

# Andrey G Starikov

## List of Publications by Year in descending order

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152  
papers

1,653  
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docs citations

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times ranked

1198  
citing authors

#	ARTICLE	IF	CITATIONS
1	Computational search for redox isomerism in Ge and Sn bis-chelates with $\hat{\pm}$ -diimine ligands. Mendeleev Communications, 2022, 32, 49-51.	0.6	1
2	Spin transitions in ferric catecholate complexes mediated by outer-sphere counteranions. Dalton Transactions, 2022, 51, 10909-10919.	1.6	4
3	Structural Changes in Five-coordinate Bromido-bis(o-aminobenzo-semiquinonato)iron(III) Complex: Spin-Crossover or Ligand-Metal Antiferromagnetic Interactions?. European Journal of Inorganic Chemistry, 2021, 2021, 756-762.	1.0	1
4	DFT computational insight into the mechanism of the monomer-trimer isomerism of Ni(II) bis-acetylacetonate. Inorganica Chimica Acta, 2021, 517, 120183.	1.2	3
5	Computational insight into magnetic behaviour of cobalt tris(2-pyridylmethyl)amine complexes with dioxolenes incorporating stable radicals. Chemical Physics Letters, 2021, 762, 138128.	1.2	11
6	Synthesis, Structure and Redox Properties of Cu(II) Chelate Complexes on the Basis of 2-(Hydroxyphenyl)-1-H-benzo[d]imidazol-1-yl Phenol Ligands. European Journal of Inorganic Chemistry, 2021, 2021, 2055-2062.	1.0	4
7	XAS Diagnostic of the Photoactive State in Co(II) Azobenzene Complex in Organic Solvents. ChemistrySelect, 2021, 6, 7087-7092.	0.7	0
8	Dithiolate and Catecholate Binding of Copper by the $OO\frac{1}{4}SS$ Bifunctional Ligand: Regioselectivity and Regioisomeric Transformations. European Journal of Inorganic Chemistry, 2021, 2021, 3292-3300.	1.0	3
9	Acene-Linked Zethrenes and Bisphenalenyls: A DFT Search for Organic Tetraradicals. Journal of Physical Chemistry A, 2021, 125, 6562-6570.	1.1	13
10	Spin-State-Switching Rearrangements of Bis(dioxolene)-Bridged CrCo Complexes: A DFT Study. European Journal of Inorganic Chemistry, 2021, 2021, 4113-4121.	1.0	2
11	Structure, spectral-luminescent and ionochromic properties of hydroxyaryl(hetaryl)idene azomethine imines. Journal of Molecular Structure, 2020, 1199, 127013.	1.8	3
12	o-Quinone phenalenyl derivatives as expedient ligands for the design of magnetically active metal complexes: A computational study. Chemical Physics Letters, 2020, 740, 137073.	1.2	6
13	Electronic structure and magnetic properties of the triangular nanographenes with radical substituents: a DFT study. Physical Chemistry Chemical Physics, 2020, 22, 1288-1298.	1.3	18
14	Heterospin magnetically active bimetallic Fe and Co complexes of aldiminato-functionalized catechol: a DFT study. Structural Chemistry, 2020, 31, 37-46.	1.0	5
15	Immobilization of UiO-67 with photochromic spiropyrans: a quantum chemical study. Journal of Molecular Modeling, 2020, 26, 212.	0.8	2
16	Low-valent oligogermanium amidophenolate complex comprising a unique Ge <sub>4</sub> chain. Mendeleev Communications, 2020, 30, 205-208.	0.6	18
17	UiO-66 type MOFs with mixed-linkers - 1,4-Benzenedicarboxylate and 1,4-naphthalenedicarboxylate: Effect of the modulator and post-synthetic exchange. Microporous and Mesoporous Materials, 2020, 305, 110324.	2.2	33
18	Some aspects of the formation and structural features of low nuclearity heterometallic carboxylates. Pure and Applied Chemistry, 2020, 92, 1093-1110.	0.9	21

