

Michael G Richmond

List of Publications by Year in descending order

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| # | ARTICLE | IF | CITATIONS |
|----|---|-----|-----------|
| 1 | Stereochemical control of the diphosphine and alkyne ligands in triruthenium clusters: the effect of reversible CO loss/addition on the ligand distribution in [Ru ₃ (μ ₃ -1,2-PhCCPh){μ ₃ -1,2-Ph ₂ PCH(Me)PPh ₂ }(CO) _{7,8}]. <i>Journal of Organometallic Chemistry</i> , 2022, , 122337. | 1.8 | 1 |
| 2 | Biomimics of [FeFe]-hydrogenases incorporating redox-active ligands: synthesis, redox properties and spectroelectrochemistry of diiron-dithiolate complexes with ferrocenyl-diphosphines as Fe ₄ S ₄ surrogates. <i>Dalton Transactions</i> , 2022, 51, 9748-9769. | 3.3 | 11 |
| 3 | Bimodal substitution behavior in the reaction of N,N TM -diisopropylformamidine with [Os ₃ (CO) ₁₀ (NCMe) ₂]: Kinetics and molecular structures of the formamidinate-substituted clusters HO ₃ (CO) ₉ [1/4-C(O)NPr C(H)NPr], HO ₃ (CO) ₁₀ [1/4-NPr C(H)NPr], and HO ₃ (CO) ₉ [1/4-3-NPr C(H)NPr]. <i>Journal of Organometallic Chemistry</i> , 2021, 936, 121709. | 1.8 | 0 |
| 4 | Proton reduction by phosphinidene-capped triiron clusters. <i>Journal of Organometallic Chemistry</i> , 2021, 943, 121816. | 1.8 | 0 |
| 5 | Ligand coordination in [Re ₂ (CO) ₉ (NCMe)] and [H ₃ Re ₃ (CO) ₁₁ (NCMe)] by triphenylantimony: Reactivity studies and Sb-Ph bond cleavage to give new antimony-containing di- and trirhenium complexes. <i>Journal of Organometallic Chemistry</i> , 2021, 953, 122034. | 1.8 | 2 |
| 6 | Reactions of [Ru ₃ (CO) ₁₂] with thiosaccharin: Synthesis and structure of di-, tri-, tetra- and penta-ruthenium complexes containing a thiosaccharinate ligand(s). <i>Journal of Organometallic Chemistry</i> , 2020, 906, 121048. | 1.8 | 7 |
| 7 | Reactions of triosmium and triruthenium clusters with 2-ethynylpyridine: new modes for alkyne C-C bond coupling and C-H bond activation. <i>RSC Advances</i> , 2020, 10, 30671-30682. | 3.6 | 6 |
| 8 | Electrocatalytic proton-reduction behaviour of telluride-capped triiron clusters: tuning of overpotentials and stabilization of redox states relative to lighter chalcogenide analogues. <i>Dalton Transactions</i> , 2020, 49, 7133-7143. | 3.3 | 5 |
| 9 | Thermolysis of [HO ₃ (CO) ₈ {μ ₃ -Ph ₂ PCH ₂ P(Ph)C ₆ H ₄ }] : New Os ₂ - and Os ₃ - cluster products based on multiple C-H bond activation of the bis(diphenylphosphino)methane ligand. <i>Inorganica Chimica Acta</i> , 2020, 510, 119733. | 2.4 | 4 |
| 10 | Electron Transfer Mediated by Iron Carbonyl Clusters Enhance Light-Driven Hydrogen Evolution in Water by Quantum Dots. <i>ChemSusChem</i> , 2020, 13, 3252-3260. | 6.8 | 7 |
| 11 | Facile Os-Os bond cleavage in the reactions of [Os ₃ (CO) ₁₀ (NCMe) ₂] and [Os ₃ (CO) ₁₀ (1/4-H) ₂] with tetramethylthiuram disulfide (tmtsd): Syntheses and crystal structures of new polynuclear osmium carbonyl complexes containing a dimethyldithiocarbamate ligand(s). <i>Journal of Organometallic Chemistry</i> , 2020, 911, 121133. | 1.8 | 7 |
