# Lai-Sheng Wang

# List of Publications by Year in Descending Order

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160 34,299 101 479 h-index g-index citations papers 36,946 489 7.6 7.45 avg, IF L-index ext. citations ext. papers

#	Paper	IF	Citations
479	Probing the Nature of the Transition-Metal-Boron Bonds and Novel Aromaticity in Small Metal-Doped Boron Clusters Using Photoelectron Spectroscopy <i>Annual Review of Physical Chemistry</i> , <b>2022</b> ,	15.7	1
478	AuB: an Au-borozene complex Chemical Communications, 2022,	5.8	1
477	Observation of Core-Excited Dipole-Bound States <i>Journal of Physical Chemistry Letters</i> , <b>2022</b> , 2124-212	<b>28</b> .4	3
476	A Heteroleptic Gold Hydride Nanocluster for Efficient and Selective Electrocatalytic Reduction of CO to CO <i>Journal of the American Chemical Society</i> , <b>2022</b> ,	16.4	9
475	Probing copper-boron interactions in the Cu2B8Ibimetallic cluster. <i>Journal of Vacuum Science and Technology A: Vacuum, Surfaces and Films</i> , <b>2022</b> , 40, 042201	2.9	
474	Monovalent lanthanide(I) in borozene complexes. <i>Nature Communications</i> , <b>2021</b> , 12, 6467	17.4	3
473	Photodetachment spectroscopy and resonant photoelectron imaging of cryogenically cooled 1-pyrenolate. <i>Journal of Chemical Physics</i> , <b>2021</b> , 154, 094308	3.9	8
472	Expanded Inverse-Sandwich Complexes of Lanthanum Borides: LaB and LaB. <i>Journal of Physical Chemistry A</i> , <b>2021</b> , 125, 2622-2630	2.8	10
471	How O-Binding Affects Structural Evolution of Medium Even-Sized Gold Clusters Au ( = 20-34). Journal of Physical Chemistry Letters, <b>2021</b> , 12, 3560-3570	6.4	3
470	Probing the Dipole-Bound State in the 9-Phenanthrolate Anion by Photodetachment Spectroscopy, Resonant Two-Photon Photoelectron Imaging, and Resonant Photoelectron Spectroscopy. <i>Journal of Physical Chemistry A</i> , <b>2021</b> , 125, 2967-2976	2.8	7
469	Double FAromaticity in a Planar Zinc-Doped Gold Cluster: AuZn. <i>Journal of Physical Chemistry A</i> , <b>2021</b> , 125, 4606-4613	2.8	9
468	The synthesis and characterization of a new diphosphine-protected gold hydride nanocluster. Journal of Chemical Physics, <b>2021</b> , 155, 034307	3.9	3
467	The Synthesis, Bonding, and Transformation of a Ligand-Protected Gold Nanohydride Cluster. <i>Angewandte Chemie - International Edition</i> , <b>2021</b> , 60, 2424-2430	16.4	16
466	The Synthesis, Bonding, and Transformation of a Ligand-Protected Gold Nanohydride Cluster. <i>Angewandte Chemie</i> , <b>2021</b> , 133, 2454-2460	3.6	4
465	B: a bilayer boron cluster. <i>Nanoscale</i> , <b>2021</b> , 13, 3868-3876	7.7	16
464	Photoelectron Spectroscopy of Size-Selected Bismuth-Boron Clusters: BiB ( = 6-8). <i>Journal of Physical Chemistry A</i> , <b>2021</b> , 125, 6751-6760	2.8	6
463	Observation of a dipole-bound excited state in 4-ethynylphenoxide and comparison with the quadrupole-bound excited state in the isoelectronic 4-cyanophenoxide. <i>Journal of Chemical Physics</i> , <b>2021</b> , 155, 124305	3.9	4

462	Transition-metal-like bonding behaviors of a boron atom in a boron-cluster boronyl complex [(B)-B-BO]. <i>Chemical Science</i> , <b>2021</b> , 12, 8157-8164		4
461	The nature of the chemical bonding in 5d transition-metal diatomic borides MB (M = Ir, Pt, Au).  Journal of Chemical Physics, <b>2020</b> , 152, 174301  3.9	,	6
460	Observation of Transition-Metal-Boron Triple Bonds in IrB O and ReB O. <i>Angewandte Chemie - International Edition</i> , <b>2020</b> , 59, 15260-15265	· .	3
459	Observation of Transition-MetalBoron Triple Bonds in IrB2Oland ReB2OllAngewandte Chemie, 3.6		
458	Spherical trihedral metallo-borospherenes. <i>Nature Communications</i> , <b>2020</b> , 11, 2766		20
457	MnB: An Open-Shell Metallaboron Analog of 3d Metallabenzenes. <i>Journal of Physical Chemistry A</i> , <b>2.8</b>		7
456	High-Resolution Photoelectron Imaging and Photodetachment Spectroscopy of Cryogenically Cooled IO. <i>Journal of Physical Chemistry A</i> , <b>2020</b> , 124, 5720-5726		1
455	Photodetachment spectroscopy and resonant photoelectron imaging of the 2-naphthoxide anion via dipole-bound excited states. <i>Journal of Chemical Physics</i> , <b>2020</b> , 152, 214307		6
454	High-resolution photoelectron imaging of MnB : Probing the bonding between the aromatic B cluster and 3d transition metals. <i>Journal of Chemical Physics</i> , <b>2020</b> , 152, 244306		2
453	Observation of MBius Aromatic Planar Metallaborocycles. <i>Journal of the American Chemical Society</i> , <b>2020</b> , 142, 3356-3360		13
452	Observation of Four-Fold Boron-Metal Bonds in RhB(BO) and RhB. <i>Journal of Physical Chemistry Letters</i> , <b>2020</b> , 11, 659-663		21
45 <sup>1</sup>	Observation of Backbonding in a Boronyl-Coordinated Transition Metal Complex TaBO. <i>Journal of Physical Chemistry A</i> , <b>2020</b> , 124, 10001-10007		
450	Observation of a Symmetry-Forbidden Excited Quadrupole-Bound State. <i>Journal of the American Chemical Society</i> , <b>2020</b> , 142, 20240-20246		6
449	Observation of a Type Dipole-Bound State in Molecular Anions. <i>Physical Review Letters</i> , <b>2020</b> , 125, 073003		16
448	Halogen effects on the electronic and optical properties of Au13 nanoclusters. <i>Nanoscale Advances</i> , 2020, 2, 4902-4907		8
447	Polarization of Valence Orbitals by the Intramolecular Electric Field from a Diffuse Dipole-Bound Electron. <i>Journal of Physical Chemistry Letters</i> , <b>2020</b> , 11, 7914-7919		11
446	Probing the electronic structure of the CoB16Idrum complex: Unusual oxidation state of CoIII Chinese Journal of Chemical Physics, <b>2019</b> , 32, 241-247	,	3
445	Probing the coupling of a dipole-bound electron with a molecular core. <i>Chemical Science</i> , <b>2019</b> , 10, 1386 <u>4.3</u> 9	1	13

444	Facile Synthesis of Unsolvated Alkali Metal Octahydrotriborate Salts MB3H8 (M=K, Rb, and Cs), Mechanisms of Formation, and the Crystal Structure of KB3H8. <i>Angewandte Chemie</i> , <b>2019</b> , 131, 2746-27	730 <sup>6</sup>	11
443	Facile Synthesis of Unsolvated Alkali Metal Octahydrotriborate Salts MB H (M=K, Rb, and Cs), Mechanisms of Formation, and the Crystal Structure of KB H. <i>Angewandte Chemie - International Edition</i> , <b>2019</b> , 58, 2720-2724	16.4	29
442	[La([]B)La] (= 7-9): a new class of inverse sandwich complexes. <i>Chemical Science</i> , <b>2019</b> , 10, 2534-2542	9.4	42
441	Double- and multi-slit interference in photodetachment from nanometer organic molecular anions. Journal of Chemical Physics, <b>2019</b> , 150, 244302	3.9	
440	Re[] B and Re[] B: New Members of the Transition-Metal-Centered Borometallic Molecular Wheel Family. <i>Journal of Physical Chemistry A</i> , <b>2019</b> , 123, 5317-5324	2.8	23
439	LaB: an inverse triple-decker lanthanide boron cluster. <i>Chemical Communications</i> , <b>2019</b> , 55, 7864-7867	5.8	25
438	Probing the structures and bonding of size-selected boron and doped-boron clusters. <i>Chemical Society Reviews</i> , <b>2019</b> , 48, 3550-3591	58.5	90
437	B and B: chiral quasi-planar boron clusters. <i>Nanoscale</i> , <b>2019</b> , 11, 9698-9704	7.7	18
436	High-Resolution Photoelectron Imaging of IrB: Observation of a EAromatic B Ring Coordinated to a Transition Metal. <i>Angewandte Chemie - International Edition</i> , <b>2019</b> , 58, 8877-8881	16.4	19
435	High-Resolution Photoelectron Imaging of IrB3ElObservation of a EAromatic B3+ Ring Coordinated to a Transition Metal. <i>Angewandte Chemie</i> , <b>2019</b> , 131, 8969-8973	3.6	5
434	Tautomer-Specific Resonant Photoelectron Imaging of Deprotonated Cytosine Anions. <i>Angewandte Chemie - International Edition</i> , <b>2019</b> , 58, 7856-7860	16.4	7
433	Tautomer-Specific Resonant Photoelectron Imaging of Deprotonated Cytosine Anions. <i>Angewandte Chemie</i> , <b>2019</b> , 131, 7938-7942	3.6	
432	Au: The Smallest Gold Cluster with the High-Symmetry Icosahedral Core Au. <i>Journal of Physical Chemistry Letters</i> , <b>2019</b> , 10, 1820-1827	6.4	12
431	Resonant Two-Photon Photoelectron Imaging and Intersystem Crossing from Excited Dipole-Bound States of Cold Anions. <i>Journal of Physical Chemistry Letters</i> , <b>2019</b> , 10, 4339-4344	6.4	19
430	Probing the Critical Dipole Moment To Support Excited Dipole-Bound States in Valence-Bound Anions. <i>Journal of Physical Chemistry Letters</i> , <b>2019</b> , 10, 6472-6477	6.4	30
429	ReB: A Metallaboron Analog of Metallabenzenes. <i>Journal of the American Chemical Society</i> , <b>2019</b> , 141, 17854-17860	16.4	12
428	High resolution photoelectron imaging of boron-bismuth binary clusters: BiB (n = 2-4). <i>Journal of Chemical Physics</i> , <b>2019</b> , 150, 064304	3.9	7
427	Planar B and B clusters with double-hexagonal vacancies. <i>Nanoscale</i> , <b>2019</b> , 11, 23286-23295	7.7	29

#### (2017-2019)

426	High-resolution photoelectron imaging and resonant photoelectron spectroscopy noncovalently bound excited states of cryogenically cooled anions. <i>Chemical Science</i> , <b>2019</b> , 10, 9409-9423	9.4	31	
425	Lanthanides with Unusually Low Oxidation States in the PrB and PrB Boride Clusters. <i>Inorganic Chemistry</i> , <b>2019</b> , 58, 411-418	5.1	23	
424	Structural Evolution of Gold-Doped Bismuth Clusters AuBin $\mathbb{Q}$ n = 4 $\mathbb{B}$ ). <i>Journal of Physical Chemistry C</i> , <b>2018</b> , 122, 6947-6954	3.8	13	
423	A high-resolution photoelectron imaging and theoretical study of CP and CP. <i>Journal of Chemical Physics</i> , <b>2018</b> , 148, 044301	3.9	7	
422	[(CpM)BH] (MI Zr or Hf): early transition metal 'guarded' heptaborane with strong covalent and electrostatic bonding. <i>Chemical Science</i> , <b>2018</b> , 9, 1976-1981	9.4	14	
421	Toward Solution Syntheses of the Tetrahedral Au Pyramid and Atomically Precise Gold Nanoclusters with Uncoordinated Sites. <i>Accounts of Chemical Research</i> , <b>2018</b> , 51, 2159-2168	24.3	48	
420	Observation of highly stable and symmetric lanthanide octa-boron inverse sandwich complexes. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , <b>2018</b> , 115, E6972-E6977	<b>,</b> 11.5	59	
419	Elucidation of the Formation Mechanisms of the Octahydrotriborate Anion (BH) through the Nucleophilicity of the B-H Bond. <i>Journal of the American Chemical Society</i> , <b>2018</b> , 140, 6718-6726	16.4	49	
418	Probing the interaction between the encapsulated water molecule and the fullerene cages in HO@C and HO@CN. <i>Chemical Science</i> , <b>2018</b> , 9, 5666-5671	9.4	13	
417	Recent Progress on the investigations of boron clusters and boron-based materials (I): borophene. <i>Scientia Sinica Chimica</i> , <b>2018</b> , 48, 98-107	1.6	10	
416	Dipole-bound excited states and resonant photoelectron imaging of phenoxide and thiophenoxide anions. <i>Journal of Chemical Physics</i> , <b>2018</b> , 149, 164301	3.9	22	
415	Probing the structures and bonding of auropolyynes, Au-(C?C) -Au ( = 1-3), using high-resolution photoelectron imaging. <i>Journal of Chemical Physics</i> , <b>2018</b> , 149, 144307	3.9	7	
414	Determination of CO Adsorption Sites on Gold Clusters Au (n = 21-25): A Size Region That Bridges the Pyramidal and Core-Shell Structures. <i>Journal of Physical Chemistry Letters</i> , <b>2018</b> , 9, 5430-5439	6.4	5	
413	Di-niobium gold clusters: Multiply-bonded Nb2 dimer coordinated equatorially by Au atoms. <i>International Journal of Mass Spectrometry</i> , <b>2018</b> , 434, 7-16	1.9	1	
412	Photodetachment spectroscopy and resonant photoelectron imaging of cryogenically-cooled deprotonated 2-hydroxypyrimidine anions. <i>Journal of Molecular Spectroscopy</i> , <b>2017</b> , 332, 86-93	1.3	12	
411	B26[]The smallest planar boron cluster with a hexagonal vacancy and a complicated potential landscape. <i>Chemical Physics Letters</i> , <b>2017</b> , 683, 336-341	2.5	33	
410	Planar B and B clusters with a double-hexagonal vacancy: molecular motifs for borophenes. <i>Nanoscale</i> , <b>2017</b> , 9, 4550-4557	7.7	61	
409	Conformation-selective resonant photoelectron imaging from dipole-bound states of cold 3-hydroxyphenoxide. <i>Journal of Chemical Physics</i> , <b>2017</b> , 147, 013910	3.9	20	

