

Janusz Rak

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135
papers

2,624
citations

30
h-index

45
g-index

145
ext. papers

2,808
ext. citations

4.7
avg. IF

4.76
L-index

#	Paper	IF	Citations
135	Low-energy tautomers and conformers of neutral and protonated arginine. <i>Journal of the American Chemical Society</i> , 2001 , 123, 11695-707	16.4	128
134	Calculation of quantum-mechanical descriptors for QSPR at the DFT level: is it necessary?. <i>Journal of Chemical Information and Modeling</i> , 2008 , 48, 1174-80	6.1	104
133	Consequences of proton transfer in guanidine. <i>Journal of Physical Organic Chemistry</i> , 2003 , 16, 91-106	2.1	95
132	AT base pair anions versus (9-methyl-A)(1-methyl-T) base pair anions. <i>Journal of the American Chemical Society</i> , 2005 , 127, 6443-50	16.4	82
131	Stabilization of very rare tautomers of uracil by an excess electron. <i>Physical Chemistry Chemical Physics</i> , 2005 , 7, 2116-25	3.6	72
130	Quasidegeneracy of zwitterionic and canonical tautomers of arginine solvated by an excess electron. <i>Journal of the American Chemical Society</i> , 2001 , 123, 11073-4	16.4	61
129	Intermolecular proton transfer in anionic complexes of uracil with alcohols. <i>Journal of Physical Chemistry B</i> , 2005 , 109, 13383-91	3.4	55
128	Toward an Understanding of the Chemiluminescence Accompanying the Reaction of 9-Carboxy-10-methylacridinium Phenyl Ester with Hydrogen Peroxide. <i>Journal of Organic Chemistry</i> , 1999 , 64, 3002-3008	4.2	55
127	Properties of Closed-Shell, Octahedral, Multiply-Charged Hexafluorometallates MF ₆ ³⁻ , M = Sc, Y, La, ZrF ₆ ²⁻ , and TaF ₆ ⁻ . <i>Journal of the American Chemical Society</i> , 1996 , 118, 1173-1180	16.4	55
126	Photoelectron spectroscopy of adiabatically bound valence anions of rare tautomers of the nucleic acid bases. <i>Journal of Chemical Physics</i> , 2007 , 127, 174309	3.9	54
125	Mechanisms of Damage to DNA Labeled with Electrophilic Nucleobases Induced by Ionizing or UV Radiation. <i>Journal of Physical Chemistry B</i> , 2015 , 119, 8227-38	3.4	53
124	Excess Electron Attachment Induces Barrier-Free Proton Transfer in Binary Complexes of Uracil with H ₂ Se and H ₂ S but Not with H ₂ O. <i>Journal of Physical Chemistry B</i> , 2003 , 107, 7889-7895	3.4	53
123	Barrier-free intermolecular proton transfer induced by excess electron attachment to the complex of alanine with uracil. <i>Journal of Chemical Physics</i> , 2004 , 120, 6064-71	3.9	52
122	Fundamental mechanisms of DNA radiosensitization: damage induced by low-energy electrons in brominated oligonucleotide trimers. <i>Journal of Physical Chemistry B</i> , 2012 , 116, 9676-82	3.4	51
121	An ab initio study of the betaine anion-dipole-bound anionic state of a model zwitterion system. <i>Journal of Chemical Physics</i> , 2001 , 114, 10673-10681	3.9	48
120	Barrier-free proton transfer in anionic complex of thymine with glycine. <i>Physical Chemistry Chemical Physics</i> , 2004 , 6, 4351-4357	3.6	47
119	Computational Study of Hydrogen-Bonded Complexes between the Most Stable Tautomers of Glycine and Uracil. <i>Journal of Physical Chemistry A</i> , 2002 , 106, 7423-7433	2.8	47

