

Alireza Ariafard

List of Publications by Year in descending order

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155
papers

5,244
citations

136740

32
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163
all docs

163
docs citations

163
times ranked

5878
citing authors

#	ARTICLE	IF	CITATIONS
1	Degradation of chlorophenols by means of advanced oxidation processes: a general review. <i>Applied Catalysis B: Environmental</i> , 2004, 47, 219-256.	10.8	1,874
2	Connecting Binuclear Pd(III) and Mononuclear Pd(IV) Chemistry by Pd-Pd Bond Cleavage. <i>Journal of the American Chemical Society</i> , 2012, 134, 12002-12009.	6.6	148
3	Understanding the Relative Easiness of Oxidative Addition of Aryl and Alkyl Halides to Palladium(0). <i>Organometallics</i> , 2006, 25, 4030-4033.	1.1	140
4	Acyl Migration versus Epoxidation in Gold Catalysis: Facile, Switchable, and Atom-Economic Synthesis of Acylindoles and Quinoline Derivatives. <i>Angewandte Chemie - International Edition</i> , 2020, 59, 471-478.	7.2	99
5	η^2 -Heteroatom versus η^2 -Hydrogen Elimination: A Theoretical Study. <i>Organometallics</i> , 2006, 25, 812-819.	1.1	87
6	Nickel(IV)-Catalyzed C-H Trifluoromethylation of (Hetero)arenes. <i>Journal of the American Chemical Society</i> , 2019, 141, 12872-12879.	6.6	82
7	DFT Studies on the Mechanism of Allylative Dearomatization Catalyzed by Palladium. <i>Journal of the American Chemical Society</i> , 2006, 128, 13010-13016.	6.6	76
8	Subtle Balance of Ligand Steric Effects in Stille Transmetalation. <i>Journal of the American Chemical Society</i> , 2009, 131, 13981-13991.	6.6	76
9	Borane-Catalyzed Stereoselective C-H Insertion, Cyclopropanation, and Ring-Opening Reactions. <i>Chem</i> , 2020, 6, 2364-2381.	5.8	70
10	A Theoretical Study on the Protodeauration Step of the Gold(I)-Catalyzed Organic Reactions. <i>Organometallics</i> , 2015, 34, 3186-3195.	1.1	68
11	Effect of the Leaving Ligand X on Transmetalation of Organostannanes (vinylSnR ₃) with LnPd(Ar)(X) in Stille Cross-Coupling Reactions. A Density Functional Theory Study. <i>Organometallics</i> , 2006, 25, 5788-5794.	1.1	67
12	DFT Study on the Mechanism of the Activation and Cleavage of CO ₂ by (NHC)CuEPPh ₃ (E = Si, Ge, Sn). <i>Organometallics</i> , 2011, 30, 1340-1349.	1.1	66
13	Highly Selective Liquid-Phase Oxidation of Cyclohexane to KA Oil over Ti-MWW Catalyst: Evidence of Formation of Oxy Radicals. <i>ACS Catalysis</i> , 2014, 4, 53-62.	5.5	66
14	Oxidatively Induced C-H Activation at High Valent Nickel. <i>Journal of the American Chemical Society</i> , 2017, 139, 6058-6061.	6.6	62
15	Experimental and Computational Studies of High-Valent Nickel and Palladium Complexes. <i>Organometallics</i> , 2017, 36, 4382-4393.	1.1	55
16	Mechanism of Pd-Catalyzed Ar-Ar Bond Formation Involving Ligand-Directed C-H Arylation and Diaryliodonium Oxidants: Computational Studies of Orthopalladation at Binuclear Pd(II) Centers, Oxidation To Form Binuclear Palladium(III) Species, and Ar-Ar Reductive Coupling. <i>Organometallics</i> , 2013, 32, 544-555.	1.1	52
17	Theoretical studies of the oxidative addition of PhBr to Pd(PX ₃) ₂ and Pd(X ₂ PCH ₂ CH ₂ PX ₂) (X=Me, H, Cl). <i>Journal of Organometallic Chemistry</i> , 2007, 692, 3984-3993.	0.8	48
18	A Density Functional Theory (DFT) Mechanistic Study of Gold(I)-Catalyzed Alkynylation of the Indole and Pyrrole Substrates, Using a Hypervalent Iodine Reagent. <i>ACS Catalysis</i> , 2014, 4, 2896-2907.	5.5	47

