

Kenneth G Caulton

List of Publications by Year in descending order

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#	ARTICLE	IF	CITATIONS
1	Anion metathesis and chlorination of late transition metal pincer complexes: Comparing Co, Rh and Zn. <i>Inorganica Chimica Acta</i> , 2021, 516, 120118.	1.2	3
2	A proton-responsive ligand becomes a dimetal linker for multisubstrate assembly <i>via</i> nitrate deoxygenation. <i>Chemical Communications</i> , 2021, 57, 2780-2783.	2.2	10
3	Back donation, intramolecular electron transfer and N=O bond scission targeting nitrogen oxyanion reduction: how can a metal complex assist?. <i>Dalton Transactions</i> , 2021, 50, 2149-2157.	1.6	1
4	A redox cascade of NO ⁺ complexes: Structures and nitrogen deoxygenation thermodynamics. <i>Polyhedron</i> , 2021, 200, 115119.	1.0	2
5	A Redox-Active Tetrazine-Based Pincer Ligand for the Reduction of NO ⁻ Oxyanions Using a Redox-Inert Metal. <i>Chemistry - A European Journal</i> , 2021, 27, 11676-11681.	1.7	7
6	Pincers with diverse donors and their interconversion: application to Ni(II). <i>Zeitschrift Fur Anorganische Und Allgemeine Chemie</i> , 2021, 647, 1524-1529.	0.6	2
7	Nickel-mediated N=N bond formation and N ₂ O liberation <i>via</i> nitrogen oxyanion reduction. <i>Chemical Science</i> , 2021, 12, 10664-10672.	3.7	8
8	An Integrated View of Nitrogen Oxyanion Deoxygenation in Solution Chemistry and Electrospray Ion Production. <i>Inorganic Chemistry</i> , 2021, 60, 17241-17248.	1.9	3
9	A Dimeric Chromium(II) Pincer as an Electron Shuttle for N=N Bond Scission. <i>Chemistry - A European Journal</i> , 2020, 26, 13915-13926.	1.7	3
10	Unusual Dinitrogen Binding and Electron Storage in Dinuclear Iron Complexes. <i>Journal of the American Chemical Society</i> , 2020, 142, 8147-8159.	6.6	24
11	A reagent for heteroatom borylation, including iron mediated reductive deoxygenation of nitrate yielding a dinitrosyl complex. <i>Dalton Transactions</i> , 2020, 49, 1681-1687.	1.6	15
12	The Influence of Nucleophilic and Redox Pincer Character as well as Alkali Metals on the Capture of Oxygen Substrates: The Case of Chromium(II). <i>Chemistry - A European Journal</i> , 2020, 26, 9547-9555.	1.7	4
13	Nitrogen oxyanion reduction by Co(ⁱⁱ) augmented by a proton responsive ligand: recruiting multiple metals. <i>Dalton Transactions</i> , 2020, 49, 7891-7896.	1.6	13
14	[Cr(pincer ²⁺)] ₂ as an electron shuttle for reductively promoted hydrazine disproportionation. <i>Dalton Transactions</i> , 2019, 48, 11642-11646.	1.6	5
15	Gross rearrangement of Fe(II) aggregate structure by replacement of two H ⁺ by two Li ⁺ : Visualizing altered structure of acid versus conjugate base. <i>Polyhedron</i> , 2019, 174, 114152.	1.0	2
16	Selective deoxygenation of nitrate to nitrosyl using trivalent chromium and the Mashima reagent: reductive silylation. <i>Chemical Science</i> , 2019, 10, 475-479.	3.7	25
17	Gauging the Redox Non-Innocence of a Highly π -Acidic Bis-Tetrazine Pincer Ligand. <i>European Journal of Inorganic Chemistry</i> , 2019, 2019, 2535-2542.	1.0	2
