

# Steve Scheiner

## List of Publications by Year in descending order

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316  
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#	ARTICLE	IF	CITATIONS
1	Promotion of TH3 (T = Si and Ge) group transfer within a tetrel bond by a cation-π interaction. <i>Physical Chemistry Chemical Physics</i> , 2022, 24, 1113-1119.	1.3	3
2	Influence of Substituents in the Benzene Ring on the Halogen Bond of Iodobenzene with Ammonia. <i>ChemPhysChem</i> , 2022, 23, .	1.0	19
3	Principles Guiding the Square Bonding Motif Containing a Pair of Chalcogen Bonds between Chalcogenadiazoles. <i>Journal of Physical Chemistry A</i> , 2022, 126, 1194-1203.	1.1	13
4	Search for an exothermic halogen bond between anions. <i>Physical Chemistry Chemical Physics</i> , 2022, 24, 6964-6972.	1.3	3
5	Characterization of Type I and II Interactions between Halogen Atoms. <i>Crystal Growth and Design</i> , 2022, 22, 2692-2702.	1.4	16
6	The Role of Hydrogen Bonds in Interactions between [PdCl <sub>4</sub> ] <sup>2-</sup> Dianions in Crystal. <i>Molecules</i> , 2022, 27, 2144.	1.7	4
7	Competition between Intra and Intermolecular Pnicogen Bonds. Complexes between Naphthalene Derivatives and Neutral or Anionic Bases. <i>ChemPhysChem</i> , 2022, , .	1.0	4
8	Resonance-assisted intramolecular triel bonds. <i>Physical Chemistry Chemical Physics</i> , 2022, 24, 15015-15024.	1.3	6
9	Experimental and Theoretical Evidence of a Pb·····Pb Ditetrel Bond Without a π-Hole. <i>ChemPhysChem</i> , 2022, 23, .	1.0	4
10	Various Sorts of Chalcogen Bonds Formed by an Aromatic System. <i>Journal of Physical Chemistry A</i> , 2022, 126, 4025-4035.	1.1	9
11	Carbon as an electron donor atom. <i>Polyhedron</i> , 2021, 193, 114905.	1.0	8
12	Unusual substituent effects in the Tr·····Te triel bond. <i>International Journal of Quantum Chemistry</i> , 2021, 121, e26526.	1.0	6
13	Comparison of Bifurcated Halogen with Hydrogen Bonds. <i>Molecules</i> , 2021, 26, 350.	1.7	12
14	Experimental and Theoretical Studies of Dimers Stabilized by Two Chalcogen Bonds in the Presence of a N·····N Pnicogen Bond. <i>Journal of Physical Chemistry A</i> , 2021, 125, 657-668.	1.1	14
15	Origins and properties of the tetrel bond. <i>Physical Chemistry Chemical Physics</i> , 2021, 23, 5702-5717.	1.3	88
16	Weak π-Hole Triel Bond between C <sub>5</sub> H <sub>5</sub> Tr (Tr=B, Al, Ga) and Haloethyne: Substituent and Cooperativity Effects. <i>ChemPhysChem</i> , 2021, 22, 481-487.	1.0	15
17	Molecular Recognition. <i>ChemPhysChem</i> , 2021, 22, 433-434.	1.0	4
18	Relative Strengths of a Pnicogen and a Tetrel Bond and Their Mutual Effects upon One Another. <i>Journal of Physical Chemistry A</i> , 2021, 125, 2631-2641.	1.1	13

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19	Noncovalent Bonds through Sigma and Pi-Hole Located on the Same Molecule. Guiding Principles and Comparisons. <i>Molecules</i> , 2021, 26, 1740.	1.7	32
20	Competition between Inter and Intramolecular Tetrel Bonds: Theoretical Studies Complemented by CSD Survey. <i>ChemPhysChem</i> , 2021, 22, 924-934.	1.0	7
21	Crystallographic and Theoretical Evidences of Anion...Anion Interaction. <i>ChemPhysChem</i> , 2021, 22, 818-821.	1.0	25
