

# Stephen Neidle

## List of Publications by Year in descending order

Source: <https://exaly.com/author-pdf/7478947/publications.pdf>

Version: 2024-02-01

381  
papers

34,077  
citations

3159

92  
h-index

4432

172  
g-index

406  
all docs

406  
docs citations

406  
times ranked

15276  
citing authors

#	ARTICLE	IF	CITATIONS
1	Non-standard and higher-order DNA structures: DNAâ€™DNA recognition. , 2022, , 109-190.		1
2	Principles of small moleculeâ€™DNA recognition. , 2022, , 191-286.		1
3	Structured Waters Mediate Small Molecule Binding to G-Quadruplex Nucleic Acids. <i>Pharmaceuticals</i> , 2022, 15, 7.	3.8	19
4	The mechanism of resistance in <i>Escherichia coli</i> to ridinilazole and other antibacterial head-to-head bis-benzimidazole compounds. <i>Medicinal Chemistry Research</i> , 2022, 31, 1176-1191.	2.4	1
5	Beyond the double helix: DNA structural diversity and the PDB. <i>Journal of Biological Chemistry</i> , 2021, 296, 100553.	3.4	25
6	Targeting the ALS/FTD-associated A-DNA kink with anthracene-based metal complex causes DNA backbone straightening and groove contraction. <i>Nucleic Acids Research</i> , 2021, 49, 9526-9538.	14.5	5
7	Water spines and networks in G-quadruplex structures. <i>Nucleic Acids Research</i> , 2021, 49, 519-528.	14.5	27
8	Asymmetrically Substituted Quadruplex-Binding Naphthalene Diimide Showing Potent Activity in Pancreatic Cancer Models. <i>ACS Medicinal Chemistry Letters</i> , 2020, 11, 1634-1644.	2.8	26
9	Challenges in developing small-molecule quadruplex therapeutics. <i>Annual Reports in Medicinal Chemistry</i> , 2020, , 517-546.	0.9	4
10	A G-Quadruplex-Binding Small Molecule and the HDAC Inhibitor SAHA (Vorinostat) Act Synergistically in Gemcitabine-Sensitive and Resistant Pancreatic Cancer Cells. <i>Molecules</i> , 2020, 25, 5407.	3.8	7
11	A G-quadruplex-binding compound shows potent activity in human gemcitabine-resistant pancreatic cancer cells. <i>Scientific Reports</i> , 2020, 10, 12192.	3.3	18
12	Substituted Naphthalenediimide Compounds Bind Selectively to Two Human Quadruplex Structures with Parallel Topology. <i>ACS Medicinal Chemistry Letters</i> , 2020, 11, 991-999.	2.8	16
13	Comment on â€œDark Nudges and Sludge in Big Alcohol: Behavioral Economics, Cognitive Biases, and Alcohol Industry Corporate Social Responsibilityâ€. <i>Milbank Quarterly</i> , 2020, 98, E1-E4.	4.4	2
14	Hierarchical Nanotube Selfâ€™Assembly of DNA Minor Grooveâ€™Binding Ligand DB921 via Alkali Halide Triggering. <i>Macromolecular Symposia</i> , 2019, 386, 1800243.	0.7	0
15	Polymorphic G:C mismatches act as hotspots for inducing right-handed Z DNA by DNA intercalation. <i>Nucleic Acids Research</i> , 2019, 47, 8899-8912.	14.5	16
16	Combining 1,3â€™Ditriazolylbenzene and Quinoline to Discover a New Gâ€™Quadruplexâ€™Interactive Small Molecule Active against Cancer Stemâ€™Like Cells. <i>ChemMedChem</i> , 2019, 14, 1325-1328.	3.2	13
17	The Targeting of Quadruplex Nucleic Acids in Human Cancers. <i>Proceedings (mdpi)</i> , 2019, 22, .	0.2	0
18	Dynamic self-assembly of DNA minor groove-binding ligand DB921 into nanotubes triggered by an alkali halide. <i>Nanoscale</i> , 2018, 10, 5550-5558.	5.6	6

#	ARTICLE	IF	CITATIONS
19	Accuracy of alcohol and breast cancer risk information on Drinkaware™'s website. <i>Drug and Alcohol Review</i> , 2018, 37, 304-306.	2.1	4
20	Targeting Multiple Effector Pathways in Pancreatic Ductal Adenocarcinoma with a G-Quadruplex-Binding Small Molecule. <i>Journal of Medicinal Chemistry</i> , 2018, 61, 2500-2517.	6.4	114
21	A naphthalene diimide G-quadruplex ligand inhibits cell growth and down-regulates BCL-2 expression in an imatinib-resistant gastrointestinal cancer cell line. <i>Bioorganic and Medicinal Chemistry</i> , 2018, 26, 2958-2964.	3.0	23
22	G-Quadruplex-binding small molecules ameliorate <i>C9orf72</i> <i>FTD</i> / <i>ALS</i> pathology <i>in vitro</i> and <i>in vivo</i> . <i>EMBO Molecular Medicine</i> , 2018, 10, 22-31.	6.9	178
23	Quadruplex nucleic acids as targets for anticancer therapeutics. <i>Nature Reviews Chemistry</i> , 2017, 1, .	30.2	357
24	Exploring the Dynamics of Propeller Loops in Human Telomeric DNA Quadruplexes Using Atomistic Simulations. <i>Journal of Chemical Theory and Computation</i> , 2017, 13, 2458-2480.	5.3	39
25	Induced-Fit Recognition of CCG Trinucleotide Repeats by a Nickel-Chromomycin Complex Resulting in Large-Scale DNA Deformation. <i>Angewandte Chemie - International Edition</i> , 2017, 56, 8761-8765.	13.8	30
26	Folding of guanine quadruplex molecules: "funnel-like mechanism or kinetic partitioning? An overview from MD simulation studies. <i>Biochimica Et Biophysica Acta - General Subjects</i> , 2017, 1861, 1246-1263.	2.4	89
27	Induced-Fit Recognition of CCG Trinucleotide Repeats by a Nickel-Chromomycin Complex Resulting in Large-Scale DNA Deformation. <i>Angewandte Chemie</i> , 2017, 129, 8887-8891.	2.0	0
28	Targeting Promoter Quadruplex Nucleic Acids for Cancer Therapy. , 2017, , 308-340.		3
29	Toward the Development of Specific G-Quadruplex Binders: Synthesis, Biophysical, and Biological Studies of New Hydrazone Derivatives. <i>Journal of Medicinal Chemistry</i> , 2016, 59, 5706-5720.	6.4	51
30	Preface. <i>Bioorganic and Medicinal Chemistry</i> , 2016, 24, 4767.	3.0	0
31	Molecular mechanisms and therapeutic strategies in amyotrophic lateral sclerosis caused by <i>C9orf72</i> mutations. <i>Lancet, The</i> , 2016, 387, S13.	13.7	0
32	Quadruplex Nucleic Acids as Novel Therapeutic Targets. <i>Journal of Medicinal Chemistry</i> , 2016, 59, 5987-6011.	6.4	481
33	Can We Execute Reliable MM-PBSA Free Energy Computations of Relative Stabilities of Different Guanine Quadruplex Folds?. <i>Journal of Physical Chemistry B</i> , 2016, 120, 2899-2912.	2.6	32
34	Structural Insights into the Quadruplex-Duplex Interface Formed from a Telomeric Repeat: A Potential Molecular Target. <i>Journal of the American Chemical Society</i> , 2016, 138, 1226-1233.	13.7	56
35	A Personal History of Quadruplex Small Molecule Targeting. <i>Chemical Record</i> , 2015, 15, 691-710.	5.8	11
36	Flexibility and structural conservation in a c-KIT G-quadruplex. <i>Nucleic Acids Research</i> , 2015, 43, 629-644.	14.5	63

#	ARTICLE	IF	CITATIONS
37	Loop flexibility in human telomeric quadruplex small-molecule complexes. <i>Nucleic Acids Research</i> , 2015, 43, 4785-4799.	14.5	42
38	The discovery of a novel antibiotic for the treatment of <i>Clostridium difficile</i> infections: a story of an effective academic–industrial partnership. <i>MedChemComm</i> , 2015, 6, 1420-1426.	3.4	22
39	A G-quadruplex-binding compound showing anti-tumour activity in an in vivo model for pancreatic cancer. <i>Scientific Reports</i> , 2015, 5, 11385.	3.3	95
40	Synthesis and biological evaluation of hybrid acridine-HSP90 ligand conjugates as telomerase inhibitors. <i>Organic and Biomolecular Chemistry</i> , 2015, 13, 8500-8504.	2.8	12
41	Atomic Force Microscopy and Voltammetric Investigation of Quadruplex Formation between a Triazole-Acridine Conjugate and Guanine-Containing Repeat DNA Sequences. <i>Analytical Chemistry</i> , 2015, 87, 6141-6149.	6.5	15
42	KRAS oncogene repression in colon cancer cell lines by G-quadruplex binding indolo[3,2-c]quinolines. <i>Scientific Reports</i> , 2015, 5, 9696.	3.3	74
43	G-quadruplexes: Emerging roles in neurodegenerative diseases and the non-coding transcriptome. <i>FEBS Letters</i> , 2015, 589, 1653-1668.	2.8	185
44	Macrocyclic naphthalene diimides as G-quadruplex binders. <i>Bioorganic and Medicinal Chemistry</i> , 2015, 23, 3819-3830.	3.0	34
45	Indolo[3,2-a]quinoline G-quadruplex Stabilizers: a Structural Analysis of Binding to the Human Telomeric G-quadruplex. <i>ChemMedChem</i> , 2015, 10, 836-849.	3.2	24
46	Extended molecular dynamics of a c-kit promoter quadruplex. <i>Nucleic Acids Research</i> , 2015, 43, 8673-8693.	14.5	49
47	Triazole-linked phenyl derivatives: Redox mechanisms and in situ electrochemical evaluation of interaction with dsDNA. <i>Bioelectrochemistry</i> , 2015, 101, 97-105.	4.6	2
48	Targeting KRAS Oncogene in Colon Cancer Cells with 7-Carboxylate Indolo[3,2-b]quinoline Tri-Alkylamine Derivatives. <i>PLoS ONE</i> , 2015, 10, e0126891.	2.5	41
49	Structure-Dependent Binding of Arylimidamides to the DNA Minor Groove. <i>ChemBioChem</i> , 2014, 15, 68-79.	2.6	20
50	Discovery of new G-quadruplex binding chemotypes. <i>Chemical Communications</i> , 2014, 50, 960-963.	4.1	24
51	Small-molecule quadruplex-targeted drug discovery. <i>Bioorganic and Medicinal Chemistry Letters</i> , 2014, 24, 2602-2612.	2.2	165
52	Targeting a c-MYC G-quadruplex DNA with a fragment library. <i>Chemical Communications</i> , 2014, 50, 1704-1707.	4.1	49
53	Preface. <i>Bioorganic and Medicinal Chemistry</i> , 2014, 22, 4355.	3.0	0
54	Structural Basis for the Identification of an i-Motif Tetraplex Core with a Parallel-Duplex Junction as a Structural Motif in CCG Triplet Repeats. <i>Angewandte Chemie - International Edition</i> , 2014, 53, 10682-10686.	13.8	30

