

Lee Brammer

List of Publications by Year in descending order

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#	ARTICLE	IF	CITATIONS
1	Multi-stimulus linear negative expansion of a breathing $M(O_2CR)_4$ -node MOF. Faraday Discussions, 2021, 225, 133-151.	3.2	2
2	Post-Synthetic Modification Unlocks a 2D-to-3D Switch in MOF Breathing Response: A Single-Crystal Diffraction Mapping Study. Angewandte Chemie, 2021, 133, 18064-18068.	2.0	1
3	Post-Synthetic Modification Unlocks a 2D-to-3D Switch in MOF Breathing Response: A Single-Crystal Diffraction Mapping Study. Angewandte Chemie - International Edition, 2021, 60, 17920-17924.	13.8	13
4	Increasing Alkyl Chain Length in a Series of Layered Metal-Organic Frameworks Aids Ultrasonic Exfoliation to Form Nanosheets. Inorganic Chemistry, 2019, 58, 10837-10845.	4.0	23
5	Benchmarking of Halogen Bond Strength in Solution with Nickel Fluorides: Bromine versus Iodine and Perfluoroaryl versus Perfluoroalkyl Donors. Chemistry - A European Journal, 2019, 25, 9237-9241.	3.3	13
6	Encapsulation of Crabtree's Catalyst in Sulfonated MIL-101(Cr): Enhancement of Stability and Selectivity between Competing Reaction Pathways by the MOF Chemical Microenvironment. Angewandte Chemie, 2018, 130, 4622-4627.	2.0	7
7	Encapsulation of Crabtree's Catalyst in Sulfonated MIL-101(Cr): Enhancement of Stability and Selectivity between Competing Reaction Pathways by the MOF Chemical Microenvironment. Angewandte Chemie - International Edition, 2018, 57, 4532-4537.	13.8	52
8	Fe(III) Protoporphyrin IX Encapsulated in a Zinc Metal-Organic Framework Shows Dramatically Enhanced Peroxidatic Activity. Inorganic Chemistry, 2018, 57, 1171-1183.	4.0	15
9	Self-complementary nickel halides enable multifaceted comparisons of intermolecular halogen bonds: fluoride ligands vs. other halides. Chemical Science, 2018, 9, 3767-3781.	7.4	27
10	Cocrystals of spironolactone and griseofulvin based on an in silico screening method. CrystEngComm, 2017, 19, 3592-3599.	2.6	39
11	Hydrogen bonding vs. halogen bonding: the solvent decides. Chemical Science, 2017, 8, 5392-5398.	7.4	176
12	Solvent-switchable continuous-breathing behaviour in a diamondoid metal-organic framework and its influence on CO ₂ versus CH ₄ selectivity. Nature Chemistry, 2017, 9, 882-889.	13.6	293
13	Arene guest selectivity and pore flexibility in a metal-organic framework with semi-fluorinated channel walls. Philosophical Transactions Series A, Mathematical, Physical, and Engineering Sciences, 2017, 375, 20160031.	3.4	5
14	Halogen bonding, chalcogen bonding, pnictogen bonding, tetrel bonding: origins, current status and discussion. Faraday Discussions, 2017, 203, 485-507.	3.2	145
15	Coordination change, lability and hemilability in metal-organic frameworks. Chemical Society Reviews, 2017, 46, 5444-5462.	38.1	216
16	Highly selective detection of Hg ²⁺ and MeHgI by di-pyridin-2-yl-[4-(2-pyridin-4-yl-vinyl)-phenyl]-amine and its zinc coordination polymer. Inorganic Chemistry Frontiers, 2016, 3, 1297-1305.	6.0	56
17	Arene Selectivity by a Flexible Coordination Polymer Host. Chemistry - A European Journal, 2016, 22, 13120-13126.	3.3	17
18	Encapsulation of an organometallic cationic catalyst by direct exchange into an anionic MOF. Chemical Science, 2016, 7, 2037-2050.	7.4	57