18	Reductive Silylation Using a Bis-silylated Diaza-2,5-cyclohexadiene. <i>Chemistry - A European Journal</i> , 2019, 25, 8105-8111.	1.7	9

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19	Reactivity of an Unusual Divalent Chromium Aggregate Supported by a Multifunctional Bis(pyrazolate) Pincer Ligand. <i>European Journal of Inorganic Chemistry</i> , 2019, 2019, 1932-1940.	1.0	9
20	A bis- π -Pyrazolate Pincer on Reduced Cr Deoxygenates CO ₂ : Selective Capture of the Derived Oxide by Cr(II). <i>Chemistry - A European Journal</i> , 2019, 25, 7935-7940.	1.7	9
21	Multi-electron Reduction Capacity and Multiple Binding Pockets in Metal-Organic Redox Assembly at Surfaces. <i>Chemistry - A European Journal</i> , 2019, 25, 5565-5573.	1.7	7
22	Di- and trivalent chromium bis(pyrazol-3-yl)pyridine pincer complexes with good leaving groups. <i>Inorganica Chimica Acta</i> , 2019, 486, 483-491.	1.2	11
23	A new face for bis(pyrazol-3-yl)pyridine: Incompatible geometric preferences dictates unprecedented pincer ligand connectivity. <i>Inorganica Chimica Acta</i> , 2019, 485, 54-57.	1.2	4
24	A Multifunctional Pincer Ligand for Cobalt-Promoted Oxidation by N ₂ O. <i>Chemistry - A European Journal</i> , 2018, 24, 5962-5966.	1.7	15
25	Dehydrohalogenation of proton responsive complexes: versatile aggregation <i>via</i> pyrazolate pincer ligand arms. <i>Dalton Transactions</i> , 2018, 47, 2052-2060.	1.6	16
26	Redox-active ligand controlled selectivity of vanadium oxidation on Au(100). <i>Chemical Science</i> , 2018, 9, 1674-1685.	3.7	24
27	A substituent-tolerant synthetic approach to N/P-loaded-heteroarenes. <i>Dalton Transactions</i> , 2018, 47, 5938-5942.	1.6	12
28	A PNNH Pincer Ligand Allows Access to Monovalent Iron. <i>Chemistry - A European Journal</i> , 2018, 24, 1330-1341.	1.7	22
29	A new access route to dimetal sandwich complexes, including a radical anion. <i>Chemical Communications</i> , 2018, 54, 12397-12399.	2.2	5
30	Electrophile Recruitment as a Structural Element in Bis-Pyrazolate Pyridine Complex Aggregation. <i>European Journal of Inorganic Chemistry</i> , 2018, 2018, 5160-5166.	1.0	4
31	Seeking Redox Activity in a Tetrazinyl Pincer Ligand: Installing Zerovalent Cr and Mo. <i>Inorganic Chemistry</i> , 2018, 57, 12671-12682.	1.9	7
32	Probing Redox Noninnocence of Copper and Zinc Bis-pyridylpyrrolides. <i>European Journal of Inorganic Chemistry</i> , 2018, 2018, 4893-4904.	1.0	4
33	A multimetal-ligand cooperative approach to CO ₂ activation. <i>Inorganica Chimica Acta</i> , 2018, 483, 510-515.	1.2	2
34	Redox Isomeric Surface Structures Are Preferred over Odd-electron Pt(1+). <i>Chemistry - A European Journal</i> , 2018, 24, 15852-15858.	1.7	7
35	Electron and Oxygen Atom Transfer Chemistry of Co(II) in a Proton Responsive, Redox Active Ligand Environment. <i>Inorganic Chemistry</i> , 2018, 57, 6176-6185.	1.9	9
36	A Multifunctional Pincer Ligand Supports Unsaturated Cobalt: Five Functionalities in One Pincer. <i>Chemistry - A European Journal</i> , 2017, 23, 8039-8050.	1.7	26

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37	Deprotonation, Chloride Abstraction, and Dehydrohalogenation as Synthetic Routes to Bis(π-Pyrazolate Pyridyl Iron(II) Complexes. <i>European Journal of Inorganic Chemistry</i> , 2017, 2017, 3999-4012.	1.0	19
