Vladimir I Polshakov

List of Publications by Year in descending order

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95 papers 1,209 citations

³⁶¹⁴¹³
20
h-index

454955 30 g-index

112 all docs

 $\begin{array}{c} 112 \\ \\ \text{docs citations} \end{array}$

112 times ranked

1194 citing authors

#	Article	IF	CITATIONS
1	Biological evaluation and spectral characterization of a novel tetracenomycin X congener. Biochimie, 2022, 192, 63-71.	2.6	5
2	Telomere length regulation by Rif1 protein from Hansenula polymorpha. ELife, 2022, 11, .	6.0	1
3	Analogs of S-Adenosyl-L-Methionine in Studies of Methyltransferases. Molecular Biology, 2022, 56, 229-250.	1.3	15
4	Antifungal Thiazolidines: Synthesis and Biological Evaluation of Mycosidine Congeners. Pharmaceuticals, 2022, 15, 563.	3.8	10
5	Structural basis for interaction between CLAMP and MSL2 proteins involved in the specific recruitment of the dosage compensation complex in <i>Drosophila</i> . Nucleic Acids Research, 2022, 50, 6521-6531.	14.5	4
6	NMR resonance assignment and backbone dynamics of a C-terminal domain homolog of orange carotenoid protein. Biomolecular NMR Assignments, 2021, 15, 17-23.	0.8	4
7	Synthesis and Biological Evaluation of PSMA Ligands with Aromatic Residues and Fluorescent Conjugates Based on Them. Journal of Medicinal Chemistry, 2021, 64, 4532-4552.	6.4	19
8	Synthesis, Characterization, and Preclinical Evaluation of a Small-Molecule Prostate-Specific Membrane Antigen-Targeted Monomethyl Auristatin E Conjugate. Journal of Medicinal Chemistry, 2021, 64, 17123-17145.	6.4	12
9	Responses of DNA Mismatch Repair Proteins to a Stable G-Quadruplex Embedded into a DNA Duplex Structure. International Journal of Molecular Sciences, 2020, 21, 8773.	4.1	12
10	Backbone resonance assignment and dynamics of 110 kDa hexameric inorganic pyrophosphatase from Mycobacterium tuberculosis. Biomolecular NMR Assignments, 2020, 14, 281-287.	0.8	1
11	Polypeptide-Based Molecular Platform and Its Docetaxel/Sulfo-Cy5-Containing Conjugate for Targeted Delivery to Prostate Specific Membrane Antigen. Molecules, 2020, 25, 5784.	3.8	13
12	Interplay of Pyrrolidine Units with Homo/Hetero Chirality and CF3–Aryl Substituents on Secondary Structures of β-Proline Tripeptides in Solution. Journal of Organic Chemistry, 2020, 85, 8865-8871.	3.2	1
13	Tetracenomycin X inhibits translation by binding within the ribosomal exit tunnel. Nature Chemical Biology, 2020, 16, 1071-1077.	8.0	43
14	Insights into the structure and function of Est3 from the Hansenula polymorpha telomerase. Scientific Reports, 2020, 10, 11109.	3.3	4
15	Williams–Beuren syndromeâ€related methyltransferase WBSCR27: cofactor binding and cleavage. FEBS Journal, 2020, 287, 5375-5393.	4.7	6
16	Thiazolidine-2,4-dione in benzoylation reaction. Chemistry of Heterocyclic Compounds, 2019, 55, 178-183.	1.2	3
17	Nybomycin-producing Streptomyces isolated from carpenter ant Camponotus vagus. Biochimie, 2019, 160, 93-99.	2.6	25
18	NMR screening and studies of target – ligand interactions. Russian Chemical Reviews, 2019, 88, 59-98.	6.5	7