408	PrB : A Praseodymium-Doped Boron Cluster with a Pr Center Coordinated by a Doubly Aromatic Planar BB Ligand. <i>Angewandte Chemie - International Edition</i> , <b>2017</b> , 56, 6916-6920	16.4	46
407	PrB7🛮A Praseodymium-Doped Boron Cluster with a PrII Center Coordinated by a Doubly Aromatic Planar 🗗 -B73 Ligand. <i>Angewandte Chemie</i> , <b>2017</b> , 129, 7020-7024	3.6	12
406	Bismuth-Boron Multiple Bonding in BiB O and Bi B. <i>Angewandte Chemie - International Edition</i> , <b>2017</b> , 56, 9551-9555	16.4	19
405	Recent progresses of global minimum searches of nanoclusters with a constrained Basin-Hopping algorithm in the TGMin program. <i>Computational and Theoretical Chemistry</i> , <b>2017</b> , 1107, 57-65	2	59
404	Observation of a metal-centered B-Ta@B tubular molecular rotor and a perfect Ta@B boron drum with the record coordination number of twenty. <i>Chemical Communications</i> , <b>2017</b> , 53, 1587-1590	5.8	90
403	Probing the Structures of Neutral B11 and B12 Using High-Resolution Photoelectron Imaging of B11 and B12 Dournal of Physical Chemistry C, <b>2017</b> , 121, 10752-10759	3.8	14
402	From planar boron clusters to borophenes and metalloborophenes. <i>Nature Reviews Chemistry</i> , <b>2017</b> , 1,	34.6	118
401	Nb[] Au: a molecular wheel with a short Nb 00000000000000000000000000000000000	9.4	12
400	Observation of Excited Quadrupole-Bound States in Cold Anions. <i>Physical Review Letters</i> , <b>2017</b> , 119, 023002	7.4	28
399	B99BMD B99BAPONDatic Plainal Science <b>2017</b> WITT TEXAGONAL Vacancy. European Journal of Inorganic Chemistry, <b>2017</b> , 2017, 4546-4551	2.3	28
398	High-Resolution Photoelectron Imaging of Cryogenically-Cooled CN and (CN) Azafullerene Anions. Journal of Physical Chemistry Letters, <b>2017</b> , 8, 6220-6225	6.4	4
397	Bismuth <b>B</b> oron Multiple Bonding in BiB2O[and Bi2B[] <i>Angewandte Chemie</i> , <b>2017</b> , 129, 9679-9683	3.6	4
396	Resonant photoelectron imaging of deprotonated uracil anion via vibrational levels of a dipole-bound excited state. <i>Chemical Physics</i> , <b>2017</b> , 482, 374-383	2.3	24
395	Probing the Structural Evolution of GoldAluminum Bimetallic Clusters (Au2AlnIn = 311) Using Photoelectron Spectroscopy and Theoretical Calculations. <i>Journal of Physical Chemistry C</i> , <b>2017</b> , 121, 18234-18243	3.8	17
394	Bond-bending isomerism of Aul: competition between covalent bonding and aurophilicity. <i>Chemical Science</i> , <b>2016</b> , 7, 475-481	9.4	14
393	Structural Evolution of Core-Shell Gold Nanoclusters: Au (n = 42-50). ACS Nano, <b>2016</b> , 10, 10013-10022	16.7	32
392	Polymorphism of Phosphine-Protected Gold Nanoclusters: Synthesis and Characterization of a New 22-Gold-Atom Cluster. <i>Small</i> , <b>2016</b> , 12, 2518-25	11	27
391	All-Metal Antiaromaticity in Sb4 -Type Lanthanocene Anions. <i>Angewandte Chemie - International Edition</i> , <b>2016</b> , 55, 5531-5	16.4	46 

# (2016-2016)

390	Beyond organic chemistry: aromaticity in atomic clusters. <i>Physical Chemistry Chemical Physics</i> , <b>2016</b> , 18, 11589-605	3.6	88
389	Probing the Electronic Structure and Chemical Bonding of Mono-Uranium Oxides with Different Oxidation States: $UOx(-)$ and $UOx$ (x = 3-5). <i>Journal of Physical Chemistry A</i> , <b>2016</b> , 120, 1084-96	2.8	22
388	Photoelectron spectroscopy of size-selected boron clusters: from planar structures to borophenes and borospherenes. <i>International Reviews in Physical Chemistry</i> , <b>2016</b> , 35, 69-142	7	195
387	Photoelectron Spectroscopy of BiAu(-) and BiBO(-): Further Evidence of the Analogy between Au and Boronyl. <i>Journal of Physical Chemistry B</i> , <b>2016</b> , 120, 1635-40	3.4	14
386	Frontispiz: The Planar CoB18[Cluster as a Motif for Metallo-Borophenes. <i>Angewandte Chemie</i> , <b>2016</b> , 128,	3.6	1
385	The Planar CoB18 (-) Cluster as a Motif for Metallo-Borophenes. <i>Angewandte Chemie - International Edition</i> , <b>2016</b> , 55, 7358-63	16.4	71
384	Hollow Gold Cages and Their Topological Relationship to Dual Fullerenes. <i>Chemistry - A European Journal</i> , <b>2016</b> , 22, 8823-34	4.8	13
383	Hollow Gold Cages and Their Topological Relationship to Dual Fullerenes. <i>Chemistry - A European Journal</i> , <b>2016</b> , 22, 8709-8709	4.8	
382	A combined photoelectron spectroscopy and relativistic ab initio studies of the electronic structures of UFO and UFO(-). <i>Journal of Chemical Physics</i> , <b>2016</b> , 144, 084309	3.9	4
381	Probing the electronic structure and Au <b>L</b> chemical bonding in AuCnland AuCnHl(n = 2, 4, and 6) using high-resolution photoelectron spectroscopy. <i>Journal of Chemical Physics</i> , <b>2016</b> , 145, 064304	3.9	15
380	Observation and characterization of the smallest borospherene, B28(-) and B28. <i>Journal of Chemical Physics</i> , <b>2016</b> , 144, 064307	3.9	119
379	Time-resolved photoelectron spectroscopy of a dinuclear Pt(II) complex: Tunneling autodetachment from both singlet and triplet excited states of a molecular dianion. <i>Journal of Chemical Physics</i> , <b>2016</b> , 144, 054305	3.9	13
378	Manganese-centered tubular boron cluster - MnB16 (-): A new class of transition-metal molecules. Journal of Chemical Physics, <b>2016</b> , 144, 154310	3.9	84
377	Probing the structures of gold-aluminum alloy clusters AuxAly(-): a joint experimental and theoretical study. <i>Nanoscale</i> , <b>2016</b> , 8, 9805-14	7.7	22
376	Catalyst design based on agostic interactions: synthesis, characterization, and catalytic activity of bis(pyrazolyl)borate copper complexes. <i>Dalton Transactions</i> , <b>2016</b> , 45, 10194-9	4.3	13
375	Second-Order Nonlinear Optical Scattering Properties of Phosphine-Protected Au20 Clusters. <i>Industrial &amp; Engineering Chemistry Research</i> , <b>2016</b> , 55, 10500-10506	3.9	12
374	Diphosphine-Protected Au Nanoclusters on Oxide Supports Are Active for Gas-Phase Catalysis without Ligand Removal. <i>Nano Letters</i> , <b>2016</b> , 16, 6560-6567	11.5	70
373	Competition between quasi-planar and cage-like structures in the B cluster: photoelectron spectroscopy and ab initio calculations. <i>Physical Chemistry Chemical Physics</i> , <b>2016</b> , 18, 29147-29155	3.6	71

372	Competition between drum and quasi-planar structures in RhB: motifs for metallo-boronanotubes and metallo-borophenes. <i>Chemical Science</i> , <b>2016</b> , 7, 7020-7027	9.4	78
371	On the gold-ligand covalency in linear [AuX2](-) complexes. <i>Dalton Transactions</i> , <b>2015</b> , 44, 5535-46	4.3	20
370	Photoelectron spectroscopy and theoretical studies of gaseous uranium hexachlorides in different oxidation states: UCl6 (q-) ( $q = 0-2$ ). <i>Journal of Chemical Physics</i> , <b>2015</b> , 142, 134308	3.9	29
369	Conformation-Selective Resonant Photoelectron Spectroscopy via Dipole-Bound States of Cold Anions. <i>Journal of Physical Chemistry Letters</i> , <b>2015</b> , 6, 2153-7	6.4	22
368	Cobalt-centred boron molecular drums with the highest coordination number in the CoB16- cluster. <i>Nature Communications</i> , <b>2015</b> , 6, 8654	17.4	147
367	Communication: Observation of dipole-bound state and high-resolution photoelectron imaging of cold acetate anions. <i>Journal of Chemical Physics</i> , <b>2015</b> , 142, 091103	3.9	22
366	Vibrational state-selective autodetachment photoelectron spectroscopy from dipole-bound states of cold 2-hydroxyphenoxide: o-HO(C6H4)O(-). <i>Journal of Chemical Physics</i> , <b>2015</b> , 142, 124309	3.9	27
365	Perspective: Electrospray photoelectron spectroscopy: From multiply-charged anions to ultracold anions. <i>Journal of Chemical Physics</i> , <b>2015</b> , 143, 040901	3.9	55
364	B27(-): Appearance of the smallest planar boron cluster containing a hexagonal vacancy. <i>Journal of Chemical Physics</i> , <b>2015</b> , 142, 204305	3.9	49
363	Probing the vibrational spectroscopy of the deprotonated thymine radical by photodetachment and state-selective autodetachment photoelectron spectroscopy dipole-bound states. <i>Chemical Science</i> , <b>2015</b> , 6, 3129-3138	9.4	29
362	Experimental and theoretical evidence of an axially chiral borospherene. ACS Nano, 2015, 9, 754-60	16.7	195
361	Vibrational State-Selective Resonant Two-Photon Photoelectron Spectroscopy of AuS(-) via a Spin-Forbidden Excited State. <i>Journal of Physical Chemistry Letters</i> , <b>2015</b> , 6, 637-42	6.4	19
360	Communication: Vibrationally resolved photoelectron spectroscopy of the tetracyanoquinodimethane (TCNQ) anion and accurate determination of the electron affinity of TCNQ. <i>Journal of Chemical Physics</i> , <b>2015</b> , 143, 221102	3.9	28
359	[BIP: a quasiplanar chiral boron cluster. <i>Angewandte Chemie - International Edition</i> , <b>2014</b> , 53, 5540-5	16.4	116
358	Strong electron correlation in UO2(-): a photoelectron spectroscopy and relativistic quantum chemistry study. <i>Journal of Chemical Physics</i> , <b>2014</b> , 140, 094306	3.9	23
357	Planar hexagonal B(36) as a potential basis for extended single-atom layer boron sheets. <i>Nature Communications</i> , <b>2014</b> , 5, 3113	17.4	503
356	Synthesis and structure determination of a new Au(20) nanocluster protected by tripodal tetraphosphine ligands. <i>Inorganic Chemistry</i> , <b>2014</b> , 53, 3932-4	5.1	68
355	High-resolution photoelectron imaging of cold Clanions and accurate determination of the electron affinity of Cli Journal of Chemical Physics, <b>2014</b> , 140, 224315	3.9	73

