# Borys OÅ›miaÅ‚owski

List of Publications by Year
in descending order



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|---|---|---|---|
| 109 papers | 1,964 citations | 304602 22 h-index | 289141 40 g-index |
| 115 all docs | 115 docs citations | 115 times ranked | 1808 citing authors |

| # | Article | IF | Citations |
|---|---|---|---|
| 1 | Tautomeric Equilibria in Relation to Pi-Electron Delocalization. Chemical Reviews, 2005, 105, 3561-3612. | 23.0 | 298 |
| 2 | GIAO/DFT calculated chemical shifts of tautomeric species. 2-Phenacylpyridines and (Z)-2-(2-hydroxy-2-phenylvinyl)pyridines. Magnetic Resonance in Chemistry, 2001, 39, 334-340. | 1.1 | 166 |
| 3 | The experimental studies on the determination of the ground and excited state dipole moments of some hemicyanine dyes. Spectrochimica Acta - Part A: Molecular and Biomolecular Spectroscopy, 2006, 63, 524-531. | 2.0 | 75 |
| 4 | Substituent and temperature controlled tautomerism: multinuclear magnetic resonance, X-ray, and theoretical studies on 2-phenacylquinolines. Perkin Transactions II RSC, 2000, , 1259-1266. | 1.1 | 68 |
| 5 | Substituent and temperature controlled tautomerism of 2-phenacylpyridine: the hydrogen bond as a configurational lock of (Zâ€Š)-2-(2-hydroxy-2-phenylvinyl)pyridine. Perkin Transactions II RSC, 2000, , 2185-2191. | 1.1 | 60 |
| 6 | Tautomeric Equilibria and Pi Electron Delocalization for Some MonohydroxyarenesQuantum Chemical Studies. Journal of Organic Chemistry, 2006, 71, 3727-3736. | 1.7 | 48 |
| 7 | Self-Organization of 2-Acylaminopyridines in the Solid State and in Solution. Journal of Physical Chemistry A, 2010, 114, 10421-10426. | 1.1 | 48 |
| 8 | Influence of Substituent and Benzoannulation on Photophysical Properties of 1-Benzoylmethyleneisoquinoline Difluoroborates. Journal of Organic Chemistry, 2015, 80, 2072-2080. | 1.7 | 47 |
| 9 | The Influence of the Ï€-Conjugated Spacer on Photophysical Properties of Difluoroboranyls Derived from Amides Carrying a Donor Group. Journal of Organic Chemistry, 2016, 81, 2280-2292. | 1.7 | 45 |
| 10 | Substituent effects on the photophysical properties of fluorescent 2-benzoylmethylenequinoline difluoroboranes: A combined experimental and quantum chemical study. Dyes and Pigments, 2013, 99, 957-965. | 2.0 | 42 |
| 11 | Toward Fully Nonempirical Simulations of Optical Band Shapes of Molecules in Solution: A Case Study of Heterocyclic Ketoimine Difluoroborates. Journal of Physical Chemistry A, 2015, 119, 5145-5152. | 1.1 | 39 |
| 12 | Influence of Bond Fixation in Benzo-AnnulatedN-Salicylideneanilines and Theirortho-C(O)X Derivatives (X = CH3, NH2, OCH3) on Tautomeric Equilibria in Solution. Journal of Organic Chemistry, 2007, 72, 5598-5607. | 1.7 | 38 |
| 13 | Photophysical Properties of Phenacylphenantridine Difluoroboranyls: Effect of Substituent and Double Benzannulation. Journal of Organic Chemistry, 2017, 82, 1529-1537. | 1.7 | 37 |
| 14 | Tautomeric equilibria, H-bonding and Ï€-electron delocalization ino-nitrosophenol. A B3LYP/6-311â€‰+â€‰G(2df,2p) study. Journal of Physical Organic Chemistry, 2005, 18, 892-897. | 0.9 | 34 |