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19	Structure and function of the N-terminal domain of the yeast telomerase reverse transcriptase. Nucleic Acids Research, 2018, 46, 1525-1540.	14.5	19
20	Chemical shift assignments and the secondary structure of the Est3 telomerase subunit in the yeast Hansenula polymorpha. Biomolecular NMR Assignments, 2018, 12, 57-62.	0.8	3
21	Theoretical and NMR Conformational Studies of \hat{l}^2 -Proline Oligopeptides With Alternating Chirality of Pyrrolidine Units. Frontiers in Chemistry, 2018, 6, 91.	3.6	8
22	Enalaprilat Inhibits Zinc-Dependent Oligomerization of Metal-Binding Domain of Amyloid-beta Isoforms and Protects Human Neuroblastoma Cells from Toxic Action of these Isoforms. Molecular Biology, 2018, 52, 590-597.	1.3	9
23	NMR assignments of the WBSCR27 protein related to Williams-Beuren syndrome. Biomolecular NMR Assignments, 2018, 12, 303-308.	0.8	3
24	A Binuclear Zinc Interaction Fold Discovered in the Homodimer of Alzheimer's Amyloidâ€Î² Fragment with Taiwanese Mutation D7H. Angewandte Chemie - International Edition, 2017, 56, 11734-11739.	13.8	25
25	Frontispiece: A Binuclear Zinc Interaction Fold Discovered in the Homodimer of Alzheimer's Amyloidâ€Î² Fragment with Taiwanese Mutation D7H. Angewandte Chemie - International Edition, 2017, 56, .	13.8	0
26	Frontispiz: A Binuclear Zinc Interaction Fold Discovered in the Homodimer of Alzheimer's Amyloidâ $\in \hat{I}^2$ Fragment with Taiwanese Mutation D7H. Angewandte Chemie, 2017, 129, .	2.0	0
27	A Binuclear Zinc Interaction Fold Discovered in the Homodimer of Alzheimer's Amyloidâ€Î² Fragment with Taiwanese Mutation D7H. Angewandte Chemie, 2017, 129, 11896-11901.	2.0	5
28	Reinvestigation of dimerization of Z-N-alkylarylmethylideneindoxyls upon exposure to UV-vis radiation. Russian Chemical Bulletin, 2017, 66, 350-354.	1.5	2
29	Versatility of the green microalga cell vacuole function as revealed by analytical transmission electron microscopy. Protoplasma, 2017, 254, 1323-1340.	2.1	49
30	Intermolecular interactions in rifabutin–2-hydroxypropyl-β-cyclodextrin–water solutions, according to solubility data. Russian Journal of Physical Chemistry A, 2016, 90, 983-989.	0.6	1
31	Control of Azomethine Cycloaddition Stereochemistry by CF ₃ Group: Structural Diversity of Fluorinated β-Proline Dimers. Organic Letters, 2016, 18, 4698-4701.	4.6	8
32	"Suppressor factor―of neutrophils: A short story of a long-term misconception. Biochemistry (Moscow), 2016, 81, 1284-1292.	1.5	0
33	Interplay of histidine residues of the Alzheimer's disease Aβ peptide governs its Zn-induced oligomerization. Scientific Reports, 2016, 6, 21734.	3.3	81
34	New conjugates of polyene macrolide amphotericin B with benzoxaboroles: synthesis and properties. Journal of Antibiotics, 2016, 69, 549-560.	2.0	24
35	NMR assignments of the N-terminal domain of Ogataea polymorpha telomerase reverse transcriptase. Biomolecular NMR Assignments, 2016, 10, 183-187.	0.8	5
36	The QTL within the H2 Complex Involved in the Control of Tuberculosis Infection in Mice Is the Classical Class II H2-Ab1 Gene. PLoS Genetics, 2015, 11, e1005672.	3.5	24

