

# Glenn E M Maguire

## List of Publications by Year in descending order

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229  
papers

4,082  
citations

172457  
29  
h-index

138484  
58  
g-index

244  
all docs

244  
docs citations

244  
times ranked

4404  
citing authors

#	ARTICLE	IF	CITATIONS
1	Molecular fluorescent signalling with “fluor“spacer“receptor™ systems: approaches to sensing and switching devices via supramolecular photophysics. <i>Chemical Society Reviews</i> , 1992, 21, 187-195.	38.1	573
2	Fluorescent PET (photoinduced electron transfer) sensors. <i>Topics in Current Chemistry</i> , 1993, , 223-264.	4.0	369
3	The role of nanotechnology in the treatment of viral infections. <i>Therapeutic Advances in Infectious Disease</i> , 2017, 4, 105-131.	1.8	233
4	Current trends in computer aided drug design and a highlight of drugs discovered via computational techniques: A review. <i>European Journal of Medicinal Chemistry</i> , 2021, 224, 113705.	5.5	229
5	“Off“on“ fluorescent sensors for physiological levels of magnesium ions based on photoinduced electron transfer (PET), which also behave as photoionic OR logic gates. <i>Journal of the Chemical Society Chemical Communications</i> , 1994, .	2.0	127
6	The Current Status of Heterogeneous Palladium Catalysed Heck and Suzuki Cross-Coupling Reactions. <i>Molecules</i> , 2018, 23, 1676.	3.8	123
7	Nanotechnology and the Treatment of HIV Infection. <i>Viruses</i> , 2012, 4, 488-520.	3.3	106
8	Enantioselective Organocatalyzed Transformations of $\beta$ -Ketoesters. <i>Chemical Reviews</i> , 2016, 116, 9375-9437.	47.7	105
9	Integrated Approach to Structure-Based Enzymatic Drug Design: Molecular Modeling, Spectroscopy, and Experimental Bioactivity. <i>Chemical Reviews</i> , 2014, 114, 493-537.	47.7	100
10	Fluorescent signalling of the brain neurotransmitter $\gamma$ -aminobutyric acid and related amino acid zwitterions. <i>Chemical Communications</i> , 1996, , 2191-2192.	4.1	96
11	Hydaphile Channels: Structural and Fluorescent Probes of Position and Function in a Phospholipid Bilayer. <i>Journal of the American Chemical Society</i> , 1999, 121, 9043-9052.	13.7	67
12	Neutral Molecule Receptor Systems Using Ferrocene's “Atomic Ball Bearing“Character as the Flexible Element. <i>Journal of the American Chemical Society</i> , 1997, 119, 1609-1618.	13.7	60
13	The effect of N-methylation of amino acids (Ac-X-OMe) on solubility and conformation: a DFT study. <i>Organic and Biomolecular Chemistry</i> , 2015, 13, 9993-10006.	2.8	55
14	Structural and functional features of enzymes of <i>Mycobacterium tuberculosis</i> peptidoglycan biosynthesis as targets for drug development. <i>Tuberculosis</i> , 2015, 95, 95-111.	1.9	54
15	Anticancer activity of ruthenium(II) arene complexes bearing 1,2,3,4-tetrahydroisoquinoline amino alcohol ligands. <i>European Journal of Medicinal Chemistry</i> , 2013, 66, 407-414.	5.5	51
16	Planar Bilayer Conductance and Fluorescence Studies Confirm the Function and Location of a Synthetic, Sodium-Ion-Conducting Channel in a Phospholipid Bilayer Membrane. <i>Journal of the American Chemical Society</i> , 1997, 119, 9061-9062.	13.7	45
17	Computational model for the acylation step of the $\beta$ -lactam ring: Potential application for l,d-transpeptidase 2 in <i>mycobacterium tuberculosis</i> . <i>Journal of Molecular Structure</i> , 2017, 1128, 94-102.	3.6	41
18	<scp>DFT</scp> study of the acid-catalyzed esterification reaction mechanism of methanol with carboxylic acid and its halide derivatives. <i>International Journal of Quantum Chemistry</i> , 2018, 118, e25497.	2.0	41

