Vladislav A Blatov

List of Publications by Year
in descending order

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216 papers	13,086 citations	44069 48 h-index	23533 111 g-index
226 all docs	226 docs citations	226 times ranked	7456 citing authors

#	Article	IF	Citations
1	Applied Topological Analysis of Crystal Structures with the Program Package ToposPro. Crystal Growth and Design, 2014, 14, 3576-3586.	3.0	2,448
2	Interpenetrating metalâ€“organic and inorganic 3D networks: a computer-aided systematic investigation. Part I. Analysis of the Cambridge structural database. CrystEngComm, 2004, 6, 377-395.	2.6	1,116
3	TOPOS3.2: a new version of the program package for multipurpose crystal-chemical analysis. Journal of Applied Crystallography, 2000, 33, 1193-1193.	4.5	964
4	Vertex-, face-, point-, SchlÃfli-, and Delaney-symbols in nets, polyhedra and tilings: recommended terminology. CrystEngComm, 2010, 12, 44-48.	2.6	694
5	Underlying nets in three-periodic coordination polymers: topology, taxonomy and prediction from a computer-aided analysis of the Cambridge Structural Database. CrystEngComm, 2011, 13, 3947.	2.6	626
6	Nanocluster analysis of intermetallic structures with the program package TOPOS. Structural Chemistry, 2012, 23, 955-963.	2.0	488
7	Interpenetrating metal-organic and inorganic 3D networks: a computer-aided systematic investigation. Part II [1]. Analysis of the Inorganic Crystal Structure Database (ICSD). Journal of Solid State Chemistry, 2005, 178, 2452-2474.	2.9	335
8	Entangled Two-Dimensional Coordination Networks: A General Survey. Chemical Reviews, 2014, 114, 7557-7580.	47.7	253
9	Voronoiâ€“dirichlet polyhedra in crystal chemistry: theory and applications. Crystallography Reviews, 2004, 10, 249-318.	1.5	236
10	Interpenetrated Three-Dimensional Networks of Hydrogen-Bonded Organic Species: A Systematic Analysis of the Cambridge Structural Database. Crystal Growth and Design, 2008, 8, 519-539.	3.0	232
11	High-nuclearity cobalt coordination clusters: Synthetic, topological and magnetic aspects. Coordination Chemistry Reviews, 2012, 256, 1246-1278.	18.8	204
12	Two metalâ€“organic frameworks with unique high-connected binodal network topologies: synthesis, structures, and catalytic properties. CrystEngComm, 2012, 14, 4210.	2.6	196
13	Three-periodic nets and tilings: natural tilings for nets. Acta Crystallographica Section A: Foundations and Advances, 2007, 63, 418-425.	0.3	188
14	Topological relations between three-periodic nets. II. Binodal nets. Acta Crystallographica Section A: Foundations and Advances, 2009, 65, 202-212.	0.3	172
15	Interpenetrated three-dimensional hydrogen-bonded networks from metalâ€“organic molecular and one- or two-dimensional polymeric motifs. CrystEngComm, 2008, 10, 1822.	2.6	160
16	The Aluminum-Ion Battery: A Sustainable and Seminal Concept?. Frontiers in Chemistry, 2019, 7, 268.	3.6	155
17	A luminescent zinc(<scp>ii</scp>) coordination polymer with unusual (3,4,4)-coordinated self-catenated 3D network for selective detection of nitroaromatics and ferric and chromate ions: a versatile luminescent sensor. Dalton Transactions, 2018, 47, 6189-6198.	3.3	147
18	Syntheses, Topological Structures, and Photoluminescences of Six New Zn(II) Coordination Polymers Based on Mixed Tripodal Imidazole Ligand and Varied Polycarboxylates. Crystal Growth and Design, 2013, 13, 1277-1289.	3.0	143

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19	Topological Motifs in Cyanometallates: From Building Units to Three-Periodic Frameworks. Chemical Reviews, 2015, 115, 12286-12319.	47.7	128
20	Crystal space analysis by means of Voronoiâ€“Dirichlet polyhedra. Acta Crystallographica Section A: Foundations and Advances, 1995, 51, 909-916.	0.3	123
21	Topology of 2-Periodic Coordination Networks: Toward Expert Systems in Crystal Design. Crystal Growth and Design, 2013, 13, 1655-1664.	3.0	119