| 15 | Topology-Driven Physicochemical Properties of Ï€-Î•lectron Systems. 1. Does the Clar Rule Work in Cyclic Ï€-Electron Systems with the Intramolecular Hydrogen or Lithium Bond?â€. Journal of Organic Chemistry, 2006, 71, 7678-7682. | 1.7 | 34 |
| 16 | Tautomeric preferences of phthalones and related compounds. Tetrahedron, 2007, 63, 9172-9178. | 1.0 | 34 |
| 17 | (1Z,3Z)-1,4-Di(pyridin-2-yl)buta-1,3-diene-2,3-diol:Â The Planar Highly Conjugated Symmetrical Enediol with Multiple Intramolecular Hydrogen Bonds. Journal of Organic Chemistry, 2002, 67, 3339-3345. | 1.7 | 32 |
| 18 | 2-Acylamino-6-pyridones: Breaking of an Intramolecular Hydrogen Bond by Self-association and Complexation with Double and Triple Hydrogen Bonding Counterparts. Uncommon Steric Effect on Intermolecular Interactions. Journal of Organic Chemistry, 2012, 77, 1653-1662. | 1.7 | 28 |

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| 19 | Effect of Ï€-Electron Delocalization on Tautomeric Equilibria â€“ Benzoannulated 2-Phenacylpyridines. European Journal of Organic Chemistry, 2006, 2006, 2817-2824. | 1.2 | 25 |
| 20 | Complexation of 2,6-Bis(acylamino)pyridines with Dipyridin-2-ylamine and 4,4-Dimethylpiperidine-2,6-dione. Journal of Physical Chemistry A, 2010, 114, 12881-12887. | 1.1 | 25 |
| 21 | <sup>15</sup>N NMR Studies of tautomerism. International Reviews in Physical Chemistry, 2012, 31, 567-629. | 0.9 | 25 |
| 22 | NMR and quantum chemical studies on association of 2,6-bis(acylamino)pyridines with selected imides and 2,2â€²-dipyridylamine. Structural Chemistry, 2010, 21, 1061-1067. | 1.0 | 23 |
| 23 | Substituent Effect in 2-Benzoylmethylenequinoline Difluoroborates Exhibiting Through-Space Couplings. Multinuclear Magnetic Resonance, X-ray Diffraction, and Computational Study. Journal of Physical Chemistry A, 2013, 117, 252-256. | 1.1 | 22 |
| 24 | Synthesis and Photophysical Properties of Novel Donorâ€“Acceptor N-(Pyridin-2-yl)-Substituted Benzo(thio)amides and Their Difluoroboranyl Derivatives. Journal of Physical Chemistry A, 2016, 120, 4116-4123. | 1.1 | 22 |
| 25 | Influence of the Nature of the Amino Group in Highly Fluorescent Difluoroborates Exhibiting Intramolecular Charge Transfer. Journal of Organic Chemistry, 2018, 83, 7779-7788. | 1.7 | 22 |
| 26 | Identity Double-Proton Transfer in (3Z)-3-Hydroxy-1,4-di(quinolin-2-yl)but-3-en-2-one. Chemistry - A European Journal, 2003, 9, 2710-2716. | 1.7 | 20 |
| 27 | 13C-NMR Based Evaluation of the Electronic and Steric Interactions in Aromatic Amines. International Journal of Molecular Sciences, 2005, 6, 52-62. | 1.8 | 20 |
| 28 | N-methyl-1,2-dihydro-2-benzoylmethylenequinolines: configurational dissimilarity with unmethylated congeners. Journal of Molecular Structure, 2000, 525, 233-239. | 1.8 | 19 |
| 29 | 4-Fluoroanilines: synthesis and decomposition. Journal of Fluorine Chemistry, 2001, 111, 1-10. | 0.9 | 19 |