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19	Synthesis and evaluation of SQ109 analogues as potential anti-tuberculosis candidates. European Journal of Medicinal Chemistry, 2010, 45, 2075-2079.	5.5	39
20	Mechanistic investigation of the uncatalyzed esterification reaction of acetic acid and acid halides with methanol: a DFT study. Journal of Molecular Modeling, 2016, 22, 235.	1.8	39
21	Synthesis and structural studies of pentacycloundecane-based HIV-1 PR inhibitors: A hybrid 2D NMR and docking/QM/MM/MD approach. European Journal of Medicinal Chemistry, 2011, 46, 3976-3985.	5.5	38
22	Synthesis of chiral pentacyclo-undecane ligands and their use in the enantioselective alkylation of benzaldehyde with diethylzinc. Tetrahedron: Asymmetry, 2004, 15, 2661-2666.	1.8	37
23	Differential flap dynamics in $\text{I}_{\text{L}}$ - $\text{d}_{\text{L}}$ -transpeptidase2 from mycobacterium tuberculosis revealed by molecular dynamics. Molecular BioSystems, 2017, 13, 1223-1234.	2.9	36
24	Pentacyclo-undecane derived cyclic tetra-amines: Synthesis and evaluation as potent anti-tuberculosis agents. European Journal of Medicinal Chemistry, 2009, 44, 4297-4305.	5.5	34
25	Synthesis and Screening of C <sup>1</sup> -Substituted Tetrahydroisoquinoline Derivatives for Asymmetric Transfer Hydrogenation Reactions. European Journal of Organic Chemistry, 2010, 2010, 972-980.	2.4	33
26	Pentacycloundecane-based inhibitors of wild-type C-South African HIV-protease. Bioorganic and Medicinal Chemistry Letters, 2011, 21, 2274-2277.	2.2	32
27	Comparison of the Molecular Dynamics and Calculated Binding Free Energies for Nine FDA-Approved HIV-1 PR Drugs Against Subtype B and C-CA HIV PR. Chemical Biology and Drug Design, 2013, 81, 208-218.	3.2	32
28	Novel tetrahydroisoquinoline based organocatalysts for asymmetric Diels-Alder reactions: insight into the catalytic mode using ROESY NMR and DFT studies. Tetrahedron: Asymmetry, 2010, 21, 2859-2867.	1.8	30
29	Design and study of peptide-based inhibitors of amylin cytotoxicity. Bioorganic and Medicinal Chemistry Letters, 2010, 20, 1360-1362.	2.2	29
30	Novel polycyclic cage-1,2-diamines as potential anti-tuberculosis agents. European Journal of Medicinal Chemistry, 2012, 54, 1-9.	5.5	28
31	Synthesis of tetrahydroisoquinoline-diamine ligands and their application in asymmetric transfer hydrogenation. Tetrahedron: Asymmetry, 2010, 21, 679-687.	1.8	26
32	Total synthesis of a depsidomycin analogue by convergent solid-phase peptide synthesis and macrolactonization strategy for antitubercular activity. Journal of Peptide Science, 2011, 17, 683-689.	1.4	26
33	Synthesis of tetrahydroisoquinoline (TIQ)-oxazoline ligands and their application in enantioselective Henry reactions. Tetrahedron: Asymmetry, 2010, 21, 846-852.	1.8	25
34	Visualization of Time-Dependent Distribution of Rifampicin in Rat Brain Using MALDI MSI and Quantitative LCMS/MS. Assay and Drug Development Technologies, 2015, 13, 277-284.	1.2	25
35	Pore formation in phospholipid bilayers by amphiphilic cavitands. Organic and Biomolecular Chemistry, 2011, 9, 4498.	2.8	24
36	Binding Free Energy Calculations of Nine FDA-Approved Protease Inhibitors Against HIV-1 Subtype C I36T-T Containing 100 Amino Acids Per Monomer. Chemical Biology and Drug Design, 2016, 87, 487-498.	3.2	23

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37	A comparative modeling and molecular docking study on <i>Mycobacterium tuberculosis</i> targets involved in peptidoglycan biosynthesis. <i>Journal of Biomolecular Structure and Dynamics</i> , 2016, 34, 2399-2417.	3.5	23
38	Crystal, spectroscopic and quantum mechanics studies of Schiff bases derived from 4-nitrocinnamaldehyde. <i>Scientific Reports</i> , 2021, 11, 8151.	3.3	23
39	Tetrahydroisoquinoline-based <i>N</i> -Oxides as Chiral Organocatalysts for the Asymmetric Allylation of Aldehydes. <i>European Journal of Organic Chemistry</i> , 2011, 2011, 6923-6932.	2.4	22
40	Pentacycloundecane derived hydroxy acid peptides: A new class of irreversible non-scissile ether bridged type isoster as potential HIV-1 wild type C-SA protease inhibitors. <i>Bioorganic Chemistry</i> , 2012, 40, 19-29.	4.1	22
41	Purification and characterization of naturally occurring HIV-1 (South African subtype C) protease mutants from inclusion bodies. <i>Protein Expression and Purification</i> , 2016, 122, 90-96.	1.3	22
42	Fluorescent PET(Photoinduced Electron Transfer) Sensors for Calcium Ions. Extension to Multiple Fluorophores and Virtual Spacers. <i>Chemistry Letters</i> , 1995, 24, 125-126.	1.3	21
43	Iridium-catalyzed asymmetric hydrogenation of olefins using TiQ phosphine-oxazoline ligands. <i>Tetrahedron: Asymmetry</i> , 2010, 21, 2295-2301.	1.8	21
44	Determination of the antitubercular drug PA-824 in rat plasma, lung and brain tissues by liquid chromatography tandem mass spectrometry: Application to a pharmacokinetic study. <i>Journal of Chromatography B: Analytical Technologies in the Biomedical and Life Sciences</i> , 2015, 988, 187-194.	2.3	21
45	A tris(macrocycle) that exhibits H-bond-induced blockage of the cation channel fraction in a phospholipid bilayer. <i>Chemical Communications</i> , 1996, , 2147.	4.1	19
46	Synthesis, 2D-NMR and molecular modelling studies of pentacycloundecane lactam-peptides and peptoids as potential HIV-1 wild type C-SA protease inhibitors. <i>Journal of Enzyme Inhibition and Medicinal Chemistry</i> , 2013, 28, 78-88.	5.2	19
47	Simulating the inhibition reaction of <i>Mycobacterium tuberculosis</i> <i>L</i> - <i>d</i> -transpeptidase 2 by carbapenems. <i>Chemical Communications</i> , 2015, 51, 12560-12562.	4.1	19
48	Targeting the cell wall of <i>Mycobacterium tuberculosis</i> : a molecular modeling investigation of the interaction of imipenem and meropenem with <i>L</i> - <i>D</i> -transpeptidase 2. <i>Journal of Biomolecular Structure and Dynamics</i> , 2016, 34, 304-317.	3.5	18
49	Lansoprazole-sulfide, pharmacokinetics of this promising anti-tuberculous agent. <i>Biomedical Chromatography</i> , 2017, 31, e4035.	1.7	18
50	Synthesis of chiral pentacyclo-undecane macrocycles and their use in enantioselective Michael addition reactions. <i>Tetrahedron: Asymmetry</i> , 2004, 15, 3775-3781.	1.8	17
51	Microwave assisted SPPS of amylin and its toxicity of the pure product to RIN-5F cells. <i>Biopolymers</i> , 2010, 94, 323-330.	2.4	17
52	Microwave-Assisted Synthesis of Guanidine Organocatalysts Bearing a Tetrahydroisoquinoline Framework and Their Evaluation in Michael Addition Reactions. <i>European Journal of Organic Chemistry</i> , 2012, 2012, 3331-3337.	2.4	16
53	Synthesis and molecular modelling studies of novel carbapeptide analogs for inhibition of HIV-1 protease. <i>European Journal of Medicinal Chemistry</i> , 2012, 53, 13-21.	5.5	16
54	Synthesis, screening and computational investigation of pentacycloundecane-peptoids as potent CSA-HIV PR inhibitors. <i>European Journal of Medicinal Chemistry</i> , 2012, 57, 459-467.	5.5	15