#	ARTICLE	IF	CITATIONS
55	Small-molecule G-quadruplex interactions: Systematic exploration of conformational space using multiple molecular dynamics. <i>Biopolymers</i> , 2013, 99, n/a-n/a.	2.4	29
56	Structure-Based Design and Evaluation of Naphthalene Diimide G-Quadruplex Ligands As Telomere Targeting Agents in Pancreatic Cancer Cells. <i>Journal of Medicinal Chemistry</i> , 2013, 56, 2959-2974.	6.4	163
57	Synthesis, G-Quadruplex Stabilisation, Docking Studies, and Effect on Cancer Cells of Indolo[3,2-a]quinolines with One, Two, or Three Basic Side Chains. <i>ChemMedChem</i> , 2013, 8, 1648-1661.	3.2	39
58	The influence of positional isomerism on G-quadruplex binding and anti-proliferative activity of tetra-substituted naphthalene diimide compounds. <i>Bioorganic and Medicinal Chemistry</i> , 2013, 21, 6162-6170.	3.0	17
59	A new plant-derived antibacterial is an inhibitor of efflux pumps in <i>Staphylococcus aureus</i> . <i>International Journal of Antimicrobial Agents</i> , 2013, 42, 513-518.	2.5	62
60	Antibacterial activity of head-to-head bis-benzimidazoles. <i>International Journal of Antimicrobial Agents</i> , 2013, 42, 361-366.	2.5	27
61	Mechanism of the Antiproliferative Activity of Some Naphthalene Diimide G-Quadruplex Ligands. <i>Molecular Pharmacology</i> , 2013, 83, 470-480.	2.3	29
62	Small-molecule Binding to the DNA Minor Groove Is Mediated by a Conserved Water Cluster. <i>Journal of the American Chemical Society</i> , 2013, 135, 1369-1377.	13.7	68
63	Optimization of anti-proliferative activity using a screening approach with a series of bis-heterocyclic G-quadruplex ligands. <i>Bioorganic and Medicinal Chemistry Letters</i> , 2013, 23, 5351-5355.	2.2	7
64	Observation of unphosphorylated STAT3 core protein binding to target DNA by PEMSA and X-ray crystallography. <i>FEBS Letters</i> , 2013, 587, 833-839.	2.8	60
65	Downregulation of Androgen Receptor Transcription by Promoter G-Quadruplex Stabilization as a Potential Alternative Treatment for Castrate-Resistant Prostate Cancer. <i>Biochemistry</i> , 2013, 52, 1429-1436.	2.5	23
66	Crystal Structure of a Promoter Sequence in the B-raf Gene Reveals an Intertwined Dimer Quadruplex. <i>Journal of the American Chemical Society</i> , 2013, 135, 19319-19329.	13.7	45
67	Conformational dynamics of the human propeller telomeric DNA quadruplex on a microsecond time scale. <i>Nucleic Acids Research</i> , 2013, 41, 2723-2735.	14.5	70
68	Inhibition of the hypoxia-inducible factor pathway by a G-quadruplex binding small molecule. <i>Scientific Reports</i> , 2013, 3, 2799.	3.3	35
69	Thioester derivatives of the natural product psammaplin A as potent histone deacetylase inhibitors. <i>Beilstein Journal of Organic Chemistry</i> , 2013, 9, 81-88.	2.2	28
70	Xanthene and Xanthone Derivatives as G-Quadruplex Stabilizing Ligands. <i>Molecules</i> , 2013, 18, 13446-13470.	3.8	14
71	Developing and paying for medicines for orphan indications in oncology: utilitarian regulation vs equitable care?. <i>British Journal of Cancer</i> , 2012, 106, 14-17.	6.4	23
72	Reply: Comment on "Developing and paying for medicines for orphan indications in oncology: utilitarian regulation vs equitable care?". <i>British Journal of Cancer</i> , 2012, 107, 584-584.	6.4	0

#	ARTICLE	IF	CITATIONS
73	Defining the Mechanism of Action and Enzymatic Selectivity of Psammaplin A against Its Epigenetic Targets. <i>Journal of Medicinal Chemistry</i> , 2012, 55, 1731-1750.	6.4	89
74	Molecular Basis of Structure–Activity Relationships between Salphen Metal Complexes and Human Telomeric DNA Quadruplexes. <i>Journal of Medicinal Chemistry</i> , 2012, 55, 209-222.	6.4	196
75	Crystal structure of a c-kit promoter quadruplex reveals the structural role of metal ions and water molecules in maintaining loop conformation. <i>Nucleic Acids Research</i> , 2012, 40, 4691-4700.	14.5	117
76	A novel series of G-quadruplex ligands with selectivity for HIF-expressing osteosarcoma and renal cancer cell lines. <i>Bioorganic and Medicinal Chemistry Letters</i> , 2012, 22, 5984-5988.	2.2	30
77	Crystallography of DNA and RNA G-Quadruplex Nucleic Acids and Their Ligand Complexes. <i>Current Protocols in Nucleic Acid Chemistry</i> , 2012, 50, Unit17.6.	0.5	36
78	Into the minor groove. <i>Nature Chemistry</i> , 2012, 4, 594-595.	13.6	21
79	Molecular Dynamics and Force Field Based Methods for Studying Quadruplex Nucleic Acids. <i>RSC Biomolecular Sciences</i> , 2012, , 33-52.	0.4	4
80	Sequences in the HSP90 promoter form G-quadruplex structures with selectivity for disubstituted phenyl bis-oxazole derivatives. <i>Bioorganic and Medicinal Chemistry Letters</i> , 2012, 22, 5930-5935.	2.2	26
81	Bioactive Compounds from <i>Carissa spinarum</i> . <i>Phytotherapy Research</i> , 2012, 26, 1496-1499.	5.8	32
82	Molecular Dynamics Studies of the STAT3 Homodimer:DNA Complex: Relationships between STAT3 Mutations and Protein–DNA Recognition. <i>Journal of Chemical Information and Modeling</i> , 2012, 52, 1179-1192.	5.4	21
83	Symmetric Bis-benzimidazoles Are Potent Anti-Staphylococcal Agents with Dual Inhibitory Mechanisms against DNA Gyrase. <i>Biochemistry</i> , 2012, 51, 5860-5871.	2.5	26
84	Structural Basis for Telomeric G-Quadruplex Targeting by Naphthalene Diimide Ligands. <i>Journal of the American Chemical Society</i> , 2012, 134, 2723-2731.	13.7	213
85	Synthesis of Small Molecules Targeting Multiple DNA Structures using Click Chemistry. <i>ChemMedChem</i> , 2012, 7, 792-804.	3.2	21
86	Identification of novel telomeric G-quadruplex-targeting chemical scaffolds through screening of three NCI libraries. <i>Bioorganic and Medicinal Chemistry Letters</i> , 2012, 22, 3006-3010.	2.2	29
87	Structural Basis of Telomeric RNA Quadruplex–Acridine Ligand Recognition. <i>Journal of the American Chemical Society</i> , 2011, 133, 2721-2728.	13.7	125
88	Water-Mediated Binding of Agents that Target the DNA Minor Groove. <i>Journal of the American Chemical Society</i> , 2011, 133, 10171-10183.	13.7	60
89	A structural analysis of G-quadruplex/ligand interactions. <i>Biochimie</i> , 2011, 93, 1239-1251.	2.6	123
90	Surface area accessibility and the preferred topology of telomeric DNA quadruplex–ligand complexes. <i>Biochimie</i> , 2011, 93, 1275-1279.	2.6	13