| 12 | Asymmetric hydrogenation of an α -unsaturated carboxylic acid catalyzed by intact chiral transition metal carbonyl clusters - diastereomeric control of enantioselectivity. <i>Dalton Transactions</i> , 2020, 49, 4244-4256. | 3.3 | 4 |
| 13 | A new synthetic route for the preparation of [Os ₃ (CO) ₁₀ (1/4-OH)(1/4-H)] and its reaction with bis(diphenylphosphino)methane (dppm): syntheses and X-ray structures of two isomers of [Os ₃ (CO) ₈ (1/4-OH)(1/4-H)(1/4-dppm)] and [Os ₃ (CO) ₇ (1/4-3-CO)(1/4-3-O)(1/4-dppm)]. <i>RSC Advances</i> , 2020, 10, 44688-44711. | 3.6 | 4 |
| 14 | Reactions of [Os ₃ (CO) ₁₀ (1/4-dppm)] and [HO ₃ (CO) ₈ {1/4-3-Ph ₂ PCH ₂ P(Ph)C ₆ H ₄ }] with Bu ₃ GeH: Ge-H and Ge-C bond cleavage at triosmium centers. <i>Journal of Organometallic Chemistry</i> , 2019, 898, 120862. | 1.8 | 7 |
| 15 | New molecular architectures containing low-valent cluster centres with di- and trimetalated 2-vinylpyrazine ligands: synthesis and molecular structures of Ru ₅ (CO) ₁₅ (1/4-5-C ₄ H ₂ N ₂ CHiC)(1/4-H) ₂ and Ru ₈ (CO) ₂₄ (1/4-7-C ₄ H ₂ N ₂ CHiC)(1/4-H) ₃ . <i>RSC Advances</i> , 2019, 9, 21025-21030. | 3.6 | 2 |
| 16 | Polyhedral Flexibility in the Sulfido-Capped Cluster H ₂ Ru ₃ (CO) ₉ (1/4-3-S) on Reaction with 2-(Diphenylphosphino)thioanisole (PS) and Reversible Tripodal Rotation of the Chelated PS Ligand in H ₂ Ru ₃ (CO) ₇ (1/4-3-PS)(1/4-3-S). <i>Organometallics</i> , 2019, 38, 2472-2484. | 2.3 | 2 |
| 17 | Models of the iron-only hydrogenase enzyme: structure, electrochemistry and catalytic activity of Fe ₂ (CO) ₃ (1/4-dithiolate)(1/4-1 ^{sup} ,1 ^{sup} -2 ^{sup} -triphos). <i>Dalton Transactions</i> , 2019, 48, 6174-6190. | 3.3 | 31 |
| 18 | Reactivity of [Mo(CO) ₃ (NCMe) ₃] towards pyrimidine-2-thiol (pymSH) and thiophenol (PhSH) in the presence of phosphine auxiliaries: Synthesis of mono- and dinuclear complexes bearing 1 ² and μ ₂ -pymS coordination motifs. <i>Polyhedron</i> , 2019, 164, 55-63. | 2.2 | 4 |

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|----|---|-----|-----------|
| 19 | Chalcogenide-capped triiron clusters $[\text{Fe}_3(\text{CO})_9(\eta^3\text{-E})_2]$, $[\text{Fe}_3(\text{CO})_7(\eta^3\text{-CO})(\eta^3\text{-E})(\eta^4\text{-dppm})]$ and $[\text{Fe}_3(\text{CO})_7(\eta^3\text{-E})_2(\eta^4\text{-dppm})]$ (E = S, Se) as proton-reduction catalysts. <i>Journal of Organometallic Chemistry</i> , 2019, 880, 213-222. | 1.8 | 6 |
| 20 | Activation of thiosaccharin at a polynuclear osmium cluster. <i>Journal of Organometallic Chemistry</i> , 2019, 880, 223-231. | 1.8 | 4 |
| 21 | Hydrogenase biomimics containing redox-active ligands: $\text{Fe}_2(\text{CO})_4(\eta^3\text{-edt})(\eta^2\text{-bpcd})$ with electron-acceptor 4,5-bis(diphenylphosphino)-4-cyclopenten-1,3-dione (bpcd) as a potential $[\text{Fe}_4\text{S}_4\text{H}]$ surrogate. <i>Dalton Transactions</i> , 2019, 48, 6051-6060. | 3.3 | 31 |