| 30 | Application of spectroscopic and theoretical methods in the studies of photoisomerization and photophysical properties of the pushâ€”pull styryl-benzimidazole dyes. Photochemical and Photobiological Sciences, 2016, 15, 117-128. | 1.6 | 19 |
| 31 | Quantifying the Performances of DFT for Predicting Vibrationally Resolved Optical Spectra: Asymmetric Fluoroborate Dyes as Working Examples. Journal of Chemical Theory and Computation, 2017, 13, 4347-4356. | 2.3 | 18 |
| 32 | 2-Acylamino- and 2,4-Bis(acylamino)pyrimidines as Supramolecular Synthons Analyzed by Multiple Noncovalent Interactions. DFT, X-ray Diffraction, and NMR Spectral Studies. Journal of Organic Chemistry, 2012, 77, 9609-9619. | 1.7 | 17 |
| 33 | Association of N-(Pyridin-2-yl),Nâ€²-substituted Ureas with 2-Amino-1,8-naphthyridines and Benzoates: NMR and Quantum Chemical Studies of the Substituent Effect on Complexation. Journal of Organic Chemistry, 2013, 78, 7582-7593. | 1.7 | 17 |
| 34 | Two-photon absorption of BF<sub>2</sub>-carrying compounds: insights from theory and experiment. Physical Chemistry Chemical Physics, 2017, 19, 5705-5708. | 1.3 | 17 |
| 35 | Intermolecular steric hindrance in 7-acylamino-[1H]-2-oxo-1,8-naphthyridines: NMR, ESI-MS, IR, and DFT calculation studies. Structural Chemistry, 2011, 22, 1143-1151. | 1.0 | 14 |
| 36 | Controlling Two-Photon Action Cross Section by Changing a Single Heteroatom Position in Fluorescent Dyes. Journal of Physical Chemistry Letters, 2020, 11, 5920-5925. | 2.1 | 14 |

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| 37 | Molecular Decoration of Ceramic Supports for Highly Effective Enzyme Immobilizationâ€"Material Approach. Materials, 2021, 14, 201. | 1.3 | 14 |
| 38 | Long-range substituent and temperature effect on prototropic tautomerism in 2-(acylmethyl)quinolines. Journal of Physical Organic Chemistry, 2001, 14, 201-204. | 0.9 | 13 |
| 39 | Predominance of 2-arylhydrazones of 1,3-diphenylpropane-1,2,3-trione over its proton-transfer products. Journal of Physical Organic Chemistry, 2001, 14, 797-803. | 0.9 | 13 |
| 40 | Benchmarking Density Functional Approximations for Excited-State Properties of Fluorescent Dyes. Molecules, 2021, 26, 7434. | 1.7 | 13 |
| 41 | NMR crystallography of 2-acylamino-6-[1H]-pyridones: Solid-state NMR, GIPAW computational, and single crystal X-ray diffraction studies. Journal of Molecular Structure, 2011, 1006, 678-683. | 1.8 | 12 |
| 42 | NMR spectral assignment of substituted salicylaldoximes by inverse pulse techniques withz-gradient selection: correlation of NMR parameters with substituent constants. Magnetic Resonance in Chemistry, 1997, 35, 778-784. | 1.1 | 11 |
| 43 | Bischromophoric styrylpyridinium dyes. Spectrochimica Acta - Part A: Molecular and Biomolecular Spectroscopy, 2007, 67, 306-315. | 2.0 | 11 |
| 44 | DFT studies on tautomeric preferences. Part 3: Proton transfer in 2-(8-acylquinolin-2-yl)-1,3-diones. Structural Chemistry, 2010, 21, 1037-1041. | 1.0 | 11 |
