

Peter C Healy

List of Publications by Year in descending order

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71102

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4601
citing authors

#	ARTICLE	IF	CITATIONS
1	Synthesis and structural characterization of some 1:1 and 1:2 adducts of silver(I) salts with hindered 2022, 535, 120857.	2.4	0
2	Synthesis and structural characterization of some 1:1 adducts of silver(I) salts with (hindered) PR3 bases (R=Åphenyl, o-tolyl, cyclohexyl). <i>Inorganica Chimica Acta</i> , 2022, 536, 120895.	2.4	0
3	Styracifoline from the Vietnamese Plant <i>Desmodium styracifolium</i> : A Potential Inhibitor of Diabetes-Related and Thrombosis-Based Proteins. <i>ACS Omega</i> , 2021, 6, 23211-23221.	3.5	5
4	Solid-State NMR, X-Ray Diffraction, and Theoretical Studies of Neutral Mononuclear Molecular Bis(triphenylphosphine)silver(I) Mono-Carboxylate and -Nitrate Systems. <i>Australian Journal of Chemistry</i> , 2020, 73, 556.	0.9	0
5	Synthesis and Structures of Bis- and Tris-(triphenylarsine)gold(I) Iodides. <i>Australian Journal of Chemistry</i> , 2020, 73, 497.	0.9	0
6	Facile amidinations of 2-aminophenylboronic acid promoted by boronate ester formation. <i>Organic and Biomolecular Chemistry</i> , 2019, 17, 803-806.	2.8	6
7	Synthesis of spirocyclic orthoesters by σ -anomalous TM rhodium(ii)-catalysed intramolecular C-H insertions. <i>Organic and Biomolecular Chemistry</i> , 2018, 16, 256-261.	2.8	8
8	5,6,7,3,4,5-Hexamethoxyflavone from the Australian plant <i>Eremophila debilis</i> (Myoporaceae). <i>FÅ-toterapÅ-Åç</i> , 2018, 126, 90-92.	2.2	3
9	Discovery of new nanomolar inhibitors of GPa: Extension of 2-oxo-1,2-dihydropyridinyl-3-yl amide-based GPa inhibitors. <i>European Journal of Medicinal Chemistry</i> , 2017, 127, 341-356.	5.5	11
10	Crystallization-induced amide bond formation creates a boron-centered spirocyclic system. <i>Heterocyclic Communications</i> , 2017, 23, 167-169.	1.2	4
11	Attempted Synthesis and Unexpected β -Fragmentation of a Hindered β -Keto Nitroxide. <i>Australian Journal of Chemistry</i> , 2017, 70, 1106.	0.9	1
12	Bioactive Dihydro- β -agarofuran Sesquiterpenoids from the Australian Rainforest Plant <i>Maytenus bilocularis</i> . <i>Journal of Natural Products</i> , 2016, 79, 1445-1453.	3.0	33
13	Nucleophilic substitution reactions of $[(\text{I}^5\text{-Cp}^*)\text{Ru}(\text{I}^6\text{-C}_6\text{H}_5\text{CO}_2\text{H})]^+$: Synthesis, characterization and cytotoxicity of organoruthenium ester and amide complexes. <i>Journal of Organometallic Chemistry</i> , 2016, 819, 1-10.	1.8	7
14	2-Oxo-1,2-dihydropyridinyl-3-yl amide-based GPa inhibitors: Design, synthesis and structure-activity relationship study. <i>European Journal of Medicinal Chemistry</i> , 2016, 111, 1-14.	5.5	8
15	Synthesis of two chiral octahydroindole scaffolds for drug discovery. <i>Tetrahedron</i> , 2016, 72, 1225-1228.	1.9	2
16	Antibacterial and antifungal screening of natural products sourced from Australian fungi and characterisation of pestalactams DÅF. <i>Phytochemistry</i> , 2016, 124, 79-85.	2.9	21