| 22 | Synthesis and redox properties of fac-BrRe(CO) ₃ [1,2-(PPh ₂) ₂ -closo-1,2-C ₂ B ₁₀ H ₁₀]: The first structurally characterized rhenium carbonyl containing a carboranyl-based diphosphine ligand. <i>Journal of Molecular Structure</i> , 2018, 1156, 397-402. | 3.6 | 2 |
| 23 | Synthesis and molecular structures of the 52-electron triiron telluride clusters $[\text{Fe}_3(\text{CO})_8(\eta^3\text{-Te})_2(\eta^2\text{-diphosphine})]$ - Electrochemical properties and activity as proton reduction catalysts. <i>Journal of Organometallic Chemistry</i> , 2018, 867, 381-390. | 1.8 | 8 |
| 24 | Experimental and computational preference for phosphine regioselectivity and stereoselective tripod rotation in $\text{H}_3\text{Os}(\text{CO})_8(\text{PPh})_2(\eta^3\text{-1,2-N,C-}\eta^5\text{-C}_7\text{H}_7\text{Cp})$. <i>RSC Advances</i> , 2018, 8, 32672-32683. | 3.6 | 10 |
| 25 | Hydrogenase Biomimetics with Redox-Active Ligands: Synthesis, Structure, and Electrocatalytic Studies on $[\text{Fe}_2(\text{CO})_4(\eta^2\text{-dppn})(\eta^2\text{-edt})]$ (edt = Ethanedithiolate; dppn =) Tj ETQq1 1 0.784314 rgBT /Overlock 10 2f 50 4971d (1,8- bis | | |
| 26 | Electrocatalytic proton reduction by thiolate-capped triiron clusters $[\text{Fe}_3(\text{CO})_9(\eta^3\text{-SR})(\eta^3\text{-H})]$ (R = iPr, tBu). <i>Inorganica Chimica Acta</i> , 2018, 480, 47-53. | 2.4 | 13 |
| 27 | Synthesis of the labile rhenium(I) complexes fac-Re(CO) ₃ (L)($\eta^2\text{-O,O-FcC(O)CHC(O)Me}$) (where Fc =) Tj ETQq1 1 0.784314 rgBT /Overlock 10 2f 50 4971d <i>Journal of Organometallic Chemistry</i> , 2018, 874, 87-100. | 1.8 | 3 |
| 28 | Mixed-valence dimolybdenum complexes containing hard oxo and soft carbonyl ligands: synthesis, structure, and electrochemistry of $\text{Mo}_2(\text{O})(\text{CO})_2(\eta^3\text{-S}(\text{CH}_2)_n\text{S})(\eta^2\text{-diphosphine})$. <i>Dalton Transactions</i> , 2018, 47, 10102-10112. | 3.3 | 3 |
| 29 | A new diphosphine-carbonyl complex of ruthenium: an efficient precursor for C=C and C=N bond coupling catalysis. <i>Dalton Transactions</i> , 2018, 47, 10264-10272. | 3.3 | 11 |
| 30 | Diphosphine-induced thiolate-bridge scission of $[\text{Re}(\text{CO})_3(\eta^3\text{-S,N-thpymS})_2]$ (thpymS =) Tj ETQq0 0 0 rgBT /Overlock 10 Tf 50 3071d isomers of $[\text{Re}(\text{CO})_3(\eta^3\text{-S,N-thpymS})_2(\eta^3\text{-1,1'-dppe})]$. <i>Journal of Organometallic Chemistry</i> , 2018, 871, 167-177. | 1.8 | 4 |
| 31 | Reversible C-H bond activation at a triosmium centre: A comparative study of the reactivity of unsaturated triosmium clusters $\text{Os}_3(\text{CO})_8(\eta^3\text{-dppm})(\eta^3\text{-H})_2$ and $\text{Os}_3(\text{CO})_8(\eta^3\text{-dppf})(\eta^3\text{-H})_2$ with activated alkynes. <i>Journal of Organometallic Chemistry</i> , 2017, 836-837, 68-80. | 1.8 | 7 |
| 32 | Mixed main group transition metal clusters: Reactions of $[\text{Ru}_3(\text{CO})_{10}(\eta^3\text{-dppm})]$ with Ph_3SnH . <i>Journal of Organometallic Chemistry</i> , 2017, 840, 47-55. | 1.8 | 8 |
