

Mark A Fox

List of Publications by Year in descending order

Source: <https://exaly.com/author-pdf/2430399/publications.pdf>

Version: 2024-02-01

183
papers

6,762
citations

57681

46
h-index

97045

71
g-index

188
all docs

188
docs citations

188
times ranked

5665
citing authors

#	ARTICLE	IF	CITATIONS
1	Thermodynamic equilibrium between locally excited and charge-transfer states through thermally activated charge transfer in 1-(pyren-2-yl)- <i>o</i> -carborane. <i>Chemical Science</i> , 2022, 13, 5205-5219.	3.7	20
2	Cyclophane Molecules Exhibiting Thermally Activated Delayed Fluorescence: Linking Donor Units to Influence Molecular Conformation. <i>Journal of Organic Chemistry</i> , 2021, 86, 429-445.	1.7	13
3	Carborane photochromism: a fatigue resistant carborane switch. <i>Chemical Communications</i> , 2021, 57, 9466-9469.	2.2	6
4	Vibrational Damping Reveals Vibronic Coupling in Thermally Activated Delayed Fluorescence Materials. <i>Chemistry of Materials</i> , 2021, 33, 3066-3080.	3.2	47
5	Nonlinear optical properties of meso-Tetra(fluorenyl)porphyrins peripherally functionalized with one to four ruthenium alkynyl substituents. <i>Dyes and Pigments</i> , 2021, 188, 109155.	2.0	15
6	Structure and hydration of polyvinylpyrrolidone-hydrogen peroxide. <i>Chemical Communications</i> , 2021, 58, 80-83.	2.2	4
7	Exploiting trifluoromethyl substituents for tuning orbital character of singlet and triplet states to increase the rate of thermally activated delayed fluorescence. <i>Materials Chemistry Frontiers</i> , 2020, 4, 3602-3615.	3.2	35
8	HFO as a CF ₃ -Building Block: Synthesis and Chemistry of CF ₃ -Nones. <i>European Journal of Organic Chemistry</i> , 2020, 2020, 6236-6244.	1.2	14
9	Metallacarborane Assemblies as Effective Antimicrobial Agents, Including a Highly Potent Anti-MRSA Agent. <i>Organometallics</i> , 2020, 39, 4253-4264.	1.1	14
10	Unusual dual-emissive heteroleptic iridium complexes incorporating TADF cyclometalating ligands. <i>Dalton Transactions</i> , 2020, 49, 2190-2208.	1.6	19
11	Fluorenylporphyrins functionalized by electrochromic ruthenium units as redox-triggered fluorescence switches. <i>Dalton Transactions</i> , 2019, 48, 11897-11911.	1.6	5
12	Unravelling the Complexities of Pseudocontact Shift Analysis in Lanthanide Coordination Complexes of Differing Symmetry. <i>Angewandte Chemie - International Edition</i> , 2019, 58, 10290-10294.	7.2	16
13	Unravelling the Complexities of Pseudocontact Shift Analysis in Lanthanide Coordination Complexes of Differing Symmetry. <i>Angewandte Chemie</i> , 2019, 131, 10396-10400.	1.6	7
14	Impact of Methoxy Substituents on Thermally Activated Delayed Fluorescence and Room-Temperature Phosphorescence in All-Organic Donor-Acceptor Systems. <i>Journal of Organic Chemistry</i> , 2019, 84, 3801-3816.	1.7	43
15	Synthetic, Structural, and Computational Studies on Heavier Tetragen and Chalcogen Triazenide Complexes. <i>Inorganic Chemistry</i> , 2019, 58, 16660-16666.	1.9	6
16	Monitoring of the ADP/ATP Ratio by Induced Circularly Polarised Europium Luminescence. <i>Angewandte Chemie - International Edition</i> , 2018, 57, 7488-7492.	7.2	74
17	Enhanced selectivity for Mg ²⁺ with a phosphinate-based chelate: APDAP versus APTRA. <i>Dalton Transactions</i> , 2018, 47, 1879-1887.	1.6	11
18	Monitoring of the ADP/ATP Ratio by Induced Circularly Polarised Europium Luminescence. <i>Angewandte Chemie</i> , 2018, 130, 7610-7614.	1.6	26