22	Anion...Anion Interactions in Aerogen-Bonded Complexes. Influence of Solvent Environment. <i>Molecules</i> , 2021, 26, 2116.	1.7	13
23	Fabricating Flexible Packaging Batteries in General Chemistry Laboratories. <i>Journal of Chemical Education</i> , 2021, 98, 2471-2475.	1.1	2
24	Proximity Effects of Substituents on Halogen Bond Strength. <i>Journal of Physical Chemistry A</i> , 2021, 125, 5069-5077.	1.1	17
25	Diboron Bonds Between BX <sub>3</sub> (X=H, F, CH <sub>3</sub> ) and BYZ <sub>2</sub> (Y=H, F; Z=CO, N <sub>2</sub> , CNH). <i>ChemPhysChem</i> , 2021, 22, 1461-1469.	1.0	4
26	Probing the Hydrogen-Bonding Environment of Individual Bases in DNA Duplexes with Isotope-Edited Infrared Spectroscopy. <i>Journal of Physical Chemistry B</i> , 2021, 125, 7613-7627.	1.2	9
27	Dissection of the Origin of $\sigma$ -Holes and the Noncovalent Bonds in Which They Engage. <i>Journal of Physical Chemistry A</i> , 2021, 125, 6514-6528.	1.1	21
28	Enhancement of the Tetrel Bond by the Effects of Substituents, Cooperativity, and Electric Field: Transition from Noncovalent to Covalent Bond. <i>ChemPhysChem</i> , 2021, 22, 2305-2312.	1.0	6
29	Partial transfer of bridging atom in halogen-bonded complexes. <i>Computational and Theoretical Chemistry</i> , 2021, 1204, 113398.	1.1	2
30	Anion...anion and anion...neutral triel bonds. <i>Physical Chemistry Chemical Physics</i> , 2021, 23, 4818-4828.	1.3	19
31	Participation of S and Se in hydrogen and chalcogen bonds. <i>CrystEngComm</i> , 2021, 23, 6821-6837.	1.3	29
32	Noncovalent bond between tetrel $\sigma$ -hole and hydride. <i>Physical Chemistry Chemical Physics</i> , 2021, 23, 10536-10544.	1.3	2
33	Anion...anion (MX <sub>3</sub> <sup>+</sup> ) <sub>2</sub> dimers (M = Zn, Cd, Hg; X = Cl, Br, I) in different environments. <i>Physical Chemistry Chemical Physics</i> , 2021, 23, 13853-13861.	1.3	16
34	Competition between a Tetrel and Halogen Bond to a Common Lewis Acid. <i>Journal of Physical Chemistry A</i> , 2021, 125, 308-316.	1.1	14
35	Ability of Lewis Acids with Shallow $\sigma$ -Holes to Engage in Chalcogen Bonds in Different Environments. <i>Molecules</i> , 2021, 26, 6394.	1.7	9
36	Anatomy of $\sigma$ -hole bonds: Linear systems. <i>Journal of Chemical Physics</i> , 2021, 155, 174302.	1.2	5

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37	Trial bonds within anion-anion complexes. <i>Physical Chemistry Chemical Physics</i> , 2021, 23, 25097-25106.	1.3	6
38	On the Ability of Nitrogen to Serve as an Electron Acceptor in a Pnictogen Bond. <i>Journal of Physical Chemistry A</i> , 2021, 125, 10419-10427.	1.1	14
39	Maximal occupation by bases of $\pi$ -hole bands surrounding linear molecules. <i>Journal of Computational Chemistry</i> , 2021, , .	1.5	2
40	Experimental and theoretical evidence of attractive interactions between dianions: $[\text{PdCl}_4]^{2-} \cdots [\text{PdCl}_4]^{2-}$ . <i>Chemical Communications</i> , 2021, 57, 13305-13308.	2.2	7
41	Structures and energetics of clusters surrounding diatomic anions stabilized by hydrogen, halogen, and other noncovalent bonds. <i>Chemical Physics</i> , 2020, 530, 110590.	0.9	15
42	The Hydrogen Bond: A Hundred Years and Counting. <i>Journal of the Indian Institute of Science</i> , 2020, 100, 61-76.	0.9	34
