

Sergiy V Rosokha

List of Publications by Year in descending order

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3809
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#	ARTICLE	IF	CITATIONS
1	Editorial: Advanced Research in Halogen Bonding. <i>Crystals</i> , 2022, 12, 133.	1.0	1
2	From weak to strong interactions: structural and electron topology analysis of the continuum from the supramolecular chalcogen bonding to covalent bonds. <i>Physical Chemistry Chemical Physics</i> , 2022, 24, 8251-8259.	1.3	15
3	Solvent and Ionic Atmosphere Effects in Anion π - π Interactions: Complexes of Halide Anions with <i>p</i> -Benzoquinones. <i>Journal of Physical Chemistry A</i> , 2022, 126, 4255-4263.	1.1	3
4	“Anti-electrostatic” halogen bonding in solution. <i>Chemical Science</i> , 2021, 12, 8246-8251.	3.7	20
5	“Anti-electrostatic” Halogen Bonding between Ions of Like Charge. <i>Chemistry - A European Journal</i> , 2021, 27, 16530-16542.	1.7	24
6	Examining a Transition from Supramolecular Halogen Bonding to Covalent Bonds: Topological Analysis of Electron Densities and Energies in the Complexes of Bromosubstituted Electrophiles. <i>ACS Omega</i> , 2021, 6, 23588-23597.	1.6	14
7	Structures, Multicenter π -Bonding, and Spin Equilibria in the Mixed-Valence Trimers of Tetramethyltetrathiafulvalene Cation-Radicals. <i>Crystal Growth and Design</i> , 2021, 21, 7257-7268.	1.4	6
8	Frontispiece: “Anti-electrostatic” Halogen Bonding between Ions of Like Charge. <i>Chemistry - A European Journal</i> , 2021, 27, .	1.7	0
9	Effects of structural variations on π -dimer formation: long-distance multicenter bonding of cation-radicals of tetrathiafulvalene analogues. <i>Physical Chemistry Chemical Physics</i> , 2020, 22, 25054-25065.	1.3	8
10	Efficient energy transfer in phenyl-ethynyl-linked asymmetric BODIPY dimers. <i>Tetrahedron</i> , 2020, 76, 131515.	1.0	4
11	Halogen Bonding in the Complexes of Brominated Electrophiles with Chloride Anions: From a Weak Supramolecular Interaction to a Covalent Br \cdots Cl Bond. <i>Crystals</i> , 2020, 10, 1075.	1.0	5
12	Halogen Bonding Between Anions: Association of Anion Radicals of Tetraiodo- <i>p</i> -benzoquinone with Iodide Anions. <i>Angewandte Chemie - International Edition</i> , 2020, 59, 17197-17201.	7.2	13
13	Halogen Bonding Between Anions: Association of Anion Radicals of Tetraiodo- <i>p</i> -benzoquinone with Iodide Anions. <i>Angewandte Chemie</i> , 2020, 132, 17350-17354.	1.6	4
14	Diversity and uniformity in anion π complexes of thiocyanate with aromatic, olefinic and quinoidal π -acceptors. <i>Dalton Transactions</i> , 2020, 49, 8734-8743.	1.6	19
15	Intermolecular Interactions between Halogen-Substituted <i>p</i> -Benzoquinones and Halide Anions: Anion π Complexes versus Halogen Bonding. <i>ChemPlusChem</i> , 2020, 85, 441-449.	1.3	8
16	Molecular Bases for Anesthetic Agents: Halothane as a Halogen- and Hydrogen-Bond Donor. <i>Angewandte Chemie</i> , 2019, 131, 12586-12589.	1.6	4
17	Complexes of Diiodine with Heteroaromatic <i>N</i> -Oxides: Effects of Halogen-Bond Acceptors in Halogen Bonding. <i>Journal of Physical Chemistry A</i> , 2019, 123, 7113-7123.	1.1	21
18	Molecular Bases for Anesthetic Agents: Halothane as a Halogen- and Hydrogen-Bond Donor. <i>Angewandte Chemie - International Edition</i> , 2019, 58, 12456-12459.	7.2	10