#	ARTICLE	IF	CITATIONS
19	Role of the Diphosphine Chelate in Emissive, Charge-Neutral Iridium(III) Complexes. <i>Chemistry - A European Journal</i> , 2018, 24, 624-635.	1.7	12
20	Selective signalling of glyphosate in water using europium luminescence. <i>Dalton Transactions</i> , 2018, 47, 16145-16154.	1.6	18
21	Bond Rotations and Heteroatom Effects in Donor-Acceptor Donor Molecules: Implications for Thermally Activated Delayed Fluorescence and Room Temperature Phosphorescence. <i>Journal of Organic Chemistry</i> , 2018, 83, 14431-14442.	1.7	61
22	Luminescent Pt(II) complexes featuring imidazolylidene-pyridylidene and dianionic bipyrazolate: from fundamentals to OLED fabrications. <i>Journal of Materials Chemistry C</i> , 2017, 5, 1420-1435.	2.7	37
23	New Blatter-type radicals from a bench-stable carbene. <i>Nature Communications</i> , 2017, 8, 15088.	5.8	36
24	Bis-Tridentate Ir(III) Metal Phosphors for Efficient Deep-Blue Organic Light-Emitting Diodes. <i>Advanced Materials</i> , 2017, 29, 1702464.	11.1	117
25	The contributions of molecular vibrations and higher triplet levels to the intersystem crossing mechanism in metal-free organic emitters. <i>Journal of Materials Chemistry C</i> , 2017, 5, 6269-6280.	2.7	83
26	Geometries of 11-Vertex Carborane Monoanion Radicals with 2 <i>n</i> +3 Skeletal Electron Counts. <i>European Journal of Inorganic Chemistry</i> , 2017, 2017, 4568-4574.	1.0	3
27	Pyridylpyrazole N ^N ligands combined with sulfonyl-functionalised cyclometalating ligands for blue-emitting iridium(III) complexes and solution-processable PhOLEDs. <i>Dalton Transactions</i> , 2017, 46, 10996-11007.	1.6	17
28	Bright green PhOLEDs using cyclometalated diiridium(III) complexes with bridging oxamidato ligands as phosphorescent dopants. <i>Journal of Materials Chemistry C</i> , 2017, 5, 6777-6789.	2.7	30
29	Biotransformation of fluorophenyl pyridine carboxylic acids by the model fungus <i>Cunninghamella elegans</i> . <i>Xenobiotica</i> , 2017, 47, 763-770.	0.5	10
30	Platinum(II) complexes of some unsymmetrical diphosphenes. <i>Journal of Organometallic Chemistry</i> , 2017, 830, 113-119.	0.8	2
31	Tethered N-heterocyclic carbene-carboranes: unique ligands that exhibit unprecedented and versatile coordination modes at rhodium. <i>Chemical Communications</i> , 2016, 52, 6443-6446.	2.2	32
32	Induced europium CPL for the selective signalling of phosphorylated amino-acids and O-phosphorylated hexapeptides. <i>Dalton Transactions</i> , 2016, 45, 8355-8366.	1.6	41
33	Sulfonyl-Substituted Heteroleptic Cyclometalated Iridium(III) Complexes as Blue Emitters for Solution-Processable Phosphorescent Organic Light-Emitting Diodes. <i>Inorganic Chemistry</i> , 2016, 55, 8612-8627.	1.9	32
34	Substituent Effects on the Fluorescence Properties of ortho-Carboranes: Unusual Emission Behaviour in C(2-Pyridyl)-ortho-carboranes. <i>European Journal of Inorganic Chemistry</i> , 2016, 2016, 403-412.	1.0	46
35	The Role of Local Triplet Excited States and Relative Orientation in Thermally Activated Delayed Fluorescence: Photophysics and Devices. <i>Advanced Science</i> , 2016, 3, 1600080.	5.6	403
36	Bis-Tridentate Iridium(III) Phosphors Bearing Functional 2-Phenyl-6-(imidazol-2-ylidene)pyridine and 2-(Pyrazol-3-yl)-6-phenylpyridine Chelates for Efficient OLEDs. <i>Organometallics</i> , 2016, 35, 1813-1824.	1.1	63

#	ARTICLE	IF	CITATIONS
37	Low-melting molecular complexes: Part VII. 2,3-, 2,5- and 3,4-hexanediones and their molecular complexes with chloroform. <i>Structural Chemistry</i> , 2016, 27, 9-15.	1.0	1
38	Challenging lanthanide relaxation theory: erbium and thulium complexes that show NMR relaxation rates faster than dysprosium and terbium analogues. <i>Physical Chemistry Chemical Physics</i> , 2015, 17, 16507-16511.	1.3	19
39	Oligo(<i>p</i> -phenyleneethynylene) (OPE) Molecular Wires: Synthesis and Length Dependence of Photoinduced Charge Transfer in OPEs with Triarylamine and Diaryloxadiazole End Groups. <i>Chemistry - A European Journal</i> , 2015, 21, 3997-4007.	1.7	33
40	Syntheses and Structures of Buta-1,3-Diynyl Complexes from σ -Complex Cross-Coupling Reactions. <i>Organometallics</i> , 2015, 34, 2395-2405.	1.1	16
41	Why are the $\{Cu_4N_4\}$ rings in copper phosphinimide clusters $[Cu_{1/4}NiPR_3]_4$ ($R = NMe_3$ or Ph) planar?. <i>Dalton Transactions</i> , 2015, 44, 5611-5619.	1.6	11
42	Chiral probe development for circularly polarised luminescence: comparative study of structural factors determining the degree of induced CPL with four heptacoordinate europium complexes. <i>Dalton Transactions</i> , 2015, 44, 14937-14951.	1.6	33
43	Syntheses and reductions of C-dimesitylboryl-1,2-dicarba-closo-dodecaboranes. <i>Dalton Transactions</i> , 2015, 44, 9766-9781.	1.6	53
44	Near infrared-emitting tris-bidentate Os(ii) phosphors: control of excited state characteristics and fabrication of OLEDs. <i>Journal of Materials Chemistry C</i> , 2015, 3, 4910-4920.	2.7	52
45	New donor-acceptor conjugates based on a trifluorenylporphyrin linked to a redox-switchable ruthenium unit. <i>Dalton Transactions</i> , 2015, 44, 9470-9485.	1.6	16
46	Ir(III)-Based Phosphors with Bipyrazolate Ancillaries; Rational Design, Photophysics, and Applications in Organic Light-Emitting Diodes. <i>Inorganic Chemistry</i> , 2015, 54, 10811-10821.	1.9	36
47	An Experimental and Computational Approach to Understanding the Reactions of Acyl Nitroso Compounds in [4 + 2] Cycloadditions. <i>Journal of Organic Chemistry</i> , 2015, 80, 9518-9534.	1.7	18
48	Crystal Structures of the Carborane Dianions $[1,4-(PhCB_{10}H_{10}C)_2C_6H_4]^{2-}$ and $[1,4-(PhCB_{10}H_{10}C)_2C_6F_4]^{2-}$ and the Stabilizing Role of the <i>para</i> -Phenylene Unit on $2\%+3$ Skeletal Electron Clusters. <i>Angewandte Chemie - International Edition</i> , 2014, 53, 3702-3705.	7.2	28
49	On the ambiguity of 1,3,2-benzodiazaboroles as donor/acceptor functionalities in luminescent molecules. <i>Dalton Transactions</i> , 2014, 43, 3347-3363.	1.6	16
50	Electronic Structure and Charge Transport Properties of a Series of 3,6-(Diphenyl)- <i>s</i> -tetrazine Derivatives: Are They Suitable Candidates for Molecular Electronics?. <i>Journal of Physical Chemistry C</i> , 2014, 118, 26427-26439.	1.5	13
51	Bimetallic Cyclometalated Iridium(III) Diastereomers with Non-Innocent Bridging Ligands for High-Efficiency Phosphorescent OLEDs. <i>Angewandte Chemie - International Edition</i> , 2014, 53, 11616-11619.	7.2	65
52	Ultrafast Dynamics and Computational Studies on Diaminodicyanoquinodimethanes (DADQs). <i>Journal of Physical Chemistry B</i> , 2014, 118, 6815-6828.	1.2	9
53	Mixed-Valence Ruthenium Complexes Rotating through a Conformational Robin-Day Continuum. <i>Chemistry - A European Journal</i> , 2014, 20, 6895-6908.	1.7	76
54	Studies on bis(1-ortho-carboranyl)benzenes and bis(1-ortho-carboranyl)biphenyls. <i>Tetrahedron</i> , 2014, 70, 5182-5189.	1.0	16

