

Stanislav Zalis

List of Publications by Year in descending order

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186
papers

7,201
citations

41344

49
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82547

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times ranked

5653
citing authors

#	ARTICLE	IF	CITATIONS
1	Excitation-Wavelength-Dependent Photophysics of d ⁸ -Di-isocyanide Complexes. <i>Inorganic Chemistry</i> , 2022, 61, 2745-2759.	4.0	5
2	Structural and Oxidation State Alternatives in Platinum and Palladium Complexes of a Redox-Active Amidinato Ligand. <i>Chemistry - A European Journal</i> , 2021, 27, 3374-3381.	3.3	1
3	Variable electronic structure and spin distribution in bis(2,2'-bipyridine)-metal complexes (M = Ru or Tl). <i>Chemical Communications</i> , 2021, 2021, 10784-10785.	3.3	5
4	Photoinduced hole hopping through tryptophans in proteins. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2021, 118, .	7.1	13
5	Optical and Infrared Spectroelectrochemical Studies of CN-Substituted Bipyridyl Complexes of Ruthenium(II). <i>Inorganic Chemistry</i> , 2021, 60, 3514-3523.	4.0	4
6	Analysis of Multiple Redox Sites in Complexes [M(C ₅ Me ₅)(Q)(NO)] _n , M=Ru or Os, Q= o-Quinones. <i>Zeitschrift Fur Anorganische Und Allgemeine Chemie</i> , 2021, 647, 867-875.	1.2	0
7	NO and NO ₂ as non-innocent ligands: A comparison. <i>Coordination Chemistry Reviews</i> , 2020, 404, 2131-14.	18.8	23
8	Oxidation State Assignments in the Organoplatinum One-Electron Redox Series [(N ⁺)PtMes ₂] _n , n = +, 0, -2. <i>European Journal of Inorganic Chemistry</i> , 2020, 2020, 2435-2443.	2.0	6
9	Analysis of a Diimine-Organonickel Redox Series. <i>European Journal of Inorganic Chemistry</i> , 2020, 2020, 3010-3015.	2.0	2
10	Time-Resolved Femtosecond Stimulated Raman Spectra and DFT Anharmonic Vibrational Analysis of an Electronically Excited Rhenium Photosensitizer. <i>Journal of Physical Chemistry A</i> , 2020, 124, 1253-1265.	2.5	13
11	Hole Hopping Across a Protein-Protein Interface. <i>Journal of Physical Chemistry B</i> , 2019, 123, 1578-1591.	2.6	8
12	Two Tryptophans Are Better Than One in Accelerating Electron Flow through a Protein. <i>ACS Central Science</i> , 2019, 5, 192-200.	11.3	28
13	Lipophilicity of acetylcholine and related ions examined by ion transfer voltammetry at a polarized room-temperature ionic liquid membrane. <i>Journal of Electroanalytical Chemistry</i> , 2018, 815, 183-188.	3.8	9
14	At the Borderline between Metal-Metal Mixed Valency and a Radical Bridge Situation: Four Charge States of a Diruthenium Complex with a Redox-Active Bis(mer-tridentate) Ligand. <i>Inorganic Chemistry</i> , 2018, 57, 3983-3992.	4.0	10
15	Charge and spin coupling in copper compounds with hemilabile noninnocent ligands - Ambivalence in three dimensions. <i>Coordination Chemistry Reviews</i> , 2018, 355, 173-179.	18.8	12
16	Photophysical Heavy-Atom Effect in Iodinated Metalloporphyrins: Spin-Orbit Coupling and Density of States. <i>Journal of Physical Chemistry A</i> , 2018, 122, 7256-7266.	2.5	22
17	Hybrid Diimine/Bis(chalcogenoether) Ligands for Copper(I) and Copper(II) Complexes. <i>Zeitschrift Fur Anorganische Und Allgemeine Chemie</i> , 2018, 644, 661-670.	1.2	6
18	The BIAN ligand 1,2-bis[(2,6-diisopropylphenyl)imino]acenaphthene: An electron sponge or a normal diimine ligand?. <i>Inorganica Chimica Acta</i> , 2017, 455, 540-548.	2.4	9