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55	Pentacycloundecaneâ€¢diolâ€¢Based HIVâ€¢1 Protease Inhibitors: Biological Screening, 2<scp>Dâ€...</scp>NMR, and Molecular Simulation Studies. <i>ChemMedChem</i> , 2012, 7, 1009-1019.	3.2	15
56	Molecular Modeling of <i>T.Ârangeli, T.Âbrucei gambiense</i> and <i>T.Âevansi</i> Sialidases in Complex with the DANA Inhibitor. <i>Chemical Biology and Drug Design</i> , 2012, 80, 114-120.	3.2	15
57	Catalytic asymmetric carbonâ€¢carbon bond forming reactions catalyzed by tetrahydroisoquinoline (TIQ) N,Nâ€¢2-dioxide ligands. <i>Tetrahedron: Asymmetry</i> , 2013, 24, 191-195.	1.8	15
58	Development and validation of a liquid chromatographyâ€¢tandem mass spectrometry (LCâ€¢MS/MS) method for the quantification of tigecycline in rat brain tissues. <i>Biomedical Chromatography</i> , 2016, 30, 837-845.	1.7	15
59	Unraveling the concerted catalytic mechanism of the human immunodeficiency virus type 1 (HIV-1) protease: a hybrid QM/MM study. <i>Structural Chemistry</i> , 2019, 30, 409-417.	2.0	15
60	Proline N-oxides: modulators of the 3D conformation of linear peptides through â€œNO-turnsâ€¢ Organic and Biomolecular Chemistry, 2014, 12, 4479.	2.8	14
61	Potential inhibition of HIV-1 encapsidation by oligoribonucleotide&ndash;dendrimer nanoparticle complexes. <i>International Journal of Nanomedicine</i> , 2017, Volume 12, 317-325.	6.7	14
62	The Impact of Active Site Mutations of South African <scp>HIV PR</scp> on Drug Resistance: Insight from Molecular Dynamics Simulations, Binding Free Energy and Perâ€¢Residue Footprints. <i>Chemical Biology and Drug Design</i> , 2014, 83, 472-481.	3.2	13
63	Pentacycloundecane lactam vs lactone norstatine type protease HIV inhibitors: binding energy calculations and DFT study. <i>Journal of Biomedical Science</i> , 2015, 22, 15.	7.0	13
64	PdCuCeOâ€¢TPAB: a new catalytic system for quasi-heterogeneous Suzukiâ€¢Miyaura cross-coupling reactions under ligand-free conditions in water. <i>New Journal of Chemistry</i> , 2017, 41, 13560-13566.	2.8	13
65	The catalytic role of water in the binding site of L,D-transpeptidase 2 within acylation mechanism: A QM/MM (ONIOM) modelling. <i>Tuberculosis</i> , 2018, 113, 222-230.	1.9	13
66	Identification of potent L,D-transpeptidase 5 inhibitors for <i>Mycobacterium tuberculosis</i> as potential anti-TB leads: virtual screening and molecular dynamics simulations. <i>Journal of Molecular Modeling</i> , 2019, 25, 328.	1.8	13
67	The Driving Force for the Acylation of $\beta^2$ â€¢Lactam Antibiotics by L,Dâ€¢Transpeptidase 2: Quantum Mechanics/Molecular Mechanics (QM/MM) Study. <i>ChemPhysChem</i> , 2019, 20, 1126-1134.	2.1	13
68	Structure and Function of L,D- and D,D-Transpeptidase Family Enzymes from <i>Mycobacterium tuberculosis</i> . <i>Current Medicinal Chemistry</i> , 2020, 27, 3250-3267.	2.4	13
69	Intracellular localization of gold nanoparticles with targeted delivery in MT-4 lymphocytes. <i>Advances in Natural Sciences: Nanoscience and Nanotechnology</i> , 2016, 7, 045013.	1.5	12
70	Investigation of the binding free energies of FDA approved drugs against subtype B and C-SA HIV PR: ONIOM approach. <i>Journal of Molecular Graphics and Modelling</i> , 2017, 76, 77-85.	2.4	12
71	Inhibition mechanism of L,D-transpeptidase 5 in presence of the $\beta^2$ -lactams using ONIOM method. <i>Journal of Molecular Graphics and Modelling</i> , 2019, 87, 204-210.	2.4	12
72	Development and Evaluation of Peptide-Functionalized Gold Nanoparticles for HIV Integrase Inhibition. <i>International Journal of Peptide Research and Therapeutics</i> , 2019, 25, 311-322.	1.9	12