43	Tuning the Competition between Hydrogen and Tetrel Bonds by a Magnesium Bond. <i>ChemPhysChem</i> , 2020, 21, 212-219.	1.0	28
44	Coordination of anions by noncovalently bonded $f$ -hole ligands. <i>Coordination Chemistry Reviews</i> , 2020, 405, 213136.	9.5	66
45	Versatility of the Cyano Group in Intermolecular Interactions. <i>Molecules</i> , 2020, 25, 4495.	1.7	8
46	Understanding noncovalent bonds and their controlling forces. <i>Journal of Chemical Physics</i> , 2020, 153, 140901.	1.2	46
47	The balance between side-chain and backbone-driven association in folding of the $\alpha$ -helical influenza A transmembrane peptide. <i>Journal of Computational Chemistry</i> , 2020, 41, 2177-2188.	1.5	3
48	Noncovalent Bonds between Tetrel Atoms. <i>ChemPhysChem</i> , 2020, 21, 1934-1944.	1.0	24
49	F-Halogen Bond: Conditions for Its Existence. <i>Journal of Physical Chemistry A</i> , 2020, 124, 7290-7299.	1.1	17
50	Relationships between Bond Strength and Spectroscopic Quantities in H-Bonds and Related Halogen, Chalcogen, and Pnictogen Bonds. <i>Journal of Physical Chemistry A</i> , 2020, 124, 7716-7725.	1.1	16
51	Effect of carbon hybridization in $\text{C}=\text{F}$ bond as an electron donor in trial bonds. <i>Journal of Chemical Physics</i> , 2020, 153, 074304.	1.2	6
52	Pnictogen Bonds Pairing Anionic Lewis Acid with Neutral and Anionic Bases. <i>Journal of Physical Chemistry A</i> , 2020, 124, 4998-5006.	1.1	24
53	Complexes of $\text{HArF}$ and $\text{AuX}$ ( $X = \text{F}, \text{Cl}, \text{Br}, \text{I}$ ). Comparison of $\text{H} \cdots \text{F}$ bonds, halogen bonds, $\text{F} \cdots \text{F}$ shared bonds and covalent bonds. <i>Applied Organometallic Chemistry</i> , 2020, 34, e5891.	1.7	6
54	Coordination of a Central Atom by Multiple Intramolecular Pnictogen Bonds. <i>Inorganic Chemistry</i> , 2020, 59, 9315-9324.	1.9	19

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55	The ditetrel bond: noncovalent bond between neutral tetrel atoms. <i>Physical Chemistry Chemical Physics</i> , 2020, 22, 16606-16614.	1.3	29
56	How Many Pnictogen Bonds can be Formed to a Central Atom Simultaneously?. <i>Journal of Physical Chemistry A</i> , 2020, 124, 2046-2056.	1.1	29
57	Xe $\pi$ -chalcogen aerogen bond. Effect of substituents and size of chalcogen atom. <i>Physical Chemistry Chemical Physics</i> , 2020, 22, 4115-4121.	1.3	11
58	Competition between Intra and Intermolecular Triel Bonds. Complexes between Naphthalene Derivatives and Neutral or Anionic Lewis Bases. <i>Molecules</i> , 2020, 25, 635.	1.7	20
59	On the Stability of Interactions between Pairs of Anions $\pi$ -Complexes of $MCl_3$ ( $M=Be, Mg, Ca, Sr, Ba$ ) with Pyridine and $CN$ . <i>ChemPhysChem</i> , 2020, 21, 870-877.	1.0	25
60	Anion $\pi$ -Anion Attraction in Complexes of $MCl_3$ ( $M=Zn, Cd, Hg$ ) with $CN$ . <i>ChemPhysChem</i> , 2020, 21, 1119-1125.	1.0	31
61	Effects of Halogen, Chalcogen, Pnictogen, and Tetrel Bonds on IR and NMR Spectra. <i>Molecules</i> , 2019, 24, 2822.	1.7	41
62	Violation of Electrostatic Rules: Shifting the Balance between Pnictogen Bonds and Lone Pair $\pi$ -Interactions Tuned by Substituents. <i>Journal of Physical Chemistry A</i> , 2019, 123, 7288-7295.	1.1	11
63	The ability of a tetrel bond to transition a neutral amino acid into a zwitterion. <i>Chemical Physics Letters</i> , 2019, 731, 136584.	1.2	9
