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	Computational design of materials for metal-ion batteries. , 2023, , 404-429.		4
2	Topological methods for analysis and design of coordination polymers. Russian Chemical Reviews, 2022, 91, .	6.5	17
3	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
4	Discovery of Electrides in Electronâ€Rich Nonâ€Electride Materials via Energy Modification of Interstitial Electrons. Advanced Functional Materials, 2022, 32, .	14.9	8
5	Polymorphism and topological features of compounds with the general formula <mml:math xmlns:mml="http://www.w3.org/1998/Math/MathML" altimg="si4.svg"><mml:mrow><mml:msubsup><mml:mi mathvariant="bold-italic">A</mml:mi><mml:mrow><mml:mn>1</mml:mn><mml:mo>â^'</mml:mo><mml:mi mathvariant="bold-italic">x</mml:mi></mml:mrow><mml:mo>+</mml:mo></mml:msubsup><mml:msubsup><mml:mi mathvariant="bold-italic">B</mml:mi></mml:mi>	2.9	3
6	Topological transformations in metalâ€“organic frameworks: a prospective design route?. CrystEngComm, 2022, 24, 2914-2924.	2.6	1
7	Metal-organic frameworks as the basis for new-generation functional materials. Russian Chemical Reviews, 2022, 91, .	6.5	8
8	Mining Knowledge from Crystal Structures: Oxidation States of Oxygen-Coordinated Metal Atoms in Ionic and Coordination Compounds. Journal of Chemical Information and Modeling, 2022, 62, 2332-2340.	5.4	4
9	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
10	Generating triply periodic surfaces from crystal structures: the tiling approach and its application to zeolites. Acta Crystallographica Section A: Foundations and Advances, 2022, 78, 327-336.	0.1	3
11	Topology: ToposPro. , 2021, , 389-412.		23
12	CrystalGrower: a generic computer program for Monte Carlo modelling of crystal growth. Chemical Science, 2021, 12, 1126-1146.	7.4	18
13	Discovery of intrinsic two-dimensional antiferromagnets from transition-metal borides. Nanoscale, 2021, 13, 8254-8263.	5.6	31
14	Simplify to understand: how to elucidate crystal structures?. Structural Chemistry, 2021, 32, 507-519.	2.0	39
15	Crystal and electronic structure engineering of tin monoxide by external pressure. Journal of Advanced Ceramics, 2021, 10, 565-577.	17.4	11
16	B₅N₃ and B₇N₅ Monolayers with High Carrier Mobility and Excellent Optical Performance. Journal of Physical Chemistry Letters, 2021, 12, 4823-4832.	4.6	18
17	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
18	High-throughput systematic topological generation of low-energy carbon allotropes. Npj Computational Materials, 2021, 7, .	8.7	14

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19	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
20	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
21	Combined DFT and geometricalâ€“topological analysis of Li-ion conductivity in complex hydrides. Inorganic Chemistry Frontiers, 2020, 7, 3115-3125.	6.0	17
22	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
23	From Simple to Complex: Design of Inorganic Crystal Structures with a Topologically Extended Zintlâ€“Klemm Concept. Journal of Physical Chemistry Letters, 2020, 11, 8114-8120.	4.6	1
24	Monolayer selfâ€•organization of cyclodextrins on carbon surface. Journal of the Chinese Chemical Society, 2020, 67, 1778-1782.	1.4	1
25	Nanoporous materials with predicted zeolite topologies. RSC Advances, 2020, 10, 17760-17767.	3.6	15
26	The CSD and knowledge databases: from answers to questions. CrystEngComm, 2020, 22, 7298-7307.	2.6	18
27	From clusters to crystals: scale chemistry of intermetallics. Structural Chemistry, 2019, 30, 2015-2027.	2.0	32
28	A universal algorithm for finding the shortest distance between systems of points. Acta Crystallographica Section A: Foundations and Advances, 2019, 75, 827-832.	0.1	3
29	Anisotropy of Elastic Properties of Metalâ€“Organic Frameworks and the Breathing Phenomenon. Journal of Physical Chemistry C, 2019, 123, 24651-24658.	3.1	18
30	Construction of (3,8)-connected three-dimensional cobalt(II) and copper(II) coordination polymers with 1,3-bis[(1,2,4-triazol-4-yl)methyl]benzene and benzene-1,3,5-tricarboxylate ligands. Acta Crystallographica Section C, Structural Chemistry, 2019, 75, 960-968.	0.5	2
31	Sulfurâ€•and Seleniumâ€Containing Compounds Potentially Exhibiting Al Ion Conductivity. Chemistry - A European Journal, 2019, 25, 8623-8629.	3.3	4
32	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
33	The Aluminum-Ion Battery: A Sustainable and Seminal Concept?. Frontiers in Chemistry, 2019, 7, 268.	3.6	155
34	Network topological model of reconstructive solid-state transformations. Scientific Reports, 2019, 9, 6007.	3.3	21
35	Topological Databases: Why Do We Need Them for Design of Coordination Polymers?. Crystal Growth and Design, 2019, 19, 2604-2614.	3.0	72
36	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