17	Entonalactams AÅC: Isoindolinone derivatives from an Australian rainforest fungus belonging to the genus <i>Entonaema</i> . <i>Phytochemistry</i> , 2015, 117, 10-16.	2.9	21
18	Denhaminols AÅH, Dihydro- β -agarofurans from the Endemic Australian Rainforest Plant <i>Denhamia celastroides</i> . <i>Journal of Natural Products</i> , 2015, 78, 111-119.	3.0	21

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19	Design and Synthesis of a Screening Library Using the Natural Product Scaffold 3-Chloro-4-hydroxyphenylacetic Acid. <i>Journal of Natural Products</i> , 2015, 78, 914-918.	3.0	10
20	Synthesis and antimalarial evaluation of amide and urea derivatives based on the thiaplakortone A natural product scaffold. <i>Organic and Biomolecular Chemistry</i> , 2015, 13, 1558-1570.	2.8	25
21	Solving the Supply of Resveratrol Tetramers from Papua New Guinean Rainforest <i>Anisoptera</i> Species That Inhibit Bacterial Type III Secretion Systems. <i>Journal of Natural Products</i> , 2014, 77, 2633-2640.	3.0	16
22	A versatile synthesis of α -tafuramycin A a potent anticancer and parasite attenuating agent. <i>Organic and Biomolecular Chemistry</i> , 2014, 12, 4260-4264.	2.8	6
23	Isomorphism in the structural chemistry of two-coordinate adducts of diphenyl(2-formylphenyl)phosphine and triphenylphosphine with gold(I) halides. <i>Journal of Molecular Structure</i> , 2014, 1072, 253-259.	3.6	20
24	Oligo-nuclear silver thiocyanate complexes with monodentate tertiary phosphine ligands, including novel α -cubane TM and α -step TM tetramer forms of AgSCN \cdot PR ₃ (1 \cdot) ₄ . <i>Dalton Transactions</i> , 2013, 42, 277.	3.3	24
25	The glycogen phosphorylase inhibitor 2-(3-benzylamino-2-oxo-1,2-dihydropyridin-1-yl)-N-(3,4-dichlorobenzyl)acetamide. <i>Acta Crystallographica Section C: Crystal Structure Communications</i> , 2013, 69, 1408-1410.	0.4	0
26	3-Benzhydryl-1,3,4-thiadiazole-2(3H)-thione. <i>Acta Crystallographica Section E: Structure Reports Online</i> , 2013, 69, o1407-o1407.	0.2	0
27	1-[4-Chloro-3-(trifluoromethyl)phenyl]-4-phenyl-1 <i>H</i> -1,2,3-triazole. <i>Acta Crystallographica Section E: Structure Reports Online</i> , 2012, 68, o3159-o3159.	0.2	1
28	Bis[2-(diphenylphosphanyl) ¹⁸ P]benzaldehyde]iodidogold(I). <i>Acta Crystallographica Section E: Structure Reports Online</i> , 2012, 68, m473-m473.	0.2	0
29	Methyl 3-[4-(4-nitrobenzyloxy)phenyl]propanoate. <i>Acta Crystallographica Section E: Structure Reports Online</i> , 2012, 68, o1997-o1998.	0.2	0
30	Dicyclohexylammonium (S)-2-azido-3-phenylpropanoate. <i>Acta Crystallographica Section E: Structure Reports Online</i> , 2012, 68, o2050-o2050.	0.2	0
31	Synthesis and Structural Studies of Tetrahedral [M(Ph ₂ P(CH=CH)PPh ₂) ₂]BPh ₄ and [M(Ph ₂ P(CH ₂) _n PPh ₂) ₂]BPh ₄ (n = 2, 3) Complexes for M = Copper(I), Silver(I), and Gold(I). <i>Australian Journal of Chemistry</i> , 2012, 65, 811.	0.9	5
32	Mechanochemical and solution synthesis, X-ray structure and IR and ³¹ P solid state NMR spectroscopic studies of copper(I) thiocyanate adducts with bulky monodentate tertiary phosphine ligands. <i>Dalton Transactions</i> , 2012, 41, 7513.	3.3	32