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19	Theoretical Investigation into the Mechanism of Reductive Elimination from Bimetallic Palladium Complexes. <i>Inorganic Chemistry</i> , 2011, 50, 6449-6457.	1.9	46
20	Understanding the Highly Regioselective Cyanothiolation of 1-Alkynes Catalyzed by Palladium Phosphine Complexes. <i>Organometallics</i> , 2008, 27, 246-253.	1.1	43
21	Cleavage of Carbon Dioxide by an Iridium-Supported Fischer Carbene. A DFT Investigation. <i>Journal of the American Chemical Society</i> , 2009, 131, 5800-5808.	6.6	43
22	Experimental and Theoretical Investigation into the Gold-Catalyzed Reactivity of Cyclopropenylmethyl Acetates. <i>Organic Letters</i> , 2010, 12, 4768-4771.	2.4	43
23	Isomerism of Cp-Containing Transition Metal Allyl Complexes. <i>Organometallics</i> , 2005, 24, 680-686.	1.1	39
24	Theoretical Investigation into the Mechanism of Au(I)-Catalyzed Reaction of Alcohols with 1,5 Enynes. <i>Journal of the American Chemical Society</i> , 2012, 134, 16882-16890.	6.6	39
25	DFT Studies on the Carboxylation of the C-H Bond of Heteroarenes by Copper(I) Complexes. <i>Organometallics</i> , 2011, 30, 6218-6224.	1.1	38
26	Activation of CS ₂ and CS by ML ₃ Complexes. <i>Journal of the American Chemical Society</i> , 2008, 130, 11928-11938.	6.6	37
27	In-depth insight into the electronic and steric effects of phosphine ligands on the mechanism of the R ² R reductive elimination from (PR ₃) ₂ PdR ₂ . <i>Journal of Organometallic Chemistry</i> , 2009, 694, 2075-2084.	0.8	37
28	Ligand Effects in Bimetallic High Oxidation State Palladium Systems. <i>Inorganic Chemistry</i> , 2010, 49, 11249-11253.	1.9	37
29	Aryl-Fluoride Bond-Forming Reductive Elimination from Nickel(IV) Centers. <i>Journal of the American Chemical Society</i> , 2019, 141, 13261-13267.	6.6	37
30	Mechanism of Endo-Exo Interconversion in η^3 -Allyl Cp Complexes: A Longstanding Unresolved Issue. <i>Organometallics</i> , 2005, 24, 2241-2244.	1.1	36
31	Formation of Ethane from Mono-Methyl Palladium(II) Complexes. <i>Journal of the American Chemical Society</i> , 2014, 136, 8237-8242.	6.6	35
32	Chiral Brønsted Acid Catalyzed Enantioselective Dehydrative Nazarov-Type Electrocyclization of Aryl and 2-Thienyl Vinyl Alcohols. <i>Journal of the American Chemical Society</i> , 2018, 140, 5834-5841.	6.6	33
33	Gas-Phase Ion-Molecule Reactions of Copper Hydride Anions [CuH ₂] ⁻ and [Cu ₂ H ₃] ⁻ . <i>Inorganic Chemistry</i> , 2017, 56, 2387-2399.	1.9	32
34	Triarylborane-Catalyzed Alkenylation Reactions of Aryl Esters with Diazo Compounds. <i>Angewandte Chemie - International Edition</i> , 2020, 59, 15492-15496.	7.2	32
35	Theoretical Investigation into the Mechanism of 3d-GMP Oxidation by [Pt ^{IV} Cl ₄](dach)]. <i>Inorganic Chemistry</i> , 2013, 52, 707-717.	1.9	31
36	Insight into the Mechanism of R ² R Reductive Elimination from the Six-Coordinate d ⁶ Complexes L ₂ Pt(R) ₄ (R = vinyl, Me). <i>Organometallics</i> , 2011, 30, 422-432.	1.1	29