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19	Electronic structures and photophysics of d8-d8 complexes. <i>Coordination Chemistry Reviews</i> , 2017, 345, 297-317.	18.8	70
20	Electronic Structures of Reduced and Superreduced Ir ₂ (1,8-diisocyanomenthane) ₄ Complexes. <i>Inorganic Chemistry</i> , 2017, 56, 2874-2883.	4.0	5
21	Probing the Intramolecular Metal-Selenoether Interaction in a Bis(iminosemiquinone)copper(II) Compound. <i>Zeitschrift Fur Anorganische Und Allgemeine Chemie</i> , 2017, 643, 1621-1627.	1.2	9
22	Metal carbonyl complexes of potentially ambidentate 2,1,3-benzothiadiazole and 2,1,3-benzoselenadiazole acceptors. <i>Zeitschrift Fur Naturforschung - Section B Journal of Chemical Sciences</i> , 2017, 72, 839-846.	0.7	7
23	Ultrafast Wiggling and Jiggling: Ir ₂ (1,8-diisocyanomenthane) ₄ Complexes. <i>Journal of Physical Chemistry A</i> , 2017, 121, 9275-9283.	2.5	4
24	Reduced and Superreduced Diplatinum Complexes. <i>Journal of the American Chemical Society</i> , 2016, 138, 5699-5705.	13.7	9
25	Extreme Basicity of Biguanide Drugs in Aqueous Solutions: Ion Transfer Voltammetry and DFT Calculations. <i>Journal of Physical Chemistry A</i> , 2016, 120, 7344-7350.	2.5	20
26	Different manifestations of enhanced Ir-acceptor ligation at every redox level of [Os(9-OP) ₂] ⁿ , n = 2+, +, 0, - (9-OP ⁻ = 9-oxidophenalenone and L =) <i>Journal of Physical Chemistry A</i> , 2016, 120, 7344-7350.	2.5	20
27	Complexation of the strontium cation with a macrocyclic lactam receptor: Experimental and theoretical study. <i>Journal of Molecular Liquids</i> , 2016, 214, 171-174.	4.9	2
28	Thermally Tunable Dual Emission of the d ⁸ Dimer [Pt ₂ (1/4-P ₂ O ₅ (BF ₂) ₂) ₄] ⁴⁺ . <i>Inorganic Chemistry</i> , 2016, 55, 2441-2449.	4.1	42
29	Luminescent Diamond Nanoparticles: Physical, Chemical and Biological Aspects of the Phenomenon. <i>Journal of Nanoscience and Nanotechnology</i> , 2015, 15, 1000-1005.	0.9	3
30	Metal-Induced Chelating N ₂ -Bis(4-dimethylaminophenyl)acetamidinyl Radical: A New Chromophore for the Near-Infrared Region. <i>Chemistry - A European Journal</i> , 2015, 21, 12275-12278.	3.3	12
31	Metal-Induced Thiophene Ring Opening and C-C Bond Formation To Produce Unique Hexa-1,3,5-trienediyl-Coupled Non-Innocent Ligand Chelates. <i>Chemistry - A European Journal</i> , 2015, 21, 15163-15166.	3.3	7
32	Electron-Transfer Acceleration Investigated by Time Resolved Infrared Spectroscopy. <i>Accounts of Chemical Research</i> , 2015, 48, 868-876.	15.6	44
33	Ruthenium Styryl Complexes with Ligands Derived from 2-Hydroxy- and 2-Mercaptopyridine and 2-Hydroxy- and 2-Mercaptoquinoline. <i>Organometallics</i> , 2015, 34, 3611-3628.	2.3	20
34	Spin-Orbit TDDFT Electronic Structure of Diplatinum(II,II) Complexes. <i>Inorganic Chemistry</i> , 2015, 54, 3491-3500.	4.0	35
35	Vinyl Ruthenium-Modified Biphenyl and 2,2'-Bipyridines. <i>Inorganic Chemistry</i> , 2015, 54, 3387-3402.	4.0	32
36	Anharmonicity Effects in IR Spectra of [Re(X)(CO) ₃ (1±-diimine)] (1±-diimine = 2,2'-bipyridine or) <i>Journal of Physical Chemistry A</i> , 2015, 119, 10137-10146.	2.5	19