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19	The chemical and electrochemical reduction of heteroligand o-semiquinonato-formazanato cobalt complexes. <i>Inorganica Chimica Acta</i> , 2019, 489, 1-7.	1.2	19
20	Tetrahedral nickel(ii) and cobalt(ii) bis-o-iminobenzosemiquinonates. <i>Dalton Transactions</i> , 2019, 48, 10723-10732.	1.6	20
21	Rational Design of Electronically Labile Dinuclear Fe and Co complexes with 1,10-Phenanthroline-5,6-diimine: A DFT study. <i>Journal of Computational Chemistry</i> , 2019, 40, 2284-2292.	1.5	5
22	Metal-ligand ferromagnetic exchange interactions in heteroligand bis-o-semiquinonato nickel complexes with 2,2'-dipyridine and 1,10-phenanthroline. <i>Polyhedron</i> , 2019, 158, 262-269.	1.0	14
23	Insight into the Electron Density Distribution in an O,N-Heterocyclic Stannylene by High-Resolution X-ray Diffraction Analysis. <i>European Journal of Inorganic Chemistry</i> , 2019, 2019, 875-884.	1.0	22
24	Synthesis and redox activity of the ruthenium complexes based on 9-hydroxy-2,4,6,8-tetra-(tert.-butyl)phenoxazin-1-one ligands. <i>Inorganica Chimica Acta</i> , 2019, 484, 430-436.	1.2	4
25	A DFT computational study of the magnetic behaviour of cobalt dioxolene complexes of tetraazamacrocyclic ligands. <i>Computational and Theoretical Chemistry</i> , 2018, 1124, 15-22.	1.1	36
26	Computational insight into magnetic behavior and properties of the transition metal complexes with redox-active ligands: a DFT approach. <i>Pure and Applied Chemistry</i> , 2018, 90, 811-824.	0.9	41
27	Pentacoordinated chloro-bis-o-iminosemiquinonato Mn and Fe complexes. <i>Journal of Molecular Structure</i> , 2018, 1165, 51-61.	1.8	20
28	Dispersion interactions in oligomerization of metal diketonates: a DFT evaluation. <i>Chemical Papers</i> , 2018, 72, 829-839.	1.0	9
29	The structurally variable network of spin couplings and migrating paramagnetic centers in binuclear o-quinone Co <sup>II</sup> complexes with biradical acene linkers: a computational DFT study. <i>Dalton Transactions</i> , 2018, 47, 15948-15956.	1.6	10
30	Computational modeling of mixed-ligand cobalt complexes with o-quinone derivative of corannulene. <i>Russian Chemical Bulletin</i> , 2018, 67, 1978-1984.	0.4	1
31	Cobalt complexes with hemilabile o-quinone-iminobenzoquinonate ligands: a novel example of redox-induced electron transfer. <i>Dalton Transactions</i> , 2018, 47, 15049-15060.	1.6	33
32	Photochromic Properties and Surface Enhanced Raman Scattering Spectra of Indoline Spiropyran in Silver-Based Nanocomposite Films. <i>Optics and Spectroscopy (English Translation of Optika i Spektroskopiya)</i> , 2018, 44, 1011-1017.	1.0	1
33	Magnetic Properties of Acenes and Their o-Quinone Derivatives: Computer Simulation. <i>Doklady Chemistry</i> , 2018, 478, 21-25.	0.2	8
34	Quantum-Chemical Modeling of B32 Complexes with Nitrogen: Endo or Exo?. <i>Russian Journal of Inorganic Chemistry</i> , 2018, 63, 902-905.	0.3	0
35	Valence-tautomeric adducts of Co(II) diketonates based on annelated di-o-quinones: Computer simulation. <i>Russian Journal of Coordination Chemistry/Koordinatsionnaya Khimiya</i> , 2017, 43, 197-205.	0.3	8
36	Quantum chemical study of the adducts of azomethine cobalt complexes with acenaphthene-1,2-diimines. <i>Russian Journal of General Chemistry</i> , 2017, 87, 98-106.	0.3	5

