

Douglas H Turner

List of Publications by Year in descending order

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156
papers

15,733
citations

27035

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120
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201
all docs

201
docs citations

201
times ranked

10185
citing authors

#	ARTICLE	IF	CITATIONS
1	Expanded sequence dependence of thermodynamic parameters improves prediction of RNA secondary structure. <i>Journal of Molecular Biology</i> , 1999, 288, 911-940.	2.0	3,486
2	Incorporating chemical modification constraints into a dynamic programming algorithm for prediction of RNA secondary structure. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2004, 101, 7287-7292.	3.3	1,332
3	Thermodynamic Parameters for an Expanded Nearest-Neighbor Model for Formation of RNA Duplexes with Watson-Crick Base Pairs. <i>Biochemistry</i> , 1998, 37, 14719-14735.	1.2	1,055
4	NNDB: the nearest neighbor parameter database for predicting stability of nucleic acid secondary structure. <i>Nucleic Acids Research</i> , 2010, 38, D280-D282.	6.5	444
5	Dynalign: an algorithm for finding the secondary structure common to two RNA sequences. <i>Journal of Molecular Biology</i> , 2002, 317, 191-203.	2.0	340
6	Prediction of RNA secondary structure by free energy minimization. <i>Current Opinion in Structural Biology</i> , 2006, 16, 270-278.	2.6	339
7	Investigation of the Structural Basis for Thermodynamic Stabilities of Tandem GU Mismatches: A Solution Structure of (rGAGGUCUC) ₂ by Two-Dimensional NMR and Simulated Annealing. <i>Biochemistry</i> , 1996, 35, 14077-14089.	1.2	333
8	Microarrays for identifying binding sites and probing structure of RNAs. <i>Nucleic Acids Research</i> , 2015, 43, 1-12.	6.5	250
9	Predicting oligonucleotide affinity to nucleic acid targets. <i>Rna</i> , 1999, 5, 1458-1469.	1.6	228
10	Expanded CUG repeat RNAs form hairpins that activate the double-stranded RNA-dependent protein kinase PKR. <i>Rna</i> , 2000, 6, 79-87.	1.6	225
11	The contribution of pseudouridine to stabilities and structure of RNAs. <i>Nucleic Acids Research</i> , 2014, 42, 3492-3501.	6.5	177
12	Laser temperature-jump, spectroscopic, and thermodynamic study of salt effects on duplex formation by dGCATGC. <i>Biochemistry</i> , 1989, 28, 4283-4291.	1.2	176
13	Structure of (rGGCGAGCC) ₂ in solution from NMR and restrained molecular dynamics. <i>Biochemistry</i> , 1993, 32, 12612-12623.	1.2	175
14	Thermodynamics of Single Mismatches in RNA Duplexes. <i>Biochemistry</i> , 1999, 38, 14214-14223.	1.2	166
15	Stabilities of consecutive A.cntdot.C, C.cntdot.C, G.cntdot.G, U.cntdot.C, and U.cntdot.U mismatches in RNA internal loops: evidence for stable hydrogen-bonded U.cntdot.U and C.cntdot.C+ pairs. <i>Biochemistry</i> , 1991, 30, 8242-8251.	1.2	164
16	Thermodynamic and spectroscopic study of bulge loops in oligoribonucleotides. <i>Biochemistry</i> , 1990, 29, 278-285.	1.2	161
17	Free energy increments for hydrogen bonds in nucleic acid base pairs. <i>Journal of the American Chemical Society</i> , 1987, 109, 3783-3785.	6.6	158
18	Reparameterization of RNA \ddagger Torsion Parameters for the AMBER Force Field and Comparison to NMR Spectra for Cytidine and Uridine. <i>Journal of Chemical Theory and Computation</i> , 2010, 6, 1520-1531.	2.3	155

