

David R Manke

List of Publications by Year in descending order

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80
papers

906
citations

623574

14
h-index

477173

29
g-index

82
all docs

82
docs citations

82
times ranked

1057
citing authors

#	ARTICLE	IF	CITATIONS
1	Molecular Chemistry of Consequence to Renewable Energy. <i>Inorganic Chemistry</i> , 2005, 44, 6879-6892.	1.9	200
2	Oxygen and hydrogen photocatalysis by two-electron mixed-valence coordination compounds. <i>Coordination Chemistry Reviews</i> , 2005, 249, 1316-1326.	9.5	103
3	Cobalt(II) Coordination Polymer as a Precatalyst for Selective Hydroboration of Aldehydes, Ketones, and Imines. <i>Journal of Organic Chemistry</i> , 2018, 83, 9442-9448.	1.7	79
4	Homochiral Metal-Organic Frameworks Based on Transition Metal Bisphosphonates. <i>Chemistry of Materials</i> , 2002, 14, 3866-3874.	3.2	72
5	Synthesis of functional bisphosphonates via new palladium-catalyzed bis-hydrophosphorylation reactions. <i>Tetrahedron Letters</i> , 2000, 41, 151-154.	0.7	53
6	Eclipsed M ₂ X ₆ Compounds Exhibiting Very Short Metal-Metal Triple Bonds. <i>Inorganic Chemistry</i> , 2004, 43, 3618-3624.	1.9	34
7	Bis(alkylamido)phenylborane Complexes of Zirconium, Hafnium, and Vanadium. <i>Inorganic Chemistry</i> , 2003, 42, 4431-4436.	1.9	31
8	Carbon dioxide reduction by early metal compounds: A propensity for oxygen atom transfer. <i>Inorganica Chimica Acta</i> , 2011, 369, 203-214.	1.2	24
9	Active Metabolite of Aeruginascin (4-Hydroxy- <i>N,N,N</i> -trimethyltryptamine): Synthesis, Structure, and Serotonergic Binding Affinity. <i>ACS Omega</i> , 2020, 5, 16940-16943.	1.6	21
10	Bis(4-acetoxy- <i>N,N</i> -dimethyltryptammonium) fumarate: a new crystalline form of psilocetin, an alternative to psilocybin as a psilocin prodrug. <i>Acta Crystallographica Section E: Crystallographic Communications</i> , 2019, 75, 900-902.	0.2	18
11	Titanium bis(alkylamido)phenylborane complexes. <i>Inorganica Chimica Acta</i> , 2003, 345, 235-240.	1.2	16
12	Genetic Survey of <i>Psilocybe</i> Natural Products. <i>ChemBioChem</i> , 2022, 23, .	1.3	16
13	Aryl Oxazoline Chelates of First-Row Transition Metals: Structures of $\{[\text{C}(\text{N}(\text{O})\text{C}_6\text{H}_4)_2\text{CMe}_2(\text{COCH}_2\text{CMe}_2)]\text{FeCl}(\text{p})$ and		