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37	New fluorogenic chemosensors derived from benzimidazole. <i>Chemistry of Heterocyclic Compounds</i> , 2017, 53, 179-185.	0.6	4
38	Light-controlled spin-state-switching rearrangements of transition metal complexes with photochromic ligands. <i>Pure and Applied Chemistry</i> , 2017, 89, 985-1005.	0.9	9
39	Dinuclear adducts of di-o-iminoquinone ligands with Coll diketonates: computational insights into two-step valence tautomeric rearrangements. <i>Journal of Molecular Modeling</i> , 2017, 23, 307.	0.8	2
40	Quantum-chemical study of (Z)-6,8-di-tert-butyl-N-(4-methoxyphenyl)-3-((4-methoxyphenyl)imino)-3H-phenoxazine-2-amine complexation with cobalt bis(chelate)s. <i>Doklady Chemistry</i> , 2017, 476, 215-218.	0.2	2
41	Computational modeling of mixed-ligand cobalt diketonate complexes with pyrene-4,5,9,10-tetraimine. <i>Doklady Chemistry</i> , 2017, 475, 168-172.	0.2	1
42	Quantum chemical modeling of valence tautomeric adducts of Coll bischelates with pyrene-4,5-diimines. <i>Russian Chemical Bulletin</i> , 2017, 66, 208-221.	0.4	10
43	Dual magnetic behavior of Co(II) and Fe(II) bis(chelate) adducts with Di-o-diiminobenzoquinone: Quantum chemical modeling. <i>Russian Journal of Coordination Chemistry/Koordinatsionnaya Khimiya</i> , 2017, 43, 718-726.	0.3	7
44	Quantum chemical study of binuclear adducts of cobalt azomethine complexes with pyrene-4,5,9,10-tetraimine. <i>Russian Chemical Bulletin</i> , 2017, 66, 1543-1549.	0.4	8
45	Protonated paramagnetic redox forms of di-o-quinone bridged with p-phenylene-extended TTF: A EPR spectroscopy study. <i>Beilstein Journal of Organic Chemistry</i> , 2016, 12, 2450-2456.	1.3	10
46	Quantum chemical modeling of magnetically bistable metal coordination compounds. Synchronization of spin crossover, valence tautomerism and charge transfer induced spin transition mechanisms. <i>Dalton Transactions</i> , 2016, 45, 12103-12113.	1.6	25
47	Synthesis and structure of 1-[[3-hydroxybenzo[b]thiophen-2-yl)methylidene]-3-oxo-5-phenyl-1-pyrazolidinium-2-ide. <i>Doklady Chemistry</i> , 2016, 471, 311-313.	0.2	3
48	Benzenoid-quinoid tautomerism of azomethines and their structural analogs 56. Azomethine imines, derivatives of salicylic and 2-hydroxynaphthoic aldehydes. <i>Russian Chemical Bulletin</i> , 2016, 65, 648-653.	0.4	4
49	Computational modeling of spin crossover phenomenon in adducts of iron bis-chelates with o-diiminobenzoquinones. <i>Russian Chemical Bulletin</i> , 2016, 65, 1464-1472.	0.4	7
50	Synthesis and studies of new photochromic spiropyrans containing a formylcoumarin fragment. <i>Russian Chemical Bulletin</i> , 2016, 65, 944-951.	0.4	8
51	Recognition of S $\cdots$ Cl Chalcogen Bonding in Metal-Bound Alkylthiocyanates. <i>Crystal Growth and Design</i> , 2016, 16, 2979-2987.	1.4	22
52	Adducts of tetracoordinate cobalt(II) complexes and 1-(pyridin-2-yl)methanimine: Computational search for valence tautomeric systems. <i>Russian Journal of General Chemistry</i> , 2016, 86, 859-864.	0.3	5
53	Quantum-chemical study of spin crossover in cobalt complexes with an o-benzoquinone ligand. <i>Doklady Chemistry</i> , 2016, 467, 83-87.	0.2	24
54	Supermolecular design: From molecules to solid states. <i>International Journal of Quantum Chemistry</i> , 2016, 116, 259-264.	1.0	5