38	Ligand Design toward Multifunctional Substrate Reductive Transformations. <i>Inorganic Chemistry</i> , 2017, 56, 9505-9514.	1.9	21
39	Coordination and electronic characteristics of a nitrogen heterocycle pincer ligand. <i>Inorganica Chimica Acta</i> , 2016, 451, 82-91.	1.2	27
40	Tetrazine Assists Reduction of Water by Phosphines: Application in the Mitsunobu Reaction. <i>Chemistry - A European Journal</i> , 2016, 22, 13985-13998.	1.7	16
41	Two- and Three-Electron Oxidation of Single-Site Vanadium Centers at Surfaces by Ligand Design. <i>Journal of the American Chemical Society</i> , 2015, 137, 7898-7902.	6.6	37
42	A <i>cis</i> -Divacant Octahedral and Mononuclear Iron(IV) Imide. <i>Angewandte Chemie - International Edition</i> , 2014, 53, 14139-14143.	7.2	74
43	Multiplying the electron storage capacity of a bis-tetrazine pincer ligand. <i>Dalton Transactions</i> , 2014, 43, 6513-6524.	1.6	39
44	In search of redox non-innocence between a tetrazine pincer ligand and monovalent copper. <i>Dalton Transactions</i> , 2014, 43, 7958-7963.	1.6	11
45	Mechanistic Understanding of a Silver Pyridylpyrrolide as a Catalyst for 3 + 2 Cyclization of a Nitrile with Diazo Ester. <i>Organometallics</i> , 2014, 33, 1544-1552.	1.1	13
46	Transition Metal Chlorides Are Lewis Acids toward Terminal Chloride Attached to Late Transition Metals. <i>Inorganic Chemistry</i> , 2014, 53, 3307-3310.	1.9	17
47	Chemical Implications of Incompatible Ligand versus Metal Coordination Geometry Preferences. <i>Inorganic Chemistry</i> , 2014, 53, 3039-3047.	1.9	11
48	Probing the Steric and Electronic Characteristics of a New Bis-Pyrrolide Pincer Ligand. <i>Inorganic Chemistry</i> , 2014, 53, 1361-1369.	1.9	46
49	A tale of hydrogen abstraction, initially detected via X-ray diffraction. <i>Acta Crystallographica Section C, Structural Chemistry</i> , 2014, 70, 250-255.	0.2	1
50	Synthesis and Oxidative Reactivity of 2,2'-Pyridylpyrrolide Complexes of Ni(II). <i>Inorganic Chemistry</i> , 2013, 52, 9511-9521.	1.9	18
51	Reactivity of the terminal oxo species ((tBu ₂ PCH ₂ SiMe ₂) ₂ N)RhO. <i>Dalton Transactions</i> , 2013, 42, 6745.	1.6	20
52	New Approaches to Functionalizing Metal-Coordinated N ₂ . <i>Angewandte Chemie - International Edition</i> , 2013, 52, 4726-4732.	7.2	33
53	2,2'-Pyridylpyrrolide Ligand Redistribution Following Reduction. <i>Inorganic Chemistry</i> , 2013, 52, 5611-5619.	1.9	20
54	Bäcklund Reactions Catalyzed by a Silver(I) Pyridylpyrrolide: Understanding Arene C-C Insertion Selectivity. <i>Organometallics</i> , 2013, 32, 3185-3191.	1.1	33

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55	Probing the redox non-innocence of dinuclear, three-coordinate Co(II) indigo complexes: not simply η^2 -diketiminato variants. <i>Chemical Communications</i> , 2012, 48, 11082.	2.2	25
56	An iridium π -pyridylpyrrolide complex exhibiting reversible binding of H ₂ . <i>Dalton Transactions</i> , 2012, 41, 9619.	1.6	15