| 33 | The reaction of $\text{Os}_3(\text{CO})_{12}$ with triphos $\{\text{MeC}(\text{CH}_2\text{PPh}_2)_3\}$: A case of multiple C-P and C-H bond activations. <i>Journal of Organometallic Chemistry</i> , 2017, 849-850, 125-129. | 1.8 | 0 |
| 34 | Reactions of $\text{Ru}_3(\text{CO})_{10}(\eta^3\text{-dppm})$ with Ph_3GeH : Ge-H and Ge-C bond cleavage in Ph_3GeH at triruthenium clusters. <i>Journal of Organometallic Chemistry</i> , 2017, 843, 75-86. | 1.8 | 12 |
| 35 | Reactions of the face-capped benzothiazolate-substituted clusters $\text{Os}_3(\text{CO})_9(\eta^3\text{-1,2-C}_7\text{H}_3\text{NSR})(\eta^3\text{-H})$ (R = H, Me) with PPh_3 : Kinetic formation of $\text{Os}_3(\text{CO})_9(\text{PPh}_3)(\eta^3\text{-1,2-C}_7\text{H}_3\text{NSR})(\eta^3\text{-H})$ and thermally induced ligand isomerization. <i>Journal of Organometallic Chemistry</i> , 2017, 849-850, 337-349. | 1.8 | 4 |
| 36 | Alkyne activation and polyhedral reorganization in benzothiazolate-capped osmium clusters on reaction with diethyl acetylenedicarboxylate (DEAD) and ethyl propiolate. <i>Dalton Transactions</i> , 2017, 46, 13597-13609. | 3.3 | 2 |

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|----|--|-----|-----------|
| 37 | Ambidentate Ligand Reactivity with the Rhenium(I) Compounds [BrRe(CO) ₄] ₂ and <i>cis</i> -BrRe(CO) ₄ L: A Kinetic and Mechanistic Study. <i>European Journal of Inorganic Chemistry</i> , 2017, 2017, 3990-3998. | 2.0 | 3 |
| 38 | Diphosphine-bridged digold(I) compounds: Structural and computational studies on the aurophilic interaction in Au ₂ Cl ₂ ($\frac{1}{4}$ -bpcd) and Au ₂ Cl ₂ ($\frac{1}{4}$ -bmi). <i>Journal of Molecular Structure</i> , 2017, 1129, 188-194. | 3.6 | 3 |
| 39 | Microwave-induced dpmm ligand substitution in triosmium clusters: Structural and DFT evaluation of Os ₃ clusters containing multiply activated dpmm ligands through cyclometalation, ortho metalation, and P C bond cleavage. <i>Journal of Organometallic Chemistry</i> , 2016, 813, 15-25. | 1.8 | 6 |
| 40 | Biomimetics of the [FeFe]-hydrogenase enzyme: Identification of kinetically favoured apical-basal [Fe ₂ (CO) ₄ ($\frac{1}{4}$ -H){ η^2 -Ph ₂ PC(Me) ₂ PPh ₂ }($\frac{1}{4}$ -pdt)] ⁺ as a proton-reduction catalyst. <i>Journal of Organometallic Chemistry</i> , 2016, 812, 247-258. | 1.8 | 54 |
| 41 | Hydrogenase biomimetics with redox-active ligands: Electrocatalytic proton reduction by [Fe ₂ (CO) ₄ (η^2 -diamine)($\frac{1}{4}$ -edt)] (diamine = 2,2'-bipy, 1,10-phen). <i>Polyhedron</i> , 2016, 116, 127-135. | 2.2 | 36 |
| 42 | Syntheses and Characterization of Tantalum Alkyl Imides and Amide Imides. DFT Studies of Unusual η^3 -SiMe ₃ Abstraction by an Amide Ligand. <i>Organometallics</i> , 2015, 34, 5687-5696. | 2.3 | 13 |
| 43 | Phenazine-substituted polynuclear osmium clusters: Synthesis and DFT evaluation of the C-metalated derivatives Os ₃ (CO) ₉ ($\frac{1}{4}$ 3,1-2-C ₁₂ H ₇ N ₂)($\frac{1}{4}$ -H) and Os ₃ (CO) ₉ ($\frac{1}{4}$ 3,1-2-C ₁₂ H ₆ N ₂)($\frac{1}{4}$ -H) ₂ . <i>Journal of Organometallic Chemistry</i> , 2015, 779, 21-29. | | 7 |