#	ARTICLE	IF	CITATIONS
55	A novel, efficient synthesis of N-aryl pyrroles via reaction of 1-boronodienes with aryl nitroso compounds. <i>Chemical Communications</i> , 2013, 49, 5414.	2.2	26
56	Electrochemical and spectroelectrochemical studies of C-benzodiazaborolyl-ortho-carboranes. <i>Dalton Transactions</i> , 2013, 42, 2266-2281.	1.6	87
57	C,Câ€²-Bis(benzodiazaborolyl)dicarba-closo-dodecaboranes: Synthesis, structures, photophysics and electrochemistry. <i>Dalton Transactions</i> , 2013, 42, 10982.	1.6	70
58	¹⁹ F and ¹³ C GIAO-NMR chemical shifts for the identification of perfluoro-quinoline and -isoquinoline derivatives. <i>Journal of Fluorine Chemistry</i> , 2013, 155, 62-71.	0.9	12
59	Remarkable cage deboronation and rearrangement of a closo-1,12-dicarbododecaborane to form a neutral nido-7,9-dicarbaundecaborane. <i>Chemical Communications</i> , 2013, 49, 3312.	2.2	7
60	Some Ruthenium Derivatives of Penta-1,4-diyne-3-one. <i>Organometallics</i> , 2013, 32, 3286-3299.	1.1	37
61	Carboranes as model hypercarbon systems; structural and bonding patterns in selected isoelectronic closo-borane and carborane systems; [B _n H _n] ²⁺ , [1-CBnâ€² ⁺ 1Hn] ⁺ and 1, <i>n</i> -C ₂ Bnâ€² ⁺ 2Hn (<i>n</i> = 5, 6, 7, 10 or 12). <i>Journal of Organometallic Chemistry</i> , 2012, 721-722, 130-136.	0.8	7
62	peri-Dimethylamino substituent effects on proton transfer at carbon in Î±-naphthylacetate esters: a model for mandelate racemase. <i>Organic and Biomolecular Chemistry</i> , 2012, 10, 590-596.	1.5	2
63	Reactions of 4-substituted tetrafluoropyridine derivatives with sulfur nucleophiles: S _N Ar and annelation processes. <i>Journal of Fluorine Chemistry</i> , 2012, 143, 148-154.	0.9	10
64	Diazaborolyl-boryl pushâ€“pull systems with ethynyleneâ€“arylene bridges as â€“turn-onâ€™ fluoride sensors. <i>Dalton Transactions</i> , 2012, 41, 10328.	1.6	23
65	Synthesis and Characterization of Dithia[3.3]paracyclophane-Bridged Binuclear Ruthenium Vinyl and Alkynyl Complexes. <i>Organometallics</i> , 2012, 31, 5321-5333.	1.1	43
66	Substitution of Tetracyanoethene by Ethynylâ€“Metal Complexes Gives Tricyanovinylethynyl (Tricyanobutenynyl) Derivatives: Syntheses, Protonation, and Addition of Metalâ€“Ligand Fragments. <i>Organometallics</i> , 2012, 31, 2639-2657.	1.1	14
67	Phosphineâ€“alkene ligand-mediated alkylâ€“alkyl and alkylâ€“halide elimination processes from palladium(II). <i>Chemical Communications</i> , 2012, 48, 10413.	2.2	12
68	Colour tuning of blue electroluminescence using bipolar carbazoleâ€“oxadiazole molecules in single-active-layer organic light emitting devices (OLEDs). <i>Journal of Materials Chemistry</i> , 2012, 22, 11816.	6.7	79
69	Platinum(II) Complexes of Cyclic Triphosphenium Ions: a ³¹ P NMR Spectroscopic and Computational Study. <i>Inorganic Chemistry</i> , 2012, 51, 9799-9808.	1.9	15
70	Syntheses, structures and redox properties of tris(pyrazolyl)borate-capped ruthenium vinyl complexes. <i>Journal of Organometallic Chemistry</i> , 2012, 721-722, 173-185.	0.8	4
71	Colour tuning from green to red by substituent effects in phosphorescent tris-cyclometalated iridium(III) complexes of carbazole-based ligands: synthetic, photophysical, computational and high efficiency OLED studies. <i>Journal of Materials Chemistry</i> , 2012, 22, 6419.	6.7	96
72	Dinuclear iridium(III) complexes of cyclometalated fluorenylpyridine ligands as phosphorescent dopants for efficient solution-processed OLEDs. <i>Journal of Materials Chemistry</i> , 2012, 22, 13529.	6.7	41