22	Topological relations between three-dimensional periodic nets. I. Uninodal nets. Acta Crystallographica Section A: Foundations and Advances, 2007, 63, 329-343.	0.3	109
23	The Zeolite Conundrum: Why Are There so Many Hypothetical Zeolites and so Few Observed? A Possible Answer from the Zeolite-Type Frameworks Perceived As Packings of Tiles. Chemistry of Materials, 2013, 25, 412-424.	6.7	90
24	Predicting crystal growth via a unified kinetic three-dimensional partition model. Nature, 2017, 544, 456-459.	27.8	88
25	Structures and Properties of Spherical 90â€Vertex Fullereneâ€Like Nanoballs. Chemistry - A European Journal, 2010, 16, 2092-2107.	3.3	87
26	An unprecedented â€œstronglyâ€•self-catenated MOF containing inclined catenated honeycomb-like units. Dalton Transactions, 2016, 45, 2426-2429.	3.3	87
27	Crystal chemistry of zirconosilicates and their analogs: topological classification of MT frameworks and suprapolyhedral invariants. Acta Crystallographica Section B: Structural Science, 2002, 58, 198-218.	1.8	83
28	Topology of molecular packings in organic crystals. Acta Crystallographica Section B: Structural Science, 2000, 56, 1035-1045.	1.8	82
29	Natural Tilings for Zeolite-Type Frameworks. Journal of Physical Chemistry C, 2010, 114, 10160-10170.	3.1	82
30	A method for topological analysis of high nuclearity coordination clusters and its application to Mn coordination compounds. Dalton Transactions, 2012, 41, 4634.	3.3	80
31	Migration maps of Li+ cations in oxygen-containing compounds. Solid State Ionics, 2008, 179, 2248-2254.	2.7	79
32	Three-dimensional hydrogen-bonded frameworks in organic crystals: a topological study. Acta Crystallographica Section B: Structural Science, 2007, 63, 791-802.	1.8	72
33	Topological Databases: Why Do We Need Them for Design of Coordination Polymers?. Crystal Growth and Design, 2019, 19, 2604-2614.	3.0	72
34	A comparative crystallochemical analysis of binary compounds and simple anhydrous salts containing pyramidal anions LO3 (L = S, Se, Te, Cl, Br, I). Acta Crystallographica Section B: Structural Science, 2006, 62, 457-466.	1.8	71
35	Analysis of migration paths in fast-ion conductors with Voronoiâ€“Dirichlet partition. Acta Crystallographica Section B: Structural Science, 2006, 62, 1010-1018.	1.8	68
36	Nanocluster Model of Intermetallic Compounds with Giant Unit Cells: Î², Î²â€²-Mg₂Al₃ Polymorphs. Inorganic Chemistry, 2010, 49, 1811-1818.	4.0	68

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37	Molecular coordination numbers in crystal structures of organic compounds. Acta Crystallographica Section B: Structural Science, 2000, 56, 501-511.	1.8	67
38	Topology-based crystal structure generator. Computer Physics Communications, 2019, 236, 1-7.	7.5	67
39	Analysis of voids in crystal structures: the methods of `dual' crystal chemistry. Acta Crystallographica Section A: Foundations and Advances, 2003, 59, 34-44.	0.3	66
40	A topological method for the classification of entanglements in crystal networksA preliminary account of this work was presented at the workshop `Topological dynamics in physics and biology' held in Pisa, 12â€“13 July 2011.. Acta Crystallographica Section A: Foundations and Advances, 2012, 68, 484-493.	0.3	66
41	Deconstruction of Crystalline Networks into Underlying Nets: Relevance for Terminology Guidelines and Crystallographic Databases. Crystal Growth and Design, 2018, 18, 3411-3418.	3.0	65
42	Search for isotypism in crystal structures by means ofÂtheÂgraph theory. Acta Crystallographica Section A: Foundations and Advances, 2000, 56, 178-188.	0.3	63
43	A Water-Stable Cl@Ag₁₄ Cluster Based Metalâ€“Organic Open Framework for Dichromate Trapping and Bacterial Inhibition. Inorganic Chemistry, 2017, 56, 11891-11899.	4.0	60
44	Intermetallic compounds of the NaCd2 family perceived as assemblies of nanoclusters. Structural Chemistry, 2009, 20, 975-982.	2.0	57
45	Î³-Brass Polyhedral Core in Intermetallics: The Nanocluster Model. Inorganic Chemistry, 2013, 52, 13094-13107.	4.0	57