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37	Theoretical study of chromium and nickel-related luminescence centers in molecular-sized nanodiamonds. <i>Diamond and Related Materials</i> , 2015, 58, 122-128.	3.9	3
38	Electronic Excited States of Tungsten(0) Arylisocyanides. <i>Inorganic Chemistry</i> , 2015, 54, 8518-8528.	4.0	34
39	Oxidation and reduction response of π -diimine complexes with tricarbonylrhenium halides and pseudohalides. <i>Journal of Organometallic Chemistry</i> , 2014, 751, 678-685.	1.8	12
40	Correlated Coordination and Redox Activity of a Hemilabile Noninnocent Ligand in Nickel Complexes. <i>Chemistry - A European Journal</i> , 2014, 20, 5414-5422.	3.3	50
41	RuII(π -diimine) or RuIII(π -diimine π)? Structural, Spectroscopic, and Theoretical Evidence for the Stabilization of a Prominent Metal-to-Ligand Charge-Transfer Excited-State Configuration in the Ground State. <i>European Journal of Inorganic Chemistry</i> , 2014, 2014, 110-119.	2.0	23
42	Theoretical Predictions of Redox Potentials of Fischer-Type Chromium Aminocarbene Complexes. <i>Organometallics</i> , 2014, 33, 4964-4972.	2.3	16
43	Magnetical and Optical Properties of Nanodiamonds Can Be Tuned by Particles Surface Chemistry: Theoretical and Experimental Study. <i>Journal of Physical Chemistry C</i> , 2014, 118, 25245-25252.	3.1	25
44	Structural flexibility of 2-hetaryl chromium aminocarbene complexes: Experimental and theoretical evidence. <i>Inorganica Chimica Acta</i> , 2014, 421, 439-445.	2.4	4
45	Photophysics of Singlet and Triplet Intraligand Excited States in [ReCl(CO) ₃ (1-(2-pyridyl)-imidazo[1,5- π]pyridine)] Complexes. <i>Journal of the American Chemical Society</i> , 2014, 136, 5963-5973.	13.7	64
46	Structure and Spectroelectrochemical Response of Arene π -Ruthenium and Arene π -Osmium Complexes with Potentially Hemilabile Noninnocent Ligands. <i>Organometallics</i> , 2014, 33, 4973-4985.	2.3	44
47	Inhibitory Effect of Water on the Oxygen Reduction Catalyzed by Cobalt(II) Tetraphenylporphyrin. <i>Journal of Physical Chemistry A</i> , 2014, 118, 2018-2028.	2.5	16
48	Non π -innocent Redox Behavior of Amidinato Ligands: Spectroscopic Evidence for Amidinyl Complexes. <i>Zeitschrift Fur Anorganische Und Allgemeine Chemie</i> , 2014, 640, 2781-2787.	1.2	9
49	Origin of the correlation between the standard Gibbs energies of ion transfer from water to a hydrophobic ionic liquid and to a molecular solvent. <i>Electrochimica Acta</i> , 2013, 87, 591-598.	5.2	8
50	Complexation of Eu ³⁺ with a macrocyclic lactam receptor: Experimental and theoretical study. <i>Journal of Molecular Structure</i> , 2013, 1038, 216-219.	3.6	2
51	Reversible and pH dependent photophysical properties of mixed-ligand Ru(II) complexes containing 2,2 π -bipyridine and nitrosobarbiturate: Experimental and theoretical approach. <i>Inorganica Chimica Acta</i> , 2013, 404, 123-130.	2.4	0
52	Electronic Excitations in Fischer-Type Cr and W Aminocarbene Complexes: A Combined ab Initio and Experimental Study. <i>Journal of Physical Chemistry A</i> , 2013, 117, 11456-11463.	2.5	11
53	Mechanistic model of the oxygen reduction catalyzed by a metal-free porphyrin in one- and two-phase liquid systems. <i>Electrochimica Acta</i> , 2013, 110, 816-821.	5.2	11
54	Discovering More Non π -innocence: Triazenido versus Triazenyl Radical Ligand Function, and a Comment on [NO ₂] ²⁻ as a π -Suspect π -Ligand. <i>Angewandte Chemie - International Edition</i> , 2013, 52, 4673-4675.	13.8	19