#	ARTICLE	IF	CITATIONS
73	Luminescence Properties of <i>ortho</i> -Carboranes as Donor-Acceptor Systems. <i>Chemistry - A European Journal</i> , 2012, 18, 8347-8357.	1.7	151
74	The synthesis, molecular and electronic structure of cyanovinylidene complexes. <i>Inorganica Chimica Acta</i> , 2012, 380, 358-371.	1.2	14
75	Experimental and Theoretical Studies on Organic π -Systems Containing Three-coordinate Boron Moieties as both Donor and Acceptor. <i>Chemistry - A European Journal</i> , 2012, 18, 1369-1382.	1.7	80
76	Simultaneous Bridge-Localized and Mixed-Valence Character in Diruthenium Radical Cations Featuring Diethynylaromatic Bridging Ligands. <i>Journal of the American Chemical Society</i> , 2011, 133, 18433-18446.	6.6	138
77	Molybdenum Complexes of <i>ortho</i> -Bis(ethynyl)carboranes: Design, Synthesis, and Study of a Weakly Coupled Mixed-Valence Compound. <i>Organometallics</i> , 2011, 30, 884-894.	1.1	29
78	Ligand redox non-innocent behaviour in ruthenium complexes of ethynyl tolans. <i>Inorganica Chimica Acta</i> , 2011, 374, 461-471.	1.2	16
79	Highly Efficient, Solution-Processed, Single-Layer, Electrophosphorescent Diodes and the Effect of Molecular Dipole Moment. <i>Advanced Functional Materials</i> , 2011, 21, 2376-2382.	7.8	66
80	1,2-Carbagerma-closo-dodecaborate as a Germanium Ligand in Coordination Chemistry - Synthesis, Structure and Reactivity. <i>European Journal of Inorganic Chemistry</i> , 2011, 2011, 3349-3356.	1.0	7
81	Luminescent Platinum(II) Complexes Containing Cyclometallated Diaryl Ketimine Ligands: Synthesis, Photophysical and Computational Properties. <i>European Journal of Inorganic Chemistry</i> , 2010, 2010, 1963-1972.	1.0	25
82	Deboronation and Deprotonation of <i>ortho</i> -Carborane with N-Heterocyclic Carbenes. <i>Chemistry - A European Journal</i> , 2010, 16, 10644-10648.	1.7	47
83	Tuning the Intramolecular Charge Transfer Emission from Deep Blue to Green in Ambipolar Systems Based on Dibenzothiophene <i>ortho</i> -Dioxide by Manipulation of Conjugation and Strength of the Electron Donor Units. <i>Journal of Organic Chemistry</i> , 2010, 75, 6771-6781.	1.7	114
84	Some reactions of an η^3 -tetracyanobutadienyl-ruthenium complex. <i>Dalton Transactions</i> , 2010, 39, 3759.	1.6	21
85	Syntheses and molecular structures of some tricobaltcarbonylclusters containing 2,4,6-trimethyl-1,3,5-trithiane. <i>Dalton Transactions</i> , 2010, 39, 1222-1234.	1.6	11
86	The electronic structures of diruthenium complexes containing an oligo(phenylene ethynylene) bridging ligand, and some related molecular structures. <i>Dalton Transactions</i> , 2010, 39, 11605.	1.6	20
87	From Cyclic Iminophosphanes to π -Conjugated Materials. <i>Angewandte Chemie - International Edition</i> , 2009, 48, 9109-9113.	7.2	12
88	Structural, spectroscopic, electrochemical and computational studies of C, π -diaryl- <i>ortho</i> -carboranes, 1-(4-XC ₆ H ₄)-2-Ph-1,2-C ₂ B ₁₀ H ₁₀ (X = H, F, OMe, NMe ₂ , NH ₂ , OH and O ⁻). <i>Journal of Solid State Electrochemistry</i> , 2009, 13, 1483-1495.	1.2	44
89	New synthetic and structural studies on nitroso- <i>ortho</i> -carboranes RCB ₁₀ H ₁₀ CNO and bis(<i>ortho</i> -carboranyl)amines (RCB ₁₀ H ₁₀ C) ₂ NH (R=Ph or Me). <i>Polyhedron</i> , 2009, 28, 789-795.	1.0	16
90	Trends in <i>ortho</i> -carboranes 1-X-2-R-1,2-C ₂ B ₁₀ H ₁₀ (R=Ph, Me) bearing an exo-CN-bonded substituent group (X=NO, NNR ⁻ or NHR ⁻). <i>Polyhedron</i> , 2009, 28, 2359-2370.	1.0	32

#	ARTICLE	IF	CITATIONS
91	A simple synthesis of trans-RuCl(CCR)(dppe) ₂ complexes and representative molecular structures. Journal of Organometallic Chemistry, 2009, 694, 2350-2358.	0.8	67
92	The syntheses and structures of mono- and di-bromovinylidenes. Journal of Organometallic Chemistry, 2009, 694, 4042-4048.	0.8	7
93	Noninnocent Ligand Behavior in Diruthenium Complexes Containing a 1,3-Diethynylbenzene Bridge. Organometallics, 2009, 28, 5266-5269.	1.1	66
94	Transition metal alkynyl complexes by transmetallation from Au(Ci€r,CAr)(PPh₃)₃ (Ar =) Tj ETQqO 0 0 rgBT /Overlock 10 Tf 50 610-620.	1.6	74
95	Synthetic, structural, photophysical and computational studies of i€-conjugated bis- and tris-1,3,2-benzodiazaboroles and related bis(boryl) dithiophenes. Dalton Transactions, 2009, , 1339.	1.6	79
96	Synthetic, structural, photophysical and computational studies on 2-arylethynyl-1,3,2-diazaboroles. Dalton Transactions, 2009, , 2823.	1.6	49
97	DFT studies of the i€-donor/i€-acceptor properties of [SnCB10H11]â€“ and its relationship to [SnCl3]â€“, CO, PF3, [SnB11H11]2â€“, SnC2B9H11, and related SnC2BnHn+2 compounds. Canadian Journal of Chemistry, 2009, 87, 63-71.	0.6	10
98	Experimental and computed dipole moments in donorâ€“bridgeâ€“acceptor systems with p-phenylene and p-carboranediyl bridges. Collection of Czechoslovak Chemical Communications, 2009, 74, 131-146.	1.0	8
99	The preparation and characterisation of ruthenium cyanovinylidene complexes. Dalton Transactions, 2008, , 433-436.	1.6	16
100	Ruthenium Complexes of <i>C,C</i>-Bis(ethynyl)carboranes:â€“ An Investigation of Electronic Interactions Mediated by Spherical Pseudo-aromatic Spacers. Journal of the American Chemical Society, 2008, 130, 3566-3578.	6.6	116
101	Some transition metal complexes derived from mono- and di-ethynyl perfluorobenzenes. Dalton Transactions, 2008, , 6763.	1.6	63
102	Quenched gas-phase reactions of tetraborane(10), B₄H₁₀, with substituted alkynes: new nido-dicarbapentaboranes and arachno-monocarbapentaboranes. Dalton Transactions, 2008, , 676-684.	1.6	3
103	N-Phosphino-amidines and -guanidines: synthesis, structure and P,N-chelate chemistry. Dalton Transactions, 2008, , 1043.	1.6	11
104	Facile photoinduced charge separation through a cyanoacetylide bridge in a heterobimetallic Fe(ii)â€“Re(i) complex. Chemical Communications, 2008, , 5845.	2.2	13
105	Carborane radical anions: spectroscopic and electronic properties of a carborane radical anion with a 2n + 3 skeletal electron count. Chemical Communications, 2007, , 2372.	2.2	61
106	Spectroscopic properties and electronic structures of 17-electron half-sandwich ruthenium acetylide complexes, [Ru(CCAr)(L2)Cpâ€“] ⁺ (Ar=phenyl, p-tolyl, 1-naphthyl, 9-anthryl; L2=(PPh3) ₂ , Cpâ€“=Cp; L2=dppe;) Tj ETQqO 0 0 rgBT /Over	1.6	74
107	Elemental fluorine. Journal of Fluorine Chemistry, 2007, 128, 29-33.	0.9	59
108	Synthetic and structural studies on C-ethynyl- and C-bromo-carboranes. Dalton Transactions, 2006, , 3544.	1.6	29