118	Electron-induced elimination of the bromide anion from brominated nucleobases. A computational study. <i>Journal of Physical Chemistry B</i> , 2012 , 116, 5612-9	3.4	46
117	Excess Electron Attachment Induces Barrier-Free Proton Transfer in Anionic Complexes of Thymine and Uracil with Formic Acid. <i>Journal of Physical Chemistry B</i> , 2004 , 108, 6919-6921	3.4	44
116	How to Find Out Whether a 5-Substituted Uracil Could Be a Potential DNA Radiosensitizer. <i>Journal of Physical Chemistry Letters</i> , 2013 , 4, 2853-2857	6.4	43
115	On the unusual stability of valence anions of thymine based on very rare tautomers: A computational study. <i>Journal of Physical Chemistry B</i> , 2006 , 110, 24696-707	3.4	41
114	Hartree-Fock and Density Functional Methods and IR and NMR Spectroscopies in the Examination of Tautomerism and Features of Neutral 9-Acridinamine in Gaseous and Condensed Media. <i>Journal of Physical Chemistry A</i> , 1997 , 101, 283-292	2.8	38
113	TG-FTIR, DSC and quantum chemical studies of the thermal decomposition of quaternary methylammonium halides. <i>Chemical Physics</i> , 2006 , 324, 425-437	2.3	38
112	Valence anions in complexes of adenine and 9-methyladenine with formic acid: stabilization by intermolecular proton transfer. <i>Journal of the American Chemical Society</i> , 2007 , 129, 1216-24	16.4	37
111	Effects of intra base-pairs flexibility on hole transfer coupling in DNA. <i>Chemical Physics Letters</i> , 2006 , 429, 546-550	2.5	34
110	The origin of luminescence accompanying electrochemical reduction or chemical decomposition of peroxydisulfates. <i>Journal of Luminescence</i> , 2003 , 105, 27-34	3.8	34
109	Interaction with glycine increases stability of a mutagenic tautomer of uracil. A density functional theory study. <i>Journal of the American Chemical Society</i> , 2005 , 127, 2238-48	16.4	33
108	Valence anions of 9-methylguanine-1-methylcytosine complexes. Computational and photoelectron spectroscopy studies. <i>Journal of the American Chemical Society</i> , 2009 , 131, 2663-9	16.4	32
107	Stabilization of very rare tautomers of 1-methylcytosine by an excess electron. <i>Journal of Physical Chemistry A</i> , 2005 , 109, 11495-503	2.8	32
106	A first-principles study of electron attachment to the fully hydrated bromonucleobases. <i>Chemical Physics Letters</i> , 2014 , 595-596, 133-137	2.5	31
105	Low-energy-barrier proton transfer induced by electron attachment to the guanine...cytosine base pair. <i>ChemPhysChem</i> , 2010 , 11, 880-8	3.2	30
104	Effect of Hydrogen Bonding on Barrier-Free Proton Transfer in Anionic Complexes of Uracil with Weak Acids: (UH ⁻ CN) versus (UH ⁻ 2S) <i>Israel Journal of Chemistry</i> , 2004 , 44, 157-170	3.4	28
103	Valence anion of thymine in the DNA pi-stack. <i>Journal of the American Chemical Society</i> , 2008 , 130, 15683-7.4	6.4	26
102	Findings on the electron-attachment-induced abasic site in a DNA double helix. <i>Angewandte Chemie - International Edition</i> , 2007 , 46, 3479-81	16.4	26
101	Effect of proton transfer on the electronic coupling in DNA. <i>Chemical Physics</i> , 2006 , 325, 567-574	2.3	24