46	A method for hierarchical comparative analysis of crystal structures. Acta Crystallographica Section A: Foundations and Advances, 2006, 62, 356-364.	0.3	54
47	How 2-periodic coordination networks are interweaved: entanglement isomerism and polymorphism. CrystEngComm, 2017, 19, 1993-2006.	2.6	51
48	Crystallochemical tools in the search for cathode materials of rechargeable Na-ion batteries and analysis of their transport properties. Solid State Ionics, 2018, 314, 129-140.	2.7	51
49	Multilevel topological description of molecular packings in 1,2-benzothiazines. CrystEngComm, 2014, 16, 1963-1970.	2.6	44
50	A Possible Route toward Expert Systems in Supramolecular Chemistry: 2-Periodic H-Bond Patterns in Molecular Crystals. Crystal Growth and Design, 2014, 14, 1938-1949.	3.0	44
51	Synthesis, structure, topology and magnetic properties of cobalt(ii) coordination polymers with 2-nitrobiphenyl-4,4â€²-dicarboxylic acid and bis(pyridyl) ligands. Dalton Transactions, 2012, 41, 14316.	3.3	43
52	A method for topological analysis of rod packings. Structural Chemistry, 2016, 27, 1605-1611.	2.0	43
53	New Types of Multishell Nanoclusters with a Frankâ€“Kasper Polyhedral Core in Intermetallics. Inorganic Chemistry, 2011, 50, 5714-5724.	4.0	39
54	Simplify to understand: how to elucidate crystal structures?. Structural Chemistry, 2021, 32, 507-519.	2.0	39

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55	Structures of the ZrZn₂₂ family: suprapolyhedral nanoclusters, methods of self-assembly and superstructural ordering. Acta Crystallographica Section B: Structural Science, 2009, 65, 300-307.	1.8	37
56	On the Way to New Possible Naâ€Ion Conductors: The Voronoiâ€“Dirichlet Approach, Data Mining and Symmetry Considerations in Ternary Na Oxides. Chemistry - A European Journal, 2015, 21, 16601-16608.	3.3	37
57	Crystal Structure and Li-Ion Transport in Li₂CoPO₄F High-Voltage Cathode Material for Li-Ion Batteries. Journal of Physical Chemistry C, 2017, 121, 3194-3202.	3.1	37
58	High-throughput search for potential potassium ion conductors: A combination of geometrical-topological and density functional theory approaches. Solid State Ionics, 2018, 326, 188-199.	2.7	37
59	Sonochemical synthesis and characterization of four nanostructural nickel coordination polymers and photocatalytic degradation of methylene blue. Ultrasonics Sonochemistry, 2019, 56, 213-228.	8.2	36
60	A Collection of Topological Types of Nanoclusters and Its Application to Icosahedron-Based Intermetallics. Inorganic Chemistry, 2015, 54, 6616-6630.	4.0	35
61	A series of Cd(<scp>ii</scp>) coordination polymers based on flexible bis(triazole) and multicarboxylate ligands: topological diversity, entanglement and properties. CrystEngComm, 2017, 19, 5797-5808.	2.6	34
62	Topology versus porosity: what can reticular chemistry tell us about free space in metalâ€“organic frameworks?. Chemical Communications, 2020, 56, 9616-9619.	4.1	34
63	Analysis of ion-migration paths in inorganic frameworks by means of tilings and Voronoiâ€“Dirichlet partition: a comparison. Acta Crystallographica Section B: Structural Science, 2009, 65, 426-434.	1.8	32
64	From clusters to crystals: scale chemistry of intermetallics. Structural Chemistry, 2019, 30, 2015-2027.	2.0	32
65	Packing topology in crystals of proteins and small molecules: a comparison. Scientific Reports, 2017, 7, 13209.	3.3	31
66	Discovery of intrinsic two-dimensional antiferromagnets from transition-metal borides. Nanoscale, 2021, 13, 8254-8263.	5.6	31
67	Lithium-cation conductivity and crystal structure of lithium diphosphate. Journal of Solid State Chemistry, 2014, 211, 170-175.	2.9	30
68	Topology of Intermetallic Structures: From Statistics to Rational Design. Accounts of Chemical Research, 2018, 51, 21-30.	15.6	30
69	Crystal chemistry of orthosilicates and their analogs: the classification by topological types of suprapolyhedral structural units. Acta Crystallographica Section B: Structural Science, 2002, 58, 948-964.	1.8	27