#	ARTICLE	IF	CITATIONS
109	Preparative and structural studies on sulfur-linked carborane icosahedra: 2-Phenyl-ortho-carboranyl-sulfur systems (2-Ph-1,2-C ₂ B ₁₀ H ₁₀) ₂ X (X=S, S ₂ or SO), and ortho-carboran-di-yl systems (1,2-C ₂ B ₁₀ H ₁₀ Y) ₂ (Y=S or SO). <i>Polyhedron</i> , 2006, 25, 300-306.	1.0	15
110	Improved syntheses of bis(ethynyl)-para-carboranes, 1,12-(RCC)2-1,12-C ₂ B ₁₀ H ₁₀ and 1,10-(RCC)2-1,10-C ₂ B ₈ H ₈ (R=H or Me ₃ Si). <i>Journal of Organometallic Chemistry</i> , 2006, 691, 3889-3894.	0.8	12
111	Reactions of Icosahedral Carboranes with Iminotris(dimethylamino)Phosphorane HNP(NMe ₂) ₃ : a Deboronation Intermediate nido-C ₂ B ₁₀ H ₁₂ ·N(H)P(NMe ₂) ₃ , Deboronation Reactions and Hydrogen-bonded Cluso-carborane Systems. <i>Journal of Cluster Science</i> , 2006, 17, 119-137.	1.7	22
112	Electronic interactions in bridged bis(cluster) assemblies – a comparison of para-CB ₁₀ H ₁₀ C, para-C ₆ H ₄ and C ₄ bridges. <i>Comptes Rendus Chimie</i> , 2005, 8, 1883-1896.	0.2	16
113	Versatile Gas/Liquid Microreactors for Industry. <i>Chemical Engineering and Technology</i> , 2005, 28, 344-352.	0.9	31
114	Elemental fluorine : Part 18. Selective direct fluorination of 1,3-ketoesters and 1,3-diketones using gas/liquid microreactor technology. <i>Lab on A Chip</i> , 2005, 5, 1132.	3.1	70
115	Gas-phase electron diffraction studies of the icosahedral carbaboranes, ortho-, meta- and para-C ₂ B ₁₀ H ₁₂ . <i>Dalton Transactions</i> , 2005, , 1310.	1.6	47
116	Elemental fluorine : Part 16. Versatile thin-film gas-liquid multi-channel microreactors for effective scale-out. <i>Lab on A Chip</i> , 2005, 5, 191-198.	3.1	107
117	Evolving Patterns in Boron Cluster Chemistry. <i>ChemInform</i> , 2004, 35, no.	0.1	0
118	Two Contrasting Ethynyl Hydroboration Pathways in the Formation of a Novel Tris-hydroboration Product from Reaction of Dimesitylborane with 2,5-Diethynylpyridine.. <i>ChemInform</i> , 2004, 35, no.	0.1	0
119	The synthesis and molecular and crystal structures of 1-methyl-2-carboxy-1,2-dicarba-closo-dodecaborane(12), 1-phenyl-2-carboxy-1,2-dicarba-closo-dodecaborane(12) and 1-phenyl-2-benzoyl-1,2-dicarba-closo-dodecaborane(12). <i>Polyhedron</i> , 2004, 23, 629-636.	1.0	25
120	Cage C...X interactions in solid-state structures of icosahedral carboranes. <i>Coordination Chemistry Reviews</i> , 2004, 248, 457-476.	9.5	140
121	Two contrasting ethynyl hydroboration pathways in the formation of a novel tris-hydroboration product from reaction of dimesitylborane with 2,5-diethynylpyridine Electronic supplementary information (ESI) available: NMR data for 4, rotatable 3-D molecular structure diagrams of optimised geometries in CHIME format and energy data for optimised geometries. See http://www.rsc.org/supplata/d3/ft/2004/216250h/ . <i>Chemical Communications</i> , 2004, , 702.	2.2	32
122	Gas-Phase Electron Diffraction Studies on Two 11-Vertex Dicarboranes, closo-2,3-C ₂ B ₉ H ₁₁ and nido-2,9-C ₂ B ₉ H ₁₃ . <i>Inorganic Chemistry</i> , 2004, 43, 5387-5392.	1.9	23
123	Sulfur, tin and gold derivatives of 1-(2-pyridyl)-ortho-carborane, 1-R-2-X-1,2-C ₂ B ₁₀ H ₁₀ (R = 2-pyridyl, X =) <i>Tj ETQq1 1 0,784314</i>	1.6	37
124	Exo-π-bonding to an ortho-carborane hypercarbon atom: systematic icosahedral cage distortions reflected in the structures of the fluoro-, hydroxy- and amino-carboranes, 1-X-2-Ph-1,2-C ₂ B ₁₀ H ₁₀ (X = F,) <i>Tj ETQq1 1 0,60 rgBT 10</i>	1.6	10
125	Synthesis and crystal structure of an assembly of three ortho-carborane cages linked via para-phenylene units: effect of aryl orientation on cage C-C bond lengths in C-aryl-ortho-carboranes. <i>Applied Organometallic Chemistry</i> , 2003, 17, 499-508.	1.7	50
126	Big macrocyclic assemblies of carboranes (big MACs): synthesis and crystal structure of a macrocyclic assembly of four carboranes containing alternate ortho- and meta-carborane icosahedra linked by para-phenylene units. <i>Journal of Organometallic Chemistry</i> , 2003, 680, 155-164.	0.8	32