64	Comparison of halogen with proton transfer. Symmetric and asymmetric systems. <i>Chemical Physics Letters</i> , 2019, 731, 136593.	1.2	4
65	Theoretical Studies of IR and NMR Spectral Changes Induced by Sigma-Hole Hydrogen, Halogen, Chalcogen, Pnictogen, and Tetrel Bonds in a Model Protein Environment. <i>Molecules</i> , 2019, 24, 3329.	1.7	35
66	Chalcogen bonding of two ligands to hypervalent $YF_4$ ( $Y = S, Se, Te, Po$ ). <i>Physical Chemistry Chemical Physics</i> , 2019, 21, 20829-20839.	1.3	27
67	Dual Geometry Schemes in Tetrel Bonds: Complexes between $TF_4$ ( $T = Si, Ge, Sn$ ) and Pyridine Derivatives. <i>Molecules</i> , 2019, 24, 376.	1.7	28
68	Switchable Aromaticity in an Isostructural Mn Phthalocyanine Series Isolated in Five Separate Redox States. <i>Journal of the American Chemical Society</i> , 2019, 141, 2604-2613.	6.6	28
69	On the ability of pnictogen atoms to engage in both $\pi$ and $\pi$ -hole complexes. Heterodimers of $ZF_2C_6H_5$ ( $Z=P, As, Sb, Bi$ ) and $NH_3$ . <i>Journal of Molecular Modeling</i> , 2019, 25, 152.	0.8	29
70	Computational Insights into $Mg\pi$ -Cl Complex Electrolytes for Rechargeable Magnesium Batteries. <i>Batteries and Supercaps</i> , 2019, 2, 792-800.	2.4	16
71	Comparison between Hydrogen and Halogen Bonds in Complexes of $6\pi$ -OX $\pi$ -Fulvene with Pnictogen and Chalcogen Electron Donors. <i>ChemPhysChem</i> , 2019, 20, 1978-1984.	1.0	16
72	Forty years of progress in the study of the hydrogen bond. <i>Structural Chemistry</i> , 2019, 30, 1119-1128.	1.0	39

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73	Interactions of (MY) <sub>6</sub> (M = Zn, Cd; Y = O, S, Se) quantum dots with N-bases. <i>Structural Chemistry</i> , 2019, 30, 1003-1014.	1.0	2
74	Structures of clusters surrounding ions stabilized by hydrogen, halogen, chalcogen, and pnictogen bonds. <i>Chemical Physics</i> , 2019, 524, 55-62.	0.9	13
75	Influence of monomer deformation on the competition between two types of $\sigma$ -holes in tetrel bonds. <i>Physical Chemistry Chemical Physics</i> , 2019, 21, 10336-10346.	1.3	20
76	On the capability of metal-halogen groups to participate in halogen bonds. <i>CrystEngComm</i> , 2019, 21, 2875-2883.	1.3	18
77	Optical Stability of 1,1'-Binaphthyl Derivatives. <i>ACS Omega</i> , 2019, 4, 6044-6049.	1.6	11
78	Differential Binding of Tetrel-Bonding Bipodal Receptors to Monatomic and Polyatomic Anions. <i>Molecules</i> , 2019, 24, 227.	1.7	21
79	Structural and Functional Characterization of Sulfonium Carbon-Oxygen Hydrogen Bonding in the Deoxyamino Sugar Methyltransferase TylM1. <i>Biochemistry</i> , 2019, 58, 2152-2159.	1.2	0
80	Hexacoordinated Tetrel-Bonded Complexes between TF <sub>4</sub> (T=Si, Ge, Sn, Pb) and NCH: Competition between $\sigma$ - and $\pi$ -holes. <i>ChemPhysChem</i> , 2019, 20, 959-966.	1.0	25
81	Comparison of $\sigma$ -hole and $\pi$ -hole tetrel bonds in complexes of borazine with TH <sub>3</sub> F and F <sub>2</sub> TO/H <sub>2</sub> TO (T=C, Si, Ge). <i>International Journal of Quantum Chemistry</i> , 2019, 119, e25910.	1.0	19
82	Definition of the chalcogen bond (IUPAC Recommendations 2019). <i>Pure and Applied Chemistry</i> , 2019, 91, 1889-1892.	0.9	322