33	Selective, Cytotoxic Organoruthenium(II) Full π -Sandwich Complexes: A Structural, Computational and In Vitro Biological Study. <i>Chemistry - an Asian Journal</i> , 2012, 7, 112-121.	3.3	31
34	Solution and mechanochemical syntheses, and spectroscopic and structural studies in the silver(I) (bi)-carbonate: triphenylphosphine system. <i>Dalton Transactions</i> , 2011, 40, 7210.	3.3	32
35	Mono- and 1,1 α ² -Disubstituted Organoruthenium Cyclopentadiene Complexes: Synthesis, Structural Characterization, and Antitumoral Evaluation. <i>Organometallics</i> , 2011, 30, 1395-1403.	2.3	17
36	Structural Systematics of the Anhydrous 1:1 Proton-Transfer Compounds of 3,5-Dinitrosalicylic Acid with Aniline and Monosubstituted Anilines. <i>Journal of Chemical Crystallography</i> , 2011, 41, 1649-1662.	1.1	25

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37	Vibrational and crystal structure analysis of a phenylenedioxydiacetic acid derivative. <i>Journal of Molecular Structure</i> , 2011, 987, 25-33.	3.6	1
38	Arylation of [6,6]-spiroacetal enol ethers: reactivity and rearrangement. <i>Tetrahedron Letters</i> , 2011, 52, 1070-1073.	1.4	6
39	(η^6 -Isopropyl η^5 -phenylcarbamate)(η^5 -pentamethylcyclopentadienyl)ruthenium(II) tetraphenylborate acetone monosolvate. <i>Acta Crystallographica Section E: Structure Reports Online</i> , 2011, 67, m1231-m1231.	0.2	2
40	Structural and Spectroscopic Characterisation of Linearly Coordinated Gold(I) Tribenzylphosphane Complexes. <i>European Journal of Inorganic Chemistry</i> , 2010, 2010, 2044-2053.	2.0	7
41	Synthesis, structure and cytotoxicity studies of four-coordinate bis (cis- η^6 -bis(diphenylphosphino)ethene) gold(I) complexes, [Au(dppe) $_2$] $_2$ X. <i>Journal of Inorganic Biochemistry</i> , 2010, 104, 625-631.	3.5	14
42	Synthesis of N-Propargyl Iminosugar Scaffolds for Compound Library Generation using Click Chemistry. <i>Australian Journal of Chemistry</i> , 2010, 63, 821.	0.9	10
43	Bis(cis-bis(diphenylphosphino)ethene)copper(I) dichloridocuprate(I). <i>Acta Crystallographica Section E: Structure Reports Online</i> , 2010, 66, m493-m494.	0.2	5
44	(η^5 -Pentamethylcyclopentadienyl)(η^6 -4-phenylbutan-2-one)ruthenium(II) tetraphenylborate. <i>Acta Crystallographica Section E: Structure Reports Online</i> , 2010, 66, m1564-m1564.	0.2	0
45	N-Benzyl-2-(3-chloro-4-hydroxyphenyl)acetamide. <i>Acta Crystallographica Section E: Structure Reports Online</i> , 2010, 66, o2521-o2521.	0.2	3
46	Synthesis, Structure, and Selective Cytotoxicity of Organometallic Cp*Ru O-Alkyl-N-phenylcarbamate Sandwich Complexes. <i>Australian Journal of Chemistry</i> , 2010, 63, 245.	0.9	13
47	Solid-State ^{65}Cu and ^{31}P NMR Spectroscopy of Bis(triphenylphosphine) Copper Species. <i>Journal of Physical Chemistry C</i> , 2010, 114, 7949-7962.	3.1	21
48	Pestalactams $\text{A}\hat{=}\text{C}$: novel caprolactams from the endophytic fungus <i>Pestalotiopsis</i> sp.. <i>Organic and Biomolecular Chemistry</i> , 2010, 8, 1785.	2.8	48
49	Synthesis, Spectroscopic Characterization, and Cytotoxic Evaluation of Pentasubstituted Ruthenocenyl Esters. <i>Organometallics</i> , 2010, 29, 6237-6244.	2.3	17
50	2-Propynyl 2-hydroxybenzoate. <i>Acta Crystallographica Section E: Structure Reports Online</i> , 2010, 66, o226-o227.	0.2	2
