

M P Evstigneev

List of Publications by Year in descending order

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148
papers

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docs citations

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times ranked

1733
citing authors

#	ARTICLE	IF	CITATIONS
1	A Novel Water-Soluble C60 Fullerene-Based Nano-Platform Enhances Efficiency of Anticancer Chemotherapy. , 2022, , 59-93.		0
2	Interceptor potential of C60 fullerene aqueous solution: a comparative analysis using the example of the antitumor antibiotic mitoxantrone. European Biophysics Journal, 2022, 51, 297-307.	1.2	1
3	Design, characterization and mechanical properties of new Na ⁺ , CO ₃ ²⁻ -apatite/alginate/C60 fullerene hybrid biocomposites. Journal of the Korean Ceramic Society, 2021, 58, 422-429.	1.1	5
4	Comparative analysis of self-aggregation of liquid crystalline Pt(II) complexes in solution and in neat films. Journal of Organometallic Chemistry, 2021, 938, 121750.	0.8	1
5	Study of aggregation of O ^N N ^O Pt(II) complexes in solution. Journal of Molecular Liquids, 2021, 334, 116062.	2.3	2
6	Antitumor efficiency of the natural alkaloid berberine complexed with C60 fullerene in Lewis lung carcinoma in vitro and in vivo. Cancer Nanotechnology, 2021, 12, .	1.9	10
7	Nanocomplex of Berberine with C60 Fullerene Is a Potent Suppressor of Lewis Lung Carcinoma Cells Invasion In Vitro and Metastatic Activity In Vivo. Materials, 2021, 14, 6114.	1.3	5
8	The interplay of enthalpic/entropic factors in nanoparticles' aggregation in solution: The case of fullerene C60. Journal of Molecular Liquids, 2020, 318, 114043.	2.3	4
9	Studies on potential interaction between cinacalcet hydrochloride and diclofenac sodium. Biophysical Chemistry, 2020, 266, 106460.	1.5	0
10	Fractal C60 fullerene aggregation: Equilibrium thermodynamics approach. Chemical Physics Letters, 2020, 742, 137161.	1.2	7
11	Single-walled carbon nanotubes loaded hydroxyapatite- α -alginate beads with enhanced mechanical properties and sustained drug release ability. Progress in Biomaterials, 2020, 9, 1-14.	1.8	14
12	A Novel Nanoconjugate of Landomycin A with C60 Fullerene for Cancer Targeted Therapy: In Vitro Studies. Cellular and Molecular Bioengineering, 2019, 12, 41-51.	1.0	19
13	The energetics of small molecules binding with nucleic acids. Journal of Chemical Thermodynamics, 2019, 139, 105887.	1.0	1
14	In vitro study of the anticancer activity of various doxorubicin-containing dispersions. BiolImpacts, 2019, 9, 57-63.	0.7	17
15	C60 Fullerene as an Effective Nanoplatfrom of Alkaloid Berberine Delivery into Leukemic Cells. Pharmaceutics, 2019, 11, 586.	2.0	29
16	Tuning the Aggregation of N ¹ N ¹ C Pt(II) Complexes by Varying the Aliphatic Side Chain and Auxiliary Halide Ligand: ¹ H and ¹⁹⁵ Pt NMR Investigation. European Journal of Inorganic Chemistry, 2019, 2019, 4122-4128.	1.0	11
17	Interaction of pseudoephedrine and azithromycin with losartan: Spectroscopic, dissolution and permeation studies. Spectrochimica Acta - Part A: Molecular and Biomolecular Spectroscopy, 2019, 221, 117194.	2.0	2
18	The theory of interceptor-protector action of DNA binding drugs. Progress in Biophysics and Molecular Biology, 2019, 149, 131-146.	1.4	4