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73	Synthesis and NMR assignment of pentacycloundecane precursors of potential pharmaceutical agents. Magnetic Resonance in Chemistry, 2010, 48, 249-255.	1.9	11
74	SQ109 analogues as potential antimicrobial candidates. Medicinal Chemistry Research, 2011, 20, 1394-1401.	2.4	11
75	Organocatalytic Mannich Reactions on a Carbapenem Core – Synthesis of Mannich Bases and Bicyclic Diazanonanes. European Journal of Organic Chemistry, 2014, 2014, 2253-2260.	2.4	11
76	Preparation of Enantiomerically Pure <i>C</i> <sub>4</sub> –Symmetric Tetramethoxyresorcarenes by Using the (–)-S-Phenylethyl Isocyanate Chiral Auxiliary. European Journal of Organic Chemistry, 2014, 2014, 4600-4609.	2.4	11
77	An insight to the molecular interactions of the FDA approved HIV-PR drugs against L38L <sup>T</sup> N <sup>T</sup> L PR mutant. Journal of Computer-Aided Molecular Design, 2018, 32, 459-471.	2.9	11
78	NMR elucidation of some pentacycloundecane derived ligands. Structural Chemistry, 2007, 18, 633-639.	2.0	10
79	NMR elucidation of some ligands derived from the pentacycloundecane skeleton. Structural Chemistry, 2008, 19, 429-434.	2.0	10
80	I36T <sup>T</sup> mutation in South African subtype C (C-SA) HIV-1 protease significantly alters protease-drug interactions. Biological Chemistry, 2017, 398, 1109-1117.	2.5	10
81	Molecular insight on the non-covalent interactions between carbapenems and l,d-transpeptidase 2 from Mycobacterium tuberculosis: ONIOM study. Journal of Computer-Aided Molecular Design, 2018, 32, 687-701.	2.9	10
82	Concerted hydrolysis mechanism of HIV-1 natural substrate against subtypes B and C-SA PR: insight through molecular dynamics and hybrid QM/MM studies. Physical Chemistry Chemical Physics, 2020, 22, 2530-2539.	2.8	10
83	A Review of Cyanoacetyl Indoles (CAIs): Versatile Starting Materials in Organic Synthesis.. Current Organic Synthesis, 2013, 10, 737-750.	1.3	10
84	A Comparison of Phenyl- and Naphthal-substituted tris(Macrocycle), Cation-conducting Channels to Assess the Effect of Extended Aromaticity. Tetrahedron Letters, 1997, 38, 6339-6342.	1.4	9
85	Linear and cyclic glycopeptide as HIV protease inhibitors. European Journal of Medicinal Chemistry, 2013, 60, 144-154.	5.5	9
86	From Recognition to Reaction Mechanism: An Overview on the Interactions between HIV-1 Protease and its Natural Targets. Current Medicinal Chemistry, 2020, 27, 2514-2549.	2.4	9
87	Synthesis and NMR elucidation of novel pentacyclo-undecane diamine ligands. Structural Chemistry, 2009, 20, 1067-1076.	2.0	8
88	Novel Linear Diamine Disubstituted Polycyclic Cage <sup>TM</sup> Derivatives as Potential Antimycobacterial Candidates. Chemical Biology and Drug Design, 2011, 78, 1022-1030.	3.2	8
89	Optically active diaryl tetrahydroisoquinoline derivatives. Acta Crystallographica Section C: Crystal Structure Communications, 2011, 67, o100-o103.	0.4	8
90	Asymmetric conjugate addition of thioglycolate to a range of chalcones using tetrahydroisoquinoline (TIQ) N,N <sup>2</sup> -dioxide ligands. Tetrahedron: Asymmetry, 2012, 23, 616-622.	1.8	8

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91	Interaction of <i>&lt;math&gt;\text{i}\alpha\beta&lt;/math&gt;</i> -Amyloid Interactions with Peptide Functionalized Gold Nanoparticles. <i>Journal of Nanoscience and Nanotechnology</i> , 2012, 12, 2179-2184.	0.9	8
92	Rapid and widespread distribution of doxycycline in rat brain: a mass spectrometric imaging study. <i>Xenobiotica</i> , 2016, 46, 385-392.	1.1	8
93	Pd <sub>0.09</sub> Ce <sub>0.91</sub> O <sub>2-<math>\delta</math></sub> : A sustainable ionic solid-solution precatalyst for heterogeneous, ligand free Heck coupling reactions. <i>Molecular Catalysis</i> , 2017, 443, 60-68.	2.0	8
94	Microwave-assisted synthesis of a new series of resorcin[4]arene cavitand-capped porphyrin capsules. <i>Organic and Biomolecular Chemistry</i> , 2009, 7, 3958.	2.8	7
95	<i>&lt;math&gt;\text{In Vitro}&lt;/math&gt;</i> Antifungal and Antibacterial Activities of Pentacycloundecane Tetra- $\epsilon$ Amines. <i>Chemical Biology and Drug Design</i> , 2011, 77, 295-299.	3.2	7
96	Enzymatic Activation of a Peptide Functionalised Gold Nanoparticle System for Prodrug Delivery. <i>Journal of Nanoscience and Nanotechnology</i> , 2011, 11, 3075-3083.	0.9	6
97	(S)-N-Benzyl-2-methyl-1,2,3,4-tetrahydroisoquinoline-3-carboxamide. <i>Acta Crystallographica Section E: Structure Reports Online</i> , 2011, 67, o67-o67.	0.2	6
98	Bisthiourea: thermal and structural investigation. <i>Journal of Thermal Analysis and Calorimetry</i> , 2013, 111, 597-603.	3.6	6
99	Inhibition of <i>&lt;math&gt;\text{i}Mycobacterium tuberculosis&lt;/math&gt;</i> L,D $\alpha$ -Transpeptidase 5 by Carbapenems: MD and QM/MM Mechanistic Studies. <i>ChemistrySelect</i> , 2018, 3, 13603-13612.	1.5	6
100	Exploring the concerted mechanistic pathway for HIV-1 PR $\alpha$ "substrate revealed by umbrella sampling simulation. <i>Journal of Biomolecular Structure and Dynamics</i> , 2022, 40, 1736-1747.	3.5	6
101	Microwave-assisted synthesis of <i>&lt;math&gt;\text{meso}&lt;/math&gt;</i> -carboxyalkyl-BODIPYs and an application to fluorescence imaging. <i>Organic and Biomolecular Chemistry</i> , 2020, 18, 7876-7883.	2.8	6
102	(1R,3S)-Methyl 2-benzyl-6,7-dimethoxy-1-phenyl-1,2,3,4-tetrahydroisoquinoline-3-carboxylate. <i>Acta Crystallographica Section E: Structure Reports Online</i> , 2009, 65, o3278-o3278.	0.2	6
103	Synthesis, Characterization and Biocompatibility of a Multifunctional Gold Nanoparticle System for the Delivery of Single-Stranded RNA to Lymphocytes. <i>South African Journal of Chemistry</i> , 2018, 71, 1-14.	0.6	6
104	NMR elucidation of a novel ( <i>&lt;math&gt;\text{i}S&lt;/math&gt;</i> ) $\alpha$ -pentacyclo $\alpha$ undecane bis(4-phenyloxazoline) ligand and related derivatives. <i>Magnetic Resonance in Chemistry</i> , 2008, 46, 1089-1095.	1.9	5
105	(1S,3S)-Methyl 2-benzyl-6,7-dimethoxy-1-phenyl-1,2,3,4-tetrahydroisoquinoline-3-carboxylate. <i>Acta Crystallographica Section E: Structure Reports Online</i> , 2011, 67, o1403-o1403.	0.2	5
106	L-Proline organocatalyzed Michael synthesis of monobactam and carbapenem $\beta$ -lactam cores. <i>Tetrahedron: Asymmetry</i> , 2014, 25, 969-973.	1.8	5
107	Synthesis and structural elucidation of a novel polymorph of alcaftadine. <i>Spectrochimica Acta - Part A: Molecular and Biomolecular Spectroscopy</i> , 2015, 142, 311-319.	3.9	5
108	Theoretical Model for HIV-1 PR That Accounts for Substrate Recognition and Preferential Cleavage of Natural Substrates. <i>Journal of Physical Chemistry B</i> , 2019, 123, 6389-6400.	2.6	5