51	(η^6 -Benzophenone)(η^5 -pentamethylcyclopentadienyl)ruthenium(II) tetraphenylborate. <i>Acta Crystallographica Section E: Structure Reports Online</i> , 2009, 65, m1486-m1486.	0.2	2
52	Novel organometallic cationic ruthenium(II) pentamethylcyclopentadienyl benzenesulfonamide complexes targeted to inhibit carbonic anhydrase. <i>Journal of Biological Inorganic Chemistry</i> , 2009, 14, 935-945.	2.6	33
53	Isolation, structure elucidation and cytotoxic evaluation of endiandrin B from the Australian rainforest plant <i>Endiandra anthropogorum</i> . <i>Bioorganic and Medicinal Chemistry</i> , 2009, 17, 1387-1392.	3.0	26
54	Glycogen phosphorylase inhibitory effects of 2-oxo-1,2-dihydropyridin-3-yl amide derivatives. <i>Bioorganic and Medicinal Chemistry</i> , 2009, 17, 4724-4733.	3.0	9

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55	1 : 2 Adducts of copper(i) halides with 1,2-bis(di-2-pyridylphosphino)ethane: solid state and solution structural studies and antitumour activity. Dalton Transactions, 2009, , 10861.	3.3	28
56	Bis[<i>cis</i> -bis(diphenylphosphino)ethene- λ^2 P, λ^2]copper(I) tetrafluoridoborate ethanol solvate. Acta Crystallographica Section E: Structure Reports Online, 2009, 65, m500-m501.	0.2	1
57	4-(Benzylideneamino)benzenesulfonamide. Acta Crystallographica Section E: Structure Reports Online, 2009, 65, o2087-o2087.	0.2	11
58	Structural and NMR Studies of the Hexameric Copper(I) Complex with <i>cis</i> -N-butyl- λ^2 -ethoxycarbonyl- λ^2 -thioureate. Zeitschrift Fur Anorganische Und Allgemeine Chemie, 2008, 634, 1537-1541.	1.2	11
59	Evaluation of novel Hyphodermin derivatives as Glycogen Phosphorylase a inhibitors. Bioorganic and Medicinal Chemistry, 2008, 16, 6172-6178.	3.0	12
60	The isolation, structure determination and cytotoxicity of the new fungal metabolite, trichodermamide C. Bioorganic and Medicinal Chemistry Letters, 2008, 18, 2836-2839.	2.2	46
61	Structural forms in complexes of 2,9-dimethyl-1,10-phenanthroline with simple salts of copper(I) and other univalent "closed shell" species. Inorganica Chimica Acta, 2008, 361, 2365-2374.	2.4	11
62	Selective Cytotoxic Ru(II) Arene Cp* Complex Salts [R-PhRuCp*] ⁺ X ⁻ for X = BF ₄ ⁻ , PF ₆ ⁻ , and BPh ₄ ⁻ . Inorganic Chemistry, 2008, 47, 8589-8591.	4.0	60
63	Total Synthesis of (λ^2)-Hyphodermins A and D. Journal of Organic Chemistry, 2008, 73, 3435-3440.	3.2	9
64	Structural, 197Au Mössbauer and solid state 31P CP/MAS NMR studies on bis (cis-bis(diphenylphosphino)ethylene) gold(i) complexes [Au(dppey)2]X for X = PF ₆ , I. Dalton Transactions, 2008, , 3723.	3.3	13
65	Diammonium biphenyl-4,4'-disulfonate. Acta Crystallographica Section E: Structure Reports Online, 2008, 64, o47-o47.	0.2	1
66	Bis(pyridine- λ^2 -N)bis(triphenylphosphine- λ^2 -P)copper(I) tetrafluoridoborate. Acta Crystallographica Section E: Structure Reports Online, 2008, 64, m607-m607.	0.2	2
67	2-(3-Chloro-4-hydroxyphenyl)-N-(3,4-dimethoxyphenethyl)acetamide. Acta Crystallographica Section E: Structure Reports Online, 2008, 64, o1057-o1057.	0.2	2
