

Gernot Frenking

List of Publications by Year in descending order

Source: <https://exaly.com/author-pdf/7989546/publications.pdf>

Version: 2024-02-01

294
papers

26,235
citations

3721

89
h-index

8370

147
g-index

319
all docs

319
docs citations

319
times ranked

8462
citing authors

#	ARTICLE	IF	CITATIONS
1	The strength of a chemical bond. <i>International Journal of Quantum Chemistry</i> , 2022, 122, e26773.	1.0	29
2	Neutral and charged group 13-16 homologs of carbon EL2 (E = B, Al, Ga, In, Tl; Si, Ge, Sn, Pb; N, P, As, Sb, Bi; O, S, Se, Te) _{0.4} . <i>Advances in Inorganic Chemistry</i> , 2022, , 247-299.		3
3	Bonding analysis of the C ₂ precursor Me ₃ C-C ₂ -I(Ph)FBF ₃ (E=Al, Si, Ge). <i>Pure and Applied Chemistry</i> , 2022, 94, 1-10.		1
4	Isolation of Stable Borepin Radicals and Anions. <i>Angewandte Chemie - International Edition</i> , 2022, 61, .	7.2	15
5	Isolation of Stable Borepin Radicals and Anions. <i>Angewandte Chemie</i> , 2022, 134, .	1.6	5
6	π-Back-Donation from a Beryllium Dibromide Fragment at the Expense of Its σ Strength. <i>Inorganic Chemistry</i> , 2022, 61, 700-705.	1.9	14
7	Carbodiphosphorane-Stabilized Parent Dioxophosphorane: A Valuable Synthetic HO ₂ P Source. <i>Journal of the American Chemical Society</i> , 2022, 144, 7357-7365.	6.6	7
8	Generation and Characterization of the Charge-Transferred Diradical Complex CaCO ₂ with an Open-Shell Singlet Ground State. <i>Journal of the American Chemical Society</i> , 2022, 144, 8355-8361.	6.6	8
9	Clarifying notes on the bonding analysis adopted by the energy decomposition analysis. <i>Physical Chemistry Chemical Physics</i> , 2022, 24, 15726-15735.	1.3	25
10	The nature of the polar covalent bond. <i>Journal of Chemical Physics</i> , 2022, 157, .	1.2	15
11	Triple bonding between beryllium and nitrogen in HNB ₂ CO. <i>Chemical Communications</i> , 2022, 58, 8532-8535.	2.2	5
12	CO ₂ -Induced Dinitrogen Fixation and Cleavage Mediated by Boron. <i>Chemistry - A European Journal</i> , 2021, 27, 2131-2137.	1.7	20
13	Isolable dicarbon stabilized by a single phosphine ligand. <i>Nature Chemistry</i> , 2021, 13, 89-93.	6.6	15
14	Generation and Characterization of the C ₃ O ₂ ⁻ Anion with an Unexpected Unsymmetrical Structure. <i>Angewandte Chemie</i> , 2021, 133, 4568-4573.	1.6	0
15	Generation and Characterization of the C ₃ O ₂ ⁻ Anion with an Unexpected Unsymmetrical Structure. <i>Angewandte Chemie - International Edition</i> , 2021, 60, 4518-4523.	7.2	7
16	Carbodicarbene Bismaalkene Cations: Unravelling the Complexities of Carbene versus Carbene in Heavy Pnictogen Chemistry. <i>Angewandte Chemie - International Edition</i> , 2021, 60, 6682-6690.	7.2	40
17	Carbodicarbene Bismaalkene Cations: Unravelling the Complexities of Carbene versus Carbene in Heavy Pnictogen Chemistry. <i>Angewandte Chemie</i> , 2021, 133, 6756-6764.	1.6	15
18	Generation and Identification of the Linear OCBNO and OBNCO Molecules with 24 Valence Electrons. <i>Chemistry - A European Journal</i> , 2021, 27, 412-418.	1.7	8

#	ARTICLE	IF	CITATIONS
19	Donor-Stabilized Antimony(I) and Bismuth(I) Ions: Heavier Valence Isoelectronic Analogues of Carbones. <i>Journal of the American Chemical Society</i> , 2021, 143, 1301-1306.	6.6	40
20	Dinitrogen complexation and reduction at low-valent calcium. <i>Science</i> , 2021, 371, 1125-1128.	6.0	131
21	Chemical Bonding in Homoleptic Carbonyl Cations $[M\{Fe(CO)_5\}_2]^{+}$ (M=Cu, Ag, Au). <i>Chemistry - A European Journal</i> , 2021, 27, 6936-6944.	1.7	13
22	Metal-CO Bonding in Mononuclear Transition Metal Carbonyl Complexes. <i>Jacs Au</i> , 2021, 1, 623-645.	3.6	57
23	Highly Coordinated Heteronuclear Calcium-Iron Carbonyl Cation Complexes $[CaFe(CO)_n]^{+}$ ($n=5-12$) with $d\pi$ Bonding. <i>Angewandte Chemie</i> , 2021, 133, 13984-13989.	1.6	0
24	Bonding in $M(NHBMe)_2$ and $M[Mn(CO)_5]_2$ complexes (M=Zn, Cd, Hg; $NHBMe=(HCNMe)_2B$): divalent group 12 metals with zero oxidation state. <i>Theoretical Chemistry Accounts</i> , 2021, 140, 1.	0.5	4
25	Highly Coordinated Heteronuclear Calcium-Iron Carbonyl Cation Complexes $[CaFe(CO)_n]^{+}$ ($n=5-12$) with $d\pi$ Bonding. <i>Angewandte Chemie - International Edition</i> , 2021, 60, 13865-13870.	7.2	18
26	Isolation of a Uranium(III)-Carbon Multiple Bond Complex. <i>Chemistry - A European Journal</i> , 2021, 27, 10006-10011.	1.7	12
27	Transition-Metal Chemistry of the Heavier Alkaline Earth Atoms Ca, Sr, and Ba. <i>Accounts of Chemical Research</i> , 2021, 54, 3071-3082.	7.6	39
28	A Critical Look at Linus Pauling's Influence on the Understanding of Chemical Bonding. <i>Molecules</i> , 2021, 26, 4695.	1.7	7
29	Covalent Bonding Between Be^{+} and CO_2 in $BeOCO^{+}$ with a Surprisingly High Antisymmetric OCO Stretching Vibration. <i>Journal of the American Chemical Society</i> , 2021, 143, 14300-14305.	6.6	10
30	Revisiting the Bonding Scenario of Two Donor Ligand Stabilized C_2 Species. <i>Journal of Physical Chemistry A</i> , 2021, 125, 291-301.	1.1	16
31	Persistent Borafluorene Radicals. <i>Angewandte Chemie - International Edition</i> , 2020, 59, 3850-3854.	7.2	70
32	Persistent Borafluorene Radicals. <i>Angewandte Chemie</i> , 2020, 132, 3878-3882.	1.6	19
33	Quadruple bonding of bare group-13 atoms in transition metal complexes. <i>Dalton Transactions</i> , 2020, 49, 14815-14825.	1.6	7
34	The bonding situation in heterometallic carbonyl complexes. <i>Dalton Transactions</i> , 2020, 49, 16762-16771.	1.6	3
35	The Valence Orbitals of the Alkaline Earth Atoms. <i>Chemistry - A European Journal</i> , 2020, 26, 14194-14210.	1.7	39
36	A diradical based on odd-electron f -bonds. <i>Nature Communications</i> , 2020, 11, 3441.	5.8	22