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109	Tetrakis- $\frac{1}{4}$ -L-alanine- $\text{^{19}O}_2$ -bis[tetraaquaterbium(III)] hexaperchlorate. Acta Crystallographica Section E: Structure Reports Online, 2010, 66, m193-m194.	0.2	5
110	{(1R,3S)-2-Benzyl-6,7-dimethoxy-1-phenyl-1,2,3,4-tetrahydroisoquinolin-3-yl}diphenylmethanol. Acta Crystallographica Section E: Structure Reports Online, 2010, 66, o638-o638.	0.2	5
111	7,11,15,28-Tetrakis(bromomethyl)-1,21,23,25-tetrapentylresorcin[4]arene cavitand 0.415-hydrate. Acta Crystallographica Section E: Structure Reports Online, 2007, 63, o4345-o4345.	0.2	4
112	Redetermination of chlorido(2,2â€¢:6â€¢,2â€¢â€¢-terpyridine-â€¢3N,Nâ€¢,Nâ€¢â€¢)gold(I) dichloride trihydrate at 173 K. Acta Crystallographica Section E: Structure Reports Online, 2008, 64, m1240-m1240.	0.2	4
113	1,1â€¢-(Propane-1,3-diyl)bis(3-phenylurea). Acta Crystallographica Section E: Structure Reports Online, 2011, 67, o2552-o2552.	0.2	4
114	(2-(Benzod[d]thiazol-2-yl-methoxy)-5-chlorophenyl)(phenyl)methanone. Acta Crystallographica Section E: Structure Reports Online, 2012, 68, o3125-o3125.	0.2	4
115	Novel PCU cage diol peptides as potential targets against wild-type CSA HIV-1 protease: synthesis, biological screening and molecular modelling studies. Medicinal Chemistry Research, 2013, 22, 3918-3933.	2.4	4
116	[2-(1,3-Benzothiazol-2-ylmethoxy)-5-bromophenyl](4-chlorophenyl)methanone. Acta Crystallographica Section E: Structure Reports Online, 2013, 69, o70-o70.	0.2	4
117	Isolation, characterization and x-ray structure determination of the schiff base ligand: 5-methyl-2-phenyl-4-[phenyl-(4-phenyl-thiazol-2-ylamino)-methylene]-2,4-dihydro-pyrazol-3-one. South African Journal of Chemistry, 2015, 68, .	0.6	4
118	Kinetic and thermodynamic characterisation of HIV-protease inhibitors against E35Dâ†‘Gâ†‘S mutant in the South African HIV-1 subtype C protease. Journal of Enzyme Inhibition and Medicinal Chemistry, 2019, 34, 1451-1456.	5.2	4
119	Serendipitous discovery of new pentacycloundecane molecules. Journal of Molecular Structure, 2020, 1204, 127497.	3.6	4
120	Ethyl 4-(4-chlorophenyl)-6-methyl-2-thioxo-1,2,3,4-tetrahydropyrimidine-5-carboxylate. Acta Crystallographica Section E: Structure Reports Online, 2009, 65, o2518-o2518.	0.2	4
121	Pdo.o2Ceo.98O2: a copper- and ligand-free quasi-heterogeneous catalyst for aquacatalytic Sonogashira cross-coupling reaction. Journal of the South African Institute of Mining and Metallurgy, 2017, 117, 955-962.	0.5	4
122	Experimental measurement of kinetic parameters using quantum plasmonic sensing. Journal of Applied Physics, 2022, 131, 084402.	2.5	4
123	Synthesis, Crystal structure, photoluminescence properties and quantum mechanics studies of two schiff bases of 2-amino-p-cresol. Journal of Molecular Structure, 2022, 1262, 133046.	3.6	4
124	5,11,17,23-Tetramethyl-2,8,14,20-tetrakis(2-phenylethyl)-4,6,10,12,16,18,22,24-octahydroxycalix[4]arene methanol pentasolvate 0.10-hydrate. Acta Crystallographica Section E: Structure Reports Online, 2007, 63, o4346-o4346.	0.2	3
125	7,11,15,28-Tetramethyl-1,21,23,25-tetrakis(2-phenylethyl)resorcin[4]arene ethyl acetate clathrate. Acta Crystallographica Section E: Structure Reports Online, 2008, 64, o98-o98.	0.2	3
126	(2 <i>i</i> R, <i>i</i> S)-Methyl 2-hydroxy-3-(4-methylbenzenesulfonamido)-3-phenylpropanoate. Acta Crystallographica Section E: Structure Reports Online, 2010, 66, o3279-o3280.	0.2	3