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19	Innentitelbild: Molecular Bases for Anesthetic Agents: Halothane as a Halogen- and Hydrogen-Bond Donor (Angew. Chem. 36/2019). Angewandte Chemie, 2019, 131, 12436-12436.	1.6	0
20	Anion-Complexes of Halides with <i>p</i> -Benzoquinones: Structures, Thermodynamics, and Criteria of Charge Transfer to Electron Transfer Transition. Journal of the American Chemical Society, 2019, 141, 9338-9348.	6.6	52
21	Structural preferences in strong anion- and halogen-bonded complexes: σ - and π -holes vs. frontier orbitals interaction. New Journal of Chemistry, 2018, 42, 10572-10583.	1.4	19
22	Effects of Supramolecular Architecture on Halogen Bonding between Diiodine and Heteroaromatic N-Oxides. Crystal Growth and Design, 2018, 18, 1198-1207.	1.4	22
23	Continuum of covalent to intermolecular bonding in the halogen-bonded complexes of 1,4-diazabicyclo[2.2.2]octane with bromine-containing electrophiles. Chemical Communications, 2018, 54, 8060-8063.	2.2	29
24	Resolving the halogen vs. hydrogen bonding dichotomy in solutions: intermolecular complexes of trihalomethanes with halide and pseudohalide anions. Physical Chemistry Chemical Physics, 2018, 20, 21999-22007.	1.3	16
25	Anion- interaction in metal-organic networks formed by metal halides and tetracyanopyrazine. Journal of Molecular Structure, 2017, 1138, 129-135.	1.8	16
26	Electron-transfer reactions of halogenated electrophiles: a different look into the nature of halogen bonding. Faraday Discussions, 2017, 203, 315-332.	1.6	22
27	The halogen bond in solution: general discussion. Faraday Discussions, 2017, 203, 347-370.	1.6	5
28	Computational approaches and sigma-hole interactions: general discussion. Faraday Discussions, 2017, 203, 131-163.	1.6	17
29	Solid-state chemistry and applications: general discussion. Faraday Discussions, 2017, 203, 459-483.	1.6	2
30	From single-point to three-point halogen bonding between zinc(II) tetrathiocyanate and tetrabromomethane. CrystEngComm, 2016, 18, 488-495.	1.3	8
31	Mechanism and Thermodynamics of Reductive Cleavage of Carbon-Halogen Bonds in the Polybrominated Aliphatic Electrophiles. Journal of Physical Chemistry A, 2016, 120, 1706-1715.	1.1	13
32	From charge transfer to electron transfer in halogen-bonded complexes of electrophilic bromocarbons with halide anions. Physical Chemistry Chemical Physics, 2015, 17, 4989-4999.	1.3	21
33	Interplay of Halogen and Charge-Transfer Bondings in Intermolecular Associates of Bromo- or Iododinitrobenzene with Tetramethyl- <i>p</i> -phenylenediamine. Journal of Physical Chemistry A, 2015, 119, 3833-3842.	1.1	23
34	Halogen bond-assisted electron transfer reactions of aliphatic bromosubstituted electrophiles. Physical Chemistry Chemical Physics, 2014, 16, 1809-1813.	1.3	30
35	Halogen bonding of electrophilic bromocarbons with pseudohalide anions. Physical Chemistry Chemical Physics, 2014, 16, 12968-12979.	1.3	40
36	-Bonded molecular wires: self-assembly of mixed-valence cation-radical stacks within the nanochannels formed by inert tetrakis[3,5-bis(trifluoromethyl)phenyl]borate anions. CrystEngComm, 2013, 15, 10638.	1.3	22