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55	Re and Br X-ray Absorption Near-Edge Structure Study of the Ground and Excited States of [ReBr(CO) ₃ (bpy)] Interpreted by DFT and TD-DFT Calculations. <i>Inorganic Chemistry</i> , 2013, 52, 5775-5785.	4.0	43
56	Identifying Intermediates of Sequential Electron and Hydrogen Loss from a Dicarbonylcobalt Hydride Complex. <i>Angewandte Chemie - International Edition</i> , 2013, 52, 6781-6784.	13.8	30
57	Ruthenium Stilbenyl and Diruthenium Distyrylene Complexes: Aspects of Electron Delocalization and Electrocatalyzed Isomerization of the <i>Z</i> -Isomer. <i>Journal of the American Chemical Society</i> , 2012, 134, 16671-16692.	13.7	89
58	UV-vis-NIR and EPR characterisation of the redox series [MQ ₃] ^{2+,+,0,âˆ’,2âˆ’} , M = Ru or Os, and Q = o-quinone derivative. <i>Dalton Transactions</i> , 2012, 41, 8913.	3.3	16
59	Competitive inhibition of a metal-free porphyrin oxygen-reduction catalyst by water. <i>Chemical Communications</i> , 2012, 48, 4094.	4.1	8
60	Solar Cell Sensitizer Models [Ru(bpy-R) ₂ (NCS) ₂] Probed by Spectroelectrochemistry. <i>Inorganic Chemistry</i> , 2012, 51, 2097-2104.	4.0	36
61	Spinâ€œOrbit Treatment of UVâ€œvis Absorption Spectra and Photophysics of Rhenium(I) Carbonylâ€œBipyridine Complexes: MS-CASPT2 and TD-DFT Analysis. <i>Journal of Physical Chemistry A</i> , 2012, 116, 11319-11329.	2.5	74
62	Structural Reassessment of [W(CO) ₅ (TCNE)]: N (Îf) Coordination Instead of an Olefin (Î€) Complex. <i>Organometallics</i> , 2012, 31, 6305-6311.	2.3	3
63	Charged, but Found â€œNot Guiltyâ€œ Innocence of the Suspect Bridging Ligands [RO(O)CNC(O)OR] ^{2â€œ} = L ^{2â€œ} in [(acac) ₂ Ru(Î¼ ⁴ -L)Ru(acac) ₂] ⁺ⁿ , <i>n</i> = +,0,â€œ,2â€œ. <i>Inorganic Chemistry</i> , 2012, 51, 9273-9281.	4.0	34
64	(Î±-Diimine)tricarbonylhalorhenium complexes: the oxidation side. <i>Dalton Transactions</i> , 2012, 41, 1013-1019.	3.3	13
65	Filling Gaps in the Series of Noninnocent Hetero-1,3-diene Chelate Ligands: Ruthenium Complexes of Redox-Active Î±-Azocarbonyl and Î±-Azothiocarbonyl Ligands RNNC(Râ€²) ₂ E, E = O or S. <i>Inorganic Chemistry</i> , 2012, 51, 6237-6244.	4.0	20
66	Metal(IV) Complexes [M(L _{N,O,S}) ₂] ⁺ⁿ (M = Ru, Os) of a Redox-Active <i>o</i> -Amidophenolate Ligand (L _{N,O,S}) ^{2â€œ} with Coordinating Thioether Appendix. <i>European Journal of Inorganic Chemistry</i> , 2012, 2012, 3569-3576.	2.0	24
67	Charge transfer in porphyrinâ€œcalixarene complexes: ultrafast kinetics, cyclic voltammetry, and DFT calculations. <i>Physical Chemistry Chemical Physics</i> , 2011, 13, 6947.	2.8	19
68	Ultrafast Excited-State Dynamics of Rhenium(I) Photosensitizers [Re(Cl)(CO) ₃ (N,N)] and [Re(imidazole)(CO) ₃ (N,N)] ⁺ⁿ : Diimine Effects. <i>Inorganic Chemistry</i> , 2011, 50, 2932-2943.	4.0	171
69	Fine tuning of the catalytic effect of a metal-free porphyrin on the homogeneous oxygenreduction. <i>Chemical Communications</i> , 2011, 47, 5446-5448.	4.1	31
70	Photoinduced Intramolecular Tryptophan Oxidation and Excited-State Behavior of [Re(L-AA)(CO) ₃ (Î±-diimine)] ⁺ⁿ (L = Pyridine or Imidazole, AA = Tryptophan,) <i>J. Phys. Chem. B</i> , 2011, 115, 10710-10716.	4.0	10
71	Electron delocalization in vinyl ruthenium substituted cyclophanes: Assessment of the through-space and the through-bond pathways. <i>Journal of Organometallic Chemistry</i> , 2011, 696, 3186-3197.	1.8	43
72	Tuning of nanodiamond particles' optical properties by structural defects and surface modifications: DFT modelling. <i>Journal of Materials Chemistry</i> , 2011, 21, 18248.	6.7	26