83	Carbene triel bonds between TrR <sub>3</sub> (Tr=B, Al) and N-heterocyclic carbenes. <i>International Journal of Quantum Chemistry</i> , 2019, 119, e25867.	1.0	27
84	Dependence of NMR chemical shifts upon CH bond lengths of a methyl group involved in a tetrel bond. <i>Chemical Physics Letters</i> , 2019, 714, 61-64.	1.2	16
85	Implications of monomer deformation for tetrel and pnictogen bonds. <i>Physical Chemistry Chemical Physics</i> , 2018, 20, 8832-8841.	1.3	67
86	Steric Crowding in Tetrel Bonds. <i>Journal of Physical Chemistry A</i> , 2018, 122, 2550-2562.	1.1	55
87	Halogen, Chalcogen, and Pnictogen Bonding Involving Hypervalent Atoms. <i>Chemistry - A European Journal</i> , 2018, 24, 8167-8177.	1.7	68
88	Effect of Magnesium Bond on the Competition Between Hydrogen and Halogen Bonds and the Induction of Proton and Halogen Transfer. <i>ChemPhysChem</i> , 2018, 19, 1456-1464.	1.0	11
89	Aerogen bonds formed between AeOF <sub>2</sub> (Ae = Kr, Xe) and diazines: comparisons between $\sigma$ -hole and $\pi$ -hole complexes. <i>Physical Chemistry Chemical Physics</i> , 2018, 20, 4676-4687.	1.3	36
90	Comparison of Various Means of Evaluating Molecular Electrostatic Potentials for Noncovalent Interactions. <i>Journal of Computational Chemistry</i> , 2018, 39, 500-510.	1.5	27

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91	The $\pi$ - $\pi$ Tetrel Bond and its Influence on Hydrogen Bonding and Proton Transfer. <i>ChemPhysChem</i> , 2018, 19, 736-743.	1.0	46
92	Crystallographic and Computational Characterization of Methyl Tetrel Bonding in S-Adenosylmethionine-Dependent Methyltransferases. <i>Molecules</i> , 2018, 23, 2965.	1.7	29
93	Ability of IR and NMR Spectral Data to Distinguish between a Tetrel Bond and a Hydrogen Bond. <i>Journal of Physical Chemistry A</i> , 2018, 122, 7852-7862.	1.1	31
94	Triel-Bonded Complexes between $TrR_3$ ( $Tr=B, Al, Ga; R=H, F, Cl, Br, CH_3$ ) and Pyrazine. <i>ChemPhysChem</i> , 2018, 19, 3122-3133.	1.0	25
95	Tetrel Bonding as a Vehicle for Strong and Selective Anion Binding. <i>Molecules</i> , 2018, 23, 1147.	1.7	39
96	Comparative Strengths of Tetrel, Pnicogen, Chalcogen, and Halogen Bonds and Contributing Factors. <i>Molecules</i> , 2018, 23, 1681.	1.7	69
97	Comparison between Tetrel Bonded Complexes Stabilized by $\pi$ and $\pi$ Hole Interactions. <i>Molecules</i> , 2018, 23, 1416.	1.7	45
98	Water-Mediated Carbon-Oxygen Hydrogen Bonding Facilitates <i>S</i> -Adenosylmethionine Recognition in the Reactivation Domain of Cobalamin-Dependent Methionine Synthase. <i>Biochemistry</i> , 2018, 57, 3733-3740.	1.2	16
99	Regium bonds between $M_n$ clusters ( $M = Cu, Ag, Au$ and $n = 2-6$ ) and nucleophiles $NH_3$ and HCN. <i>Physical Chemistry Chemical Physics</i> , 2018, 20, 22498-22509.	1.3	46
100	Comparison of tetrel bonds in neutral and protonated complexes of pyridine $T_3$ and furan $T_3$ ( $T = C, Si, \text{ and } Ge$ ) with $NH_3$ . <i>Physical Chemistry Chemical Physics</i> , 2017, 19, 5550-5559.	1.3	98
101	Assembly of Effective Halide Receptors from Components. Comparing Hydrogen, Halogen, and Tetrel Bonds. <i>Journal of Physical Chemistry A</i> , 2017, 121, 3606-3615.	1.1	56
102	Comparison of halide receptors based on H, halogen, chalcogen, pnicogen, and tetrel bonds. <i>Faraday Discussions</i> , 2017, 203, 213-226.	1.6	57