| 45 | Tuning the hydrogen-bonding strength in 2,6-bis(cycloalkylcarbonylamino)pyridine assemblies by variable flexibility. Association constants measured by hydrogen-bonded vs. non-hydrogen-bonded protons. Supramolecular Chemistry, 2011, 23, 579-586. | 1.5 | 11 |
| 46 | The influence of secondary interactions on complex stability and double proton transfer reaction in 2-[1H]-pyridone/2-hydroxypyridine dimers. Journal of Molecular Modeling, 2011, 17, 2491-2500. | 0.8 | 11 |
| 47 | The influence of CH bond polarization on the self-association of 2-acylaminopyrimidines by NH/CHÂ·Â·Â·O/N interactions: XRD, NMR, DFT, and AIM study. Structural Chemistry, 2013, 24, 2203-2209. | 1.0 | 11 |
| 48 | Substituent effects in hydrogen bonding: DFT and QTAIM studies on acids and carboxylates complexes with formamide. Journal of Molecular Modeling, 2014, 20, 2356. | 0.8 | 11 |
| 49 | Tuning the Electronic Properties of the Dative Nâ†’B Bond with Associated Oâ†’B Interaction: Electron Localizability Indicator from Xâ€Ray Wavefunction Refinement. ChemPhysChem, 2016, 17, 2395-2406. | 1.0 | 11 |
| 50 | The fabrication, characterization, and pervaporation performance of poly(ether-block-amide) membranes blended with 4-(trifluoromethyl)-N(pyridine-2-yl)benzamide and 4-(dimethylamino)-N(pyridine-2-yl)benzamide fillers. Separation and Purification Technology, 2021, 268, 118707. | 3.9 | 11 |
| 51 | Electron ionization mass spectra and tautomerism of substituted 2â€phenacylquinolines. Rapid Communications in Mass Spectrometry, 2009, 23, 1075-1084. | 0.7 | 10 |
| 52 | Systematic investigation of 2,7-dihydroxy-1,8-naphthyridine dimerization â€" secondary interactions and tautomeric preferences calculations. Computational and Theoretical Chemistry, 2009, 908, 92-101. | 1.5 | 10 |
| 53 | Spectral and physicochemical properties of difluoroboranyls containing N,N-dimethylamino group studied by solvatochromic methods. Spectrochimica Acta - Part A: Molecular and Biomolecular Spectroscopy, 2018, 199, 86-95. | 2.0 | 10 |
| 54 | Difluoroboranyl derivatives as efficient panchromatic photoinitiators in radical polymerization reactions. Polymer Bulletin, 2018, 75, 3267-3281. | 1.7 | 10 |

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| 55 | Proton transfer reaction and intermolecular interactions in associates of 2,5-dihydroxy-1,8-naphthyridine. Journal of Molecular Modeling, 2012, 18, 1633-1644. | 0.8 | 9 |
| 56 | The impact of the heteroatom in a five-membered ring on the photophysical properties of difluoroborates. Dyes and Pigments, 2019, 170, 107481. | 2.0 | 9 |
| 57 | SYNTHESIS AND NMR SPECTRA OF 2-METHYL-2-QUINOLIN-2-YL-PROPIOPHENONES. Heterocyclic Communications, 1999, 5, . | 0.6 | 8 |
| 58 | Electron Ionization Mass Spectra and Tautomerism of 2-Phenacylpyridines. European Journal of Mass Spectrometry, 2006, 12, 25-29. | 0.5 | 8 |
| 59 | Secondary interactions as driving force in heterocomplex formation of 2,7-disubstituted-1,8-naphthyridines: Quantum chemical, NMR and mass spectral investigations. Journal of Molecular Structure, 2009, 931, 60-67. | 1.8 | 8 |