68	(λ^5 -Pentamethylcyclopentadienyl)(λ^6 -p-toluenesulfonamide)ruthenium(II) tetraphenylborate. Acta Crystallographica Section E: Structure Reports Online, 2008, 64, m1568-m1568.	0.2	2
69	N-Cinnamoyl-L-phenylalanine methyl ester. Acta Crystallographica Section E: Structure Reports Online, 2008, 64, o139-o139.	0.2	0
70	1,4-Bis(piperidin-1-ylcarbonyl)benzene. Acta Crystallographica Section E: Structure Reports Online, 2008, 64, o31-o31.	0.2	0
71	λ^2 -Azido-5-cholestene. Acta Crystallographica Section E: Structure Reports Online, 2008, 64, o1738-o1738.	0.2	2
72	<i>cis</i> -tert-Butyl 2-(3-acetylamino-2-oxo-1,2-dihydro-1-pyridyl)acetate. Acta Crystallographica Section E: Structure Reports Online, 2008, 64, o2492-o2493.	0.2	1

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91	Mixed 3,5-dinitrosalicylate(2 σ) and partial 3,5-dinitrosalicylate(1 σ) and picrate(1 σ) species in tris(piperidinium) bis(3,5-dinitrosalicylate) picrate monohydrate. Acta Crystallographica Section E: Structure Reports Online, 2006, 62, o610-o613.	0.2	4
92	(Δ)-1-Hydroxy-6,6-dimethyl-1H,6H-naphtho[1,2-c]furan-3,9-trione. Acta Crystallographica Section E: Structure Reports Online, 2006, 62, o1440-o1442.	0.2	1
93	A dianionic 5-sulfonatosalicylate species in the proton-transfer compound bis(benzylaminium) 3-carboxylato-4-hydroxybenzenesulfonate. Acta Crystallographica Section E: Structure Reports Online, 2006, 62, o1863-o1865.	0.2	3
94	3-Methoxyanilinium 3-carboxy-4-hydroxybenzenesulfonate dihydrate. Acta Crystallographica Section E: Structure Reports Online, 2006, 62, o2313-o2315.	0.2	1
95	Ethyl (3-nitro-2-oxo-1,2-dihydropyridin-1-yl)acetate. Acta Crystallographica Section E: Structure Reports Online, 2006, 62, o2714-o2716.	0.2	1
96	Ethylenediaminium pyridine-2,5-dicarboxylate dihydrate. Acta Crystallographica Section E: Structure Reports Online, 2006, 62, o3124-o3126.	0.2	1
97	2-Phenylquinoline 1-oxide. Acta Crystallographica Section E: Structure Reports Online, 2006, 62, o3805-o3807.	0.2	1
98	The 1:1 proton-transfer compound of 8-quinolinol (oxine) with pyrazine-2,3-dicarboxylic acid: 8-hydroxyquinolinium 3-carboxypyrazine-2-carboxylate dihydrate. Acta Crystallographica Section E: Structure Reports Online, 2006, 62, o5089-o5091.	0.2	4
99	Anhydrous quinolinium 2,4,6-trinitrobenzenesulfonate. Acta Crystallographica Section E: Structure Reports Online, 2006, 62, o5510-o5512.	0.2	1
100	Strychnine-8-ammonio-2-naphthalenesulfonate-water (1/1/3.5): The first structure of a strychnine or brucine compound with a zwitterionic species. Journal of Chemical Crystallography, 2006, 36, 805-811.	1.1	3
101	Hydrogen bonding in proton-transfer compounds of 5-sulfosalicylic acid with ortho-substituted monocyclic heteroaromatic Lewis bases. Journal of Chemical Crystallography, 2006, 36, 841-849.	1.1	20
102	Molecular Recognition in Proton-Transfer Compounds of Brucine with Achiral Substituted Salicylic Acid Analogues. Australian Journal of Chemistry, 2006, 59, 320.	0.9	19
103	The isolation and synthesis of 3-chloro-4-hydroxyphenylacetamide produced by a plant-associated microfungus of the genus Xylaria. Tetrahedron Letters, 2005, 46, 919-921.	1.4	30