| 44 | Electrocatalytic proton reduction catalysed by the low-valent tetrairon-oxo cluster [Fe ₄ (CO) ₁₀ (η^2 -dppn)($\frac{1}{4}$ -O)] ²⁺ [dppn = 1,1'-bis(diphenylphosphino)naphthalene]. <i>Dalton Transactions</i> , 2015, 44, 5160-5169. | 3.3 | 11 |
| 45 | Synthesis of the Stereoisomeric Clusters 1,2-Os ₃ (CO) ₁₀ (trans-dpmn) and 1,2-Os ₃ (CO) ₁₀ (cis-dpmn) [where dpmn = 2,3-bis(diphenylphosphinomethyl)-5-norbornene]: DFT Evaluation of the Isomeric Clusters 1,2-Os ₃ (CO) ₁₀ (dpmn) and Isomer-Dependent Diphosphine Ligand Activation. <i>Journal of Cluster Science</i> , 2015, 26, 93-109. | 3.3 | 2 |
| 46 | Synthesis, structure and bonding of new mono- and dinuclear molybdenum complexes containing pyridine-2-thiolate (pyS) and different P-donors. <i>Inorganica Chimica Acta</i> , 2015, 434, 150-157. | 2.4 | 13 |
| 47 | Nonheme Fe(IV) Oxo Complexes of Two New Pentadentate Ligands and Their Hydrogen-Atom and Oxygen-Atom Transfer Reactions. <i>Inorganic Chemistry</i> , 2015, 54, 7152-7164. | 4.0 | 63 |
| 48 | Reaction of ethyl (2Z)-cyano-6-methoxyquinolin-2(1H)-ylidene-ethanoate (L) with rhenium carbonyls: Structural and computational studies on the rhenium(I) compound cis-BrRe(CO) ₄ L. <i>Polyhedron</i> , 2015, 94, 83-89. | 2.2 | 5 |
| 49 | New azido-substituted tantalum compounds: syntheses and DFT examination of nitrogen-rich mono-, di-, and trinuclear tantalum(V) compounds. <i>Dalton Transactions</i> , 2014, 43, 3453. | 3.3 | 5 |
| 50 | A comparative study of the reactivity of the lightly stabilized cluster [Os ₃ (CO) ₈ ($\frac{1}{4}$ 3-Ph ₂ PCH ₂ P(Ph)C ₆ H ₄)($\frac{1}{4}$ -H)] towards tri(2-thienyl)-, tri(2-furyl)- and triphenyl-phosphine. <i>Journal of Organometallic Chemistry</i> , 2014, 751, 399-411. | 1.8 | 9 |
| 51 | Backbone modified small bite-angle diphosphines: Synthesis, structure, and DFT evaluation of the thermal activation products based on Os ₃ (CO) ₁₀ ($\frac{1}{4}$ -Ph ₂ PC(Me) ₂ PPh ₂). <i>Journal of Organometallic Chemistry</i> , 2014, 750, 49-58. | 1.8 | 9 |
| 52 | DFT examination of rare η^3 -SiMe ₃ abstraction in Ta(NMe ₂) ₄ [N(SiMe ₃) ₂]: formation of the imide compound Ta(η^3 -SiMe ₃)(NMe ₂) ₃ and its trapping to give guanidinate imides. <i>Dalton Transactions</i> , 2014, 43, 12390-12395. | 3.3 | 2 |
| 53 | Unusual chemical transformations of acetone thiosemicarbazone mediated by ruthenium: C-H bond activation, thiolation, and C-N bond cleavage. <i>RSC Advances</i> , 2014, 4, 1432-1440. | 3.6 | 10 |
| 54 | Synthesis and Characterization of Dimethylbis(2-pyridyl)borate Nickel(II) Complexes: Unimolecular Square-Planar to Square-Planar Rotation around Nickel(II). <i>Organometallics</i> , 2014, 33, 2019-2026. | 2.3 | 8 |

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|----|--|-----|-----------|