354	The B35 cluster with a double-hexagonal vacancy: a new and more flexible structural motif for borophene. <i>Journal of the American Chemical Society</i> , <b>2014</b> , 136, 12257-60	16.4	250
353	A photoelectron spectroscopy and ab initio study of the structures and chemical bonding of the B25(-) cluster. <i>Journal of Chemical Physics</i> , <b>2014</b> , 141, 034303	3.9	54
352	Assessment of Quantum Mechanical Methods for Copper and Iron Complexes by Photoelectron Spectroscopy. <i>Journal of Chemical Theory and Computation</i> , <b>2014</b> , 10, 1283-1291	6.4	16
351	Controlling gold nanoclusters by diphospine ligands. <i>Journal of the American Chemical Society</i> , <b>2014</b> , 136, 92-5	16.4	187
350	[B30]EA Quasiplanar Chiral Boron Cluster. <i>Angewandte Chemie</i> , <b>2014</b> , 126, 5646-5651	3.6	28
349	Observation of an all-boron fullerene. <i>Nature Chemistry</i> , <b>2014</b> , 6, 727-31	17.6	590
348	Complexes between planar boron clusters and transition metals: a photoelectron spectroscopy and ab initio study of CoB12(-) and RhB12(-). <i>Journal of Physical Chemistry A</i> , <b>2014</b> , 118, 8098-105	2.8	111
347	Electronic structure and chemical bonding of a highly stable and aromatic auro-aluminum oxide cluster. <i>Journal of Physical Chemistry A</i> , <b>2014</b> , 118, 5204-11	2.8	7
346	Isomerism and structural fluxionality in the Au26 and Au26(-) nanoclusters. ACS Nano, 2014, 8, 7413-22	16.7	40
345	Understanding boron through size-selected clusters: structure, chemical bonding, and fluxionality. <i>Accounts of Chemical Research</i> , <b>2014</b> , 47, 1349-58	24.3	382
344	Boronyl chemistry: the BO group as a new ligand in gas-phase clusters and synthetic compounds. <i>Accounts of Chemical Research</i> , <b>2014</b> , 47, 2435-45	24.3	64
343	Hexagonal Bipyramidal [Ta2B6] 100 Clusters: B6 Rings as Structural Motifs. <i>Angewandte Chemie</i> , <b>2014</b> , 126, 1312-1316	3.6	26
342	Vibrational Spectroscopy of the Dehydrogenated Uracil Radical by Autodetachment of Dipole-Bound Excited States of Cold Anions. <i>Angewandte Chemie</i> , <b>2014</b> , 126, 2496-2500	3.6	3
341	Hexagonal bipyramidal [Ta(2)B(6)](-/0) clusters: B(6) rings as structural motifs. <i>Angewandte Chemie - International Edition</i> , <b>2014</b> , 53, 1288-92	16.4	38
340	Vibrational spectroscopy of the dehydrogenated uracil radical by autodetachment of dipole-bound excited states of cold anions. <i>Angewandte Chemie - International Edition</i> , <b>2014</b> , 53, 2464-8	16.4	61
339	Probing the electronic and vibrational structure of Au2Al2(-) and Au2Al2 using photoelectron spectroscopy and high resolution photoelectron imaging. <i>Journal of Chemical Physics</i> , <b>2014</b> , 141, 22430	93.9	8
338	High resolution photoelectron imaging of UO(-) and UO2(-) and the low-lying electronic states and vibrational frequencies of UO and UO2. <i>Journal of Chemical Physics</i> , <b>2014</b> , 141, 244302	3.9	12
337	The design and construction of a high-resolution velocity-map imaging apparatus for photoelectron spectroscopy studies of size-selected clusters. <i>Review of Scientific Instruments</i> , <b>2014</b> , 85, 083106	1.7	107

336	Probing the electronic structure and Au-C chemical bonding in AuC2(-) and AuC2 using high-resolution photoelectron spectroscopy. <i>Journal of Chemical Physics</i> , <b>2014</b> , 140, 084303	3.9	24
335	Probing the nature of gold-carbon bonding in gold-alkynyl complexes. <i>Nature Communications</i> , <b>2013</b> , 4, 2223	17.4	44
334	Electron tunneling from electronically excited states of isolated bisdisulizole-derived trianion chromophores following UV absorption. <i>Physical Chemistry Chemical Physics</i> , <b>2013</b> , 15, 6726-36	3.6	15
333	Transition-metal-centered monocyclic boron wheel clusters (MI Bn): a new class of aromatic borometallic compounds. <i>Accounts of Chemical Research</i> , <b>2013</b> , 46, 350-8	24.3	184
332	Photoelectron spectroscopy of aromatic compound clusters of the B12 all-boron benzene: B12Au-and B12(BO) <i>Physical Chemistry Chemical Physics</i> , <b>2013</b> , 15, 9646-53	3.6	39
331	Geometric and electronic factors in the rational design of transition-metal-centered boron molecular wheels. <i>Journal of Chemical Physics</i> , <b>2013</b> , 138, 134315	3.9	53
330	Probing the electronic structure and chemical bonding in tricoordinate uranyl complexes UO2X3- (X = F, Cl, Br, I): competition between Coulomb repulsion and U-X bonding. <i>Inorganic Chemistry</i> , <b>2013</b> , 52, 6617-26	5.1	46
329	Monohafnium oxide clusters $HfO(n)$ - and $HfO(n)$ ( $n = 1-6$ ): oxygen radicals, superoxides, peroxides, diradicals, and triradicals. <i>Journal of Physical Chemistry A</i> , <b>2013</b> , 117, 1042-52	2.8	20
328	High resolution photoelectron imaging of Au2(-). Journal of Chemical Physics, 2013, 138, 184304	3.9	41
327	On the structures and bonding in boron-gold alloy clusters: B6Au(n)- and B6Au(n) (n = 1-3). <i>Journal of Chemical Physics</i> , <b>2013</b> , 138, 084306	3.9	21
326	A photoelectron spectroscopy and density functional study of di-tantalum boride clusters: $Ta2B(x)-(x=2-5)$ . <i>Journal of Chemical Physics</i> , <b>2013</b> , 138, 034308	3.9	27
325	Pi and sigma double conjugations in boronyl polyboroene nanoribbons: B(n)(BO)2- and B(n)(BO)2 (n = 5-12). <i>Journal of Chemical Physics</i> , <b>2013</b> , 139, 174301	3.9	35
324	Resonant photoelectron spectroscopy of Au2(-) via a Feshbach state using high-resolution photoelectron imaging. <i>Journal of Chemical Physics</i> , <b>2013</b> , 139, 194306	3.9	10
323	Observation of linear to planar structural transition in sulfur-doped gold clusters: $Au(x)S-(x = 2-5)$ . <i>Journal of Chemical Physics</i> , <b>2013</b> , 138, 174303	3.9	21
322	On the way to the highest coordination number in the planar metal-centred aromatic Ta $\Box$ B10-cluster: evolution of the structures of TaB(n)- (n = 3-8). <i>Journal of Chemical Physics</i> , <b>2013</b> , 139, 104312	3.9	48
321	Probing the electronic structures of low oxidation-state uranium fluoride molecules UF(x)- ( $x = 2-4$ ). <i>Journal of Chemical Physics</i> , <b>2013</b> , 139, 244303	3.9	13
320	Photoelectron spectroscopy of boron-gold alloy clusters and boron boronyl clusters: B3Au(n)(-) and B3(BO)n(-) (n = 1, 2). <i>Journal of Chemical Physics</i> , <b>2013</b> , 139, 044308	3.9	26
319	A combined photoelectron spectroscopy and ab initio study of the quasi-planar B24(-) cluster. <i>Journal of Chemical Physics</i> , <b>2013</b> , 139, 144307	3.9	111

318	Observation of Mode-Specific Vibrational Autodetachment from Dipole-Bound States of Cold Anions. <i>Angewandte Chemie</i> , <b>2013</b> , 125, 9146-9149	3.6	7
317	Observation of mode-specific vibrational autodetachment from dipole-bound states of cold anions. <i>Angewandte Chemie - International Edition</i> , <b>2013</b> , 52, 8976-9	16.4	84
316	Communication: Vibrational spectroscopy of Au4 from high resolution photoelectron imaging. Journal of Chemical Physics, <b>2013</b> , 139, 021106	3.9	36
315	Observation of the highest coordination number in planar species: decacoordinated Ta[] B10(-) and Nb[] B10(-) anions. <i>Angewandte Chemie - International Edition</i> , <b>2012</b> , 51, 2101-5	16.4	160
314	The electronic structure and chemical bonding in gold dihydride: AuH2[and AuH2. <i>Chemical Science</i> , <b>2012</b> , 3, 3286	9.4	47
313	On the electronic structure and conflicting d-orbital aromaticity in the Re3O3læluster. <i>RSC Advances</i> , <b>2012</b> , 2, 2707	3.7	15
312	Observation and investigation of the uranyl tetrafluoride dianion (UO2F42Dand its solvation complexes with water and acetonitrile. <i>Chemical Science</i> , <b>2012</b> , 3, 1137	9.4	41
311	Geometrical requirements for transition-metal-centered aromatic boron wheels: the case of VB10(-). <i>Physical Chemistry Chemical Physics</i> , <b>2012</b> , 14, 13663-9	3.6	43
310	Structural and electronic properties of reduced transition metal oxide clusters, M4O10 and M4O10-(M = Cr, W), from photoelectron spectroscopy and quantum chemical calculations. <i>Journal of Physical Chemistry A</i> , <b>2012</b> , 116, 5256-71	2.8	33
309	A photoelectron spectroscopy and ab initio study of B21-: negatively charged boron clusters continue to be planar at 21. <i>Journal of Chemical Physics</i> , <b>2012</b> , 136, 104310	3.9	112
308	Probing the structures of neutral boron clusters using infrared/vacuum ultraviolet two color ionization: B11, B16, and B17. <i>Journal of Chemical Physics</i> , <b>2012</b> , 137, 014317	3.9	94
307	Photoelectron Spectroscopy of Palladium(I) Dimers with Bridging Allyl Ligands. <i>Organometallics</i> , <b>2012</b> , 31, 8571-8576	3.8	4
306	Probing the structures and chemical bonding of boron-boronyl clusters using photoelectron spectroscopy and computational chemistry: B4(BO)(n)- (n = 1-3). <i>Journal of Chemical Physics</i> , <b>2012</b> , 137, 044307	3.9	23
305	Elongation of planar boron clusters by hydrogenation: boron analogues of polyenes. <i>Journal of the American Chemical Society</i> , <b>2012</b> , 134, 13228-31	16.4	67
304	Experimental and computational evidence of octa- and nona-coordinated planar iron-doped boron clusters: Fe[] B8[and Fe[] B9[]Journal of Organometallic Chemistry, 2012, 721-722, 148-154	2.3	68
303	Transition-metal-centered nine-membered boron rings: M(c)B9 and M(c)B9(-) (M = Rh, Ir). <i>Journal of the American Chemical Society</i> , <b>2012</b> , 134, 165-8	16.4	132
302	B22- and B23-: all-boron analogues of anthracene and phenanthrene. <i>Journal of the American Chemical Society</i> , <b>2012</b> , 134, 18065-73	16.4	172
301	Unraveling the mechanisms of O2 activation by size-selected gold clusters: transition from superoxo to peroxo chemisorption. <i>Journal of the American Chemical Society</i> , <b>2012</b> , 134, 9438-45	16.4	129