57	Systematics and Future Projections Concerning Redox-Noninnocent Amide/Imine Ligands. <i>European Journal of Inorganic Chemistry</i> , 2012, 2012, 435-443.	1.0	115
58	An evaluation of monovalent osmium supported by the PNP ligand environment. <i>Dalton Transactions</i> , 2011, 40, 1105-1110.	1.6	16
59	Assessment of the Electronic Structure of η^2 -Pyridylpyrrolides as Ligands. <i>Inorganic Chemistry</i> , 2011, 50, 8121-8131.	1.9	44
60	Three-Coordinate Ni(II): Tracing the Origin of an Unusual, Facile Si ⁺ C(sp ³) Bond Cleavage in [(⁺ Bu ₂ PCH ₂ SiMe ₂) ₂ N]Ni ⁺ . <i>Journal of the American Chemical Society</i> , 2011, 133, 2571-2582.	6.6	26
61	Mechanism of N/O Bond Scission of N ₂ O by an Unsaturated Rhodium Transient. <i>Journal of the American Chemical Society</i> , 2011, 133, 12576-12583.	6.6	26
62	Selectivity of attack on a Si ⁺ C(sp ³) sigma bond coordinated to Ni(II). <i>Inorganica Chimica Acta</i> , 2011, 374, 79-87.	1.2	9
63	Evaluating the reducing power of 3-coordinate T-shaped Ni(II). <i>Inorganica Chimica Acta</i> , 2011, 369, 49-54.	1.2	6
64	Redox and Lewis Acid Reactivity of Unsaturated Os(II). <i>European Journal of Inorganic Chemistry</i> , 2010, 2010, 4790-4800.	1.0	26
65	Ligand Influence on Metal Aggregation: a Unique Bonding Mode for Pyridylpyrrolides. <i>Inorganic Chemistry</i> , 2010, 49, 7626-7628.	1.9	18
66	Mechanism of Heterolysis of H ₂ by an Unsaturated d ⁸ Nickel Center: via Tetravalent Nickel?. <i>Journal of the American Chemical Society</i> , 2010, 132, 910-911.	6.6	76
67	Spin State, Structure, and Reactivity of Terminal Oxo and Dioxygen Complexes of the (PNP)Rh Moiety. <i>Chemistry - A European Journal</i> , 2008, 14, 7680-7686.	1.7	51
68	Reactivity of the Radical NO with a Masked Form of 14 Valence Electron (PNP)Rh: Forming Rh(0, I or II)?. <i>European Journal of Inorganic Chemistry</i> , 2008, 2008, 4704-4709.	1.0	8
69	[Ni{N(SiMe ₂ CH ₂ P(⁺ Bu ₂) ₂) ₂ }] ⁺ : Direct Observation of Transannular Si ⁺ C(sp ³) Bond Coordination. <i>Angewandte Chemie - International Edition</i> , 2008, 47, 9112-9114.	7.2	31
70	Mechanism of alkyne conversion to carbyne by 14- or 16-electron Os(H)2ClL2X (L=PiPr3; X=OTf or Tf) ETQq0 0 0 rgBT /Overlock 10 Tf 5	0.8	12
71	Coupling of terminal alkynes by RuHXL2 (X=Cl or N(SiMe3)2, L=PiPr3). <i>Journal of Organometallic Chemistry</i> , 2008, 693, 1664-1673.	0.8	19
72	Nitrogen-Ligated Iron Complexes: Photolytic Approach to the FeN ⁺ Moiety. <i>Inorganic Chemistry</i> , 2008, 47, 5129-5135.	1.9	18

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73	[(tBu ₂ PCH ₂ SiMe ₂) ₂ N]Rh: Rapidly Reversible H ^α -C(sp ³) and H ^β -C(sp ²) Bond Cleavage by Rhodium(I). <i>Organometallics</i> , 2008, 27, 166-168.	1.1	57
74	Redox Chemistry of the Triplet Complex (PNP)Co ^I . <i>Journal of the American Chemical Society</i> , 2008, 130, 4262-4276.	6.6	61
75	Influence of the d-Electron Count on CO Binding by Three-Coordinate [(t ⁺ Bu) ₂ PCH ₂ SiMe ₂] ₂ N]Fe, -Co, and -Ni. <i>Inorganic Chemistry</i> , 2008, 47, 407-409.	1.9	78
76	Bis(1,1,1,3,3,3-Hexamethyldisilazanato)-Bis(Tetrahydrofuran)Barium. <i>Inorganic Syntheses</i> , 2007, , 8-10.	0.3	12
