

Oleg V Shishkin

List of Publications by Year in descending order

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2,718
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159585

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46
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125
all docs

125
docs citations

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2361
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#	ARTICLE	IF	CITATIONS
1	Microwave-Assisted Three-Component Synthesis of 7-Aryl-2-alkylthio-4,7-dihydro-1,2,4-triazolo[1,5-a]-pyrimidine-6-carboxamides and Their Selective Reduction. <i>ACS Combinatorial Science</i> , 2006, 8, 427-434.	3.3	114
2	Synthesis and imido-group exchange reactions of tert-butylimidotitanium complexes. <i>Journal of the Chemical Society Dalton Transactions</i> , 1997, , 1549-1558.	1.1	109
3	OH \cdots F and halogen \cdots F interactions as driving forces in the crystal organisations of tri-bromo and tri-iodo trityl alcohols. <i>CrystEngComm</i> , 2008, 10, 715.	2.6	87
4	Electron density distribution in stacked benzene dimers: A new approach towards the estimation of stacking interaction energies. <i>Journal of Chemical Physics</i> , 2005, 122, 144104.	3.0	85
5	Diastereoselective enolate chemistry using atropisomeric amide. <i>Tetrahedron Letters</i> , 1996, 37, 7607-7610.	1.4	82
6	One-Pot, Multicomponent Route to Pyrazoloquinolizinones. <i>Organic Letters</i> , 2007, 9, 1691-1694.	4.6	80
7	Supramolecular architecture of crystals of fused hydrocarbons based on topology of intermolecular interactions. <i>CrystEngComm</i> , 2012, 14, 1795.	2.6	80
8	Cyclocondensation reactions of 5-aminopyrazoles, pyruvic acids and aldehydes. Multicomponent approaches to pyrazolopyridines and related products. <i>Tetrahedron</i> , 2007, 63, 1229-1242.	1.9	72
9	Role of supramolecular synthons in the formation of the supramolecular architecture of molecular crystals revisited from an energetic viewpoint. <i>Physical Chemistry Chemical Physics</i> , 2014, 16, 6773.	2.8	67
10	Structural nonrigidity of nucleic acid bases. Post-Hartree-Fock ab initio study. <i>International Journal of Quantum Chemistry</i> , 2000, 80, 1116-1124.	2.0	62
11	Intramolecular Hydrogen Bonds in Canonical 2-Deoxyribonucleotides: An Atoms in Molecules Study. <i>Journal of Physical Chemistry B</i> , 2006, 110, 4413-4422.	2.6	62
12	Multicomponent cyclocondensation reactions of aminoazoles, arylpyruvic acids and aldehydes with controlled chemoselectivity. <i>Tetrahedron</i> , 2008, 64, 11041-11049.	1.9	59
13	Unexpected alternative direction of a Biginelli-like multicomponent reaction with 3-amino-1,2,4-triazole as the urea component. <i>Tetrahedron Letters</i> , 2010, 51, 2095-2098.	1.4	50
14	Conformational Analysis of Canonical 2-Deoxyribonucleotides. 2. Purine Nucleotides. <i>Journal of Biomolecular Structure and Dynamics</i> , 2004, 22, 227-243.	3.5	48
15	Cyclobutane-Derived Diamines: Synthesis and Molecular Structure. <i>Journal of Organic Chemistry</i> , 2010, 75, 5941-5952.	3.2	48
16	Dotting the i's in three-component Biginelli-like condensations using 3-amino-1,2,4-triazole as a 1,3-binucleophile. <i>RSC Advances</i> , 2012, 2, 6719.	3.6	48
17	Three-Component Procedure for the Synthesis of 5-Aryl-5,8-dihydroazolo[1,5-a]pyrimidine-7-carboxylic Acids. <i>Synthesis</i> , 2005, 2005, 2597-2601.	2.3	46
18	Heterobinuclear Zn \cdots Ln and Ni \cdots Ln Complexes with Schiff-Base and Carbacylamidophosphate Ligands: Synthesis, Crystal Structures, and Catalytic Activity. <i>European Journal of Inorganic Chemistry</i> , 2014, 2014, 3720-3730.	2.0	46