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37	DFT mechanistic investigation into phenol dearomatization mediated by an iodine (<sc>iii</sc>) reagent. <i>Organic and Biomolecular Chemistry</i> , 2019, 17, 3521-3528.	1.5	29
38	Hydrogen peroxide decomposition over La _{0.9} Sr _{0.1} Ni ¹⁺ Cr O ₃ perovskites. <i>Catalysis Communications</i> , 2003, 4, 561-566.	1.6	28
39	Mechanistic Elucidation of the Arylation of Non-Spectator <i>N</i> -Heterocyclic Carbenes at Copper Using a Combined Experimental and Computational Approach. <i>Organometallics</i> , 2015, 34, 3497-3507.	1.1	28
40	Gold-Catalyzed Regiospecific Annulation of Unsymmetrically Substituted 1,5-Diynes for the Precise Synthesis of Bispentalenes. <i>Chemistry - A European Journal</i> , 2019, 25, 12180-12186.	1.7	28
41	Dual Gold-Catalyzed Cycloaromatization of Unconjugated (<i>E</i>)-Enediynes. <i>Angewandte Chemie - International Edition</i> , 2019, 58, 2114-2119.	7.2	28
42	Copper-catalysed synthesis of $\hat{\text{I}}$ -alkylidene cyclic carbonates from propargylic alcohols and CO ₂ . <i>Green Chemistry</i> , 2021, 23, 889-897.	4.6	28
43	Site-Selective C(sp ³) ₃ -C(sp ³) ₃ -C(sp ³) ₃ -C(sp ³) ₂ Cross-Coupling Reactions Using Frustrated Lewis Pairs. <i>Journal of the American Chemical Society</i> , 2021, 143, 4451-4464.	6.6	28
44	Assembly of Cyclometalated Platinum(II) Complexes via 1,1-Bis(diphenylphosphino)ferrocene Ligand: Kinetics and Mechanisms. <i>Organometallics</i> , 2011, 30, 1466-1477.	1.1	27
45	Theoretical Studies of Cycloaddition Reactions of Cationic Aluminum $\hat{\text{I}}$ -Diketiminato Alkyl Complexes with Alkenes and Alkynes. <i>Organometallics</i> , 2005, 24, 5140-5146.	1.1	26
46	Synthesis of Double-End-Capped Polyethylene by a Cationic Tris(pyrazolyl)borate Zirconium Benzyl Complex. <i>Organometallics</i> , 2008, 27, 5867-5875.	1.1	26
47	Synthesis, structure and gas-phase reactivity of the mixed silver hydride borohydride nanocluster [Ag ₃ ($\hat{\text{I}}$ ₃ -H)($\hat{\text{I}}$ ₃ -BH ₄)L ³ Ph ₃]BF ₄ . <i>Journal of the American Chemical Society</i> , 2016, 138, 1261-1269.	12.4	26
48	Computational study of C(sp ³) $\hat{\text{C}}$ -O bond formation at a Pd ^{IV} centre. <i>Dalton Transactions</i> , 2017, 46, 3742-3748.	1.6	25
49	Gold-catalyzed domino cyclization $\hat{\text{C}}$ -alkynylation reactions with EBX reagents: new insights into the reaction mechanism. <i>Dalton Transactions</i> , 2017, 46, 12257-12262.	1.6	25
50	Nickel(II/IV) Manifold Enables Room-Temperature C(sp ³) $\hat{\text{C}}$ -H Functionalization. <i>Journal of the American Chemical Society</i> , 2019, 141, 19513-19520.	6.6	25
51	Decarboxylative-Coupling of Allyl Acetate Catalyzed by Group 10 Organometallics, [(phen)M(CH ₃) ₃] ⁺ . <i>Journal of Organic Chemistry</i> , 2014, 79, 12056-12069.	1.7	24
52	A one-pot route to thioamides discovered by gas-phase studies: palladium-mediated CO ₂ extrusion followed by insertion of isothiocyanates. <i>Chemical Communications</i> , 2017, 53, 3854-3857.	2.2	24
53	Computational Study of Intramolecular Arene Palladation at a Palladium(IV) Center. <i>Organometallics</i> , 2015, 34, 1085-1090.	1.1	23
54	Density Functional Theory Study on the Mechanism of the Reductive Cleavage of CO ₂ by a Bis- $\hat{\text{I}}$ -Diketoiminatodiiron Dinitrogen Complex. <i>Inorganic Chemistry</i> , 2010, 49, 7773-7782.	1.9	22