77	Reactivity of ⁺ OsO ₄ with an osmium polyhydride: Reductive elimination and reductive nitrosylation on the path from odd- to even-electron molecules. <i>New Journal of Chemistry</i> , 2007, 31, 838-840.	1.4	6
78	Exploring the Reactivity of Four-Coordinate PNPCoX with Access to Three-Coordinate Spin Triplet PNPCo. <i>Inorganic Chemistry</i> , 2007, 46, 10321-10334.	1.9	49
79	Facile Hydrogenation of N ₂ O by an Operationally Unsaturated Osmium Polyhydride. <i>Journal of the American Chemical Society</i> , 2007, 129, 8706-8707.	6.6	59
80	Metal-Dependent Reactivity Differences for Transients Formed By Flash Photolysis of (PNP)M(CO), M = Co and Rh. <i>Journal of the American Chemical Society</i> , 2007, 129, 15430-15431.	6.6	8
81	Influence of Chelate Substituents on the Structure and Spin State of Unsaturated [N(SiMe ₂ CH ₂ PtBu ₂) ₂]Ru ^{II} X. <i>Inorganic Chemistry</i> , 2007, 46, 4612-4616.	1.9	17
82	2-Electron reduction of N ₂ O yields a new ligand binding mode: N-Nitrosoimide. <i>Polyhedron</i> , 2007, 26, 4731-4736.	1.0	9
83	NO binds to unsaturated Os(IV) polyhydrides as a redox reagent. <i>Journal of Organometallic Chemistry</i> , 2007, 692, 3121-3132.	0.8	1
84	Reducing Power of Three-Coordinate Cobalt(I). <i>Journal of the American Chemical Society</i> , 2006, 128, 4248-4249.	6.6	50
85	Triple Benzylic Dehydrogenation by Osmium in an Amide Ligand Environment. <i>Organometallics</i> , 2006, 25, 802-804.	1.1	46
86	Cleavage of F ^α -C(sp ²) bonds by MHR(CO)(PtBu ₂ Me) ₂ (M=Os and Ru; R=H, CH ₃ or Aryl): Product dependence on M and R. <i>Polyhedron</i> , 2006, 25, 459-468.	1.0	36
87	N ₂ Provides Insight into the Mechanism of H ^α -C(sp ³) Bond Cleavage. <i>Journal of the American Chemical Society</i> , 2005, 127, 16780-16781.	6.6	34
88	A Facile Approach to a d ⁴ Ru ^{II} -N ³ Moiety. <i>Journal of the American Chemical Society</i> , 2005, 127, 5330-5331.	6.6	75
89	New d ⁴ dihydrides of Ru(IV) and Os(IV) with σ -donor ligands: M(H) ₂ (chelate)(PiPr ₃) ₂ with chelate = ortho-XYC ₆ H ₄ with X, Y = O, NR; R = H or CH ₃ . <i>New Journal of Chemistry</i> , 2005, 29, 193-204.	1.4	21
90	Room-temperature hydrosilylation of the CF bond of vinyl fluoride catalyzed by osmium hydrides. <i>Journal of Molecular Catalysis A</i> , 2004, 224, 125-131.	4.8	20

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91	Facile Insertion of Terminal Acetylenes into the Rull π NR ₂ Bond of a 14-Valence-Electron Complex. <i>Organometallics</i> , 2004, 23, 4814-4816.	1.1	18
92	Reactivity of the Hydrido/Nitrosyl Radical MHCl(NO)(CO)(PiPr ₃) ₂ , M = Ru, Os. <i>Inorganic Chemistry</i> , 2004, 43, 351-360.	1.9	38
93	Hydrocarbyl Ligand π -Tuning of the PtII/IV Redox Potential. <i>Inorganic Chemistry</i> , 2004, 43, 3642-3646.	1.9	17
94	Terminal Acetylenes React to Increase Unsaturation in [(tBu ₂ PCH ₂ SiMe ₂) ₂ N]Re(H) ₄ . <i>Organometallics</i> , 2004, 23, 4934-4943.	1.1	16
95	Design and Synthesis of Tridentate Facially Chelating Ligands of the [2.n.1]-(2,6)-Pyridinophane Family. <i>Journal of Organic Chemistry</i> , 2003, 68, 4806-4814.	1.7	19
96	Double C(sp ³) dehydrogenation as a route to coordinated Arduengo carbenes: experiment and computation on comparative π -acidity. <i>New Journal of Chemistry</i> , 2003, 27, 1446-1450.	1.4	36