70	New knowledge and tools for crystal design: local coordination versus overall network topology and much more. CrystEngComm, 2015, 17, 2913-2924.	2.6	27
71	Two d10 metal coordination polymers as dual functional luminescent probes for sensing of Fe3+ ions and acetylacetone with high selectivity and sensitivity. Journal of Solid State Chemistry, 2020, 289, 121460.	2.9	27
72	Crystal Structures of Inorganic Oxoacid Salts Perceived as Cation Arrays: A Periodic-Graph Approach. Structure and Bonding, 2011, , 31-66.	1.0	25

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73	Analysis of environment of alkali atoms in oxygen-containing compounds with Voronoi-Dirichlet polyhedra. Zeitschrift Fur Kristallographie - Crystalline Materials, 1998, 213, 202-209.	0.8	23
74	Methods for topological analysis of atomic nets. Journal of Structural Chemistry, 2009, 50, 160-167.	1.0	23
75	A â€œStronglyâ€•Self-Catenated Metalâ€“Organic Framework with the Highest Topological Density among 3,4-Coordinated Nets. Inorganic Chemistry, 2013, 52, 10732-10734.	4.0	23
76	An unusual (3,10)-coordinated 3D network coordination polymer as a potential luminescent sensor for detection of nitroaromatics and ferric ion. Journal of Luminescence, 2018, 199, 126-132.	3.1	23
77	Topology: ToposPro. , 2021, , 389-412.		23
78	Orthotetrahedral crystal structures My(TO4)z (T = Si, Ge, P, As, S, Se, Cl, Br, I): geometrical-topological analysis, quasi-binary representation, and comparison with the AyXz compounds by the method of coordination sequences. Zeitschrift Fur Kristallographie - Crystalline Materials, 2004, 219, 468-478.	0.8	21
79	Predicting New Zeolites: A Combination of Thermodynamic and Kinetic Factors. Chemistry of Materials, 2018, 30, 2829-2837.	6.7	21
80	Network topological model of reconstructive solid-state transformations. Scientific Reports, 2019, 9, 6007.	3.3	21
81	Comparative topological analysis of simple anhydrous borates, carbonates and nitrates. Zeitschrift Fur Kristallographie - Crystalline Materials, 2002, 217, .	0.8	20
82	Sizes of molecules in organic crystals: the Voronoiâ€“Dirichlet approach. Acta Crystallographica Section B: Structural Science, 2004, 60, 447-452.	1.8	20
83	Structural chemistry of metal microclusters: Questions and answers. Glass Physics and Chemistry, 2009, 35, 1-12.	0.7	20
84	New method for computer analysis of complex intermetallic compounds and nanocluster model of the samson phase Cd3Cu4. Crystallography Reports, 2010, 55, 1100-1105.	0.6	20
85	A Database of Topological Representations of Polynuclear Nickel Compounds. European Journal of Inorganic Chemistry, 2013, 2013, 520-526.	2.0	20
86	Computer-Aided Modeling of Aluminophosphate Zeolites As Packings of Building Units. Journal of Physical Chemistry C, 2012, 116, 6734-6744.	3.1	19
87	Topological analysis of ionic packings in crystal structures of inorganic sulfides: the method of coordination sequences. Zeitschrift Fur Kristallographie - Crystalline Materials, 2001, 216, 165-171.	0.8	19
88	Conduction mechanism in the low-temperature phase of KAlO2. Inorganic Materials, 2010, 46, 1234-1241.	0.8	18
89	Anisotropy of Elastic Properties of Metalâ€“Organic Frameworks and the Breathing Phenomenon. Journal of Physical Chemistry C, 2019, 123, 24651-24658.	3.1	18
90	The CSD and knowledge databases: from answers to questions. CrystEngComm, 2020, 22, 7298-7307.	2.6	18

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91	CrystalGrower: a generic computer program for Monte Carlo modelling of crystal growth. Chemical Science, 2021, 12, 1126-1146.	7.4	18
92	B₅N₃ and B₇N₅ Monolayers with High Carrier Mobility and Excellent Optical Performance. Journal of Physical Chemistry Letters, 2021, 12, 4823-4832.	4.6	18
93	Order and Topology in Systems with Many Particles. Acta Crystallographica Section A: Foundations and Advances, 1997, 53, 144-160.	0.3	17