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19	Polymer-supported RNA synthesis and its application to test the nearest-neighbor model for duplex stability. <i>Biochemistry</i> , 1986, 25, 7840-7846.	1.2	148
20	A Periodic Table of Tandem Mismatches in RNA. <i>Biochemistry</i> , 1995, 34, 3204-3211.	1.2	140
21	Folding and Finding RNA Secondary Structure. <i>Cold Spring Harbor Perspectives in Biology</i> , 2010, 2, a003665-a003665.	2.3	136
22	Stability of XGCGCp, GCGCYp, and XGCGCYp helices: an empirical estimate of the energetics of hydrogen bonds in nucleic acids. <i>Biochemistry</i> , 1986, 25, 3214-3219.	1.2	134
23	Effects of 3' dangling end stacking on the stability of GGCC and CCGG double helices. <i>Biochemistry</i> , 1983, 22, 6198-6206.	1.2	131
24	Comparison of binding of mixed ribose-deoxyribose analogs of CUCU to a ribozyme and to GGAGAA by equilibrium dialysis: evidence for ribozyme specific interactions with 2'-hydroxy groups. <i>Biochemistry</i> , 1991, 30, 10632-10640.	1.2	129
25	Contributions of dangling end stacking and terminal base-pair formation to the stabilities of XGGCCp, XCCGGp, XGGCCYp, and XCCGGYp helices. <i>Biochemistry</i> , 1985, 24, 4533-4539.	1.2	128
26	Nearest-neighbor parameters for G.cntdot.U mismatches: 5'GU3'/3'UG5' is destabilizing in the contexts CGUG/GUGC, UGUA/AUGU, and AGUU/UUGA but stabilizing in GGUC/CUGC. <i>Biochemistry</i> , 1991, 30, 11124-11132.	1.2	128
27	Thermodynamics of RNA-RNA Duplexes with 2- or 4-Thiouridines: Implications for Antisense Design and Targeting a Group I Intron. <i>Biochemistry</i> , 1999, 38, 16655-16662.	1.2	118
28	Effects of GA mismatches on the structure and thermodynamics of RNA internal loops. <i>Biochemistry</i> , 1990, 29, 8813-8819.	1.2	117
29	Solution Structure of (rGCGGACGC) ₂ by Two-Dimensional NMR and the Iterative Relaxation Matrix Approach. <i>Biochemistry</i> , 1996, 35, 9677-9689.	1.2	108
30	The influence of locked nucleic acid residues on the thermodynamic properties of 2'-O-methyl RNA/RNA heteroduplexes. <i>Nucleic Acids Research</i> , 2005, 33, 5082-5093.	6.5	104
31	Laser temperature-jump study of stacking in adenylic acid polymers. <i>Biochemistry</i> , 1979, 18, 5757-5762.	1.2	102
32	Measuring the thermodynamics of RNA secondary structure formation. , 1997, 44, 309-319.		100
33	Stacking in RNA: NMR of Four Tetramers Benchmark Molecular Dynamics. <i>Journal of Chemical Theory and Computation</i> , 2015, 11, 2729-2742.	2.3	99
34	Conformational ensembles of RNA oligonucleotides from integrating NMR and molecular simulations. <i>Science Advances</i> , 2018, 4, eaar8521.	4.7	99
35	Thermodynamic study of internal loops in oligoribonucleotides: symmetric loops are more stable than asymmetric loops. <i>Biochemistry</i> , 1991, 30, 6428-6436.	1.2	98
36	Benchmarking AMBER Force Fields for RNA: Comparisons to NMR Spectra for Single-Stranded r(GACC) Are Improved by Revised \ddagger Torsions. <i>Journal of Physical Chemistry B</i> , 2011, 115, 9261-9270.	1.2	95

