . Materials Research Letters, 2021, 9, 112-118.	4.1	17
8	Where is the unpaired transition metal in substoichiometric diboride line compounds?. Acta Materialia, 2021, 204, 116510.	3.8	21
9	In-plane ordered quaternary M <sub>4/3</sub> A <sub>2</sub> M <sub>2/3</sub> B <sub>2</sub> phases (i-MAB): electronic structure and mechanical properties from first-principles calculations. Journal of Physics Condensed Matter, 2021, 33, 255402.	0.7	4
10	Boridene: Two-dimensional Mo <sub>4/3</sub> B <sub>2-x</sub> with ordered metal vacancies obtained by chemical exfoliation. Science, 2021, 373, 801-805.	6.0	126
11	Out-of-Plane Ordered Laminate Borides and Their 2D Ti-Based Derivative from Chemical Exfoliation. Advanced Materials, 2021, 33, e2008361.	11.1	14
12	Predictions of attainable compositions of layered quaternary i-MAB phases and solid solution MAB phases. Nanoscale, 2021, 13, 18311-18321.	2.8	12
13	Phase Stability of Nanolaminated Epitaxial (Cr <sub>1-x</sub> Fe <sub>x</sub> ) <sub>2</sub> AlC MAX Phase Thin Films on MgO(111) and Al <sub>2</sub> O <sub>3</sub> (0001) for Use as Conductive Coatings. ACS Applied Nano Materials, 2021, 4, 13761-13770.	2.4	6
14	Predictive theoretical screening of phase stability for chemical order and disorder in quaternary 312 and 413 MAX phases. Nanoscale, 2020, 12, 785-794.	2.8	56
15	Ta-based 413 and 211 MAX phase solid solutions with Hf and Nb. Journal of the European Ceramic Society, 2020, 40, 1829-1838.	2.8	31
16	Theoretical Prediction and Experimental Verification of the Chemically Ordered Atomic-Laminate i-MAX Phases (Cr <sub>2/3</sub> Sc <sub>1/3</sub> ) <sub>2</sub> GaC and (Mn <sub>2/3</sub> Sc <sub>1/3</sub> ) <sub>2</sub> GaC. Crystal Growth and Design, 2020, 20, 55-61.	1.4	16
17	Theoretical Prediction and Synthesis of a Family of Atomic Laminate Metal Borides with In-Plane Chemical Ordering. Journal of the American Chemical Society, 2020, 142, 18583-18591.	6.6	55
18	Impact of strain, pressure, and electron correlation on magnetism and crystal structure of Mn <sub>2</sub> GaC from first-principles. Scientific Reports, 2020, 10, 11384.	1.6	13

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19	Single Crystal Growth and Structural Characterization of Theoretically Predicted Nanolaminates $M_2Al_2C_3$ , Where M = Sc and Er. <i>Crystal Growth and Design</i> , 2020, 20, 7640-7646.	1.4	3
20	Chemically stable new MAX phase $V_2SnC$ : a damage and radiation tolerant TBC material. <i>RSC Advances</i> , 2020, 10, 43783-43798.	1.7	34
21	Compatibility of $Zr_2AlC$ MAX phase-based ceramics with oxygen-poor, static liquid lead-bismuth eutectic. <i>Corrosion Science</i> , 2020, 171, 108704.	3.0	24
22	Possible monoclinic distortion of $Mo_2GaC$ under high pressure. <i>Journal of Applied Physics</i> , 2020, 127, 145103.	1.1	2
23	In- and Out-of-Plane Ordered MAX Phases and Their MXene Derivatives. , 2019, , 37-52.		9
24	Theoretical Analysis, Synthesis, and Characterization of 2D $W_{1.33}C$ (MXene) with Ordered Vacancies. <i>ACS Applied Nano Materials</i> , 2019, 2, 6209-6219.	2.4	37
25	Atomically Layered and Ordered Rare-Earth <i>MAX</i> Phases: A New Class of Magnetic Quaternary Compounds. <i>Chemistry of Materials</i> , 2019, 31, 2476-2485.	3.2	89
26	Transmorphic epitaxial growth of AlN nucleation layers on SiC substrates for high-breakdown thin GaN transistors. <i>Applied Physics Letters</i> , 2019, 115, .	1.5	25
27	A Tungsten-Based Nanolaminated Ternary Carbide: $(W,Ti)_4C_4$ . <i>Inorganic Chemistry</i> , 2019, 58, 1100-1106.	1.9	9