94	The natural tiling approach to cation conductivity in KAlO₂ polymorphs. Acta Crystallographica Section B: Structural Science, 2012, 68, 356-363.	1.8	17
95	Combined DFT and geometricalâ€“topological analysis of Li-ion conductivity in complex hydrides. Inorganic Chemistry Frontiers, 2020, 7, 3115-3125.	6.0	17
96	Topological methods for analysis and design of coordination polymers. Russian Chemical Reviews, 2022, 91, .	6.5	17
97	Analysis of Lanthanide Ï€ Complexes in Terms of Voronoiâ€“Dirichlet Polyhedra. Russian Journal of Coordination Chemistry/Koordinatsionnaya Khimiya, 2000, 26, 847-856.	1.0	16
98	Local Coordination versus Overall Topology in Crystal Structures: Deriving Knowledge from Crystallographic Databases. Crystal Growth and Design, 2017, 17, 774-785.	3.0	16
99	Ionic Conductivity in Ti-Doped KFeO₂: Experiment and Mathematical Modeling. Journal of Physical Chemistry C, 2017, 121, 21128-21135.	3.1	16
100	Identification of solid oxygenâ€containing Naâ€electrolytes: An assessment based on crystallographic and economic parameters. Crystal Research and Technology, 2017, 52, 1600223.	1.3	16
101	Unique self-catenated 3D Cd(II)-MOF with a rare (3,3,9)-connected underlying network exhibiting high photocatalytic activities. Inorganic Chemistry Communication, 2019, 102, 126-129.	3.9	16
102	Topological representations of crystal structures: generation, analysis and implementation in the TopCryst system. Science and Technology of Advanced Materials Methods, 2022, 2, 250-265.	1.3	16
103	Computer simulation of the self-assembly of paulingite crystal structure from suprapolyhedral nanocluster precursors K6, K16, and K20. Crystallography Reports, 2011, 56, 75-83.	0.6	15
104	Nanoporous materials with predicted zeolite topologies. RSC Advances, 2020, 10, 17760-17767.	3.6	15
105	Study of rare-earth Ï€-complexes by means of Voronoiâ€“Dirichlet polyhedra. Acta Crystallographica Section B: Structural Science, 2001, 57, 261-270.	1.8	14
106	Symmetry and topology codes of cluster self-assembly for icosahedral structures of the NaZn13-cF112 and TRB66-cF1944 family. Glass Physics and Chemistry, 2015, 41, 341-351.	0.7	14
107	Knowledge-Based Approaches to H-Bonding Patterns in Heterocycle-1-Carbohydrazoneamides. Crystal Growth and Design, 2016, 16, 6354-6362.	3.0	14
108	Natural tilings and free space in zeolites: models, statistics, correlations, prediction. Zeitschrift Fur Kristallographie - Crystalline Materials, 2019, 234, 421-436.	0.8	14

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109	High-throughput systematic topological generation of low-energy carbon allotropes. Npj Computational Materials, 2021, 7, .	8.7	14
110	Theoretical crystal chemistry of phosphates: Topological analysis and classification of suprapolyhedral ensembles of orthophosphates and their analogues M x (TO4)y. Crystallography Reports, 2004, 49, 327-342.	0.6	13
111	Vacancy Ordering as a Driving Factor for Structural Changes in Ternary Germanides: The New R2Zn1â€“xGe6Series of Polar Intermetallics (R = Rare-Earth Metal). Inorganic Chemistry, 2015, 54, 2411-2424.	4.0	13
112	Molecular coordination numbers and crystal structure of simple substances. Computational and Theoretical Chemistry, 1999, 489, 225-236.	1.5	12
113	An unusual (4,6)-coordinated copper(II) coordination polymer: High efficient degradation of organic dyes under visible light irradiation and electrochemical properties. Polyhedron, 2018, 148, 81-87.	2.2	12
114	Computational Search for Novel Zn-Ion Conductorsâ€”A Crystallochemical, Bond Valence, and Density Functional Study. Journal of Physical Chemistry C, 2021, 125, 17590-17599.	3.1	12
115	Analysis of the environment of beryllium, magnesium and alkaline earth atoms in oxygen-containing compounds. Acta Crystallographica Section B: Structural Science, 1999, 55, 139-146.	1.8	11
116	Structure-forming components in crystals of ternary and quaternary 3d-metal complex fluorides. Acta Crystallographica Section B: Structural Science, 2003, 59, 361-377.	1.8	11