103	The halogen bond in solution: general discussion. <i>Faraday Discussions</i> , 2017, 203, 347-370.	1.6	5
104	Computational approaches and sigma-hole interactions: general discussion. <i>Faraday Discussions</i> , 2017, 203, 131-163.	1.6	17
105	Can HCCH/HBNH Break B-N/C-C Bonds of Single-Wall BN/Carbon Nanotubes at Their Surface?. <i>Journal of Physical Chemistry C</i> , 2017, 121, 26044-26053.	1.5	0
106	Systematic Elucidation of Factors That Influence the Strength of Tetrel Bonds. <i>Journal of Physical Chemistry A</i> , 2017, 121, 5561-5568.	1.1	108
107	Monitoring the Charge Distribution during Proton and Sodium Ion Conduction along Chains of Water Molecules and Protein Residues. <i>Israel Journal of Chemistry</i> , 2017, 57, 385-392.	1.0	2
108	Halogen Bonds Formed between Substituted Imidazoliums and N Bases of Varying N-Hybridization. <i>Molecules</i> , 2017, 22, 1634.	1.7	18

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109	Assessment of the Presence and Strength of H-Bonds by Means of Corrected NMR. <i>Molecules</i> , 2016, 21, 1426.	1.7	30
110	Hâ€bonding and stacking interactions between chloroquine and temozolomide. <i>International Journal of Quantum Chemistry</i> , 2016, 116, 1196-1204.	1.0	7
111	Segmentation and additive approach: A reliable technique to study noncovalent interactions of large molecules at the surface of singleâ€wall carbon nanotubes. <i>Journal of Computational Chemistry</i> , 2016, 37, 1953-1961.	1.5	0
112	Enhancing the Reduction Potential of Quinones via Complex Formation. <i>Journal of Organic Chemistry</i> , 2016, 81, 4316-4324.	1.7	9
113	Torsional and Electronic Factors Control the CâˆHâ€...â€...O Interaction. <i>Chemistry - A European Journal</i> , 2016, 22, 16513-16521.	1.7	18
114	Highly Selective Halide Receptors Based on Chalcogen, Pnicogen, and Tetrel Bonds. <i>Chemistry - A European Journal</i> , 2016, 22, 18850-18858.	1.7	98
115	Interactions of Nucleic Acid Bases with Temozolomide. Stacked, Perpendicular, and Coplanar Heterodimers. <i>Journal of Physical Chemistry B</i> , 2016, 120, 9347-9361.	1.2	10
116	Effects of Angular Deformation on the Energetics of the S<sub>N</sub>2 Reaction. <i>European Journal of Organic Chemistry</i> , 2016, 2016, 3964-3968.	1.2	1
117	Interactions between temozolomide and quercetin. <i>Structural Chemistry</i> , 2016, 27, 1577-1588.	1.0	7
118	Interpretation of Spectroscopic Markers of Hydrogen Bonds. <i>ChemPhysChem</i> , 2016, 17, 2263-2271.	1.0	17
119	NXâ€Y halogen bonds. Comparison with NHâ€Y H-bonds and CXâ€Y halogen bonds. <i>Physical Chemistry Chemical Physics</i> , 2016, 18, 18015-18023.	1.3	17
120	Building a Better Halide Receptor: Optimum Choice of Spacer, Binding Unit, and Halosubstitution. <i>ChemPhysChem</i> , 2016, 17, 836-844.	1.0	15
121	Hydrogen bonded and stacked geometries of the temozolomide dimer. <i>Journal of Molecular Modeling</i> , 2016, 22, 77.	0.8	13
122	Catalysis of the Aza-Dielsâ€Alder Reaction by Hydrogen and Halogen Bonds. <i>Journal of Organic Chemistry</i> , 2016, 81, 2589-2597.	1.7	38
123	Sulfurâ€Oxygen Chalcogen Bonding Mediates AdoMet Recognition in the Lysine Methyltransferase SET7/9. <i>ACS Chemical Biology</i> , 2016, 11, 748-754.	1.6	93