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37	Experimental and Computational Probes of the Nature of Halogen Bonding: Complexes of Bromine-Containing Molecules with Bromide Anions. <i>Chemistry - A European Journal</i> , 2013, 19, 8774-8788.	1.7	109
38	Substituent-Induced Switch of the Role of Charge-Transfer Complexes in the Diels-Alder Reactions of <i>o</i> -Chloranil and Styrenes. <i>Journal of Organic Chemistry</i> , 2012, 77, 5971-5981.	1.7	10
39	Hybrid Network Formation via Halogen Bonding of the Neutral Bromo-Substituted Organic Molecules with Anionic Metal-Bromide Complexes. <i>Crystal Growth and Design</i> , 2012, 12, 4149-4156.	1.4	32
40	Intermolecular π -dimer of oxoverdazyl radicals with long-distance multicenter (2e/8c) bonding via nitrogen atoms. <i>Journal of Physical Organic Chemistry</i> , 2010, 23, 395-399.	0.9	8
41	One- and two-dimensional coordination networks of the tetracyanoethylene anion-radicals with potassium counter-ions. <i>Polyhedron</i> , 2009, 28, 4136-4140.	1.0	7
42	Unusual structural effects of intermolecular π -bonding in the tetracyanopyrazine (ion-radical) dimer. <i>New Journal of Chemistry</i> , 2009, 33, 545-553.	1.4	26
43	Lewis acid effects on donor-acceptor associations and redox reactions: ternary complexes of heteroaromatic N-oxides with boron trifluoride and organic donors. <i>New Journal of Chemistry</i> , 2009, 33, 2317.	1.4	15
44	Trimorphism of a model carcinogen 4-nitroquinoline-N-oxide. <i>CrystEngComm</i> , 2009, 11, 2400.	1.3	7
45	Counter-ion modulation of long-distance π -bonding of the open-shell p-benzoquinone anions. <i>Physical Chemistry Chemical Physics</i> , 2009, 11, 324-332.	1.3	28
46	Spectroscopic and Electrochemical Evaluation of Salt Effects on Electron-Transfer Equilibria between Donor/Acceptor and Ion-Radical Pairs in Organic Solvents. <i>ChemPhysChem</i> , 2008, 9, 2406-2413.	1.0	6
47	The Spectral Elucidation versus the X-ray Structure of the Critical Precursor Complex in Bimolecular Electron Transfers: Application of Experimental/Theoretical Solvent Probes to Ion-Radical (Redox) Dyads. <i>Journal of the American Chemical Society</i> , 2008, 130, 1944-1952.	6.6	35
48	Fresh Look at Electron-Transfer Mechanisms via the Donor/Acceptor Bindings in the Critical Encounter Complex. <i>Accounts of Chemical Research</i> , 2008, 41, 641-653.	7.6	359
49	Halogen-bonded assembly of hybrid inorganic/organic 3D-networks from dibromocuprate salts and tetrabromomethane. <i>Chemical Communications</i> , 2007, , 3383.	2.2	31
50	Reversible Interchange of Charge-Transfer versus Electron-Transfer States in Organic Electron Transfer via Cross-Exchanges between Diamagnetic (Donor/Acceptor) Dyads. <i>Journal of Physical Chemistry B</i> , 2007, 111, 6655-6666.	1.2	33
51	Molecular and Electronic Structures of the Long-Bonded π -Dimers of Tetrathiafulvalene Cation-Radical in Intermolecular Electron Transfer and in (Solid-State) Conductivity. <i>Journal of the American Chemical Society</i> , 2007, 129, 828-838.	6.6	173
52	Continuum of Outer- and Inner-Sphere Mechanisms for Organic Electron Transfer. Steric Modulation of the Precursor Complex in Paramagnetic (Ion-Radical) Self-Exchanges. <i>Journal of the American Chemical Society</i> , 2007, 129, 3683-3697.	6.6	115
53	Tris(thianthrene)(2+) bis(dodecamethylcarba-closo-dodecaborate) dichloromethane tetrasolvate: a crossed triple-decker π -trimer dication. <i>Acta Crystallographica Section C: Crystal Structure Communications</i> , 2007, 63, o347-o349.	0.4	8
54	X-ray Structures and Electronic Spectra of the π -Halogen Complexes between Halogen Donors and Acceptors with π -Receptors. , 2007, , 137-160.		37