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19	Hetero-association models of non-covalent molecular complexation. <i>Physical Chemistry Chemical Physics</i> , 2019, 21, 7717-7731.	1.3	10
20	General up-scaled model of ligand binding with C60 fullerene clusters in aqueous solution. <i>Chemical Physics Letters</i> , 2019, 721, 22-26.	1.2	10
21	Does C60 fullerene act as a transporter of small aromatic molecules?. <i>Colloids and Surfaces B: Biointerfaces</i> , 2018, 164, 134-143.	2.5	34
22	Self-Organization of Pristine C60 Fullerene and its Complexes with Chemotherapy Drugs in Aqueous Solution as Promising Anticancer Agents. <i>Springer Proceedings in Physics</i> , 2018, , 3-22.	0.1	5
23	Study of the properties of doxorubicin-resistant cells affected by acute leucosis. <i>Journal of Bioenergetics and Biomembranes</i> , 2018, 50, 53-58.	1.0	1
24	Generalized shape-independent approach to studying molecular hetero-assembly in solution using NMR diffusometry. <i>Journal of Molecular Liquids</i> , 2018, 265, 88-95.	2.3	4
25	Determination of the equilibrium constant of C ₆₀ fullerene binding with drug molecules. <i>Physical Chemistry Chemical Physics</i> , 2017, 19, 6777-6784.	1.3	11
26	Measuring Self-Association of Pt Complexes by ¹⁹⁵ Pt NMR. <i>ChemistrySelect</i> , 2017, 2, 3353-3355.	0.7	5
27	Comparative Analysis of the Antineoplastic Activity of C60 Fullerene with 5-Fluorouracil and Pyrrole Derivative In Vivo. <i>Nanoscale Research Letters</i> , 2017, 12, 8.	3.1	28
28	C60 fullerene enhances cisplatin anticancer activity and overcomes tumor cell drug resistance. <i>Nano Research</i> , 2017, 10, 652-671.	5.8	61
29	Molecular Modeling-Based Energy Analysis of Dimeric Binding of Ligands to the Minor DNA Groove. <i>Biophysics (Russian Federation)</i> , 2017, 62, 876-884.	0.2	0
30	A nanocomplex of C ₆₀ fullerene with cisplatin: design, characterization and toxicity. <i>Beilstein Journal of Nanotechnology</i> , 2017, 8, 1494-1501.	1.5	41
31	Effect of Nicotinamide on the Photolysis of Riboflavin in Aqueous Solution. <i>Scientia Pharmaceutica</i> , 2016, 84, 289-303.	0.7	2
32	Biophysical characterization of the complexation of C ₆₀ fullerene with doxorubicin in a prokaryotic model. <i>Materialwissenschaft Und Werkstofftechnik</i> , 2016, 47, 92-97.	0.5	12
33	Clustering of hydrochloric acid on the surface of C ₆₀ /C ₇₀ fullerite and its composites with nanosilica. <i>Materialwissenschaft Und Werkstofftechnik</i> , 2016, 47, 172-179.	0.5	1
34	Interaction of C ₆₀ fullerene complexed to cisplatin with model bilipid membranes and its uptake by HeLa cells. <i>Materialwissenschaft Und Werkstofftechnik</i> , 2016, 47, 105-111.	0.5	2
35	Conformational, IR spectroscopic and electronic properties of conium alkaloids and their adducts with C60 fullerene. <i>Journal of Molecular Structure</i> , 2016, 1118, 167-171.	1.8	2
36	Physiographic factors of seasonal distribution of linear trends in air temperature on the Azov-Black sea coast. <i>Russian Meteorology and Hydrology</i> , 2016, 41, 19-27.	0.2	3