#	ARTICLE	IF	CITATIONS
37	d σ -d π Dative Bonding Between Iron and the Alkaline-Earth Metals Calcium, Strontium, and Barium. <i>Angewandte Chemie</i> , 2020, 132, 14723-14728.	1.6	7
38	Stabilization of Linear C ₃ by Two Donor Ligands: A Theoretical Study of L ₂ C ₃ L (L=PPh ₃ , NHC Me, cAAC) <i>J. Phys. Chem. Lett.</i> , 2020, 11, 1000-1004.	1.7	21
39	Carbones and Carbon Atom as Ligands in Transition Metal Complexes. <i>Molecules</i> , 2020, 25, 4943.	1.7	43
40	Group-6 Hexacarbonyls as Ligands for the Silver Cation: Syntheses, Characterization, and Analysis of the Bonding Compared with the Isolectronic Group-5 Hexacarbonylates. <i>Chemistry - A European Journal</i> , 2020, 26, 17203-17211.	1.7	16
41	Beryllium Atom Mediated Dinitrogen Activation via Coupling with Carbon Monoxide. <i>Angewandte Chemie</i> , 2020, 132, 18358-18364.	1.6	3
42	Comment on "Topological Analysis of the Electron Density in the Carbonyl Complexes M(CO) ₈ (M = Ca, Sr, Ba)" <i>Organometallics</i> , 2020, 39, 2956-2958.	1.1	6
43	Di-ortho-beryllated Carbodiphosphorane: A Compound with a Metal-Carbon Double Bond to an Element of the s-Block. <i>Organometallics</i> , 2020, 39, 3224-3231.	1.1	44
44	d σ -d π Dative Bonding Between Iron and the Alkaline-Earth Metals Calcium, Strontium, and Barium. <i>Angewandte Chemie - International Edition</i> , 2020, 59, 14615-14620.	7.2	30
45	Synthesis and characterization of heterometallic complexes involving coinage metals and isoelectronic Fe(CO) ₅ , [Mn(CO) ₅] ⁺ and [Fe(CO) ₄ CN] ⁺ ligands. <i>Dalton Transactions</i> , 2020, 49, 8566-8581.	1.6	13
46	Alkaline Earth Metals Activate N ₂ and CO in Cubic Complexes Just Like Transition Metals: A Conceptual Density Functional Theory and Energy Decomposition Analysis Study. <i>Chemistry - A European Journal</i> , 2020, 26, 12785-12793.	1.7	20
47	Filling a Gap: The Coordinatively Saturated Group-4 Carbonyl Complexes TM(CO) ₈ (TM=Zr) <i>J. Phys. Chem. Lett.</i> , 2020, 11, 1000-1004.	1.7	21
48	Side-On Bonded Beryllium Dinitrogen Complexes. <i>Angewandte Chemie - International Edition</i> , 2020, 59, 10603-10609.	7.2	51
49	Beryllium Atom Mediated Dinitrogen Activation via Coupling with Carbon Monoxide. <i>Angewandte Chemie - International Edition</i> , 2020, 59, 18201-18207.	7.2	29
50	Side-On Bonded Beryllium Dinitrogen Complexes. <i>Angewandte Chemie</i> , 2020, 132, 10690-10696.	1.6	13
51	Double donation in trigonal planar iron-carbodiphosphorane complexes - a concise study on their spectroscopic and electronic properties. <i>Dalton Transactions</i> , 2020, 49, 2537-2546.	1.6	20
52	Comment on "Revisiting π backbonding: the influence of d orbitals on metal-CO bonds and ligand red shifts" by D. Koch, Y. Chen, P. Golub and S. Manzhos, <i>Phys. Chem. Chem. Phys.</i> , 2019, 21, 20814. <i>Physical Chemistry Chemical Physics</i> , 2020, 22, 5377-5379.	1.3	7
53	A Stable, Crystalline Beryllium Radical Cation. <i>Journal of the American Chemical Society</i> , 2020, 142, 4560-4564.	6.6	80
54	Isolable cyclic radical cations of heavy main-group elements. <i>Chemical Communications</i> , 2020, 56, 2167-2170.	2.2	21

#	ARTICLE	IF	CITATIONS
55	Bonding Analysis of the Shortest Bond between Two Atoms Heavier than Hydrogen and Helium: O ₂ ²⁺ . <i>Journal of Physical Chemistry A</i> , 2020, 124, 1087-1092.	1.1	12
56	Comment on "Realization of Lewis Basic Sodium Anion in the NaBH ₃ ⁺ Cluster". <i>Angewandte Chemie</i> , 2020, 132, 8836-8839.	1.6	16
57	Comment on "Realization of Lewis Basic Sodium Anion in the NaBH ₃ ⁺ Cluster". <i>Angewandte Chemie - International Edition</i> , 2020, 59, 8756-8759.	7.2	43
58	Response to Comment on "Observation of alkaline earth complexes M(CO) ₈ (M = Ca, Sr, Ba)". <i>Journal of Physical Chemistry A</i> , 2020, 124, 1087-1092.	6.0	54
59	Octa-coordinated alkaline earth metal dinitrogen complexes M(N ₂) ₈ (M=Ca, Sr, Ba). <i>Nature Communications</i> , 2019, 10, 3375.	5.8	79
60	Nine questions on energy decomposition analysis. <i>Journal of Computational Chemistry</i> , 2019, 40, 2248-2283.	1.5	113
61	Chemical Bonding and Bonding Models of Main-Group Compounds. <i>Chemical Reviews</i> , 2019, 119, 8781-8845.	23.0	232
62	Octacarbonyl Anion Complexes of Actinides [An(CO) ₈] ⁻ (An=Th, U) and the Role of f Orbitals in Metal-Ligand Bonding. <i>Chemistry - A European Journal</i> , 2019, 25, 11772-11784.	1.7	38
63	Transition-Metal Chemistry of Alkaline-Earth Elements: The Trisbenzene Complexes M(Bz) ₃ (M=Ca, Sr, Ba). <i>Angewandte Chemie - International Edition</i> , 2019, 58, 17365-17374.	7.2	82
64	Transition-Metal Chemistry of Alkaline-Earth Elements: The Trisbenzene Complexes M(Bz) ₃ (M=Ca, Sr, Ba). <i>Angewandte Chemie</i> , 2019, 131, 17526-17535.	1.6	28
65	Alkali Metal Covalent Bonding in Nickel Carbonyl Complexes ENi(CO) ₃ ⁺ . <i>Angewandte Chemie</i> , 2019, 131, 1746-1752.	1.6	53
66	Synthesis of cAAC stabilized biradical of Me ₂ Si• and Me ₂ SiCl• monoradical from Me ₂ SiCl ₂ an important feedstock material. <i>Chemical Communications</i> , 2019, 55, 4534-4537.	2.2	9
67	Chemical bonding in the hexamethylbenzene-SO ₂ ⁺ dication. <i>Theoretical Chemistry Accounts</i> , 2019, 138, 1.	0.5	20
68	Bent Phosphaallenes With "Hidden" Lone Pairs as Ligands. <i>Chemistry - A European Journal</i> , 2019, 25, 7912-7920.	1.7	2
69	An Experimental and Theoretical Study of the Structures and Properties of [CDP ⁺ Me] ⁺ Ni(CO) ₃ and [Ni ₂ (CO) ₄](μ ₂ -CO)(μ ₂ -CDP ⁺ Me)]. <i>European Journal of Inorganic Chemistry</i> , 2019, 2019, 4546-4554.	1.0	13
70	Cerium-carbon dative interactions supported by carbodiphosphorane. <i>Dalton Transactions</i> , 2019, 48, 16108-16114.	1.6	20
71	The Diels-Alder Reaction from the EDA-NOCV Perspective: A Re-Examination of the Frontier Molecular Orbital Model. <i>European Journal of Organic Chemistry</i> , 2019, 2019, 478-485.	1.2	10
72	Octacarbonyl Anion Complexes of the Late Lanthanides Ln(CO) ₈ ⁻ (Ln=Tm, Yb). <i>Journal of Physical Chemistry A</i> , 2020, 124, 1087-1092.	1.7	38