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55	A Mechanistic Investigation of the Gold(III)-Catalyzed Hydrofurylation of C=C Multiple Bonds. <i>Journal of the American Chemical Society</i> , 2016, 138, 14599-14608.	6.6	22
56	DFT Mechanistic Investigation into BF ₃ -Catalyzed Alcohol Oxidation by a Hypervalent Iodine(III) Compound. <i>ACS Catalysis</i> , 2019, 9, 6510-6521.	5.5	22
57	Density functional studies of influences of Ni triad metals and solvents on oxidative addition of MeI to [M(CH ₃) ₂ (NH ₃) ₂] complexes and C=C reductive elimination from [M(CH ₃) ₃ (NH ₃) ₂] complexes. <i>Journal of Organometallic Chemistry</i> , 2011, 696, 3351-3358.	0.8	21
58	Theoretical Study on the Ring-Opening Reactions of Cyclopropenes Mediated by a Au ^I Complex. <i>Journal of Organic Chemistry</i> , 2013, 78, 9553-9559.	1.7	21
59	Theoretical Investigation into the Mechanism of Cyanomethylation of Aldehydes Catalyzed by a Nickel Pincer Complex in the Absence of Base Additives. <i>ACS Catalysis</i> , 2016, 6, 60-68.	5.5	21
60	Computational Mechanistic Study of Palladium(II)-Catalyzed Carboxyalkynylation of an Olefin Using an Iodine(III) Oxidant Reagent. <i>Organometallics</i> , 2014, 33, 7318-7324.	1.1	20
61	A transition-metal-free fast track to flavones and 3-arylcoumarins. <i>Chemical Communications</i> , 2017, 53, 10676-10679.	2.2	20
62	Synthetic and computational studies of the palladium(IV) system Pd(alkyl)(aryl)(alkynyl)(bidentate)(triflate) exhibiting selectivity in C=C reductive elimination. <i>Dalton Transactions</i> , 2012, 41, 11820.	1.6	19
63	Experimental and computational evidence for KOt-Bu-promoted synthesis of oxopyrazino[1,2-a]indoles. <i>RSC Advances</i> , 2015, 5, 101353-101361.	1.7	19
64	Total Synthesis of (±)-Dihydroisobamamol. <i>Advanced Synthesis and Catalysis</i> , 2017, 359, 866-874.	2.1	19
65	A computational mechanistic investigation into the reduction of Pt(IV) prodrugs with two axial chlorides by biological reductants. <i>Chemical Communications</i> , 2017, 53, 1413-1416.	2.2	19
66	Gold Catalyzed Cyclopropanation/[5+3] Cycloaddition of 1,4,9- and 1,4,10-Allenynes to Bicyclo[3.3.1]nonane Derivatives. <i>Advanced Synthesis and Catalysis</i> , 2020, 362, 1084-1095.	2.1	19
67	In-depth insight into metal-alkene bonding interactions. <i>Inorganica Chimica Acta</i> , 2006, 359, 3527-3534.	1.2	18
68	The different roles of a cationic gold(I) complex in catalysing hydroarylation of alkynes and alkenes with a heterocycle. <i>Chemical Communications</i> , 2016, 52, 9422-9425.	2.2	18
69	Different Selectivities in the Insertions into C(sp ²)-H Bonds: Benzofulvenes by Dual Gold Catalysis Competition Experiments. <i>Chemistry - A European Journal</i> , 2018, 24, 10766-10772.	1.7	18
70	Mechanistic investigation into phenol oxidation by IBX elucidated by DFT calculations. <i>Organic and Biomolecular Chemistry</i> , 2020, 18, 1117-1129.	1.5	18
71	Tris(pentafluorophenyl)borane-catalyzed Carbenium Ion Generation and Autocatalytic Pyrazole Synthesis: A Computational and Experimental Study. <i>Angewandte Chemie - International Edition</i> , 2021, 60, 24395-24399.	7.2	18
72	Reduction of a platinum(IV) prodrug model by sulfur containing biological reductants: computational mechanistic elucidation. <i>Chemical Communications</i> , 2018, 54, 10491-10494.	2.2	17