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37	Solvent effects on the dynamics of (dG-dC) ₃ . <i>Biopolymers</i> , 1983, 22, 1107-1131.	1.2	92
38	Sequence dependence of stability for coaxial stacking of RNA helices with Watson-Crick base paired interfaces. <i>Biochemistry</i> , 1994, 33, 12715-12719.	1.2	89
39	Thermodynamics of Three-Way Multibranch Loops in RNA. <i>Biochemistry</i> , 2001, 40, 6971-6981.	1.2	88
40	The Stability and Structure of Tandem GA Mismatches in RNA Depend on Closing Base Pairs. <i>Biochemistry</i> , 1994, 33, 11349-11354.	1.2	87
41	Optical Melting Measurements of Nucleic Acid Thermodynamics. <i>Methods in Enzymology</i> , 2009, 468, 371-387.	0.4	86
42	Identification of potential conserved RNA secondary structure throughout influenza A coding regions. <i>Rna</i> , 2011, 17, 991-1011.	1.6	85
43	Thermodynamics of Nonsymmetric Tandem Mismatches Adjacent to G-C Base Pairs in RNA. <i>Biochemistry</i> , 1997, 36, 12486-12497.	1.2	84
44	Solution Structure of (rGGCAGGCC) ₂ by Two-Dimensional NMR and the Iterative Relaxation Matrix Approach. <i>Biochemistry</i> , 1997, 36, 4449-4460.	1.2	82
45	Testing the Nearest Neighbor Model for Canonical RNA Base Pairs: Revision of GU Parameters. <i>Biochemistry</i> , 2012, 51, 3508-3522.	1.2	80
46	Thermodynamics of unpaired terminal nucleotides on short RNA helices correlates with stacking at helix termini in larger RNAs. <i>Journal of Molecular Biology</i> , 1999, 290, 967-982.	2.0	79
47	The Nuclear Magnetic Resonance of CCCC RNA Reveals a Right-Handed Helix, and Revised Parameters for AMBER Force Field Torsions Improve Structural Predictions from Molecular Dynamics. <i>Biochemistry</i> , 2013, 52, 996-1010.	1.2	78
48	Solvent effects on the kinetics and thermodynamics of stacking in poly(cytidylic acid). <i>Biochemistry</i> , 1981, 20, 1419-1426.	1.2	77
49	A model for the stabilities of RNA hairpins based on a study of the sequence dependence of stability for hairpins of six nucleotides. <i>Biochemistry</i> , 1994, 33, 14289-14296.	1.2	76
50	Effects of substrate structure on the kinetics of circle opening reactions of the self-splicing intervening sequence from <i>Tetrahymena thermophila</i> : evidence for substrate and Mg ²⁺ binding interactions. <i>Nucleic Acids Research</i> , 1989, 17, 355-371.	6.5	74
51	Factors Affecting the Thermodynamic Stability of Small Asymmetric Internal Loops in RNA. <i>Biochemistry</i> , 2000, 39, 9257-9274.	1.2	73
52	Experimentally Derived Nearest-Neighbor Parameters for the Stability of RNA Three- and Four-Way Multibranch Loops. <i>Biochemistry</i> , 2002, 41, 869-880.	1.2	73
53	Structure determination of noncanonical RNA motifs guided by ¹ H NMR chemical shifts. <i>Nature Methods</i> , 2014, 11, 413-416.	9.0	72
54	Long-Range Cooperativity in Molecular Recognition of RNA by Oligodeoxynucleotides with Multiple C5-(1-Propynyl) Pyrimidines. <i>Journal of the American Chemical Society</i> , 2001, 123, 4107-4118.	6.6	69