100	Theoretical studies on the prototropic tautomerism, structure, and features of acridine and 9-acridinamine free bases and their protonated forms. <i>Journal of Organic Chemistry</i> , 1992 , 57, 3720-3725 ^{4.2}	24
99	5-Thiocyanato-2-Deoxyuridine as a possible radiosensitizer: electron-induced formation of uracil-C5-thiyl radical and its dimerization. <i>Physical Chemistry Chemical Physics</i> , 2015 , 17, 16907-16	3.6 23
98	Anion of the formic acid dimer as a model for intermolecular proton transfer induced by a pi* excess electron. <i>Journal of Chemical Physics</i> , 2005 , 122, 204304	3.9 23
97	Electron-induced degradation of 8-bromo-2-Deoxyadenosine 3',5'-Diphosphate, a DNA radiosensitizing nucleotide. <i>Journal of Physical Chemistry B</i> , 2013 , 117, 8681-8	3.4 22
96	Single strand break in DNA coupled to the O-P bond cleavage. A computational study. <i>Journal of Physical Chemistry B</i> , 2011 , 115, 1911-7	3.4 22
95	The Effect of Pyrimidine Bases on the Hole-Transfer Coupling in DNA. <i>Journal of Physical Chemistry B</i> , 2002 , 106, 7919-7926	3.4 22
94	Splitting of Cyclobutane-Type Uracil Dimer Cation Radicals. Hartree-Fock, MP2, and Density Functional Studies. <i>Journal of Physical Chemistry A</i> , 1998 , 102, 7168-7175	2.8 21
93	Electron stimulated desorption of anions from native and brominated single stranded oligonucleotide trimers. <i>Journal of Chemical Physics</i> , 2012 , 136, 075101	3.9 20
92	The effect of two- and three-body interactions in ArnCO ₂ (n=1,2) on the asymmetric stretching CO ₂ coordinate: An ab initio study. <i>Journal of Chemical Physics</i> , 1997 , 106, 10215-10221	3.9 20
91	Origins and modeling of many-body exchange effects in van der Waals clusters. <i>Journal of Chemical Physics</i> , 1997 , 106, 3301-3310	3.9 20
90	Barrier-free proton transfer in the valence anion of 2-Deoxyadenosine-5-Monophosphate. II. A computational study. <i>Journal of Chemical Physics</i> , 2008 , 128, 044315	3.9 20
89	Intermolecular proton transfer induced by excess electron attachment to adenine(formic acid) _n (n=2, 3) hydrogen-bonded complexes. <i>Chemical Physics</i> , 2007 , 342, 215-222	2.3 20
88	Effect of Proton Transfer on the Anionic and Cationic Pathways of Pyrimidine Photodimer Cleavage. A Computational Study. <i>Journal of Physical Chemistry A</i> , 1999 , 103, 3569-3574	2.8 20
87	Theoretical Studies on the Structure, Stability, Ability To Undergo Internal Transformations, and Tautomerization, as Well as Reactivity, of H ₂ PPH ₂ and HPPH ₃ Molecules. <i>Journal of the American Chemical Society</i> , 1995 , 117, 2638-2648	16.4 20
86	TG-FTIR, DSC, and quantum-chemical studies on the thermal decomposition of quaternary ethylammonium halides. <i>Journal of Physical Chemistry A</i> , 2006 , 110, 5066-74	2.8 18
85	5-Selenocyanatouracil: A Potential Hypoxic Radiosensitizer. Electron Attachment Induced Formation of Selenium Centered Radical. <i>Journal of Physical Chemistry B</i> , 2017 , 121, 6139-6147	3.4 17
84	Electrophilic 5-Substituted Uracils as Potential Radiosensitizers: A Density Functional Theory Study. <i>ChemPhysChem</i> , 2016 , 17, 2572-8	3.2 17
83	Photoinduced single strand breaks and intrastrand cross-links in an oligonucleotide labeled with 5-bromouracil. <i>Journal of Physical Chemistry B</i> , 2014 , 118, 5009-16	3.4 17