#	ARTICLE	IF	CITATIONS
91	The triazatruxene derivative azatrux binds to the parallel form of the human telomeric G-quadruplex under molecular crowding conditions: Biophysical and molecular modeling studies. <i>Biochimie</i> , 2011, 93, 1318-1327.	2.6	37
92	N-Cyclic Bay-Substituted Perylene G-Quadruplex Ligands Have Selective Antiproliferative Effects on Cancer Cells and Induce Telomere Damage. <i>Journal of Medicinal Chemistry</i> , 2011, 54, 1140-1156.	6.4	51
93	Fluorine in medicinal chemistry: $\hat{2}$ -fluorination of peripheral pyrrolidines attached to acridine ligands affects their interactions with G-quadruplex DNA. <i>Organic and Biomolecular Chemistry</i> , 2011, 9, 1328.	2.8	65
94	Targeting G-quadruplexes in gene promoters: a novel anticancer strategy?. <i>Nature Reviews Drug Discovery</i> , 2011, 10, 261-275.	46.4	1,447
95	Targeting pancreatic cancer with a G-quadruplex ligand. <i>Bioorganic and Medicinal Chemistry</i> , 2011, 19, 7151-7157.	3.0	58
96	On the function of the internal cavity of histone deacetylase protein 8: R37 is a crucial residue for catalysis. <i>Bioorganic and Medicinal Chemistry Letters</i> , 2011, 21, 2129-2132.	2.2	36
97	Mapping the sequences of potential guanine quadruplex motifs. <i>Nucleic Acids Research</i> , 2011, 39, 4917-4927.	14.5	29
98	Rational Design of Acridine-Based Ligands with Selectivity for Human Telomeric Quadruplexes. <i>Journal of the American Chemical Society</i> , 2010, 132, 12263-12272.	13.7	98
99	Tetrasubstituted naphthalene diimide ligands with selectivity for telomeric G-quadruplexes and cancer cells. <i>Bioorganic and Medicinal Chemistry Letters</i> , 2010, 20, 6459-6463.	2.2	93
100	C-11 diamino cryptolepine derivatives NSC748392, NSC748393, and NSC748394: Anticancer profile and G-quadruplex stabilization. <i>Bioorganic and Medicinal Chemistry Letters</i> , 2010, 20, 7042-7045.	2.2	26
101	A novel small-molecule inhibitor of IL-6 signalling. <i>Bioorganic and Medicinal Chemistry Letters</i> , 2010, 20, 7029-7032.	2.2	16
102	Human telomeric G-quadruplex: The current status of telomeric G-quadruplexes as therapeutic targets in human cancer. <i>FEBS Journal</i> , 2010, 277, 1118-1125.	4.7	481
103	A crystallographic and modelling study of a human telomeric RNA (TERRA) quadruplex. <i>Nucleic Acids Research</i> , 2010, 38, 5569-5580.	14.5	213
104	Electrospray Mass Spectrometry of Telomeric RNA (TERRA) Reveals the Formation of Stable Multimeric G-Quadruplex Structures. <i>Journal of the American Chemical Society</i> , 2010, 132, 9328-9334.	13.7	124
105	Targeting the c-Myc Promoter G-quadruplexes with 6-Substituted Indenoisoquinolines. <i>ACS Medicinal Chemistry Letters</i> , 2010, 1, 306-310.	2.8	67
106	Molecular Modeling on Inhibitor Complexes and Active-Site Dynamics of Cytochrome P450 C17, a Target for Prostate Cancer Therapy. <i>Journal of Molecular Biology</i> , 2010, 400, 1078-1098.	4.2	25
107	A click chemistry approach to C3 symmetric, G-quadruplex stabilising ligands. <i>Organic and Biomolecular Chemistry</i> , 2010, 8, 2926.	2.8	28
108	Structure-Activity Relationships of Monomeric C2-Aryl Pyrrolo[2,1- <i>b</i> ][1,4]benzodiazepine (PBD) Antitumor Agents. <i>Journal of Medicinal Chemistry</i> , 2010, 53, 2927-2941.	6.4	39

#	ARTICLE	IF	CITATIONS
109	The structures of quadruplex nucleic acids and their drug complexes. <i>Current Opinion in Structural Biology</i> , 2009, 19, 239-250.	5.7	407
110	G-quadruplex compounds and cis-platin act synergistically to inhibit cancer cell growth in vitro and in vivo. <i>Biochemical Pharmacology</i> , 2009, 78, 115-122.	4.4	34
111	Design, synthesis and evaluation of 4,5-di-substituted acridone ligands with high G-quadruplex affinity and selectivity, together with low toxicity to normal cells. <i>Bioorganic and Medicinal Chemistry Letters</i> , 2009, 19, 5109-5113.	2.2	33
112	G-quadruplex nucleic acids as therapeutic targets. <i>Current Opinion in Chemical Biology</i> , 2009, 13, 345-353.	6.1	532
113	Selective G-quadruplex ligands: The significant role of side chain charge density in a series of perylene derivatives. <i>Bioorganic and Medicinal Chemistry Letters</i> , 2009, 19, 3903-3908.	2.2	24
114	A G-Rich Sequence within the <i>c-kit</i> Oncogene Promoter Forms a Parallel G-Quadruplex Having Asymmetric G-Tetrad Dynamics. <i>Journal of the American Chemical Society</i> , 2009, 131, 13399-13409.	13.7	195
115	Targeting Human Gastrointestinal Stromal Tumor Cells with a Quadruplex-Binding Small Molecule. <i>Journal of Medicinal Chemistry</i> , 2009, 52, 3774-3783.	6.4	126
116	Shedding Light on the Interaction between TMPyP4 and Human Telomeric Quadruplexes. <i>Journal of Physical Chemistry B</i> , 2009, 113, 14779-14786.	2.6	145
117	Bioactive Pyridine- <i>N</i> -oxide Disulfides from <i>Allium stipitatum</i> . <i>Journal of Natural Products</i> , 2009, 72, 360-365.	3.0	103
118	A Role for Water Molecules in DNA-Ligand Minor Groove Recognition. <i>Accounts of Chemical Research</i> , 2009, 42, 11-21.	15.6	119
119	Selectivity in Ligand Recognition of G-Quadruplex Loops. <i>Biochemistry</i> , 2009, 48, 1675-1680.	2.5	114
120	Recognition and discrimination of DNA quadruplexes by acridine-peptide conjugates. <i>Organic and Biomolecular Chemistry</i> , 2009, 7, 76-84.	2.8	60
121	A molecular model for drug binding to tandem repeats of telomeric G-quadruplexes. <i>Biochemical Society Transactions</i> , 2009, 37, 583-588.	3.4	30
122	Amide bond direction modulates G-quadruplex recognition and telomerase inhibition by 2,6 and 2,7 bis-substituted anthracenedione derivatives. <i>Bioorganic and Medicinal Chemistry</i> , 2008, 16, 354-361.	3.0	31
123	Tri- and tetra-substituted naphthalene diimides as potent G-quadruplex ligands. <i>Bioorganic and Medicinal Chemistry Letters</i> , 2008, 18, 1668-1673.	2.2	128
124	TRAP-LIG, a modified telomere repeat amplification protocol assay to quantitate telomerase inhibition by small molecules. <i>Analytical Biochemistry</i> , 2008, 380, 99-105.	2.4	101
125	Effects of Metal Coordination Geometry on Stabilization of Human Telomeric Quadruplex DNA by Square-Planar and Square-Pyramidal Metal Complexes. <i>Inorganic Chemistry</i> , 2008, 47, 11910-11919.	4.0	126
126	Quadruplex DNA crystal structures and drug design. <i>Biochimie</i> , 2008, 90, 1184-1196.	2.6	147



#	ARTICLE	IF	CITATIONS
127	Rational Design of Substituted Diarylureas: A Scaffold for Binding to G-Quadruplex Motifs. <i>Journal of Medicinal Chemistry</i> , 2008, 51, 7751-7767.	6.4	61
128	Aminoacyl- $\gamma$ -Anthraquinone Conjugates as Telomerase Inhibitors: Synthesis, Biophysical and Biological Evaluation. <i>Journal of Medicinal Chemistry</i> , 2008, 51, 5566-5574.	6.4	58
129	Topology Conservation and Loop Flexibility in Quadruplex-Drug Recognition: Crystal Structures of Inter- and Intramolecular Telomeric DNA Quadruplex-Drug Complexes. <i>Journal of Molecular Biology</i> , 2008, 381, 1145-1156.	4.2	165
130	Molecular Dynamics and Principal Components Analysis of Human Telomeric Quadruplex Multimers. <i>Biophysical Journal</i> , 2008, 95, 296-311.	0.5	189
131	Structural Basis of DNA Quadruplex Recognition by an Acridine Drug. <i>Journal of the American Chemical Society</i> , 2008, 130, 6722-6724.	13.7	295
132	The relationship of potential G-quadruplex sequences in cis-upstream regions of the human genome to SP1-binding elements. <i>Nucleic Acids Research</i> , 2008, 36, 2700-2704.	14.5	66
133	Targeting telomerase and telomeres: a click chemistry approach towards highly selective G-quadruplex ligands. <i>Molecular BioSystems</i> , 2008, 4, 629.	2.9	36
134	The Building-Blocks of DNA and RNA. , 2008, , 20-37.		11
135	DNA Structure as Observed in Fibers and Crystals. , 2008, , 38-80.		8
136	High-resolution crystal structure of the intramolecular d(TpA) thymine-adenine photoadduct and its mechanistic implications. <i>Nucleic Acids Research</i> , 2007, 35, 1048-1053.	14.5	25
137	Observation of the Coexistence of Sodium and Calcium Ions in a DNA G-Quadruplex Ion Channel. <i>Journal of the American Chemical Society</i> , 2007, 129, 10106-10107.	13.7	67
138	Induced Fit Conformational Changes of a $\alpha$ -Reversed Amidine-Heterocycle: Optimized Interactions in a DNA Minor Groove Complex. <i>Journal of the American Chemical Society</i> , 2007, 129, 5688-5698.	13.7	47
139	Structural Basis for Binding of Porphyrin to Human Telomeres,. <i>Biochemistry</i> , 2007, 46, 2390-2397.	2.5	303
140	Sequence occurrence and structural uniqueness of a G-quadruplex in the human c-kit promoter. <i>Nucleic Acids Research</i> , 2007, 35, 5799-5808.	14.5	132
141	Structure of an Unprecedented G-Quadruplex Scaffold in the Human c-kit Promoter. <i>Journal of the American Chemical Society</i> , 2007, 129, 4386-4392.	13.7	418
142	Chemical Variation of Natural-Product-Like Scaffolds: Design, Synthesis, and Biological Activity of Fused Bicyclic Acetal Derivatives. <i>Angewandte Chemie - International Edition</i> , 2007, 46, 2493-2496.	13.8	51
143	Structure-based design of benzylamino-acridine compounds as G-quadruplex DNA telomere targeting agents. <i>Bioorganic and Medicinal Chemistry Letters</i> , 2007, 17, 2293-2298.	2.2	65
144	Structure-specific recognition of quadruplex DNA by organic cations: Influence of shape, substituents and charge. <i>Biophysical Chemistry</i> , 2007, 126, 140-153.	2.8	182

