

George I Makhatadze

List of Publications by Year in descending order

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96
papers

7,339
citations

66343

42
h-index

54911

84
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120
all docs

120
docs citations

120
times ranked

5975
citing authors

#	ARTICLE	IF	CITATIONS
1	From Protein Design to the Energy Landscape of a Cold Unfolding Protein. <i>Journal of Physical Chemistry B</i> , 2022, 126, 1212-1231.	2.6	3
2	Effects of Hydrostatic Pressure on the Thermodynamics of CspB-Bs Interactions with the ssDNA Template. <i>Biochemistry</i> , 2021, 60, 3086-3097.	2.5	2
3	Evolutionary conservation and structural localizations suggest a physical trace of metabolism's progressive geochronological emergence. <i>Journal of Biomolecular Structure and Dynamics</i> , 2020, 38, 3700-3719.	3.5	3
4	Protein adaptation to high hydrostatic pressure: Computational analysis of the structural proteome. <i>Proteins: Structure, Function and Bioinformatics</i> , 2020, 88, 584-592.	2.6	12
5	Sequence-independent recognition of the amyloid structural motif by GFP protein family. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2020, 117, 22122-22127.	7.1	7
6	Strategy for Stabilization of CutA1 Proteins Due to Ion-Ion Interactions at Temperatures of over 100 Å°C. <i>Biochemistry</i> , 2018, 57, 2649-2656.	2.5	7
7	Predictive Model of Linear Antimicrobial Peptides Active against Gram-Negative Bacteria. <i>Journal of Chemical Information and Modeling</i> , 2018, 58, 1141-1151.	5.4	57
8	Evidence for the principle of minimal frustration in the evolution of protein folding landscapes. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2017, 114, E1627-E1632.	7.1	59
9	Reply to Candell et al.: Evidence for evolutionary conservation of folding kinetics in the thioredoxin protein family. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2017, 114, E4124-E4124.	7.1	2
10	Non-linearity of the collagen triple helix in solution and implications for collagen function. <i>Biochemical Journal</i> , 2017, 474, 2203-2217.	3.7	21
11	Linking computation and experiments to study the role of charge-charge interactions in protein folding and stability. <i>Physical Biology</i> , 2017, 14, 013002.	1.8	8
12	Molecular determinant of the effects of hydrostatic pressure on protein folding stability. <i>Nature Communications</i> , 2017, 8, 14561.	12.8	75
13	Molecular Determinants of Temperature Dependence of Protein Volume Change upon Unfolding. <i>Journal of Physical Chemistry B</i> , 2017, 121, 8300-8310.	2.6	13
14	Putting the Piezolyte Hypothesis under Pressure. <i>Biophysical Journal</i> , 2017, 113, 974-977.	0.5	13
15	Intramolecular diffusion controls aggregation of the PAPf39 peptide. <i>Biophysical Chemistry</i> , 2016, 216, 37-43.	2.8	11
16	Influence of surface charge, binding site residues and glycosylation on <i>Thielavia terrestris</i> cutinase biochemical characteristics. <i>Applied Microbiology and Biotechnology</i> , 2016, 100, 4435-4446.	3.6	25
17	Applications of pressure perturbation calorimetry to study factors contributing to the volume changes upon protein unfolding. <i>Biochimica Et Biophysica Acta - General Subjects</i> , 2016, 1860, 1036-1042.	2.4	5
18	Gallic Acid Is an Antagonist of Semen Amyloid Fibrils That Enhance HIV-1 Infection. <i>Journal of Biological Chemistry</i> , 2016, 291, 14045-14055.	3.4	12