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19	Conformational flexibility of pyrimidine ring in adenine and related compounds. <i>Chemical Physics Letters</i> , 2000, 330, 603-611.	2.6	43
20	A density functional theory study of vibrational coupling between ribose and base rings of nucleic acids with ribosyl guanosine as a model system. <i>Journal of Chemical Physics</i> , 2000, 113, 5986-5990.	3.0	42
21	Structural non-rigidity of six-membered aromatic rings. <i>Journal of Molecular Structure</i> , 2002, 616, 159-166.	3.6	42
22	Conformational Analysis of Canonical 2-Deoxyribonucleotides. 1. Pyrimidine Nucleotides. <i>Journal of Biomolecular Structure and Dynamics</i> , 2004, 21, 537-553.	3.5	42
23	Layered crystal structure of bicyclic aziridines as revealed by analysis of intermolecular interactions energy. <i>CrystEngComm</i> , 2010, 12, 1816.	2.6	39
24	Switchable selectivity in multicomponent heterocyclizations of acetoacetamides, aldehydes, and 3-amino-1,2,4-triazoles/5-aminopyrazoles. <i>Tetrahedron</i> , 2011, 67, 9389-9400.	1.9	39
25	Imidotitanium Tris(pyrazolyl)hydroborates: Synthesis, Solution Dynamics, and Solid-State Structure. <i>Inorganic Chemistry</i> , 1996, 35, 1006-1012.	4.0	38
26	Analysis of the crystal structure of two polymorphic modifications of 3,4-diamino-1,2,4-triazole based on the energy of the intermolecular interactions. <i>CrystEngComm</i> , 2010, 12, 909-916.	2.6	37
27	Columnar supramolecular architecture of crystals of 2-(4-iodophenyl)-1,10-phenanthroline derived from values of intermolecular interaction energy. <i>CrystEngComm</i> , 2011, 13, 800-805.	2.6	36
28	Azatriquinanes: Synthesis, Structure, and Reactivity. <i>Journal of Organic Chemistry</i> , 1998, 63, 6016-6020.	3.2	34
29	Study of the Chemoselectivity of Multicomponent Heterocyclizations Involving 3-Amino-1,2,4-triazole and Pyruvic Acids as Key Reagents, and Biological Activity of the Reaction Products. <i>European Journal of Organic Chemistry</i> , 2015, 2015, 4481-4492.	2.4	33
30	Diversity oriented heterocyclizations of pyruvic acids, aldehydes and 5-amino-N-aryl-1H-pyrazole-4-carboxamides: catalytic and temperature control of chemoselectivity. <i>Molecular Diversity</i> , 2010, 14, 523-531.	3.9	31
31	Structure and hydrogen bonding in polyhydrated complexes of guanine. <i>Structural Chemistry</i> , 2008, 19, 171-180.	2.0	30
32	Novel type of mixed $\text{O} \cdots \text{H} \cdots \text{N} / \text{O} \cdots \text{H} \cdots \text{N}$ hydrogen bonds: monohydrate of pyridine. <i>Structural Chemistry</i> , 2009, 20, 37-41.	2.0	30
33	Features of switchable multicomponent heterocyclizations of salicylic aldehydes and 5-aminopyrazoles with pyruvic acids and antimicrobial activity of the reaction products. <i>Tetrahedron</i> , 2013, 69, 9261-9269.	1.9	30
34	Supramolecular architecture of molecular crystals possessing shearing mechanical properties: columns versus layers. <i>CrystEngComm</i> , 2013, 15, 160-167.	2.6	29
35	Hydration of nucleic acid bases: a Car Parrinello molecular dynamics approach. <i>Physical Chemistry Chemical Physics</i> , 2010, 12, 3363.	2.8	28
36	Insights into the crystal packing of phosphorylporphyrins based on the topology of their intermolecular interaction energies. <i>CrystEngComm</i> , 2014, 16, 10428-10438.	2.6	28