#	ARTICLE	IF	CITATIONS
145	Mechanism of acridine-based telomerase inhibition and telomere shortening. <i>Biochemical Pharmacology</i> , 2007, 74, 679-689.	4.4	126
146	Trisubstituted Acridines as G-quadruplex Telomere Targeting Agents. Effects of Extensions of the 3,6- and 9-Side Chains on Quadruplex Binding, Telomerase Activity, and Cell Proliferation. <i>Journal of Medicinal Chemistry</i> , 2006, 49, 582-599.	6.4	192
147	Topology Variation and Loop Structural Homology in Crystal and Simulated Structures of a Bimolecular DNA Quadruplex. <i>Journal of the American Chemical Society</i> , 2006, 128, 5480-5487.	13.7	61
148	Stabilization of G-Quadruplex DNA by Highly Selective Ligands via Click Chemistry. <i>Journal of the American Chemical Society</i> , 2006, 128, 15972-15973.	13.7	212
149	Virtual Screening of DNA Minor Groove Binders. <i>Journal of Medicinal Chemistry</i> , 2006, 49, 4232-4238.	6.4	43
150	Stabilization of G-Quadruplex DNA and Inhibition of Telomerase Activity by Square-Planar Nickel(II) Complexes. <i>Journal of the American Chemical Society</i> , 2006, 128, 5992-5993.	13.7	314
151	Discovery of G-quadruplex stabilizing ligands through direct ELISA of a one-bead-one-compound library. <i>Organic and Biomolecular Chemistry</i> , 2006, 4, 4364.	2.8	15
152	New mustard-linked 2-aryl-bis-benzimidazoles with anti-proliferative activity. <i>Organic and Biomolecular Chemistry</i> , 2006, 4, 1305.	2.8	41
153	A Conserved Quadruplex Motif Located in a Transcription Activation Site of the Human c-kit Oncogene. <i>Biochemistry</i> , 2006, 45, 7854-7860.	2.5	370
154	Quadruplex DNA: sequence, topology and structure. <i>Nucleic Acids Research</i> , 2006, 34, 5402-5415.	14.5	2,049
155	Natural and synthetic G-quadruplex interactive berberine derivatives. <i>Bioorganic and Medicinal Chemistry Letters</i> , 2006, 16, 1707-1711.	2.2	202
156	Synthesis of distamycin A polyamides targeting G-quadruplex DNA. <i>Organic and Biomolecular Chemistry</i> , 2006, 4, 3479.	2.8	45
157	Targeting the DNA minor groove with fused ring dicationic compounds: Comparison of in silico screening and a high-resolution crystal structure. <i>Bioorganic and Medicinal Chemistry Letters</i> , 2006, 16, 15-19.	2.2	11
158	Predictive modelling of topology and loop variations in dimeric DNA quadruplex structures. <i>Nucleic Acids Research</i> , 2006, 34, 2117-2127.	14.5	62
159	Structural Bioinformatics in Cancer. , 2006, , 127-140.		0
160	Truncated azinomycin analogues intercalate into DNA. <i>Bioorganic and Medicinal Chemistry Letters</i> , 2005, 15, 653-656.	2.2	19
161	Chemical approaches to the discovery and development of cancer therapies. <i>Nature Reviews Cancer</i> , 2005, 5, 285-296.	28.4	205
162	Highly prevalent putative quadruplex sequence motifs in human DNA. <i>Nucleic Acids Research</i> , 2005, 33, 2901-2907.	14.5	872

#	ARTICLE	IF	CITATIONS
163	Putative DNA Quadruplex Formation within the Human c-kit Oncogene. <i>Journal of the American Chemical Society</i> , 2005, 127, 10584-10589.	13.7	526
164	Out-of-Shape DNA Minor Groove Binders: Induced Fit Interactions of Heterocyclic Dications with the DNA Minor Groove. <i>Biochemistry</i> , 2005, 44, 14701-14708.	2.5	59
165	The G-Quadruplex-Interactive Molecule BRACO-19 Inhibits Tumor Growth, Consistent with Telomere Targeting and Interference with Telomerase Function. <i>Cancer Research</i> , 2005, 65, 1489-1496.	0.9	517
166	The role of chemical sciences in 21st century cancer drug discovery. <i>Discovery Medicine</i> , 2005, 5, 450-4.	0.5	0
167	Synthesis, biophysical and biological evaluation of 3,6-bis-amidoacridines with extended 9-anilino substituents as potent G-quadruplex-binding telomerase inhibitors. <i>Bioorganic and Medicinal Chemistry Letters</i> , 2004, 14, 4347-4351.	2.2	64
168	Evaluation of by disubstituted acridone derivatives as telomerase inhibitors: the importance of G-quadruplex binding. <i>Bioorganic and Medicinal Chemistry Letters</i> , 2004, 14, 5845-5849.	2.2	74
169	Tetrapeptides induce selective recognition for G-quadruplexes when conjugated to a DNA-binding platform. <i>Organic and Biomolecular Chemistry</i> , 2004, 2, 2925.	2.8	36
170	Thiophene-Based Diamidine Forms a $\pi$ - $\pi$ Stacked $\pi$ -AT Binding Minor Groove Agent. <i>Journal of the American Chemical Society</i> , 2004, 126, 13659-13669.	13.7	82
171	Synthesis and evaluation of analogues of 10H-indolo[3,2-b]quinoline as G-quadruplex stabilising ligands and potential inhibitors of the enzyme telomerase. <i>Organic and Biomolecular Chemistry</i> , 2004, 2, 981.	2.8	106
172	Loop-Length-Dependent Folding of G-Quadruplexes. <i>Journal of the American Chemical Society</i> , 2004, 126, 16405-16415.	13.7	428
173	Characterization of a Novel DNA Minor-Groove Complex. <i>Biophysical Journal</i> , 2004, 86, 1028-1041.	0.5	88
174	A G-quadruplex telomere targeting agent produces p16-associated senescence and chromosomal fusions in human prostate cancer cells. <i>Molecular Cancer Therapeutics</i> , 2004, 3, 1201-6.	4.1	112
175	The structure of telomeric DNA. <i>Current Opinion in Structural Biology</i> , 2003, 13, 275-283.	5.7	420
176	DNA sequence specificity of triplex-binding ligands. <i>FEBS Journal</i> , 2003, 270, 4982-4992.	0.2	19
177	Studies on the Nitroreductase Prodrug-Activating System. Crystal Structures of Complexes with the Inhibitor Dicoumarol and Dinitrobenzamide Prodrugs and of the Enzyme Active Form. <i>Journal of Medicinal Chemistry</i> , 2003, 46, 4009-4020.	6.4	107
178	Trisubstituted Acridine Derivatives as Potent and Selective Telomerase Inhibitors. <i>Journal of Medicinal Chemistry</i> , 2003, 46, 4463-4476.	6.4	173
179	Sequence-specific minor groove binding by bis-benzimidazoles: water molecules in ligand recognition. <i>Nucleic Acids Research</i> , 2003, 31, 1514-1524.	14.5	59
180	Structure of a G-quadruplex-Ligand Complex. <i>Journal of Molecular Biology</i> , 2003, 326, 117-125.	4.2	258

#	ARTICLE	IF	CITATIONS
181	Mechanistic and anti-proliferative studies of two novel, biologically active bis-benzimidazoles. <i>European Journal of Cancer</i> , 2003, 39, 2548-2555.	2.8	41
182	Structure of the First Parallel DNA Quadruplex-Drug Complex. <i>Journal of the American Chemical Society</i> , 2003, 125, 4066-4067.	13.7	212
183	Acquired Cellular Resistance to Flavopiridol in a Human Colon Carcinoma Cell Line Involves Up-Regulation of the Telomerase Catalytic Subunit and Telomere Elongation. Sensitivity of Resistant Cells to Combination Treatment with a Telomerase Inhibitor. <i>Molecular Pharmacology</i> , 2003, 64, 1101-1108.	2.3	37
184	The Design of G-quadruplex Ligands as Telomerase Inhibitors. <i>Mini-Reviews in Medicinal Chemistry</i> , 2003, 3, 11-21.	2.4	107
185	Fundamentals of DNA and RNA structure. <i>Methods of Biochemical Analysis</i> , 2003, 44, 41-73.	0.2	0
186	Telomerase inhibitors in cancer therapy: current status and future directions. <i>Current Opinion in Investigational Drugs</i> , 2003, 4, 675-85.	2.3	8
187	Protein and drug interactions in the minor groove of DNA. <i>Nucleic Acids Research</i> , 2002, 30, 1182-1191.	14.5	68
188	Strong Binding in the DNA Minor Groove by an Aromatic Diamidine with a Shape That Does Not Match the Curvature of the Groove. <i>Journal of the American Chemical Society</i> , 2002, 124, 13680-13681.	13.7	71
189	Solid-Phase Synthesis of Symmetrical 3,6-Bispeptide- $\alpha$ -Acridone Conjugates. <i>Organic Letters</i> , 2002, 4, 2509-2512.	4.6	15
190	Crystal Structure of the Potassium Form of an <i>Oxytricha nova</i> G-quadruplex. <i>Journal of Molecular Biology</i> , 2002, 320, 189-200.	4.2	317
191	A G-Quadruplex-Interactive Potent Small-Molecule Inhibitor of Telomerase Exhibiting in Vitro and in Vivo Antitumor Activity. <i>Molecular Pharmacology</i> , 2002, 61, 1154-1162.	2.3	280
192	Synthesis of a novel dimeric bis-benzimidazole with site-selective DNA-binding properties. <i>Tetrahedron Letters</i> , 2002, 43, 7239-7241.	1.4	15
193	Crystal structure of parallel quadruplexes from human telomeric DNA. <i>Nature</i> , 2002, 417, 876-880.	27.8	1,889
194	Telomere maintenance as a target for anticancer drug discovery. <i>Nature Reviews Drug Discovery</i> , 2002, 1, 383-393.	46.4	577
195	A New Class of Symmetric Bisbenzimidazole-Based DNA Minor Groove-Binding Agents Showing Antitumor Activity. <i>Journal of Medicinal Chemistry</i> , 2001, 44, 138-144.	6.4	234
196	A standard reference frame for the description of nucleic acid base-pair geometry 1 Edited by P. E. Wright 2 This is a document of the Nomenclature Committee of IUBMB (NC-IUBMB)/IUPAC-IUBMB Joint Commission on Biochemical Nomenclature (JCBN), whose members are R. Cammack (chairman), A. Bairoch, H.M. Berman, S. Boyce, C.R. Cantor, K. Elliott, D. Horton, M. Kanehisa, A. Kotyk, G.P. Moss, N. Sharon and K.F. Tipton.. <i>Journal of Molecular Biology</i> , 2001, 313, 229-237.	4.2	533
197	Structure-based design of selective and potent G quadruplex-mediated telomerase inhibitors. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2001, 98, 4844-4849.	7.1	455
198	DNA minor-groove recognition by small molecules (up to 2000). <i>Natural Product Reports</i> , 2001, 18, 291-309.	10.3	455