300	Probing the electronic properties and structural evolution of anionic gold clusters in the gas phase. <i>Nanoscale</i> , <b>2012</b> , 4, 4038-53	7.7	88
299	Probing the electronic structure and chemical bonding of the "staple" motifs of thiolate gold nanoparticles: Au(SCH3)2- and Au2(SCH3)3 <i>Physical Chemistry Chemical Physics</i> , <b>2012</b> , 14, 9323-9	3.6	39
298	Observation of the Highest Coordination Number in Planar Species: Decacoordinated Tall B10land Nbll B10lanions. <i>Angewandte Chemie</i> , <b>2012</b> , 124, 2143-2147	3.6	34
297	Note: Photoelectron spectroscopy of cold UF5(-). <i>Journal of Chemical Physics</i> , <b>2012</b> , 137, 116101	3.9	17
296	Resonant tunneling through the repulsive Coulomb barrier of a quadruply charged molecular anion. <i>Physical Review A</i> , <b>2012</b> , 85,	2.6	23
295	Photoelectron spectroscopy and the electronic structure of the uranyl tetrachloride dianion: UO2C(4(2-). <i>Journal of Chemical Physics</i> , <b>2012</b> , 137, 064315	3.9	41
294	Photoelectron spectroscopy and theoretical studies of UF5(-) and UF6(-). <i>Journal of Chemical Physics</i> , <b>2012</b> , 136, 194304	3.9	18
293	Photoelectron spectroscopy and ab initio study of boron-carbon mixed clusters: CB9- and C2B8 <i>Journal of Chemical Physics</i> , <b>2012</b> , 137, 234306	3.9	18
292	Stoichiometric and oxygen-rich M2O(n)- and M2O(n) (M = Nb, Ta; n = 5-7) clusters: molecular models for oxygen radicals, diradicals, and superoxides. <i>Journal of the American Chemical Society</i> , <b>2011</b> , 133, 3085-94	16.4	46
291	All-boron analogues of aromatic hydrocarbons: B17- and B18 <i>Journal of Chemical Physics</i> , <b>2011</b> , 134, 224304	3.9	242
290	Valence isoelectronic substitution in the B8(-) and B9(-) molecular wheels by an Al dopant atom: umbrella-like structures of AlB7(-) and AlB8(-). <i>Journal of Chemical Physics</i> , <b>2011</b> , 135, 104301	3.9	62
289	Molecular wheel to monocyclic ring transition in boron-carbon mixed clusters C2B6? and C3B5?. <i>Physical Chemistry Chemical Physics</i> , <b>2011</b> , 13, 8805-10	3.6	31
288	On the electronic structure of mono-rhenium oxide clusters: ReOn- and ReOn (n = 3, 4). Chemical Physics Letters, <b>2011</b> , 512, 49-53	2.5	14
287	Bridging 🛘 -BO in B2(BO)3(-) and B3(BO)3(-) clusters: boronyl analogs of boranes. <i>ChemPhysChem</i> , <b>2011</b> , 12, 2549-53	3.2	34
286	Aromatic Metal-Centered Monocyclic Boron Rings: Co[] B8[and Ru[] B9[]Angewandte Chemie, <b>2011</b> , 123, 9506-9509	3.6	27
285	Aromatic metal-centered monocyclic boron rings: Co B8- and Ru B9 <i>Angewandte Chemie - International Edition</i> , <b>2011</b> , 50, 9334-7	16.4	151
284	The mixed cyanide halide Au(I) complexes, [XAuCN][[X = F, Cl, Br, and I): evolution from ionic to covalent bonding. <i>Chemical Science</i> , <b>2011</b> , 2, 2101	9.4	37
283	Aluminum avoids the central position in AlB9- and AlB10-: photoelectron spectroscopy and ab initio study. <i>Journal of Physical Chemistry A</i> , <b>2011</b> , 115, 10391-7	2.8	40

282	Chemisorption-Induced 2DBDØD Structural Transitions in Gold Heptamer: (CO)nAu7[[n = 1]]). Journal of Physical Chemistry Letters, <b>2011</b> , 2, 2288-2293	6.4	26
281	Planarization of B7- and B12- clusters by isoelectronic substitution: AlB6- and AlB11 <i>Journal of the American Chemical Society</i> , <b>2011</b> , 133, 8646-53	16.4	66
280	Structure evolution of gold cluster anions between the planar and cage structures by isoelectronic substitution: Au(n)- (n = 13-15) and MAu(n)- (n = 12-14; M = Ag, Cu). <i>Journal of Chemical Physics</i> , <b>2011</b> , 134, 054306	3.9	44
279	A concentric planar doubly Earomatic BII cluster. <i>Nature Chemistry</i> , <b>2010</b> , 2, 202-6	17.6	424
278	Guiding electron emissions by excess negative charges in multiply charged molecular anions. <i>Physical Review Letters</i> , <b>2010</b> , 105, 263001	7.4	13
277	Isomer identification and resolution in small gold clusters. <i>Journal of Chemical Physics</i> , <b>2010</b> , 132, 05430	<b>15</b> .9	78
276	On the analogy of B-BO and B-Au chemical bonding in B11O- and B10Au- clusters. <i>Journal of Physical Chemistry A</i> , <b>2010</b> , 114, 12155-61	2.8	69
275	Covalent gold. Physical Chemistry Chemical Physics, <b>2010</b> , 12, 8694-705	3.6	94
274	On the electronic and structural properties of tri-niobium oxide clusters Nb3O(n)- (n = 3-8): photoelectron spectroscopy and density functional calculations. <i>Journal of Physical Chemistry A</i> , <b>2010</b> , 114, 5958-66	2.8	39
273	Probing the interactions of O(2) with small gold cluster anions (Au(n)(-), $n = 1-7$ ): chemisorption vs physisorption. <i>Journal of the American Chemical Society</i> , <b>2010</b> , 132, 4344-51	16.4	112
272	Photoelectron spectroscopy of C60Fn- and C60Fm2- (n = 17, 33, 35, 43, 45, 47; m = 34, 46) in the gas phase and the generation and characterization of C1-C60F47- and D2-C60F44 in solution. <i>Journal of Physical Chemistry A</i> , <b>2010</b> , 114, 1756-65	2.8	15
271	Vibrationally resolved photoelectron spectroscopy of di-gold carbonyl clusters Au2(CO)n- (n = 1-3): experiment and theory. <i>Journal of Physical Chemistry A</i> , <b>2010</b> , 114, 1247-54	2.8	45
270	Probing the structural evolution of medium-sized gold clusters: Au(n)(-) (n = 27-35). <i>Journal of the American Chemical Society</i> , <b>2010</b> , 132, 6596-605	16.4	111
269	Photoelectron imaging of doubly charged anions, (-)O2C(CH2)nCO2(-) (n = 2-8): observation of near 0 eV electrons due to secondary dissociative autodetachment. <i>Journal of Physical Chemistry A</i> , <b>2010</b> , 114, 4524-30	2.8	16
268	Photoelectron imaging and spectroscopy of MI(2)(-) (M = Cs, Cu, Au): evolution from ionic to covalent bonding. <i>Journal of Physical Chemistry A</i> , <b>2010</b> , 114, 11244-51	2.8	50
267	Planar to linear structural transition in small boron-carbon mixed clusters: $C(x)B(5-x)-(x = 1-5)$ .  Journal of the American Chemical Society, <b>2010</b> , 132, 14104-12	16.4	33
266	Observation of earlier two-to-three dimensional structural transition in gold cluster anions by isoelectronic substitution: MAu(n)(-) (n=8-11; M=Ag,Cu). <i>Journal of Chemical Physics</i> , <b>2010</b> , 132, 114306	3.9	72
265	Stepwise hydration of the cyanide anion: a temperature-controlled photoelectron spectroscopy and ab initio computational study of CN-(H2O)n, n=2-5. <i>Journal of Chemical Physics</i> , <b>2010</b> , 132, 124306	3.9	23

264	Probing the electronic structure of early transition metal oxide clusters: Molecular models towards mechanistic insights into oxide surfaces and catalysis. <i>Chemical Physics Letters</i> , <b>2010</b> , 500, 185-195	2.5	90
263	Magnetic doping of the golden cage cluster M@Au16[[M=Fe,Co,Ni). <i>Physical Review B</i> , <b>2009</b> , 79,	3.3	82
262	Probing the 2D to 3D structural transition in gold cluster anions using argon tagging. <i>Physical Review Letters</i> , <b>2009</b> , 102, 153401	7.4	145
261	Experimental and theoretical investigation of three-dimensional nitrogen-doped aluminum clusters Al8N- and Al8N. <i>Journal of Chemical Physics</i> , <b>2009</b> , 130, 134303	3.9	11
260	Tuning the electronic properties of the golden buckyball by endohedral doping: M@Au16(-) (M = Ag,Zn,In). <i>Journal of Chemical Physics</i> , <b>2009</b> , 130, 051101	3.9	61
259	The [(AL2O3)2]- anion cluster: electron localization-delocalization isomerism. <i>ChemPhysChem</i> , <b>2009</b> , 10, 2410-3	3.2	25
258	Diversity of Functionalized Germanium Zintl Clusters: Syntheses and Theoretical Studies of [Ge9PdPPh3]3[and [Ni@(Ge9PdPPh3)]2[] Journal of Cluster Science, <b>2009</b> , 20, 601-609	3	29
257	Microsolvation of the acetate anion [CH3CO2-(H2O)n, n= 1B]: A photoelectron spectroscopy and ab initio computational study. <i>Chemical Physics Letters</i> , <b>2009</b> , 477, 41-44	2.5	14
256	Photoelectron spectroscopy of cold hydrated sulfate clusters, SO4(2-)(H2O)n (n = 4-7): temperature-dependent isomer populations. <i>Journal of Physical Chemistry A</i> , <b>2009</b> , 113, 5567-76	2.8	43
255	Are carboxyl groups the most acidic sites in amino acids? Gas-phase acidities, photoelectron spectra, and computations on tyrosine, p-hydroxybenzoic acid, and their conjugate bases. <i>Journal of the American Chemical Society</i> , <b>2009</b> , 131, 1174-81	16.4	62
254	Structural and electronic properties of reduced transition metal oxide clusters, M3O8 and M3O8-(M = Cr, W), from photoelectron spectroscopy and quantum chemical calculations. <i>Journal of Physical Chemistry A</i> , <b>2009</b> , 113, 11273-88	2.8	55
253	Structural transition of gold nanoclusters: from the golden cage to the golden pyramid. <i>ACS Nano</i> , <b>2009</b> , 3, 1225-30	16.7	99
252	Photoelectron angular distribution and molecular structure in multiply charged anions. <i>Journal of Physical Chemistry A</i> , <b>2009</b> , 113, 945-8	2.8	11
251	Structural evolution, sequential oxidation, and chemical bonding in tritantalum oxide clusters: $Ta(3)O(n)(-)$ and $Ta(3)O(n)$ (n = 1-8). <i>Journal of Physical Chemistry A</i> , <b>2009</b> , 113, 9804-13	2.8	44
250	Observation of a remarkable temperature effect in the hydrogen bonding structure and dynamics of the CN(-)(H2O) cluster. <i>Journal of Physical Chemistry A</i> , <b>2009</b> , 113, 9579-84	2.8	10
249	Evidence of significant covalent bonding in Au(CN)(2)(-). <i>Journal of the American Chemical Society</i> , <b>2009</b> , 131, 16368-70	16.4	137
248	Detecting weak interactions between Au- and gas molecules: a photoelectron spectroscopic and ab initio study. <i>Journal of the American Chemical Society</i> , <b>2009</b> , 131, 9484-5	16.4	41
247	Structural evolution of doped gold clusters: $MAu(x)(-)$ (M = Si, Ge, Sn; x = 5-8). Journal of the American Chemical Society, <b>2009</b> , 131, 3396-404	16.4	87

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246	Investigating the weak to evaluate the strong: an experimental determination of the electron binding energy of carborane anions and the gas phase acidity of carborane acids. <i>Journal of the American Chemical Society</i> , <b>2009</b> , 131, 18050-1	16.4	29
245	Au10-: isomerism and structure-dependent O2 reactivity. <i>Physical Chemistry Chemical Physics</i> , <b>2009</b> , 11, 2663-7	3.6	52
244	Probing the electronic and structural properties of the niobium trimer cluster and its mono- and dioxides: Nb3On- and Nb3On (n = 0-2). <i>Journal of Physical Chemistry A</i> , <b>2009</b> , 113, 3866-75	2.8	52
243	Photoelectron spectroscopy of multiply charged anions. <i>Annual Review of Physical Chemistry</i> , <b>2009</b> , 60, 105-26	15.7	108
242	Probing the electronic stability of multiply charged anions: sulfonated pyrene tri- and tetraanions. Journal of the American Chemical Society, <b>2009</b> , 131, 9836-42	16.4	14
241	Experimental and theoretical investigations of CB8-: towards rational design of hypercoordinated planar chemical species. <i>Physical Chemistry Chemical Physics</i> , <b>2009</b> , 11, 9840-9	3.6	80
240	CO chemisorption on the surfaces of the golden cages. <i>Journal of Chemical Physics</i> , <b>2009</b> , 131, 234305	3.9	30
239	Relativistic effects and the unique low-symmetry structures of gold nanoclusters. <i>ACS Nano</i> , <b>2008</b> , 2, 897-904	16.7	119
238	A photoelectron spectroscopic and theoretical study of B16- and B16(2-): an all-boron naphthalene. <i>Journal of the American Chemical Society</i> , <b>2008</b> , 130, 7244-6	16.4	231
237	Probing the electronic and structural properties of chromium oxide clusters (CrO3)n(-) and (CrO3)n (n = 1-5): photoelectron spectroscopy and density functional calculations. <i>Journal of the American Chemical Society</i> , <b>2008</b> , 130, 5167-77	16.4	92
236	Carbon avoids hypercoordination in CB6(-), CB6(2-), and C2B5(-) planar carbon-boron clusters. <i>Journal of the American Chemical Society</i> , <b>2008</b> , 130, 9248-50	16.4	138
235	B2(BO)2(2-)-diboronyl diborene: a linear molecule with a triple boron-boron bond. <i>Journal of the American Chemical Society</i> , <b>2008</b> , 130, 2573-9	16.4	142
234	Probing the electronic and structural properties of doped aluminum clusters: MAl12- (M=Li, Cu, and Au). <i>Journal of Chemical Physics</i> , <b>2008</b> , 128, 024305	3.9	42
233	High resolution and low-temperature photoelectron spectroscopy of an oxygen-linked fullerene dimer dianion: C(120)O(2-). <i>Journal of Chemical Physics</i> , <b>2008</b> , 128, 114307	3.9	14
232	Aromaticity and antiaromaticity in transition-metal systems. <i>Physical Chemistry Chemical Physics</i> , <b>2008</b> , 10, 257-67	3.6	166
231	Photoelectron spectroscopy and Ab initio study of the structure and bonding of Al7N- and Al7N. <i>Journal of Physical Chemistry A</i> , <b>2008</b> , 112, 1873-9	2.8	25
230	On the electronic structure and chemical bonding in the tantalum trimer cluster. <i>Journal of Physical Chemistry A</i> , <b>2008</b> , 112, 10962-7	2.8	43
229	Observation of entropic effect on conformation changes of complex systems under well-controlled temperature conditions. <i>Journal of Physical Chemistry A</i> , <b>2008</b> , 112, 172-5	2.8	27