#	ARTICLE	IF	CITATIONS
73	The Lewis electron-pair bonding model: the physical background, one century later. <i>Nature Reviews Chemistry</i> , 2019, 3, 35-47.	13.8	52
74	The Lewis electron-pair bonding model: modern energy decomposition analysis. <i>Nature Reviews Chemistry</i> , 2019, 3, 48-63.	13.8	197
75	Dative versus electron-sharing bonding in <i>N</i> -imides and phosphane imides R_3ENX and relative energies of the $R_2EN(X)R$ isomers (E = N, P; R = H, Cl, Me, Ph; X = H, F, Cl). <i>Molecular Physics</i> , 2019, 117, 1306-1314.	0.8	14
76	Isolation of Transient Acyclic Germanium(I) Radicals Stabilized by Cyclic Alkyl(amino) Carbenes. <i>Journal of the American Chemical Society</i> , 2019, 141, 1908-1912.	6.6	27
77	Dative and electron-sharing bonding in transition metal compounds. <i>Journal of Computational Chemistry</i> , 2019, 40, 247-264.	1.5	74
78	Synthesis and Reactivity Studies of Amido-Substituted Germanium(I)/Tin(I) Dimers and Clusters. <i>Chemistry - A European Journal</i> , 2019, 25, 2773-2785.	1.7	46
79	Alkali Metal Covalent Bonding in Nickel Carbonyl Complexes $ENi(CO)_3^+$. <i>Angewandte Chemie - International Edition</i> , 2019, 58, 1732-1738.	7.2	62
80	Barium as Honorary Transition Metal in Action: Experimental and Theoretical Study of $Ba(CO)^+$ and $Ba(CO)^+$. <i>Angewandte Chemie</i> , 2018, 130, 4038-4044.	1.6	16
81	Electronic Structure and Bonding Situation in M_2O_2 (M = Be, Mg, Ca) Rhombic Clusters. <i>Journal of Physical Chemistry A</i> , 2018, 122, 2816-2822.	1.1	34
82	Octacarbonyl Anion Complexes of Group Three Transition Metals $[TM(CO)_8]^-$ (TM=Sc, Y, La) and the 18-Electron Rule. <i>Angewandte Chemie</i> , 2018, 130, 6344-6349.	1.6	10
83	Dative and Electron-Sharing Bonding in C_2F_4 . <i>Chemistry - A European Journal</i> , 2018, 24, 9083-9089.	1.7	73
84	Octacarbonyl Anion Complexes of Group Three Transition Metals $[TM(CO)_8]^-$ (TM=Sc, Y, La) and the 18-Electron Rule. <i>Angewandte Chemie - International Edition</i> , 2018, 57, 6236-6241.	7.2	49
85	Dative versus electron-sharing bonding in N-oxides and phosphane oxides R_3EO and relative energies of the R_2EOR isomers (E = N, P; R = H, F, Cl, Me, Ph). A theoretical study. <i>Physical Chemistry Chemical Physics</i> , 2018, 20, 11856-11866.	1.3	32
86	Barium as Honorary Transition Metal in Action: Experimental and Theoretical Study of $Ba(CO)^+$ and $Ba(CO)^+$. <i>Angewandte Chemie - International Edition</i> , 2018, 57, 3974-3980.	7.2	60
87	Photoinduced Sulfur-Nitrogen Bond Rotation and Thermal Nitrogen Inversion in Heterocumulene OSNSO. <i>Journal of the American Chemical Society</i> , 2018, 140, 1231-1234.	6.6	9
88	Relativistic Effects on Donor-Acceptor Interactions in Coinage Metal Carbonyl Complexes $[TM(CO)_n]^+$ (TM=Cu, Ag, Au; $n=1, 2$). <i>Chemistry - A European Journal</i> , 2018, 24, 11675-11682.	1.7	16
89	A Route to Base Coordinate Silicon Difluoride and the Silicon Trifluoride Radical. <i>Chemistry - A European Journal</i> , 2018, 24, 1264-1268.	1.7	24
90	Energy decomposition analysis. <i>Wiley Interdisciplinary Reviews: Computational Molecular Science</i> , 2018, 8, e1345.	6.2	369

#	ARTICLE	IF	CITATIONS
91	Berichtigung: Barium as Honorary Transition Metal in Action: Experimental and Theoretical Study of Ba(CO) ₈ and Ba(CO) ₉ . <i>Angewandte Chemie</i> , 2018, 130, 15856-15857.	1.6	0
92	Double dative bond between divalent carbon(0) and uranium. <i>Nature Communications</i> , 2018, 9, 4997.	5.8	63
93	Vinyltrifluoroborate Complexes of Silver Supported by N-Heterocyclic Carbenes. <i>European Journal of Inorganic Chemistry</i> , 2018, 2018, 4142-4152.	1.0	11
94	Suppressed Phosphine Dissociation by Polarization Effects on the Donor-Acceptor Bonds in [Ni(PEt ₃) ₄ ·n(ECp*)] (E = Al, Ga). <i>Inorganic Chemistry</i> , 2018, 57, 12657-12664.	1.9	15
95	Buckyball Difluoride F ₂ @C ₆₀ A Single-Molecule Crystal. <i>Angewandte Chemie - International Edition</i> , 2018, 57, 13931-13934.	7.2	28
96	Buckyball Difluoride F ₂ @C ₆₀ A Single-Molecule Crystal. <i>Angewandte Chemie</i> , 2018, 130, 14127-14130.	1.6	3
97	Observation of alkaline earth complexes M(CO) ₈ (M = Ca, Sr, or Ba) that mimic transition metals. <i>Science</i> , 2018, 361, 912-916.	6.0	207
98	Comparison of Two Phosphinidenes Binding to Silicon(IV)dichloride as well as to Silylene. <i>Journal of the American Chemical Society</i> , 2018, 140, 9409-9412.	6.6	42
99	Anion stabilised hypercloso-hexaalane Al ₆ H ₆ . <i>Nature Communications</i> , 2018, 9, 3079.	5.8	39
100	Bonding in Binuclear Carbonyl Complexes M ₂ (CO) ₉ (M = Fe, Ru, Os). <i>Inorganic Chemistry</i> , 2018, 57, 7780-7791.	1.9	50
101	Heterocumulene Sulfinyl Radical OCNSO. <i>Angewandte Chemie - International Edition</i> , 2017, 56, 2140-2144.	7.2	17
102	Heterocumulene Sulfinyl Radical OCNSO. <i>Angewandte Chemie</i> , 2017, 129, 2172-2176.	1.6	5
103	Aromaticity, the Hückel 4n+2 Rule and Magnetic Current. <i>ChemistrySelect</i> , 2017, 2, 863-870.	0.7	66
104	The Bonding Situation in Metalated Ylides. <i>Chemistry - A European Journal</i> , 2017, 23, 4422-4434.	1.7	92
105	Ein stabiles neutrales Radikal in der Koordinationssphäre des Aluminiums. <i>Angewandte Chemie</i> , 2017, 129, 407-411.	1.6	23
106	Ligand-Supported E ₃ Clusters (E=Si-Sn). <i>Chemistry - A European Journal</i> , 2017, 23, 7463-7473.	1.7	11
107	(L) ₂ C ₂ P ₂ : Dicarbondiphosphide Stabilized by N-Heterocyclic Carbenes or Cyclic Diamido Carbenes. <i>Angewandte Chemie - International Edition</i> , 2017, 56, 5744-5749.	7.2	102
108	(L) ₂ C ₂ P ₂ : Dicarbondiphosphide Stabilized by N-Heterocyclic Carbenes or Cyclic Diamido Carbenes. <i>Angewandte Chemie</i> , 2017, 129, 5838-5843.	1.6	55