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37	Natural tilings and free space in zeolites: models, statistics, correlations, prediction. Zeitschrift Fur Kristallographie - Crystalline Materials, 2019, 234, 421-436.	0.8	14
38	Topology-based crystal structure generator. Computer Physics Communications, 2019, 236, 1-7.	7.5	67
39	Rules for designing rod metalâ€“organic frameworks: a topological approach. Acta Crystallographica Section A: Foundations and Advances, 2019, 75, e513-e513.	0.1	0
40	Crystallochemical computational tools, web services, databases, and approaches for joining researchers over the world. Acta Crystallographica Section A: Foundations and Advances, 2019, 75, e722-e722.	0.1	0
41	Predicting New Zeolites: A Combination of Thermodynamic and Kinetic Factors. Chemistry of Materials, 2018, 30, 2829-2837.	6.7	21
42	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
43	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
44	Topology of Intermetallic Structures: From Statistics to Rational Design. Accounts of Chemical Research, 2018, 51, 21-30.	15.6	30
45	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
46	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
47	4. Battery Materials. , 2018, , 75-260.		0
48	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
49	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
50	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
51	Nanocluster model and its application for crystal structure prediction of complex intermetallics. Acta Crystallographica Section A: Foundations and Advances, 2018, 74, e54-e54.	0.1	0
52	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
53	Modeling of self-organization in crystal-forming systems: Symmetry and topology codes of cluster self-assembly of icosahedral structure Sc12B185C9 (P6/mmm, hP212). Glass Physics and Chemistry, 2017, 43, 115-124.	0.7	0
54	Predicting crystal growth via a unified kinetic three-dimensional partition model. Nature, 2017, 544, 456-459.	27.8	88

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55	Topological methods for complex intermetallics. Zeitschrift Fur Kristallographie - Crystalline Materials, 2017, 232, 497-506.	0.8	11
56	How 2-periodic coordination networks are interweaved: entanglement isomerism and polymorphism. CrystEngComm, 2017, 19, 1993-2006.	2.6	51
57	Local Coordination versus Overall Topology in Crystal Structures: Deriving Knowledge from Crystallographic Databases. Crystal Growth and Design, 2017, 17, 774-785.	3.0	16
58	Packing topology in crystals of proteins and small molecules: a comparison. Scientific Reports, 2017, 7, 13209.	3.3	31
59	Ionic Conductivity in Ti-Doped KFeO₂: Experiment and Mathematical Modeling. Journal of Physical Chemistry C, 2017, 121, 21128-21135.	3.1	16
60	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
61	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
62	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
63	A fascinating building unit: Mackay cluster in intermetallics. Structural Chemistry, 2017, 28, 133-140.	2.0	11
64	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
65	Topological approach for the design of new materials. Acta Crystallographica Section A: Foundations and Advances, 2017, 73, C21-C21.	0.1	0
66	Multilevel topological analysis in application to design of coordination networks. Acta Crystallographica Section A: Foundations and Advances, 2017, 73, C73-C73.	0.1	0
67	Assessment of potential Al ion conductors from large crystallographic databases. Acta Crystallographica Section A: Foundations and Advances, 2017, 73, C328-C328.	0.1	0
68	Two-shell nanoclusters in intermetallic compounds: beyond the icosahedral core. Acta Crystallographica Section A: Foundations and Advances, 2016, 72, s316-s316.	0.1	0
69	Bergman, Bergman-based and 63-atom nanoclusters in intermetallics. Structural Chemistry, 2016, 27, 1685-1692.	2.0	5
70	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
71	Knowledge-Based Approaches to H-Bonding Patterns in Heterocycle-1-Carbohydrazoneamides. Crystal Growth and Design, 2016, 16, 6354-6362.	3.0	14
72	Topological collections and knowledge bases for applications in crystal engineering of coordination compounds. Acta Crystallographica Section A: Foundations and Advances, 2016, 72, s165-s165.	0.1	0