104	Synthesis and structure-activity relationships of novel benzene sulfonamides with potent binding affinity for bovine carbonic anhydrase II. Bioorganic and Medicinal Chemistry Letters, 2005, 15, 5429-5433.	2.2	17
105	Layered structures in proton-transfer compounds of 5-sulfosalicylic acid with the aromatic polyamines 2,6-diaminopyridine and 1,4-phenylenediamine. Acta Crystallographica Section C: Crystal Structure Communications, 2005, 61, o555-o558.	0.4	9
106	Aurophilic and hydrogen-bonding interactions in [1-(diphenylphosphino)-4-(diphenylphosphinoyl)butane- η^2 P]iodogold(I) monohydrate. Acta Crystallographica Section E: Structure Reports Online, 2005, 61, m30-m32.	0.2	0
107	N-(2-Nitrophenylsulfonyl)glycine methyl ester. Acta Crystallographica Section E: Structure Reports Online, 2005, 61, o323-o325.	0.2	1
108	Methyl 2,3,4-tri-O-acetyl-1-azido-1-deoxy- β -D-glucopyranuronate at room temperature. Acta Crystallographica Section E: Structure Reports Online, 2005, 61, o738-o740.	0.2	1

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109	Cytosinium 3,5-dinitrosalicylate. <i>Acta Crystallographica Section E: Structure Reports Online</i> , 2005, 61, o746-o748.	0.2	4
110	N,N-É²-Ethylenedisuccinimide. <i>Acta Crystallographica Section E: Structure Reports Online</i> , 2005, 61, o1072-o1073.	0.2	0
111	Succinimido 4-(N-maleimidomethyl)cyclohexanecarboxylate. <i>Acta Crystallographica Section E: Structure Reports Online</i> , 2005, 61, o1203-o1204.	0.2	1
112	Tetrakis[4-(trifluoromethyl)phenyl]stannane. <i>Acta Crystallographica Section E: Structure Reports Online</i> , 2005, 61, m956-m957.	0.2	2
113	(Z,Z)-3,6-Bis[(3,4,5-trimethoxyphenyl)methylene]piperazine-2,5-dione. <i>Acta Crystallographica Section E: Structure Reports Online</i> , 2005, 61, o1443-o1445.	0.2	0
114	N-(2-Nitrobenzenesulfonyl)-L-alanine methyl ester. <i>Acta Crystallographica Section E: Structure Reports Online</i> , 2005, 61, o1665-o1667.	0.2	1
115	1-(4-Bromophenyl)-2-(2-propenylsulfonyl)ethanone. <i>Acta Crystallographica Section E: Structure Reports Online</i> , 2005, 61, o1668-o1669.	0.2	0
116	Brucinium 2-carboxy-6-nitrophthalate dihydrate: the 1:1 proton-transfer compound of brucine with 3-nitrophthalic acid. <i>Acta Crystallographica Section E: Structure Reports Online</i> , 2005, 61, o2008-o2011.	0.2	5
117	Brucinium toluene-4-sulfonate trihydrate at 130 K. <i>Acta Crystallographica Section E: Structure Reports Online</i> , 2005, 61, o2646-o2648.	0.2	12
118	Methyl 3,4-O-dibenzoyl-2-O-benzyl-1-thio-Î²-L-fucopyranoside. <i>Acta Crystallographica Section E: Structure Reports Online</i> , 2005, 61, o2899-o2901.	0.2	0
119	4-(2-Thienyl)-1H-pyrrole-2-carbaldehyde. <i>Acta Crystallographica Section E: Structure Reports Online</i> , 2005, 61, o3401-o3402.	0.2	2
120	Diversion from Bicyclo[4.2.0]octanol Formation Through the Use of Vinyl Electrophiles. <i>Synthesis</i> , 2005, 2005, 2220-2226.	2.3	3
121	5-Nitrosalicylic Acid and its Proton-Transfer Compounds with Aliphatic Lewis Bases. <i>Australian Journal of Chemistry</i> , 2005, 58, 47.	0.9	58
122	Structural and solid state 31P NMR studies of the four-coordinate copper(i) complexes [Cu(PPh3)3X] and [Cu(PPh3)3(CH3CN)]X. <i>Dalton Transactions</i> , 2005, , 2547.	3.3	32