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55	Quantum chemical simulation of trans- and cis-isomers of bis-chelate azomethine complexes of Ni(II), Pd(II), and Pt(II) with the MN <sub>2</sub> Y <sub>2</sub> (Y = O, S, Se) coordination core. <i>Journal of Structural Chemistry</i> , 2016, 57, 431-436.	0.3	4
56	Photochromism of novel [1]benzothien-2-yl fulgides. <i>Tetrahedron</i> , 2016, 72, 5776-5782.	1.0	9
57	Quantum chemical study of the self-assembly of tetrathiocalix[4]arenes and their oxygen analogs functionalized by hydrazide groups. <i>Russian Chemical Bulletin</i> , 2016, 65, 47-53.	0.4	0
58	Computational modeling of LD LISC and LIESST rearrangements of a Fe(II) complex with phenanthroline modified by photochromic chromene. <i>Doklady Chemistry</i> , 2016, 468, 152-155.	0.2	6
59	Quantum chemical modeling of pyrene-4,5-dione adducts with cobalt diketonates. <i>Computational and Theoretical Chemistry</i> , 2016, 1076, 74-80.	1.1	13
60	EPR spectroscopy study of di-o-quinone bridged by $\pi$ -extended TTF: redox behavior and binding modes as a ligand. <i>New Journal of Chemistry</i> , 2016, 40, 1244-1249.	1.4	7
61	Quantum-chemical study of manganese(II) diketonate adducts with diimine. <i>Doklady Chemistry</i> , 2015, 463, 211-214.	0.2	2
62	Structure and magnetic properties of bis-o-benzosemiquinonato zinc complexes. <i>Polyhedron</i> , 2015, 102, 715-721.	1.0	13
63	Germanium, carbon-germanium, and silicon-germanium triangulenes. <i>Journal of Computational Chemistry</i> , 2015, 36, 2193-2199.	1.5	8
64	Indirect Magnetic Exchange between $\pi$ -Iminosemiquinonate Ligands Controlled by Apical Substituent in Pentacoordinated Gallium(III) Complexes. <i>Inorganic Chemistry</i> , 2015, 54, 6090-6099.	1.9	28
65	The Pearson's HSAB principle in the quantum-chemical model of formation of the MN <sub>2</sub> O <sub>2</sub> or MN <sub>2</sub> S <sub>2</sub> coordination node in the bischelates of Be(II) and Hg(II) with ambidentate ligands based on azomethines and their cyclic analogs. <i>Russian Journal of General Chemistry</i> , 2015, 85, 2629-2633.	0.3	3
66	Quantum chemical study of photomagnetic properties of Ni(II) monochelates with chromenes. <i>Doklady Chemistry</i> , 2015, 462, 118-122.	0.2	4
67	Quantum-chemical model for the formation of the coordination mode structure of Be(II), Ni(II), Pd(II), Pt(II), and Hg(II) bis(chelate) complexes with polydentate azomethine ligands. <i>Russian Journal of Coordination Chemistry/Koordinatsionnaya Khimiya</i> , 2015, 41, 421-427.	0.3	5
68	Theoretical modeling of photocontrolled spin crossover in Fe(II) phenanthroline complexes. <i>Doklady Chemistry</i> , 2015, 460, 5-9.	0.2	10
69	Theoretical modeling of valence tautomeric dinuclear cobalt complexes. Adducts of Co <sup>II</sup> diketonates with cyclic redox-active tetraone ligands. <i>Dalton Transactions</i> , 2015, 44, 17819-17828.	1.6	21
70	Photoswitchable dihetarylethene chemosensors for the selective "naked-eye" detection of fluoride anions. <i>Tetrahedron</i> , 2015, 71, 8817-8822.	1.0	15
71	Tetranuclear Cu(II) and Ni(II) complexes with 1,3,5-triketone ligands: A quantum-chemical simulation of exchange interactions. <i>Russian Journal of Coordination Chemistry/Koordinatsionnaya Khimiya</i> , 2015, 41, 487-495.	0.3	11
72	Quantum-chemical simulation of structure of bischelate Ni(II) complexes based on cyclic analogs of azomethines. <i>Russian Journal of General Chemistry</i> , 2015, 85, 1698-1705.	0.3	4