#	ARTICLE	IF	CITATIONS
199	Stabilisation of TG- and AG-containing antiparallel DNA triplexes by triplex-binding ligands. <i>Nucleic Acids Research</i> , 2001, 29, 1935-1942.	14.5	27
200	A novel inhibitor of human telomerase derived from 10H-indolo[3,2-b]quinoline. <i>Bioorganic and Medicinal Chemistry Letters</i> , 2000, 10, 2063-2066.	2.2	111
201	DNA minor groove interactions and the biological activity of 2,5-bis-[4-(N-alkylamidino)phenyl] furans. <i>Bioorganic and Medicinal Chemistry Letters</i> , 2000, 10, 2593-2597.	2.2	33
202	A DNA-porphyrin minor-groove complex at atomic resolution: The structural consequences of porphyrin ruffling. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2000, 97, 9476-9481.	7.1	68
203	Stability of DNA Triplexes on Shuttle Vector Plasmids in the Replication Pool in Mammalian Cells. <i>Journal of Biological Chemistry</i> , 2000, 275, 39117-39124.	3.4	7
204	A thermodynamic and structural analysis of DNA minor-groove complex formation 1 Edited by I. Tinoco. <i>Journal of Molecular Biology</i> , 2000, 300, 321-337.	4.2	130
205	A new crystal form for the dodecamer C-G-C-G-A-A-T-T-C-G-C-G: symmetry effects on sequence-dependent DNA structure. <i>Journal of Molecular Biology</i> , 2000, 300, 551-561.	4.2	50
206	Structural Characterization of a Guanine <sup>+</sup> Quadruplex Ligand Complex. <i>Biochemistry</i> , 2000, 39, 13422-13432.	2.5	100
207	Crystal Structure of FMN-Dependent Nitroreductase from <i>Escherichia coli</i> B: A Prodrug-Activating Enzyme. <i>Journal of Medicinal Chemistry</i> , 2000, 43, 3624-3631.	6.4	119
208	DNA minor groove recognition of a non-self-complementary AT-rich sequence by a tris-benzimidazole ligand. <i>Nucleic Acids Research</i> , 1999, 27, 2691-2698.	14.5	35
209	Structure of the oligonucleotide d(CGTATATACG) as a site-specific complex with nickel ions. <i>Nucleic Acids Research</i> , 1999, 27, 1593-1599.	14.5	65
210	Human telomerase inhibition by substituted acridine derivatives. <i>Bioorganic and Medicinal Chemistry Letters</i> , 1999, 9, 2463-2468.	2.2	164
211	2,7-Disubstituted Amidofluorenone Derivatives as Inhibitors of Human Telomerase. <i>Journal of Medicinal Chemistry</i> , 1999, 42, 2679-2684.	6.4	212
212	Stabilization of DNA triple helices by a series of mono- and disubstituted amidoanthraquinones. <i>FEBS Journal</i> , 1999, 263, 817-825.	0.2	27
213	Symmetric bis-benzimidazoles: new sequence-selective DNA-binding molecules. <i>Chemical Communications</i> , 1999, , 929-930.	4.1	44
214	Molecular Modeling Studies on G-Quadruplex Complexes of Telomerase Inhibitors: Structure-Activity Relationships. <i>Journal of Medicinal Chemistry</i> , 1999, 42, 4538-4546.	6.4	144
215	Crystal Structure of Human DT-diaphorase: A Model for Interaction with the Cytotoxic Prodrug 5-(Aziridin-1-yl)-2,4-dinitrobenzamide (CB1954). <i>Journal of Medicinal Chemistry</i> , 1999, 42, 4325-4330.	6.4	50
216	Telomerase as an anti-cancer target: current status and future prospects. <i>Anti-cancer Drug Design</i> , 1999, 14, 341-7.	0.3	15

#	ARTICLE	IF	CITATIONS
217	Design, synthesis and evaluation of human telomerase inhibitors based upon a tetracyclic structural motif. <i>Anti-cancer Drug Design</i> , 1999, 14, 373-82.	0.3	1
218	Human Telomerase Inhibition by Regioisomeric Disubstituted Amidoanthracene-9,10-diones. <i>Journal of Medicinal Chemistry</i> , 1998, 41, 4873-4884.	6.4	167
219	New insights into sequence-dependent DNA structure. <i>Nature Structural Biology</i> , 1998, 5, 754-756.	9.7	18
220	Structure of the DNA Decamer d(GGCAATTGCG) Contains both Major- and Minor-Groove Binding GÂ·(GÂ·C) Base Triplets. <i>Acta Crystallographica Section D: Biological Crystallography</i> , 1998, 54, 577-583.	2.5	3
221	1,4- and 2,6-Disubstituted Amidoanthracene-9,10-dione Derivatives as Inhibitors of Human Telomerase. <i>Journal of Medicinal Chemistry</i> , 1998, 41, 3253-3260.	6.4	162
222	Anthracene-9,10-diones as Potential Anticancer Agents:Â Bacterial Mutation Studies of Amido-Substituted Derivatives Reveal an Unexpected Lack of Mutagenicity. <i>Journal of Medicinal Chemistry</i> , 1998, 41, 3748-3752.	6.4	33
223	Visualisation of extensive water ribbons and networks in a DNA minor-groove drug complex. <i>Nucleic Acids Research</i> , 1998, 26, 2873-2878.	14.5	27
224	An Integrated Sequence-Structure Database incorporating matching mRNA sequence, amino acid sequence and protein three-dimensional structure data. <i>Nucleic Acids Research</i> , 1998, 26, 327-331.	14.5	24
225	Structure of a Bis-Amidinium Derivative of Hoechst 33258 Complexed to Dodecanucleotide D(CGCGAATTCGCC)2: The Role of Hydrogen Bonding in Minor Groove Drug-DNA Recognition. <i>Nucleic Acids Research</i> , 1997, 25, 1510-1515.	14.5	34
226	Recognition of GC base pairs by triplex forming oligonucleotides containing nucleosides derived from 2-aminopyridine. <i>Nucleic Acids Research</i> , 1997, 25, 4891-4898.	14.5	48
227	Sequence-dependent crossed helix packing in the crystal structure of a B-DNA decamer yields a detailed model for the holliday junction 1 Edited by K. Nagai. <i>Journal of Molecular Biology</i> , 1997, 269, 827-841.	4.2	29
228	A model for the [C+ -Gâ·...C]n triple helix derived from observation of the C+ -Gâ·...C base triplet in a crystal structure. <i>FEBS Letters</i> , 1997, 416, 86-89.	2.8	11
229	Inhibition of Human Telomerase by a G-Quadruplex-Interactive Compound. <i>Journal of Medicinal Chemistry</i> , 1997, 40, 2113-2116.	6.4	763
230	Cytotoxicity of bis(phenylamidinium)furan alkyl derivatives in human tumour cell lines: Relation to DNA minor groove binding. <i>Bioorganic and Medicinal Chemistry Letters</i> , 1997, 7, 1403-1408.	2.2	36
231	Mercaptoacyl matrix metalloproteinase inhibitors: The effect of substitution at the mercaptoacyl moiety. <i>Bioorganic and Medicinal Chemistry Letters</i> , 1997, 7, 2765-2770.	2.2	21
232	Structure of the A-DNA Octamer d(GGCATGCC). <i>Acta Crystallographica Section D: Biological Crystallography</i> , 1997, 53, 269-273.	2.5	2
233	Conformations of the sugar-phosphate backbone in helical DNA crystal structures. , 1997, 42, 113-124.		198
234	Homology modelling of the enzyme P450 17 alpha-hydroxylase/17,20-lyase--a target for prostate cancer chemotherapy--from the crystal structure of P450BM-3. <i>Anti-cancer Drug Design</i> , 1997, 12, 113-23.	0.3	2