| 55 | Diastereomeric control of enantioselectivity: evidence for metal cluster catalysis. <i>Chemical Communications</i> , 2014, 50, 7705-7708. | 4.1 | 7 |
| 56 | Biinspired Hydrogenase Models: The Mixed-Valence Triiron Complex $[\text{Fe}_3(\text{CO})_7(\eta^4\text{-edt})_2]$ and Phosphine Derivatives $[\text{Fe}_3(\text{CO})_7\text{P}(\text{PPh}_3)_2(\eta^4\text{-edt})_2]$ ($x = 1, 2$) and $[\text{Fe}_3(\text{CO})_5(\eta^5\text{-diphosphine})(\eta^4\text{-edt})_2]$ as Proton Reduction Catalysts. <i>Organometallics</i> , 2014, 33, 1356-1366. | 2.3 | 22 |
| 57 | Synthesis of $[\text{Ru}_3(\text{CO})_9(\eta^4\text{-dppf})\{\text{P}(\text{C}_4\text{H}_3\text{E})_3\}]$ ($\text{E} = \text{O}, \text{S}$) and thermally induced cyclometalation to form $[(\eta^4\text{-H})\text{Ru}_3(\text{CO})_7(\eta^4\text{-dppf})\{\eta^4\text{-}(\text{C}_4\text{H}_3\text{E})_2\text{P}(\text{C}_4\text{H}_2\text{E})\}]$ ($\text{dppf} = 1,1'\text{-bis}(\text{diphenylphosphino})\text{ferrocene}$). <i>Journal of Organometallic Chemistry</i> , 2014, 760, 231-239. | 1.8 | 13 |
| 58 | Cycloruthenation of <i>N</i> -(Naphthyl)salicylaldimine and Related Ligands: Utilization of the Ru-C Bond in Catalytic Transfer Hydrogenation. <i>European Journal of Inorganic Chemistry</i> , 2014, 2014, 4600-4610. | 2.0 | 27 |
| 59 | Bimetallic osmium-tin complexes: Stannylene and hydrostannylene clusters upon addition of Ph_3SnH to unsaturated triosmium clusters $[(\eta^4\text{-H})_2\text{Os}_3(\text{CO})_8(\eta^4\text{-diphosphine})]$ (diphosphine = <i>dppm</i> , <i>dppf</i>). <i>Inorganica Chimica Acta</i> , 2014, 409, 320-329. | 2.4 | 21 |
| 60 | Synthesis, structure, and conformational dynamics of rhodium and iridium complexes of dimethylbis(2-pyridyl)borate. <i>Polyhedron</i> , 2014, 84, 24-31. | 2.2 | 8 |
| 61 | Experimental and computational studies on the reaction of silanes with the diphosphine-bridged triruthenium clusters $\text{Ru}_3(\text{CO})_{10}(\eta^4\text{-dppf})$, $\text{Ru}_3(\text{CO})_{10}(\eta^4\text{-dppm})$ and $\text{Ru}_3(\text{CO})_9\{\eta^4\text{-}(\text{PPhCH}_2\text{PPh}(\text{C}_6\text{H}_4))\}$. <i>Journal of Organometallic Chemistry</i> , 2014, 767, 185-195. | 1.8 | 9 |
| 62 | Models of the iron-only hydrogenase: a comparison of chelate and bridge isomers of $\text{Fe}_2(\text{CO})_4\{\text{Ph}_2\text{PN}(\text{R})\text{PPh}_2\}(\eta^4\text{-pdt})$ as proton-reduction catalysts. <i>Dalton Transactions</i> , 2013, 42, 6775. | 3.3 | 111 |
| 63 | Reaction of 4-(2,2-dimethylhydrazino)dimethylhydrazone-3(Z)-penten-2-one with $\text{BrRe}(\text{CO})_5$ and $\text{fac-BrRe}(\text{CO})_3(\text{THF})_2$: Synthesis, structural characterization, and DFT examination of the η^2 -diketimine-substituted compound $\text{fac-BrRe}(\text{CO})_3[(\text{Me}_2\text{NNCMe})_2\text{CH}_2]$. <i>Journal of Organometallic Chemistry</i> , 2013, 748, 56-62. | 1.8 | 10 |