228	Probing the electronic structure and chemical bonding of gold oxides and sulfides in AuOn(-) and AuSn(-) (n = 1, 2). <i>Journal of the American Chemical Society</i> , <b>2008</b> , 130, 9156-67	16.4	67
227	Observation of H2 aggregation onto a doubly charged anion in a temperature-controlled ion trap. <i>Journal of Physical Chemistry A</i> , <b>2008</b> , 112, 13271-4	2.8	22
226	Chemisorption-induced Structural Changes and Transition from Chemisorption to Physisorption in Au6(CO)n $\mathbb{I}$ (n = 4 $\mathbb{B}$ ). <i>Journal of Physical Chemistry C</i> , <b>2008</b> , 112, 11920-11928	3.8	49
225	Stable icosahedral hollow cage clusters: stannaspherene () and plumbaspherene (). <i>International Reviews in Physical Chemistry</i> , <b>2008</b> , 27, 139-166	7	22
224	Low-lying isomers of the B9(-) boron cluster: the planar molecular wheel versus three-dimensional structures. <i>Journal of Chemical Physics</i> , <b>2008</b> , 129, 024302	3.9	71
223	Development of a low-temperature photoelectron spectroscopy instrument using an electrospray ion source and a cryogenically controlled ion trap. <i>Review of Scientific Instruments</i> , <b>2008</b> , 79, 073108	1.7	241
222	Photoelectron spectroscopy of anions at 118.2 nm: observation of high electron binding energies in superhalogens MCl4- (M=Sc, Y, La). <i>Journal of Chemical Physics</i> , <b>2008</b> , 128, 201102	3.9	72
221	Boronyls as key structural units in boron oxide clusters: B(BO)2- and B(BO)3 <i>Journal of the American Chemical Society</i> , <b>2007</b> , 129, 9254-5	16.4	91
220	On the chemical bonding of gold in auro-boron oxide clusters AunBO- (n = 1-3). <i>Journal of Physical Chemistry A</i> , <b>2007</b> , 111, 1648-58	2.8	41
219	Pd(2)@Sn(18)(4-): fusion of two endohedral stannaspherenes. <i>Journal of the American Chemical Society</i> , <b>2007</b> , 129, 9560-1	16.4	109
218	Microsolvation of the dicyanamide anion: $[N(CN)(2)(-)](H(2)O)n$ (n = 0-12). Journal of Physical Chemistry A, <b>2007</b> , 111, 7719-25	2.8	17
217	Photoelectron Spectroscopy of Singly and Doubly Charged Higher Fullerenes at Low Temperatures: C76-, C78-, C84- and C762-, C782-, C842- <i>Journal of Physical Chemistry C</i> , <b>2007</b> , 111, 17	684 <sup>8</sup> 17	68 <sup>8</sup>
216	Probing the electronic structure of early transition-metal oxide clusters: polyhedral cages of $(V2O5)n(-)$ (n = 2-4) and $(M2O5(2)(-)$ (M = Nb, Ta). <i>Journal of the American Chemical Society</i> , <b>2007</b> , 129, 13270-6	16.4	105
215	Probing the structure and bonding in Al6N- and Al6N by photoelectron spectroscopy and ab initio calculations. <i>Journal of Physical Chemistry A</i> , <b>2007</b> , 111, 34-41	2.8	15
214	Structural Transitions from Pyramidal to Fused Planar to Tubular to Core/Shell Compact in Gold Clusters: Aun- (n = 21🗹5). <i>Journal of Physical Chemistry C</i> , <b>2007</b> , 111, 4190-4198	3.8	85
213	A photoelectron spectroscopic and computational study of sodium auride clusters, NanAun- (n = 1-3). <i>Journal of Physical Chemistry A</i> , <b>2007</b> , 111, 7555-61	2.8	15
212	Doping golden buckyballs:Cu@Au16- and Cu@Au17- cluster anions. <i>Angewandte Chemie - International Edition</i> , <b>2007</b> , 46, 2915-8	16.4	106
211	CB7-: experimental and theoretical evidence against hypercoordinate planar carbon. <i>Angewandte Chemie - International Edition</i> , <b>2007</b> , 46, 4550-3	16.4	108

# (2006-2007)

210	Electrospray ionization photoelectron spectroscopy: Probing the electronic structure of inorganic metal complexes in the gas-phase. <i>Coordination Chemistry Reviews</i> , <b>2007</b> , 251, 474-491	23.2	32
209	Electronic structure and fragmentation properties of [Fe4S4(SEt)4½(SSEt)x]2[International Journal of Mass Spectrometry, <b>2007</b> , 263, 260-266	1.9	6
208	Vibrationally resolved photoelectron spectroscopy of BO- and BO2-: a joint experimental and theoretical study. <i>Journal of Physical Chemistry A</i> , <b>2007</b> , 111, 1030-5	2.8	151
207	Doping the golden cage Au16(-) with Si, Ge, and Sn. <i>Journal of the American Chemical Society</i> , <b>2007</b> , 129, 15136-7	16.4	87
206	Probing the electronic structure and band gap evolution of titanium oxide clusters ( $TiO(2)$ )(n)(-) (n = 1-10) using photoelectron spectroscopy. <i>Journal of the American Chemical Society</i> , <b>2007</b> , 129, 3022-6	16.4	159
205	Au34-: A Fluxional CoreBhell Cluster. <i>Journal of Physical Chemistry C</i> , <b>2007</b> , 111, 8228-8232	3.8	98
204	Formation of monodisperse (WO3)3 clusters on TiO2(110). <i>Angewandte Chemie - International Edition</i> , <b>2006</b> , 45, 4786-9	16.4	88
203	Planar nitrogen-doped aluminum clusters AlxN- (x=3-5). <i>Journal of Chemical Physics</i> , <b>2006</b> , 125, 124305	3.9	30
202	Probing the electronic properties of dichromium oxide clusters Cr2On- (n=1-7) using photoelectron spectroscopy. <i>Journal of Chemical Physics</i> , <b>2006</b> , 125, 164315	3.9	42
201	Evidence of hollow golden cages. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , <b>2006</b> , 103, 8326-30	11.5	345
200	On the structure and chemical bonding of Si6(2-) and Si6(2-) in NaSi6(-) upon Na+ coordination. Journal of Chemical Physics, <b>2006</b> , 124, 124305	3.9	28
199	Direct experimental probe of the on-site Coulomb repulsion in the doubly charged fullerene anion C70 2 <i>Physical Review Letters</i> , <b>2006</b> , 96, 143002	7.4	66
198	First steps towards dissolution of NaSO4- by water. <i>Physical Chemistry Chemical Physics</i> , <b>2006</b> , 8, 4294-6	3.6	19
197	Gold apes hydrogen. The structure and bonding in the planar B7Au2- and B7Au2 clusters. <i>Journal of Physical Chemistry A</i> , <b>2006</b> , 110, 1689-93	2.8	108
196	Photoelectron spectroscopy of the bis(dithiolene) anions [M(mnt)2]n- (M = Fe - Zn; n = 1, 2): changes in electronic structure with variation of metal center and with oxidation. <i>Inorganic Chemistry</i> , <b>2006</b> , 45, 5841-51	5.1	16
195	Probing the Electronic Structure of Fe?S Clusters: Ubiquitous Electron Transfer Centers in Metalloproteins Using Anion Photoelectron Spectroscopy in the Gas Phase <b>2006</b> , 63-117		3
194	Sn12(2-): stannaspherene. <i>Journal of the American Chemical Society</i> , <b>2006</b> , 128, 8390-1	16.4	140
193	Gold as hydrogen: structural and electronic properties and chemical bonding in Si3Au3(+0-) and comparisons to Si3H3(+0-). <i>Journal of Chemical Physics</i> , <b>2006</b> , 125, 133204	3.9	71

Facile syntheses of monodisperse ultrasmall Au clusters. Journal of Physical Chemistry B, 2006, 110, 2141628 192 124 Pb12 2-: plumbaspherene. Journal of Physical Chemistry A, 2006, 110, 10169-72 2.8 191 105 Determination of the electron affinity of the acetyloxyl radical (CH3COO) by low-temperature anion photoelectron spectroscopy and ab initio calculations. Journal of Physical Chemistry A, 2006, 2.8 190 47 110, 5047-50 Structural evolution of anionic silicon clusters SiN (20. Journal of Physical Chemistry A, 2006, 110, 908-122.8 189 69 Probing the intrinsic electronic structure of the bis(dithiolene) anions [M(mnt)2]2- and [M(mnt)2]1-188 (M = Ni, Pd, Pt; mnt = 1,2-S2C2(CN)2) in the gas phase by photoelectron spectroscopy. Journal of 16.4 28 the American Chemical Society, 2006, 128, 4282-91 Experimental and computational studies of alkali-metal coinage-metal clusters. Journal of Physical 2.8 187 67 Chemistry A, 2006, 110, 4244-50 Observation of cysteine thiolate and -S...H-O intermolecular hydrogen bond. Journal of Physical 186 2.8 54 Chemistry A, 2006, 110, 12603-6 Observation of triatomic species with conflicting aromaticity: AlSi2- and AlGe2-. Journal of Physical 185 10 3.4 Chemistry B, 2006, 110, 9743-6 Photoelectron spectroscopy of free multiply charged Keggin anions alpha-[PM12O40]3- (M = Mo, 2.8 184 26 W) in the gas phase. Journal of Physical Chemistry A, 2006, 110, 10737-41 Low-temperature photoelectron spectroscopy of aliphatic dicarboxylate monoanions, 183 HO2C(CH2(nCO2- (n = 1-10)): hydrogen bond induced cyclization and strain energies. Journal of 2.8 12 Physical Chemistry A, 2006, 110, 7801-5 Theoretical probing of deltahedral closo-auroboranes B(x)Au(x)2-(x = 5-12). Inorganic Chemistry, 182 5.1 28 2006, 45, 5269-71 On the structure and chemical bonding of tri-tungsten oxide clusters W3On- and W3On (n=7-10): W3O8 as a potential molecular model for O-deficient defect sites in tungsten oxides. Journal of 181 2.8 75 Physical Chemistry A, 2006, 110, 85-92 All-boron aromatic clusters as potential new inorganic ligands and building blocks in chemistry. 180 23.2 509 Coordination Chemistry Reviews, 2006, 250, 2811-2866 Free tetra- and hexa-coordinated platinum-cyanide dianions, and: A combined photodetachment 179 2.3 photoelectron spectroscopic and theoretical study. Chemical Physics, 2006, 329, 230-238 MX3(-) superhalogens (M = Be, Mg, Ca; X = Cl, Br): a photoelectron spectroscopic and ab initio 178 2.8 142 theoretical study. Journal of Physical Chemistry A, 2005, 109, 11560-7 Photoelectron spectroscopy of doubly and singly charged group VIB dimetalate anions: M2O72-, 2.8 70 MM'O72-, and M2O7- (M, M' = Cr, Mo, W). Journal of Physical Chemistry A, 2005, 109, 10512-20 Electronic structure of the hydroxo and methoxo oxometalate anions MO3(OH)- and MO3(OCH3)-2.8 176 15 (M = Cr, Mo, and W). Journal of Physical Chemistry A, 2005, 109, 11771-80 Cu3C4-: a new sandwich molecule with two revolving C2(2-) units. Journal of Physical Chemistry A, 2.8 27 **2005**, 109, 562-70