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37	Investigation of topology of intermolecular interactions in the benzene-acetylene co-crystal by different theoretical methods. <i>Structural Chemistry</i> , 2014, 25, 1547-1552.	2.0	28
38	An Entry into Hexahydro-2H-thieno[2,3-c]pyrrole 1,1-Dioxide Derivatives. <i>Journal of Organic Chemistry</i> , 2011, 76, 7010-7016.	3.2	26
39	Low-melting molecular complexes. Halogen bonds in molecular complexes of bromoform. <i>CrystEngComm</i> , 2012, 14, 8222.	2.6	26
40	The C-Cl interactions inside supramolecular nanotubes of hexaethynylhexamethoxy[6]pericyclyne. <i>Physical Chemistry Chemical Physics</i> , 2011, 13, 6837.	2.8	24
41	Exploiting morph-DAST mediated ring-expansion of substituted cyclic β^2 -amino alcohols for the preparation of cyclic fluorinated amino acids. Synthesis of 5-fluoromethylproline and 5-fluoropipecolic acid. <i>Tetrahedron</i> , 2011, 67, 3091-3097.	1.9	24
42	Dalton communications. Exchange of organoimido groups at a mononuclear titanium centre and a crystallographic evaluation of the relative structural influences of the NBut, NC ₆ H ₄ Me-4 and NC ₆ H ₄ NO ₂ -4 ligands. <i>Journal of the Chemical Society Dalton Transactions</i> , 1995, , 3743.	1.1	23
43	Dependence of Deformability of Geometries and Characteristics of Intramolecular Hydrogen Bonds in Canonical 2-Deoxyribonucleotides on DNA Conformations. <i>Journal of Biomolecular Structure and Dynamics</i> , 2009, 26, 653-661.	3.5	23
44	Stable N-Heterocyclic Carbenes: Alkyl-N-phosphanylbenzimidazol-2-ylidenes. <i>European Journal of Organic Chemistry</i> , 2012, 2012, 4018-4033.	2.4	22
45	Intra- and Intermolecular Interactions in the Crystals of 3,4-Diamino-1,2,4-triazole and Its 5-Methyl Derivative. Experimental and Theoretical Investigations of Charge Density Distribution. <i>Journal of Physical Chemistry A</i> , 2011, 115, 8550-8562.	2.5	21
46	Influence of Deuteration and Fluorination on the Supramolecular Architecture of Pyridine N-Oxide Crystals. <i>ChemPhysChem</i> , 2013, 14, 847-856.	2.1	21
47	Multicomponent Cyclocondensations of β -Ketosulfones with Aldehydes and Aminoazole Building Blocks. <i>Heterocycles</i> , 2007, 73, 469.	0.7	20
48	Synthesis of 2- and 3-trifluoromethylmorpholines: useful building blocks for drug discovery. <i>Tetrahedron</i> , 2013, 69, 3796-3804.	1.9	20
49	Remarkably strong polarization of amidine fragment in the crystals of 1-imino-1H-isoindol-3-amine. <i>Structural Chemistry</i> , 2013, 24, 1089-1097.	2.0	20
50	Role of CHF ₂ - and CF ₃ -Substituents on Molecular Arrangement in the Solid State: Experimental and Theoretical Crystal Structure Analysis of CH ₃ /CHF ₂ /CF ₃ -Substituted Benzene. <i>Crystal Growth and Design</i> , 2014, 14, 3124-3130.	3.0	20
51	3-Benzyl-3-azabicyclo[3.1.1]heptan-6-one: A Promising Building Block for Medicinal Chemistry. <i>Organic Letters</i> , 2010, 12, 4372-4375.	4.6	19
52	Partially hydrogenated 2-amino[1,2,4]triazolo[1,5-a]pyrimidines as synthons for the preparation of polycondensed heterocycles: reaction with chlorocarboxylic acid chlorides. <i>Tetrahedron</i> , 2014, 70, 684-701.	1.9	19
53	Dibutylphosphinoylmethoxythiacalix[4]arenes. Synthesis, structure, americium, europium and technetium extraction. <i>Supramolecular Chemistry</i> , 2014, 26, 864-872.	1.2	19
54	Azatriquinane, azatriquinacene, and a remarkable dimerization product. <i>Tetrahedron Letters</i> , 1996, 37, 131-134.	1.4	18