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55	Replacement of the Conserved G.cntdot.U with a G-C Pair at the Cleavage Site of the Tetrahymena Ribozyme Decreases Binding, Reactivity, and Fidelity. <i>Biochemistry</i> , 1994, 33, 13856-13863.	1.2	66
56	Structural Features of a Six-Nucleotide RNA Hairpin Loop Found in Ribosomal RNA. <i>Biochemistry</i> , 1996, 35, 6539-6548.	1.2	66
57	Revision of AMBER Torsional Parameters for RNA Improves Free Energy Predictions for Tetramer Duplexes with GC and iGiC Base Pairs. <i>Journal of Chemical Theory and Computation</i> , 2012, 8, 172-181.	2.3	65
58	NMR Structures of r(GCAGGCGUGC) ₂ and Determinants of Stability for Single Guanosine-Guanosine Base Pairs. <i>Biochemistry</i> , 2000, 39, 11748-11762.	1.2	61
59	Oligonucleotide directed misfolding of RNA inhibits <i>Candida albicans</i> group I intron splicing. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2002, 99, 11091-11096.	3.3	59
60	Nuclear Magnetic Resonance Spectroscopy and Molecular Modeling Reveal That Different Hydrogen Bonding Patterns Are Possible for G-U Pairs: One Hydrogen Bond for Each G-U Pair in r(GCGUGCC) ₂ and Two for Each G-U Pair in r(GAGUGUC) ₂ . <i>Biochemistry</i> , 2000, 39, 8970-8982.	1.2	56
61	Melting and chemical modification of a cyclized self-splicing group I intron: similarity of structures in 1 M sodium, in 10 mM magnesium and in the presence of substrate. <i>Biochemistry</i> , 1990, 29, 10147-10158.	1.2	55
62	The time dependence of chemical modification reveals slow steps in the folding of a group I ribozyme. <i>Biochemistry</i> , 1995, 34, 6504-6512.	1.2	55
63	Investigation of the Structural Basis for Thermodynamic Stabilities of Tandem GU Wobble Pairs: NMR Structures of (rGGAGUUC) ₂ and (rGGAUGUC) ₂ . <i>Biochemistry</i> , 1997, 36, 8030-8038.	1.2	55
64	Thermodynamic Studies of RNA Stability. <i>Journal of Biomolecular Structure and Dynamics</i> , 1984, 1, 1229-1242.	2.0	52
65	Thermodynamics of Coaxially Stacked Helices with GA and CC Mismatches. <i>Biochemistry</i> , 1996, 35, 13753-13761.	1.2	52
66	The Determination of RNA Folding Nearest Neighbor Parameters. <i>Methods in Molecular Biology</i> , 2014, 1097, 45-70.	0.4	52
67	In vivo analysis of influenza A mRNA secondary structures identifies critical regulatory motifs. <i>Nucleic Acids Research</i> , 2019, 47, 7003-7017.	6.5	51
68	RNA Challenges for Computational Chemists. <i>Biochemistry</i> , 2005, 44, 13225-13234.	1.2	47
69	Sequence dependence for the energetics of terminal mismatches in ribooligonucleotides. <i>Biochemistry</i> , 1987, 26, 4559-4562.	1.2	46
70	Stability and Structure of RNA Duplexes Containing Isoguanosine and Isocytidine. <i>Journal of the American Chemical Society</i> , 2001, 123, 1267-1274.	6.6	46
71	An RNA Molecular Switch: Intrinsic Flexibility of 23S rRNA Helices 40 and 68 5'-UAA/5'-GAN Internal Loops Studied by Molecular Dynamics Methods. <i>Journal of Chemical Theory and Computation</i> , 2010, 6, 910-929.	2.3	46
72	NMR-Assisted Prediction of RNA Secondary Structure: Identification of a Probable Pseudoknot in the Coding Region of an R2 Retrotransposon. <i>Journal of the American Chemical Society</i> , 2008, 130, 10233-10239.	6.6	45