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73	Gold-Catalyzed [5,5]-Rearrangement. ACS Catalysis, 2021, 11, 6510-6518.	5.5	17
74	Factors Dictating Carbene Formation at (PNP)Ir. Organometallics, 2010, 29, 4239-4250.	1.1	16
75	Synthesis of Amidines by Palladium-Mediated CO ₂ Extrusion Followed by Insertion of Carbodiimides: Translating Mechanistic Studies to Develop a One-Pot Method. Organometallics, 2019, 38, 424-435.	1.1	16
76	Structure and Bonding of d ⁸ Allyl Complexes M(η -3-allyl)L ₃ (M = Co, Rh, Ir; L = Phosphine or Carbonyl). Organometallics, 2005, 24, 3800-3806.	1.1	14
77	Theoretical Investigation into the Palladium-Catalyzed Silaboration of Pyridines. Organometallics, 2012, 31, 1680-1687.	1.1	14
78	Mechanistic Insights into the Oxidative Coupling of N-Heterocyclic Carbenes within the Coordination Sphere of Copper Complexes. Chemistry - A European Journal, 2014, 20, 12729-12733.	1.7	14
79	A Computational Mechanistic Investigation into Reduction of Gold(III) Complexes by Amino Acid Glycine: A New Variant for Amine Oxidation. Chemistry - A European Journal, 2018, 24, 8361-8368.	1.7	14
80	A Molecular Orbital Rationalization of Ligand Effects in N ₂ Activation. Chemistry - A European Journal, 2008, 14, 6119-6124.	1.7	13
81	Reactivity of CO ₂ towards Mo[N(R)Ph] ₃ . Dalton Transactions, 2009, , 9266.	1.6	13
82	Tuning the Laplaza-Cummins 3-coordinate M[N(R)Ph] ₃ catalyst to activate and cleave CO ₂ . Dalton Transactions, 2011, 40, 5569.	1.6	12
83	Proton supplier role of binuclear gold complexes in promoting hydrofunctionalisation of nonactivated alkenes. Catalysis Science and Technology, 2019, 9, 1420-1426.	2.1	11
84	The role of hypervalent iodine(ⁱⁱⁱ) reagents in promoting alkoxylation of unactivated C(sp ³)-H bonds catalyzed by palladium(ⁱⁱ) complexes. Chemical Science, 2021, 12, 7185-7195.	3.7	11
85	Understanding the Influence of Donor-Acceptor Diazo Compounds on the Catalyst Efficiency of B(C ₆ F ₅) ₃ Towards Carbene Formation. Chemistry - A European Journal, 2022, 28, .	1.7	11
86	Titanium-mediated rearrangement of cyclopropenylmethyl acetates to (E)-halodienes. Organic and Biomolecular Chemistry, 2011, 9, 3359.	1.5	10
87	Density Functional Theory Studies on the Oxidation of 5 ² -dGMP and 5 ² -dAMP by a Platinum(IV) Complex. Inorganic Chemistry, 2012, 51, 8002-8013.	1.9	10
88	Theoretical rationalisation for the mechanism of N-heterocyclic carbene-halide reductive elimination at Cu ^{III} , Ag ^{III} and Au ^{III} . Chemical Communications, 2016, 52, 5057-5060.	2.2	10
89	A Modified Cationic Mechanism for PdCl ₂ -Catalyzed Transformation of a Homoallylic Alcohol to an Allyl Ether. Organometallics, 2019, 38, 2953-2962.	1.1	10
90	Disclosure of Some Obscure Mechanistic Aspects of the Copper-Catalyzed Click Reactions Involving N ₂ Elimination Promoted by the Use of Electron-Deficient Azides from a DFT Perspective. Organometallics, 2019, 38, 256-267.	1.1	10