#	ARTICLE	IF	CITATIONS
235	Recent developments in triple-helix regulation of gene expression. <i>Anti-cancer Drug Design</i> , 1997, 12, 433-42.	0.3	30
236	Targeting the Minor Groove of DNA: Crystal Structures of Two Complexes between Furan Derivatives of Berenil and the DNA Dodecamer d(CGCGAATTCGCG) <sub>2</sub> . <i>Journal of Medicinal Chemistry</i> , 1996, 39, 4554-4562.	6.4	92
237	The High Resolution Crystal Structure of the DNA Decamer d(AGGCATGCCT). <i>Journal of Molecular Biology</i> , 1996, 256, 340-351.	4.2	36
238	Non-random usage of "degenerate" codons is related to protein three-dimensional structure. <i>FEBS Letters</i> , 1996, 399, 78-82.	2.8	88
239	Homologs of Idoxifene: Variation of Estrogen Receptor Binding and Calmodulin Antagonism with Chain Length. <i>Journal of Medicinal Chemistry</i> , 1996, 39, 999-1004.	6.4	19
240	Conformational flexibility and its potential influence on metabolic reactions for the anti-tumour compound a-ammine-bd-bis(butyrate)-cf-dichloro-e-cyclohexylamine platinum(IV). <i>Bioorganic and Medicinal Chemistry Letters</i> , 1996, 6, 421-426.	2.2	6
241	Synthesis of Sequence-Selective C8-Linked Pyrrolo[2,1-c][1,4]benzodiazepine DNA Interstrand Cross-Linking Agents. <i>Journal of Organic Chemistry</i> , 1996, 61, 8141-8147.	3.2	108
242	Designer DNA-Binding Drugs: The Crystal Structure of a Meta-Hydroxy Analogue of Hoechst 33258 Bound to d(CGCGAATTCGCG) <sub>2</sub> . <i>Nucleic Acids Research</i> , 1996, 24, 4882-4889.	14.5	58
243	Efficient triple helix formation by oligodeoxyribonucleotides containing alpha- or beta-2-amino-5-(2-deoxy-D-ribofuranosyl) pyridine residues. <i>Nucleic Acids Research</i> , 1996, 24, 4176-4184.	14.5	50
244	Equilibrium and Kinetic Measurements Reveal Rapidly Reversible Binding of Ras to Raf. <i>Journal of Biological Chemistry</i> , 1996, 271, 6713-6719.	3.4	52
245	Crystal structure of a DNA decamer showing a novel pseudo four-way helix-helix junction.. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 1995, 92, 10767-10771.	7.1	36
246	A molecular anchor for stabilizing triple-helical DNA.. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 1995, 92, 7887-7891.	7.1	75
247	Perturbations in DNA structure upon interaction with porphyrins revealed by chemical probes, DNA footprinting and molecular modelling. <i>Bioorganic and Medicinal Chemistry</i> , 1995, 3, 671-677.	3.0	25
248	Active-site conformation of 17-(3-pyridyl)androsta-5,16-dien-3 <sup>ol</sup> , a potent inhibitor of the P450 enzyme C17 $\alpha$ -hydroxylase/C17-20 lyase. <i>Bioorganic and Medicinal Chemistry Letters</i> , 1995, 5, 1125-1130.	2.2	11
249	Sequence-Dependent Drug Binding to the Minor Groove of DNA: Crystal Structure of the DNA Dodecamer d(CGCAAATTTGCG) <sub>2</sub> Complexed with Propamidine. <i>Journal of Medicinal Chemistry</i> , 1995, 38, 2317-2325.	6.4	68
250	Characteristics of triplex-directed photoadduct formation by psoralen-linked oligodeoxynucleotides. <i>Nucleic Acids Research</i> , 1995, 23, 4283-4289.	14.5	19
251	An approach to protein homology modelling based on an ensemble of NMR structures: application to the Sox-5 HMG-box protein. <i>Protein Engineering, Design and Selection</i> , 1995, 8, 615-625.	2.1	13
252	Analysis of van der Waals and Electrostatic Contributions in the Interactions of Minor Groove Binding Benzimidazoles with DNA. <i>Journal of the American Chemical Society</i> , 1995, 117, 4716-4717.	13.7	84

#	ARTICLE	IF	CITATIONS
253	Inhibition of aromatase expression by a psoralen-linked triplex-forming oligonucleotide targeted to a coding sequence. <i>FEBS Letters</i> , 1995, 372, 222-228.	2.8	20
254	Detection and kinetic studies of triplex formation by oligodeoxynucleotides using real-time biomolecular interaction analysis (BIA). <i>Nucleic Acids Research</i> , 1995, 23, 3627-3632.	14.5	75
255	Rationally Designed Analogs of Tamoxifen with Improved Calmodulin Antagonism. <i>Journal of Medicinal Chemistry</i> , 1995, 38, 241-248.	6.4	38
256	Variability in DNA minor groove width recognised by ligand binding: the crystal structure of a bis-benzimidazole compound bound to the DNA duplex d(CGCGAATTCGCG) <sub>2</sub> . <i>Nucleic Acids Research</i> , 1995, 23, 3678-3684.	14.5	44
257	Sequence-dependent effects in drug-DNA interaction: the crystal structure of Hoechst 33258 bound to the d(CGCAAATTTGCG) <sub>2</sub> duplex. <i>Nucleic Acids Research</i> , 1994, 22, 1607-1612.	14.5	113
258	Crystal Structure of gamma-Oxapentamidine Complexed with d(CGCGAATTCGCG) <sub>2</sub> . The Effects of Drug Structural Change on DNA Minor-Groove Recognition. <i>FEBS Journal</i> , 1994, 226, 953-961.	0.2	23
259	A homology-based molecular model of the proline-rich homeodomain protein Prh, from haematopoietic cells. <i>FEBS Letters</i> , 1994, 345, 93-98.	2.8	5
260	Antisense Nucleic Acids and Proteins: Fundamentals and Applications. <i>International Journal of Biological Macromolecules</i> , 1994, 16, 110.	7.5	0
261	Nucleic acid targeted drug design C. L. Propst and T. J. Perun (editors), Dekker, New York, 1992. Pages xiii + 619. \$165.00. ISBN 0-8247-8662-9.. <i>Talanta</i> , 1994, 41, 166-166.	5.5	0
262	Structure of a Covalent DNA Minor Groove Adduct with a Pyrrolobenzodiazepine Dimer: Evidence for Sequence-Specific Interstrand Crosslinking. <i>Journal of Medicinal Chemistry</i> , 1994, 37, 4529-4537.	6.4	87
263	"...the tyranny of the lattice...".. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 1994, 91, 3579-3583.	7.1	194
264	NMR and molecular modeling studies of the interaction of berenil and pentamidine with d(CGCAAATTTGCG) <sub>2</sub> . <i>FEBS Journal</i> , 1993, 213, 1175-1184.	0.2	63
265	Factors involved in guanine/cytosine (G/C) selectivity of DNA-binding drugs: a molecular modelling study of ligand interactions with a tyrT DNA sequence. <i>Journal of the Chemical Society, Faraday Transactions</i> , 1993, 89, 2651.	1.7	5
266	A detailed molecular model for human aromatase. <i>Journal of Steroid Biochemistry and Molecular Biology</i> , 1993, 44, 399-407.	2.5	90
267	Crystal structure of an oligonucleotide duplex containing G.C base pairs: influence of mispairing on DNA backbone conformation.. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 1993, 90, 804-808.	7.1	93
268	Solution and Solid State Structure of 2- $\beta$ -BIS-(O-Trityl)-3-Oximinouridine. <i>Nucleosides &amp; Nucleotides</i> , 1993, 12, 605-614.	0.5	2
269	DNA minor groove recognition properties of pentamidine and its analogs: a molecular modeling study. <i>Molecular Pharmacology</i> , 1993, 43, 982-8.	2.3	15
270	Conformation and Dynamics of Drug-DNA Intercalation. <i>Journal of Biomolecular Structure and Dynamics</i> , 1992, 10, 97-139.	3.5	33



#	ARTICLE	IF	CITATIONS
271	Prediction of the structure of the Y+Â·Râˆ™Â·R+type DNA triple helix by molecular modelling. <i>Nucleic Acids Research</i> , 1992, 20, 6535-6541.	14.5	35
272	Crystal structure of a pentamidine-oligonucleotide complex: implications for DNA-binding properties. <i>Biochemistry</i> , 1992, 31, 7104-7109.	2.5	177
273	Rational design of a highly efficient irreversible DNA interstrand cross-linking agent based on the pyrrolobenzodiazepine ring system. <i>Journal of the American Chemical Society</i> , 1992, 114, 4939-4941.	13.7	147
274	Structural properties of four isomeric C2'/C3' modified uridines. <i>Journal of the American Chemical Society</i> , 1992, 114, 2687-2696.	13.7	9
275	A molecular-modeling study of the interactions between the antiestrogen drug tamoxifen and several derivatives, and the calcium-binding protein calmodulin. <i>Journal of Medicinal Chemistry</i> , 1992, 35, 2753-2761.	6.4	23
276	Anthracene-9,10-diones as potential anticancer agents. Synthesis, DNA-binding, and biological studies on a series of 2,6-disubstituted derivatives. <i>Journal of Medicinal Chemistry</i> , 1992, 35, 1418-1429.	6.4	83
277	Minor-groove width and accessibility in B-DNA drug and protein complexes. <i>FEBS Letters</i> , 1992, 298, 97-99.	2.8	37
278	Molecular dynamics simulation of the DNA triplex d(TC)5 Â· d(GA)5 Â· d(C+T)5. <i>Journal of Molecular Biology</i> , 1992, 223, 519-529.	4.2	40
279	Crystal structure of a berenil-d(CGCAAATTTGCCG) complex. <i>Journal of Molecular Biology</i> , 1992, 226, 481-490.	4.2	103
280	Molecular structure of the B-DNA dodecamer d(CGCAAATTTGCCG) <sub>2</sub> An examination of propeller twist and minor-groove water structure at 2Â·2Âˆ resolution. <i>Journal of Molecular Biology</i> , 1992, 226, 1161-1173.	4.2	137
281	Derivatives of 2â€²,3â€²-dithiouridine and [1-Î²-D-(2,3-dithioxylofuranosyl)]uracil. <i>Tetrahedron Letters</i> , 1992, 33, 8151-8154.	1.4	8
282	Comparative structural studies of [3.1.0]-fused 2',3'-modified .beta.-D-nucleosides by x-ray crystallography, NMR spectroscopy, and molecular mechanics calculations. <i>Journal of Organic Chemistry</i> , 1991, 56, 6884-6892.	3.2	32
283	[22] Molecular modeling to study DNA intercalation by anti-tumor drugs. <i>Methods in Enzymology</i> , 1991, 203, 433-458.	1.0	13
284	Molecular Dynamics Simulations of Dinucleoside and Dinucleoside-Drug Crystal Hydrates. <i>Journal of Biomolecular Structure and Dynamics</i> , 1991, 9, 363-386.	3.5	9
285	Crystal structure and sequence-dependent conformation of the A.C mispaired oligonucleotide d(CGCAAGCTGGCG).. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 1990, 87, 6693-6697.	7.1	70
286	X-ray structural studies on three analogues of the Î±-anomer of the antitumour antibiotic showdomycin. Different ring-puckering effects of hydroxyl sugar substituents in lyxo and arabino configurations. <i>Acta Crystallographica Section B: Structural Science</i> , 1990, 46, 426-431.	1.8	0
287	Structure and conformational features of 9-(4-diethylaminophenyl)acridine. <i>Acta Crystallographica Section B: Structural Science</i> , 1990, 46, 860-862.	1.8	2
288	Crystal structure of a berenil-dodecanucleotide complex: the role of water in sequence-specific ligand binding.. <i>EMBO Journal</i> , 1990, 9, 1329-1334.	7.8	105