#	ARTICLE	IF	CITATIONS
109	Helium Accepts Back-Donation In Highly Polar Complexes: New Insights into the Weak Chemical Bond. <i>Journal of Physical Chemistry Letters</i> , 2017, 8, 3334-3340.	2.1	24
110	Normal-to-abnormal rearrangement of an N-heterocyclic carbene with a silylene transition metal complex. <i>Dalton Transactions</i> , 2017, 46, 7791-7799.	1.6	32
111	Carbene stabilized interconnected bis-germylene and its silicon analogue with small methyl substituents. <i>Dalton Transactions</i> , 2017, 46, 7947-7952.	1.6	23
112	Frontispiece: Ligand-Supported E ₃ Clusters (E=Si-Sn). <i>Chemistry - A European Journal</i> , 2017, 23, .	1.7	0
113	Dative bonding in main group compounds. <i>Coordination Chemistry Reviews</i> , 2017, 344, 163-204.	9.5	174
114	Dicarbonyls of Carbon and Methylidyne Cations. <i>Journal of Physical Chemistry A</i> , 2017, 121, 2903-2910.	1.1	5
115	An Electrophilic Carbene-Anchored Silylene-Phosphinidene. <i>Angewandte Chemie - International Edition</i> , 2017, 56, 4219-4223.	7.2	54
116	The aromaticity of dicupra[10]annulenes. <i>Physical Chemistry Chemical Physics</i> , 2017, 19, 9669-9675.	1.3	33
117	An Electrophilic Carbene-Anchored Silylene-Phosphinidene. <i>Angewandte Chemie</i> , 2017, 129, 4283-4287.	1.6	27
118	Carbenes as Ligands in Novel Main-Group Compounds E[C(NHC) ₂] ₂ (E=Be, Tl, Pb, Bi, Sn, Sb, Te, Se, S, O, N, C). <i>Chemistry - A European Journal</i> , 2017, 23, 17222-17226.	1.7	85
119	A Very Short Be-Be Distance but No Bond: Synthesis and Bonding Analysis of Ng-Be ₂ O ₂ -Ng (Ng, Ng=Ne, Ar, Kr, Xe). <i>Chemistry - A European Journal</i> , 2017, 23, 2035-2039.	7.2	46
120	NHC-Stabilised Acetylene-How Far Can the Analogy Be Pushed?. <i>Chemistry - A European Journal</i> , 2017, 23, 2926-2934.	1.7	65
121	A Stable Neutral Radical in the Coordination Sphere of Aluminum. <i>Angewandte Chemie - International Edition</i> , 2017, 56, 397-400.	7.2	56
122	Heterobimetallic Complexes Featuring Fe(CO) ₅ as a Ligand on Gold. <i>Chemistry - A European Journal</i> , 2017, 23, 17222-17226.	1.7	18
123	The trans Effect in Palladium Phosphine Sulfonate Complexes. <i>Journal of Physical Chemistry A</i> , 2017, 121, 7709-7716.	1.1	13
124	Organosilicon Radicals with Si-H and Si-Me Bonds from Commodity Precursors. <i>Journal of the American Chemical Society</i> , 2017, 139, 11028-11031.	6.6	25
125	Parent Thioketene S ₂ O ₂ CCSO: Gas-Phase Generation, Structure, and Bonding Analysis. <i>Chemistry - A European Journal</i> , 2017, 23, 16566-16573.	1.7	39
126	The hypothiocyanite radical OSCN and its isomers. <i>Physical Chemistry Chemical Physics</i> , 2017, 19, 16713-16720.	1.3	6

#	ARTICLE	IF	CITATIONS
127	InnenrÄ¼cktitelbild: Heterocumulene Sulfinyl Radical OCNSO (Angew. Chem. 8/2017). Angewandte Chemie, 2017, 129, 2253-2253.	1.6	0
128	Carbodicarbenes: Unexpected ĩ€-Accepting Ability during Reactivity with Small Molecules. Journal of the American Chemical Society, 2017, 139, 12830-12836.	6.6	57
129	A C(sp ²)ĤH Dehydrogenation of Heteroarenes and Arenes by a Functionalized Aluminum Hydride. Chemistry - A European Journal, 2017, 23, 13633-13637.	1.7	28
130	A C ₂ Fragment as FourĖElectron ĩf Donor. Zeitschrift Fur Anorganische Und Allgemeine Chemie, 2017, 643, 1096-1099.	0.6	20
131	Covalent Bonding and Charge Shift Bonds: Comment on ĖThe CarbonĖNitrogen Bonds in Ammonium Compounds Are Charge Shift BondsĖ. Chemistry - A European Journal, 2017, 23, 18320-18324.	1.7	5
132	An acyclic zincagermylene: rapid activation of dihydrogen at sub-ambient temperature. Chemical Communications, 2017, 53, 12692-12695.	2.2	35
133	Structural Exploration of Phantom Oligoguanidine from Asymmetric Diamine and Guanidine Hydrochloride. Macromolecular Chemistry and Physics, 2016, 217, 1834-1841.	1.1	0
134	Proton Affinities of Cationic Carbene Adducts [AC(PPh ₃) ₂] ⁺ (A=Halogen, Hydrogen, Methyl) and Unusual Electronic Structures of the Cations and Dications [AC(H)(PPh ₃) ₂] ²⁺ . Chemistry - A European Journal, 2016, 22, 8536-8546.	1.7	27
135	The [B ₃ (NN) ₃] ⁺ and [B ₃ (CO) ₃] ⁺ Complexes Featuring the Smallest ĖAromatic Species B ₃ ⁺ . Angewandte Chemie - International Edition, 2016, 55, 2078-2082.	7.2	64
136	Unusually Short BeĖBe Distances with and without a Bond in Be ₂ F ₂ and in the Molecular Discuses Be ₂ B ₈ and Be ₂ B ₇ ⁺ . Angewandte Chemie - International Edition, 2016, 55, 7841-7846.	7.2	60
137	The Chemical Bond in C ₂ . Chemistry - A European Journal, 2016, 22, 4100-4108.	1.7	75
138	The [B ₃ (NN) ₃] ⁺ and [B ₃ (CO) ₃] ⁺ Complexes Featuring the Smallest ĖAromatic Species B ₃ ⁺ . Angewandte Chemie, 2016, 128, 2118-2122.	1.6	24
139	Comment on ĖThe Quadruple Bonding in C ₂ Reproduces the Properties of the MoleculeĖ. Chemistry - A European Journal, 2016, 22, 18975-18976.	1.7	33
140	The Structure of the Carbene Stabilized Si ₂ H ₂ May Be Equally Well Described with Coordinate Bonds as with Classical Double Bonds. Journal of the American Chemical Society, 2016, 138, 10429-10432.	6.6	105
141	Bonding analysis of ylidone complexes EL ₂ (E = CĖPb) with phosphine and carbene ligands L. Canadian Journal of Chemistry, 2016, 94, 1006-1014.	0.6	19
142	New Avenues in sĖBlock Chemistry: Beryllium(0) Complexes. Angewandte Chemie - International Edition, 2016, 55, 13380-13382.	7.2	27
143	Neue Wege in der sĖBlockĖChemie Ė Komplexe mit Beryllium in der Oxidationsstufe Null. Angewandte Chemie, 2016, 128, 13576-13578.	1.6	8
144	A Triatomic Silicon(0) Cluster Stabilized by a Cyclic Alkyl(amino) Carbene. Angewandte Chemie, 2016, 128, 3210-3213.	1.6	25

#	ARTICLE	IF	CITATIONS
145	Unusually Short Be ⁺ Be Distances with and without a Bond in Be ₂ F ₂ and in the Molecular Discs Be ₂ B ₈ and Be ₂ B ₇ ⁺ . <i>Angewandte Chemie</i> , 2016, 128, 7972-7977.	1.6	34
146	Observation of Main-Group Tricarbonyls [B(CO) ₃] and [C(CO) ₃] ⁺ Featuring a Tilted One-Electron Donor Carbonyl Ligand. <i>Chemistry - A European Journal</i> , 2016, 22, 2376-2385.	1.7	23
147	Zn ⁻ Zn interactions at nickel and palladium centers. <i>Chemical Science</i> , 2016, 7, 6413-6421.	3.7	30
148	A Triatomic Silicon(0) Cluster Stabilized by a Cyclic Alkyl(amino) Carbene. <i>Angewandte Chemie - International Edition</i> , 2016, 55, 3158-3161.	7.2	54
149	Donor-acceptor bonding in novel low-coordinated compounds of boron and group-14 atoms C ⁺ Sn. <i>Chemical Society Reviews</i> , 2016, 45, 1129-1144.	18.7	162
150	Gilbert Lewis and the Model of Dative Bonding. <i>Structure and Bonding</i> , 2016, , 131-156.	1.0	12
151	Reaction Mechanism of the Symmetry-Forbidden [2+2] Addition of Ethylene and Acetylene to Amido-Substituted Digermynes and Distannynes Ph ₂ Ni ₂ NPh ₂ , (E=Ge, Sn): A Theoretical Study. <i>Chemistry - A European Journal</i> , 2015, 21, 12405-12413.	1.7	27
152	Stabilization of Heterodiatomic SiC Through Ligand Donation: Theoretical Investigation of SiC(L) ₂ (L=NHC ⁺ Me, CAAC ⁺ Me, PMe ₃). <i>Angewandte Chemie - International Edition</i> , 2015, 54, 12319-12324.	7.2	102
153	Stabilisierung von heterodiatomarem SiC durch Donorliganden - theoretische Untersuchung von SiC(L) ₂ (L=NHC ⁺ Me, CAAC ⁺ Me, PMe ₃). <i>Angewandte Chemie</i> , 2015, 127, 12494-12500.	1.6	44
154	Formation and Characterization of the Boron Dicarboxyl Complex [B(CO) ₂] ⁺ . <i>Angewandte Chemie - International Edition</i> , 2015, 54, 11078-11083.	7.2	107
155	Carbon Monoxide Bonding With BeO and BeCO ₃ : Surprisingly High CO Stretching Frequency of OBeCO ₃ . <i>Angewandte Chemie - International Edition</i> , 2015, 54, 124-128.	7.2	70
156	No Need for a Re-examination of the Electrostatic Notation of the Hydrogen Bonding: A Comment. <i>Angewandte Chemie - International Edition</i> , 2015, 54, 2596-2599.	7.2	78
157	The Fate of NHC-Stabilized Dicarboxyl. <i>Chemistry - A European Journal</i> , 2015, 21, 3377-3386.	1.7	42
158	The f-Aromatic Clusters [Zn ₃] ⁺ and [Zn ₂ Cu]: Embryonic Brass. <i>Angewandte Chemie - International Edition</i> , 2015, 54, 4370-4374.	7.2	72
159	Reactivity of Amido-Digermynes, LGeGeL (L = Bulky Amide), toward Olefins and Related Molecules: Facile Reduction, C-H Activation, and Reversible Cycloaddition of Unsaturated Substrates. <i>Organometallics</i> , 2015, 34, 3175-3185.	1.1	60
160	The boron-boron triple bond in NHC ⁺ B ⁻ B ⁺ NHC. <i>Chemical Science</i> , 2015, 6, 4089-4094.	3.7	73
161	Bonding situation in silicon complexes [(L) ₂ (Si)] and [(L) ₂ (Si)] with NHC and cAAC ligands. <i>Journal of Organometallic Chemistry</i> , 2015, 792, 139-148.	0.8	33
162	Two-coordinate group 14 element hydrides as reagents for the facile, and sometimes reversible, hydrogermylation/hydrostannylation of unactivated alkenes and alkynes. <i>Chemical Science</i> , 2015, 6, 7249-7257.	3.7	69