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19	Coordination Polymer Flexibility Leads to Polymorphism and Enables a Crystalline Solid–Vapour Reaction: A Multi–technique Mechanistic Study. <i>Chemistry - A European Journal</i> , 2015, 21, 8799-8811.	3.3	25
20	The Contrasting Character of Early and Late Transition Metal Fluorides as Hydrogen Bond Acceptors. <i>Journal of the American Chemical Society</i> , 2015, 137, 11820-11831.	13.7	29
21	Solvent-vapour-assisted pathways and the role of pre-organization in solid-state transformations of coordination polymers. <i>IUCr</i> , 2015, 2, 188-197.	2.2	10
22	Metal Hydrides Form Halogen Bonds: Measurement of Energetics of Binding. <i>Journal of the American Chemical Society</i> , 2014, 136, 1288-1291.	13.7	35
23	Highly fluorinated naphthalenes and bifurcated C–H⋯F⋯C hydrogen bonding. <i>CrystEngComm</i> , 2014, 16, 9711-9720.	2.6	21
24	A solvent-resistant halogen bond. <i>Chemical Science</i> , 2014, 5, 4179-4183.	7.4	122
25	Crystallographic studies of gas sorption in metal–organic frameworks. <i>Acta Crystallographica Section B: Structural Science, Crystal Engineering and Materials</i> , 2014, 70, 404-422.	1.1	79
26	Chemical transformations of a crystalline coordination polymer: a multi-stage solid–vapour reaction manifold. <i>Chemical Science</i> , 2013, 4, 696-708.	7.4	35
27	Zippering and Unzipping of a Paddlewheel Metal–Organic Framework to Enable Two–Step Synthetic and Structural Transformation. <i>Chemistry - A European Journal</i> , 2013, 19, 3552-3557.	3.3	28
28	Persistent C–I⋯ halogen-bonded layer motifs involving 4-iodobenzoate paddlewheel units, Cu ₂ (4- <i>l</i> bz) ₄ (L) ₂ . <i>CrystEngComm</i> , 2013, 15, 3160.	2.6	18
29	Coordination chemistry meets halogen bonding and hydrogen bonding: building networks from 3-iodobenzoate paddlewheel units [Cu ₂ (3- <i>l</i> bz) ₄ (L) ₂]. <i>CrystEngComm</i> , 2013, 15, 3151.	2.6	29
30	One-dimensional organization of free radicals via halogen bonding. <i>CrystEngComm</i> , 2012, 14, 6381.	2.6	30
31	Cyanometallates as Halogen Bond Acceptors. <i>Crystal Growth and Design</i> , 2012, 12, 205-216.	3.0	81
32	Tuning the magneto-structural properties of non-porous coordination polymers by HCl chemisorption. <i>Nature Communications</i> , 2012, 3, 828.	12.8	99
33	Diiodoacetylene: compact, strong ditopic halogen bond donor. <i>CrystEngComm</i> , 2012, 14, 3033.	2.6	60
34	Synthesis and polymorphism of (4-ClpyH) ₂ [CuCl ₄]: solid–gas and solid–solid reactions. <i>CrystEngComm</i> , 2011, 13, 3189-3196.	2.6	38
35	Different structural destinations: comparing reactions of [CuBr ₂ (3-Brpy) ₂] crystals with HBr and HCl gas. <i>CrystEngComm</i> , 2011, 13, 4400.	2.6	22
36	Energetics of Halogen Bonding of Group 10 Metal Fluoride Complexes. <i>Journal of the American Chemical Society</i> , 2011, 133, 14338-14348.	13.7	64

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37	Mechanistic Insights into a Gas-Solid Reaction in Molecular Crystals: The Role of Hydrogen Bonding. <i>Angewandte Chemie - International Edition</i> , 2010, 49, 8892-8896.	13.8	59
38	Effects of halogen bonding in ferromagnetic chains based on Co(ii) coordination polymers. <i>CrystEngComm</i> , 2010, 12, 2339.	2.6	43
39	Rational Modification of the Hierarchy of Intermolecular Interactions in Molecular Crystal Structures by Using Tunable Halogen Bonds. <i>Chemistry - A European Journal</i> , 2009, 15, 7554-7568.	3.3	164
40	Halogen Bonded Supramolecular Assemblies of [Ru(bipy)(CN) ₄] ²⁻ Anions and N-Methyl-Halopyridinium Cations in the Solid State and in Solution. <i>Inorganic Chemistry</i> , 2009, 48, 1666-1677.	4.0	86
41	Ligand Substitution within Nonporous Crystals of a Coordination Polymer: Elimination from and Insertion into Ag-O Bonds by Alcohol Molecules in a Solid-Vapor Reaction. <i>Angewandte Chemie - International Edition</i> , 2008, 47, 1693-1697.	13.8	65
42	Noncovalent Interactions under Extreme Conditions: High-Pressure and Low-Temperature Diffraction Studies of the Isostructural Metal-Organic Networks (4-Chloropyridinium) ₂ [CoX ₄] (X = Cl, Br). <i>Journal of the American Chemical Society</i> , 2008, 130, 9058-9071.	13.7	82
43	Competition between coordination network and halogen bond network formation: towards halogen-bond functionalised network materials using copper-iodobenzoate units. <i>CrystEngComm</i> , 2008, 10, 1335.	2.6	34
44	Combining metals with halogen bonds. <i>CrystEngComm</i> , 2008, 10, 1712.	2.6	300
45	Metal Fluorides Form Strong Hydrogen Bonds and Halogen Bonds: Measuring Interaction Enthalpies and Entropies in Solution. <i>Journal of the American Chemical Society</i> , 2008, 130, 7842-7844.	13.7	143
46	Ligand flexibility and framework rearrangement in a new family of porous metal-organic frameworks. <i>Chemical Communications</i> , 2007, , 1532-1534.	4.1	73
47	Reversible Gas Uptake by a Nonporous Crystalline Solid Involving Multiple Changes in Covalent Bonding. <i>Journal of the American Chemical Society</i> , 2007, 129, 15606-15614.	13.7	82
48	Metal-Organic Halogen-Bonded Network Formation in MX ₂ (4-halopyridine) ₂ Complexes (M = Pd, Pt; X = Cl, I). <i>Journal of the American Chemical Society</i> , 2007, 129, 15606-15614.	13.7	82
49	Solvent hydrolysis leads to an unusual Cu(ii) metal-organic framework. <i>CrystEngComm</i> , 2006, 8, 473.	2.6	50
50	Unexpected structural homologies involving hydrogen-bonded and halogen-bonded networks in halopyridinium halometallate salts. <i>CrystEngComm</i> , 2006, 8, 425.	2.6	51
51	Designing Intermolecular Interactions between Halogenated Peripheries of Inorganic and Organic Molecules: Electrostatically Directed Metal-Organic Halogen Bonds. <i>Angewandte Chemie - International Edition</i> , 2006, 45, 435-440.	13.8	152
52	New trends in crystal engineering. <i>CrystEngComm</i> , 2005, 7, 1.	2.6	412
53	Supramolecular Chemistry of Halogens: Complementary Features of Inorganic (MX) and Organic (C-X) Halogens Applied to Metal-Organic Halogen Bond Formation. <i>Journal of the American Chemical Society</i> , 2005, 127, 5979-5989.	13.7	365
54	Halometallate and halide ions: nucleophiles in competition for hydrogen bond and halogen bond formation in halopyridinium salts of mixed halide-halometallate anions. <i>CrystEngComm</i> , 2005, 7, 350.	2.6	75