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91	Iron Triflate Salts as Highly Active Catalysts for the Solvent-Free Oxidation of Cyclohexane. <i>European Journal of Organic Chemistry</i> , 2020, 2020, 3552-3559.	1.2	10
92	Copper(I)-catalysed site-selective C(sp ³)-H bond chlorination of ketones, (E)-enones and alkylbenzenes by dichloramine-T. <i>Nature Communications</i> , 2021, 12, 4065.	5.8	10
93	Borane Catalyzed Selective Diazo Cross-Coupling Towards Pyrazoles. <i>Advanced Synthesis and Catalysis</i> , 2022, 364, 773-780.	2.1	10
94	Density functional theory study on structural isomers and bonding of model complexes M(CO) ₅ (BH ₃ -PH ₃) (M=Cr, Mo, W) and W(CO) ₅ (BH ₃ -AH ₃) (A=N, P, As, Sb). <i>Journal of Organometallic Chemistry</i> , 2005, 690, 1147-1156.	0.8	9
95	Theoretical Investigation into the Mechanism of Ethylene Polymerization by a Cationic Dinuclear Aluminum Complex: A Longstanding Unresolved Issue. <i>Organometallics</i> , 2013, 32, 1687-1693.	1.1	9
96	Computational Study of Carbostannylation Implicating Bimetallic Catalysis Involving σ -Au ^I -Vinyl-Pd ^{II} -Species. <i>ACS Catalysis</i> , 2014, 4, 860-869.	5.5	9
97	Prying open a Reactive Site for Allylic Arylation by Phosphine-Ligated Geminally Diaurated Aryl Complexes. <i>Organometallics</i> , 2015, 34, 3255-3263.	1.1	9
98	Mechanistic Elucidation of Gold(I)-Catalyzed Oxidation of a Propargylic Alcohol by a N-Oxide in the Presence of an Imine Using DFT Calculations. <i>Organometallics</i> , 2019, 38, 489-497.	1.1	9
99	Chiral Gold Complex Catalyzed Cycloisomerization/Regio- and Enantioselective Nitroso-Diels-Alder Reaction of 1,6-Diyne Esters with Nitrosobenzenes. <i>ACS Catalysis</i> , 2022, 12, 7288-7299.	5.5	9
100	Theoretical study on interaction of different coordination modes of BH ₄ ligand with transition metal in [TM(BH ₄)(CO) ₄] ⁻ (TM=Cr, Mo). <i>Journal of Organometallic Chemistry</i> , 2005, 690, 84-95.	0.8	8
101	Mechanistic Studies of Ligand Fluxionality in [M(η -5-Cp)(η -1-Cp)(L) ₂] _n . <i>Journal of Physical Chemistry A</i> , 2009, 113, 2982-2989.	1.1	8
102	Gas-Phase and Computational Study of Identical Nickel- and Palladium-Mediated Organic Transformations Where Mechanisms Proceeding via M ^{II} or M ^{IV} Oxidation States Are Determined by Ancillary Ligands. <i>Journal of the American Chemical Society</i> , 2015, 137, 13588-13593.	6.6	8
103	DFT-Based Comparison between Mechanistic Aspects of Amine and Alcohol Oxidation Mediated by IBX. <i>Journal of Organic Chemistry</i> , 2020, 85, 515-525.	1.7	8
104	Triarylborane catalysed <i>N</i> -alkylation of amines with aryl esters. <i>Catalysis Science and Technology</i> , 2020, 10, 7523-7530.	2.1	8
105	Bidentate Nitrogen-Ligated I(V) Reagents, Bi(<i>N</i>)-HVLs: Preparation, Stability, Structure, and Reactivity. <i>Journal of Organic Chemistry</i> , 2021, 86, 6566-6576.	1.7	8
106	DFT study of metal-tetrahydroborato ligand interactions in [Ti(CO) ₄ (BH ₄)] ⁻ . <i>Computational and Theoretical Chemistry</i> , 2003, 625, 305-314.	1.5	7
107	DFT studies of structural preference of coordinated ethylene in W(CO) ₃ (PX ₃) ₂ (CH ₂ -...-CH ₂) (X=H, CH ₃ , F,) Tj ETQg1.1 0.784314 rgB /	0.8	7
108	Scission of Carbon Monoxide Using TaR ₃ , R=(<i>t</i> -Bu)Ph or OSi(<i>t</i> -Bu) ₃ : A DFT Investigation. <i>Chemistry - A European Journal</i> , 2010, 16, 8117-8132.	1.7	7