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55	Reactions of 4,7-dihydro-1,2,4-triazolo[1,5-a]pyrimidines with α,β -unsaturated carbonyl compounds. <i>Journal of Heterocyclic Chemistry</i> , 2003, 40, 1081-1086.	2.6	18
56	Cyclecondensation of 3-amino-1,2,4-triazole with substituted methyl cinnamates. <i>Journal of Heterocyclic Chemistry</i> , 1999, 36, 205-208.	2.6	17
57	Reactions of <i>N</i> -arylmaleimides with 3-amino-1,2,4-triazole and 2-aminobenzimidazole. <i>Journal of Heterocyclic Chemistry</i> , 2011, 48, 888-895.	2.6	17
58	Controlled Switching of Multicomponent Heterocyclizations of 5-Amino- <i>N</i> -arylpiperazine-4-carboxamides, 1,3-Cyclohexanediones, and Aldehydes. <i>European Journal of Organic Chemistry</i> , 2012, 2012, 5515-5524.	2.6	17
59	New tricks of well-known aminoazoles in isocyanide-based multicomponent reactions and antibacterial activity of the compounds synthesized. <i>Beilstein Journal of Organic Chemistry</i> , 2017, 13, 1050-1063.	2.2	17
60	Opposite charges assisted extra strong C-H \cdots O hydrogen bond in protonated 2-deoxyadenosine monophosphate. <i>Chemical Physics Letters</i> , 2008, 452, 198-205.	2.6	14
61	Conjugation and Hyperconjugation in Conformational Analysis of Cyclohexene Derivatives Containing an Exocyclic Double Bond. <i>Journal of Physical Chemistry A</i> , 2008, 112, 7080-7089.	2.5	14
62	Role of different molecular fragments in formation of the supramolecular architecture of the crystal of 1,1-dioxo-tetrahydro-1H-6-thiopyran-3-one. <i>CrystEngComm</i> , 2012, 14, 8698.	2.6	14
63	A Convenient Route to α -alkyl- α -trifluoromethyl-1,2,3-triazole-4-carboxylic Acids Employing a Diazo Transfer Reaction. <i>European Journal of Organic Chemistry</i> , 2013, 2013, 2891-2897.	2.4	14
64	Synthesis of structurally constrained 4-quinazolinone derivatives with a tetrahedral C-2 atom present in three rings. <i>Tetrahedron</i> , 2009, 65, 8582-8586.	1.9	13
65	The unexpected influence of aryl substituents in <i>N</i> -aryl-3-oxobutanamides on the behavior of their multicomponent reactions with 5-amino-3-methylisoxazole and salicylaldehyde. <i>Beilstein Journal of Organic Chemistry</i> , 2014, 10, 3019-3030.	2.2	13
66	Dynamical Nonplanarity of Benzene. Evidences from the Car-Parrinello Molecular Dynamics Study. <i>Journal of Physical Chemistry Letters</i> , 2011, 2, 2881-2884.	4.6	12
67	Nucleic Acid Bases in Anionic 2-Deoxyribonucleotides: A DFT/B3LYP Study of Structures, Relative Stability, and Proton Affinities. <i>Journal of Physical Chemistry B</i> , 2013, 117, 2841-2849.	2.6	11
68	Acceptor properties of amino groups in aminobenzene crystals: study from the energetic viewpoint. <i>CrystEngComm</i> , 2017, 19, 6274-6288.	2.6	11
69	Estimating stacking interaction energy using atom in molecules properties: Homodimers of benzene and pyridine. <i>International Journal of Quantum Chemistry</i> , 2012, 112, 3008-3017.	2.0	10
70	1-Amino-4,4-difluorocyclohexanecarboxylic acid as a promising building block for drug discovery: design, synthesis and characterization. <i>Tetrahedron</i> , 2013, 69, 4066-4075.	1.9	10
71	Interactions of Water with Mono- and Diamino Derivatives of N,N-Dimethyluracil. <i>Journal of Physical Chemistry A</i> , 2002, 106, 7828-7833.	2.5	9
72	A novel synthesis and transformations of isothiochroman 2,2-dioxide. <i>Tetrahedron Letters</i> , 2012, 53, 4296-4299.	1.4	9