97	Conversion of Ethylene to Hydride and Ethylidyne by an Amido Rhenium Polyhydride. <i>Organometallics</i> , 2003, 22, 2539-2541.	1.1	42
98	N π -PtIV π -H/N π -H π -PtII intramolecular redox equilibrium in a product of H π -C(sp ²) cleavage and unusual alkane/arene C π -H bond selectivity of ([2.1.1]pyridinophane)PtII(CH ₃)+Electronic supplementary information (ESI) available: experimental details. See http://www.rsc.org/suppdata/cc/b2/b207797n/ . <i>Chemical Communications</i> , 2003, , 358-359.	2.2	29
99	Amido/phosphine pincer hydrides of ruthenium. <i>New Journal of Chemistry</i> , 2003, 27, 263-273.	1.4	25
100	π -Donor olefin substituents alter olefin binding to CpFe(CO) ₂ +. <i>New Journal of Chemistry</i> , 2003, 27, 1769-1774.	1.4	19
101	Evaluation of energies of isomeric SO ₂ complexesElectronic supplementary information (ESI) available: DFT geometry optimized structures of isomeric RuCl(CO)(PMe ₃) ₂ HSO ₂ . See http://www.rsc.org/suppdata/nj/b2/b212025a/ . <i>New Journal of Chemistry</i> , 2003, 27, 680-683.	1.4	5
102	[2.1.1]-(2,6)-Pyridinophane(L)-controlled alkane C π -H bond cleavage: (L)PtMe ₂ H ⁺ as a precursor to the geometrically π -tense π -transient (L)PtMe+Electronic supplementary information (ESI) available: complete experimental and characterization data. See http://www.rsc.org/suppdata/nj/b3/b302055j/ . <i>New Journal of Chemistry</i> , 2003, 27, 665-667.	1.4	30
103	Facile C(sp ²)/O ₂ CR bond cleavage by Ru or Os. <i>New Journal of Chemistry</i> , 2003, 27, 1451-1462.	1.4	22
104	Four-Coordinate, Planar Ru(II). A Triplet State as a Response to a 14-Valence Electron Configuration. <i>Journal of the American Chemical Society</i> , 2003, 125, 8426-8427.	6.6	91
105	Assessing isomeric structures of pincer-ligated ruthenium and osmium polyhydrides using density functional calculations. <i>Molecular Physics</i> , 2002, 100, 385-395.	0.8	3
106	Conventional Lithium Bases as Unconventional Sources of Methyl Anion: π -Facile Me π -Si and Me π -C Bond Cleavage in RLi, R ₂ NLi, and BR ₄ - by an Electrophilic Osmium Dihydride. <i>Organometallics</i> , 2002, 21, 4030-4049.	1.1	7
107	An Electron-Excessive Nitrosyl Complex: π -Reactivity of a Ligand-Centered Radical Leading to Coordinated HNO. <i>Inorganic Chemistry</i> , 2002, 41, 4087-4089.	1.9	24
108	[n.1.1]-(2,6)-Pyridinophanes: A New Ligand Type Imposing Unusual Metal Coordination Geometries. <i>Inorganic Chemistry</i> , 2002, 41, 6867-6874.	1.9	16

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109	Electrophilic (Li+) acceleration of C-H reductive elimination and oxidative addition reactions of Os(II)/Os(0) nitrosyl complexes. Electronic supplementary information (ESI) available: full experimental and computational details. See http://www.rsc.org/suppdata/nj/b2/b200180m/ . <i>New Journal of Chemistry</i> , 2002, 26, 498-502.	1.4	1