82	Benign decay vs. photolysis in the photophysics and photochemistry of 5-bromouracil. A computational study. <i>Journal of Physical Chemistry A</i> , 2009 , 113, 5489-95	2.8	17
81	Theoretical studies on structure, thermochemistry, vibrational spectroscopy, and other features of ZrX ₂ B (X=F,Cl,Br,I): Coulombic energy in inorganic and organic hexahalogenozirconates. <i>Journal of Chemical Physics</i> , 1994 , 100, 5810-5820	3.9	17
80	Is 9-acridinamine anion a dispersion-bound anion?. <i>Journal of Chemical Physics</i> , 2001 , 115, 11193-11199	3.9	16
79	Electron induced single strand break and cyclization: a DFT study on the radiosensitization mechanism of the nucleotide of 8-bromoguanine. <i>Physical Chemistry Chemical Physics</i> , 2014 , 16, 6568-74	3.6	15
78	Visible-Light Photocatalytic Activity of Ionic Liquid TiO ₂ Spheres: Effect of the Ionic Liquid Anion Structure. <i>ChemCatChem</i> , 2017 , 9, 4377-4388	5.2	14
77	Electron-induced single strand break in the nucleotide of 5- and 6-bromouridine. A DFT study. <i>Chemical Physics Letters</i> , 2014 , 612, 289-294	2.5	14
76	An ESR and DFT study of hydration of the 2'-deoxyuridine-5-yl radical: a possible hydroxyl radical intermediate. <i>Chemical Communications</i> , 2014 , 50, 14605-8	5.8	14
75	Stable Valence Anions of Nucleic Acid Bases and DNA Strand Breaks Induced by Low Energy Electrons. <i>Challenges and Advances in Computational Chemistry and Physics</i> , 2008 , 619-667	0.7	14
74	An ab initio study of (H ₃ B←NH ₃) ⁺ dipole-bound anion supported by the dative charge-transfer bond in the neutral host. <i>Journal of Chemical Physics</i> , 2000 , 113, 8961-8968	3.9	14
73	UV-Induced Strand Breaks in Double-Stranded DNA Labeled with 5-Bromouracil: Frank or Secondary?. <i>Journal of Physical Chemistry Letters</i> , 2013 , 4, 4014-4018	6.4	13
72	Electron-Induced Dissociation of the Potential Radiosensitizer 5-Selenocyanato-2'-deoxyuridine. <i>Journal of Physical Chemistry B</i> , 2019 , 123, 1274-1282	3.4	12
71	5-Selenocyanato and 5-trifluoromethanesulfonyl derivatives of 2'-deoxyuridine: synthesis, radiation and computational chemistry as well as cytotoxicity.. <i>RSC Advances</i> , 2018 , 8, 21378-21388	3.7	12
70	Structure, Properties, Thermodynamics, and Isomerization Ability of 9-Acridinones. <i>Journal of Physical Chemistry A</i> , 2002 , 106, 3957-3963	2.8	12
69	Theoretical Studies on the Structure, Thermochemistry, Vibrational Spectroscopy, and Other Features of HfX ₆ ²⁻ (X = F, Cl, Br, I). Electrostatic Energy in Hexahalogenohafnates. <i>Inorganic Chemistry</i> , 1994 , 33, 6187-6193	5.1	12
68	5-Bromo-2'-deoxycytidine-a potential DNA photosensitizer. <i>Organic and Biomolecular Chemistry</i> , 2016 , 14, 9312-9321	3.9	11
67	Stability of the valence anion of cytosine is governed by nucleobases sequence in the double stranded DNA pi-stack: A computational study. <i>Journal of Chemical Physics</i> , 2009 , 131, 085103	3.9	11
66	NMR and DFT investigations of the substituent and solvent effect on amino-imino tautomerism in acridin-9-amines substituted at the exocyclic nitrogen atom. <i>Journal of Physical Organic Chemistry</i> , 2005 , 18, 870-879	2.1	11
65	PCR synthesis of double stranded DNA labeled with 5-bromouridine. A step towards finding a bromonucleoside for clinical trials. <i>Journal of Pharmaceutical and Biomedical Analysis</i> , 2011 , 56, 671-7	3.5	10