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37	Development of heterogeneous preparation with inulinase for tubular reactor systems. <i>Journal of Molecular Catalysis B: Enzymatic</i> , 2016, 129, 1-5.	1.8	6
38	Study of the complexation between Landomycin A and C60 fullerene in aqueous solution. <i>RSC Advances</i> , 2016, 6, 81231-81236.	1.7	12
39	Does fullerene C60 affect cell membranes integrity?. <i>New Biotechnology</i> , 2016, 33, S146-S147.	2.4	0
40	Optimal experiment design: Link between the concentration and the accuracy of estimation of aggregation parameters. <i>Chemical Physics Letters</i> , 2016, 664, 133-137.	1.2	0
41	Interrelation of Entropic Contributors to π -Stacking in Solution. <i>International Journal of Thermophysics</i> , 2016, 37, 1.	1.0	1
42	Hidden entropic contribution in the thermodynamics of molecular complexation. <i>Physical Chemistry Chemical Physics</i> , 2016, 18, 7617-7625.	1.3	5
43	Interaction of C 60 fullerene complexed to doxorubicin with model bilipid membranes and its uptake by HeLa cells. <i>Materials Science and Engineering C</i> , 2016, 59, 398-403.	3.8	33
44	Complex of C60 Fullerene with Doxorubicin as a Promising Agent in Antitumor Therapy. <i>Nanoscale Research Letters</i> , 2015, 10, 499.	3.1	57
45	Application of C₆₀ Fullerene-Doxorubicin Complex for Tumor Cell Treatment <i>in vitro</i> and <i>in vivo</i>. <i>Journal of Biomedical Nanotechnology</i> , 2015, 11, 1139-1152.	0.5	83
46	Entropic binding mode preference in cooperative homo-dimeric drug-DNA recognition. <i>Chemical Physics Letters</i> , 2015, 624, 12-14.	1.2	4
47	Structural organization of C60 fullerene, doxorubicin, and their complex in physiological solution as promising antitumor agents. <i>Journal of Nanoparticle Research</i> , 2015, 17, 1.	0.8	49
48	PMR Characterization of the Water Structure in Tibetan Milk Mushroom <i>Zooglea</i> : Influence of Medium Hydration and Hydrophobicity. <i>Journal of Applied Spectroscopy</i> , 2015, 82, 353-359.	0.3	1
49	C60 fullerene affects elastic properties and osmoregulation reactions of human lymphocytes. <i>European Biophysics Journal</i> , 2015, 44, 493-498.	1.2	7
50	Shape-independent model (SHIM) approach for studying aggregation by NMR diffusometry. <i>Journal of Chemical Physics</i> , 2015, 142, 104202.	1.2	9
51	Structural self-organization of C₆₀ and cisplatin in physiological solution. <i>Physical Chemistry Chemical Physics</i> , 2015, 17, 26084-26092.	1.3	40
52	Structural Features of Highly Stable Reproducible C₆₀ Fullerene Aqueous Colloid Solution Probed by Various Techniques. <i>Fullerenes Nanotubes and Carbon Nanostructures</i> , 2015, 23, 530-534.	1.0	103
53	Tumor-Inhibitory Effect of C60 Fullerene Complex with Doxorubicin. <i>Nanomedicine and Nanobiology</i> , 2015, 2, 49-53.	0.4	5
54	Thermal analysis of ligand-DNA interaction: determination of binding parameters. <i>AIMS Biophysics</i> , 2015, 2, 423-440.	0.3	8