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73	Adducts of manganese diketonates with redox-active ligands: Computational modeling of valence tautomeric systems. <i>Computational and Theoretical Chemistry</i> , 2015, 1070, 132-142.	1.1	3
74	Intramolecular spin state switching mechanisms of transition metal complexes. <i>Russian Chemical Bulletin</i> , 2015, 64, 475-497.	0.4	28
75	Valence tautomeric dinuclear adducts of Co(II) diketonates with redox-active diquinones for the design of spin qubits: computational modeling. <i>Dalton Transactions</i> , 2015, 44, 1982-1991.	1.6	33
76	Spin crossover in monoadducts of Co(Salen) with pyridine and imidazole: a quantum chemical study. <i>Structural Chemistry</i> , 2014, 25, 1865-1871.	1.0	12
77	Cobalt diketonate adducts with redox-active diiminosuccinonitriles. <i>Mendeleev Communications</i> , 2014, 24, 329-331.	0.6	16
78	Theoretical DFT modeling of the structure of nickel(II) bis(chelate) complexes based on aromatic azomethines. <i>Doklady Chemistry</i> , 2014, 458, 181-184.	0.2	5
79	Sn(IV) complexes with bi- and tridentate phenoxazin-1-one ligands: Synthesis, structure and magnetic properties. <i>Inorganica Chimica Acta</i> , 2014, 418, 66-72.	1.2	17
80	A biradical chelate Zn(II) complex with phenoxazin-1-one ligands. <i>Inorganica Chimica Acta</i> , 2014, 410, 144-149.	1.2	12
81	Self-association of $\alpha$ -tocopherol: a computer simulation. <i>Russian Chemical Bulletin</i> , 2014, 63, 54-59.	0.4	1
82	Compactly Fused Quinone-Extended Tetrathiafulvalene-Quinone Triad: a Redox-Amphoteric Ligand. <i>European Journal of Organic Chemistry</i> , 2014, 2014, 4571-4576.	1.2	19
83	Structure of 4-methyl-N-[2-[2-alkylamino-5-nitrophenyliminomethyl]phenyl]benzenesulfonamides. <i>Crystallography Reports</i> , 2013, 58, 437-441.	0.1	4
84	Hybrid carbon-silicon triangulenes. <i>Doklady Chemistry</i> , 2013, 448, 23-28.	0.2	8
85	Computational modeling of chelating properties of quinoline spiropyrans. <i>Doklady Chemistry</i> , 2013, 453, 263-267.	0.2	5
86	DFT Computational Design of a Ligand-Driven Light-Induced Mechanism for Spin-State Switching. <i>European Journal of Inorganic Chemistry</i> , 2013, 2013, 4203-4219.	1.0	10
87	Adducts of cobalt(II) bis(salicylaldiminates) and redox-active phenoxazin-1-one: synthesis, structure, and magnetic properties. <i>Russian Chemical Bulletin</i> , 2013, 62, 1744-1751.	0.4	34
88	X-ray diffraction, magnetochemical, and quantum chemical study of the structure and properties of binuclear copper(II) complexes. <i>Russian Journal of General Chemistry</i> , 2012, 82, 1770-1776.	0.3	4
89	Effective pH sensors based on 1-(anthracen-9-ylmethyl)-1H-benzimidazol-2-amine. <i>Chemistry of Heterocyclic Compounds</i> , 2012, 47, 1230-1236.	0.6	6
90	Valence tautomeric complexes of cobalt diketonates with Diimines: A quantum-chemical study. <i>Doklady Chemistry</i> , 2011, 440, 289-293.	0.2	15

