

# Janusz Rak

## List of Publications by Year in descending order

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145  
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docs citations

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times ranked

1930  
citing authors

#	ARTICLE	IF	CITATIONS
1	Low-Energy Tautomers and Conformers of Neutral and Protonated Arginine. <i>Journal of the American Chemical Society</i> , 2001, 123, 11695-11707.	13.7	133
2	Calculation of Quantum-Mechanical Descriptors for QSPR at the DFT Level: Is It Necessary?. <i>Journal of Chemical Information and Modeling</i> , 2008, 48, 1174-1180.	5.4	110
3	Consequences of proton transfer in guanidine. <i>Journal of Physical Organic Chemistry</i> , 2003, 16, 91-106.	1.9	109
4	AT Base Pair Anions versus (9-Methyl-A)(1-Methyl-T) Base Pair Anions. <i>Journal of the American Chemical Society</i> , 2005, 127, 6443-6450.	13.7	84
5	Stabilization of very rare tautomers of uracil by an excess electron. <i>Physical Chemistry Chemical Physics</i> , 2005, 7, 2116.	2.8	73
6	Mechanisms of Damage to DNA Labeled with Electrophilic Nucleobases Induced by Ionizing or UV Radiation. <i>Journal of Physical Chemistry B</i> , 2015, 119, 8227-8238.	2.6	73
7	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.	3.2	67
8	Quasidegeneracy of Zwitterionic and Canonical Tautomers of Arginine Solvated by an Excess Electron. <i>Journal of the American Chemical Society</i> , 2001, 123, 11073-11074.	13.7	64
9	Properties of Closed-Shell, Octahedral, Multiply-Charged Hexafluorometallates MF <sub>6</sub> <sup>3-</sup> , M = Sc, Y, La, ZrF <sub>6</sub> <sup>2-</sup> , and TaF <sub>6</sub> <sup>-</sup> . <i>Journal of the American Chemical Society</i> , 1996, 118, 1173-1180.	13.7	59
10	Photoelectron spectroscopy of adiabatically bound valence anions of rare tautomers of the nucleic acid bases. <i>Journal of Chemical Physics</i> , 2007, 127, 174309.	3.0	59
11	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.	4.6	59
12	Fundamental Mechanisms of DNA Radiosensitization: Damage Induced by Low-Energy Electrons in Brominated Oligonucleotide Trimers. <i>Journal of Physical Chemistry B</i> , 2012, 116, 9676-9682.	2.6	57
13	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, 6064-6071.	3.0	55
14	Intermolecular Proton Transfer in Anionic Complexes of Uracil with Alcohols. <i>Journal of Physical Chemistry B</i> , 2005, 109, 13383-13391.	2.6	55
15	Excess Electron Attachment Induces Barrier-Free Proton Transfer in Binary Complexes of Uracil with H <sub>2</sub> Se and H <sub>2</sub> S but Not with H <sub>2</sub> O. <i>Journal of Physical Chemistry B</i> , 2003, 107, 7889-7895.	2.6	53
16	Electron-Induced Elimination of the Bromide Anion from Brominated Nucleobases. A Computational Study. <i>Journal of Physical Chemistry B</i> , 2012, 116, 5612-5619.	2.6	52
17	An ab initio study of the betaine anion's dipole-bound anionic state of a model zwitterion system. <i>Journal of Chemical Physics</i> , 2001, 114, 10673-10681.	3.0	49
18	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.	2.5	49