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109	Oxo iron(IV) as an oxidative active intermediate of p-chlorophenol in the Fenton reaction: a DFT study. <i>Physical Chemistry Chemical Physics</i> , 2012, 14, 3766.	1.3	7
110	Understanding the mechanism of CuI-catalyzed N-H carboxylation of heterocyclic rings with CO ₂ from a theoretical point of view. <i>Journal of Organometallic Chemistry</i> , 2013, 748, 89-97.	0.8	7
111	Dual Gold-Catalyzed Cycloaromatization of Unconjugated (E)-Enediyne. <i>Angewandte Chemie</i> , 2019, 131, 2136-2141.	1.6	7
112	Theoretical studies of rotational barriers of dithiocarbamate ligands in the square planar complexes TM(L)(L ²)(H ₂ dtc) (TM=Ir, Rh). <i>Computational and Theoretical Chemistry</i> , 2003, 636, 49-56.	1.5	6
113	Does type of phosphine affect rotational barrier of vinylidene in the complexes OsHCl(CCH ₂)(L) ₂ (L=phosphine)? <i>Inorganic Chemistry Communication</i> , 2004, 7, 999-1002.	1.8	6
114	Sulfur Dioxide Activation: A Theoretical Investigation into Dual S-O Bond Cleavage by Three-Coordinate Molybdenum(III) Complexes. <i>Inorganic Chemistry</i> , 2015, 54, 534-543.	1.9	6
115	η ³ -Alkenyl endo-palladacycle formation via regiospecific functionalisation of an unreactive NHC-tethered C(sp ²)-H bond. <i>Chemical Communications</i> , 2015, 51, 5513-5515.	2.2	6
116	Computational study of selectivity in the [Pt ^{II} Cl ₄] ²⁻ -catalysed arylation of arenes by diaryliodonium reagents: arene activation at Pt ^{IV} centres. <i>Dalton Transactions</i> , 2017, 46, 15480-15486.	1.6	6
117	Substituent effects in the decarboxylation reactions of coordinated arylcarboxylates in dinuclear copper complexes, [(napy) ₂ Cu ₂ (O) ₂ CC ₆ H ₄ X] ⁺ . <i>European Journal of Mass Spectrometry</i> , 2017, 23, 351-358.	0.5	6
118	Nazarov cyclisations initiated by DDQ-oxidised pentadienyl ether: a mechanistic investigation from the DFT perspective. <i>Organic and Biomolecular Chemistry</i> , 2018, 16, 9021-9029.	1.5	6
119	Accessing Chelating Extended Linker Bis(NHC) Palladium(II) Complexes: Sterically Triggered Divergent Reaction Pathways. <i>Organometallics</i> , 2019, 38, 3032-3038.	1.1	6
120	Computational Analysis of Mesomerism in <i>para</i> -Substituted <i>mer</i> -NCN Pincer Platinum(II) Complexes, Including its Relationships with Hammett <i>p</i> -Substituent Parameters. <i>Chemistry - A European Journal</i> , 2020, 26, 15629-15635.	1.7	6
121	Gold-Catalyzed Annulation of 1,8-Dialkynyl naphthalenes: Synthesis and Photoelectric Properties of Indenophenylene-Based Derivatives. <i>Chemistry - A European Journal</i> , 2021, 27, 3552-3559.	1.7	6
122	Hydroalkylation of Alkenes with 1,3-Diketones via Gold(III) or Silver(I) Catalysis: Divergent Mechanistic Pathways Revealed by a DFT-Based Investigation. <i>ACS Catalysis</i> , 2021, 11, 5795-5807.	5.5	6
123	Oxidation of Electron-Deficient Phenols Mediated by Hypervalent Iodine(V) Reagents: Fundamental Mechanistic Features Revealed by a Density Functional Theory-Based Investigation. <i>Journal of Organic Chemistry</i> , 2021, 86, 12237-12246.	1.7	6
124	Exploring Cyclization Strategies to Access Stemona Alkaloids: Subtle Effects Influencing Reactivity in Intramolecular Michael Additions. <i>Organic Letters</i> , 2021, 23, 8494-8498.	2.4	6
125	Stability of Heavier Group 14 Analogues of Vinylidene Complexes: A Theoretical Study. <i>Organometallics</i> , 2005, 24, 6283-6286.	1.1	5
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