#	ARTICLE	IF	CITATIONS
127	Dimesitylborane monomer-dimer equilibrium in solution, and the solid-state structure of the dimer by single crystal neutron and X-ray diffraction. <i>Journal of Organometallic Chemistry</i> , 2003, 680, 165-172.	0.8	47
128	9,12-Diiodo-1,2-dicarba-closo-dodecaborane(12). <i>Acta Crystallographica Section C: Crystal Structure Communications</i> , 2003, 59, o74-o76.	0.4	17
129	Synthesis and characterisation of some new boron compounds containing the 2,4,6-(CF ₃) ₃ C ₆ H ₂ (fluoromes = Ar), 2,6-(CF ₃) ₂ C ₆ H ₃ (fluoroxyl = Ar ^ε), or 2,4-(CF ₃) ₂ C ₆ H ₃ (Ar ^ε) ligands. <i>Dalton Transactions</i> , 2003, , 4395-4405.		79
130	Alkynyl Gold(I) Rigid-Rod Molecules from 1,12-Bis(ethynyl)-1,12-dicarba-closo-dodecaborane(12). <i>Organometallics</i> , 2003, 22, 4792-4797.	1.1	44
131	Intra- and inter-molecular carboranyl C ^δ -H ^δ : N hydrogen bonds in pyridyl-containing ortho-carboranes Electronic supplementary information (ESI) available: rotatable 3-D molecular structure diagrams of experimental structures of 1a ^ε and of MP2/6-31G* optimised geometries 1a ^ε and 7b in CHIME format. Computed GIAO NMR data for 1b ^ε and 4c. See http://www.rsc.org/suppdata/dt/b2/b209931d/ . <i>Dalton Transactions</i> , 2003, , 475-482.	1.6	90
132	Evolving patterns in boron cluster chemistry. <i>Pure and Applied Chemistry</i> , 2003, 75, 1315-1323.	0.9	74
133	Unexpected formation of new fluoroboranes from the reaction of NMe ₄ B ₃ H ₈ with BF ₃ and MeC≡CH: exo-2-FB ₄ H ₉ and trans-MeCH ₂ CHBF ₂ . <i>Chemical Communications</i> , 2002, , 2052-2053.	2.2	3
134	Solution and Solid-State Structure of the Anion [Ag ₂ {closo-CB ₁₁ H ₁₂ } ₄] ²⁻ . <i>Inorganic Chemistry</i> , 2002, 41, 4567-4573.	1.9	32
135	Synthesis of a new boron carbonitride with a B ₄ C-like structure from the thermolysis of N-alkylated borazines. <i>Chemical Communications</i> , 2002, , 718-719.	2.2	16
136	Model compounds and monomers for phenylene ether carboranylene ketone (PECK) polymer synthesis: preparation and characterization of boron-arylated ortho-carboranes bearing carboxyphenyl, phenoxyphenyl or benzoylphenyl substituents. <i>Journal of Materials Chemistry</i> , 2002, 12, 1301-1306.	6.7	62
137	A new nido-5-vertex cluster, phosphacarba-nido-pentaborane, 2-tBu-1,2-PCB ₃ H ₅ Electronic supplementary information (ESI) available: rotatable 3-D molecular structure diagrams of MP2-optimised geometries for 2-tBu-1,2-PCB ₃ H ₅ , 1-tBu-2,1-PCB ₃ H ₅ and P ₄ in CHIME format. See http://www.rsc.org/suppdata/cc/b2/b204409a/ . <i>Chemical Communications</i> , 2002, , 1448-1449.	2.2	21
138	A convenient cyanide-free one-pot synthesis of nido-Me ₃ N-7-CB ₁₀ H ₁₂ and nido-7-CB ₁₀ H ₁₃ ⁺ . <i>Dalton Transactions RSC</i> , 2002, , 2624.	2.3	25
139	Do the discrete dianions C ₂ B ₉ H ₁₁ ²⁻ exist? Characterisation of alkali metal salts of the 11-vertex nido dicarboranes, C ₂ B ₉ H ₁₁ ²⁻ , in solution. <i>Dalton Transactions RSC</i> , 2002, , 2009.	2.3	20
140	Cage-closing reactions of the nido-carborane anion 7,9-C ₂ B ₉ H ₁₂ ⁻ and derivatives; formation of neutral 11-vertex carboranes by acidification. <i>Dalton Transactions RSC</i> , 2002, , 3505.	2.3	21
141	Crystal and molecular structures of the nido-carborane anions, 7,9- and 2,9-C ₂ B ₉ H ₁₂ ⁻ . <i>Dalton Transactions RSC</i> , 2002, , 2132.	2.3	41
142	Halogenation of Tris(amido)tantalacarboranes with Dihalomethanes CH ₂ X ₂ (X = Cl, Br). <i>Collection of Czechoslovak Chemical Communications</i> , 2002, 67, 791-807.	1.0	4
143	The Molecular Structure of (PSH ⁺)(nido-7,8-C ₂ B ₉ H ₁₂ ⁻) Determined by Neutron Diffraction (PS = Proton) <i>J. Chem. Soc. Dalton Trans.</i> 1997, 1, 107-114	1.9	44
144	Synthesis of isomeric B-methylated tantalum carboranes, (Me ₂ N) ₃ TaC ₂ B ₉ H ₁₀ Me ^ε . <i>Dalton Transactions RSC</i> , 2001, , 2263-2269.	2.3	21