110	Geminal dehydrogenation of ether and amine C(sp ³)H ₂ groups by electron-rich Ru(II) and Os. Electronic supplementary information (ESI) available: crystallographic data, fractional coordinates and isotropic thermal parameters, anisotropic thermal parameters, and bond distances and angles. See http://www.rsc.org/suppdata/nj/b2/b200168n/ . <i>New Journal of Chemistry</i> , 2002, 26, 687-700.	1.4	57
111	Evolving approaches to integration of computational with experimental chemistry: an example from osmium hydrido (η ¹ -η ²)-nitrosyl/nitroxyl chemistry. <i>New Journal of Chemistry</i> , 2002, 26, 1267-1269.	1.4	3
112	Title is missing!. <i>Angewandte Chemie</i> , 2002, 114, 4276-4278.	1.6	1
113	A comprehensive view of C-H addition across the RC≡CH bond: frustration culminating in ultimate union. <i>New Journal of Chemistry</i> , 2001, 25, 1244-1255.	1.4	56
114	C≡DO (DO = Ï€-donor, F) Cleavage in H ₂ CCH(DO) by (Cp ₂ ZrHCl) _n : Mechanism, Agostic Fluorines, and a Carbene of Zr(IV). <i>Journal of the American Chemical Society</i> , 2001, 123, 603-611.	6.6	88
115	Facile C(sp ²)/OR Bond Cleavage by Ru or Os. <i>Inorganic Chemistry</i> , 2001, 40, 6610-6621.	1.9	23
116	Os(H) ₂ NO(PiPr ₃) ₂ : Mechanism and Energetics of a Room Temperature Reversible Ethyl/Hydridoethylene Equilibrium and Contrasting Double Insertion of Acetylene. <i>Organometallics</i> , 2001, 20, 2040-2046.	1.1	17
117	Factors governing the equilibrium between metal-alkyl, alkylidene and alkylidyne: MCX ₂ R, Xi-, M...CXR and X ₂ M†CR. <i>Journal of Organometallic Chemistry</i> , 2001, 617-618, 56-64.	0.8	72
118	Carbene transposition involving double dehydrogenation of an sp ³ carbon. <i>Chemical Communications</i> , 2001, , 1158-1159.	2.2	22
119	Facile and Reversible Cleavage of C-F Bonds. Contrasting Thermodynamic Selectivity for RuCF ₂ H vs FOsCFH. <i>Journal of the American Chemical Society</i> , 2000, 122, 8916-8931.	6.6	99
120	Coordinated carbenes from electron-rich olefins on RuHCl(PPr ₃) ₂ . <i>New Journal of Chemistry</i> , 2000, 24, 9-26.	1.4	87
121	Structural Distortions in mer-M(H) ₃ (NO)L ₂ (M = Ru, Os) and Their Influence on Intramolecular Fluxionality and Quantum Exchange Coupling. <i>Inorganic Chemistry</i> , 2000, 39, 1919-1932.	1.9	23
122	Fate of CH ₂ CH(E) in the Presence of Unsaturated Ru(X)(H)L ₂ q+(X = Cl, q= 0; X = CO, q= 1): Highly Sensitive to X and E. <i>Organometallics</i> , 2000, 19, 2291-2298.	1.1	17
123	Cleavage of H-C(sp ²) and C(sp ²)-X Bonds (X = Alkyl, Aryl, OR, NR ₂): Facile Decarbonylation, Isonitrile Abstraction, or Dehydrogenation of Aldehydes, Esters, Amides, Amines, and Imines by [RuHCl(PiPr ₃) ₂] ₂ . <i>Organometallics</i> , 2000, 19, 3569-3578.	1.1	58
124	A New Class of Electron-Rich Unsaturated Molecules: Ru ₂ HnX _{4-n} (PiPr ₃) ₄ , X = Anion. <i>Inorganic Chemistry</i> , 2000, 39, 3757-3764.	1.9	28
125	Reactivity of [RuHCl(PiPr ₃) ₂] ₂ with Functionalized Vinyl Substrates. The H ₂ Ligand as a Sensitive Probe of Electronic Structure. <i>Inorganic Chemistry</i> , 2000, 39, 3749-3756.	1.9	43
126	A 14-Electron Ruthenium(II) Hydride, [RuH(CO)(PtBu ₂ Me) ₂]BAR ₄ (Ar = 3,5-(C ₆ H ₃)(CF ₃) ₂): Synthesis, Structure, and Reactivity toward Alkenes and Oxygen Ligands. <i>Organometallics</i> , 2000, 19, 2281-2290.	1.1	59

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