#### (2005-2005)

174	Interior and interfacial aqueous solvation of benzene dicarboxylate dianions and their methylated analogues: A combined molecular dynamics and photoelectron spectroscopy study. <i>Journal of Physical Chemistry A</i> , <b>2005</b> , 109, 5042-9	2.8	20
173	Temperature-dependent photoelectron spectroscopy of methyl benzoate anions: observation of steric effect in o-methyl benzoate. <i>Journal of Physical Chemistry A</i> , <b>2005</b> , 109, 11395-400	2.8	20
172	Electronic structure and intrinsic redox properties of [2Fe-2S]+ clusters with tri- and tetracoordinate iron sites. <i>Inorganic Chemistry</i> , <b>2005</b> , 44, 1202-4	5.1	14
171	Gold as hydrogen. An experimental and theoretical study of the structures and bonding in disilicon gold clusters $Si2Au(n)$ - and $Si2Au(n)$ ( $n = 2$ and 4) and comparisons to $Si2H2$ and $Si2H4$ . <i>Journal of Physical Chemistry A</i> , <b>2005</b> , 109, 4366-74	2.8	96
170	Electronic and structural evolution and chemical bonding in ditungsten oxide clusters: W2O(n)- and W2O(n) (n = 1-6). <i>Journal of Physical Chemistry A</i> , <b>2005</b> , 109, 6019-30	2.8	64
169	Unique CO chemisorption properties of gold hexamer: Au6(CO)n- (n = 0-3). <i>Journal of the American Chemical Society</i> , <b>2005</b> , 127, 12098-106	16.4	96
168	Probing the electronic structure of [2Fe-2S] clusters with three coordinate iron sites by use of photoelectron spectroscopy. <i>Journal of Physical Chemistry A</i> , <b>2005</b> , 109, 1815-20	2.8	14
167	Photoelectron spectroscopy and electronic structures of fullerene oxides: C60Ox- ( $x = 1-3$ ). <i>Journal of Physical Chemistry A</i> , <b>2005</b> , 109, 11089-92	2.8	7
166	All-metal aromaticity and antiaromaticity. Chemical Reviews, 2005, 105, 3716-57	68.1	457
165	Probing the low-barrier hydrogen bond in hydrogen maleate in the gas phase: a photoelectron spectroscopy and ab initio study. <i>Journal of Physical Chemistry A</i> , <b>2005</b> , 109, 10633-7	2.8	34
164	Chemical bonding in Si5(2-) and NaSi5(-) via photoelectron spectroscopy and ab initio calculations. Journal of Physical Chemistry A, <b>2005</b> , 109, 11385-94	2.8	26
163	Synthesis of the H-cluster framework of iron-only hydrogenase. <i>Nature</i> , <b>2005</b> , 433, 610-3	50.4	467
162	Observation of weak C-HO hydrogen bonding to unactivated alkanes. <i>Angewandte Chemie - International Edition</i> , <b>2005</b> , 44, 4968-72	16.4	46
161	Experimental and theoretical investigation of the electronic and geometrical structures of the Au32 cluster. <i>Angewandte Chemie - International Edition</i> , <b>2005</b> , 44, 7119-23	16.4	124
160	Chemisorption sites of CO on small gold clusters and transitions from chemisorption to physisorption. <i>Journal of Chemical Physics</i> , <b>2005</b> , 122, 51101	3.9	87
159	Photoelectron spectroscopy and ab initio study of the doubly antiaromatic B(6) (2-) dianion in the LiB(6) (-) cluster. <i>Journal of Chemical Physics</i> , <b>2005</b> , 122, 54313	3.9	92
158	Magnetic properties in transition-metal-doped gold clusters: M@Au6 (M = Ti, V, Cr). <i>Physical Review Letters</i> , <b>2005</b> , 95, 253401	7.4	149
157	Planar-to-tubular structural transition in boron clusters: B20 as the embryo of single-walled boron nanotubes. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , <b>2005</b> , 102, 961-4	11.5	428

156	Observation of Au2H- impurity in pure gold clusters and implications for the anomalous Au-Au distances in gold nanowires. <i>Journal of Chemical Physics</i> , <b>2004</b> , 121, 8231-6	3.9	43
155	Direct experimental observation of the low ionization potentials of guanine in free oligonucleotides by using photoelectron spectroscopy. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , <b>2004</b> , 101, 17588-92	11.5	124
154	[SiAu4]: Aurosilane. Angewandte Chemie - International Edition, 2004, 43, 2125-9	16.4	120
153	Multiple aromaticity and antiaromaticity in silicon clusters. ChemPhysChem, 2004, 5, 1885-91	3.2	46
152	Terminal ligand influence on the electronic structure and intrinsic redox properties of the [Fe4S4]2+ cubane clusters. <i>Inorganic Chemistry</i> , <b>2004</b> , 43, 3647-55	5.1	15
151	Electronic structure and chemical bonding in MO(n)- and MO(n) clusters (M = Mo, W; n = 3-5): a photoelectron spectroscopy and ab initio study. <i>Journal of the American Chemical Society</i> , <b>2004</b> , 126, 16134-41	16.4	104
150	Mechanistic Insight into the Symmetric Fission of [4Fe\( \Pi \)S] Analogue Complexes and Implications for Cluster Conversions in Iron\( \Pi \) ulfur Proteins. Journal of Physical Chemistry A, 2004, 108, 6750-6757	2.8	23
149	Solvent-mediated folding of a doubly charged anion. <i>Journal of the American Chemical Society</i> , <b>2004</b> , 126, 876-83	16.4	39
148	Solvation of the Azide Anion (N3-) in Water Clusters and Aqueous Interfaces: A Combined Investigation by Photoelectron Spectroscopy, Density Functional Calculations, and Molecular Dynamics Simulations <i>Journal of Physical Chemistry A</i> , <b>2004</b> , 108, 7820-7826	2.8	44
147	Bulk versus interfacial aqueous solvation of dicarboxylate dianions. <i>Journal of the American Chemical Society</i> , <b>2004</b> , 126, 11691-8	16.4	53
146	Direct measurement of the hydrogen-bonding effect on the intrinsic redox potentials of [4Fe-4S] cubane complexes. <i>Journal of the American Chemical Society</i> , <b>2004</b> , 126, 15790-4	16.4	46
145	Photoelectron spectroscopy of the doubly-charged anions [MIVO(mnt)2]2- (M = Mo, W; mnt = S2C2(CN)2(2-): access to the ground and excited states of the [MVO(mnt)2]- anion. <i>Journal of the American Chemical Society</i> , <b>2004</b> , 126, 5119-29	16.4	23
144	Sequential oxidation of the cubane [4Fe4S] cluster from [4Fe4S](-) to [4Fe4S](3+) in Fe(4)S(4)L(n)(-) complexes. <i>Journal of the American Chemical Society</i> , <b>2004</b> , 126, 8413-20	16.4	11
143	Electronic Structure, Isomerism, and Chemical Bonding in B7- and B7. <i>Journal of Physical Chemistry A</i> , <b>2004</b> , 108, 3509-3517	2.8	177
142	Molecular wheel B8(2-) as a new inorganic ligand. photoelectron spectroscopy and ab initio characterization of LiB8(-). <i>Inorganic Chemistry</i> , <b>2004</b> , 43, 3552-4	5.1	135
141	Toward the Solution Synthesis of the Tetrahedral Au20 Cluster. <i>Journal of Physical Chemistry B</i> , <b>2004</b> , 108, 12259-12263	3.4	96
140	Photoelectron Spectroscopy of Free Polyoxoanions Mo6O192- and W6O192- in the Gas Phase. Journal of Physical Chemistry A, <b>2004</b> , 108, 10089-10093	2.8	50
139	Competition between linear and cyclic structures in monochromium carbide clusters CrCn- and CrCn (n = 2-8): a photoelectron spectroscopy and density functional study. <i>Journal of Chemical Physics</i> , <b>2004</b> , 120, 8996-9008	3.9	57

138	Icosahedral gold cage clusters: M@Au12- (M=V, Nb, and Ta). <i>Journal of Chemical Physics</i> , <b>2004</b> , 121, 836	93794	127
137	Structure of the Na(x)Cl(x+1) (-) (x=1-4) clusters via ab initio genetic algorithm and photoelectron spectroscopy. <i>Journal of Chemical Physics</i> , <b>2004</b> , 121, 5709-19	3.9	247
136	From helical nanowires, nanocrosses to aligned micro-carbon fibers. <i>Materials Research Society Symposia Proceedings</i> , <b>2003</b> , 776, 721		
135	Au20: a tetrahedral cluster. <i>Science</i> , <b>2003</b> , 299, 864-7	33.3	990
134	Photoelectron spectroscopy of TinItlusters (n=11130). Journal of Chemical Physics, 2003, 118, 2108-2115	3.9	40
133	Hepta- and Octacoordinate Boron in Molecular Wheels of Eight- and Nine-Atom Boron Clusters: Observation and Confirmation. <i>Angewandte Chemie</i> , <b>2003</b> , 115, 6186-6190	3.6	65
132	Au20: A Tetrahedral Cluster <i>ChemInform</i> , <b>2003</b> , 34, no		3
131	Hepta- and octacoordinate boron in molecular wheels of eight- and nine-atom boron clusters: observation and confirmation. <i>Angewandte Chemie - International Edition</i> , <b>2003</b> , 42, 6004-8	16.4	419
130	Gold dichloride and gold dibromide with gold atoms in three different oxidation states. <i>Angewandte Chemie - International Edition</i> , <b>2003</b> , 42, 311-4	16.4	54
129	Collision-induced symmetric fission of doubly-charged cubelike [Fe4S4X4]2lælusters. <i>International Journal of Mass Spectrometry</i> , <b>2003</b> , 228, 797-805	1.9	19
128	Hydrocarbon analogues of boron clustersplanarity, aromaticity and antiaromaticity. <i>Nature Materials</i> , <b>2003</b> , 2, 827-33	27	567
127	Structure and Bonding in B6- and B6: Planarity and Antiaromaticity. <i>Journal of Physical Chemistry A</i> , <b>2003</b> , 107, 1359-1369	2.8	175
126	Photoelectron Spectroscopy and ab Initio Study of B3- and B4- Anions and Their Neutrals. <i>Journal of Physical Chemistry A</i> , <b>2003</b> , 107, 9319-9328	2.8	169
125	Probing the Electronic Structure of the Di-Iron Subsite of [Fe]-Hydrogenase: A Photoelectron Spectroscopic Study of Fe(I)He(I) Model Complexes. <i>Journal of Physical Chemistry A</i> , <b>2003</b> , 107, 4612-46	1 <b>3</b> .8	28
124	On the Electronic and Atomic Structures of Small AuN- (N = 4114) Clusters: A Photoelectron Spectroscopy and Density-Functional Study. <i>Journal of Physical Chemistry A</i> , <b>2003</b> , 107, 6168-6175	2.8	572
123	Photodetachment of zwitterions: probing intramolecular coulomb repulsion and attraction in the gas phase using pyridinium dicarboxylate anions. <i>Journal of the American Chemical Society</i> , <b>2003</b> , 125, 296-304	16.4	41
122	Photodetachment of zwitterions: probing intramolecular Coulomb repulsion and attraction in the gas phase using mono-decarboxylated pyridinium dicarboxylates. Implications on the mechanism of orotidine 5'-monophosphate decarboxylase. <i>Journal of the American Chemical Society</i> , <b>2003</b> , 125, 6814-2	16.4 2 <b>6</b>	12
121	Electronic and Structural Evolution of Monoiron Sulfur Clusters, FeSn- and FeSn (n = 1 <b>B</b> ), from Anion Photoelectron Spectroscopy. <i>Journal of Physical Chemistry A</i> , <b>2003</b> , 107, 2821-2828	2.8	38