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91	Valence tautomerism of a manganese complex with phenoxybenzoquinone imine ligands: A quantum-chemical study. Doklady Chemistry, 2011, 441, 365-370.	0.2	12
92	Synthesis, Molecular and Electronic Structures of Six-Coordinate Transition Metal (Mn, Fe, Co, Ni) Tj ETQq0 0 0 rgBT /Overlock 10 Tf 50 2011, 50, 7022-7032.	1.9	41
93	Carbon and silicon triangulenes: searching for molecular magnets. Russian Chemical Bulletin, 2011, 60, 1517-1524.	0.4	11
94	Structure and stability of the mixed polymolecular complexes of nitrogen and carbon nonoxide: A quantum chemical study. Russian Journal of General Chemistry, 2011, 81, 807-818.	0.3	5
95	Structure of indoline spiropyran containing a fused coumarin fragment: Quantum-chemical investigation. Russian Journal of Organic Chemistry, 2011, 47, 1742-1745.	0.3	3
96	Quantum-chemical modeling of metal coordination compounds with photoswitchable magnetic properties controlled by ligand rearrangements. Theoretical and Experimental Chemistry, 2011, 46, 363-370.	0.2	7
97	Metal complexes with azomethines containing the isomeric E-Z azo fragments. Russian Journal of Coordination Chemistry/Koordinatsionnaya Khimiya, 2010, 36, 479-489.	0.3	17
98	Quantum chemical study of pyridine addition to Ni(II) $\hat{1}^2$ -diketonate complexes. Russian Journal of Coordination Chemistry/Koordinatsionnaya Khimiya, 2010, 36, 597-604.	0.3	3
99	Quantum chemical calculations for the geometry and intramolecular rearrangements of a model nickel(ii) o-semiquinone complex with a PCP pincer ligand. Russian Chemical Bulletin, 2010, 59, 1110-1115.	0.4	3
100	Sandwich Compounds of Transition Metals with Cyclopolyenes and Isolobal Boron Analogues. Chemistry - A European Journal, 2010, 16, 2272-2281.	1.7	15
101	Quantum-chemical study of valence tautomerism of a cobalt complex with phenoxybenzoquinone imine. Doklady Chemistry, 2010, 435, 319-323.	0.2	27
102	A Quantum Chemical Study of Bis-(iminoquinonephenolate) Zn(II) Complexes. Journal of Physical Chemistry A, 2010, 114, 7780-7785.	1.1	36
103	Effect of ligand environment on the mechanism of enantiomerization of Bell, ZnII, and CuII bischelate complexes. Russian Chemical Bulletin, 2009, 58, 513-521.	0.4	2
104	$\hat{1}^2$ -Complexes of transition metal tricarbonyls with cyclopolyenes and their boron analogs. Russian Chemical Bulletin, 2009, 58, 691-705.	0.4	5
105	Novel structural motif for stabilization of polyacene systems. Doklady Physical Chemistry, 2009, 425, 77-80.	0.2	0
106	Theoretical modeling of enantiomerization mechanisms of tetrahedral bis-( $\hat{1}^2$ -diiminato) Ni(II) complexes. Computational and Theoretical Chemistry, 2009, 895, 138-141.	1.5	20
107	Theoretical modeling of the molecular and crystal structures and a square-planar to tetrahedral conformational rearrangement of trans-planar bis(N-methylsalicylaldiminato)nickel(II). Mendeleev Communications, 2009, 19, 64-66.	0.6	4
108	Influence of structural factors on the magnetic properties of the binuclear copper complexes with salicylaldehyde hydrazone and bis(hydrazone)-2,6-diformylphenol: Quantum-chemical calculations. Russian Journal of Coordination Chemistry/Koordinatsionnaya Khimiya, 2009, 35, 616-620.	0.3	14

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109	Quantum-chemical simulation of the mechanisms of stereoisomerism of tetracoordinated Ni(II) complexes. Russian Journal of General Chemistry, 2009, 79, 1793-1801.	0.3	2
110	Quantum-chemical study of intramolecular spin-forbidden rearrangements of the transition metal chelate complexes. Russian Journal of General Chemistry, 2009, 79, 2792-2806.	0.3	8
111	Cooperative effects in polymolecular nitrogen clusters. Russian Chemical Bulletin, 2008, 57, 2037-2044.	0.4	3
112	Theoretical modeling of the square-planar to tetrahedral isomerization of bis-chelate nickel(II) complexes. Chemical Physics Letters, 2008, 459, 27-32.	1.2	27
113	Tautomeric crown-containing chemosensors for alkali-earth metal cations. Tetrahedron, 2008, 64, 3160-3167.	1.0	33
114	Quantum-chemical study of endohedral fullerenes. Russian Journal of General Chemistry, 2008, 78, 793-810.	0.3	12
115	Steric and electronic structure of complexes of pyrylium and thiopyrylium cations with borabenzene anion. Russian Journal of General Chemistry, 2008, 78, 1354-1360.	0.3	0
116	Extended organoboron structures containing several planar tetracoordinate carbon atoms. Doklady Chemistry, 2008, 419, 101-107.	0.2	9
117	Hypercoordination of first-row elements in heteroanalogues of prismanes and propellanes. Doklady Chemistry, 2008, 422, 255-259.	0.2	3
118	Hexacoordinated carbon and nitrogen atoms in extended organoboron cage structures. Doklady Chemistry, 2007, 416, 235-240.	0.2	2
119	Hydrogen bond in FHâ€¦FM (M = Li, Na, K) dimers: Nonempirical calculations. Russian Journal of Physical Chemistry A, 2007, 81, 1100-1103.	0.1	3
120	Effect of the counterion on the steric and electronic structure of pyrylium cation. Russian Journal of General Chemistry, 2007, 77, 1373-1385.	0.3	9
121	Hypercoordinated carbon in endohedral hydrocarbon cage complexes C@C20H 20 4âˆ™ and C@C20H20 Â·Li4. Doklady Chemistry, 2006, 407, 47-50.	0.2	11
122	Geometry and electronic structure of lithium sandwich complexes [(OLIGOCENE)2] n+1Li n+1: A quantum-chemical study. Doklady Chemistry, 2006, 409, 113-116.	0.2	2
123	Photoisomerization of quinolin-2-yl derivatives of Î²-tropolone. Russian Chemical Bulletin, 2006, 55, 484-491.	0.4	3
124	New method for the synthesis of Î²-tropolones: Structures of condensation products of o-quinones with 2-methylquinolines and the mechanism of their formation. Russian Chemical Bulletin, 2006, 55, 2032-2055.	0.4	22
125	Sandwich compounds with central hypercoordinate carbon, nitrogen, and oxygen: A quantum-chemical study. Heteroatom Chemistry, 2006, 17, 464-474.	0.4	12
126	Synthesis and structure of N-arylimines of Î²-tellurocyclohexenals with the intramolecular coordination Nâ†’Te bonds. Journal of Organometallic Chemistry, 2005, 690, 103-116.	0.8	22