117	Structure of a new high-pressureâ€“high-temperature modification of antimony(III) oxide, Î³-Sb₂O₃, from high-resolution synchrotron powder diffraction data. Acta Crystallographica Section B: Structural Science, 2012, 68, 1-7.	1.8	11
118	New types of two-layer nanoclusters with an icosahedral core. Glass Physics and Chemistry, 2013, 39, 229-234.	0.7	11
119	Molecular packings and specific-bonding patterns in sulfonamides. New Journal of Chemistry, 2014, 38, 4099-4106.	2.8	11
120	Topological methods for complex intermetallics. Zeitschrift Fur Kristallographie - Crystalline Materials, 2017, 232, 497-506.	0.8	11
121	A fascinating building unit: Mackay cluster in intermetallics. Structural Chemistry, 2017, 28, 133-140.	2.0	11
122	Crystal and electronic structure engineering of tin monoxide by external pressure. Journal of Advanced Ceramics, 2021, 10, 565-577.	17.4	11
123	Visualization and Quantification of Geometric Diversity in Metalâ€“Organic Frameworks. Chemistry of Materials, 0, , .	6.7	11
124	Methods of crystallochemical analysis of supramolecular complexes by means of Voronoiâ€“Dirichlet polyhedra: a study of cucurbituril hostâ€“guest compounds. Acta Crystallographica Section B: Structural Science, 2004, 60, 350-357.	1.8	10
125	Assembly models for zeolite crystal structures according to the data of topological analysis by the tiling method. Crystallography Reports, 2012, 57, 875-884.	0.6	10
126	Specific features of the crystal structure of polymorphous modifications of KFeO2 and their correlation with ionic conductivity. Physics of the Solid State, 2013, 55, 1050-1056.	0.6	10

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127	Construction of five zinc coordination polymers with 4-substituted bis(trizole) and multicarboxylate ligands: Syntheses, structures and properties. Polyhedron, 2018, 155, 223-231.	2.2	10
128	Cluster self-organization of intermetallic systems: Quasi-spherical nanocluster precursors with internal Friauf polyhedra (A-172) and icosahedra (B-137) in the Li19Na8Ba15 (hP842) crystal structure. Crystallography Reports, 2010, 55, 1093-1099.	0.6	9
129	Structure and chemical composition of the new zeolite ISC-1 from the data of nanocluster modeling. Glass Physics and Chemistry, 2010, 36, 663-672.	0.7	9
130	Interpenetration of three-periodic networks in crystal structures: Description and classification methods, geometrical-topological conditions of implementation. Journal of Structural Chemistry, 2014, 55, 1308-1325.	1.0	9
131	Topos: Program package for topological analysis of crystal structures. Journal of Structural Chemistry, 1994, 34, 820-822.	1.0	8
132	A new set of molecular descriptors. Acta Crystallographica Section B: Structural Science, 2002, 58, 219-226.	1.8	8
133	Theoretical crystal chemistry of M x (TO4)y sulfates and selenates: Topological analysis and classification of suprapolyhedral invariants. Crystallography Reports, 2006, 51, 366-378.	0.6	8
134	Intentional selection of coordination compounds with the required thermochemical properties on the basis of the cambridge bank of structural data. Russian Journal of Physical Chemistry A, 2012, 86, 1340-1351.	0.6	8
135	Two new zinc(ii) coordination complexes with helix characteristics showing both interpenetration and self-catenation features: a platform for the synthesis of chiral and catenated structures assembled by length-modulated dicarboxylates. Dalton Transactions, 2014, 43, 15151-15158.	3.3	8
136	Combinatorial and topological modeling of cluster self-assembly of the crystal structure of zeolites. Crystallography Reports, 2015, 60, 453-465.	0.6	8
137	The role of local heteropolyhedral substitutions in the stoichiometry, topological characteristics and ion-migration paths in the eudialyte-related structures: a quantitative analysis. Acta Crystallographica Section B: Structural Science, Crystal Engineering and Materials, 2022, 78, 80-90.	1.1	8