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19	Barrier-free proton transfer in anionic complex of thymine with glycine. <i>Physical Chemistry Chemical Physics</i> , 2004, 6, 4351-4357.	2.8	49
20	The origin of luminescence accompanying electrochemical reduction or chemical decomposition of peroxydisulfates. <i>Journal of Luminescence</i> , 2003, 105, 27-34.	3.1	44
21	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.	2.6	44
22	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, 24696-24707.	2.6	44
23	TG-FTIR, DSC and quantum chemical studies of the thermal decomposition of quaternary methylammonium halides. <i>Chemical Physics</i> , 2006, 324, 425-437.	1.9	40
24	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.	2.5	39
25	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, 1216-1224.	13.7	37
26	Effects of intra base-pairs flexibility on hole transfer coupling in DNA. <i>Chemical Physics Letters</i> , 2006, 429, 546-550.	2.6	35
27	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, 2238-2248.	13.7	34
28	Valence Anions of 9-Methylguanine-1-Methylcytosine Complexes. Computational and Photoelectron Spectroscopy Studies. <i>Journal of the American Chemical Society</i> , 2009, 131, 2663-2669.	13.7	33
29	Stabilization of Very Rare Tautomers of 1-Methylcytosine by an Excess Electron. <i>Journal of Physical Chemistry A</i> , 2005, 109, 11495-11503.	2.5	32
30	A first-principles study of electron attachment to the fully hydrated bromonucleobases. <i>Chemical Physics Letters</i> , 2014, 595-596, 133-137.	2.6	32
31	Low-Energy Barrier Proton Transfer Induced by Electron Attachment to the Guanine...Cytosine Base Pair. <i>ChemPhysChem</i> , 2010, 11, 880-888.	2.1	31
32	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, 16907-16916.	2.8	29
33	Effect of Hydrogen Bonding on Barrier-Free Proton Transfer in Anionic Complexes of Uracil with Weak Acids: (Uracil-HCN) versus (Uracil-H <sub>2</sub> S). <i>Israel Journal of Chemistry</i> , 2004, 44, 157-170.	2.3	28
34	Findings on the Electron-Attachment-Induced Abasic Site in a DNA Double Helix. <i>Angewandte Chemie - International Edition</i> , 2007, 46, 3479-3481.	13.8	27
35	Effect of proton transfer on the electronic coupling in DNA. <i>Chemical Physics</i> , 2006, 325, 567-574.	1.9	26
36	Valence Anion of Thymine in the DNA $\pi$ -Stack. <i>Journal of the American Chemical Society</i> , 2008, 130, 15683-15687.	13.7	26

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37	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.	3.2	25
38	The Effect of Pyrimidine Bases on the Hole-Transfer Coupling in DNA. <i>Journal of Physical Chemistry B</i> , 2002, 106, 7919-7926.	2.6	24
39	2,6-diaminopurine promotes repair of DNA lesions under prebiotic conditions. <i>Nature Communications</i> , 2021, 12, 3018.	12.8	24
40	Theoretical Studies on the Structure, Stability, Ability To Undergo Internal Transformations, and Tautomerization, as Well as Reactivity, of H <sub>2</sub> PPH <sub>2</sub> and HPPH <sub>3</sub> Molecules. <i>Journal of the American Chemical Society</i> , 1995, 117, 2638-2648.	13.7	23
41	Anion of the formic acid dimer as a model for intermolecular proton transfer induced by a $\dot{\text{I}}\text{C}^*$ excess electron. <i>Journal of Chemical Physics</i> , 2005, 122, 204304.	3.0	23
42	Single Strand Break in DNA Coupled to the O <sup>3</sup> P Bond Cleavage. A Computational Study. <i>Journal of Physical Chemistry B</i> , 2011, 115, 1911-1917.	2.6	23
43	Electron-Induced Degradation of 8-Bromo-2'-deoxyadenosine 3',5'-Diphosphate, a DNA Radiosensitizing Nucleotide. <i>Journal of Physical Chemistry B</i> , 2013, 117, 8681-8688.	2.6	23
44	The effect of two- and three-body interactions in Ar <sub>n</sub> CO <sub>2</sub> (n=1,2) on the asymmetric stretching CO <sub>2</sub> coordinate: An ab initio study. <i>Journal of Chemical Physics</i> , 1997, 106, 10215-10221.	3.0	22
45	Electron stimulated desorption of anions from native and brominated single stranded oligonucleotide trimers. <i>Journal of Chemical Physics</i> , 2012, 136, 075101.	3.0	22
46	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.	2.6	22
47	Electron-Induced Dissociation of the Potential Radiosensitizer 5-Selenocyanato-2'-deoxyuridine. <i>Journal of Physical Chemistry B</i> , 2019, 123, 1274-1282.	2.6	22
48	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.	2.5	21
49	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.	2.5	21
50	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.	3.0	21
51	Visible-Light Photocatalytic Activity of Ionic Liquid TiO <sub>2</sub> Spheres: Effect of the Ionic Liquid's Anion Structure. <i>ChemCatChem</i> , 2017, 9, 4377-4388.	3.7	21
52	Origins and modeling of many-body exchange effects in van der Waals clusters. <i>Journal of Chemical Physics</i> , 1997, 106, 3301-3310.	3.0	20
53	Intermolecular proton transfer induced by excess electron attachment to adenine(formic acid) <sub>n</sub> (n=2, 3). <i>Journal of Chemical Physics</i> , 2019, 150, 074314.	1.9	20
54	Electrophilic 5-Substituted Uracils as Potential Radiosensitizers: A Density Functional Theory Study. <i>ChemPhysChem</i> , 2016, 17, 2572-2578.	2.1	20