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73	Complementary hydrogen bonding of a carboxylato-barbiturate with urea and acetamide: Experimental and theoretical approach. <i>Spectrochimica Acta - Part A: Molecular and Biomolecular Spectroscopy</i> , 2011, 83, 532-539.	3.9	2
74	Origin of electronic absorption spectra of MLCT-excited and one-electron reduced 2,2'-bipyridine and 1,10-phenanthroline complexes. <i>Inorganica Chimica Acta</i> , 2011, 374, 578-585.	2.4	67
75	Vinyl-ruthenium entities as markers for intramolecular electron transfer processes. <i>Inorganica Chimica Acta</i> , 2011, 374, 36-50.	2.4	61
76	Phototriggering Electron Flow through Re ^I -modified <i>Pseudomonas aeruginosa</i> Azurins. <i>Chemistry - A European Journal</i> , 2011, 17, 5350-5361.	3.3	51
77	Relativistic effects in spectroscopy and photophysics of heavy-metal complexes illustrated by spin-orbit calculations of [Re(imidazole)(CO) ₃ (phen)] ⁺ . <i>Coordination Chemistry Reviews</i> , 2011, 255, 975-989.	18.8	90
78	Fischer aminocarbene complexes of chromium and iron: Anomalous electrochemical reduction of p-carbonyl substituted derivatives. <i>Electrochimica Acta</i> , 2011, 56, 6853-6859.	5.2	27
79	The 1,4-diazabutadiene/1,2-enediamido non-innocent ligand system in the formation of iridaheteroaromatic compounds: Spectroelectrochemistry and electronic structure. <i>Journal of Organometallic Chemistry</i> , 2010, 695, 1052-1058.	1.8	35
80	Quantum chemical interpretation of redox properties of ruthenium complexes with vinyl and TCNX type non-innocent ligands. <i>Coordination Chemistry Reviews</i> , 2010, 254, 1383-1396.	18.8	93
81	Synthesis and electrochemical study of iron, chromium and tungsten aminocarbenes: Role of ligand structure and central metal nature. <i>Electrochimica Acta</i> , 2010, 55, 8341-8351.	5.2	31
82	Oxygen Reduction Catalyzed by a Fluorinated Tetraphenylporphyrin Free Base at Liquid/Liquid Interfaces. <i>Journal of the American Chemical Society</i> , 2010, 132, 13733-13741.	13.7	80
83	Charge Transport in DNA Oligonucleotides with Various Base-Pairing Patterns. <i>Journal of Physical Chemistry B</i> , 2010, 114, 5196-5205.	2.6	34
84	Ultrafast Excited-State Dynamics of [Re(L)(CO) ₃ (bpy)] ⁿ⁺ Complexes: Involvement of the Solvent. <i>Journal of Physical Chemistry A</i> , 2010, 114, 6361-6369.	2.5	118
85	How to elucidate and control the redox sequence in vinylbenzoate and vinylpyridine bridged diruthenium complexes. <i>Dalton Transactions</i> , 2010, 39, 8000.	3.3	27
86	Ligand-Centred Reactivity of Bis(picoyl)amine Iridium: Sequential Deprotonation, Oxidation and Oxygenation of a Non-Innocent Ligand. <i>Chemistry - A European Journal</i> , 2009, 15, 11878-11889.	3.3	60
87	Boron Atoms as Spin Carriers in Two- and Three-Dimensional Systems. <i>Angewandte Chemie - International Edition</i> , 2009, 48, 5082-5091.	13.8	109
88	An Odd-Electron Complex [Ru ^{III} (NO ⁺)(Q ⁿ⁺)(terpy)] ²⁺ with Two Prototypical Non-Innocent Ligands. <i>Angewandte Chemie - International Edition</i> , 2009, 48, 4242-4245.	13.8	53
89	Evidence for the dimer-of-(mixed-valent dimers) configuration in tetranuclear {(1/4-TCNX)[Ru(NH ₃) ₅] ₄ } ⁸⁺ , TCNX=ATCNE and TCNQ, from DFT calculations. <i>Monatshefte für Chemie</i> , 2009, 140, 765-773.	1.8	24
90	Silver(I), Copper(I) and Copper(II) Complex of the New N-Se-Chelate Ligand 2-Phenylselenomethyl-1H-benzimidazole: Electrochemistry and Structure. <i>Zeitschrift Für Anorganische Und Allgemeine Chemie</i> , 2009, 635, 1001-1007.	1.2	8