#	ARTICLE	IF	CITATIONS
145	Electrochemical evidence for electronic interactions through the para-carborane skeleton in the novel tricluster $[\{Co_2C_2(SiMe_3)(CO)_4(dppm)\}_2(\mu-CB_{10}H_{10}C)]$. <i>Chemical Communications</i> , 2001, , 1610-1611.	2.2	24
146	Phosphine promoted substituent redistribution reactions of B-chlorocatechol borane: molecular structures of ClBcat, BrBcat and $\mu-CB_{10}H_{10}C$ (cat = 1,2-O ₂ C ₆ H ₄ ; L = PMe ₃ , PEt ₃ , PBut ₃ , PCy ₃ , NEt ₃). <i>Dalton Transactions RSC</i> , 2001, , 1201-1209.	2.2	61
147	Empirical and ab Initio Energy/Architectural Patterns for 73 nido-6-vertex Carborane Isomers, from B ₆ H ₉ -to C ₄ B ₂ H ₆ . <i>Inorganic Chemistry</i> , 2001, 40, 1790-1801.	1.9	32
148	1,2-Ph ₂ -9-I-1,2-closo-C ₂ B ₁₀ H ₉ . <i>Acta Crystallographica Section C: Crystal Structure Communications</i> , 2000, 56, 487-488.	0.4	12
149	Synthesis and structure of 1,12-diethynyl-para-carborane. <i>Journal of Organometallic Chemistry</i> , 2000, 610, 20-24.	0.8	23
150	Gas-phase reactions of nido-1-methylpentaborane with propyne and 2-pentyne. Formation of B-alkyl nido and closo carbaboranes. <i>Journal of Organometallic Chemistry</i> , 2000, 614-615, 262-268.	0.8	11
151	Syntheses and reactions of some new C-pentafluorophenyl and tetrafluorophenylene carborane systems. <i>Journal of Organometallic Chemistry</i> , 2000, 597, 157-163.	0.8	15
152	Why are B ₂ O ₂ rings rare?. <i>Chemical Communications</i> , 2000, , 2217-2218.	2.2	15
153	First structural characterisation of a 2,1,12-MC ₂ B ₉ metallacarborane, [2,2,2-(NMe ₂) ₃ -closo-2,1,12-TaC ₂ B ₉ H ₁₁]. Trends in boron NMR shifts on replacing a {BH} vertex with a metal {ML _n } vertex in icosahedral carboranes. <i>Dalton Transactions RSC</i> , 2000, , 3519-3525.	2.3	19
154	Deboronation of 9-substituted-ortho- and -meta-carboranes. <i>Journal of Organometallic Chemistry</i> , 1999, 573, 279-291.	0.8	56
155	Deboronation of ortho-carborane by an iminophosphorane: crystal structures of the novel carborane adduct nido-C ₂ B ₁₀ H ₁₂ A-HNP(NMe ₂) ₃ and the borenium salt [(Me ₂ N) ₃ PNHBNP(NMe ₂) ₃]O ₂ +(C ₂ B ₉ H ₁₂ A) ⁺ . <i>Chemical Communications</i> , 1999, , 1649-1650.	2.2	65
156	Gas-Phase Flash Reactions of Diborane, Triborane Carbonyl and Tetraborane with Alkynes. <i>Collection of Czechoslovak Chemical Communications</i> , 1999, 64, 806-818.	1.0	12
157	Studies on tetraborane(8) carbonyl, B ₄ H ₈ A-CO: its isomeric composition in the gas phase and in solution, and its reactions with alkenes. Dedicated to Professor Ken Wade on the occasion of his 65th birthday in recognition of his outstanding contributions to the chemistry of the boranes, organometallics, and Main Group chemistry generally.1. <i>Journal of Organometallic Chemistry</i> , 1998, 550, 207-212.	0.8	6
158	Six-vertex nido-carborane structures with unusual CHB bridges or endo-CH hydrogens. Dedicated to Professor Ken Wade on the occasion of his 65th birthday in recognition of his outstanding contributions to organometallic and inorganic chemistry.1. <i>Journal of Organometallic Chemistry</i> , 1998, 550, 331-340.	0.8	19
159	2,4-Ethanotetraborane derivatives. 3. [1] determination of the molecular structure of 2,4-(t-butylethano)tetraborane(10), 2,4-(ButCHCH ₂)B ₄ H ₈ , in the gas phase by electron diffraction. <i>Journal of Molecular Structure</i> , 1998, 445, 319-334.	1.8	11
160	Transmission of electronic effects by icosahedral carboranes; skeletal carbon-13 chemical shifts and ultraviolet-visible spectra of substituted aryl-p-carboranes (1,12-dicarba-closo-dodecaboranes). <i>Journal of the Chemical Society Dalton Transactions</i> , 1998, , 401-412.	1.1	70
161	Molecular structures of 1,12-B ₁₂ H ₁₀ (CO) ₂ and its dihydrate 1,12-B ₁₂ H ₁₀ [C(OH) ₂] ₂ a novel bis-carbene complex. <i>Chemical Communications</i> , 1998, , 2487-2488.	2.2	22
162	Gas-Phase Reaction of Tetraborane(10) and Ethyne: Molecular Structure of nido-1,2-C ₂ B ₃ H ₇ in the Gas Phase. <i>Inorganic Chemistry</i> , 1998, 37, 2166-2176.	1.9	21