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55	C60 Fullerene as Synergistic Agent in Tumor-Inhibitory Doxorubicin Treatment. <i>Drugs in R and D</i> , 2014, 14, 333-340.	1.1	69
56	General features of the energetics of complex formation between ligand and nucleic acids. <i>Biophysics (Russian Federation)</i> , 2014, 59, 546-551.	0.2	2
57	Energy analysis of non-covalent ligand binding to nucleic acids: Present and future. <i>Biophysics (Russian Federation)</i> , 2014, 59, 552-555.	0.2	2
58	On the Origin of C ₆₀ Fullerene Solubility in Aqueous Solution. <i>Langmuir</i> , 2014, 30, 3967-3970.	1.6	109
59	Effect of phosphate buffer on the complexation and photochemical interaction of riboflavin and caffeine in aqueous solution: A kinetic study. <i>Journal of Photochemistry and Photobiology A: Chemistry</i> , 2014, 273, 17-22.	2.0	12
60	Quantitative correlation of the in vitro biological effect with parameters of molecular complexation in mutagen-interceptor systems. <i>Journal of Theoretical Biology</i> , 2014, 357, 268-271.	0.8	2
61	Binding polynomial in molecular self-assembly. <i>Physical Review E</i> , 2014, 89, 062138.	0.8	4
62	Evidence of entropically driven C60 fullerene aggregation in aqueous solution. <i>Journal of Chemical Physics</i> , 2014, 140, 104909.	1.2	25
63	Theoretical Description of Metabolism Using Queueing Theory. <i>Bulletin of Mathematical Biology</i> , 2014, 76, 2238-2248.	0.9	7
64	Characterization of C ₆₀ fullerene complexation with antibiotic doxorubicin. <i>Physical Chemistry Chemical Physics</i> , 2014, 16, 23164-23172.	1.3	55
65	Hetero-association of aromatic molecules in aqueous solution. <i>International Reviews in Physical Chemistry</i> , 2014, 33, 229-273.	0.9	29
66	Interceptor effect of C60 fullerene on the in vitro action of aromatic drug molecules. <i>European Biophysics Journal</i> , 2014, 43, 265-276.	1.2	44
67	Complexation of aromatic drugs with single-walled carbon nanotubes. <i>Journal of Nanoparticle Research</i> , 2014, 16, 1.	0.8	14
68	Spectroscopic Study of Proflavine Adsorption on the Carbon Nanotube Surface. <i>Applied Spectroscopy</i> , 2014, 68, 232-237.	1.2	3
69	Dimerization Energetics of DNA Minor Groove Binders. <i>Ukrainian Journal of Physics</i> , 2014, 59, 461-472.	0.1	3
70	Indistinguishability of the models of molecular self-assembly. <i>Supramolecular Chemistry</i> , 2013, 25, 199-203.	1.5	15
71	The role of mixing entropy in molecular self-assembly. <i>Chemical Physics Letters</i> , 2013, 567, 48-49.	1.2	6
72	Complexation of C ₆₀ Fullerene with Aromatic Drugs. <i>ChemPhysChem</i> , 2013, 14, 568-578.	1.0	68

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73	On the reliability of quantitation of biological effect in drugâ€“interceptorâ€“DNA systems. European Biophysics Journal, 2013, 42, 315-319.	1.2	4
74	C60 fullerene aggregation in aqueous solution. Physical Chemistry Chemical Physics, 2013, 15, 9351.	1.3	75
75	General statistical-thermodynamical treatment of one-dimensional multicomponent molecular hetero-assembly in solution. Chemical Physics, 2013, 421, 77-83.	0.9	11
76	Thiazotropsin aggregation and its relationship to molecular recognition in the DNA minor groove. Biophysical Chemistry, 2013, 179, 1-11.	1.5	14
77	Development of an Analytical Approach to Study a Three-Component Hetero-Association by Means of Spectrophotometry. Applied Spectroscopy, 2013, 67, 29-35.	1.2	5
78	Physicochemical Mechanisms of Synergistic Biological Action of Combinations of Aromatic Heterocyclic Compounds. Organic Chemistry International, 2013, 2013, 1-10.	1.0	16
79	On the a priori possibility of the formation of hexameric mini-hairpin d(GCGAGC) in solution. Journal of Molecular Structure, 2012, 1027, 124-127.	1.8	2
80	Energy of ligand-RNA complex formation. Biophysics (Russian Federation), 2012, 57, 450-463.	0.2	5
81	Relation between the change in DNA elasticity on ligand binding and the binding energetics. Physical Review E, 2012, 86, 031919.	0.8	6
82	Energetics of ligand binding to the DNA minor groove. Physical Chemistry Chemical Physics, 2012, 14, 5588.	1.3	21
83	Random versus sequential pathway of molecular self-assembly. Physical Review E, 2012, 85, 061405.	0.8	13
84	Intermolecular hydrogen bonds in hetero-complexes of biologically active aromatic molecules probed by the methods of vibrational spectroscopy. Spectrochimica Acta - Part A: Molecular and Biomolecular Spectroscopy, 2012, 95, 224-229.	2.0	7
85	¹ H NMR study of the complexation of aromatic drugs with dimethylxanthine derivatives. Journal of Molecular Structure, 2012, 1010, 139-145.	1.8	5
86	General analysis of competitive binding in drugâ€“interceptorâ€“DNA systems. European Biophysics Journal, 2012, 41, 273-283.	1.2	16
87	Thermodynamic analysis of complex formation of ethidium bromide with DNA in water solutions. Biophysics (Russian Federation), 2011, 56, 214-219.	0.2	9
88	Contribution of enthalpy to the energetics of complex formation of aromatic ligands with DNA. Biophysics (Russian Federation), 2011, 56, 634-643.	0.2	4
89	A revised treatment of the non-electrostatic contribution to the solvation free energy of DNA-binding ligands. Journal of Molecular Liquids, 2011, 163, 178-180.	2.3	4
90	Quantification of the interceptor action of caffeine on the in vitro biological effect of the anti-tumour agent topotecan. European Biophysics Journal, 2011, 40, 969-980.	1.2	15