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91	2-Organoselenomethyl-1H-benzimidazole Complexes of Copper(II) and Copper(I). <i>Zeitschrift Fur Anorganische Und Allgemeine Chemie</i> , 2009, 635, 2177-2184.	1.2	2
92	The Semiquinone ^{•-} Ruthenium Combination as a Remarkably Invariant Feature in the Redox and Substitution Series [Ru(Q) _n (acac) _{3-n}] ⁺ , $n = 1, 2, 3$; $n = (2, 1, 0, +1, +2)$; Q = 4,6-Di-tert-butyl-N-phenyl-o-iminobenzoquinone. <i>Inorganic Chemistry</i> , 2009, 48, 11853-11864.	4.0	61
93	A Five-Center Redox System: Molecular Coupling of Two Noninnocent Imino-o-benzoquinonato-Ruthenium Functions through a π Acceptor Bridge. <i>Journal of the American Chemical Society</i> , 2009, 131, 8895-8902.	13.7	56
94	Electron Transfer Across Multiple Hydrogen Bonds: The Case of Ureapyrimidinedione-Substituted Vinyl Ruthenium and Osmium Complexes. <i>Journal of the American Chemical Society</i> , 2009, 131, 4892-4903.	13.7	53
95	Charge Delocalization in a Heterobimetallic Ferrocene ^{•-} (Vinyl)Ru(CO)Cl(PiPr ₃) ₂ System ^{•-} Dedicated to Prof. Dr. Helmut Werner on the occasion of his 75th birthday. <i>Organometallics</i> , 2009, 28, 4196-4209.	2.3	79
96	Relaxation Dynamics of <i>Pseudomonas aeruginosa</i> Re ^I (CO) ₃ (\pm -diimine)(HisX) ⁺ (X = 83, 107, 109, 124, 126)Cu ^{II} Azurins. <i>Journal of the American Chemical Society</i> , 2009, 131, 11788-11800.	13.7	55
97	Excited-state relaxation dynamics of Re(I) tricarbonyl complexes with macrocyclic phenanthroline ligands studied by time-resolved IR spectroscopy. <i>Dalton Transactions</i> , 2009, , 3941.	3.3	26
98	Effect of metal exchange (Os vs. Ru) and co-ligand variation (Cl ^{•-} vs. acac ^{•-}) on the oxidation state distribution in complexes of an o-phenylenediamido(2-quinone)diimine redox system. <i>Dalton Transactions</i> , 2009, , 7778.	3.3	20
99	Localised to intraligand charge-transfer states in cyclometalated platinum complexes: an experimental and theoretical study into the influence of electron-rich pendants and modulation of excited states by ion binding. <i>Dalton Transactions</i> , 2009, , 1728.	3.3	85
100	How do vibrations change their composition upon electronic excitation? ^{•-} EXSY-T2D-IR measurements challenge DFT calculations.. <i>Springer Series in Chemical Physics</i> , 2009, , 421-423.	0.2	0
101	Dimerization of the 2-(Methylselenomethyl)benzimidazole ^{•-} silver(I) Cation Involving a Ag ^I Interaction. <i>Zeitschrift Fur Anorganische Und Allgemeine Chemie</i> , 2008, 634, 2343-2346.	1.2	9
102	Halide Ligands ^{•-} More Than Just σ -Donors? A Structural and Spectroscopic Study of Homologous Organonickel Complexes. <i>Inorganic Chemistry</i> , 2008, 47, 11324-11333.	4.0	44
103	Redox Properties of Ruthenium Nitrosyl Porphyrin Complexes with Different Axial Ligation: Structural, Spectroelectrochemical (IR, UV-Visible, and EPR), and Theoretical Studies. <i>Inorganic Chemistry</i> , 2008, 47, 7106-7113.	4.0	50
104	Singlet Diradical Complexes of Ruthenium and Osmium: Geometrical and Electronic Structures and their Unexpected Changes on Oxidation. <i>Inorganic Chemistry</i> , 2008, 47, 1625-1633.	4.0	46
105	A tetranuclear organorhenium(I) complex of the 2,3,5,6-tetrafluoro-7,7,8,8-tetracyano-p-quinodimethane radical anion, TCNQF ₄ ^{•-} . <i>Dalton Transactions</i> , 2008, , 5749.	3.3	15
106	Stabilizing the Elusive ortho-Quinone/Copper(I) Oxidation State Combination through π - π Interaction in an Isolated Complex. <i>Journal of the American Chemical Society</i> , 2008, 130, 15230-15231.	13.7	69
107	Solvation-Driven Excited-State Dynamics of [Re(4-Et-Pyridine)(CO) ₃ (2,2'-bipyridine)] ⁺ in Imidazolium Ionic Liquids. A Time-Resolved Infrared and Phosphorescence Study. <i>Journal of Physical Chemistry A</i> , 2008, 112, 3506-3514.	2.5	31
108	Ligand-Centered Oxidations and Electron Delocalization in a Tetranuclear Complex of a Tetrador-Substituted Olefin. <i>Organometallics</i> , 2008, 27, 3321-3324.	2.3	46