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127	(R)-4-Phenyl-2-[(S)-1,2,3,4-tetrahydroisoquinolin-3-yl]-4,5-dihydro-1,3-oxazole. Acta Crystallographica Section E: Structure Reports Online, 2010, 66, o1818-o1818.	0.2	3
128	3,3â€2-Diphenyl-1,1â€2-(butane-1,4-diyl)dithiourea. Acta Crystallographica Section E: Structure Reports Online, 2011, 67, o2380-o2380.	0.2	3
129	4,5,6,10,11,12,16,17,18,22,23,24-Dodecakis[(methoxycarbonyl)methoxy]-2,8,14,20-tetrapentylresorcin[4]arene. Acta Crystallographica Section E: Structure Reports Online, 2011, 67, o3305-o3306.	0.2	3
130	<math>\text{<i>N</i>-Benzyl-5-(dimethylamino)naphthalene-1-sulfonamide. Acta Crystallographica Section E: Structure Reports Online, 2011, 67, o2458-o2459.}	0.2	3
131	1,1â€2-(Ethane-1,2-diyl)bis(3-phenylthiourea). Acta Crystallographica Section E: Structure Reports Online, 2011, 67, o2819-o2819.	0.2	3
132	Ethyl 2-(3,5-dimethyl-1,1-dioxo-2 <i>H</i> -1,1â€2-<math>\text{H</i>-1}^{\circ}\text{<sup>6</sup>},2,6-thiadiazin-4-yl)benzoate. Acta Crystallographica Section E: Structure Reports Online, 2012, 68, o2160-o2160.	0.2	3
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135	Synthesis and NMR elucidation of novel amino acid cavitand derivatives. Tetrahedron, 2014, 70, 7057-7066.	1.9	3
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141	N-(10-Bromoanthracen-9-ylmethyl)-N-[2-(5,5-dimethyl-1,3,2-dioxaborinan-2-yl)benzyl]methylamine at 240â€...K. Acta Crystallographica Section E: Structure Reports Online, 2004, 60, o571-o573.	0.2	2
142	3-(2-Amino-1,3-thiazol-4-yl)-6-bromo-2H-chromen-2-one. Acta Crystallographica Section E: Structure Reports Online, 2009, 65, o3047-o3048.	0.2	2
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144	NMR elucidation of novel SQ109 derivatives. Structural Chemistry, 2010, 21, 1203-1209.	2.0	2

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145	(1R,3S)-Methyl 6,7-dimethoxy-1-(4-methoxyphenyl)-1,2,3,4-tetrahydroisoquinoline-3-carboxylate. <i>Acta Crystallographica Section E: Structure Reports Online</i> , 2010, 66, o3105-o3105.	0.2	2
146	tert-ButylN-[(11-exo-benzyloxycarbonyl-8-oxopentacyclo[5.4.0.02,6.03,10.05,9]undecane-11-endo-yloxy)carbonylmethyl]carbamate. <i>Acta Crystallographica Section E: Structure Reports Online</i> , 2010, 66, o2607-o2608.	0.2	2
147	N,Nâ€“-[(8-endo,11-endo-Dihydroxypentacyclo[5.4.0.02,6.03,10.05,9]undecane-8,11-diyl)bis(methylenecarbonyl)]di-L-phenylalanine. <i>Acta Crystallographica Section E: Structure Reports Online</i> , 2010, 66, o2537-o2538.	0.2	2
148	(3S)-2-Benzyl-3-carboxy-1,2,3,4-tetrahydroisoquinolinium chloride monohydrate. <i>Acta Crystallographica Section E: Structure Reports Online</i> , 2011, 67, o228-o228.	0.2	2
149	3-Benzyl-5,7-dimethoxychroman-4-ol. <i>Acta Crystallographica Section E: Structure Reports Online</i> , 2011, 67, o703-o703.	0.2	2
150	Ethyl 6-methyl-2-sulfanylidene-4-[4-(trifluoromethyl)phenyl]-1,2,3,4-tetrahydropyrimidine-5-carboxylate. <i>Acta Crystallographica Section E: Structure Reports Online</i> , 2011, 67, o1559-o1560.	0.2	2
151	2-[(1R,3S)-6,7-Dimethoxy-1-phenyl-1,2,3,4-tetrahydroisoquinolin-3-yl]-4-phenyl-1,3-thiazole. <i>Acta Crystallographica Section E: Structure Reports Online</i> , 2011, 67, o2722-o2722.	0.2	2
152	4,10,16,22-Tetrakis(2-chloroacetoxy)-6,12,18,24-tetramethoxy-2,8,14,20-tetrapentylresorcin[4]arene. <i>Acta Crystallographica Section E: Structure Reports Online</i> , 2011, 67, o2907-o2907.	0.2	2
153	Ethyl 4-(1,3-benzodioxol-5-yl)-6-methyl-2-sulfanylidene-1,2,3,4-tetrahydropyrimidine-5-carboxylate. <i>Acta Crystallographica Section E: Structure Reports Online</i> , 2011, 67, o3069-o3070.	0.2	2
154	6,12,18,24-Tetramethoxy-4,10,16,22-tetrakis[(methoxycarbonyl)methoxy]-2,8,14,20-tetrakis(2-phenylethyl)resorcin[4]arene. <i>Acta Crystallographica Section E: Structure Reports Online</i> , 2012, 68, o97-o98.	0.2	2
155	2-(3,5-Dimethyl-1,1-dioxo-2H-1â†»6,2,6-thiadiazin-4-yl)benzoic acid. <i>Acta Crystallographica Section E: Structure Reports Online</i> , 2012, 68, o2859-o2859.	0.2	2
156	Synthesis of novel N-alkyl-2-(3,5-dimethyl-1,1-dioxido-2H-1,2,6-thiadiazin-4-yl)benzamides as potential antimicrobial agents. <i>Journal of Pharmacy Research</i> , 2013, 6, 355-360.	0.4	2
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158	Optimized Procedure for Recovering HIV-1 Protease (C-SA) from Inclusion Bodies. <i>Protein Journal</i> , 2019, 38, 30-36.	1.6	2
159	7,11,15,28-Tetrakis[(2-formylphenoxy)methyl]-1,21,23,25-tetramethylresorcin[4]arene cavitand ethyl acetate clathrate at 173 K. <i>Acta Crystallographica Section E: Structure Reports Online</i> , 2009, 65, o692-o693.	0.2	2
160	2-(4-Chloro-3-nitrophenyl)-4-(4-chlorophenyl)-1,3-thiazole. <i>Acta Crystallographica Section E: Structure Reports Online</i> , 2009, 65, o2611-o2612.	0.2	2
161	3-(2-Amino-1,3-thiazol-4-yl)-6-chloro-2H-chromen-2-one. <i>Acta Crystallographica Section E: Structure Reports Online</i> , 2009, 65, o3111-o3111.	0.2	2
162	Di-tert-butyl 2-(tert-butyloxycarbonylmethoxy)phenyliminodiacetate. <i>Acta Crystallographica Section E: Structure Reports Online</i> , 2003, 59, o1650-o1652.	0.2	1