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55	Water molecules insert into Nâ€”H...Clâ€”M hydrogen bonds while Mâ€”Cl...Xâ€”C halogen bonds remain intact in dihydrates of halopyridinium hexachloroplatinates. <i>Acta Crystallographica Section B: Structural Science</i> , 2004, 60, 512-519.	1.8	41
56	Developments in inorganic crystal engineering. <i>Chemical Society Reviews</i> , 2004, 33, 476.	38.1	685
57	Metals and hydrogen bonds. <i>Dalton Transactions</i> , 2003, , 3145.	3.3	359
58	Hydrogen bond patterns in aromatic and aliphatic dioximes. <i>New Journal of Chemistry</i> , 2003, 27, 1084-1094.	2.8	56
59	Involving metals in halogenâ€”halogen interactions: second-sphere Lewis acid ligands for perhalometallate ions (Mâ€”Xâ€”Xâ€”C). <i>CrystEngComm</i> , 2003, 5, 343-345.	2.6	100
60	Binding Studies on the Control of the Conformation and Self-assembly of a Calix[4]arene-dicarboxylic Acid through Hydrogen Bonding Interactions. <i>Supramolecular Chemistry</i> , 2003, 15, 385-390.	1.2	7
61	Hydrogen Bonds in Inorganic Chemistry: Application to Crystal Design. <i>Perspectives in Supramolecular Chemistry</i> , 2003, , 1-75.	0.1	22
62	Designing neutral coordination networks with the aid of hydrogen bond mimicry using silver(i) carboxylates. <i>CrystEngComm</i> , 2002, 4, 239-248.	2.6	54
63	Bridging mode flexibility of 1,3-dithiacyclohexane in silver(i) co-ordination polymers. <i>Dalton Transactions RSC</i> , 2002, , 4134.	2.3	14
64	Hydrogen bonding and perhalometallate ions: A supramolecular synthetic strategy for new inorganic materials. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2002, 99, 4956-4961.	7.1	126
65	Silver(I) carboxylates: versatile inorganic analogs of carboxylic acids for supramolecular network formation. <i>Chemical Communications</i> , 2001, , 2468-2469.	4.1	65
66	Understanding the Behavior of Halogens as Hydrogen Bond Acceptors. <i>Crystal Growth and Design</i> , 2001, 1, 277-290.	3.0	631
67	Synthesis and characterization of sterically hindered diarylsilanes containing 2,4,6-trimethylphenyl and 2,4,6-tris(trifluoromethyl)phenyl substituents. X-ray crystal structure of bis[2,4,6-tris(trifluoromethylphenyl)]fluorosilane. <i>Journal of Organometallic Chemistry</i> , 1995, 499, 89-98.	1.8	24
68	Supplement. Tables of bond lengths determined by X-ray and neutron diffraction. Part 2. Organometallic compounds and co-ordination complexes of the d- and f-block metals. <i>Journal of the Chemical Society Dalton Transactions</i> , 1989, , S1.	1.1	1,165
69	Diffraction Studies in Crystal Engineering. , 0, , 241-265.		0