28	Materials synthesis, neutron powder diffraction, and first-principles calculations of $Mo_{1.33}Ti$ . <i>Physical Review Materials</i> , 2019, 3, .	1.9	9
29	W-Based Atomic Laminates and Their 2D Derivative $W_{1.33}C$ MXene with Vacancy Ordering. <i>Advanced Materials</i> , 2018, 30, e1706409.	11.1	240
30	Origin of Chemically Ordered Atomic Laminates ( <i>MAX</i> ): Expanding the Elemental Space by a Theoretical/Experimental Approach. <i>ACS Nano</i> , 2018, 12, 7761-7770.	7.3	99
31	Theoretical Prediction and Synthesis of $(Cr_{2/3}Zr_{1/3})_2AlC$ <i>MAX</i> Phase. <i>Inorganic Chemistry</i> , 2018, 57, 6237-6244.	1.9	59
32	Electronic structure, bonding characteristics, and mechanical properties in $(W_{2/3}Sc_{1/3})_2AlC$ and $(W_{2/3}Y_{1/3})_2AlC$ <i>MAX</i> phases from first-principles calculations. <i>Journal of Physics Condensed Matter</i> , 2018, 30, 305502.	0.7	9
33	Synthesis of $Ti_3AuC_2$ , $Ti_3Au_2C_2$ and $Ti_3IrC_2$ by noble metal substitution reaction in $Ti_3SiC_2$ for high-temperature-stable Ohmic contacts to SiC. <i>Nature Materials</i> , 2017, 16, 814-818.	13.3	142
34	Two-dimensional $Mo_{1.33}C$ MXene with divacancy ordering prepared from parent 3D laminate with in-plane chemical ordering. <i>Nature Communications</i> , 2017, 8, 14949.	5.8	525
35	Theoretical stability and materials synthesis of a chemically ordered MAX phase, $Mo_2ScAlC_2$ , and its two-dimensional derivate $Mo_2ScC_2$ MXene. <i>Acta Materialia</i> , 2017, 125, 476-480.	3.8	185
36	Dataset on the structure and thermodynamic and dynamic stability of $Mo_2ScAlC_2$ from experiments and first-principles calculations. <i>Data in Brief</i> , 2017, 10, 576-582.	0.5	3

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37	Prediction and synthesis of a family of atomic laminate phases with Kagomé-like and in-plane chemical ordering. <i>Science Advances</i> , 2017, 3, e1700642.	4.7	156
38	Ti <sub>2</sub> Au <sub>2</sub> C and Ti <sub>3</sub> Au <sub>2</sub> C <sub>2</sub> formed by solid state reaction of gold with Ti <sub>2</sub> AlC and Ti <sub>3</sub> AlC <sub>2</sub> . <i>Chemical Communications</i> , 2017, 53, 9554-9557.	2.2	53
39	Benefits of oxygen incorporation in atomic laminates. <i>Journal of Physics Condensed Matter</i> , 2016, 28, 135501.	0.7	7
40	Phase stability of the nanolaminates V <sub>2</sub> Ga <sub>2</sub> C and (Mo <sub>1-x</sub> V <sub>x</sub> ) <sub>2</sub> Ga <sub>2</sub> C from first-principles calculations. <i>Physical Chemistry Chemical Physics</i> , 2016, 18, 12682-12688.	1.3	10
41	Magnetic MAX phases from theory and experiments; a review. <i>Journal of Physics Condensed Matter</i> , 2016, 28, 433003.	0.7	84
42	Long-range antiferromagnetic order in epitaxial Mn <sub>2</sub> AlC films from neutron reflectometry. <i>Physical Review B</i> , 2016, 94, .	1.1	28
43	Magnetically driven anisotropic structural changes in the atomic laminate Mn <sub>2</sub> AlC. <i>Physical Review B</i> , 2016, 94, .	1.1	44
44	Magnetic exchange interactions and critical temperature of the nanolaminate Mn <sub>2</sub> AlC from first-principles supercell methods. <i>Physical Review B</i> , 2016, 93, .	1.1	12
45	Theoretical stability, thin film synthesis and transport properties of the Mo <sub>1-x</sub> N <sub>x</sub> +1 GaC <sub>1-x</sub> N <sub>x</sub> MAX phase. <i>Physica Status Solidi - Rapid Research Letters</i> , 2015, 9, 197-201.	1.2	28