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55	TG-FTIR, DSC, and Quantum-Chemical Studies on the Thermal Decomposition of Quaternary Ethylammonium Halides. <i>Journal of Physical Chemistry A</i> , 2006, 110, 5066-5074.	2.5	19
56	Benign Decay vs. Photolysis in the Photophysics and Photochemistry of 5-Bromouracil. A Computational Study. <i>Journal of Physical Chemistry A</i> , 2009, 113, 5489-5495.	2.5	19
57	Theoretical studies on structure, thermochemistry, vibrational spectroscopy, and other features of $ZrX_2\hat{\sim}6$ (X=F, Cl, Br, I): Coulombic energy in inorganic and organic hexahalogenozirconates. <i>Journal of Chemical Physics</i> , 1994, 100, 5810-5820.	3.0	18
58	Is 9-acridinamine anion a dispersion-bound anion?. <i>Journal of Chemical Physics</i> , 2001, 115, 11193-11199.	3.0	18
59	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, 6568-6574.	2.8	17
60	Photoinduced Single Strand Breaks and Intrastrand Cross-Links in an Oligonucleotide Labeled with 5-Bromouracil. <i>Journal of Physical Chemistry B</i> , 2014, 118, 5009-5016.	2.6	17
61	5-Selenocyanato and 5-trifluoromethanesulfonyl derivatives of 2 $\hat{\sim}$ -deoxyuridine: synthesis, radiation and computational chemistry as well as cytotoxicity. <i>RSC Advances</i> , 2018, 8, 21378-21388.	3.6	16
62	Stable Valence Anions of Nucleic Acid Bases and DNA Strand Breaks Induced by Low Energy Electrons. Challenges and Advances in Computational Chemistry and Physics, 2008, , 619-667.	0.6	15
63	An ESR and DFT study of hydration of the 2 $\hat{\sim}$ -deoxyuridine-5-yl radical: a possible hydroxyl radical intermediate. <i>Chemical Communications</i> , 2014, 50, 14605-14608.	4.1	15
64	Anab initiostudy of (H3B $\hat{\dagger}$ NH3) $\hat{\sim}$ a dipole-bound anion supported by the dative charge-transfer bond in the neutral host. <i>Journal of Chemical Physics</i> , 2000, 113, 8961-8968.	3.0	14
65	Structure, Properties, Thermodynamics, and Isomerization Ability of 9-Acridinones. <i>Journal of Physical Chemistry A</i> , 2002, 106, 3957-3963.	2.5	14
66	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.	2.6	14
67	Design, synthesis and biological evaluation of betulin-3-yl 2-amino-2-deoxy- $\hat{\dagger}$ -d-glycopyranosides. <i>Bioorganic Chemistry</i> , 2020, 96, 103568.	4.1	14
68	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.	4.0	13
69	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.	4.6	13
70	5-Bromo-2 $\hat{\sim}$ -deoxycytidine $\hat{\sim}$ a potential DNA photosensitizer. <i>Organic and Biomolecular Chemistry</i> , 2016, 14, 9312-9321.	2.8	13
71	5-Iodo-4-thio-2 $\hat{\sim}$ -Deoxyuridine as a Sensitizer of X-ray Induced Cancer Cell Killing. <i>International Journal of Molecular Sciences</i> , 2019, 20, 1308.	4.1	13
72	Uracil-5-yl O-Sulfamate: An Illusive Radiosensitizer. Pitfalls in Modeling the Radiosensitizing Derivatives of Nucleobases. <i>Journal of Physical Chemistry B</i> , 2020, 124, 5600-5613.	2.6	13