#	ARTICLE	IF	CITATIONS
163	Cyclic trinuclear copper(<i>scpi</i>), silver(<i>scpi</i>), and gold(<i>scpi</i>) complexes: a theoretical insight. Dalton Transactions, 2015, 44, 377-385.	1.6	36
164	Experimental and Theoretical Studies of the Infrared Spectra and Bonding Properties of NgBeCO ₃ and a Comparison with NgBeO (Ng = He, Ne, Ar, Kr, Xe). Journal of Physical Chemistry A, 2015, 119, 2543-2552.	1.1	62
165	Analysis of the E–E Bond in Group-13 Complexes [(PMe ₃) ₂ (E ₂ Hn)] (E = B, In, n = 4, 2, 0). Croatica Chemica Acta, 2014, 87, 413-422.	0.1	8
166	Bonding Situation in Dimeric Group 15 Complexes [(NHC) ₂ (E ₂)] (E = N, Bi). Zeitschrift Fur Naturforschung - Section A Journal of Physical Sciences, 2014, 69, 385-395.	0.7	14
167	Experimental Charge Density Study of a Silylone. Angewandte Chemie - International Edition, 2014, 53, 2766-2770.	7.2	115
168	Comparative bonding analysis of N ₂ and P ₂ versus tetrahedral N ₄ and P ₄ . Theoretical Chemistry Accounts, 2014, 133, 1.	0.5	27
169	Stabilization of a Cobalt–Cobalt Bond by Two Cyclic Alkyl Amino Carbenes. Journal of the American Chemical Society, 2014, 136, 1770-1773.	6.6	55
170	Dative Bonds in Main-Group Compounds: A Case for More Arrows!. Angewandte Chemie - International Edition, 2014, 53, 6040-6046.	7.2	235
171	Reaction Mechanisms of Small-Molecule Activation by Amidoditetrylides R ₂ N–E–NR ₂ (E = Si, Ge, Sn). Inorganic Chemistry, 2014, 53, 6482-6490.	1.9	23
172	Low Coordinate Germanium(II) and Tin(II) Hydride Complexes: Efficient Catalysts for the Hydroboration of Carbonyl Compounds. Journal of the American Chemical Society, 2014, 136, 3028-3031.	6.6	290
173	New bonding modes of carbon and heavier group 14 atoms Si–Pb. Chemical Society Reviews, 2014, 43, 5106-5139.	18.7	227
174	Coinage Metals Binding as Main Group Elements: Structure and Bonding of the Carbene Complexes [TM(cAAC) ₂] and [TM(cAAC) ₂] ⁺ (TM = Cu, Ag, Au). Journal of the American Chemical Society, 2014, 136, 17123-17135.	6.6	84
175	One-Electron-Mediated Rearrangements of 2,3-Disiladibene. Journal of the American Chemical Society, 2014, 136, 8919-8922.	6.6	73
176	Beryllium chemistry the safe way: a theoretical evaluation of low oxidation state beryllium compounds. Dalton Transactions, 2013, 42, 11375.	1.6	102
177	Isolation of Neutral Mono- and Dinuclear Gold Complexes of Cyclic (Alkyl)(amino)carbenes. Angewandte Chemie - International Edition, 2013, 52, 8964-8967.	7.2	119
178	Exploiting the Twofold Donor Ability of Carbodiphosphoranes: Theoretical Studies of [(PPh ₃) ₂] ₂ Cat ⁺ EH ₂ (E ⁺ =Be, B ⁺) and [(Ph ₃ P) ₂ Ci ^{3/4} CH ₂] ₂ ⁺ . ChemPlusChem, 2013, 78, 1024-1032.	1.3	58
179	Formation of a 1,4-Diamino-2,3-disila-1,3-butadiene Derivative. Journal of the American Chemical Society, 2013, 135, 15990-15993.	6.6	49
180	Reductive elimination: a pathway to low-valent aluminium species. Chemical Communications, 2013, 49, 2858.	2.2	94

#	ARTICLE	IF	CITATIONS
181	End-On and Side-On π -Acid Ligand Adducts of Gold(I): Carbonyl, Cyanide, Isocyanide, and Cyclooctyne Gold(I) Complexes Supported by N-Heterocyclic Carbenes and Phosphines. <i>Inorganic Chemistry</i> , 2013, 52, 729-742.	1.9	69
182	Reaction Pathways for Addition of H_2 to Amido-Ditetrylynes R_2N-NR_2 (E = Si, Ge, Sn). A Theoretical Study. <i>Organometallics</i> , 2013, 32, 6666-6673.	1.1	39
183	Spectroscopic investigation of the far-infrared properties of liquid crystals. , 2013, , .		0
184	A Stable Singlet Biradicaloid Siladibene: $(L)_2Si$. <i>Angewandte Chemie - International Edition</i> , 2013, 52, 2963-2967.	7.2	246
185	Conversion of a Singlet Silylene to a stable Biradical. <i>Angewandte Chemie - International Edition</i> , 2013, 52, 1801-1805.	7.2	167
186	Dinitrogen as Double Lewis Acid: Structure and Bonding of Triphenylphosphinazine $N_2(PPH_3)_2$. <i>Angewandte Chemie - International Edition</i> , 2013, 52, 3004-3008.	7.2	76
187	Bonding Analysis of the Trimethylenemethane (TMM) Complexes $[(\eta^6-C_6H_6)M-TMM]$ (M = Fe, Ru, Os), $[(\eta^5-C_5H_5)M-TMM]$ (M = Co, Rh, Ir), and $[(\eta^4-C_4H_4)M-TMM]$ (M = Ni, Pd, Pt). <i>Organometallics</i> , 2013, 32, 1742-1751.	1.1	38
188	Bonding analysis of trimethylenemethane (TMM) complexes $[(CO)_3M-TMM]$ (M = Fe, Ru, Os, Rh+). Absence of expected bond paths. <i>Journal of Organometallic Chemistry</i> , 2013, 748, 2-7.	0.8	18
189	Critical Comments on "One Molecule, Two Atoms, Three Views, Four Bonds". <i>Angewandte Chemie - International Edition</i> , 2013, 52, 5922-5925.	7.2	53
190	Tris(alkyne) and Bis(alkyne) Complexes of Coinage Metals: Synthesis and Characterization of $(cyclooctyne)_3M$ (M = Cu, Ag) and $(cyclooctyne)_2Au$ and Coinage Metal (M = Cu, Ag, Au) Family Group Trends. <i>Organometallics</i> , 2013, 32, 3135-3144.	1.1	59
191	Complexation behavior of two-coordinated carbon compounds containing fluorenyl ligands. <i>Dalton Transactions</i> , 2013, 42, 13349.	1.6	15
192	Structure and bonding of tetrylone complexes $[(CO)_4W\{E(PPH_3)_2\}]$ (E = C, Pb). <i>Molecular Physics</i> , 2013, 111, 2640-2646.	0.8	23
193	Activation of H_2 by a Multiply Bonded Amido-Digermene: Evidence for the Formation of a Hydrido-Germene. <i>Angewandte Chemie</i> , 2013, 125, 10389-10393.	1.6	53
194	Activation of H_2 by a Multiply Bonded Amido-Digermene: Evidence for the Formation of a Hydrido-Germene. <i>Angewandte Chemie - International Edition</i> , 2013, 52, 10199-10203.	7.2	154
195	An N-heterocyclic carbene adduct of diatomic tin, $:Sn-Sn$. <i>Chemical Communications</i> , 2012, 48, 9855.	2.2	162
196	Transition-Metal Complexes of Tetrylones $[(CO)_5W\{E(PPH_3)_2\}]$ and Tetrylenes $[(CO)_5W\{NHE\}]$ (E=C, Pb): A Theoretical Study. <i>Chemistry - A European Journal</i> , 2012, 18, 12733-12748.	1.7	69
197	A Crystalline Singlet Phosphinonitrene: A Nitrogen Atom-Transfer Agent. <i>Science</i> , 2012, 337, 1526-1528.	6.0	148
198	Energy decomposition analysis. <i>Wiley Interdisciplinary Reviews: Computational Molecular Science</i> , 2012, 2, 43-62.	6.2	613