#	ARTICLE	IF	CITATIONS
289	Molecular-Mechanics Modelling of Drug-DNA Structures; the Effects of Differing Dielectric Treatment on Helix Parameters and Comparison with a Fully Solvated Structural Model. <i>Journal of Biomolecular Structure and Dynamics</i> , 1990, 8, 359-373.	3.5	55
290	Crystal and solution structures of the oligonucleotide d(ATGCGCAT) <sub>2</sub> : a combined X-ray and NMR study. <i>Nucleic Acids Research</i> , 1990, 18, 5521-5528.	14.5	38
291	Interaction of berenil with the tyrT DNA sequence studied by footprinting and molecular modelling. Implications for the design of sequence-specific DNA recognition agents. <i>Nucleic Acids Research</i> , 1990, 18, 4479-4488.	14.5	51
292	Inhibitors of the P450 enzymes aromatase and lyase. Crystallographic and molecular modeling studies suggest structural features of pyridylacetic acid derivatives responsible for differences in enzyme inhibitory activity. <i>Journal of Medicinal Chemistry</i> , 1990, 33, 3055-3060.	6.4	24
293	Crystallographic and molecular modeling studies on 3-ethyl-3-(4-pyridyl)piperidine-2,6-dione and its butyl analog, inhibitors of mammalian aromatase. Comparison with natural substrates: prediction of enantioselectivity for N-alkyl derivatives. <i>Journal of Medicinal Chemistry</i> , 1990, 33, 2673-2679.	6.4	21
294	Nucleic acids and molecular biology – Volume 3. <i>International Journal of Biological Macromolecules</i> , 1990, 12, 278-279.	7.5	0
295	Conformational effects of nucleotide exchange in ras p21 proteins as studied by fluorescence spectroscopy. <i>FEBS Letters</i> , 1990, 262, 127-130.	2.8	12
296	A molecular model for the enzyme cytochrome P45017 $\alpha$ , a major target for the chemotherapy of prostatic cancer. <i>Biochemical and Biophysical Research Communications</i> , 1990, 171, 1160-1167.	2.1	91
297	Crystal structure of a berenil-dodecanucleotide complex: the role of water in sequence-specific ligand binding. <i>EMBO Journal</i> , 1990, 9, 1329-34.	7.8	31
298	Structural studies on bio-active compounds. Part XIV. Molecular modelling of the interactions between pentamidine and DNA. <i>Anti-cancer Drug Design</i> , 1990, 5, 243-8.	0.3	9
299	The non-covalent interaction of pyrrolo[2, 1-c] [1, 4]benzodiazepine-5, 11-diones with DNA. <i>Anti-cancer Drug Design</i> , 1990, 5, 249-64.	0.3	12
300	Preliminary crystallographic data for NAD(P)H quinone reductase isolated from the Walker 256 rat carcinoma cell line. <i>Journal of Molecular Biology</i> , 1989, 205, 623-624.	4.2	6
301	The anti-conformation of 2',3'-dideoxy nucleosides may be essential for anti-HIV activity: Evidence from the crystal structure of 2',3'-dideoxy formycin A. <i>Biochemical and Biophysical Research Communications</i> , 1989, 161, 910-916.	2.1	12
302	Thermodynamic studies on the interactions of di-substituted anthraquinones with DNA. <i>Biochemical Pharmacology</i> , 1989, 38, 216-217.	4.4	4
303	Synthesis of the aromatase inhibitor 3-ethyl-3-(4-pyridyl)piperidine-2,6-dione and its enantiomers. <i>Journal of the Chemical Society Perkin Transactions 1</i> , 1989, , 196.	0.9	13
304	Molecular modelling of DNA-antitumour drug intercalation interactions: correlation of structural and energetic features with biological properties for a series of phenylquinoline-8-carboxamide compounds. <i>Molecular Pharmacology</i> , 1989, 35, 720-8.	2.3	2
305	Quinazoline antifolates inhibiting thymidylate synthase: computer modelling of the N10 substituent. <i>Anti-cancer Drug Design</i> , 1989, 3, 243-8.	0.3	4
306	Reaction between glycol benzyl ethers and thallium(III) nitrate. Synthesis of showdomycin analogues.. <i>Tetrahedron Letters</i> , 1988, 29, 1841-1844.	1.4	35

#	ARTICLE	IF	CITATIONS
307	Oxidative ring contraction of benzeneselenenate adducts of glycol ethers. synthesis of showdomycin analogues. <i>Tetrahedron Letters</i> , 1988, 29, 2711-2714.	1.4	32
308	A molecular model for proflavine-DNA intercalation. <i>Nucleic Acids Research</i> , 1988, 16, 8999-9016.	14.5	54
309	The influence of the ethyl group on the conformational flexibility of aminoglutethimide. <i>Biochemical Pharmacology</i> , 1988, 37, 143-145.	4.4	8
310	Binding of unfused aromatic cations to DNA. The influence of molecular twist on intercalation. <i>Journal of the American Chemical Society</i> , 1988, 110, 8292-8299.	13.7	52
311	Synthesis, molecular modeling, DNA binding, and antitumor properties of some substituted amidoanthraquinones. <i>Journal of Medicinal Chemistry</i> , 1988, 31, 847-857.	6.4	59
312	Glycol Ethers as Starting Materials in C-Nucleoside Synthesis. <i>Nucleosides &amp; Nucleotides</i> , 1988, 7, 609-612.	0.5	4
313	Experimental DNA-Binding and Computer Modelling Studies on an Analogue of the Anti-Tumour Drug Amsacrine. <i>Journal of Biomolecular Structure and Dynamics</i> , 1988, 6, 471-488.	3.5	12
314	Molecular modelling of the interactions of tetra-(4-N-methylpyridyl) porphin with TA and CG sites on DNA. <i>Nucleic Acids Research</i> , 1987, 15, 6553-6562.	14.5	59
315	The crystal structure of the DNA-binding drug berenil: molecular modelling studies of berenil-DNA complexes. <i>Nucleic Acids Research</i> , 1987, 15, 3469-3478.	14.5	46
316	DNA sequence preferences for an intercalating porphyrin compound revealed by footprinting. <i>Nucleic Acids Research</i> , 1987, 15, 2221-2234.	14.5	50
317	Crystallographic and Molecular Mechanics Calculations on the Anti-Tumor Drugs N-[(2-Dimethylamino)Ethyl]- and N-[(2-Dimethyl-Amino)Butyl]-9-Aminoacridine-4-Carboxamides and their Dications: Implications for Models of DNA-Binding. <i>Journal of Biomolecular Structure and Dynamics</i> , 1987, 5, 145-158.	3.5	13
318	The Influence of Intercalator Structure on DNA Binding Strength: The Importance of Side Chain Orientation. <i>Journal of Biomolecular Structure and Dynamics</i> , 1987, 5, 327-344.	3.5	3
319	Nucleic acid binding drugs. 18: Calculated charges for a DNA intercalation complex; a study of the relative influence of polarization and conformational effects. <i>International Journal of Biological Macromolecules</i> , 1987, 9, 193-196.	7.5	0
320	Conformational properties of 3-azido-2-deoxy-thymidine (AZT), an inhibitor of HIV reverse transcriptase. <i>Biochemical and Biophysical Research Communications</i> , 1987, 145, 1356-1361.	2.1	17
321	Podophyllotoxin analogues: synthesis and computer modelling investigation of structure-activity relationships. <i>Anti-cancer Drug Design</i> , 1987, 2, 247-52.	0.3	4
322	DNA sequence preferences for the anti-cancer drug mitoxantrone and related anthraquinones revealed by DNase I footprinting. <i>FEBS Letters</i> , 1986, 202, 289-294.	2.8	38
323	Origins of stereospecificity in DNA damage by anti-benzo[a]pyrene diol-epoxides. <i>FEBS Letters</i> , 1986, 209, 269-276.	2.8	11
324	Nucleic acid binding drugs—XIV. <i>Biochemical Pharmacology</i> , 1986, 35, 3915-3921.	4.4	11