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19	Modulation of folding energy landscape by charge–charge interactions: Linking experiments with computational modeling. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2015, 112, E259-66.	7.1	48
20	Thermal expansivities of peptides, polypeptides and proteins as measured by pressure perturbation calorimetry. <i>Methods</i> , 2015, 76, 61-66.	3.8	5
21	ProteinVolume: calculating molecular van der Waals and void volumes in proteins. <i>BMC Bioinformatics</i> , 2015, 16, 101.	2.6	84
22	Alterations of Nonconserved Residues Affect Protein Stability and Folding Dynamics through Charge–Charge Interactions. <i>Journal of Physical Chemistry B</i> , 2015, 119, 13103-13112.	2.6	22
23	Molecular Determinants of Expansivity of Native Globular Proteins: A Pressure Perturbation Calorimetry Study. <i>Journal of Physical Chemistry B</i> , 2014, 118, 6117-6122.	2.6	8
24	Structural Characterization of Semen Coagulum-Derived SEM1(86–107) Amyloid Fibrils That Enhance HIV-1 Infection. <i>Biochemistry</i> , 2014, 53, 3267-3277.	2.5	17
25	Structural and thermodynamic characterization of the recognition of the S100-binding peptides TRTK12 and p53 by calmodulin. <i>Protein Science</i> , 2014, 23, 1247-1261.	7.6	6
26	Novel Interactions of the TRTK12 Peptide with S100 Protein Family Members: Specificity and Thermodynamic Characterization. <i>Biochemistry</i> , 2013, 52, 5844-5856.	2.5	10
27	Bacterial expression and purification of the amyloidogenic peptide PAPf39 for multidimensional NMR spectroscopy. <i>Protein Expression and Purification</i> , 2013, 88, 196-200.	1.3	5
28	The Role of Cross-Chain Ionic Interactions for the Stability of Collagen Model Peptides. <i>Biophysical Journal</i> , 2013, 105, 1681-1688.	0.5	29
29	Backtracking due to Residual Structure in the Unfolded State Changes the Folding of the Third Fibronectin Type III Domain from Tenascin-C. <i>Journal of Physical Chemistry B</i> , 2013, 117, 800-810.	2.6	12
30	Crystal and NMR structures of a Trp-cage mini-protein benchmark for computational fold prediction. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2012, 109, 12521-12525.	7.1	31
31	Thermodynamic and Kinetic Analysis of Peptides Derived from CapZ, NDR, p53, HDM2, and HDM4 Binding to Human S100B. <i>Biochemistry</i> , 2012, 51, 7189-7201.	2.5	16
32	Core Sequence of PAPf39 Amyloid Fibrils and Mechanism of pH-Dependent Fibril Formation: The Role of Monomer Conformation. <i>Biochemistry</i> , 2012, 51, 10127-10136.	2.5	30
33	Molecular Mechanism for the Preferential Exclusion of TMAO from Protein Surfaces. <i>Journal of Physical Chemistry B</i> , 2012, 116, 12095-12104.	2.6	162
34	Electrostatic Contribution of Surface Charge Residues to the Stability of a Thermophilic Protein: Benchmarking Experimental and Predicted pKa Values. <i>PLoS ONE</i> , 2012, 7, e30296.	2.5	34
35	Enzyme Activity in the Crowded Milieu. <i>PLoS ONE</i> , 2012, 7, e39418.	2.5	54
36	Kinetic consequences of native state optimization of surface-exposed electrostatic interactions in the Fyn SH3 domain. <i>Proteins: Structure, Function and Bioinformatics</i> , 2012, 80, 858-870.	2.6	42

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37	Contribution of buried aspartic acid to the stability of the PDZ2 protein. <i>Journal of Chemical Thermodynamics</i> , 2012, 52, 64-68.	2.0	2
38	Protonation/deprotonation effects on the stability of the Trp-cage miniprotein. <i>Physical Chemistry Chemical Physics</i> , 2011, 13, 17056.	2.8	21
39	Statistical analysis of protein structures suggests that buried ionizable residues in proteins are hydrogen bonded or form salt bridges. <i>Proteins: Structure, Function and Bioinformatics</i> , 2011, 79, 2027-2032.	2.6	27
40	Energetics of charge-charge interactions between residues adjacent in sequence. <i>Proteins: Structure, Function and Bioinformatics</i> , 2011, 79, 3494-3499.	2.6	14
41	Thermodynamics of the Trp-cage miniprotein unfolding in urea. <i>Proteins: Structure, Function and Bioinformatics</i> , 2010, 78, 1376-1381.	2.6	29
42	Pressure Perturbation Calorimetry of Unfolded Proteins. <i>Journal of Physical Chemistry B</i> , 2010, 114, 16166-16170.	2.6	21
43	Rational stabilization of enzymes by computational redesign of surface charge-charge interactions. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2009, 106, 2601-2606.	7.1	201
44	Chapter 7 A Computational Approach for the Rational Design of Stable Proteins and Enzymes. <i>Methods in Enzymology</i> , 2009, 454, 175-211.	1.0	29
45	Mechanism of Fibril Formation by a 39-Residue Peptide (PAPf39) from Human Prostatic Acidic Phosphatase. <i>Biochemistry</i> , 2009, 48, 11582-11591.	2.5	33
46	Universal Convergence of the Specific Volume Changes of Globular Proteins upon Unfolding. <i>Biochemistry</i> , 2009, 48, 10846-10851.	2.5	36
47	Chapter 22 Use of Pressure Perturbation Calorimetry to Characterize the Volumetric Properties of Proteins. <i>Methods in Enzymology</i> , 2009, 466, 527-547.	1.0	31
48	Annexin I and Annexin II N-Terminal Peptides Binding to S100 Protein Family Members: Specificity and Thermodynamic Characterization. <i>Biochemistry</i> , 2009, 48, 2788-2798.	2.5	39
49	Protein Stabilization by the Rational Design of Surface Charge-Charge Interactions. <i>Methods in Molecular Biology</i> , 2009, 490, 261-283.	0.9	27
50	Cooperativity of complex salt bridges. <i>Protein Science</i> , 2008, 17, 1285-1290.	7.6	35
51	Heat capacity changes upon burial of polar and nonpolar groups in proteins. <i>Protein Science</i> , 2008, 10, 1343-1352.	7.6	120
52	Role of the Charge-Charge Interactions in Defining Stability and Halophilicity of the CspB Proteins. <i>Journal of Molecular Biology</i> , 2007, 366, 842-856.	4.2	62
53	Unfolding Thermodynamics of Trp-Cage, a 20 Residue Miniprotein, Studied by Differential Scanning Calorimetry and Circular Dichroism Spectroscopy. <i>Biochemistry</i> , 2007, 46, 2876-2880.	2.5	115
54	Computational design of the Fyn SH3 domain with increased stability through optimization of surface charge-charge interactions. <i>Protein Science</i> , 2007, 16, 2694-2702.	7.6	56