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19	Bis(4-hydroxy- <i>N</i> -isopropyl- <i>N</i> -methyltryptammonium) fumarate: a new crystalline form of miprocin. <i>Acta Crystallographica Section E: Crystallographic Communications</i> , 2020, 76, 514-517.	0.2	10
20	Synthesis, Structural Characterization, and Pharmacological Activity of Novel Quaternary Salts of 4-Substituted Tryptamines. <i>ACS Omega</i> , 2022, 7, 24888-24894.	1.6	10
21	The first-row transition-metal series of tris(ethylenediamine) diacetate complexes [<i>M</i> (en) ₃](OAc) ₂ (<i>M</i> is Mn, Fe, Co, Ni, Cu, and Zn). <i>Acta Crystallographica Section C, Structural Chemistry</i> , 2017, 73, 442-446.	0.2	8
22	The fumarate salts of the <i>N</i> -isopropyl- <i>N</i> -methyl derivatives of DMT and psilocin. <i>Acta Crystallographica Section E: Crystallographic Communications</i> , 2019, 75, 1316-1320.	0.2	8
23	6-Bromo-1H-indole-2,3-dione hemihydrate. <i>IUCrData</i> , 2016, 1, .	0.1	7
24	First-row transition metalâ€“pyridine (py)â€“sulfate [(py) _x <i>M</i>](SO ₄) complexes (<i>M</i> = Ni, Cu and Zn): crystal field theory in action. <i>Acta Crystallographica Section C, Structural Chemistry</i> , 2018, 74, 263-268.	0.2	6
25	Psilacetin derivatives: fumarate salts of the methylâ€“ethyl, methylâ€“allyl and diallyl variants of the psilocin prodrug. <i>Acta Crystallographica Section E: Crystallographic Communications</i> , 2021, 77, 101-106.	0.2	6
26	Mononuclear, dinuclear and polymeric cobalt(II) complexes built on 4-aryl-2,6-bis(2- π -pyrazinyl)pyridines. <i>Polyhedron</i> , 2017, 132, 64-69.	1.0	5
27	A tripodal aminothioether ligand scaffold: Synthesis and coordination to zirconium and hafnium. <i>Polyhedron</i> , 2017, 121, 264-268.	1.0	5
28	2,5-Dimethylbufotenine and 2,5-dimethylbufotenidine: novel derivatives of natural tryptamines found in <i>Bufo alvarius</i> toads. <i>Acta Crystallographica Section E: Crystallographic Communications</i> , 2021, 77, 190-194.	0.2	5
29	4-Bromo-1H-indole-2,3-dione. <i>IUCrData</i> , 2016, 1, .	0.1	5
30	The crystal structures of iron and cobalt pyridine (py)â€“sulfates, [Fe(SO ₄)(py) ₄] _n and [Co ₃ (SO ₄) ₃ (py) ₁₁] _n . <i>Acta Crystallographica Section E: Crystallographic Communications</i> , 2018, 74, 857-861.	0.2	5
31	The crystal structure of baeocystin. <i>Acta Crystallographica Section E: Crystallographic Communications</i> , 2022, 78, 550-553.	0.2	5
32	Facile synthesis of new divergent imidazole-containing ligands for a 1-D cobalt(II) coordination polymer. <i>Polyhedron</i> , 2017, 127, 355-360.	1.0	4
33	Crystal structure of 2,6-dimethyl-4-pyridone hemihydrate. <i>Acta Crystallographica Section E: Crystallographic Communications</i> , 2015, 71, o533-o533.	0.2	4
34	DMT analogues: <i>N</i> -ethyl- <i>N</i> -propyltryptamine and <i>N</i> -allyl- <i>N</i> -methyltryptamine as their hydrofumarate salts. <i>Acta Crystallographica Section E: Crystallographic Communications</i> , 2020, 76, 1201-1205.	0.2	4
35	5-Bromo-1H-indole-2,3-dione. <i>IUCrData</i> , 2016, 1, .	0.1	4
36	4-Iodo-1H-indole-2,3-dione. <i>IUCrData</i> , 2016, 1, .	0.1	4

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37	5-MeO-DALT: the freebase of <i>N,N</i> -diallyl-5-methoxytryptamine. IUCrData, 2020, 5, .	0.1	4
38	Designing crystalline, flexible covalent metal-organic networks through controlled ligand deprotection. CrystEngComm, 2019, 21, 4255-4257.	1.3	3
39	Quaternary tryptammonium salts: <i>N,N</i> -dimethyl- <i>N</i> -propyltryptammonium (DMPT) iodide and <i>N</i> -allyl- <i>N,N</i> -dimethyltryptammonium (DMALT) iodide. Acta Crystallographica Section E: Crystallographic Communications, 2020, 76, 1357-1360.	0.2	3
40	The hydrochloride salt of 4-hydroxy- <i>N,N</i> -di- <i>n</i> -propyltryptamine (4-HO-DPT). IUCrData, 2020, 5, .	0.1	3
41	7-Bromo-1H-indole-2,3-dione. IUCrData, 2016, 1, .	0.1	3
42	4-Chloro-1H-indole-2,3-dione. IUCrData, 2016, 1, .	0.1	3
43	More crystal field theory in action: the metal-4-picoline (pic)-sulfate [<i>M</i> (pic) ₄ SO ₄] ₄ complexes (<i>M</i> = Fe, Co, Ni, Cu, Zn, and Cd). Acta Crystallographica Section C, Structural Chemistry, 2019, 75, 568-574.	0.2	3
44	Crystal structure of serotonin. Acta Crystallographica Section E: Crystallographic Communications, 2022, 78, 365-368.	0.2	3
45	Crystal structure of 4-amino-2,6-dichlorophenol. Acta Crystallographica Section E: Crystallographic Communications, 2015, 71, o406-o406.	0.2	2
46	A modular synthesis of tris(aryl)tren ligands: Synthesis, structure and lithiation chemistry. Inorganica Chimica Acta, 2017, 461, 71-77.	1.2	2
47	Bufotenidinium iodide. IUCrData, 2021, 6, .	0.1	2
48	'Foxtrot' fumarate: a water-soluble salt of <i>N,N</i> -diallyl-5-methoxytryptamine (5-MeO-DALT). Acta Crystallographica Section E: Crystallographic Communications, 2021, 77, 416-419.	0.2	2
49	The crystalline forms of nine hydrochloride salts of substituted tryptamines. Acta Crystallographica Section C, Structural Chemistry, 2021, 77, 615-620.	0.2	2
50	Crystal structure of 3,4-dimethoxyphenol. Acta Crystallographica Section E: Crystallographic Communications, 2015, 71, o1019-o1019.	0.2	2
51	7-Iodo-1H-indole-2,3-dione. IUCrData, 2016, 1, .	0.1	2
52	4,6-Dichloro-1H-indole-2,3-dione. IUCrData, 2016, 1, .	0.1	2
53	The varied structures of cobalt(II)-pyridine (py)-sulfate: [Co(SO ₄)(py) ₄] ₂ , [Co ₂ (SO ₄) ₂ (py) ₆] ₂ , and [Co ₃ (SO ₄) ₃ (py) ₁₁] ₂ . Acta Crystallographica Section E: Crystallographic Communications, 2019, 75, 1888-1891.	0.2	2
54	A cuboidal [Cu ₄ (SO ₄) ₄] ₄ structure supported by $\hat{1}^2$ -picoline ligands. Acta Crystallographica Section E: Crystallographic Communications, 2022, 78, 108-110.	0.2	2