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164	A structural investigation of 1,2 (bis-phenoxy)ethane and its derivatives. <i>Structural Chemistry</i> , 2007, 18, 683-687.	2.0	1
165	NMR elucidation of novel ligands derived from (R)-(+)-camphor. <i>Structural Chemistry</i> , 2009, 20, 925-932.	2.0	1
166	Synthesis and NMR elucidation of novel pentacycloundecane-based peptides. <i>Magnetic Resonance in Chemistry</i> , 2010, 48, 435-442.	1.9	1
167	1,7-Dimethylpentacyclo[5.4.0.0<sup>2,6</sup>.0<sup>3,10</sup>.0<sup>5,9</sup>]undecane-8,11-dione. <i>Acta Crystallographica Section E: Structure Reports Online</i> , 2010, 66, o1901-o1902.	0.2	1
168	Methyl 1-cyclohexyl-6,7-dimethoxy-3,4-dihydroisoquinoline-3-carboxylate. <i>Acta Crystallographica Section E: Structure Reports Online</i> , 2011, 67, o883-o883.	0.2	1
169	(S)-2-Benzyl-N-(2,6-diisopropylphenyl)-1,2,3,4-tetrahydroisoquinoline-3-carboxamide. <i>Acta Crystallographica Section E: Structure Reports Online</i> , 2011, 67, o1106-o1106.	0.2	1
170	4,6,10,12,16,18,22,24-Octa-O-methyl-2,8,14,20-tetrapentylresorcin[4]arene. <i>Acta Crystallographica Section E: Structure Reports Online</i> , 2011, 67, o2565-o2565.	0.2	1
171	1-(2-Aminoethyl)-3-phenylthiourea. <i>Acta Crystallographica Section E: Structure Reports Online</i> , 2011, 67, o2621-o2621.	0.2	1
172	(S)-Methyl 3-(3,4-dimethoxyphenyl)-2-[2-(diphenylphosphanyl)benzamido]propanoate. <i>Acta Crystallographica Section E: Structure Reports Online</i> , 2011, 67, o3285-o3285.	0.2	1
173	(1R,3S)-N-Benzhydryl-2-benzyl-6,7-dimethoxy-1-phenyl-1,2,3,4-tetrahydroisoquinoline-3-carbothioamide. <i>Acta Crystallographica Section E: Structure Reports Online</i> , 2011, 67, o3441-o3441.	0.2	1
174	(S)-Benzyl 3-phenylcarbamoyl-1,2,3,4-tetrahydroisoquinoline-2-carboxylate. <i>Acta Crystallographica Section E: Structure Reports Online</i> , 2012, 68, o883-o883.	0.2	1
175	(S)-Methyl 2-benzamido-3-(3,4-dimethoxyphenyl)propanoate. <i>Acta Crystallographica Section E: Structure Reports Online</i> , 2012, 68, o177-o177.	0.2	1
176	Methyl 2-(2,3,5-trimethyl-1,1-dioxo-2H-1,6,2,6-thiadiazin-4-yl)benzoate. <i>Acta Crystallographica Section E: Structure Reports Online</i> , 2012, 68, o3360-o3360.	0.2	1
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181	cis-N-(2-Hydroxycyclohexyl)-p-toluenesulfonamide. <i>Acta Crystallographica Section E: Structure Reports Online</i> , 2010, 66, o463-o463.	0.2	1
182	(Anthracen-9-ylmethyl)diethylamine at 100°C. <i>Acta Crystallographica Section E: Structure Reports Online</i> , 2005, 61, o52-o54.	0.2	0
183	(4R,5R)-Bis(hydroxydimethylmethyl)-2,2-dimethyl-1,3-dioxolane. <i>Acta Crystallographica Section E: Structure Reports Online</i> , 2006, 62, o4339-o4341.	0.2	0
184	3-Phenyl-1,5-di-2-pyridylpentane-1,5-dione. <i>Acta Crystallographica Section E: Structure Reports Online</i> , 2007, 63, o153-o155.	0.2	0
185	(S)-(+)-2-Formamido-4-methylpentanoic acid. <i>Acta Crystallographica Section E: Structure Reports Online</i> , 2007, 63, o3445-o3445.	0.2	0
186	(1R,4S)-(â")-3,3-Ethylenedioxy-1,7,7-trimethylbicyclo[2.2.1]heptan-2-one. <i>Acta Crystallographica Section E: Structure Reports Online</i> , 2007, 63, o3765-o3765.	0.2	0
187	(S)-(+)-2-Formylamino-3-methylbutanoic acid. <i>Acta Crystallographica Section E: Structure Reports Online</i> , 2007, 63, o3912-o3912.	0.2	0
188	endo-11-(Dibenzylamino)tetracyclo[5.4.0.03,10.05,9]undecane-8-one. <i>Acta Crystallographica Section E: Structure Reports Online</i> , 2011, 67, o619-o619.	0.2	0
189	tert-Butyl 2-hydroxy-3-(4-methylbenzenesulfonamido)butanoate. <i>Acta Crystallographica Section E: Structure Reports Online</i> , 2011, 67, o648-o648.	0.2	0
190	[(1R,3S)-6,7-Dimethoxy-1-phenyl-1,2,3,4-tetrahydroisoquinolin-3-yl]methanol 2.33-hydrate. <i>Acta Crystallographica Section E: Structure Reports Online</i> , 2011, 67, o700-o700.	0.2	0