#	ARTICLE	IF	CITATIONS
199	A Boron-Boron Triple Bond. <i>Science</i> , 2012, 336, 1394-1395.	6.0	83
200	Isolable Tris(alkyne) and Bis(alkyne) Complexes of Gold(I). <i>Angewandte Chemie - International Edition</i> , 2012, 51, 3940-3943.	7.2	49
201	The Facile Reduction of Carbon Dioxide to Carbon Monoxide with an Amido-Digermine. <i>Angewandte Chemie - International Edition</i> , 2012, 51, 8611-8614.	7.2	101
202	Are They Linear, Bent, or Cyclic? Quantum Chemical Investigation of the Heavier Group 14 and Group 15 Homologues of HCN and HNC. <i>Chemistry - an Asian Journal</i> , 2012, 7, 1296-1311.	1.7	17
203	Borylene Complexes (BH) ₂ and Nitrogen Cation Complexes (N ⁺) ₂ : Isoelectronic Homologues of Carbones CL ₂ . <i>Chemistry - A European Journal</i> , 2012, 18, 5676-5692.	1.7	131
204	Building a Bridge between Coordination Compounds and Clusters: Bonding Analysis of the Icosahedral Molecules [M(ER) ₁₂] (M = Cr, Mo, W; E = Zn, Cd, Hg). <i>Journal of Physical Chemistry A</i> , 2011, 115, 12758-12768.	1.1	27
205	[Hg{C(PPh) ₃ }] ₂ and [Cu{C(PPh) ₃ }] ₂ and Comparative Theoretical Study of Carbene Complexes [M(NHC) ₂] with Carbene Complexes [M{C(PH) ₃ }] ₂ (M = Cu ⁺ , Ag ⁺) <i>Tj ETQq1 1 0.784314 rgBT /Overlock</i>	1.1	37
206	N-Heterocyclic carbenes versus transition metals for stabilizing phosphinyl radicals. <i>Chemical Science</i> , 2011, 2, 858.	3.7	108
207	Synthesis and Characterization of a Neutral Tricoordinate Organoboron Isoelectronic with Amines. <i>Science</i> , 2011, 333, 610-613.	6.0	486
208	A Digermine with a Ge-Ge Single Bond That Activates Dihydrogen in the Solid State. <i>Journal of the American Chemical Society</i> , 2011, 133, 18622-18625.	6.6	202
209	Preparation, Characterization, and Theoretical Analysis of Group 14 Element(I) Dimers: A Case Study of Magnesium(I) Compounds as Reducing Agents in Inorganic Synthesis. <i>Inorganic Chemistry</i> , 2011, 50, 12315-12325.	1.9	139
210	Cationic Gold Carbonyl Complex on a Phosphine Support. <i>Inorganic Chemistry</i> , 2011, 50, 4253-4255.	1.9	44
211	Carbodicarbenes—divalent carbon(0) compounds exhibiting carbon—carbon donor—acceptor bonds. <i>Wiley Interdisciplinary Reviews: Computational Molecular Science</i> , 2011, 1, 869-878.	6.2	65
212	Divalent Pb(0) compounds. <i>Theoretical Chemistry Accounts</i> , 2011, 129, 615-623.	0.5	44
213	The Reaction of BeCl ₂ with Carbodiphosphorane C(PPh) ₃ ₂ ; Experimental and Theoretical Studies. <i>Zeitschrift Fur Anorganische Und Allgemeine Chemie</i> , 2011, 637, 1702-1710.	0.6	46
214	1,5-Electrocyclization versus 1,5-proton shift in imidazolium allylides and 2-phosphaallylides: a DFT investigation. <i>Journal of Physical Organic Chemistry</i> , 2011, 24, 786-797.	0.9	7
215	Distinguishing Carbones from Allenes by Complexation to AuCl. <i>Chemistry - A European Journal</i> , 2011, 17, 9944-9956.	1.7	84
216	Structures and Stabilities of Group 13 Adducts [(NHC)(EX) ₃] and [(NHC) ₂ (E) ₂ X] (E=B to In; X=H, Cl; <i>n</i> =4, 2, 0); <i>Tj ETQq0 0 0 rgBT /Overlock 10 Tf</i>	1.7	133
	<i>Chemistry - A European Journal</i> , 2011, 17, 13517-13525.		

#	ARTICLE	IF	CITATIONS
217	Carbodiphosphoranes and Related Ligands. Topics in Organometallic Chemistry, 2010, , 49-92.	0.7	89
218	Carbodicarbenes and Related Divalent Carbon(0) Compounds. Chemistry - A European Journal, 2010, 16, 10160-10170.	1.7	135
219	Molecular Alloys: Experimental and Theoretical Investigations on the Substitution of Zinc by Cadmium and Mercury in the Homologous Series [Mo(Mâ€²R)â€²] and [M(Mâ€²R)â€²] (M=Pd, Pt; Mâ€²=Zn, Cd, Hg). Chemistry - A European Journal, 2010, 16, 13372-13384.	1.7	27
220	Carbodiylides C(ECp*)â€² (E=Bâ€²Ti): Another Class of Theoretically Predicted Divalent Carbon(0) Compounds. Angewandte Chemie - International Edition, 2010, 49, 7106-7110.	7.2	23
221	Isolation of crystalline carbene-stabilized P2-radical cations and P2-dications. Nature Chemistry, 2010, 2, 369-373.	6.6	282
222	Synthesis of a stable adduct of dialane(4) (Al2H4) via hydrogenation of a magnesium(I) dimer. Nature Chemistry, 2010, 2, 865-869.	6.6	221
223	A Crystalline Phosphinyl Radical Cation. Journal of the American Chemical Society, 2010, 132, 10262-10263.	6.6	185
224	Divalent carbon(0) compounds. Pure and Applied Chemistry, 2009, 81, 597-614.	0.9	211
225	Divalent Silicon(0) Compounds. Chemistry - A European Journal, 2009, 15, 3448-3456.	1.7	140
226	Divalent E(0) Compounds (E=Siâ€²Sn). Chemistry - A European Journal, 2009, 15, 8593-8604.	1.7	141
227	Carbodiphosphorane C(PPhâ€²)â€² as a Single and Twofold Lewis Base with Boranes: Synthesis, Crystal Structures and Theoretical Studies on [Hâ€²]B(C(PPhâ€²)â€²)â€² and [(1/4â€²)Hâ€²]B(C(PPhâ€²)â€²)â€². European Journal of Inorganic Chemistry, 2009, 2009, 4507-4517.	1.0	88
228	Synthesis and Ligand Properties of a Persistent, Allâ€²Carbon Fourâ€²Memberedâ€²Ring Allene. Angewandte Chemie - International Edition, 2009, 48, 4792-4795.	7.2	122
229	Nâ€²Heterocyclic Carbene Stabilized Digermanium(0). Angewandte Chemie - International Edition, 2009, 48, 9701-9704.	7.2	304
230	Molecular Alloys, Linking Organometallics with Intermetallic Humeâ€²Rothery Phases: The Highly Coordinated Transition Metal Compounds [M(ZnR)â€²] (â€² = Ni, Cu, Ag) Containing Organoâ€²Zinc Ligands. Journal of the American Chemical Society, 2009, 131, 16063-16077.	6.6	65
231	Heavy Halogen Atom Effect on â€²C NMR Chemical Shifts in Monohalo Derivatives of Cyclohexane and Pyran. Experimental and Theoretical Study. Journal of Chemical Theory and Computation, 2009, 5, 2222-2228.	2.3	32
232	Molecules with All Triple Bonds: OCBBCO, Nâ€²BBNâ€², and [OBBBBO]â€². Journal of Physical Chemistry A, 2009, 113, 11693-11698.	1.1	69
233	Tolmanâ€²s Electronic Parameters for Divalent Carbon(0) Compounds. Organometallics, 2009, 28, 3901-3905.	1.1	109
234	The Dewarâ€²Chattâ€²Duncanson model reversed â€² Bonding analysis of group-10 complexes [(PMâ€²)â€²]â€² (M = Ni, Pd, Pt; E = B, Al, Ga, In, Tl; X = H, F, Cl, Br, I) Tj ETQq 0 0 0 0 BT /Overl	0.0	0