64	Theoretical Studies on the Effect of the Medium on Tautomeric Phenomena in Neutral and Protonated Acridin-9-amine. Mechanism of Tautomerization in Neutral Entities. <i>Australian Journal of Chemistry</i> , 1997 , 50, 97	1.2	10
63	The Transformation Mechanism of 3,4,6-Tri-O-acetyl-1,5-anhydro-2-deoxy-D-arabino-hex-1-enitol in Water. <i>Journal of Organic Chemistry</i> , 1996 , 61, 2988-2994	4.2	10
62	Molecular features of thymidine analogues governing the activity of human thymidine kinase. <i>Structural Chemistry</i> , 2018 , 29, 1367-1374	1.8	9
61	The radiosensitivity of 5- and 6-bromocytidine derivatives--electron induced DNA degradation. <i>Physical Chemistry Chemical Physics</i> , 2014 , 16, 19424-8	3.6	9
60	Presolvated low energy electron attachment to peptide methyl esters in aqueous solution: C-O bond cleavage at 77 K. <i>Journal of Physical Chemistry B</i> , 2013 , 117, 2872-7	3.4	9
59	Valence anions of N-acetylproline in the gas phase: computational and anion photoelectron spectroscopic studies. <i>Journal of Chemical Physics</i> , 2011 , 135, 114301	3.9	9
58	Thermal behaviour and thermochemistry of hexachlorozirconates of mononitrogen aromatic bases. <i>Thermochimica Acta</i> , 1993 , 230, 269-292	2.9	9
57	Radiation damage to single stranded oligonucleotide trimers labelled with 5-iodopyrimidines. <i>Organic and Biomolecular Chemistry</i> , 2016 , 14, 9331-9337	3.9	9
56	The anionic (9-methyladenine)-(1-methylthymine) base pair solvated by formic acid. A computational and photoelectron spectroscopy study. <i>Journal of Physical Chemistry B</i> , 2010 , 114, 11353-62	3.4	8
55	Local excitation of the 5-bromouracil chromophore in DNA. Computational and UV spectroscopic studies. <i>Journal of Physical Chemistry B</i> , 2011 , 115, 4532-7	3.4	8
54	Photoelectron spectroscopic studies of 5-halouracil anions. <i>Journal of Chemical Physics</i> , 2011 , 134, 015101	3.1	8
53	Can an excess electron localize on a purine moiety in the adenine-thymine Watson-Crick base pair? A computational study. <i>International Journal of Quantum Chemistry</i> , 2007 , 107, 2224-2232	2.1	8
52	Theoretical Studies on the Geometry, Thermochemistry, Vibrational Spectroscopy, and Charge Distribution in TiX ₆ ²⁻ (X = F, Cl, Br, I). Coulombic Energy in hexahalogenotitanate Lattices. <i>The Journal of Physical Chemistry</i> , 1994 , 98, 6280-6286		8
51	Photoinduced electron transfer in 5-bromouracil labeled DNA. A contrathermodynamic mechanism revisited by electron transfer theories. <i>Physical Chemistry Chemical Physics</i> , 2019 , 21, 4387-4393	3.6	7
50	5-Iodo-4-thio-2-Deoxyuridine as a Sensitizer of X-ray Induced Cancer Cell Killing. <i>International Journal of Molecular Sciences</i> , 2019 , 20,	6.3	7
49	Barrier-free proton transfer induced by electron attachment to the complexes between 1-methylcytosine and formic acid. <i>Molecular Physics</i> , 2010 , 108, 2621-2631	1.7	7
48	Dominant Pathways of Adenosyl Radical-Induced DNA Damage Revealed by QM/MM Metadynamics. <i>Journal of Chemical Theory and Computation</i> , 2017 , 13, 6415-6423	6.4	6
47	Dipole-bound and dispersion-bound anions supported by the asymmetric tautomers of aminophosphine: H ₃ NPH and HNPH ₃ . <i>Chemical Physics</i> , 2002 , 279, 101-110	2.3	6