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55	Dimethyl 2,2-dinitrophenyl-4,4-dicarboxylate. Acta Crystallographica Section E: Structure Reports Online, 2014, 70, o305-o305.	0.2	1
56	5-Methoxy-N,N-di-n-propyltryptamine (5-MeO-DPT): freebase and fumarate. Acta Crystallographica Section E: Crystallographic Communications, 2021, 77, 522-526.	0.2	1
57	Di-1/4-chlorido-bis[(2,2-bipyridine-1,2,6-triazole)nickel(II)]. IUCrData, 2016, 1, o.1		1
58	2-(Phenylsulfanyl)aniline. IUCrData, 2016, 1, .	0.1	1
59	6-Fluoro-1H-indole-2,3-dione. IUCrData, 2016, 1, .	0.1	1
60	6-Iodo-1H-indole-2,3-dione. IUCrData, 2016, 1, .	0.1	1
61	6-Chloro-1H-indole-2,3-dione. IUCrData, 2016, 1, .	0.1	1
62	4,7-Dichloro-1H-indole-2,3-dione. IUCrData, 2016, 1, .	0.1	1
63	5,7-Dichloro-1H-indole-2,3-dione. IUCrData, 2016, 1, .	0.1	1
64	7-Methyl-1H-indole-2,3-dione. IUCrData, 2017, 2, .	0.1	1
65	Receptor binding profiles and behavioral effects of psilocybin analogs. FASEB Journal, 2022, 36, .	0.2	1
66	Dimethyl 2-aminobiphenyl-4,4-dicarboxylate. Acta Crystallographica Section E: Structure Reports Online, 2013, 69, o797-o797.	0.2	0
67	Dimethyl 2-nitrobiphenyl-4,4-dicarboxylate. Acta Crystallographica Section E: Structure Reports Online, 2014, 70, o371-o371.	0.2	0
68	Crystal structure of dimethyl 9H-carbazole-2,7-dicarboxylate. Acta Crystallographica Section E: Crystallographic Communications, 2015, 71, o784-o785.	0.2	0
69	4,4-Oxybis(2,6-dimethylpyridinium) bis(trifluoromethanesulfonate). Acta Crystallographica Section E: Structure Reports Online, 2013, 69, o1633-o1634.	0.2	0
70	Crystal structure of N,N,N-tris[(1,3-benzothiazol-2-yl)methyl]amine. Acta Crystallographica Section E: Crystallographic Communications, 2015, 71, o786-o787.	0.2	0
71	Crystal structure of 3-(diethylamino)phenol. Acta Crystallographica Section E: Crystallographic Communications, 2015, 71, o1075-o1075.	0.2	0
72	Bis[1/4-2-(phenylsulfanyl)anilido-1,2-N]bis[bis(tetrahydrofuran-1,3-dioxane)lithium]. IUCrData, 2016, 1, .	0.1	0

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73	2-Bromo-5-methylpyridine. IUCrData, 2016, 1, .	0.1	0
74	4-Methylcyclohexylp-toluenesulfonate. IUCrData, 2016, 1, .	0.1	0
75	1-(3,4-Dimethoxyphenyl)ethanone. IUCrData, 2016, 1, .	0.1	0
76	5,7-Dimethyl-1 <i>H</i> -indole-2,3-dione. IUCrData, 2017, 2, .	0.1	0
77	Poly[$\text{diaqua}(\frac{1}{4})\text{-}2,3\text{-dioxindoline-5-sulfonato}$]sodium]. IUCrData, 2017, 2, .	0.1	0
78	5,6-Dibromo-1 <i>H</i> -indole-2,3-dione. IUCrData, 2017, 2, .	0.1	0
79	Bis(4-hydroxy- <i>N,N</i> -di- <i>n</i> -propyltryptammonium) fumarate tetrahydrate. IUCrData, 2019, 4, .	0.1	0
80	Bis(oxotremorine) fumarate bis(fumaric acid). IUCrData, 2022, 7, .	0.1	0