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55	Calorimetric Evidence for a Two-State Unfolding of the Î²-Hairpin Peptide Trpzip4. <i>Journal of the American Chemical Society</i> , 2006, 128, 30-31.	13.7	49
56	Protein Stability and Surface Electrostatics: A Charged Relationship. <i>Biochemistry</i> , 2006, 45, 2761-2766.	2.5	285
57	How to improve nature: study of the electrostatic properties of the surface of Î±-lactalbumin. <i>Protein Engineering, Design and Selection</i> , 2005, 18, 425-433.	2.1	40
58	Enthalpy of helix-coil transition: Missing link in rationalizing the thermodynamics of helix-forming propensities of the amino acid residues. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2005, 102, 1413-1418.	7.1	52
59	Thermodynamics Of Î±-Helix Formation. <i>Advances in Protein Chemistry</i> , 2005, 72, 199-226.	4.4	42
60	Effects of Charge-to-Alanine Substitutions on the Stability of Ribosomal Protein L30e from <i>Thermococcus celer</i> . <i>Biochemistry</i> , 2005, 44, 16817-16825.	2.5	43
61	Temperature Dependence of the Thermodynamics of Helix-Coil Transition. <i>Journal of Molecular Biology</i> , 2004, 335, 1029-1037.	4.2	70
62	Mechanism of Thermostabilization in a Designed Cold Shock Protein with Optimized Surface Electrostatic Interactions. <i>Journal of Molecular Biology</i> , 2004, 336, 929-942.	4.2	74
63	Noncharged amino acid residues at the solvent-exposed positions in the middle and at the C terminus of the Î±-helix have the same helical propensity. <i>Protein Science</i> , 2003, 12, 1169-1176.	7.6	29
64	Contribution of Surface Salt Bridges to Protein Stability: Guidelines for Protein Engineering. <i>Journal of Molecular Biology</i> , 2003, 327, 1135-1148.	4.2	216
65	Isothermal Titration Calorimetry. , 2002, 173, 121-126.		24
66	The enthalpy of the alanine peptide helix measured by isothermal titration calorimetry using metal-binding to induce helix formation. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2002, 99, 1298-1302.	7.1	104
67	Differential Scanning Calorimetry. , 2002, 173, 113-119.		41
68	Conformational and thermodynamic properties of peptide binding to the human S100P protein. <i>Protein Science</i> , 2002, 11, 1367-1375.	7.6	31
69	Thermodynamic Consequences of Burial of Polar and Non-polar Amino Acid Residues in the Protein Interior. <i>Journal of Molecular Biology</i> , 2002, 320, 343-357.	4.2	110
70	Removal of surface charge-charge interactions from ubiquitin leaves the protein folded and very stable. <i>Protein Science</i> , 2002, 11, 174-177.	7.6	26
71	Molecular Characterization and Tissue Distribution of a Novel Member of the S100 Family of EF-Hand Proteins. <i>Biochemistry</i> , 2001, 40, 15538-15548.	2.5	53
72	To charge or not to charge?. <i>Trends in Biotechnology</i> , 2001, 19, 132-135.	9.3	203