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73	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.	1.9	12
74	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, 671-677.	2.8	12
75	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.	3.0	11
76	The radiosensitivity of 5- and 6-bromocytidine derivatives $\alpha$ electron induced DNA degradation. <i>Physical Chemistry Chemical Physics</i> , 2014, 16, 19424.	2.8	11
77	Radiation damage to single stranded oligonucleotide trimers labelled with 5-iodopyrimidines. <i>Organic and Biomolecular Chemistry</i> , 2016, 14, 9331-9337.	2.8	11
78	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.	3.2	10
79	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.	0.9	10
80	Molecular features of thymidine analogues governing the activity of human thymidine kinase. <i>Structural Chemistry</i> , 2018, 29, 1367-1374.	2.0	10
81	Cytotoxicity of doxorubicin conjugated with C60 fullerene. Structural and in vitro studies. <i>Structural Chemistry</i> , 2019, 30, 2327-2338.	2.0	10
82	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, 2819.	3.8	10
83	Thermal behaviour and thermochemistry of hexachlorozirconates of mononitrogen aromatic bases. <i>Thermochimica Acta</i> , 1993, 230, 269-292.	2.7	9
84	Valence anions of N-acetylproline in the gas phase: Computational and anion photoelectron spectroscopic studies. <i>Journal of Chemical Physics</i> , 2011, 135, 114301.	3.0	9
85	Photoelectron spectroscopic studies of 5-halouracil anions. <i>Journal of Chemical Physics</i> , 2011, 134, 015101.	3.0	9
86	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, 2872-2877.	2.6	9
87	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.	2.8	9
88	Modifications at the C(5) position of pyrimidine nucleosides. <i>Russian Chemical Reviews</i> , 2020, 89, 281-310.	6.5	9
89	Prototropic tautomerism in N,N-dimethyl-N <sup>2</sup> -(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.	0.9	8
90	Theoretical Studies on the Geometry, Thermochemistry, Vibrational Spectroscopy, and Charge Distribution in TiX <sub>6</sub> <sup>2-</sup> (X = F, Cl, Br, I). Coulombic Energy in hexahalogenotitanate Lattices. <i>The Journal of Physical Chemistry</i> , 1994, 98, 6280-6286.	2.9	8