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127	Octacoordinated Carbon in a Boron-Carbon Cage. Doklady Chemistry, 2005, 404, 193-198.	0.2	10
128	Planar Four-Coordinate Carbon in Star-Like Perithioannulenes $C_nLi_n$ ( $n = 3\text{--}6$ ). Russian Journal of Organic Chemistry, 2005, 41, 1289-1295.	0.3	18
129	Planar and Pyramidal Tetracoordinate Carbon in Organoboron Compounds. Journal of Organic Chemistry, 2005, 70, 6693-6704.	1.7	56
130	A hydrocarbon dication with nonplanar hexacoordinated carbon. Mendeleev Communications, 2004, 14, 47-48.	0.6	15
131	Spiroconjugation Energy of Spiroheterocyclic Structures. Doklady Chemistry, 2004, 396, 99-102.	0.2	3
132	Structure and stability of closo-hexaboranes and their heteroanalogs. Russian Chemical Bulletin, 2004, 53, 1159-1167.	0.4	12
133	Poly[n]prismanes: A Family of Stable Cage Structures with Half-Planar Carbon Centers.. ChemInform, 2004, 35, no.	0.1	0
134	Title is missing!. Russian Chemical Bulletin, 2003, 52, 519-525.	0.4	2
135	Double $\pi$ - and $\delta$ -hydrogen bonding in formic acid complexes with pyrrole and imidazole: an ab initio and density functional theory study. Mendeleev Communications, 2003, 13, 207-209.	0.6	0
136	Poly[n]prismanes: A Family of Stable Cage Structures with Half-Planar Carbon Centers. Journal of Organic Chemistry, 2003, 68, 8588-8594.	1.7	64
137	Heptacoordinated Carbon and Nitrogen in a Planar Boron Ring. Doklady Chemistry, 2002, 382, 41-45.	0.2	43
138	Induced aromaticity. Russian Chemical Bulletin, 2001, 50, 2325-2335.	0.4	8
139	Stabilization of the glycine zwitterionic form by complexation with $Na^+$ and $Cl^-$ : an ab initio study. Mendeleev Communications, 2000, 10, 43-44.	0.6	6
140	Pathways of the reactions of nucleophilic addition of $H_2O$ and $HF$ molecules to formaldehyde in the gas phase and in the complex with formic acid: ab initio calculations. Russian Chemical Bulletin, 1998, 47, 2078-2086.	0.4	4
141	Preparation and reactivity of metal-containing monomers. Russian Chemical Bulletin, 1993, 42, 66-70.	0.4	1
142	Preparation and reactivity of metal-containing monomers. 23. Transition metal complexes of methacryloylacetone. Bulletin of the Russian Academy of Sciences Division of Chemical Science, 1992, 41, 545-548.	0.0	1
143	Metal-containing monomers. Part 2*. Metal chelate monomers based on 1-phenyl-4-methylpent-4-en-1,3-dione. Transition Metal Chemistry, 1992, 17, 458-463.	0.7	7
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