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73	Energetics of Target Peptide Binding by Calmodulin Reveals Different Modes of Binding. <i>Journal of Biological Chemistry</i> , 2001, 276, 14083-14091.	3.4	107
74	Interactions of the Cold Shock Protein CspB from <i>Bacillus subtilis</i> with Single-stranded DNA. <i>Journal of Biological Chemistry</i> , 2001, 276, 15511-15518.	3.4	51
75	Major cold shock proteins, CspA from <i>Escherichia coli</i> and CspB from <i>Bacillus subtilis</i> , interact differently with single-stranded DNA templates. <i>BBA - Proteins and Proteomics</i> , 2000, 1479, 196-202.	2.1	46
76	Contribution of proton linkage to the thermodynamic stability of the major cold shock protein of <i>Escherichia coli</i> CspA. <i>Protein Science</i> , 2000, 9, 387-394.	7.6	27
77	Interactions of the Major Cold Shock Protein of <i>Bacillus subtilis</i> CspB with Single-stranded DNA Templates of Different Base Composition. <i>Journal of Biological Chemistry</i> , 1999, 274, 33601-33608.	3.4	63
78	Heat capacity change for ribonuclease A folding. <i>Protein Science</i> , 1999, 8, 1500-1504.	7.6	77
79	Cold denaturation of ubiquitin. <i>BBA - Proteins and Proteomics</i> , 1999, 1429, 384-390.	2.1	69
80	Thermodynamics of Protein Interactions with Urea and Guanidinium Hydrochloride. <i>Journal of Physical Chemistry B</i> , 1999, 103, 4781-4785.	2.6	177
81	Thermal versus Guanidine-Induced Unfolding of Ubiquitin. An Analysis in Terms of the Contributions from Charge-Charge Interactions to Protein Stability. <i>Biochemistry</i> , 1999, 38, 8138-8149.	2.5	243
82	Engineering a Thermostable Protein via Optimization of Charge-Charge Interactions on the Protein Surface. <i>Biochemistry</i> , 1999, 38, 16419-16423.	2.5	213
83	MEARA Sequence Repeat of Human CstF-64 Polyadenylation Factor Is Helical in Solution. A Spectroscopic and Calorimetric Study. <i>Biochemistry</i> , 1999, 38, 12869-12875.	2.5	44
84	Cloning, overexpression, purification, and spectroscopic characterization of human S100P. <i>Protein Science</i> , 1998, 7, 211-215.	7.6	28
85	Anion binding to the ubiquitin molecule. <i>Protein Science</i> , 1998, 7, 689-697.	7.6	118
86	Heat capacities of amino acids, peptides and proteins. <i>Biophysical Chemistry</i> , 1998, 71, 133-156.	2.8	74
87	On the entropy of protein folding. <i>Protein Science</i> , 1996, 5, 507-510.	7.6	74
88	Solvent isotope effect and protein stability. <i>Nature Structural Biology</i> , 1995, 2, 852-855.	9.7	172
89	Energetics of Protein Structure. <i>Advances in Protein Chemistry</i> , 1995, 47, 307-425.	4.4	933
90	Thermodynamics of barnase unfolding. <i>Protein Science</i> , 1994, 3, 669-676.	7.6	74

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91	Effect of pH and phosphate ions on self-association properties of the major cold-shock protein from <i>Bacillus subtilis</i> . <i>Protein Science</i> , 1994, 3, 2144-2147.	7.6	39
92	Thermodynamics of ubiquitin unfolding. <i>Proteins: Structure, Function and Bioinformatics</i> , 1994, 18, 246-253.	2.6	170
93	Energetics of ribonuclease T1 structure. <i>Biochemistry</i> , 1994, 33, 3312-3319.	2.5	81
94	Thermodynamics of bpti folding. <i>Protein Science</i> , 1993, 2, 2028-2036.	7.6	102
95	Contribution of Hydration to Protein Folding Thermodynamics. <i>Journal of Molecular Biology</i> , 1993, 232, 660-679.	4.2	406
96	Protein interactions with urea and guanidinium chloride. <i>Journal of Molecular Biology</i> , 1992, 226, 491-505.	4.2	610