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73	Novel transformations of 1H-isothiochromen-4(3H)-one 2,2-dioxide. Monatshefte für Chemie, 2013, 144, 263-271.	1.8	9
74	Isotopic Transformation Principle in Molecular Crystals. Analysis of Supramolecular Architecture of Fluorinated Benzenes and Pyridines. Crystal Growth and Design, 2018, 18, 4445-4448.	3.0	9
75	Heterocyclization of 6-hydroxyimino-6,7-dihydro-1,2,4-triazolo[1,5-a]pyrimidines into 1,2,4-triazolo[1,5-a]pyrimido[5,4-b]- and -[5,6-b]indoles. Mendeleev Communications, 2006, 16, 280-282.	1.6	8
76	Molecular and crystal structure of crown ethers containing biphenyl fragment. Journal of Molecular Structure, 2007, 832, 199-208.	3.6	8
77	Intramolecular Cyclization of <i>N</i> -Arylphosphinimidic Isocyanates – Novel Approach to a 4a,8a-dihydro-1,3,2λ ⁵ -benzodiazaphosphinin-4(3H)-one System. European Journal of Inorganic Chemistry, 2008, 2008, 3348-3352.		8
78	2-Substituted-Isoindoles: A Novel Synthetic Route and a Study of the Diels–Alder and Michael Reactions. Journal of Chemical Research, 2011, 35, 615-618.	1.3	8
79	Formylation of 4,7-dihydro-1,2,4-triazolo[1,5-a]pyrimidines Using Vilsmeier–Haack Conditions. Journal of Heterocyclic Chemistry, 2012, 49, 1019-1025.	2.6	8
80	3,3'-Dinitrophenolsulphonephthalein: an acid-base indicator dye with unusual properties. Coloration Technology, 2017, 133, 135-144.	1.5	8
81	Influence of substituents on the acceptor properties of the amino groups in the diaminobenzene analogues. CrystEngComm, 2017, 19, 7162-7176.	2.6	8
82	An alternative approach to the synthesis of 5H-chromeno[4,3-b]pyridin-5-one system using the cleavage of 5H,9H-pyrano[2',3':5,6]chromeno[4,3-b]pyridine-5,9-diones with binucleophiles. Chemistry of Heterocyclic Compounds, 2018, 54, 96-99.	1.2	8
83	The structure of cyanine dyes of tetrazoloisoindole row. 1. Bis-(1-Methyltetrazolo[5,1-a]isoindole-5)monomethyncyanine perchlorate. Journal of Molecular Structure, 2004, 707, 193-198.	3.6	7
84	Synthesis and crystal structure determination of 2,6-di-tert-butyl-4-(2,4,6-triphenylpyridinium-1-yl)phenolate and its corresponding perchlorate salt. Dyes and Pigments, 2012, 92, 1394-1399.	3.7	7
85	Upper-rim calixarene phosphines consisting of multiple lower-rim OH functional groups: synthesis and characterisation. Supramolecular Chemistry, 2014, 26, 825-835.	1.2	7
86	Synthesis and characterization of sulfolane-based amino alcohols: A combined experimental and computational study. Journal of Molecular Structure, 2018, 1157, 149-158.	3.6	7
87	Crystal, molecular structure and tautomerism of (5-methyl-1H-[1,2,4]triazol-3-ylsulfanyl)-acetic acid. Structural Chemistry, 2008, 19, 407-412.	2.0	6
88	Rotation around the glycosidic bond as driving force of proton transfer in protonated 2'-deoxyriboadenosine monophosphate (dAMP). Chemical Physics Letters, 2010, 490, 221-225.	2.6	6
89	Hexakis(dimethylformamide- \hat{O})manganese(II) $\hat{1}/4$ -oxido-bis[trichloridoferrate(III)]. Acta Crystallographica Section E: Structure Reports Online, 2011, 67, m1563-m1564.	0.2	6
90	Reaction of 1-Ethoxyisoindole with Maleimide and Its Derivatives. Heterocycles, 2012, 85, 1671.	0.7	6