120	Synthesis, Characterization, and Manipulation of Helical SiO2 Nanosprings. <i>Nano Letters</i> , <b>2003</b> , 3, 577-5	5 <b>80</b> 1.5	178
119	Probing the intrinsic electronic structure of the cubane [4Fe-4S] cluster: nature's favorite cluster for electron transfer and storage. <i>Journal of the American Chemical Society</i> , <b>2003</b> , 125, 14072-81	16.4	63
118	Combined Quantum Chemistry and Photoelectron Spectroscopy Study of the Electronic Structure and Reduction Potentials of Rubredoxin Redox Site Analogues. <i>Journal of Physical Chemistry A</i> , <b>2003</b> , 107, 2898-2907	2.8	38
117	On the Electronic Structure of [1Fe] Feß Complexes from Anionic Photoelectron Spectroscopy. Journal of Physical Chemistry A, <b>2003</b> , 107, 1703-1709	2.8	27
116	Structural and electronic properties of small titanium clusters: A density functional theory and anion photoelectron spectroscopy study. <i>Journal of Chemical Physics</i> , <b>2003</b> , 118, 2116-2123	3.9	73
115	All-metal antiaromatic molecule: rectangular Al4(4-) in the Li3Al4(-) anion. <i>Science</i> , <b>2003</b> , 300, 622-5	33.3	205
114	Structural and electronic properties of iron monoxide clusters FenO and FenO[(n=2f): A combined photoelectron spectroscopy and density functional theory study. <i>Journal of Chemical Physics</i> , <b>2003</b> , 119, 11135-11145	3.9	51
113	On the electronic structures of gaseous transition metal halide complexes, FeX4land MX3ll (M=Mn, Fe, Co, Ni, X=Cl, Br), using photoelectron spectroscopy and density functional calculations. <i>Journal of Chemical Physics</i> , <b>2003</b> , 119, 8311-8320	3.9	47
112	Photodetachment of hydrated oxalate dianions in the gas phase, C2O42[H2O)n (n=3월0): From solvated clusters to nanodroplet. <i>Journal of Chemical Physics</i> , <b>2003</b> , 119, 3631-3640	3.9	37
111	Experimental Observation and Confirmation of Icosahedral W@Au12 and Mo@Au12 Molecules. <i>Angewandte Chemie</i> , <b>2002</b> , 114, 4980-4983	3.6	49
110	Experimental observation and confirmation of icosahedral W@Au12 and Mo@Au12 molecules. <i>Angewandte Chemie - International Edition</i> , <b>2002</b> , 41, 4786-9	16.4	299
109	Photoelectron spectroscopy of pentaatomic tetracoordinate planar carbon molecules: CAl3Siland CAl3Gell <i>Chemical Physics Letters</i> , <b>2002</b> , 357, 415-419	2.5	51
108	Electronic structure and chemical boning in nonstoichiometric molecules: Al3X2[X=C,Si,Ge). A photoelectron spectroscopy and ab initio study. <i>Journal of Chemical Physics</i> , <b>2002</b> , 116, 1330-1338	3.9	29
107	Collision-induced dissociation and photodetachment of singly and doubly charged anionic polynuclear transition metal carbonyl clusters: Ru3Co(CO)13[Ru6C(CO)162[land Ru6(CO)182[l Journal of Chemical Physics, 2002, 116, 6560-6566	3.9	26
106	Electronic structure and chemical bonding of B5Dand B5 by photoelectron spectroscopy and ab initio calculations. <i>Journal of Chemical Physics</i> , <b>2002</b> , 117, 7917-7924	3.9	193
105	Electronic structure and chemical bonding of divanadium-oxide clusters (V2Ox, x=3🛭) from anion photoelectron spectroscopy. <i>Journal of Chemical Physics</i> , <b>2002</b> , 117, 7882-7888	3.9	75
104	Coulomb- and antiferromagnetic-induced fission in doubly charged cubelike fe-s clusters. <i>Physical Review Letters</i> , <b>2002</b> , 89, 163401	7.4	19
103	Evolution of the electronic properties of small Nin[n=1f]00) clusters by photoelectron spectroscopy. <i>Journal of Chemical Physics</i> , <b>2002</b> , 117, 9758-9765	3.9	28

# (2001-2002)

102	Probing solution-phase species and chemistry in the gas phase. <i>International Reviews in Physical Chemistry</i> , <b>2002</b> , 21, 473-498	7	67
101	Helical Crystalline SiC/SiO2 CoreBhell Nanowires. <i>Nano Letters</i> , <b>2002</b> , 2, 941-944	11.5	238
100	Photodetachment of Hydrated Sulfate Doubly Charged Anions: SO42-(H2O)n (n = 4個0)回 <i>Journal of Physical Chemistry A</i> , <b>2002</b> , 106, 7607-7616	2.8	69
99	Peculiar antiaromatic inorganic molecules of tetrapnictogen in Na+Pn4- (Pn = P, As, Sb) and important consequences for hydrocarbons. <i>Inorganic Chemistry</i> , <b>2002</b> , 41, 6062-70	5.1	64
98	In search of covalently bound tetra- and penta-oxygen species: a photoelectron spectroscopic and Ab initio investigation of MO4- and MO5- (M = Li, Na, K, Cs). <i>Journal of the American Chemical Society</i> , <b>2002</b> , 124, 6742-50	16.4	13
97	Al(6)(2-) - fusion of two aromatic Al(3)(-) units. A combined photoelectron spectroscopy and ab initio study of $M(+)[Al(6)(2-)]$ (M = Li, Na, K, Cu, and Au). <i>Journal of the American Chemical Society</i> , <b>2002</b> , 124, 11791-801	16.4	121
96	Probing the electronic structure of [MoOS(4)](-) centers using anionic photoelectron spectroscopy. Journal of the American Chemical Society, <b>2002</b> , 124, 10182-91	16.4	20
95	Lithium-Assisted Self-Assembly of Aluminum Carbide Nanowires and Nanoribbons. <i>Nano Letters</i> , <b>2002</b> , 2, 105-108	11.5	39
94	Photodetachment and theoretical study of free and water-solvated nitrate anions, NO3(H2O)n (n=0B). <i>Journal of Chemical Physics</i> , <b>2002</b> , 116, 561-570	3.9	69
93	Probing the Electronic Structure and Aromaticity of Pentapnictogen Cluster Anions Pn5- (Pn = P, As, Sb, and Bi) Using Photoelectron Spectroscopy and ab Initio Calculations. <i>Journal of Physical Chemistry A</i> , <b>2002</b> , 106, 5600-5606	2.8	84
92	Aromatic Mercury Clusters in Ancient Amalgams Work done at Utah State University is supported by the donors to The Petroleum Research Fund, administered by the American Chemical Society.  Work done at Iowa State University is supported by Basic Energy Sciences, the U.S. Department of	16.4	134
91	Energy. Work done at Washington State University is supported by the National Science  Photodetachment of F(H2O)n (n=14): Observation of charge-transfer states [F(H2O)n+] and the transition state of F+H2O-hydrogen abstraction reaction. Journal of Chemical Physics, 2001, 115, 2889-240, 3369-3372	892	49
90	Electronic structure of chromium oxides, CrOnland CrOn (n=18) from photoelectron spectroscopy and density functional theory calculations. <i>Journal of Chemical Physics</i> , <b>2001</b> , 115, 7935-79	944	105
89	Photoelectron spectroscopy of mono-niobium carbide clusters NbCn[h=21]): Evidence for a cyclic to linear structural transition. <i>Journal of Chemical Physics</i> , <b>2001</b> , 115, 5170-5178	3.9	33
88	The electronic structure of CuCl2 and CuBr2 from anion photoelectron spectroscopy and ab initio calculations. <i>Journal of Chemical Physics</i> , <b>2001</b> , 114, 7388-7395	3.9	51
87	Experimental and Theoretical Investigations of the Stability, Energetics, and Structures of H2PO4-, H2P2O72-, and H3P3O102-in the Gas Phase. <i>Journal of Physical Chemistry A</i> , <b>2001</b> , 105, 10468-10474	2.8	39
86	Bulk-like features in the photoemission spectra of hydrated doubly charged anion clusters. <i>Science</i> , <b>2001</b> , 294, 1322-5	33.3	171
85	Observation of all-metal aromatic molecules. <i>Science</i> , <b>2001</b> , 291, 859-61	33.3	531