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91	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.	2.0	8
92	The Anionic (9-Methyladenine) <sup>+</sup> (1-Methylthymine) Base Pair Solvated by Formic Acid. A Computational and Photoelectron Spectroscopy Study. <i>Journal of Physical Chemistry B</i> , 2010, 114, 11353-11362.	2.6	8
93	Local Excitation of the 5-Bromouracil Chromophore in DNA. Computational and UV Spectroscopic Studies. <i>Journal of Physical Chemistry B</i> , 2011, 115, 4532-4537.	2.6	8
94	Electron-Induced Decomposition of Uracil-5-yl O-(N,N-dimethylsulfamate): Role of Methylation in Molecular Stability. <i>International Journal of Molecular Sciences</i> , 2021, 22, 2344.	4.1	8
95	Development of Sulfamoylated 4-(1-Phenyl-1 <i>H</i> -1,2,3-triazol-4-yl)phenol Derivatives as Potent Steroid Sulfatase Inhibitors for Efficient Treatment of Breast Cancer. <i>Journal of Medicinal Chemistry</i> , 2022, 65, 5044-5056.	6.4	8
96	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.	0.3	7
97	Infrared and Raman spectroscopy of 9-acridinones. <i>Vibrational Spectroscopy</i> , 2001, 27, 139-152.	2.2	7
98	Dipole-bound and dispersion-bound anions supported by the asymmetric tautomers of aminophosphine: H <sub>3</sub> NPH and HNPH <sub>3</sub> . <i>Chemical Physics</i> , 2002, 279, 101-110.	1.9	7
99	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
100	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, 6923.	4.1	7
101	Radicals Formed in N-Acetylproline by Electron Attachment: Electron Spin Resonance Spectroscopy and Computational Studies. <i>Journal of Physical Chemistry B</i> , 2011, 115, 14846-14851.	2.6	6
102	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.	5.3	6
103	Theoretical and Experimental Studies on the Visible Light Activity of TiO <sub>2</sub> Modified with Halide-Based Ionic Liquids. <i>Catalysts</i> , 2020, 10, 371.	3.5	6
104	Influence of Hypoxia on Radiosensitization of Cancer Cells by 5-Bromo-2-deoxyuridine. <i>International Journal of Molecular Sciences</i> , 2022, 23, 1429.	4.1	6
105	Thermochemistry, lattice energetics and stability of hexahalogenohafnates. <i>Journal of Alloys and Compounds</i> , 1994, 210, 63-70.	5.5	5
106	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.	0.9	5
107	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.	2.8	5
108	Photoelectron spectroscopic and density functional theoretical studies of the 2-deoxycytidine homodimer radical anion. <i>Journal of Chemical Physics</i> , 2013, 139, 075101.	3.0	5

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109	5-(N-Trifluoromethylcarboxy)aminouracil as a Potential DNA Radiosensitizer and Its Radiochemical Conversion into N-Uracil-5-yloxamic Acid. <i>International Journal of Molecular Sciences</i> , 2020, 21, 6352.	4.1	5
110	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.	4.6	5
111	Low-Energy Electron Induced Reactions in Metronidazole at Different Solvation Conditions. <i>Pharmaceuticals</i> , 2022, 15, 701.	3.8	5
112	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.	5.5	4
113	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.	3.6	4
114	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.	1.5	4
115	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.	2.0	4
116	Photoelectron spectroscopy and density functional theory studies on the uridine homodimer radical anions. <i>Journal of Chemical Physics</i> , 2012, 137, 205101.	3.0	4
117	Artificial Plasmid Labeled with 5-Bromo-2-deoxyuridine: A Universal Molecular System for Strand Break Detection. <i>ChemBioChem</i> , 2014, 15, 1409-1412.	2.6	4
118	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.	3.8	4
119	Excess Electron Attachment to the Nucleoside Pair 2-Deoxyadenosine (dA)—2-Deoxythymidine (dT). <i>Journal of Physical Chemistry B</i> , 2016, 120, 4955-4962.	2.6	4
120	Quantitative assay of photoinduced DNA strand breaks by real-time PCR. <i>Journal of Pharmaceutical and Biomedical Analysis</i> , 2016, 128, 480-484.	2.8	4
121	Chemically—enzymatic synthesis of photosensitive DNA. <i>Journal of Photochemistry and Photobiology B: Biology</i> , 2017, 167, 228-235.	3.8	4
122	Guanosine Dianions Hydrated by One to Four Water Molecules. <i>Journal of Physical Chemistry Letters</i> , 2022, , 3230-3236.	4.6	4
123	Inactive-to-Active Transition of Human Thymidine Kinase 1 Revealed by Molecular Dynamics Simulations. <i>Journal of Chemical Information and Modeling</i> , 2022, 62, 142-149.	5.4	4
124	Radiosensitization of PC3 Prostate Cancer Cells by 5-Thiocyanato-2-deoxyuridine. <i>Cancers</i> , 2022, 14, 2035.	3.7	4
125	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.	2.7	3
126	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.	0.9	3



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127	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.	4.4	3
128	Photoelectron Spectroscopy and Computational Modeling of Thymidine Homodimer Anions. <i>Journal of Physical Chemistry B</i> , 2012, 116, 13975-13981.	2.6	3
129	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.	2.8	3
130	DNA Damage Radiosensitizers Geared Towards Hydrated Electrons. , 2022, , 125-169.		3
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