#	ARTICLE	IF	CITATIONS
235	Donor Acceptor Complexes of Noble Gases. <i>Journal of the American Chemical Society</i> , 2009, 131, 3942-3949.	6.6	78
236	Isolation of a C5-Deprotonated Imidazolium, a Crystalline σ -Abnormal ϵ -N-Heterocyclic Carbene. <i>Science</i> , 2009, 326, 556-559.	6.0	404
237	Exocyclic Delocalization at the Expense of Aromaticity in 3,5-bis(π -Donor) Substituted Pyrazolium Ions and Corresponding Cyclic Bent Allenes. <i>Journal of the American Chemical Society</i> , 2009, 131, 11875-11881.	6.6	119
238	Divalent Carbon(0) Chemistry, Part 1: Parent Compounds. <i>Chemistry - A European Journal</i> , 2008, 14, 3260-3272.	1.7	384
239	Divalent Carbon(0) Chemistry, Part 2: Protonation and Complexes with Main Group and Transition Metal Lewis Acids. <i>Chemistry - A European Journal</i> , 2008, 14, 3273-3289.	1.7	285
240	First and Second Proton Affinities of Carbon Bases. <i>ChemPhysChem</i> , 2008, 9, 1474-1481.	1.0	181
241	Transition Metal π -Carbon Complexes. A Theoretical Study. <i>Journal of the American Chemical Society</i> , 2007, 129, 7596-7610.	6.6	119
242	Pseudopotential Calculations of Transition Metal Compounds: Scope and Limitations. <i>Reviews in Computational Chemistry</i> , 2007, , 63-144.	1.5	100
243	Nonclassical Metal Carbonyls. <i>Progress in Inorganic Chemistry</i> , 2007, , 1-112.	3.0	94
244	Is This a Chemical Bond? A Theoretical Study of Ng_2C_{60} ($\text{Ng}=\text{He, Ne, Ar, Kr}$). <i>Tj ETQq0 0 0 rgBT / Overlock 10</i>	1.7	179
245	Reply to R��plique: A New Concept for Bonding in Carbodiphosphanes?. <i>Angewandte Chemie - International Edition</i> , 2007, 46, 2986-2987.	7.2	40
246	$\text{C}(\text{NHC})_2$: Divalent Carbon(0) Compounds with $\text{N}\pi$ -Heterocyclic Carbene Ligands – Theoretical Evidence for a Class of Molecules with Promising Chemical Properties. <i>Angewandte Chemie - International Edition</i> , 2007, 46, 8695-8698.	7.2	361
247	Electronic structure of $\text{CO}\pi$ – An exercise in modern chemical bonding theory. <i>Journal of Computational Chemistry</i> , 2007, 28, 117-126.	1.5	114
248	Unicorns in the world of chemical bonding models. <i>Journal of Computational Chemistry</i> , 2007, 28, 15-24.	1.5	198
249	The Dewar-chatt-Duncanson bonding model of transition metal-olefin complexes examined by modern quantum chemical methods. , 2007, , 111-122.		25
250	Orbital Overlap and Chemical Bonding. <i>Chemistry - A European Journal</i> , 2006, 12, 9196-9216.	1.7	292
251	σ -Naked Ca^+ and In^+ as Pure Acceptor Ligands: Structure and Bonding of $[\text{GaPt}(\text{GaCp}^*)_4][\text{BArF}]$. <i>Angewandte Chemie - International Edition</i> , 2006, 45, 5207-5210.	7.2	61
252	Carbodiphosphanes: The Chemistry of Divalent Carbon(0). <i>Angewandte Chemie - International Edition</i> , 2006, 45, 8038-8042.	7.2	378

#	ARTICLE	IF	CITATIONS
253	The Nature of the Chemical Bond Revisited: An Energy-Partitioning Analysis of Nonpolar Bonds. Chemistry - A European Journal, 2005, 11, 1813-1825.	1.7	137
254	The nature of the chemical bond in the light of an energy decomposition analysis. , 2005, , 291-372.		62
255	Why Do the Heavy-Atom Analogues of Acetylene E2H2 (E = Si~Pb) Exhibit Unusual Structures?. Journal of the American Chemical Society, 2005, 127, 6290-6299.	6.6	178
256	Chemical Bonding in Octahedral XeF6 and SF6. Australian Journal of Chemistry, 2004, 57, 1191.	0.5	18
257	The nature of the chemical bond revisited. An energy partitioning analysis of diatomic molecules E2 (E=N~Bi, F~I), CO and BF. Theoretical Chemistry Accounts, 2004, 111, 381-389.	0.5	134
258	Energy Partitioning Analysis of the Bonding in Ethylene and Acetylene Complexes of Group 6, 8, and 11 Metals: (CO)5TM~C2Hx and Cl4TM~C2Hx (TM = Cr, Mo, W), (CO)4TM~C2Hx (TM = Fe, Ru, Os), and TM+~C2Hx (TM = Cu, Ag, Au). Journal of Physical Chemistry A, 2004, 108, 3134-3142.	1.1	146
259	Towards a rigorously defined quantum chemical analysis of the chemical bond in donor~acceptor complexes. Coordination Chemistry Reviews, 2003, 238-239, 55-82.	9.5	394
260	Energy Partitioning Analysis of the Bonding in L2TM~C2H2 and L2TM~C2H4 (TM = Ni, Pd, Pt; L2 = (PH3)2, Tj ETQq 0 0 0 rgBT / Overlock	1.1	89
261	Energy decomposition analysis of the chemical bond in main group and transition metal compounds. Faraday Discussions, 2003, 124, 365-378.	1.6	138
262	Bis(benzene)chromium Is a ~Bonded Molecule and Ferrocene Is a ~Bonded Molecule. Organometallics, 2003, 22, 3304-3308.	1.1	116
263	Nature of the Metal~Ligand Bond in M(CO)5PX3 Complexes (M = Cr, Mo, W; X = H, Me, F, Cl): A Synthesis, Molecular Structure, and Quantum-Chemical Calculations. Organometallics, 2002, 21, 2921-2930.	1.1	162
264	Structures, Bond Energies, Heats of Formation, and Quantitative Bonding Analysis of Main-Group Metallocenes [E(Cp)2] (E=Be~Ba, Zn, Si~Pb) and [E(Cp)] (E=Li~Cs, Ba~Tl). Chemistry - A European Journal, 2002, 8, 4693-4707.	1.7	128
265	Energy Analysis of Metal-Ligand Bonding in Transition Metal Complexes with Terminal Group-13 Diyl Ligands (CO)4Fe-ER, Fe(EMe)5 and Ni(EMe)4 (E = B~Tl; R = Cp, N(SiH3)2, Ph, Me) Reveals Significant ~ Bonding in Homoleptical Molecules. Journal of the American Chemical Society, 2001, 123, 1683-1693.	6.6	200
266	Understanding the nature of the bonding in transition metal complexes: from Dewar's molecular orbital model to an energy partitioning analysis of the metal~ligand bond. Journal of Organometallic Chemistry, 2001, 635, 9-23.	0.8	155
267	Turning a Transition State into a Minimum~The Nature of the Bonding in Diplumbylene Compounds RPbPbR (R=H, Ar). Angewandte Chemie - International Edition, 2001, 40, 2051-2055.	7.2	27
268	Chemical bonding in mononuclear transition metal complexes with Group 13 diyl ligands ER (E=B~Tl) Part X: Theoretical studies of inorganic compounds. Coordination Chemistry Reviews, 2000, 197, 249-276.	9.5	121
269	The Nature of the Transition Metal~Carbonyl Bond and the Question about the Valence Orbitals of Transition Metals. A Bond-Energy Decomposition Analysis of TM(CO)6q (TMq = Hf2-, Ta-, W, Re+, Os2+). Tj ETQq 1 1 0 0 78431 2 rgBT / Overlock	10.7	148
270	Nature of the Chemical Bond between a Transition Metal and a Group-13 Element: A Structure and Bonding of Transition Metal Complexes with Terminal Group-13 Diyl Ligands ER (E = B to Tl; R = Cp, Tj ETQq 0 0 0 0 rgBT / Overlock 10 Tf 5		