#	ARTICLE	IF	CITATIONS
163	Convenient direct syntheses of novel fused-ring CB ₄ N ₅ systems by nitrile hydroboration. <i>Journal of the Chemical Society Dalton Transactions</i> , 1997, , 3411-3413.	1.1	5
164	Molecular Structure of Trifluorophosphine Tetraborane(8), B ₄ H ₈ PF ₃ , As Determined in the Gas Phase by Electron Diffraction and ab Initio Computations. <i>Inorganic Chemistry</i> , 1997, 36, 1048-1054.	1.9	10
165	Some Boron-Containing Ring Systems. <i>Phosphorus, Sulfur and Silicon and the Related Elements</i> , 1997, 124, 73-82.	0.8	13
166	Cage-fluorination during deboronation of meta-carboranes. <i>Polyhedron</i> , 1997, 16, 2517-2525.	1.0	35
167	Fluoride-ion deboronation of p-fluorophenyl-ortho- and -meta-carboranes. NMR evidence for the new fluoroborate, HOBHF ₂ . <i>Polyhedron</i> , 1997, 16, 2499-2507.	1.0	62
168	Gas-Phase Reaction of Tetraborane(10) with Allene: The Fluxional arachno-1-Carbapentaborane(10) Isomeric System and Derivatives 1,2- and 1,3-Me ₂ -1-CB ₄ H ₈ ; Analogies in 1-CB ₄ H ₁₀ , MeB ₅ H ₁₀ , and B ₅ H ₁₀ . <i>Angewandte Chemie International Edition in English</i> , 1997, 36, 1498-1501.	4.4	20
169	Gasphasenreaktion von Tetraboran(10) mit Allen: das fluktuerende arachno-1-Carbapentaboran(10)-Isomerensystem und die Derivate 1,2- und 1,3-Me ₂ -1-CB ₄ H ₈ ; Analogien bei 1-CB ₄ H ₁₀ , MeB ₅ H ₁₀ und B ₅ H ₁₀ . <i>Angewandte Chemie</i> , 1997, 109, 1542-1544.		5
170	Existence of C ₃ -Me ₂ -closo-1,2-C ₂ B ₃ H ₃ Refuted by the ab Initio/IGLO, GIAO-MP2/NMR Method. Attempted Repetition of the Synthesis. <i>Inorganic Chemistry</i> , 1996, 35, 6170-6178.	1.9	24
171	Gas-phase reactions of tetraborane(10) with 1-en-3-yne: syntheses of the parent tricarbahexaborane, nido-2,3,4-C ₃ B ₃ H ₇ , and its derivatives. <i>Chemical Communications</i> , 1996, , 175.	2.2	13
172	Crystallographic evidence for the diene character of C ₂ B ₁₀ H ₁₀ C ₄ H ₄ (benzocarbonae™) and a Diels-Alder reaction of its anionic nido-analogue, [C ₂ B ₉ H ₁₀ C ₄ H ₄] ⁻ : crystal structures of C ₂ B ₁₀ H ₁₀ C ₄ H ₄ and C ₂ B ₁₀ H ₁₀ C ₄ H ₆ . <i>Chemical Communications</i> , 1996, , 2033-2034.	2.2	20
173	Deboronation of C-substituted ortho- and meta-closo-carboranes using wet fluoride ion solutions. <i>Polyhedron</i> , 1996, 15, 565-571.	1.0	100
174	2,4-Ethanotetraborane Derivatives. 2. Synthesis, Characterization, and Gas-Phase Structures of 2,4-(MeCHCH ₂)B ₄ H ₈ , 2,4-(trans-MeCHCHMe)B ₄ H ₈ , and 2- and 4-Pr-2,4-(MeCHCH ₂)B ₄ H ₇ . <i>Inorganic Chemistry</i> , 1995, 34, 2841-2849.	1.9	17
175	Re-identification of the major volatile carbaboranes from the gas-phase reactions of tetraborane(10) and alkynes at 50 °C. <i>Journal of the Chemical Society Chemical Communications</i> , 1995, , 667-668.	2.0	12
176	The Structures of Alkyl Derivatives of arachno-1-CB ₄ H ₁₀ from Reactions of B ₄ H ₁₀ with Alkynes. <i>Angewandte Chemie International Edition in English</i> , 1994, 33, 2298-2300.	4.4	24
177	Die Strukturen von Alkylderivaten des arachno-1-CB ₄ H ₁₀ aus der Reaktion von B ₄ H ₁₀ mit Alkinen. <i>Angewandte Chemie</i> , 1994, 106, 2384-2386.	1.6	14
178	The true identity of the bare-carbon™ cluster, closo-C ₃ B ₅ H ₇ . <i>Journal of the Chemical Society Dalton Transactions</i> , 1994, , 3197-3198.	1.1	10
179	C-arylation and C-heteroarylation of icosahedral carboranes via their copper(I) derivatives. <i>Journal of Organometallic Chemistry</i> , 1993, 462, 19-29.	0.8	151
180	X-ray structure and bonding of 1-phenylethynyl-2-phenyl-1,2-dicarbado-decaborane(12), [1-(PhC≡C)-2-Ph-1,2-C ₂ B ₁₀ H ₁₀], a model alkyne complex containing a rich variety of carbon-carbon bond types. <i>Polyhedron</i> , 1993, 12, 2711-2717.	1.0	29

#	ARTICLE	IF	CITATIONS
181	Proposed structure of the first Hypo carbaborane, C ₃ B ₄ H ₁₂ . Polyhedron, 1993, 12, 1849-1853.	1.0	16
182	Even more reliable NMR chemical shift computations by the GIAO-MP2 method. Journal of the Chemical Society Chemical Communications, 1993, , 1766-1768.	2.0	37
183	A pentuply-bridging thiocarbonyl group: x-ray crystal structure of a salt of the 1-thio-2-phenyl-1,2-dicarbododecaborate (12) anion, [LH] ⁺ [S(Ph)C ₂ B ₁₀ H ₁₀] ⁻ (L =) Tj ETQq1 1 0.784314 rgBT /Overlock 1037 50 65		