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91	Calculation of the electrostatic charges and energies for intercalation of aromatic drug molecules with DNA. <i>International Journal of Quantum Chemistry</i> , 2011, 111, 711-721.	1.0	17
92	A novel computational approach to study ligand binding to finite lattice. <i>Biopolymers</i> , 2011, 95, 208-216.	1.2	3
93	NMR analysis of Nile Blue (C. I. Basic Blue 12) and Thionine (C. I. 52000) in solution. <i>Dyes and Pigments</i> , 2011, 88, 315-325.	2.0	15
94	Additional stabilization of hetero-complexes of aromatic molecules: H-bonds or charge-transfer?. <i>Journal of Molecular Structure</i> , 2011, 985, 403-406.	1.8	4
95	Parsing of the free energy of aromatic aromatic stacking interactions in solution. <i>Journal of Chemical Thermodynamics</i> , 2011, 43, 1424-1434.	1.0	32
96	Complete solution of the problem of one-dimensional non-covalent non-cooperative self-assembly in two-component systems. <i>Journal of Chemical Physics</i> , 2011, 134, 194902.	1.2	19
97	Does the ligand-biopolymer equilibrium binding constant depend on the number of bound ligands?. <i>Biopolymers</i> , 2010, 93, 932-935.	1.2	5
98	Hexamer oligonucleotide topology and assembly under solution phase NMR and theoretical modeling scrutiny. <i>Biopolymers</i> , 2010, 93, 1023-1038.	1.2	5
99	Relation between structure and enthalpy for stacking interactions of aromatic molecules. <i>Molecular Physics</i> , 2010, 108, 1941-1947.	0.8	8
100	Computation of energetic contribution of loss of degrees of freedom upon complexation of aromatic molecules. <i>Journal of Physical Studies</i> , 2010, 14, .	0.2	1
101	Hydration change on complexation of aromatic ligands with DNA: molecular dynamics simulations. <i>Biopolymers and Cell</i> , 2010, 26, 36-44.	0.1	8
102	Partition of thermodynamic energies of drug-DNA complexation. <i>Biopolymers</i> , 2009, 91, 773-790.	1.2	32
103	Interaction of ethidium bromide and caffeine with DNA in aqueous solution. <i>Journal of Applied Spectroscopy</i> , 2009, 76, 132-139.	0.3	40
104	Complexation of Biologically Active Aromatic Compounds with DNA in the Presence of Theophylline. <i>Journal of Biological Physics</i> , 2009, 35, 115-126.	0.7	13
105	Contribution of changes in translational, rotational, and vibrational degrees of freedom to the energy of complex formation of aromatic ligands with DNA. <i>Biophysics (Russian Federation)</i> , 2009, 54, 428-436.	0.2	5
106	Calculation of the thermodynamic potentials of changes in translational, rotational, and vibrational degrees of freedom in the dimerization of aromatic molecules. <i>Russian Journal of Physical Chemistry B</i> , 2009, 3, 707-712.	0.2	2
107	Profiles of equilibrium constants for self-association of aromatic molecules. <i>Journal of Chemical Physics</i> , 2009, 130, 165105.	1.2	27
108	Hydrophobic contribution to the free energy of complexation of aromatic ligands with DNA. <i>Biopolymers and Cell</i> , 2009, 25, 133-141.	0.1	6