| 64 | Structural and computational features of four highly polar quinolin-2(1H)-ylidene derivatives: Equilibrium preference for enaminothione, enamine, and enamino tautomeric structures. <i>Journal of Molecular Structure</i> , 2013, 1054-1055, 262-270. | 3.6 | 5 |
| 65 | 2-[(Diphenylphosphino)methyl]-6-methylpyridine (PN) coordination chemistry at triosmium clusters: Regiospecific ligand activation and DFT evaluation of the isomeric $\text{Os}_3(\text{CO})_{10}(\text{PN})$ clusters. <i>Journal of Organometallic Chemistry</i> , 2013, 744, 24-34. | 1.8 | 6 |
| 66 | Synthesis, Characterization, and Dynamic Behaviour of Triosmium Clusters Containing the Tridentate Ligand $\{\text{Ph}_2\text{PCH}_2\text{CH}_2\}_2\text{S}$ (PSP). <i>European Journal of Inorganic Chemistry</i> , 2013, 2013, 2447-2459. | 2.0 | 10 |
| 67 | DFT Investigation of the Mechanism of Phosphine-Thioether Isomerization in the Triosmium Cluster $\text{Os}_3(\text{CO})_{10}(\text{Ph}_2\text{PCH}_2\text{CH}_2\text{SMe})$: Migratory Preference for the Formation of an Edge-Bridged Thioether versus a Phosphine Moiety. <i>Organometallics</i> , 2012, 31, 6608-6613. | 2.3 | 11 |
| 68 | Phosphinoborane-induced fragmentation of the unsaturated hydride $\text{H}_2\text{Re}_2(\text{CO})_8$: X-ray structure of $\text{HRe}(\text{CO})_4(\eta^5\text{-P-Ph}_2\text{PCH}_2\text{CH}_2\text{BR}_2)$ (where $\text{BR}_2 = \text{borabicyclo}[3.3.1]\text{nonanyl}$) and DFT Evaluation of hydride versus CO coordination by the ancillary borane. <i>Journal of Organometallic Chemistry</i> , 2012, 700, 103-109. | 1.8 | 9 |
| 69 | Synthesis and Structural Characterization of the Pentanuclear Tantalum Cluster $\text{Ta}_5(\eta^5\text{-N})(\eta^4\text{-N})\text{Tj ETQq1 1 0.784314 rgBT /Over}$. <i>Journal of Chemical Crystallography</i> , 2012, 42, 916-922. | 1.1 | 4 |
| 70 | CO Substitution in $\text{HOs}_3(\text{CO})_{10}(\eta^4\text{-SC}_6\text{H}_4\text{Me-4})$ by the Diphosphine 4,5-Bis(diphenylphosphino)-4-cyclopentadiene-1,3-dione (bpcd): Structural and DFT Evaluation of the Isomeric Clusters $\text{HOs}_3(\text{CO})_8(\text{bpcd})(\eta^4\text{-SC}_6\text{H}_4\text{Me-4})$. <i>Journal of Cluster Science</i> , 2012, 23, 685-702. | 3.3 | 11 |
| 71 | Backbone Modified Small Bite-Angle Diphosphines: Synthesis, Structure and Regioselective Thermal Rearrangements of $\text{Os}_3(\text{CO})_{10}(\eta^4\text{-Ph}_2\text{PCH}(\text{Me})\text{PPh}_2)$. <i>Journal of Cluster Science</i> , 2012, 23, 781-798. | 3.3 | 11 |
| 72 | Ligand substitution in $1,2\text{-Os}_3(\text{CO})_{10}(\text{MeCN})_2$ by the diphosphine $(\text{PhO})_2\text{PN}(\text{Me})\text{N}(\text{Me})\text{P}(\text{OPh})_2$: X-ray diffraction structure of $1,2\text{-Os}_3(\text{CO})_{10}[(\text{PhO})_2\text{PN}(\text{Me})\text{N}(\text{Me})\text{P}(\text{OPh})_2]$ and DFT investigation of the isomeric $\text{Os}_3(\text{CO})_{10}[(\text{PhO})_2\text{PN}(\text{Me})\text{N}(\text{Me})\text{P}(\text{OPh})_2]$ clusters. <i>Journal of Molecular Structure</i> , 2012, 1010, 91-97. | 3.6 | 3 |

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