

Mariusz Michalczyk

List of Publications by Year in descending order

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#	ARTICLE	IF	CITATIONS
1	Implications of monomer deformation for tetrel and pnictogen bonds. <i>Physical Chemistry Chemical Physics</i> , 2018, 20, 8832-8841.	1.3	67
2	Coordination of anions by noncovalently bonded σ -hole ligands. <i>Coordination Chemistry Reviews</i> , 2020, 405, 213136.	9.5	66
3	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
4	Comparison between Tetrel Bonded Complexes Stabilized by σ and π Hole Interactions. <i>Molecules</i> , 2018, 23, 1416.	1.7	45
5	Aerogen bonds formed between $AeOF_2$ ($Ae = Kr, Xe$) and diazines: comparisons between σ -hole and π -hole complexes. <i>Physical Chemistry Chemical Physics</i> , 2018, 20, 4676-4687.	1.3	36
6	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
7	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
8	Anion-...-Anion Attraction in Complexes of MCl_3^+ ($M=Zn, Cd, Hg$) with CN^+ . <i>ChemPhysChem</i> , 2020, 21, 1119-1125.	1.0	31
9	On the ability of pnictogen atoms to engage in both σ and π -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
10	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
11	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
12	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
13	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
14	Hexacoordinated Tetrel-Bonded Complexes between TF_4 ($T=Si, Ge, Sn, Pb$) and NCH : Competition between σ and π Holes. <i>ChemPhysChem</i> , 2019, 20, 959-966.	1.0	25
15	On the Stability of Interactions between Pairs of Anions in Complexes of MCl_3^+ ($M=Be, Mg, Ca, Sr, Ba$) with Pyridine and CN^+ . <i>ChemPhysChem</i> , 2020, 21, 870-877.	1.0	25
16	Crystallographic and Theoretical Evidences of Anion-...-Anion Interaction. <i>ChemPhysChem</i> , 2021, 22, 818-821.	1.0	25
17	Noncovalent Bonds between Tetrel Atoms. <i>ChemPhysChem</i> , 2020, 21, 1934-1944.	1.0	24
18	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

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19	Influence of monomer deformation on the competition between two types of σ -holes in tetrel bonds. <i>Physical Chemistry Chemical Physics</i> , 2019, 21, 10336-10346.	1.3	20
20	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
21	Anion π -anion and anion π -neutral triel bonds. <i>Physical Chemistry Chemical Physics</i> , 2021, 23, 4818-4828.	1.3	19
22	Pnictogen bonding in pyrazine π -PnX ₅ (Pn = P, As, Sb and X = F, Cl, Br) complexes. <i>Journal of Molecular Modeling</i> , 2017, 23, 328.	0.8	18
23	Anion π -anion (MX ₃) ⁺ ₂ 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
24	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
25	Synthesis, crystal structure, DFT calculations, molecular docking study and Hirshfeld surface analysis of alkoxido-bridged dinuclear iron(III) complex. <i>Research on Chemical Intermediates</i> , 2020, 46, 4155-4171.	1.3	15
26	On the opposite trends of correlations between interaction energies and electrostatic potentials of chlorinated and methylated amine complexes stabilized by halogen bond. <i>Theoretical Chemistry Accounts</i> , 2017, 136, 1.	0.5	14
27	Experimental and Theoretical Studies of Dimers Stabilized by Two Chalcogen Bonds in the Presence of a N π -N Pnictogen Bond. <i>Journal of Physical Chemistry A</i> , 2021, 125, 657-668.	1.1	14
28	Structures of clusters surrounding ions stabilized by hydrogen, halogen, chalcogen, and pnictogen bonds. <i>Chemical Physics</i> , 2019, 524, 55-62.	0.9	13
29	Anion π -Anion Interactions in Aerogen-Bonded Complexes. Influence of Solvent Environment. <i>Molecules</i> , 2021, 26, 2116.	1.7	13
30	Nature of the interaction between ammonia derivatives and carbon disulfide. A theoretical investigation. <i>International Journal of Quantum Chemistry</i> , 2017, 117, e25369.	1.0	12
31	S π -N chalcogen bonded complexes of carbon disulfide with diazines. Theoretical study. <i>Chemical Physics</i> , 2018, 500, 37-44.	0.9	12
32	The role of hydrogen bonding in π - π stacking interactions in Ni(II) complex derived from triethanolamine: synthesis, crystal structure, antimicrobial, and DFT studies. <i>Research on Chemical Intermediates</i> , 2019, 45, 5649-5664.	1.3	11
33	Ability of Lewis Acids with Shallow σ -Holes to Engage in Chalcogen Bonds in Different Environments. <i>Molecules</i> , 2021, 26, 6394.	1.7	9
34	Synthesis, characterization, DFT optimization and anticancer evaluation of phosphane-gold(I) dithiocarbamates. <i>Journal of Molecular Structure</i> , 2020, 1218, 128486.	1.8	8
35	Competition between Inter and Intramolecular Tetrel Bonds: Theoretical Studies Complemented by CSD Survey. <i>ChemPhysChem</i> , 2021, 22, 924-934.	1.0	7
36	Experimental and theoretical evidence of attractive interactions between dianions: [PdCl ₄] ²⁻ π [PdCl ₄] ²⁻ . <i>Chemical Communications</i> , 2021, 2, 2, 13305-13308.		7

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37	Iron (III) complex exhibiting efficient catechol oxidase activity: Experimental, kinetic and theoretical approach. <i>Journal of Molecular Structure</i> , 2022, 1252, 131685.	1.8	7
38	Triel bonds within anion-anion complexes. <i>Physical Chemistry Chemical Physics</i> , 2021, 23, 25097-25106.	1.3	6
39	Theoretical modeling of argentophilic interactions in [Ag(CN) ₂] ⁻ trimer found in a copper(II) complex of cis-1,2-diaminocyclohexane (Dach), [Cu(Dach) ₂ -Ag(CN) ₂ -Cu(Dach) ₂][Ag(CN) ₂] ₃ . <i>Chemical Physics Letters</i> , 2018, 709, 11-15.	1.2	4
40	The Role of Hydrogen Bonds in Interactions between [PdCl ₄] ²⁻ Dianions in Crystal. <i>Molecules</i> , 2022, 27, 2144.	1.7	4
41	Competition between Intra and Intermolecular Pnicogen Bonds. Complexes between Naphthalene Derivatives and Neutral or Anionic Bases. <i>ChemPhysChem</i> , 2022, , .	1.0	4
42	Experimental and Theoretical Evidence of a Pb...Pb Ditetrel Bond Without a "Hole". <i>ChemPhysChem</i> , 2022, 23, .	1.0	4
43	Crystal structure and theoretical investigation of bis(cis-1,2-diaminocyclohexane)zinc(II) tetrachloridozincate(II). <i>Zeitschrift Fur Naturforschung - Section B Journal of Chemical Sciences</i> , 2017, 72, 627-630.	0.3	3
44	Interactions of (MY) ₆ (M = Zn, Cd; Y = O, S, Se) quantum dots with N-bases. <i>Structural Chemistry</i> , 2019, 30, 1003-1014.	1.0	2
45	Possible coordination modes of copper(II) atom in model silsesquioxanes complexes at various pH conditions: DFT study. <i>Chemical Physics Letters</i> , 2021, 778, 138739.	1.2	2