#	ARTICLE	IF	CITATIONS
325	Structural aspects of drug-DNA complexes: molecular modelling of intercalative interactions. <i>Drugs Under Experimental and Clinical Research</i> , 1986, 12, 455-62.	0.3	1
326	Nucleic acid binding drugs. Part XIII. Molecular motion in a drug-nucleic acid model system: thermal motion analysis of a proflavine-dinucleoside crystal structure. <i>Nucleic Acids Research</i> , 1985, 13, 5671-5684.	14.5	7
327	Variability in the Geometry of RNA Double Helices Generated from Dinucleoside Building-Blocks. <i>Journal of Biomolecular Structure and Dynamics</i> , 1985, 2, 1235-1243.	3.5	0
328	Comparative computer graphics and solution studies of the DNA interaction of substituted anthraquinones based on doxorubicin and mitoxantrone. <i>Journal of Medicinal Chemistry</i> , 1985, 28, 857-864.	6.4	98
329	Structural studies on some tamoxifen derivatives. <i>Journal of Medicinal Chemistry</i> , 1985, 28, 1497-1503.	6.4	23
330	Studies on the conformation and dynamics of the C8-substituted guanine adduct of the carcinogen acetyaminofluorene; model for a possible Z-DNA modified structure. <i>Nucleic Acids Research</i> , 1984, 12, 8219-8233.	14.5	28
331	Nucleic acid binding drugs. X. A theoretical study of proflavine intercalation into RNA and DNA fragments: comparison with crystallographic results. <i>Acta Crystallographica Section B: Structural Science</i> , 1984, 40, 424-429.	1.8	9
332	Synthesis and biophysical studies of short oligodeoxynucleotides with novel modifications: a possible approach to the problem of mixed base oligodeoxynucleotide synthesis. <i>Nucleic Acids Research</i> , 1984, 12, 7435-7453.	14.5	135
333	Molecular models for the interaction of the anti-tumour drug nogalamycin with DNA. <i>Biochemical Pharmacology</i> , 1984, 33, 2877-2880.	4.4	43
334	Structural and Sequence-Dependent Aspects of Drug Intercalation Into Nucleic Acid. <i>Critical Reviews in Biochemistry</i> , 1984, 17, 73-121.	7.5	162
335	Computer graphics in the study of drug-nucleic acid interactions. <i>Biochemical Society Transactions</i> , 1984, 12, 1008-1011.	3.4	4
336	Nucleic acid structure: New twists to left-handed DNA. <i>Nature</i> , 1983, 302, 574-574.	27.8	13
337	The structure of the ribodinucleoside monophosphate guanylyl-3',5'-cytidine as its ammonium octahydrate salt. <i>Acta Crystallographica Section B: Structural Science</i> , 1983, 39, 98-104.	1.8	22
338	Nucleic acid binding drugs. VII. Molecular-mechanics studies on the conformational properties of the anti-cancer drug daunomycin: some observations on the use of differing potential-energy functions. <i>Acta Crystallographica Section B: Structural Science</i> , 1983, 39, 114-119.	1.8	18
339	The antitumor complex ethylenediamine platinum (II) malonate: X-ray structure analysis, and studies of its stability in solution. <i>Journal of Inorganic Biochemistry</i> , 1983, 18, 213-220.	3.5	27
340	Crystal and molecular structure of three isomers of dichlorodiamminedihydroxoplatinum(IV): cis-trans isomerization on recrystallization from water. <i>Inorganic Chemistry</i> , 1983, 22, 3620-3624.	4.0	87
341	Experimental and computer graphics simulation analyses of the dna interaction of 1,8-bis-(2-diethylaminoethylamino)-anthracene-9, 10-dione, a compound modelled on doxorubicin. <i>Biochemical Pharmacology</i> , 1983, 32, 2801-2808.	4.4	16
342	X-ray crystallographic studies of nucleic acids and nucleic acid-drug complexes. <i>Progress in Biophysics and Molecular Biology</i> , 1983, 41, 43-66.	2.9	53

#	ARTICLE	IF	CITATIONS
343	Monte Carlo Studies on Water in the dCpG/Proflavin Crystal Hydrate. <i>Journal of Biomolecular Structure and Dynamics</i> , 1983, 1, 287-297.	3.5	28
344	Computer Modelling Studies of the Covalent Interactions between DNA and the Enantiomers of anti-7,8-Diol,9,10-Epoxy-Benzo[a]pyrene. <i>Journal of Biomolecular Structure and Dynamics</i> , 1983, 1, 873-881.	3.5	16
345	X-ray crystallographic analysis of (±)-7,8,9,10-tetrahydrobenzo[a]pyrene: molecular structure of a diol epoxide. <i>Carcinogenesis</i> , 1983, 4, 415-418.	2.8	8
346	Molecular modelling studies on the non-covalent intercalative interactions between DNA and the enantiomers of anti-benzo[a]pyrene 7,8-diol-9,10-epoxide. <i>Carcinogenesis</i> , 1983, 4, 211-215.	2.8	20
347	X-Ray crystallographic analysis of 3-(2-phenyl-2-bithiazole-4-carboxamido)propyl dimethylsulphonium iodide, an analogue of the DNA-binding portion of bleomycin A2. <i>Nucleic Acids Research</i> , 1982, 10, 4753-4763.	14.5	18
348	Computer graphics in rational anti-cancer drug design. <i>Biochemical Society Transactions</i> , 1982, 10, 501-501.	3.4	1
349	The structure of a pseudo intercalated complex between actinomycin and the DNA binding sequence d(GpC). <i>Nature</i> , 1982, 296, 466-469.	27.8	117
350	Molecular structure of (+/-)-7,8,9,10-tetrahydroxy-7,8,9,10-tetrahydrobenzo(a)pyrene determined by x-ray crystallography. <i>Cancer Research</i> , 1982, 42, 3766-8.	0.9	10
351	New twists to DNA and DNA-carcinogen interactions. <i>Nature</i> , 1981, 292, 292-293.	27.8	5
352	The molecular structure of (±)-7,8-dihydroxy-7,8-dihydrobenzo[a]pyrene, an early metabolite of benzo[a]pyrene. <i>Carcinogenesis</i> , 1981, 2, 533-536.	2.8	10
353	Molecular structures of three hydroxyl derivatives of the hepatic carcinogen acetylaminofluorene. Evidence for a preferred orientation of the acetyl amino side-chain. <i>Carcinogenesis</i> , 1981, 2, 901-907.	2.8	7
354	The molecular structure of (±)-8,9,10,11-tetrahydroxy-8,9,10,11-tetrahydrobenz[a]anthracene; an X-ray crystallographic analysis. <i>Carcinogenesis</i> , 1981, 2, 445-449.	2.8	0
355	The structure of the antitumor complex cis-(diammino) (1,1-cyclobutanedicarboxylato)-Pt(II): X ray and nmr studies. <i>Journal of Inorganic Biochemistry</i> , 1980, 13, 205-212.	3.5	121
356	Highly structured water network in crystals of a deoxydinucleoside-drug complex. <i>Nature</i> , 1980, 288, 129-133.	27.8	134
357	Molecular structure of (±)-7,8,9,10-tetrahydrobenzo[a]pyrene; an X-ray crystallographic study. <i>Carcinogenesis</i> , 1980, 1, 249-254.	2.8	20
358	The structure of drug-deoxydinucleoside phosphate complex; generalized conformational behavior of intercalation complexes with RNA and DNA fragments. <i>Nucleic Acids Research</i> , 1980, 8, 85-98.	14.5	146
359	3 The Molecular Basis for the Action of Some DNA-Binding Drugs. <i>Progress in Medicinal Chemistry</i> , 1979, 16, 151-221.	10.4	117
360	Molecular and crystal structure of an intercalation complex: Proflavine-cytidylyl-(3',5')-guanosine. <i>Biopolymers</i> , 1979, 18, 2405-2429.	2.4	67

#	ARTICLE	IF	CITATIONS
361	Nucleic acid statics and dynamics. Nature, 1979, 279, 474-475.	27.8	3
362	Nucleic acid binding drugs. Some conformational properties of the anti-cancer drug daunomycin and several of its derivatives:. FEBS Letters, 1979, 107, 348-354.	2.8	25
363	Crystallisation of tRNACUGLeu from Escherichia coli after purification with hydroxyapatite. Biochemical and Biophysical Research Communications, 1979, 86, 66-70.	2.1	5
364	Carcinogen binding to DNA. Nature, 1978, 276, 444-445.	27.8	2
365	A 1:2 crystalline complex of ApA:proflavine: a model for binding to single-stranded reigons in RNA. Nucleic Acids Research, 1978, 5, 4417-4422.	14.5	45
366	Drug-nucleic acid interactions: conformational flexibility at the intercalation site.. Proceedings of the National Academy of Sciences of the United States of America, 1978, 75, 828-832.	7.1	65
367	Approaches to effective anti-cancer drugs. Nature, 1977, 268, 195-196.	27.8	28
368	Structure of a dinucleoside phosphateâ€“drug complex as model for nucleic acidâ€“drug interaction. Nature, 1977, 269, 304-307.	27.8	121
369	The crystal and molecular structure of an osmium bispyridine adduct of thymine. Nucleic Acids and Protein Synthesis, 1976, 418, 226-231.	1.7	34
370	Deeper insights into tRNA. Nature, 1976, 259, 267-268.	27.8	1
371	Drug action at the molecular level. Nature, 1976, 261, 14-15.	27.8	1
372	Crystal structure of proflavine, a DNA binding agent. Nature, 1975, 253, 284-285.	27.8	16
373	Drugs and nucleic acids. Nature, 1975, 255, 190-191.	27.8	2
374	Repressors at the molecular level. Nature, 1974, 249, 212-212.	27.8	0
375	Transfer RNA revisited. Nature, 1974, 250, 699-700.	27.8	0
376	On the determination of heavy-atom positions in various elastase derivatives. Acta Crystallographica Section B: Structural Crystallography and Crystal Chemistry, 1973, 29, 2645-2647.	0.4	7
377	Swazine, a new alkaloid from compton: a chemical and x-ray crystal study of a novel spiro dilactone. Tetrahedron Letters, 1972, 13, 707-710.	1.4	5
378	X-Ray determination of the structure and absolute configuration of a novel sesquiterpenoid, melampodin. Journal of the Chemical Society Chemical Communications, 1972, , 140.	2.0	25

#	ARTICLE	IF	CITATIONS
379	Assignment of Absolute Configuration from Anomalous Dispersion by Oxygen Atoms. <i>Nature</i> , 1970, 225, 376-376.	27.8	10
380	The crystal and molecular structure of streptomycin oxime selenate. <i>Tetrahedron Letters</i> , 1968, 9, 4725-4728.	1.4	34
381	Design and Analysis of G4 Recognition Compounds. , 0, , 337-359.		0