46	Infrared and Raman spectroscopy of 9-acridinones. <i>Vibrational Spectroscopy</i> , 2001 , 27, 139-152	2.1	6
45	Prototropic tautomerism in N,N-dimethyl-N ⁹ -(1-nitro-9-acridyl)propane-1,3-diamine and its nitro isomers. Application of MNDO and PPP methods for the examination of structure and electronic absorption spectra. <i>Journal of the Chemical Society Perkin Transactions II</i> , 1990 , 1501-1508		6
44	Why Does the Type of Halogen Atom Matter for the Radiosensitizing Properties of 5-Halogen Substituted 4-Thio-2-Deoxyuridines?. <i>Molecules</i> , 2019 , 24,	4.8	5
43	Design, synthesis and biological evaluation of betulin-3-yl 2-amino-2-deoxy-β-D-glycopyranosides. <i>Bioorganic Chemistry</i> , 2020 , 96, 103568	5.1	5
42	Cytotoxicity of doxorubicin conjugated with C60 fullerene. Structural and in vitro studies. <i>Structural Chemistry</i> , 2019 , 30, 2327-2338	1.8	5
41	Photoelectron spectroscopic and density functional theoretical studies of the 2-Deoxycytidine homodimer radical anion. <i>Journal of Chemical Physics</i> , 2013 , 139, 075101	3.9	5
40	Theoretical studies on interactions between low energy electrons and protein-DNA fragments: valence anions of AT-amino acids side chain complexes. <i>Physical Chemistry Chemical Physics</i> , 2011 , 13, 19499-507	3.6	5
39	Radicals formed in N-acetylproline by electron attachment: electron spin resonance spectroscopy and computational studies. <i>Journal of Physical Chemistry B</i> , 2011 , 115, 14846-51	3.4	5
38	X-Ray, Quantum Mechanics and Density Functional Methods in the Examination of Structure and Tautomerism of N-Methyl-Substituted Acridin-9-amine Derivatives. <i>Australian Journal of Chemistry</i> , 1998 , 51, 643	1.2	5
37	Thermochemistry, lattice energetics and stability of hexahalogenohafnates. <i>Journal of Alloys and Compounds</i> , 1994 , 210, 63-70	5.7	5
36	Theoretical and Experimental Studies on the Visible Light Activity of TiO ₂ Modified with Halide-Based Ionic Liquids. <i>Catalysts</i> , 2020 , 10, 371	4	4
35	Uracil-5-yl -Sulfamate: An Illusive Radiosensitizer. Pitfalls in Modeling the Radiosensitizing Derivatives of Nucleobases. <i>Journal of Physical Chemistry B</i> , 2020 , 124, 5600-5613	3.4	4
34	Modifications at the C(5) position of pyrimidine nucleosides. <i>Russian Chemical Reviews</i> , 2020 , 89, 281-310	6.8	4
33	Quantitative assay of photoinduced DNA strand breaks by real-time PCR. <i>Journal of Pharmaceutical and Biomedical Analysis</i> , 2016 , 128, 480-484	3.5	4
32	DHPLC and MS studies of a photoinduced intrastrand cross-link in DNA labeled with 5-bromo-2-Deoxyuridine. <i>Journal of Photochemistry and Photobiology B: Biology</i> , 2014 , 130, 86-92	6.7	4
31	Artificial plasmid labeled with 5-bromo-2-Deoxyuridine: a universal molecular system for strand break detection. <i>ChemBioChem</i> , 2014 , 15, 1409-12	3.8	4
30	Photoelectron spectroscopy and density functional theory studies on the uridine homodimer radical anions. <i>Journal of Chemical Physics</i> , 2012 , 137, 205101	3.9	4
29	IR, Raman, NMR and density functional methods in the examination of tautomerism and features of N-methyl substituted 9-acridinamine derivatives. <i>Journal of Molecular Structure</i> , 1999 , 476, 45-55	3.4	4