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109	Quantitation of the molecular mechanisms of biological synergism in a mixture of DNA-acting aromatic drugs. <i>Biophysical Chemistry</i> , 2008, 132, 148-158.	1.5	21
110	Complexation of heterocyclic ligands with DNA in aqueous solution. <i>Journal of Applied Spectroscopy</i> , 2008, 75, 251-260.	0.3	24
111	Electrostatic contribution to the energy of binding of aromatic ligands with DNA. <i>Biopolymers</i> , 2008, 89, 680-690.	1.2	23
112	Structural and thermodynamic analysis of the hetero-association of theophylline with aromatic drug molecules. <i>Journal of Molecular Structure</i> , 2008, 889, 229-236.	1.8	12
113	A "microscopic" model of complexation between intercalators and DNA: Analysis of NMR spectroscopy data. <i>Biophysics (Russian Federation)</i> , 2008, 53, 38-43.	0.2	1
114	Structural Basis for the Binding Affinity of a Homologous Series of Synthetic Phenoxazone Drugs with DNA: NMR and Molecular Mechanics Analysis. <i>Journal of Biomolecular Structure and Dynamics</i> , 2007, 24, 443-453.	2.0	5
115	A method for analysis of multicomponent systems of interacting aromatic molecules in solution. <i>Journal of Chemical Physics</i> , 2007, 127, 154511.	1.2	14
116	On the origin of the decrease in stability of the DNA hairpin d(GCGAAGC) on complexation with aromatic drugs. <i>Biophysical Chemistry</i> , 2007, 129, 56-59.	1.5	7
117	Investigation of the complexation of the anti-cancer drug novantrone with the hairpin structure of the deoxyheptanucleotide 5'-d(GpCpGpApApGpC). <i>Journal of Molecular Structure</i> , 2007, 843, 78-86.	1.8	8
118	Structural and thermodynamic analysis of the conformational states of self-complementary hexanucleotides 5'-d(GCATGC) and 5'-d(GCTAGC) in Aqueous Solution. <i>Biophysics (Russian Federation)</i> , 2007, 52, 375-382.	0.2	1
119	Formation of complexes of antimicrobial agent norfloxacin with antitumor antibiotics of anthracycline series. <i>Russian Journal of Physical Chemistry A</i> , 2007, 81, 802-807.	0.1	6
120	Estimation of the mean number of molecules in associates of aromatic compounds. <i>Russian Journal of Physical Chemistry A</i> , 2007, 81, 1888-1890.	0.1	3
121	¹ H NMR study of the hetero-association of flavin-mononucleotide with mutagenic dyes: ethidium bromide and proflavine. <i>Molecular Physics</i> , 2006, 104, 647-654.	0.8	14
122	Interaction between aromatic antibiotics and vitamins: ¹ H NMR study of heteroassociation of nicotinamide and anthracycline antitumor antibiotics. <i>Biophysics (Russian Federation)</i> , 2006, 51, 176-185.	0.2	3
123	Heteroassociation of antibiotic norfloxacin with aromatic vitamins in aqueous solution. <i>Biophysics (Russian Federation)</i> , 2006, 51, 592-598.	0.2	4
124	Complexation of the quinolone antibiotic norfloxacin with DNA. <i>Molecular Biology</i> , 2006, 40, 805-810.	0.4	6
125	Complexation of norfloxacin with DNA in the presence of caffeine. <i>Biophysical Chemistry</i> , 2006, 121, 84-95.	1.5	39
126	NMR investigation of the effect of caffeine on the hetero-association of an anticancer drug with a vitamin. <i>Chemical Physics Letters</i> , 2006, 432, 248-251.	1.2	22