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73	Topological analysis of self-catenated motifs in coordination networks. Acta Crystallographica Section A: Foundations and Advances, 2016, 72, s434-s434.	0.1	0
74	Can we predict Al ion conductors? A combination of crystallographic and energetic evaluation tools. Acta Crystallographica Section A: Foundations and Advances, 2016, 72, s291-s291.	0.1	0
75	Crystal structure, microstructure and ionic conductivity of the cost-efficient sodium solid electrolyte Na5YSi4O12. Acta Crystallographica Section A: Foundations and Advances, 2016, 72, s288-s288.	0.1	0
76	Synthesis, Structures, and Properties of Cadmium(II) and Nickel(II) Coordination Polymers Based on a 4,4â€²â€²Biphenylâ€Containing Ligand and Aliphatic Carboxylic Acids. Zeitschrift Fur Anorganische Und Allgemeine Chemie, 2016, 642, 1184-1190.	1.2	2
77	Combinatorial-topological modeling of the cluster self-assembly of zeolite crystal structures: computer search for molecular templates for new zeolite ISC-2. Russian Chemical Bulletin, 2016, 65, 29-39.	1.5	2
78	A method for topological analysis of rod packings. Structural Chemistry, 2016, 27, 1605-1611.	2.0	43
79	An unprecedented â€œstronglyâ€•self-catenated MOF containing inclined catenated honeycomb-like units. Dalton Transactions, 2016, 45, 2426-2429.	3.3	87
80	The taxonomy of rod-packing coordination networksÂ(CNs). Acta Crystallographica Section A: Foundations and Advances, 2016, 72, s435-s435.	0.1	1
81	Crystallochemical analysis of ion conductivity in K+-oxygen containing inorganic compounds. Acta Crystallographica Section A: Foundations and Advances, 2016, 72, s292-s293.	0.1	0
82	SACADA - the database of three periodic carbon allotropes. Acta Crystallographica Section A: Foundations and Advances, 2015, 71, s356-s356.	0.1	0
83	On the relation between topology of halogen-bondedÂmolecular crystals and type of halogen-halogen contacts. Acta Crystallographica Section A: Foundations and Advances, 2015, 71, s454-s454.	0.1	0
84	A knowledge database for intermetallics: the collection of Topological Types of Nanoclusters. Acta Crystallographica Section A: Foundations and Advances, 2015, 71, s357-s357.	0.1	0
85	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
86	A Collection of Topological Types of Nanoclusters and Its Application to Icosahedron-Based Intermetallics. Inorganic Chemistry, 2015, 54, 6616-6630.	4.0	35
87	Rod packing nets in 3-periodic metal-organic frameworks. Acta Crystallographica Section A: Foundations and Advances, 2015, 71, s357-s357.	0.1	0
88	Structural chemistry of metal inorganic 3D MT frameworks. Templated nanoclustersâ€“precursors Ga2(PO4)2F2 and Ga2(PO4)2(HF)F2 and self-assembly of crystalline gallophosphate structures (NH4)2[Ga2(PO4)2F2](KTP-Type) and (NH4)2[Ga2(PO4)2(HF)F2] (p-KTP-Type). Glass Physics and Chemistry, 2015, 41, 561-571.	0.7	0
89	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
90	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

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91	New knowledge and tools for crystal design: local coordination versus overall network topology and much more. CrystEngComm, 2015, 17, 2913-2924.	2.6	27
92	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
93	Topological Motifs in Cyanometallates: From Building Units to Three-Periodic Frameworks. Chemical Reviews, 2015, 115, 12286-12319.	47.7	128
94	Combinatorial-topological modeling of the cluster self-assembly of crystal structures of zeolites of the GME, AFX, AFT, and ISC-2 family. Glass Physics and Chemistry, 2015, 41, 443-452.	0.7	3
95	Combinatorial and topological modeling of cluster self-assembly of the crystal structure of zeolites. Crystallography Reports, 2015, 60, 453-465.	0.6	8
96	Topological properties of crystal structures: from description to prediction. Acta Crystallographica Section A: Foundations and Advances, 2015, 71, s74-s74.	0.1	0
97	The nanocluster approach to elucidate complex intermetallics. Acta Crystallographica Section A: Foundations and Advances, 2015, 71, s54-s54.	0.1	0
98	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
99	Icosahedral nanoclusters-precursors and self-assembly of crystal structures of the WAl12 (Im-3, cI26) family and sillenite Bi12 SiO20 (I23, cI66). Glass Physics and Chemistry, 2014, 40, 591-599.	0.7	0
100	Multilevel topological description of molecular packings in 1,2-benzothiazines. CrystEngComm, 2014, 16, 1963-1970.	2.6	44
101	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
102	Molecular packings and specific-bonding patterns in sulfonamides. New Journal of Chemistry, 2014, 38, 4099-4106.	2.8	11
103	Lithium-cation conductivity and crystal structure of lithium diphosphate. Journal of Solid State Chemistry, 2014, 211, 170-175.	2.9	30
104	Structural chemistry of organo-siloxanes: Composition and structure of Si n (O,C) m (n = 2â€“21) clusters with Si-O-Si bridging bonds. Glass Physics and Chemistry, 2014, 40, 180-189.	0.7	3
105	Entangled Two-Dimensional Coordination Networks: A General Survey. Chemical Reviews, 2014, 114, 7557-7580.	47.7	253
106	Applied Topological Analysis of Crystal Structures with the Program Package ToposPro. Crystal Growth and Design, 2014, 14, 3576-3586.	3.0	2,448
107	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
108	New types of two-layer nanoclusters with an icosahedral core. Glass Physics and Chemistry, 2013, 39, 229-234.	0.7	11