#	ARTICLE	IF	CITATIONS
271	The Nature of the Bonding in Transition-Metal Compounds. <i>Chemical Reviews</i> , 2000, 100, 717-774.	23.0	1,101
272	Trends in Molecular Geometries and Bond Strengths of the Homoleptic d10 Metal Carbonyl Cations [M(CO) _n] ^{x+} (M _{x+} =Cu ⁺ , Ag ⁺ , Au ⁺ , Zn ²⁺ , Cd ²⁺ , Hg ²⁺ ; n=1-6): A Theoretical Study. <i>Chemistry - A European Journal</i> , 1999, 5, 2573-2583.	1.7	123
273	Synthesis and Structure of [Ni{Ga ⁺ C(SiMe ₃) ₃ } ₄] and Quantum-Chemical Verification of Strong π Back-Bonding in the Model Compounds [Ni(EMe) ₄] (E = B, Al, Ga, In, Tl). <i>Organometallics</i> , 1999, 18, 3778-3780.	1.1	99
274	Reaction of Carbodiphosphorane Ph ₃ PCPPH ₃ with Ni(CO) ₄ . Experimental and Theoretical Study of the Structures and Properties of (CO) ₃ NiC(PPh ₃) ₂ and (CO) ₂ NiC(PPh ₃) ₂ . <i>Organometallics</i> , 1999, 18, 619-626.	1.1	93
275	Structure and Bonding of Low-Valent (Fischer-Type) and High-Valent (Schrock-Type) Transition Metal Carbene Complexes. <i>Chemistry - A European Journal</i> , 1998, 4, 1428-1438.	1.7	142
276	Structure and Bonding of Low-Valent (Fischer-Type) and High-Valent (Schrock-Type) Transition Metal Carbyne Complexes. <i>Chemistry - A European Journal</i> , 1998, 4, 1439-1448.	1.7	100
277	Nonclassical Metal Carbonyls: Appropriate Definitions with a Theoretical Justification. <i>Angewandte Chemie - International Edition</i> , 1998, 37, 2113-2116.	7.2	156
278	Ab initio studies of transition-metal compounds: the nature of the chemical bond to a transition metal. <i>Journal of the Chemical Society Dalton Transactions</i> , 1997, , 1653-1662.	1.1	154
279	Structure and Bonding of the Isolelectronic Hexacarbonyls [Hf(CO) ₆] ²⁻ , [Ta(CO) ₆] ⁻ , W(CO) ₆ , [Re(CO) ₆] ⁺ , [Os(CO) ₆] ²⁺ , and [Ir(CO) ₆] ³⁺ : A Theoretical Study. <i>Organometallics</i> , 1997, 16, 4807-4815.	1.1	128
280	Theoretical Analysis of the Bonding between CO and Positively Charged Atoms. <i>Journal of Physical Chemistry A</i> , 1997, 101, 9551-9559.	1.1	226
281	[(η -5-Cp*)Al ₂ Fe(CO) ₄] - Synthese, Struktur, Bindungsverhältnisse. <i>Angewandte Chemie</i> , 1997, 109, 95-97.	1.6	84
282	Structure and Bonding of the Transition-Metal Carbonyl Complexes M(CO) ₅ L (M = Cr, Mo, W) and M(CO) ₃ L (M = Ni, Pd, Pt; L = CO, SiO, CS, N ₂ , NO ⁺ , CN ⁻ , NC ⁻ , HCCH, CCH ₂ , CH ₂ , CF ₂ , H ₂). <i>Organometallics</i> , 1996, 15, 105-117.	1.1	193
283	The bonding of acetylene and ethylene in high-valent and low-valent transition metal compounds. <i>Journal of Organometallic Chemistry</i> , 1996, 525, 269-278.	0.8	64
284	Theoretical Studies of Organometallic Compounds. XIV. Structure and Bonding of the Transition Metal Methyl and Phenyl Compounds MCH ₃ and MC ₆ H ₅ (M = Cu, Ag, Au) and M(CH ₃) ₂ and M(C ₆ H ₅) ₂ (M = Ti, Zr, Hf, Ta, Nb, Mo, Cr, Ru, Rh, Ir, Pt, Os, Pd, Ni, Fe, Co, Ni, Cu, Ag, Au). <i>Journal of Organometallic Chemistry</i> , 1995, 487, 101-118.	0.0	0
285	Theoretical Studies of Organometallic Compounds. XIX. Complexes of Transition Metals in High and Low Oxidation States with Side-On-Bonded π -Ligands. <i>Organometallics</i> , 1995, 14, 5325-5336.	1.1	88
286	Investigation of Donor-Acceptor Interactions: A Charge Decomposition Analysis Using Fragment Molecular Orbitals. <i>The Journal of Physical Chemistry</i> , 1995, 99, 9352-9362.	2.9	710
287	Comparative Theoretical Study of Lewis Acid-Base Complexes of BH ₃ , BF ₃ , BCl ₃ , AlCl ₃ , and SO ₂ . <i>Journal of the American Chemical Society</i> , 1994, 116, 8741-8753.	6.6	435
288	Theoretical studies of organometallic compounds. 6. Structures and bond energies of M(CO) _n ⁺ , MCN, and M(CN) ₂ ⁻ (M = silver, gold; n = 1-3). <i>Organometallics</i> , 1993, 12, 4613-4622.	1.1	85

#	ARTICLE	IF	CITATIONS
289	Light noble gas chemistry: structures, stabilities, and bonding of helium, neon, and argon compounds. Journal of the American Chemical Society, 1990, 112, 4240-4256.	6.6	103
290	Helium bonding in singly and doubly charged first-row diatomic cations HeXn+ (X = Li-Ne; n = 1,2). The Journal of Physical Chemistry, 1989, 93, 3397-3410.	2.9	99
291	Stabilities and nature of the attractive interactions in HeBeO, NeBeO, and ArBeO and a comparison with analogs NGLiF, NGBN, and NGLiH (NG = He, Ar). A theoretical investigation. Journal of the American Chemical Society, 1988, 110, 8007-8016.	6.6	158
292	Helium chemistry: theoretical predictions and experimental challenge. Journal of the American Chemical Society, 1987, 109, 5917-5934.	6.6	207
293	The Chemical Bond – an Entrance Door of Chemistry to the Neighboring Sciences and to Philosophy. Israel Journal of Chemistry, 0, , .	1.0	6
294	How to capture C2O2: Structures and bonding of neutral and charged complexes [(NHC)-C2O2-(NHC)]q (NHC = N-heterocyclic carbene; q = 0, 1+, 2+). Physical Chemistry Chemical Physics, 0, , .	1.3	0