| 60 | Conformational equilibrium in supramolecular chemistry: Dibutyltriuret case. Beilstein Journal of Organic Chemistry, 2015, 11, 2105-2116. | 1.3 | 8 |
| 61 | A detailed theoretical and experimental study on the N H, P O and C O stretching frequencies in two new phosphoric triamides and a statistical comparison with analogous structures. Polyhedron, 2019, 158, 215-224. | 1.0 | 8 |
| 62 | Structural characterization of Î²-2â€²-pyridylaminocrotonoyl-2-pyridylamide by ESI-MS, NMR, single crystal X-ray analysis and ab initio methods. Journal of Molecular Structure, 2003, 654, 61-69. | 1.8 | 7 |
| 63 | The effect of benzoannulation on the transition state and the proton transfer equilibrium in di(2-pyridyl)methane derivatives. New Journal of Chemistry, 2011, 35, 1433. | 1.4 | 7 |
| 64 | The Application of 2,6-Bis(4-Methoxybenzoyl)-Diaminopyridine in Solvent Extraction and Polymer Membrane Separation for the Recovery of Au(III), Ag(I), Pd(II) and Pt(II) Ions from Aqueous Solutions. International Journal of Molecular Sciences, 2021, 22, 9123. | 1.8 | 7 |
| 65 | Synthesis of 6-acylmethylphenanthridine enaminones. Journal of the Iranian Chemical Society, 2005, 2, 294-299. | 1.2 | 6 |
| 66 | Complex tauto- and rotamerism of 2-(R-phenyl)-1,2,3,4-tetrahydroquinazolines. Journal of Physical Organic Chemistry, 2005, 18, 737-742. | 0.9 | 6 |
| 67 | (1Z,3Z)-3-[Quinolin-2(1H)-ylidene]-1-(quinolin-2-yl)prop-1-en-2-ol: An unexpected most stable tautomer of 1,3-bis(quinolin-2-yl)acetone. Journal of Molecular Structure, 2009, 930, 78-82. | 1.8 | 6 |
| 68 | Noncovalent interactions between classical supramolecular synthons in solution: Hydrogen bonding in hindered 2-acylaminopyridine/2-pyridone associates. Journal of Molecular Structure, 2012, 1018, 84-87. | 1.8 | 6 |
| 69 | Comment on â€œNon-symmetric substituted ureas locked in an (E,Z) conformation: an unusual anion binding via supramolecular assemblyâ€•by M. Olivari, C. Caltagirone, A. Garau, F. Isaia, M. E. Light, V. Lippolis, R. Montis and M. A. Scorciapino, New J. Chem., 2013, 37, 663. New Journal of Chemistry, 2014, 38, 2701. | 1.4 | 6 |
| 70 | Design of Twoâ€Photonâ€Excited Fluorescent Dyes Containing Fluoroborylene Groups. ChemPhotoChem, 2019, 3, 719-726. | 1.5 | 6 |
| 71 | Symmetric Fluoroborate and its Boron Modification: Crystal and Electronic Structures. Crystals, 2019, 9, 662. | 1.0 | 6 |
| 72 | 2,6-Bis((benzoyl-R)amino)pyridine (R = H, 4-Me, and 4-NMe2) Derivatives for the Removal of Cu(II), Ni(II), Co(II), and Zn(II) Ions from Aqueous Solutions in Classic Solvent Extraction and a Membrane Extraction. Membranes, 2021, 11, 233. | 1.4 | 6 |

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| 73 | Substituted 2-(2-hydroxyphenyl)â€“3H-quinazolin-4-ones and their difluoroboron complexes: Synthesis and photophysical properties. Spectrochimica Acta - Part A: Molecular and Biomolecular Spectroscopy, 2021, 252, 119497. | 2.0 | 6 |
| 74 | Tailoring the nonlinear absorption of fluorescent dyes by substitution at a boron center. Journal of Materials Chemistry C, 2021, 9, 6225-6233. | 2.7 | 6 |