191	Benzyl 5-hydroxy-4-oxapentacyclo[5.4.1.02,6.03,10.08,11]dodecane-3-carboxylate. <i>Acta Crystallographica Section E: Structure Reports Online</i> , 2011, 67, o877-o877.	0.2	0
192	6,7-Dimethoxy-3-methoxycarbonyl-1-(2-methoxyphenyl)-3,4-dihydroisoquinoline 2-oxide. <i>Acta Crystallographica Section E: Structure Reports Online</i> , 2011, 67, o1352-o1352.	0.2	0
193	N-(Adamantan-1-yl)-2-chloroacetamide. <i>Acta Crystallographica Section E: Structure Reports Online</i> , 2011, 67, o1444-o1444.	0.2	0
194	(1S,3S)-Methyl 6,7-dimethoxy-1-phenyl-1,2,3,4-tetrahydroisoquinoline-3-carboxylate. <i>Acta Crystallographica Section E: Structure Reports Online</i> , 2011, 67, o1501-o1501.	0.2	0
195	(S)-4-Phenyl-2-(1,2,3,4-tetrahydroisoquinolin-3-yl)-1,3-thiazole. <i>Acta Crystallographica Section E: Structure Reports Online</i> , 2012, 68, o176-o176.	0.2	0
196	(S)-4-tert-Butyl-2-(1,2,3,4-tetrahydroisoquinolin-3-yl)-1,3-thiazole. <i>Acta Crystallographica Section E: Structure Reports Online</i> , 2012, 68, o2487-o2487.	0.2	0
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201	Crystal structure of bis(tert.-butylaminium) 4-(2-carboxylatophenyl)-3,5-dimethyl-4H-1,2,6-thiadiazin-4-ide 1,1-dioxide , C20H34N4O4S. Zeitschrift Fur Kristallographie - New Crystal Structures, 2013, 228, 469-470.	0.3	0
202	Crystal structure of (S)-tert-butyl 3-carbamothioyl-3,4-dihydroisoquinoline- 2(1H)-carboxylate, C63H86N8O9S4. Zeitschrift Fur Kristallographie - New Crystal Structures, 2013, 228, 449-450.	0.3	0
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204	Crystal structure of 2-(2,4-dioxopentan-3-yl)benzoic acid, C12H12O4. Zeitschrift Fur Kristallographie - New Crystal Structures, 2014, 229, 323-324.	0.3	0
205	Crystal structure of 2-methyl-(2,4-dioxopentan-3-yl)benzoate, C13H14O4. Zeitschrift Fur Kristallographie - New Crystal Structures, 2014, 229, 325-326.	0.3	0
206	Crystal structure of N-butyl-2-(3,5-dimethyl-1,1-dioxido-2H-1,2,6- thiadiazin-4-yl)benzamide, C16H21N3O3S. Zeitschrift Fur Kristallographie - New Crystal Structures, 2014, 229, .	0.3	0
207	Crystal structure of 2-(3-chloro-phenyl)-5-methyl-4-[1-(4-p-tolyl-thiazol-) Tj ETQq1 1 0.784314 rgBT /Overlock 10 Tf 50 427 Td (2-ylamino Crystal Structures, 2014, 229, 319-320.	0.3	0
208	Crystal structure of 5-methyl-2-phenyl-4-[1-(4-p-tolyl-thiazol-2-ylamino)-ethylidene]-2,4-dihdropyrazol-3-one, C22H20N4OS. Zeitschrift Fur Kristallographie - New Crystal Structures, 2014, 229, 313-314.	0.3	0
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221	Crystal structure of 4,10,16,22-tetrahydroxy-6,12,18,24-tetramethoxy-2,8,14,20-tetraethylphenylresorcin[4]arene "ethyl acetate (1/1), C <sub>68</sub> H <sub>72</sub> O <sub>10</sub> . Zeitschrift Fur Kristallographie - New Crystal Structures, 2017, 232, 305-307.	0.3	0
222	Crystal structure of 5,11,17,23-tetra(tert-butyl)-25,26,27,28-tetrahexoxycalix[4]arene, C <sub>68</sub> H <sub>104</sub> O <sub>4</sub> . Zeitschrift Fur Kristallographie - New Crystal Structures, 2017, 232, 397-402.	0.3	0
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226	(4-Hydroxy-3-nitrobenzyl)methylammonium chloride. Acta Crystallographica Section E: Structure Reports Online, 2008, 64, o625-o625.	0.2	0
227	7,11,15,28-Tetrabromo-1,21,23,25-tetraphenethylresorcin[4]arene cavitand "acetone "chloroform (1/1.31/0.69) "at 173 K. Acta Crystallographica Section E: Structure Reports Online, 2009, 65, o631-o632.	0.2	0
228	Crystal structure of (S)-N-benzyl-1,2,3,4-tetrahydroisoquinoline-3-carboxamide, C <sub>17</sub> H <sub>20</sub> N <sub>2</sub> O <sub>2</sub> . Zeitschrift Fur Kristallographie - New Crystal Structures, 2013, 228, 401-402.	0.3	0
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