123	Hydrogen bonding in proton-transfer compounds of 8-quinolinol (oxine) with aromatic sulfonic acids. <i>Acta Crystallographica Section C: Crystal Structure Communications</i> , 2004, 60, o600-o603.	0.4	1
124	1,2-Bis(di-4-pyridylphosphino)ethane (d4pype). <i>Acta Crystallographica Section E: Structure Reports Online</i> , 2004, 60, o43-o44.	0.2	0
125	4-Amino-2,6-dichloro-5-nitropyrimidine. <i>Acta Crystallographica Section E: Structure Reports Online</i> , 2004, 60, o241-o243.	0.2	0
126	N-Allyl-N-(2-nitrobenzenesulfonyl)-L-leucine methyl ester. <i>Acta Crystallographica Section E: Structure Reports Online</i> , 2004, 60, o383-o385.	0.2	1

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127	Dicyclohexylammonium 4-nitroanthranilate. Acta Crystallographica Section E: Structure Reports Online, 2004, 60, o684-o686.	0.2	5
128	Bis(guanidinium) 5-sulfosalicylate monohydrate. Acta Crystallographica Section E: Structure Reports Online, 2004, 60, o687-o689.	0.2	3
129	The 1/1/1 adduct hydrate of 8-hydroxy-7-iodoquinoline-5-sulfonic acid (ferron) with urea. Acta Crystallographica Section E: Structure Reports Online, 2004, 60, o1040-o1042.	0.2	0
130	Polymeric anhydrous sodium 2-aminobenzenesulfonate. Acta Crystallographica Section E: Structure Reports Online, 2004, 60, m836-m838.	0.2	2
131	(1 <i>RS</i> ,2 <i>RS</i> ,6 <i>SR</i> ,8 <i>SR</i>)-2-Methyl-8-(phenylsulfonyl)bicyclo[4.2.0]octan-1-ol. Acta Crystallographica Section E: Structure Reports Online, 2004, 60, o1151-o1153.	0.2	0
132	(1 <i>RS</i> ,2 <i>SR</i> ,6 <i>SR</i> ,8 <i>SR</i>)-2,6-Dimethyl-8-(phenylsulfonyl)bicyclo[4.2.0]octan-1-ol. Acta Crystallographica Section E: Structure Reports Online, 2004, 60, o1154-o1156.	0.2	0
133	Quinolin-8-aminium toluene-4-sulfonate. Acta Crystallographica Section E: Structure Reports Online, 2004, 60, o1257-o1259.	0.2	1
134	4-Amino-2-chloro-5-nitro-6-(propylamino)pyrimidine. Acta Crystallographica Section E: Structure Reports Online, 2004, 60, o1260-o1262.	0.2	1
135	Adenosinium 3,5-dinitrosalicylate. Acta Crystallographica Section E: Structure Reports Online, 2004, 60, o1573-o1576.	0.2	0
136	A second crystal polymorph of anilinium picrate. Acta Crystallographica Section E: Structure Reports Online, 2004, 60, o1800-o1803.	0.2	22
137	Bis(2-pyrimidinyl) disulfide dihydrate: a redetermination. Acta Crystallographica Section E: Structure Reports Online, 2004, 60, o2438-o2440.	0.2	4
138	Xanthones from a microfungus of the genus Xylaria. Phytochemistry, 2004, 65, 2373-2378.	2.9	72
139	Probing the Formation of Bicyclo[4.2.0]octan-1-ols. Journal of Organic Chemistry, 2004, 69, 5690-5698.	3.2	4
140	Evidence for Au(I)⋯Au(I) Interactions in a Sterically Congested Environment: Two-Coordinate Gold(I) Halide Phosphine Complexes. Australian Journal of Chemistry, 2004, 57, 213.	0.9	26
141	1-(Diphenylphosphino)-4-(diphenylphosphinoyl)butane. Acta Crystallographica Section E: Structure Reports Online, 2003, 59, o596-o598.	0.2	3
142	Methyl 2,5-dihydroxybenzoate. Acta Crystallographica Section E: Structure Reports Online, 2003, 59, o630-o631.	0.2	0
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