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73	Contributions of Stacking, Preorganization, and Hydrogen Bonding to the Thermodynamic Stability of Duplexes between RNA and 2'-O-Methyl RNA with Locked Nucleic Acids. <i>Biochemistry</i> , 2009, 48, 4377-4387.	1.2	43
74	Facilitating RNA Structure Prediction with Microarrays. <i>Biochemistry</i> , 2006, 45, 581-593.	1.2	42
75	APneumocystis carinii Group I Intron Ribozyme That Does Not Require 2'-OH Groups on Its 5' Exon Mimic for Binding to the Catalytic Core. <i>Biochemistry</i> , 1997, 36, 15303-15314.	1.2	41
76	Thermodynamic Stabilities of Internal Loops with GU Closing Pairs in RNA. <i>Biochemistry</i> , 2001, 40, 11509-11517.	1.2	41
77	Effects of Restrained Sampling Space and Nonplanar Amino Groups on Free-Energy Predictions for RNA with Imino and Sheared Tandem GA Base Pairs Flanked by GC, CG, iGiC or iCiG Base Pairs. <i>Journal of Chemical Theory and Computation</i> , 2009, 5, 2088-2100.	2.3	39
78	The 3' Splice Site of Influenza A Segment 7 mRNA Can Exist in Two Conformations: A Pseudoknot and a Hairpin. <i>PLoS ONE</i> , 2012, 7, e38323.	1.1	39
79	Isoenergetic penta- and hexanucleotide microarray probing and chemical mapping provide a secondary structure model for an RNA element orchestrating R2 retrotransposon protein function. <i>Nucleic Acids Research</i> , 2008, 36, 1770-1782.	6.5	37
80	Solution Structure of an RNA Internal Loop with Three Consecutive Sheared GA Pairs. <i>Biochemistry</i> , 2005, 44, 2845-2856.	1.2	36
81	Nearest neighbor parameters for Watson-Crick complementary heteroduplexes formed between 2'-O-methyl RNA and RNA oligonucleotides. <i>Nucleic Acids Research</i> , 2006, 34, 3609-3614.	6.5	36
82	GA and UA Mismatches Can Stabilize RNA Internal Loops of Three Nucleotides. <i>Biochemistry</i> , 1996, 35, 16105-16109.	1.2	35
83	Secondary Structures for 5' Regions of R2 Retrotransposon RNAs Reveal a Novel Conserved Pseudoknot and Regions that Evolve under Different Constraints. <i>Journal of Molecular Biology</i> , 2009, 390, 428-442.	2.0	35
84	A chemical synthesis of LNA-2,6-diaminopurine riboside, and the influence of 2'-O-methyl-2,6-diaminopurine and LNA-2,6-diaminopurine ribosides on the thermodynamic properties of 2'-O-methyl RNA/RNA heteroduplexes. <i>Nucleic Acids Research</i> , 2007, 35, 4055-4063.	6.5	34
85	Antisense Oligonucleotides Targeting Influenza A Segment 8 Genomic RNA Inhibit Viral Replication. <i>Nucleic Acid Therapeutics</i> , 2016, 26, 277-285.	2.0	34
86	A Mechanistic Framework for the Second Step of Splicing Catalyzed by the Tetrahymena Ribozyme. <i>Biochemistry</i> , 1996, 35, 648-658.	1.2	33
87	Factors Affecting Thermodynamic Stabilities of RNA 3' Internal Loops. <i>Biochemistry</i> , 2004, 43, 12865-12876.	1.2	33
88	NMR Structures of (rGCUGAGGCU) ₂ and (rGCGGAUGCU) ₂ : Probing the Structural Features That Shape the Thermodynamic Stability of GA Pairs. <i>Biochemistry</i> , 2007, 46, 1511-1522.	1.2	33
89	Optimization of an AMBER Force Field for the Artificial Nucleic Acid, LNA, and Benchmarking with NMR of L(CAAU). <i>Journal of Physical Chemistry B</i> , 2014, 118, 1216-1228.	1.2	32
90	Physics-based atom modeling of RNA energetics and structure. <i>Wiley Interdisciplinary Reviews RNA</i> , 2017, 8, e1422.	3.2	32

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91	The energetics of small internal loops in RNA. <i>Biopolymers</i> , 1999, 52, 157-167.	1.2	31
92	Uptake and antifungal activity of oligonucleotides in <i>Candida albicans</i> . <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2003, 100, 1530-1534.	3.3	31
93	RNA Internal Loops with Tandem AC Pairs: The Structure of the 5'GAGU/3'UAGA Loop Can Be Dramatically Different from Others, Including 5'AAGU/3'UAGA. <i>Biochemistry</i> , 2010, 49, 5817-5827.	1.2	31
94	Self-Folding of Naked Segment 8 Genomic RNA of Influenza A Virus. <i>PLoS ONE</i> , 2016, 11, e0148281.	1.1	31
95	Influenza A Virus Coding Regions Exhibit Host-Specific Global Ordered RNA Structure. <i>PLoS ONE</i> , 2012, 7, e35989.	1.1	30
96	RNA pseudoknots: folding and finding. <i>F1000 Biology Reports</i> , 2010, 2, 8.	4.0	30
97	The NMR Structure of an Internal Loop from 23S Ribosomal RNA Differs from Its Structure in Crystals of 50S Ribosomal Subunits. <i>Biochemistry</i> , 2006, 45, 11776-11789.	1.2	28
98	An Alternating Sheared AA Pair and Elements of Stability for a Single Sheared Purine-Purine Pair Flanked by Sheared GA Pairs in RNA. <i>Biochemistry</i> , 2006, 45, 6889-6903.	1.2	27
99	Consecutive GA Pairs Stabilize Medium-Size RNA Internal Loops. <i>Biochemistry</i> , 2006, 45, 4025-4043.	1.2	27
100	Proton magnetic resonance melting studies of CCGGp, CCGGAp, ACCGGp, CCGGUp, and ACCGGUp. <i>Biochemistry</i> , 1983, 22, 269-277.	1.2	26
101	A CA ⁺ Pair Adjacent to a Sheared GA or AA Pair Stabilizes Size-Symmetric RNA Internal Loops. <i>Biochemistry</i> , 2009, 48, 5738-5752.	1.2	26
102	Secondary Structure of a Conserved Domain in the Intron of Influenza A NS1 mRNA. <i>PLoS ONE</i> , 2013, 8, e70615.	1.1	26
103	Predicting the Kinetics of RNA Oligonucleotides Using Markov State Models. <i>Journal of Chemical Theory and Computation</i> , 2017, 13, 926-934.	2.3	26
104	Mutations Designed by Ensemble Defect to Misfold Conserved RNA Structures of Influenza A Segments 7 and 8 Affect Splicing and Attenuate Viral Replication in Cell Culture. <i>PLoS ONE</i> , 2016, 11, e0156906.	1.1	26
105	The Thermodynamics of 3'-Terminal Pyrene and Guanosine for the Design of Isoenergetic 2'-O-Methyl-RNA-LNA Chimeric Oligonucleotide Probes of RNA Structure. <i>Biochemistry</i> , 2008, 47, 1249-1258.	1.2	25
106	The influenza A segment 7 mRNA 3' splice site pseudoknot/hairpin family. <i>RNA Biology</i> , 2012, 9, 1305-1310.	1.5	25
107	Binding Enhancement by Tertiary Interactions and Suicide Inhibition of a <i>Candida albicans</i> Group I Intron by Phosphoramidate and 2'-O-Methyl Hexanucleotides. <i>Biochemistry</i> , 2001, 40, 6520-6526.	1.2	24
108	NMR structure of a 4-4 nucleotide RNA internal loop from an R2 retrotransposon: Identification of a three-purine-purine sheared pair motif and comparison to MC-SYM predictions. <i>Rna</i> , 2011, 17, 1664-1677.	1.6	24