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37	The English (H6R) familial Alzheimer's disease mutation facilitates zinc-induced dimerization of the amyloid- \hat{l}^2 metal-binding domain. Metallomics, 2015, 7, 422-425.	2.4	38
38	Intermolecular interactions in rifabutin—2-hydroxypropyl-β-cyclodextrin—water solutions. Russian Journal of Physical Chemistry A, 2015, 89, 797-801.	0.6	4
39	pH-responsive modulation of insulin aggregation and structural transformation of the aggregates. Biochimie, 2015, 109, 49-59.	2.6	18
40	Backbone 1H, 13C and 15N resonance assignments of the human eukaryotic release factor eRF1. Biomolecular NMR Assignments, 2015, 9, 37-42.	0.8	0
41	Structural Studies and Anticancer Activity of a Novel Class of βâ€Peptides. Chemistry - an Asian Journal, 2015, 10, 383-389.	3.3	17
42	NMR screening of potential inhibitors of methionine \hat{I}^3 -lyase from Citrobacter freundii. Molecular Biology, 2014, 48, 896-905.	1.3	4
43	Phosphorylation of Ser8 promotes zinc-induced dimerization of the amyloid- \hat{l}^2 metal-binding domain. Molecular BioSystems, 2014, 10, 2590-2596.	2.9	49
44	P4-028: ZINC-INDUCED DIMERS OF CHEMICALLY MODIFIED A \hat{I}^2 - ARE POSSIBLE AGGREGATION SEEDS. , 2014, 10 P793-P793.	0,	1
45	NMR Solution Structure of Rat Aβ(1–16): Toward Understanding the Mechanism of Rats' Resistance to Alzheimer's Disease. Biophysical Journal, 2012, 102, 136-143.	0.5	56
46	Structure and dynamics in solution of the stop codon decoding Nâ€terminal domain of the human polypeptide chain release factor eRF1. Protein Science, 2012, 21, 896-903.	7.6	3
47	NMR Structures of ApoL. caseiDihydrofolate Reductase and Its Complexes with Trimethoprim and NADPH: Contributions to Positive Cooperative Binding from Ligand-Induced Refolding, Conformational Changes, and Interligand Hydrophobic Interactions. Biochemistry, 2011, 50, 3609-3620.	2.5	24
48	Z/E(C=C)-isomerization and fluorescence modulation of imines of 7-N,N-dialkylamino-4-hydroxy-3-formylcoumarins in organic solvents. Heterocyclic Communications, 2011, ,	1.2	0
49	E/Z(C=C)-Isomerization of enamines of 3-formyl-4-hydroxycoumarin induced by organic solvents. Russian Chemical Bulletin, 2010, 59, 1605-1611.	1.5	11
50	NMR solution structure and function of the Câ€terminal domain of eukaryotic class 1 polypeptide chain release factor. FEBS Journal, 2010, 277, 2611-2627.	4.7	18
51	Optimization of the methods for small peptide solution structure determination by NMR spectroscopy. Molecular Biology, 2010, 44, 958-967.	1.3	4
52	Z/E (C=C)-isomerization and fluorescence modulation of imines of 7-N,N-dialkylamino-4-hydroxy-3-formylcoumarins in organic solvents. Heterocyclic Communications, 2010, 16, .	1.2	2
53	NMR solution structure and function of the C-terminal domain of eukaryotic class $\hat{a} \in f1$ polypeptide chain release factor. FEBS Journal, 2010, 277, 2611-2627.	4.7	12
54	Z/E(C=C)-isomerization of coumarin enamines induced by organic solvents. Mendeleev Communications, 2009, 19, 214-216.	1.6	20