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109	Nanosecond CO Photodissociation and Excited-State Character of [Ru(X)(X ^ε)(CO) ₂] (X = N, N ^ε -diisopropyl-1,4-diazabutadiene) (X = X ^ε = Cl or I; X = Me, X ^ε) Tj FTQq1 10. Calculations. <i>Inorganic Chemistry</i> , 2008, 47, 4236-4242.	4.0	110
110	Ruthenium Complexes with Vinyl, Styryl, and Vinylpyrenyl Ligands: A Case of Non-innocence in Organometallic Chemistry. <i>Journal of the American Chemical Society</i> , 2008, 130, 259-268.	13.7	111
111	Femtosecond Fluorescence and Intersystem Crossing in Rhenium(I) Carbonyl ^π Bipyridine Complexes. <i>Journal of the American Chemical Society</i> , 2008, 130, 8967-8974.	13.7	269
112	Electronic structures of two kinds of oligoborane radical anions: mixed-valence description of a reduced <i>p</i> -phenylenediborane and spin density distributions in distorted octahedral clusters [B ₆ Hal ₆] ^{•-} , Hal = Al, Br, I. <i>Main Group Chemistry</i> , 2007, 5, 267-276.	0.8	10
113	Raman Spectroscopy of Nanocrystalline Li-Ti-O Spinel and Comparative DFT Calculations on Ti ₂ O ₃ and LiTiO ₃ Clusters. <i>Collection of Czechoslovak Chemical Communications</i> , 2007, 72, 171-184.	1.0	5
114	Singlet Diradical Complexes of Chromium, Molybdenum, and Tungsten with Azo Anion Radical Ligands from M(CO) ₆ Precursors. <i>Inorganic Chemistry</i> , 2007, 46, 8584-8593.	4.0	44
115	Spectroelectrochemistry and DFT Analysis of a New {RuNO} ⁿ Redox System with Multifrequency EPR Suggesting Conformational Isomerism in the {RuNO} ⁷ State. <i>Inorganic Chemistry</i> , 2007, 46, 9254-9261.	4.0	21
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