124	Comparison of ĩ€-hole tetrel bonding with ĩf-hole halogen bonds in complexes of XCN (X = F, Cl, Br, I) and NH<sub>3</sub>. <i>Physical Chemistry Chemical Physics</i> , 2016, 18, 3581-3590.	1.3	99
125	Regioselectivity of the interaction of temozolomide with borane and boron trifluoride. <i>Structural Chemistry</i> , 2015, 26, 1359-1365.	1.0	9
126	Competitive Halide Binding by Halogen Versus Hydrogen Bonding: Bisâ€triazole Pyridinium. <i>Chemistry - A European Journal</i> , 2015, 21, 13330-13335.	1.7	33



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127	Dissection of the Factors Affecting Formation of a CH <sup>δ+</sup> ...O H-Bond. A Case Study. <i>Crystals</i> , 2015, 5, 327-345.		15
128	S <sup>δ-</sup> ...X <sup>δ+</sup> Chalcogen Bonds between SF <sub>2</sub> or SF <sub>4</sub> and C≡C Multiple Bonds. <i>Journal of Physical Chemistry A</i> , 2015, 119, 5889-5897.	1.1	40
129	Substituent Effects on the Binding of Halides by Neutral and Dicationic Bis(triazolium) Receptors. <i>Journal of Physical Chemistry A</i> , 2015, 119, 13064-13073.	1.1	35
130	B-N Bond Cleavage and BN Ring Expansion at the Surface of Boron Nitride Nanotubes by Iminoborane. <i>Journal of Physical Chemistry C</i> , 2015, 119, 3253-3259.	1.5	15
131	Intramolecular S <sup>δ-</sup> ...O Chalcogen Bond as Stabilizing Factor in Geometry of Substituted Phenyl-SF <sub>3</sub> Molecules. <i>Journal of Organic Chemistry</i> , 2015, 80, 2356-2363.	1.7	61
132	Anionic CH <sup>δ+</sup> ...X <sup>δ-</sup> Hydrogen Bonds: Origin of Their Strength, Geometry, and Other Properties. <i>Chemistry - A European Journal</i> , 2015, 21, 1474-1481.	1.7	26
133	Site and chirality selective chemical modifications of boron nitride nanotubes (BNNTs) via Lewis acid-base interactions. <i>Physical Chemistry Chemical Physics</i> , 2015, 17, 3850-3866.	1.3	20
134	Chalcogen Bonds in Complexes of SOXY (X, Y = F, Cl) with Nitrogen Bases. <i>Journal of Physical Chemistry A</i> , 2015, 119, 535-541.	1.1	58
135	Frontispiece: Anionic CH <sup>δ+</sup> ...X <sup>δ-</sup> Hydrogen Bonds: Origin of Their Strength, Geometry, and Other Properties. <i>Chemistry - A European Journal</i> , 2015, 21, n/a-n/a.	1.7	0
136	Comparison of CH <sup>δ+</sup> ...O, SH <sup>δ+</sup> ...O, Chalcogen, and Tetrel Bonds Formed by Neutral and Cationic Sulfur-Containing Compounds. <i>Journal of Physical Chemistry A</i> , 2015, 119, 9189-9199.	1.1	92
137	Structure and Properties of [8]BN-Circulenes: Inorganic Analogues of [8]Circulenes. <i>Journal of Physical Chemistry C</i> , 2015, 119, 15541-15546.	1.5	11
138	Long-range behavior of noncovalent bonds. Neutral and charged H-bonds, pnictogen, chalcogen, and halogen bonds. <i>Chemical Physics</i> , 2015, 456, 34-40.	0.9	21
139	Interactions between Thiourea and Imines. Prelude to Catalysis. <i>Journal of Organic Chemistry</i> , 2015, 80, 10334-10341.	1.7	6
140	Tetrel, chalcogen, and CH <sup>δ+</sup> ...O hydrogen bonds in complexes pairing carbonyl-containing molecules with 1, 2, and 3 molecules of CO <sub>2</sub> . <i>Journal of Chemical Physics</i> , 2015, 142, 034307.	1.2	81
141	Microsolvation of anions by molecules forming CH <sup>δ+</sup> ...X <sup>δ-</sup> hydrogen bonds. <i>Chemical Physics</i> , 2015, 463, 137-144.	0.9	6
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