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55	Interface of the interaction of the middle domain of human translation termination factor eRF1 with eukaryotic ribosomes. Molecular Biology, 2008, 42, 939-948.	1.3	3
56	Distribution coefficient of rifabutin in liposome/water system as measured by different methods. European Journal of Pharmaceutics and Biopharmaceutics, 2008, 68, 400-405.	4.3	16
57	Eukaryotic classâ \in f1 translation termination factor eRF1â \in fâ^â \in fthe NMR structure and dynamics of the middleâ \in fdomain involved in triggering ribosomeâ \in dependent peptidylâ \in tRNA hydrolysis. FEBS Journal, 2007, 274, 4223-4237.	4.7	21
58	Laser control of the structure of a photosensitive substrate for enzymatic reaction. Laser Physics, 2007, 17, 1262-1265.	1.2	1
59	Structural factors determining the binding selectivity of the antibacterial drug trimethoprim to dihydrofolate reductase. Pharmaceutical Chemistry Journal, 2007, 41, 350-353.	0.8	7
60	NMR assignments of the C-terminal domain of human polypeptide release factor eRF1. Biomolecular NMR Assignments, 2007, 1, 183-185.	0.8	6
61	Effects of Co-operative Ligand Binding on Protein Amide NH Hydrogen Exchange. Journal of Molecular Biology, 2006, 356, 886-903.	4.2	25
62	NMR Assignments of the Middle Domain of Human Polypeptide Release Factor eRF1. Journal of Biomolecular NMR, 2006, 36, 8-8.	2.8	3
63	Solution Structure of Human Dihydrofolate Reductase in its Complex with Trimethoprim and NADPH. Journal of Biomolecular NMR, 2005, 33, 69-72.	2.8	28
64	Towards understanding the origins of the different specificities of binding the reduced (NADPH) and oxidised (NADP+) forms of nicotinamide adenine dinucleotide phosphate coenzyme to dihydrofolate reductase. Journal of Molecular Structure, 2002, 602-603, 257-267.	3.6	9
65	NMR-based solution structure of the complex of Lactobacillus casei dihydrofolate reductase with trimethoprim and NADPH. Journal of Biomolecular NMR, 2002, 24, 67-70.	2.8	18
66	Dihydrofolate reductase: structural aspects of mechanisms of enzyme catalysis and inhibition. Russian Chemical Bulletin, 2001, 50, 1733-1751.	1.5	18
67	NMR Studies of Ligand Carboxylate Group Interactions with Arginine Residues in Complexes of Lactobacillus casei Dihydrofolate Reductase with Substrates and Substrate Analogues. Biochemistry, 2000, 39, 9819-9825.	2.5	14
68	Validation of a new restraint docking method for solution structure determinations of protein-ligand complexes. Journal of Biomolecular NMR, 1999, 14, 115-122.	2.8	10
69	Characterization of Rates of Ring-Flipping in Trimethoprim in Its Ternary Complexes withLactobacillus caseiDihydrofolate Reductase and Coenzyme Analoguesâ€. Biochemistry, 1999, 38, 15962-15969.	2.5	21
70	Structure and dynamics in solution of the complex of <i>lactobacillus casei</i> dihydrofolate reductase with the new lipophilic antifolate drug trimetrexate. Protein Science, 1999, 8, 467-481.	7.6	29
71	The solution structure of the complex of Lactobacillus casei dihydrofolate reductase with methotrexate. Journal of Molecular Biology, 1998, 277, 119-134.	4.2	38
72	High-resolution solution structure of human pNR-2/ps2: A single trefoil motif protein. Journal of Molecular Biology, 1997, 267, 418-432.	4.2	60