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91	Catalysis by Lithium Perchlorate Enables Double-Conjugate Addition of Electron-Deficient Maleimides to 2-Aminopyridines and 2-Aminothiazoles. <i>Synthetic Communications</i> , 2012, 42, 3304-3310.	2.1	5
92	Direct synthesis and properties of monomeric and dimeric Mn(III) salen complexes tuned by tetrahalocadmate anions. <i>Inorganic Chemistry Communication</i> , 2012, 20, 282-285.	3.9	5
93	Conformational behaviour of peptides containing a 2-pyrrolidinemethanesulfonic acid (2PyMS) residue. <i>Organic and Biomolecular Chemistry</i> , 2013, 11, 975-983.	2.8	5
94	Entropy versus aromaticity in the conformational dynamics of aromatic rings. <i>Journal of Molecular Modeling</i> , 2013, 19, 4073-4077.	1.8	5
95	Conformational behavior of peptides containing residues of 3-azetidinesulfonic (3AzeS) and 4-piperidinemethanesulfonic (4PiMS) acids. <i>Tetrahedron: Asymmetry</i> , 2014, 25, 229-237.	1.8	5
96	Revisiting tetranitrophenolsulfonephthalein. <i>Coloration Technology</i> , 2015, 131, 236-244.	1.5	5
97	Reaction of 2-Hetaryl-2-(tetrahydro-2-furanyliden)acetonitriles with 1,3-N,N-Binucleophiles. <i>Synlett</i> , 2012, 23, 2063-2068.	1.8	4
98	Binding properties and self-assembly of C _{2v} -symmetrical resorcin[4]arene tetrabenzoates. <i>Tetrahedron</i> , 2012, 68, 9429-9434.	1.9	4
99	Molecular clips based on diphenylglycoluril and benzocrown ethers: promising complexing agents for the alkali metal cations. <i>Journal of Inclusion Phenomena and Macrocyclic Chemistry</i> , 2014, 79, 343-348.	1.6	4
100	A Convenient Approach to N-(Di-tert-butylphosphanyl)- and N-(Di-tert-butylphosphoroselenoyl)formamidinium Salts: Carbene Precursors. <i>European Journal of Inorganic Chemistry</i> , 2014, 2014, 1192-1203.	2.0	4
101	Cyclic α -amino acids as precursors for synthesis of 2-amino-3-hetarylpyrrolin-4-ones and their spiro derivatives. <i>Monatshefte für Chemie</i> , 2012, 143, 779-789.	1.8	3
102	Environment-induced stabilization of hydrogen-bonded dimers in crystal of lysine (5-methyl-1H-[1,2,4]triazol-3-ylsulfanyl)-acetate. <i>Structural Chemistry</i> , 2012, 23, 581-586.	2.0	3
103	Supramolecular Architecture of Substituted Tetraphenylcarbo-benzenes from the Energetic Viewpoint. <i>ChemPhysChem</i> , 2017, 18, 2499-2508.	2.1	3
104	Functionalised Cookson's Diketones in Chlorosulfonic Acid: Towards Polysubstituted C ₃ -Trishomocubanes. <i>Journal of Chemical Research</i> , 2017, 41, 718-721.	1.3	3
105	Halogen... interactions in the complexes of fluorenonophane with haloforms. <i>Structural Chemistry</i> , 2022, 33, 257-266.	2.0	3
106	Bis[2-(2-hydroxyethyl)(methyl)amino]ethanolato}bis[3-N-methyl-2,2-azanedioldiethanolato}tetrakis(thiocyanato-N)dichromedimethylformamide tetrasolvate. <i>Acta Crystallographica Section E: Structure Reports Online</i> , 2011, 67, m1864-m1865.	0.2	2
107	Synthesis of condensed tetrahydroimidazo[1,2-a]quinazoline-1,5-dione derivatives. <i>Tetrahedron</i> , 2012, 68, 3098-3102.	1.9	2
108	Complexation of thiacalix[4]arene methylphosphonic and sulphonic acids with amino acids. <i>Supramolecular Chemistry</i> , 2013, 25, 263-268.	1.2	2

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109	Synthesis and structural investigation of new isothiochromen-4-one 2,2-dioxide derivatives. Structural Chemistry, 2016, 27, 273-280.	2.0	2
110	Weak but strong: role of weak C-H...X (X=O, N) hydrogen bonds in organization of crystals of (1S,2S,3S,4R,5R,8S)-diethyl 2,4-dicyano-3-(furan-2-yl)-8-morpholino-6-oxobicyclo[3.2.1]octane-2,4-dicarboxylate. Structural Chemistry, 2016, 27, 315-321.	2.0	2
111	Functionalized organic frameworks explored as second order NLO agents. Journal of Chemical Sciences, 2016, 128, 297-309.	1.5	2
112	Quantum delocalization of benzene in the ring puckering coordinates. International Journal of Quantum Chemistry, 2014, 114, 534-542.	2.0	1
113	Unexpected synthesis of pyrazolone derivatives. Tetrahedron, 2015, 71, 1283-1286.	1.9	1
114	Molecular Structures, Relative Stability, and Proton Affinities of Nucleotides: Broad View and Novel Findings. Challenges and Advances in Computational Chemistry and Physics, 2014, , 137-180.	0.6	0
115	Synthesis of spiro 2-(5-amino-2,3-dihydro-3-oxopyrrol-4-yl)-1,3-dialkylbenzimidazolium chlorides. Monatshefte für Chemie, 2015, 146, 931-939.	1.8	0