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109	Understanding the role of base stacking in nucleic acids. MD and QM analysis of tandem GA base pairs in RNA duplexes. <i>Physical Chemistry Chemical Physics</i> , 2012, 14, 12580.	1.3	24
110	Secondary Structure of a Conserved Domain in an Intron of Influenza A M1 mRNA. <i>Biochemistry</i> , 2014, 53, 5236-5248.	1.2	24
111	Crystal structure of a poly(rA) staggered zipper at acidic pH: evidence that adenine N1 protonation mediates parallel double helix formation. <i>Nucleic Acids Research</i> , 2016, 44, 8417-8424.	6.5	24
112	Antisense Binding Enhanced by Tertiary Interactions: Binding of Phosphorothioate and N ³ -Phosphoramidate Hexanucleotides to the Catalytic Core of a Group I Ribozyme from the Mammalian Pathogen <i>Pneumocystis carinii</i> . <i>Biochemistry</i> , 1998, 37, 9379-9385.	1.2	23
113	C5-(1-Propynyl)-2-deoxy-Pyrimidines Enhance Mismatch Penalties of DNA:RNA Duplex Formation. <i>Biochemistry</i> , 2001, 40, 12738-12745.	1.2	23
114	Recognition Elements for 5' Exon Substrate Binding to the <i>Candida albicans</i> Group I Intron. <i>Biochemistry</i> , 2001, 40, 6507-6519.	1.2	23
115	Thermodynamics of RNA Internal Loops with a Guanosine-Guanosine Pair Adjacent to Another Noncanonical Pair. <i>Biochemistry</i> , 2001, 40, 2478-2483.	1.2	23
116	Secondary structure models of the 3' untranslated regions of diverse R2 RNAs. <i>Rna</i> , 2004, 10, 978-987.	1.6	23
117	Targeting a <i>Pneumocystis carinii</i> Group I Intron with Methylphosphonate Oligonucleotides: A Backbone Charge Is Not Required for Binding or Reactivity. <i>Biochemistry</i> , 2000, 39, 6991-7000.	1.2	22
118	Structural Features and Thermodynamics of the J4/5 Loop from the <i>Candida albicans</i> and <i>Candida dubliniensis</i> Group I Introns. <i>Biochemistry</i> , 2004, 43, 15822-15837.	1.2	22
119	Improving RNA nearest neighbor parameters for helices by going beyond the two-state model. <i>Nucleic Acids Research</i> , 2018, 46, 4883-4892.	6.5	22
120	Nuclear Magnetic Resonance of Single-Stranded RNAs and DNAs of CAAU and UCAAUC as Benchmarks for Molecular Dynamics Simulations. <i>Journal of Chemical Theory and Computation</i> , 2020, 16, 1968-1984.	2.3	22
121	An Updated Recursive Algorithm for RNA Secondary Structure Prediction with Improved Thermodynamic Parameters. <i>ACS Symposium Series</i> , 1997, , 246-257.	0.5	21
122	NMR Reveals the Absence of Hydrogen Bonding in Adjacent UU and AG Mismatches in an Isolated Internal Loop from Ribosomal RNA. <i>Biochemistry</i> , 2007, 46, 12665-12678.	1.2	21
123	Chemical Synthesis of LNA-2-thiouridine and Its Influence on Stability and Selectivity of Oligonucleotide Binding to RNA. <i>Biochemistry</i> , 2009, 48, 10882-10893.	1.2	21
124	Biophysical Analysis of Influenza A Virus RNA Promoter at Physiological Temperatures. <i>Journal of Biological Chemistry</i> , 2011, 286, 22965-22970.	1.6	19
125	Inhibition of <i>Escherichia coli</i> RNase P by oligonucleotide directed misfolding of RNA. <i>Rna</i> , 2003, 9, 1437-1445.	1.6	18
126	New approaches to targeting RNA with oligonucleotides: Inhibition of group I intron self-splicing. <i>Biopolymers</i> , 2004, 73, 151-161.	1.2	18