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73	NMR Detection of Arginine-Ligand Interactions in Complexes of Lactobacillus casei Dihydrofolate Reductase. FEBS Journal, 1996, 238, 435-439.	0.2	21
74	NMR-Based Structural Studies of the pNR-2/pS2 Single Domain Trefoil Peptide. Similarities to Porcine Spasmolytic Peptide and Evidence for a Monomeric Structure. FEBS Journal, 1995, 233, 847-855.	0.2	18
75	Determination of Stereospecific Assignments, Torsion-Angle Constraints, and Rotamer Populations in Proteins Using the Program Anglesearch. Journal of Magnetic Resonance Series B, 1995, 108, 31-43.	1.6	27
76	The use of PM3 SCF MO quantum mechanical calculations to refine NMR-determined structures of complexes of antifolate drugs with dihydrofolate reductase in solution. Computational and Theoretical Chemistry, 1995, 357, 207-216.	1.5	9
77	Solution Structure of a Brodimoprim Analog in Its Complex with Lactobacillus casei Dihydrofolate Reductase. Biochemistry, 1995, 34, 11690-11702.	2.5	23
78	Solution Structure of Bound Trimethoprim in Its Complex with Lactobacillus casei Dihydrofolate Reductase. Biochemistry, 1994, 33, 12416-12426.	2.5	38
79	Synthesis of 4-nitro-1,5-diarylpyrazoles based upon?-nitroacetophenones. Pharmaceutical Chemistry Journal, 1993, 26, 889-893.	0.8	3
80	New synthesis of 2-aryl-3-hydroxy(alkoxy)-4-quinolones by ring expansion of 1-acetyl-2-arylmethylene-3-indolinones. Chemistry of Heterocyclic Compounds, 1992, 28, 234-235.	1.2	1
81	1-Acetyl-2-chloro-3-iminoindoline hydrochloride and its N-acetyl derivatives in nucleophilic substitution reactions. Chemistry of Heterocyclic Compounds, 1992, 28, 43-47.	1.2	O
82	Pyrido [2,3-d]pyrimidines 4. Synthesis and some transformations of oxo(hydroxy)pyrido [2,3-d]pyrimidines. Chemistry of Heterocyclic Compounds, 1991, 27, 538-544.	1.2	0
83	1-Acetyl-2-bromo-3-indolinone in nucleophilic substitution reactions and the synthesis of pyrrolo[3,2-b]indoles. Pharmaceutical Chemistry Journal, 1990, 24, 917-923.	0.8	2
84	Interaction of spirazidine, prospidine, and spirobromine with components of nucleic acids. Pharmaceutical Chemistry Journal, 1989, 23, 1-8.	0.8	0
85	Pharmacokinetics and metabolism of spirobromine in oral administration. Pharmaceutical Chemistry Journal, 1988, 22, 582-584.	0.8	0
86	Reactions of spirazidine in aqueous media. Pharmaceutical Chemistry Journal, 1988, 22, 871-874.	0.8	0
87	2-Allylaminothiazolin-4-one in acylation reactions. Chemistry of Heterocyclic Compounds, 1987, 23, 910-914.	1.2	1
88	Interaction of the antitumor drugs phospidine and spirobromine with nucleotides. Pharmaceutical Chemistry Journal, 1987, 21, 301-308.	0.8	1
89	Investigation of the pathways of biotransformation of spirobromine. Pharmaceutical Chemistry Journal, 1986, 20, 519-525.	0.8	0
90	Photochemical reactions of biologically important quinoxaline N-oxides. Pharmaceutical Chemistry Journal, 1986, 20, 227-235.	0.8	0

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91	Synthesis and properties of 4-0x0-3,4-dihydrobenzo[g]pteridine-2-carboxylic acid and its derivatives. Pharmaceutical Chemistry Journal, 1986, 20, 117-122.	0.8	O
92	Study of the alkylating capacity of prospidin in model reactions with organic phosphates. Pharmaceutical Chemistry Journal, 1985, 19, 376-381.	0.8	0
93	Conversions of the antineoplastic preparation prospidin in aqueous media. Pharmaceutical Chemistry Journal, 1984, 18, 521-527.	0.8	O
94	31P and 1H NMR study of reaction between hydride complexes of platinum and divalent tin halides. Bulletin of the Academy of Sciences of the USSR Division of Chemical Science, 1982, 31, 917-921.	0.0	4
95	Williams-Beuren Syndrome Related Methyltransferase WBSCR27: From Structure to Possible Function. Frontiers in Molecular Biosciences, 0, 9, .	3.5	1