138	Discovery of Electrides in Electronâ€Rich Nonâ€Electride Materials via Energy Modification of Interstitial Electrons. Advanced Functional Materials, 2022, 32, .	14.9	8
139	Metal-organic frameworks as the basis for new-generation functional materials. Russian Chemical Reviews, 2022, 91, .	6.5	8
140	Analysis of Li+ cation migration paths in oxygen-containing compounds. Russian Journal of Electrochemistry, 2009, 45, 417-428.	0.9	7
141	Cluster self-organization of crystal-forming systems: Suprapolyhedral cluster precursors and self-assembly of the icosahedral structure of ZrZn22(cF184). Crystallography Reports, 2009, 54, 548-554.	0.6	7
142	Geometric and topological analysis of icosahedral structures of samson Mg2Zn11 (cP39) Phases, K6Na15Tl18H (cP40), and Tm3In7Co9 (cP46): Nanocluster precursors, self-assembly mechanism, and superstructure ordering. Russian Journal of Inorganic Chemistry, 2011, 56, 729-737.	1.3	7
143	Modeling of self-organization processes in crystal-forming systems: Symmetry and topological codes of cluster self-assembly of a 2D layered icosahedral structure of Sc18B238 (Pbam, oP514). Glass Physics and Chemistry, 2016, 42, 221-229.	0.7	7
144	An Unprecedented 3D Self-catenated Four-coordinated Dense Net of Silver-Organic Framework. Bulletin of the Korean Chemical Society, 2013, 34, 1891-1894.	1.9	7

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145	Vaporization of Molecular Coordination Organotitanium Compounds: Development of the Structure-Thermochemical Approach with Programmed Use of the Cambridge Structural Database. Russian Journal of Coordination Chemistry/Koordinatsionnaya Khimiya, 2004, 30, 679-684.	1.0	5
146	Nanoclusters based on pentagondodecahedra with shells in the form of D32, D42, and D50 deltahedra in crystal structures of intermetallic compounds. Crystallography Reports, 2012, 57, 1-9.	0.6	5
147	On similarity of structure of icosahedral viral capsids and shells of metallic nanoclusters. Glass Physics and Chemistry, 2013, 39, 101-104.	0.7	5
148	Bergman, Bergman-based and 63-atom nanoclusters in intermetallics. Structural Chemistry, 2016, 27, 1685-1692.	2.0	5
149	Perceiving Zeolite Self-Assembly: A Combined Top-Down and Bottom-Up Approach within the Tiling Model. Journal of Physical Chemistry C, 2020, 124, 1523-1528.	3.1	5
150	A Comparative Analysis of Crystal Lattice Topology in Molybdates and Binary Compounds. Journal of Structural Chemistry, 2001, 42, 436-445.	1.0	4
151	Hierarchical crystal-chemical analysis of binary intermetallic compounds. Russian Journal of Inorganic Chemistry, 2007, 52, 1577-1585.	1.3	4
152	Computer modeling of self-assembly of the crystal structure of zeolite Na384[Al384Si384O1536] (H2O)422 (LTN, cF4080) from suprapolyhedral cluster precursors AB 2 (A-K 48, B-K 24). Crystallography Reports, 2012, 57, 360-368.	0.6	4
153	Computer simulation of the self-assembly of crystal structures of zeolites Ca64(Sr,K,Ba)48(Cu12(O,Cl))4[Si192Al192O786](H2O) n (tschoertnerite, TSC, V = 31 614 Ã...3) and Ca2K2[Al6Si6O24](H2O)10 (willhendersonite, cha, V = 804 Ã...3) from template nanocluster precursors K48 and K12. Crystallography Reports, 2013, 58, 531-540.	0.6	4
154	Modeling of self-organization processes in crystal-forming systems: Templated precursor nanoclusters T48 and the self-assembly of crystal structures of 15-crown-5, Na-FAU, 18-crown-6, Na-EMT, and Ca,Ba-TSC zeolites. Russian Journal of Inorganic Chemistry, 2015, 60, 469-482.	1.3	4
155	Symmetry and topology code of the cluster self-assembly of framework MT structures of alumophosphates AlPO4(H2O)2 (metavariscite and variscite) and Al2(PO4)2(H2O)3 (APC). Crystallography Reports, 2017, 62, 174-184.	0.6	4
156	Sulfurâ€•and Seleniumâ€Containing Compounds Potentially Exhibiting Al Ion Conductivity. Chemistry - A European Journal, 2019, 25, 8623-8629.	3.3	4
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