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109	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
110	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
111	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
112	Î³-Brass Polyhedral Core in Intermetallics: The Nanocluster Model. Inorganic Chemistry, 2013, 52, 13094-13107.	4.0	57
113	On similarity of structure of icosahedral viral capsids and shells of metallic nanoclusters. Glass Physics and Chemistry, 2013, 39, 101-104.	0.7	5
114	A Database of Topological Representations of Polynuclear Nickel Compounds. European Journal of Inorganic Chemistry, 2013, 2013, 520-526.	2.0	20
115	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
116	Topology of 2-Periodic Coordination Networks: Toward Expert Systems in Crystal Design. Crystal Growth and Design, 2013, 13, 1655-1664.	3.0	119
117	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
118	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
119	A topological study of three-dimensional hydrogen-bonded frameworks. Acta Crystallographica Section A: Foundations and Advances, 2013, 69, s484-s484.	0.3	0
120	4-150-atom Lennardâ€“Jones clusters in intermetallics. Acta Crystallographica Section A: Foundations and Advances, 2013, 69, s457-s457.	0.3	0
121	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
122	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
123	Computer-Aided Modeling of Aluminophosphate Zeolites As Packings of Building Units. Journal of Physical Chemistry C, 2012, 116, 6734-6744.	3.1	19
124	New icosahedral nanoclusters in crystal structures of intermetallic compounds: Topological types of 50-atom deltahedra D50 in samson phases Î²-Mg2Al3 and É›-Mg23Al30. Crystallography Reports, 2012, 57, 885-891.	0.6	2
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	Two metalâ€“organic frameworks with unique high-connected binodal network topologies: synthesis, structures, and catalytic properties. CrystEngComm, 2012, 14, 4210.	2.6	196

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127	Totally unimodular nets. Acta Crystallographica Section A: Foundations and Advances, 2012, 68, 286-294.	0.3	3
128	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
129	A 1D â† 3D Twoâ€Fold Interpenetration Array Formed by Hydrogenâ€Bonding Interactions. Zeitschrift Fur Anorganische Und Allgemeine Chemie, 2012, 638, 992-995.	1.2	1
130	Topological systematization of layered coordination compounds of Cu, Ag, Zn, and Cd. Russian Journal of Coordination Chemistry/Koordinatsionnaya Khimiya, 2012, 38, 309-314.	1.0	1
131	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
132	Nanocluster analysis of intermetallic structures with the program package TOPOS. Structural Chemistry, 2012, 23, 955-963.	2.0	488
133	High-nuclearity cobalt coordination clusters: Synthetic, topological and magnetic aspects. Coordination Chemistry Reviews, 2012, 256, 1246-1278.	18.8	204
134	The natural tiling approach to cation conductivity in KAlO₂ polymorphs. Acta Crystallographica Section B: Structural Science, 2012, 68, 356-363.	1.8	17
135	Mechanism of structural phase transitions in the Li4GeO4-ZnGeO4 system: Computer modeling and identification of invariant nanocluster structures in Li4GeO4, LISICON Li6Zn(GeO4)2, and Li4Zn2(GeO4)2 (Î³ phase). Russian Journal of Inorganic Chemistry, 2012, 57, 846-853.	1.3	2
136	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
137	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
138	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
139	New Types of Multishell Nanoclusters with a Frankâ€“Kasper Polyhedral Core in Intermetallics. Inorganic Chemistry, 2011, 50, 5714-5724.	4.0	39
140	Crystal Structures of Inorganic Oxoacid Salts Perceived as Cation Arrays: A Periodic-Graph Approach. Structure and Bonding, 2011, , 31-66.	1.0	25
141	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
142	Topological systematization of the framework coordination polymers formed by iron, cobalt, or nickel complexes. Russian Journal of Coordination Chemistry/Koordinatsionnaya Khimiya, 2011, 37, 81-94.	1.0	2
143	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
144	Cluster self-organization of germanate systems: Suprapolyhedral precursor nanoclusters and crystal structures re]20100419. Russian Journal of Inorganic Chemistry, 2011, 56, 1572-1578.	1.3	3

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145	Geometric and topological analysis of zeolite crystal structures by the tiling method: The model of structure of Na,K-paulingite (PAU) Na82K72[Al154Si518O1344] Â· wH2O. Russian Journal of Inorganic Chemistry, 2011, 56, 1782-1787.	1.3	2
146	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
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