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127	Stacking Effects on Local Structure in RNA: Changes in the Structure of Tandem GA Pairs when Flanking GC Pairs Are Replaced by isoG~isoC Pairs. <i>Journal of Physical Chemistry B</i> , 2007, 111, 6718-6727.	1.2	17
128	Activity of Hoechst 33258 against <i>Pneumocystis carinii</i> f. sp. muris, <i>Candida albicans</i> , and <i>Candida dubliniensis</i> . <i>Antimicrobial Agents and Chemotherapy</i> , 2005, 49, 1326-1330.	1.4	16
129	Novel Conformation of an RNA Structural Switch. <i>Biochemistry</i> , 2012, 51, 9257-9259.	1.2	16
130	Fluorescence Competition Assay Measurements of Free Energy Changes for RNA Pseudoknots. <i>Biochemistry</i> , 2010, 49, 623-634.	1.2	15
131	Structural Features of a 3' Splice Site in Influenza A. <i>Biochemistry</i> , 2015, 54, 3269-3285.	1.2	15
132	Accurate geometrical restraints for Watson-Crick base pairs. <i>Acta Crystallographica Section B: Structural Science, Crystal Engineering and Materials</i> , 2019, 75, 235-245.	0.5	14
133	Effects of Mg ²⁺ and the 2'-OH of Guanosine on Steps Required for Substrate Binding and Reactivity with the Tetrahymena Ribozyme Reveal Several Local Folding Transitions. <i>Biochemistry</i> , 1997, 36, 11131-11139.	1.2	13
134	Fundamental interactions in RNA: Questions answered and remaining. <i>Biopolymers</i> , 2013, 99, n/a-n/a.	1.2	13
135	Identification of conserved RNA secondary structures at influenza B and C splice sites reveals similarities and differences between influenza A, B, and C. <i>BMC Research Notes</i> , 2014, 7, 22.	0.6	13
136	Nuclear Magnetic Resonance-Assisted Prediction of Secondary Structure for RNA: Incorporation of Direction-Dependent Chemical Shift Constraints. <i>Biochemistry</i> , 2015, 54, 6769-6782.	1.2	13
137	Long-Range Cooperativity Due to C5-Propynylation of Oligopyrimidines Enhances Specific Recognition by Uridine of ribo-Adenosine over ribo-Guanosine. <i>Journal of the American Chemical Society</i> , 2001, 123, 9186-9187.	6.6	12
138	Nearest neighbor rules for RNA helix folding thermodynamics: improved end effects. <i>Nucleic Acids Research</i> , 2022, 50, 5251-5262.	6.5	12
139	Nuclear Magnetic Resonance Spectra and AMBER OL3 and ROC-RNA Simulations of UCUCGU Reveal Force Field Strengths and Weaknesses for Single-Stranded RNA. <i>Journal of Chemical Theory and Computation</i> , 2022, 18, 1241-1254.	2.3	11
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