| 75 | Less is more: On the effect of benzannulation on solid-state emission of difluoroborates. Journal of Materials Chemistry C, 0, , . | 2.7 | 6 |
| 76 | Predominance of resonance over polar effects on1H,13C and15N NMR substituent chemical shifts inN-arylglycines. Magnetic Resonance in Chemistry, 1998, 36, 848-854. | 1.1 | 5 |
| 77 | Predominance of inductive over resonance substituent effect on33S NMR chemical shifts of 4-substituted phenyl-4â€²-methylphenacyl sulfones. Magnetic Resonance in Chemistry, 1999, 37, 437-440. | 1.1 | 5 |
| 78 | GIAO/DFT 13C NMR Chemical Shifts of 1,3,4-Thiadiazoles. Phosphorus, Sulfur and Silicon and the Related Elements, 2007, 182, 2217-2225. | 0.8 | 5 |
| 79 | Use of time-resolved fluorescence spectroscopy to evaluate diagnostic value of collagen degradation products. Journal of Biomedical Optics, 2015, 20, 051039. | 1.4 | 5 |
| 80 | Conformational change in the association of a heterocyclic urea derivative forming two intramolecular hydrogen bonds in polar solvent. New Journal of Chemistry, 2017, 41, 1073-1081. | 1.4 | 5 |
| 81 | The trans/cis photoisomerization in hydrogen bonded complexes with stability controlled by substituent effects: 3-(6-aminopyridin-3-yl)acrylate case study. RSC Advances, 2018, 8, 23698-23710. | 1.7 | 5 |
| 82 | Synthesis and Photophysical Studies of Novel Vâ€Shaped 2,3â€Bis{5â€arylâ€2â€thienyl}(dibenzo[f,h])quinoxalines. Asian Journal of Organic Chemistry, 2020, 9, 673-681. | 1.3 | 5 |
| 83 | Two-Photon Absorption Activity of BOPHY Derivatives: Insights from Theory. Journal of Physical Chemistry A, 2021, 125, 2581-2587. | 1.1 | 5 |
| 84 | Effect of vinylene and 1,4-phenylene spacers on efficiency of the ground-state intramolecular charge-transfer in enlarged 4-dimethylamino-1-methylpyridinium cations. Structural Chemistry, 2009, 20, 655-662. | 1.0 | 4 |
| 85 | DFT studies on tautomeric preferences of 1-(pyridin-2-yl)-4-(quinolin-2-yl)butane-2,3-dione in the gas phase and in solution. Structural Chemistry, 2010, 21, 1283-1287. | 1.0 | 4 |
| 86 | Association of 2-acylaminopyridines and benzoic acids. Steric and electronic substituent effect studied by XRD, solution and solid-state NMR and calculations. Journal of Molecular Structure, 2013, 1054-1055, 157-163. | 1.8 | 4 |
| 87 | The Copper(II) Ions Solvent Extraction with a New Compound: 2,6-Bis(4-Methoxybenzoyl)-Diaminopyridine. Processes, 2019, 7, 954. | 1.3 | 4 |
| 88 | (A)symmetric chromophores based on cyano and fluorine-substituted 2,3-bis(5-arylthiophen-2-yl)quinoxalines: Synthesis, photophysical properties and application prospects. Dyes and Pigments, 2022, 204, 110434. | 2.0 | 4 |
| 89 | NMR spectral and X-ray structural investigation of 1,3-bis(2-quinolyl)-2-(p-chlorophenyl)-2-propanol. Journal of Molecular Structure, 2000, 525, 241-245. | 1.8 | 3 |
| 90 | NMR Spectra of Anilines. , 0, , 347-371. | | 3 |