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55	Quinones as Electron Acceptors. X-Ray Structures, Spectral (EPR, UV-vis) Characteristics and Electron-Transfer Reactivities of Their Reduced Anion Radicals as Separated vs Contact Ion Pairs. <i>Journal of the American Chemical Society</i> , 2006, 128, 16708-16719.	6.6	78
56	Electronic structures of intermolecular charge-transfer states in fast electron transfers with tetrathiafulvalene donor. Thermal and photoactivation of [2 + 4] cycloaddition to o-chloranil acceptor. <i>Photochemical and Photobiological Sciences</i> , 2006, 5, 914.	1.6	27
57	Very Fast Electron Migrations within p-Doped Aromatic Cofacial Arrays Leading to Three-Dimensional (Toroidal) π -Delocalization. <i>Journal of the American Chemical Society</i> , 2006, 128, 9394-9407.	6.6	78
58	The Question of Aromaticity in Open-Shell Cations and Anions as Ion-Radical Offsprings of Polycyclic Aromatic and Antiaromatic Hydrocarbons. <i>Journal of Organic Chemistry</i> , 2006, 71, 9357-9365.	1.7	34
59	Steric Modulations in the Reversible Dimerizations of Phenalenyl Radicals via Unusually Weak Carbon-Centered π - and σ -Bonds. <i>Journal of Organic Chemistry</i> , 2006, 71, 520-526.	1.7	87
60	2,3,4,5,6-Pentanitroaniline 1,2-dichloroethane disolvate: 'push-pull' deformation of aromatic rings by intramolecular charge transfer. <i>Acta Crystallographica Section C: Crystal Structure Communications</i> , 2006, 62, 0464-0466.	0.4	5
61	Mulliken-Hush elucidation of the encounter (precursor) complex in intermolecular electron transfer via self-exchange of tetracyanoethylene anion-radical. <i>Chemical Physics</i> , 2006, 324, 117-128.	0.9	28
62	Through-Space (Cofacial) π -Delocalization among Multiple Aromatic Centers: Toroidal Conjugation in Hexaphenylbenzene-like Radical Cations. <i>Angewandte Chemie - International Edition</i> , 2005, 44, 5133-5136.	7.2	78
63	Halide Recognition through Diagnostic Anion- π Interactions: Molecular Complexes of Cl ⁻ , Br ⁻ , and I ⁻ with Olefinic and Aromatic π Receptors. <i>Angewandte Chemie - International Edition</i> , 2005, 44, 2178-2178.	7.2	1
64	Separated versus Contact Ion-Pair Structures in Solution from Their Crystalline States: A Dynamic Effects on Dinitrobenzene as a Mixed-Valence Anion. <i>Journal of the American Chemical Society</i> , 2005, 127, 1797-1809.	6.6	50
65	Characterizing the Dimerizations of Phenalenyl Radicals by ab Initio Calculations and Spectroscopy: π -Bond Formation versus Resonance π -Stabilization. <i>Journal of Physical Chemistry A</i> , 2005, 109, 11261-11267.	1.1	90
66	Intermolecular Electron-Transfer Mechanisms via Quantitative Structures and Ion-Pair Equilibria for Self-Exchange of Anionic (Dinitrobenzene) Donors. <i>Journal of the American Chemical Society</i> , 2005, 127, 7411-7420.	6.6	29
67	Halide Recognition through Diagnostic Anion- π Interactions: Molecular Complexes of Cl ⁻ , Br ⁻ , and I ⁻ with Olefinic and Aromatic π Receptors. <i>Angewandte Chemie - International Edition</i> , 2004, 43, 4650-4652.	7.2	339
68	Intermolecular π -to- π Bonding between Stacked Aromatic Dyads. Experimental and Theoretical Binding Energies and Near-IR Optical Transitions for Phenalenyl Radical/Radical versus Radical/Cation Dimerizations. <i>Journal of the American Chemical Society</i> , 2004, 126, 13850-13858.	6.6	286
69	Donor-Acceptor (Electronic) Coupling in the Precursor Complex to Organic Electron Transfer: A Intermolecular and Intramolecular Self-Exchange between Phenothiazine Redox Centers. <i>Journal of the American Chemical Society</i> , 2004, 126, 1388-1401.	6.6	168
70	Isolation of the Latent Precursor Complex in Electron-Transfer Dynamics. Intermolecular Association and Self-Exchange with Acceptor Anion Radicals. <i>Journal of the American Chemical Society</i> , 2003, 125, 2559-2571.	6.6	110
71	Molecular Recognition of NO/NO ⁺ via Multicenter (Charge-Transfer) Binding to Bridged Diarene Donors. Effect of Structure on the Optical Transitions and Complexation Thermodynamics. <i>Journal of Organic Chemistry</i> , 2003, 68, 3947-3957.	1.7	27
72	Stable (Long-Bonded) Dimers via the Quantitative Self-Association of Different Cationic, Anionic, and Uncharged π -Radicals: Structures, Energetics, and Optical Transitions. <i>Journal of the American Chemical Society</i> , 2003, 125, 12161-12171.	6.6	263

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73	Charge-Transfer Mechanism for Electrophilic Aromatic Nitration and Nitrosation via the Convergence of (ab Initio) Molecular-Orbital and Marcus's Hush Theories with Experiments. <i>Journal of the American Chemical Society</i> , 2003, 125, 3273-3283.	6.6	88
74	Intervalence (Charge-Resonance) Transitions in Organic Mixed-Valence Systems. Through-Space versus Through-Bond Electron Transfer between Bridged Aromatic (Redox) Centers. <i>Journal of the American Chemical Society</i> , 2003, 125, 15950-15963.	6.6	111
75	X-ray Structure Analysis and the Intervalent Electron Transfer in Organic Mixed-Valence Crystals with Bridged Aromatic Cation Radicals. <i>Journal of the American Chemical Society</i> , 2002, 124, 843-855.	6.6	110
76	Conformation, Distance, and Connectivity Effects on Intramolecular Electron Transfer between Phenylene-Bridged Aromatic Redox Centers. <i>Journal of Physical Chemistry A</i> , 2002, 106, 2283-2292.	1.1	71
77	Strong electronic coupling in intermolecular (charge-transfer) complexes. Mechanistic relevance to thermal and optical electron transfer from aromatic donors. <i>New Journal of Chemistry</i> , 2002, 26, 851-860.	1.4	21
78	The Preorganization Step in Organic Reaction Mechanisms. Charge-Transfer Complexes as Precursors to Electrophilic Aromatic Substitutions. <i>Journal of Organic Chemistry</i> , 2002, 67, 1727-1737.	1.7	85
79	Novel Arene Receptors as Nitric Oxide (NO) Sensors. <i>Journal of the American Chemical Society</i> , 2002, 124, 5620-5621.	6.6	18
80	Mechanism of Inner-Sphere Electron Transfer via Charge-Transfer (Precursor) Complexes. Redox Energetics of Aromatic Donors with the Nitrosonium Acceptor. <i>Journal of the American Chemical Society</i> , 2001, 123, 8985-8999.	6.6	64
81	Charge-Transfer Effects on Arene Structure and Reactivity. , 0, , 435-478.		5