28	A cyclic intermediate of the splitting reaction of cyclobutane-type pyrimidine dimer cation radicals. A computational finding as challenge for experimental techniques. <i>Computational and Theoretical Chemistry</i> , 1999 , 488, 163-168		4
27	Thermal features and thermochemistry of hexachlorozirconates of aliphatic and aromatic mono-amines. Stability of hexahalogenozirconates. <i>Journal of Alloys and Compounds</i> , 1995 , 224, 1-13	5-7	4
26	Crystal Structure of 9(10-Methyl)-Acridinimine Hydriodide. Lattice Energetics of this Compound and Halide Salts of Nitrogen Organic Bases. <i>Molecular Crystals and Liquid Crystals</i> , 1996 , 276, 91-104		4
25	2,6-diaminopurine promotes repair of DNA lesions under prebiotic conditions. <i>Nature Communications</i> , 2021 , 12, 3018	17.4	4
24	Excess Electron Attachment to the Nucleoside Pair 2-Deoxyadenosine (dA)-2-Deoxythymidine (dT). <i>Journal of Physical Chemistry B</i> , 2016 , 120, 4955-62	3-4	4
23	Chemically-enzymatic synthesis of photosensitive DNA. <i>Journal of Photochemistry and Photobiology B: Biology</i> , 2017 , 167, 228-235	6-7	3
22	Photoelectron spectroscopy and computational modeling of thymidine homodimer anions. <i>Journal of Physical Chemistry B</i> , 2012 , 116, 13975-81	3-4	3
21	Enzymatic synthesis of long double-stranded DNA labeled with haloderivatives of nucleobases in a precisely pre-determined sequence. <i>BMC Biochemistry</i> , 2011 , 12, 47	4-8	3
20	Energetics of the splitting of pyrimidine photodimers induced by electron transfer to rhodium(III) complexes. A quantum chemical study. <i>International Journal of Quantum Chemistry</i> , 2000 , 77, 128-138	2-1	3
19	CGC, MS and theoretical studies on the transformation mechanism of 3,4-di-O-acetyl-1,5-anhydro-2-deoxy-D-threo-pent-1-enitol in aqueous solutions. <i>Journal of the Chemical Society Perkin Transactions II</i> , 1995 , 569-575		3
18	Thermal properties, crystal lattice energy, mechanism and energetics of the thermal decomposition of hydrochlorides of 2-amino acid esters. <i>Thermochimica Acta</i> , 1990 , 171, 253-277	2-9	3
17	The Product of Matrix Metalloproteinase Cleavage of Doxorubicin Conjugate for Anticancer Drug Delivery: Calorimetric, Spectroscopic, and Molecular Dynamics Studies on Peptide-Doxorubicin Binding to DNA. <i>International Journal of Molecular Sciences</i> , 2020 , 21,	6-3	3
16	UV-induced electron transfer between triethylamine and 5-bromo-2-deoxyuridine. A puzzle concerning the photochemical debromination of labeled DNA. <i>Journal of Pharmaceutical and Biomedical Analysis</i> , 2017 , 142, 262-269	3-5	2
15	Reactivity pattern of bromonucleosides induced by 2-hydroxypropyl radicals: photochemical, radiation chemical, and computational studies. <i>Journal of Physical Chemistry B</i> , 2015 , 119, 6545-54	3-4	2
14	The Sequence Dependence of Photoinduced Single Strand Break in 5-Bromo-2-Deoxyuridine Labeled DNA Supports That Electron Transfer Is Responsible for the Damage. <i>Journal of Physical Chemistry B</i> , 2017 , 121, 9169-9174	3-4	2
13	Valence Anions of DNA-Related Systems in the Gas Phase: Computational and Anion Photoelectron Spectroscopy Studies 2014 , 323-392		2
12	Absorption and luminescence spectroscopic analysis of tautomeric forms of protonated N,N-dimethyl-N'-(1-nitro-9-acridinyl)-1,3-propanediamine (nitracrine) and its nitro isomers in poly(vinyl alcohol) films. <i>Journal of Fluorescence</i> , 1991 , 1, 57-68	2-4	2
11	Photoelectron Spectroscopy and Theoretical Investigations of Gaseous Doubly Deprotonated 2-Deoxynucleoside 5-Monophosphate Dianions. <i>Journal of Physical Chemistry Letters</i> , 2021 , 12, 9463-9469	6-4	2

10	Unexpected photoproduct generated via the acetone-sensitized photolysis of 5-bromo-2-deoxyuridine in a water/isopropanol solution: experimental and computational studies. <i>Journal of Physical Chemistry B</i> , 2010 , 114, 16902-7	3.4	1
9	Consequences of Electron Attachment to Modified Nucleosides Incorporated into DNA 2015 , 1-22		1
8	Guanosine Dianions Hydrated by One to Four Water Molecules.. <i>Journal of Physical Chemistry Letters</i> , 2022 , 3230-3236	6.4	1
7	DNA Damage Radiosensitizers Geared Towards Hydrated Electrons 2022 , 125-169		0
6	Low-Energy Electron Induced Reactions in Metronidazole at Different Solvation Conditions. <i>Pharmaceuticals</i> , 2022 , 15, 701	5.2	0
5	Consequences of Electron Attachment to Modified Nucleosides Incorporated into DNA 2017 , 1895-1916		
4	Why Does the Type of Halogen Atom Matter for Radiosensitizing Properties of 5-Substituted 4-Thio-2-Deoxyuridines?. <i>Proceedings (mdpi)</i> , 2019 , 22, 40	0.3	
3	Theoretical Approach in Explanation of Energy Donor Properties of 1,4-Dioxane and 1,4-Dioxane-Water Complexes. <i>Zeitschrift Fur Naturforschung - Section A Journal of Physical Sciences</i> , 1988 , 43, 621-626	1.4	
2	Preliminary Observations on the Dependence of Potential Energy Surfaces on Intramolecular Degrees of Freedom 2000 , 73-82		
1	Electrophilic Properties of 2-Deoxyadenosine-Thymine Dimer: Photoelectron Spectroscopy and DFT Studies. <i>Journal of Physical Chemistry A</i> , 2021 , 125, 6591-6599	2.8	