46	Experimental and theoretical characterization of ordered MAX phases Mo <sub>2</sub> TiAlC <sub>2</sub> and Mo <sub>2</sub> Ti <sub>2</sub> AlC <sub>3</sub> . <i>Journal of Applied Physics</i> , 2015, 118, .	1.1	217
47	Influence of boron vacancies on phase stability, bonding and structure of MB <sub>2</sub> (M = Ti, Zr, Hf, V, Nb, Ta, Cr, Mo, W) with AlB <sub>2</sub> type structure. <i>Journal of Physics Condensed Matter</i> , 2015, 27, 435702.	1.1	33
48	Order and disorder in quaternary atomic laminates from first-principles calculations. <i>Physical Chemistry Chemical Physics</i> , 2015, 17, 31810-31821.	1.3	71
49	A critical evaluation of GGA+U modeling for atomic, electronic and magnetic structure of Cr <sub>2</sub> AlC, Cr <sub>2</sub> GaC and Cr <sub>2</sub> GeC. <i>Journal of Physics Condensed Matter</i> , 2015, 27, 095601.	0.7	23
50	Structural and chemical determination of the new nanolaminated carbide Mo <sub>2</sub> Ga <sub>2</sub> C from first principles and materials analysis. <i>Acta Materialia</i> , 2015, 99, 157-164.	3.8	75
51	Effect of Ti-Al cathode composition on plasma generation and plasma transport in direct current vacuum arc. <i>Journal of Applied Physics</i> , 2014, 115, .	1.1	45
52	First-principles calculations of the electronic, vibrational, and elastic properties of the magnetic laminate Mn <sub>2</sub> GaC. <i>Journal of Applied Physics</i> , 2014, 116, .	1.1	23
53	A Nanolaminated Magnetic Phase: Mn <sub>2</sub> GaC. <i>Materials Research Letters</i> , 2014, 2, 89-93.	4.1	128
54	Temperature dependent phase stability of nanolaminated ternaries from first-principles calculations. <i>Computational Materials Science</i> , 2014, 91, 251-257.	1.4	32

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55	Oxygen incorporation in TiAlC thin films studied by electron energy loss spectroscopy and ab initio calculations. Journal of Materials Science, 2013, 48, 3686-3691.	1.7	17
56	Synthesis and <i>ab initio</i> calculations of nanolaminated (Cr,Mn) $\text{AlC}$ compounds. Physical Review B, 2013, 87, .	1.1	93
57	Magnetic Self-Organized Atomic Laminate from First Principles and Thin Film Synthesis. Physical Review Letters, 2013, 110, 195502.	2.9	146
58	Phase stability of $\text{Cr}_{n+1}\text{GaC}_n$ MAX phases from first principles and $\text{Cr}_2\text{GaC}$ thin film synthesis using magnetron sputtering from elemental targets. Physica Status Solidi - Rapid Research Letters, 2013, 7, 971-974.	1.2	32
59	Characterization of plasma chemistry and ion energy in cathodic arc plasma from Ti-Si cathodes of different compositions. Journal of Applied Physics, 2013, 113, 163304.	1.1	22
60	Correlation between magnetic state and bulk modulus of $\text{Cr}_2\text{AlC}$ . Journal of Applied Physics, 2013, 113, .	1.1	50
61	Discovery of the Ternary Nanolaminated Compound $\text{Nb}_2\text{GeC}$ by a Systematic Theoretical-Experimental Approach. Physical Review Letters, 2012, 109, 035502.	2.9	73
62	Magnetic nanoscale laminates with tunable exchange coupling from first principles. Physical Review B, 2011, 84, .	1.1	75
63	Phase stability of $\text{Ti}_2\text{AlC}$ oxygen incorporation: A first-principles investigation. Physical Review B, 2010, 81, .	1.1	10
64	Oxygen incorporation in $\text{Ti}_2\text{AlC}$ : Tuning of anisotropic conductivity. Applied Physics Letters, 2010, 97, .	1.5	44
65	Stability trends of $\text{M}_2\text{AX}$ phases from first principles. Physical Review B, 2010, 81, .	1.1	10
66	Dynamics of Polymer Adsorption from Bulk Solution onto Planar Surfaces. Macromolecules, 2009, 42, 3641-3649.	2.2	39