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127	Stochastic models (cooperative and non-cooperative) for NMR analysis of the hetero-association of aromatic molecules in aqueous solution. <i>Chemical Physics</i> , 2006, 321, 25-33.	0.9	20
128	¹ H NMR determination of the self-association of an acridine homodimer and its complexation with ethidium bromide in aqueous solution. <i>Journal of Molecular Structure</i> , 2006, 784, 162-168.	1.8	4
129	Spectrophotometric investigation of the hetero-association of Caffeine and thiazine dye in aqueous solution. <i>Spectrochimica Acta - Part A: Molecular and Biomolecular Spectroscopy</i> , 2006, 64, 693-697.	2.0	32
130	Self-association of daunomycin antibiotic in various buffer solutions. <i>Russian Journal of Physical Chemistry A</i> , 2006, 80, 741-746.	0.1	17
131	Aggregation of 1,3,7-trimethylxanthine with methylene blue in aqueous solution. <i>Journal of Applied Spectroscopy</i> , 2006, 73, 171-177.	0.3	12
132	Complexation of anthracycline drugs with DNA in the presence of caffeine. <i>European Biophysics Journal</i> , 2006, 36, 1-11.	1.2	31
133	Effect of a mixture of caffeine and nicotinamide on the solubility of vitamin (B2) in aqueous solution. <i>European Journal of Pharmaceutical Sciences</i> , 2006, 28, 59-66.	1.9	51
134	Hetero-association of anticancer antibiotics in aqueous solution: NMR and molecular mechanics analysis. <i>Biophysical Chemistry</i> , 2005, 117, 111-118.	1.5	17
135	Complexation of daunomycin with a DNA oligomer in the presence of an aromatic vitamin (B2) determined by NMR spectroscopy. <i>Biophysical Chemistry</i> , 2005, 118, 118-127.	1.5	19
136	Structural and thermodynamic analysis of heteroassociation of daunomycin and flavin mononucleotide molecules in water by ¹ H NMR spectroscopy. <i>Journal of Structural Chemistry</i> , 2005, 46, 67-74.	0.3	6
137	Analysis by Means of ¹ H NMR Spectroscopy of Heteroassociation in Water Solution of Antitumor Antibiotics Daunomycin and Actinomycin D. <i>Russian Journal of Organic Chemistry</i> , 2005, 41, 1158-1164.	0.3	3
138	<title>Comparative analysis of the self-association of ethidium mono- and bis-intercalators in aqueous solution using ^1H NMR spectroscopy</title>. , 2004, , .		0
139	¹ H NMR Study of Heteroassociation of Ethidium Homodimer and Propidium Iodide in Water. <i>Journal of Structural Chemistry</i> , 2004, 45, 793-799.	0.3	1
140	¹ H NMR structural and thermodynamical analysis of the hetero-association of daunomycin and novatrone in aqueous solution. <i>Journal of Molecular Structure</i> , 2004, 701, 31-37.	1.8	17
141	¹ H NMR investigation of the self-association of ethidium homodimer and its complexation with propidium iodide in aqueous solution. <i>Journal of Molecular Structure</i> , 2004, 690, 17-24.	1.8	1
142	¹ H NMR Study of Heteroassociation of Daunomycin and Propidium Iodide in Aqueous Solution. <i>Russian Journal of Organic Chemistry</i> , 2002, 38, 1035-1041.	0.3	3
143	¹ H-NMR analysis of the interaction of antibiotic mitoxantrone with DNA in the presence of caffeine in aqueous solution. <i>Biopolymers and Cell</i> , 2002, 18, 287-296.	0.1	1
144	Self-association of the antitumour agent novatrone (mitoxantrone) and its hetero-association with caffeine. <i>Perkin Transactions II RSC</i> , 2001, , 61-67.	1.1	47

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145	¹ H NMR Analysis of Heteroassociation of Caffeine with Mitoxanthrone in Aqueous Solution. Journal of Structural Chemistry, 2001, 42, 777-783.	0.3	4
146	A general nuclear magnetic resonance analysis of hetero-association of aromatic molecules in aqueous solution. Journal of Chemical Physics, 2001, 115, 2252-2266.	1.2	25
147	A generalized statistical-thermodynamical model of hetero-association of aromatic molecules in aqueous solution for the NMR data interpretation. Biopolymers and Cell, 2001, 17, 501-511.	0.1	1
148	¹ H NMR investigation of the hetero-association of aromatic molecules in aqueous solution: factors involved in the stabilization of complexes of daunomycin and acridine drugs. Molecular Physics, 2000, 98, 1961-1971.	0.8	45