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| 91 | Two (E)-2-({[4-(dialkylamino)phenyl]imino}methyl)-4-nitrophenols. Acta Crystallographica Section C: Crystal Structure Communications, 2012, 68, o279-o282. | 0.4 | 3 |
| 92 | STABILITY OF THE DIMERS OF AZA ANALOGS OF 2-FORMYLPYRROLE. CONJUGATION VERSUS HYDROGEN BONDING. Heterocyclic Communications, 2003, 9, . | 0.6 | 2 |
| 93 | Collision induced dissociation of N -(pyridin-2-yl)-substituted benzo(thio)amides and N -(isoquinolin-1-yl)furan(thiophene)-2-carboxamides and their difluoroboranyl derivatives. International Journal of Mass Spectrometry, 2018, 428, 35-42. | 0.7 | 2 |
| 94 | Tautomeric equilibrium, proton affinity and mass spectrometry fragmentation of flexible hydrogen-bonded precursors and rigid $$hbox {N}longrightarrow hbox {BF}_2$$ fluorescent dyes. Scientific Reports, 2021, 11, 15995. | 1.6 | 2 |
| 95 | Conformational Equilibrium and Substituent Effects in Hydrogen-bonded Complexes. Current Organic Chemistry, 2018, 22, 2182-2199. | 0.9 | 2 |
| 96 | (Z)-Ethyl 2-oxo-3-(1,2-dihydroquinolin-2-ylidene)propanoate. Acta Crystallographica Section E: Structure Reports Online, 2010, 66, o1746-o1747. | 0.2 | 1 |
| 97 | N2,N2,N6,N6-Tetrakis(2,3,4,5,6-pentafluorobenzoyl)pyridine-2,6-diamine. Acta Crystallographica Section E: Structure Reports Online, 2011, 67, o3429-o3430. | 0.2 | 1 |
| 98 | Conformational and Tautomeric Control by Supramolecular Approach in Ureido-N-iso-propyl,Nâ€™-4-(3-pyridin-2-one)pyrimidine. Molecules, 2019, 24, 2491. | 1.7 | 1 |
| 99 | 2-Methyl-N-(pyrazin-2-yl)propanamideâ€“1,2,4,5-tetrafluoro-3,6-diiodobenzene (2/1). IUCrData, 2016, 1, . | 0.1 | 1 |
| 100 | N-(Pyrazin-2-yl)adamantane-1-carboxamide. IUCrData, 2016, 1, . | 0.1 | 1 |
| 101 | Effect of conjugated system extension on structural features and electron-density distribution in chargeâ€“transfer difluoroborates. Acta Crystallographica Section C, Structural Chemistry, 2021, 77, 807-813. | 0.2 | 1 |
| 102 | Supramolecular Approach to Tuning the Photophysical Properties of Quadrupolar Squaraines. Frontiers in Chemistry, 2021, 9, 800541. | 1.8 | 1 |
| 103 | Tautomeric Equilibria in Relation to Pi-Electron Delocalization. ChemInform, 2006, 37, no. | 0.1 | 0 |
| 104 | N-[2-(2,2-Dimethylpropanamido)pyrimidin-4-yl]-2,2-dimethylpropanamiden-hexane 0.25-solvate hemihydrate. Acta Crystallographica Section E: Structure Reports Online, 2013, 69, o1617-o1618. | 0.2 | 0 |
| 105 | N-(2-Benzoyl-4-chlorophenyl)-4-chlorobenzenesulfonamide. Acta Crystallographica Section E: Structure Reports Online, 2008, 64, o737-o737. | 0.2 | 0 |
| 106 | 6-Amino-2-(pivaloylamino)pyridinium benzoate. Acta Crystallographica Section E: Structure Reports Online, 2013, 69, o1483-o1484. | 0.2 | 0 |
| 107 | N,Nâ€²-Bis(pyridin-2-yl)octanediamide. IUCrData, 2016, 1, . | 0.1 | 0 |
| 108 | 2,2-Difluoro-3-(4-fluorophenyl)-2H-benzo[e][1,3,2]oxazaborinin-3-ium-2-uide. IUCrData, 2017, 2, . | 0.1 | 0 |

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| 109 | 2-[4-(Dimethylamino)phenyl]-3,3-difluoro-3H-naphtho[1,2-e][1,3,2]oxazaborinin-2-ium-3-uide. IUCrData, 2017, 2, . | 0.1 | 0 |