84 Beyond Classical Stoichiometry: Experiment and Theory. *Journal of Physical Chemistry A*, **2001**, 105, 107598107751

Vibrationally Resolved Photoelectron Spectroscopy of MgO- and ZnO- and the Low-Lying Electronic States of MgO, MgO-, and ZnO. Journal of Physical Chemistry A, 2001, 105, 5709-5718  Experimental Observation of Pentaatomic Tetracoordinate Planar Si- and Ge-Containing Molecules: MAI(4)(-) and MAI(4) The theoretical work was done at Utah State University and supported by the donors of The Petroleum Research Fund (ACS-PRF no. 35255-AC6), administered by the American Probing the electronic structure of iron clusters using photoelectron spectroscopy. Chemical Physics, 2000, 262, 53-63  PHOTODETACHMENT PHOTOELECTRON SPECTROSCOPY OF TRANSITION METAL OXIDE SPECIES. Advanced Series in Physical Chemistry, 2000, 854-957  On the origin of planarity in Al5Iand Al5 clusters: The importance of a four-center peripheral bond. Journal of Chemical Physics, 2000, 113, 5130  The electronic structure and electron affinities of higher chlorine oxide radicals ClOx (x=28) from photoelectron spectroscopy of ClOxIanions. Journal of Chemical Physics, 2000, 113, 10928-10933  Aluminum cluster anions: Photoelectron spectroscopy and ab initio simulations. Physical Review B, 2000, 62, 13216-13228  Vibrationally resolved photoelectron spectroscopy of the first row transition metal and C3 clusters: MC3I[M=Sc, V, Cr, Mn, Fe, Co, and Ni). Journal of Chemical Physics, 2000, 112, 3602-3608  The electronic structure and chemical bonding of aluminum acetylide: Al2C2 and Al2C2IIAn experimental and theoretical investigation. Journal of Chemical Physics, 2000, 113, 2671-2679  (MgO)(-)(n) (n = 1-5) clusters: multipole-bound anions and photodetachment spectroscopy. Physical Review Letters, 2000, 85, 3145-8  Probing the electronic structure of redox species and direct determination of intrinsic reorganization energies of electron transfer reactions. Journal of Chemical Physics, 2000, 112, 6959-6962  Intramolecular Coulomb repulsion and anisotropies of the repulsive Coulomb barrier in multiply charged anions. Journal of Chemical Physics, 2000, 113, 653-661		
States of MgO, MgO-, and ZnO. Journal of Physical Chemistry A, 2001, 105, 5709-5718  Experimental Observation of Pentaatomic Tetracoordinate Planar Si- and Ge-Containing Molecules: MAI(4)(4) and MAI(4) The theoretical work was done at Utah State University and supported by the donors of The Petroleum Research Fund (ACS-PRF no. 35255-AC6), administered by the American Probing the electronic structure of iron clusters using photoelectron spectroscopy. Chemical Physics, 2000, 262, 53-63  PHOTODETACHMENT PHOTOELECTRON SPECTROSCOPY OF TRANSITION METAL OXIDE SPECIES. Advanced Series in Physical Chemistry, 2000, 854-957  On the origin of planarity in AISland AIS clusters: The importance of a four-center peripheral bond. Journal of Chemical Physics, 2000, 113, 5130  The electronic structure and electron affinities of higher chlorine oxide radicals ClOx (x=28) from photoelectron spectroscopy of ClOxBnions. Journal of Chemical Physics, 2000, 113, 10928-10933  Aluminum cluster anions: Photoelectron spectroscopy and ab initio simulations. Physical Review B, 2000, 62, 13216-13228  Vibrationally resolved photoelectron spectroscopy of the first row transition metal and C3 clusters: MC3I[M=Sc, V, Cr, Mn, Fe, Co, and Ni). Journal of Chemical Physics, 2000, 112, 3602-3608  The electronic structure and chemical bonding of aluminum acetylide: Al2C2 and Al2C2BIAn experimental and theoretical investigation. Journal of Chemical Physics, 2000, 113, 2671-2679  (MgO)(-)(n) (n = 1-5) clusters: multipole-bound anions and photodetachment spectroscopy. Physical Review Letters, 2000, 85, 3145-8  Probing the electronic structure of redox species and direct determination of intrinsic reorganization energies of electron transfer reactions. Journal of Chemical Physics, 2000, 112, 6959-6962  Intramolecular Coulomb repulsion and anisotropies of the repulsive Coulomb barrier in multiply charged anions. Journal of Chemical Physics, 2000, 113, 653-661	6.4	201
MAI(4)(-) and MAI(4) The theoretical work was done at Utah State University and supported by the donors of The Petroleum Research Fund (ACS-PRF no. 35255-AC6), administered by the American  Probing the electronic structure of iron clusters using photoelectron spectroscopy. Chemical Physics, 2000, 262, 53-63  PHOTODETACHMENT PHOTOELECTRON SPECTROSCOPY OF TRANSITION METAL OXIDE SPECIES. Advanced Series in Physical Chemistry, 2000, 854-957  On the origin of planarity in AISIand AIS clusters: The importance of a four-center peripheral bond. Journal of Chemical Physics, 2000, 113, 5130  The electronic structure and electron affinities of higher chlorine oxide radicals ClOx (x=28) from photoelectron spectroscopy of ClOxlanions. Journal of Chemical Physics, 2000, 113, 10928-10933  Aluminum cluster anions: Photoelectron spectroscopy and ab initio simulations. Physical Review B, 2000, 62, 13216-13228  Vibrationally resolved photoelectron spectroscopy of the first row transition metal and C3 clusters: MC3I[M=Sc, V, Cr, Mn, Fe, Co, and Ni). Journal of Chemical Physics, 2000, 112, 3602-3608  The electronic structure and chemical bonding of aluminum acetylide: Al2C2 and Al2C2IlAn experimental and theoretical investigation. Journal of Chemical Physics, 2000, 113, 2671-2679  (MgO)(-)(n) (n = 1-5) clusters: multipole-bound anions and photodetachment spectroscopy. Physical Review Letters, 2000, 85, 3145-8  Probing the electronic structure of redox species and direct determination of intrinsic reorganization energies of electron transfer reactions. Journal of Chemical Physics, 2000, 112, 6959-6962  Intramolecular Coulomb repulsion and anisotropies of the repulsive Coulomb barrier in multiply charged anions. Journal of Chemical Physics, 2000, 113, 653-661  Vibrationally resolved photoelectron spectra of CuCNland AgCNland ab initio studies of the	2.8	49
Physics, 2000, 262, 53-63  PHOTODETACHMENT PHOTOELECTRON SPECTROSCOPY OF TRANSITION METAL OXIDE SPECIES.  Advanced Series in Physical Chemistry, 2000, 854-957  On the origin of planarity in AlSiand AlS clusters: The importance of a four-center peripheral bond.  Journal of Chemical Physics, 2000, 113, 5130  The electronic structure and electron affinities of higher chlorine oxide radicals ClOx (x=2½) from photoelectron spectroscopy of ClOxianions. Journal of Chemical Physics, 2000, 113, 10928-10933  Aluminum cluster anions: Photoelectron spectroscopy and ab initio simulations. Physical Review B, 2000, 62, 13216-13228  Vibrationally resolved photoelectron spectroscopy of the first row transition metal and C3 clusters: MC3I[M=Sc, V, Cr, Mn, Fe, Co, and Ni). Journal of Chemical Physics, 2000, 112, 3602-3608  The electronic structure and chemical bonding of aluminum acetylide: Al2C2 and Al2C2IlAn experimental and theoretical investigation. Journal of Chemical Physics, 2000, 113, 2671-2679  (MgO)(-)(n) (n = 1-5) clusters: multipole-bound anions and photodetachment spectroscopy. Physical Review Letters, 2000, 85, 3145-8  Probing the electronic structure of redox species and direct determination of intrinsic reorganization energies of electron transfer reactions. Journal of Chemical Physics, 2000, 112, 6959-6962  Intramolecular Coulomb repulsion and anisotropies of the repulsive Coulomb barrier in multiply charged anions. Journal of Chemical Physics, 2000, 113, 653-661  Vibrationally resolved photoelectron spectra of CuCNIand AgCNIand ab initio studies of the	16.4	69
PHOTODETACHMENT PHOTOELECTRON SPECTROSCOPY OF TRANSITION METAL OXIDE SPECIES.  Advanced Series in Physical Chemistry, 2000, 854-957  On the origin of planarity in AlSiand Al5 clusters: The importance of a four-center peripheral bond.  Journal of Chemical Physics, 2000, 113, 5130  The electronic structure and electron affinities of higher chlorine oxide radicals ClOx (x=2f) from photoelectron spectroscopy of ClOxianions. Journal of Chemical Physics, 2000, 113, 10928-10933  Aluminum cluster anions: Photoelectron spectroscopy and ab initio simulations. Physical Review B, 2000, 62, 13216-13228  Vibrationally resolved photoelectron spectroscopy of the first row transition metal and C3 clusters: MC3I[M=Sc, V, Cr, Mn, Fe, Co, and Ni). Journal of Chemical Physics, 2000, 112, 3602-3608  The electronic structure and chemical bonding of aluminum acetylide: Al2C2 and Al2C2ilAn experimental and theoretical investigation. Journal of Chemical Physics, 2000, 113, 2671-2679  (MgO)(-)(n) (n = 1-5) clusters: multipole-bound anions and photodetachment spectroscopy. Physical Review Letters, 2000, 85, 3145-8  Probing the electronic structure of redox species and direct determination of intrinsic reorganization energies of electron transfer reactions. Journal of Chemical Physics, 2000, 112, 6959-6962  Intramolecular Coulomb repulsion and anisotropies of the repulsive Coulomb barrier in multiply charged anions. Journal of Chemical Physics, 2000, 113, 653-661  Vibrationally resolved photoelectron spectra of CuCNIand AgCNIand ab initio studies of the	2.3	82
The electronic structure and electron affinities of higher chlorine oxide radicals ClOx (x=28) from photoelectron spectroscopy of ClOx\( \text{Laminous}\). Journal of Chemical Physics, 2000, 113, 10928-10933  Aluminum cluster anions: Photoelectron spectroscopy and ab initio simulations. Physical Review B, 2000, 62, 13216-13228  Vibrationally resolved photoelectron spectroscopy of the first row transition metal and C3 clusters: MC3[[M=Sc, V, Cr, Mn, Fe, Co, and Ni). Journal of Chemical Physics, 2000, 112, 3602-3608  The electronic structure and chemical bonding of aluminum acetylide: Al2C2 and Al2C2[IAn experimental and theoretical investigation. Journal of Chemical Physics, 2000, 113, 2671-2679  (MgO)(-)(n) (n = 1-5) clusters: multipole-bound anions and photodetachment spectroscopy. Physical Review Letters, 2000, 85, 3145-8  Probing the electronic structure of redox species and direct determination of intrinsic reorganization energies of electron transfer reactions. Journal of Chemical Physics, 2000, 112, 6959-6962  Intramolecular Coulomb repulsion and anisotropies of the repulsive Coulomb barrier in multiply charged anions. Journal of Chemical Physics, 2000, 113, 653-661  Vibrationally resolved photoelectron spectra of CuCN[and AgCN[and ab initio studies of the		5
photoelectron spectroscopy of ClOxlanions. <i>Journal of Chemical Physics</i> , <b>2000</b> , 113, 10928-10933  Aluminum cluster anions: Photoelectron spectroscopy and ab initio simulations. <i>Physical Review B</i> , <b>2000</b> , 62, 13216-13228  Vibrationally resolved photoelectron spectroscopy of the first row transition metal and C3 clusters: MC3I[M=Sc, V, Cr, Mn, Fe, Co, and Ni). <i>Journal of Chemical Physics</i> , <b>2000</b> , 112, 3602-3608  The electronic structure and chemical bonding of aluminum acetylide: Al2C2 and Al2C2ElAn experimental and theoretical investigation. <i>Journal of Chemical Physics</i> , <b>2000</b> , 113, 2671-2679  (MgO)(-)(n) (n = 1-5) clusters: multipole-bound anions and photodetachment spectroscopy. <i>Physical Review Letters</i> , <b>2000</b> , 85, 3145-8  Probing the electronic structure of redox species and direct determination of intrinsic reorganization energies of electron transfer reactions. <i>Journal of Chemical Physics</i> , <b>2000</b> , 112, 6959-6962  Intramolecular Coulomb repulsion and anisotropies of the repulsive Coulomb barrier in multiply charged anions. <i>Journal of Chemical Physics</i> , <b>2000</b> , 113, 653-661  Vibrationally resolved photoelectron spectra of CuCNIand AgCNIand ab initio studies of the	3.9	45
76 2000, 62, 13216-13228  77 Vibrationally resolved photoelectron spectroscopy of the first row transition metal and C3 clusters: MC3I[M=Sc, V, Cr, Mn, Fe, Co, and Ni). Journal of Chemical Physics, 2000, 112, 3602-3608  78 The electronic structure and chemical bonding of aluminum acetylide: Al2C2 and Al2C2IIAn experimental and theoretical investigation. Journal of Chemical Physics, 2000, 113, 2671-2679  79 (MgO)(-)(n) (n = 1-5) clusters: multipole-bound anions and photodetachment spectroscopy. Physical Review Letters, 2000, 85, 3145-8  70 Probing the electronic structure of redox species and direct determination of intrinsic reorganization energies of electron transfer reactions. Journal of Chemical Physics, 2000, 112, 6959-6962  70 Intramolecular Coulomb repulsion and anisotropies of the repulsive Coulomb barrier in multiply charged anions. Journal of Chemical Physics, 2000, 113, 653-661  70 Vibrationally resolved photoelectron spectra of CuCNIand AgCNIand ab initio studies of the	3.9	46
75 MC3I[M=Sc, V, Cr, Mn, Fe, Co, and Ni). Journal of Chemical Physics, 2000, 112, 3602-3608  76 The electronic structure and chemical bonding of aluminum acetylide: Al2C2 and Al2C2IAn experimental and theoretical investigation. Journal of Chemical Physics, 2000, 113, 2671-2679  76 (MgO)(-)(n) (n = 1-5) clusters: multipole-bound anions and photodetachment spectroscopy. Physical Review Letters, 2000, 85, 3145-8  77 Probing the electronic structure of redox species and direct determination of intrinsic reorganization energies of electron transfer reactions. Journal of Chemical Physics, 2000, 112, 6959-6962  78 Intramolecular Coulomb repulsion and anisotropies of the repulsive Coulomb barrier in multiply charged anions. Journal of Chemical Physics, 2000, 113, 653-661  78 Vibrationally resolved photoelectron spectra of CuCNIand AgCNIand ab initio studies of the	3.3	105
experimental and theoretical investigation. <i>Journal of Chemical Physics</i> , <b>2000</b> , 113, 2671-2679  (MgO)(-)(n) (n = 1-5) clusters: multipole-bound anions and photodetachment spectroscopy. <i>Physical Review Letters</i> , <b>2000</b> , 85, 3145-8  Probing the electronic structure of redox species and direct determination of intrinsic reorganization energies of electron transfer reactions. <i>Journal of Chemical Physics</i> , <b>2000</b> , 112, 6959-6962  Intramolecular Coulomb repulsion and anisotropies of the repulsive Coulomb barrier in multiply charged anions. <i>Journal of Chemical Physics</i> , <b>2000</b> , 113, 653-661  Vibrationally resolved photoelectron spectra of CuCNIand AgCNIand ab initio studies of the	3.9	53
Probing the electronic structure of redox species and direct determination of intrinsic reorganization energies of electron transfer reactions. <i>Journal of Chemical Physics</i> , <b>2000</b> , 112, 6959-6962  Intramolecular Coulomb repulsion and anisotropies of the repulsive Coulomb barrier in multiply charged anions. <i>Journal of Chemical Physics</i> , <b>2000</b> , 113, 653-661  Vibrationally resolved photoelectron spectra of CuCNIand AgCNIand ab initio studies of the	3.9	28
reorganization energies of electron transfer reactions. <i>Journal of Chemical Physics</i> , <b>2000</b> , 112, 6959-6962  Intramolecular Coulomb repulsion and anisotropies of the repulsive Coulomb barrier in multiply charged anions. <i>Journal of Chemical Physics</i> , <b>2000</b> , 113, 653-661  Vibrationally resolved photoelectron spectra of CuCNIand AgCNIand ab initio studies of the	<sup>7</sup> ·4	47
charged anions. <i>Journal of Chemical Physics</i> , <b>2000</b> , 113, 653-661  Vibrationally resolved photoelectron spectra of CuCNIand AgCNIand ab initio studies of the	i.9	42
	3.9	44
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Experimental and theoretical study of the photoelectron spectra of MnOx[k=1B) clusters. <i>Journal of Chemical Physics</i> , <b>2000</b> , 113, 1473-1483	;.9	49
Probing the Electronic Structure and MetalMetal Bond of Re2Cl82- in the Gas Phase. <i>Journal of the American Chemical Society</i> , <b>2000</b> , 122, 2096-2100	16.4	32
Probing Free Multiply Charged Anions Using Photodetachment Photoelectron Spectroscopy.  Journal of Physical Chemistry A, <b>2000</b> , 104, 1978-1990	2.8	180

#### (1999-2000)

66	Photoelectron Spectroscopy and Theoretical Calculations of SO4- and HSO4-: Confirmation of High Electron Affinities of SO4 and HSO4. <i>Journal of Physical Chemistry A</i> , <b>2000</b> , 104, 504-508	2.8	54
65	Experimental and Theoretical Investigations of the Stability of Two Small Gaseous Dicarboxylate Dianions: Acetylene Dicarboxylate and Succinate. <i>Journal of the American Chemical Society</i> , <b>2000</b> , 122, 4499-4507	16.4	37
64	Experimental Observation of Pentaatomic Tetracoordinate Planar Carbon-Containing Molecules. Journal of the American Chemical Society, <b>2000</b> , 122, 7681-7687	16.4	188
63	Photodetachment of Gaseous Multiply Charged Anions, Copper Phthalocyanine Tetrasulfonate Tetraanion: Tuning Molecular Electronic Energy Levels by Charging and Negative Electron Binding. <i>Journal of Physical Chemistry A</i> , <b>2000</b> , 104, 25-33	2.8	44
62	Photodetachment of the First Zwitterionic Anions in the Gas Phase: Probing Intramolecular Coulomb Repulsion and Attraction. <i>Journal of the American Chemical Society</i> , <b>2000</b> , 122, 8305-8306	16.4	11
61	Experimental Observation of a Very High Second Electron Affinity for ZrF6 from Photodetachment of Gaseous ZrF62- Doubly Charged Anions. <i>Journal of Physical Chemistry A</i> , <b>2000</b> , 104, 4429-4432	2.8	38
60	Photodetachment of Multiply Charged Anions: The Electronic Structure of Gaseous Square-Planar Transition Metal Complexes PtX42- ( $X = Cl$ , Br). <i>Journal of the American Chemical Society</i> , <b>2000</b> , 122, 23	39-2 <del>3</del> 4	5 <sup>45</sup>
59	Napoleon Hatl Structure of Tetraatomic Molecules. A Combined Photoelectron Spectroscopy and Ab Initio Study of CAlSi2- and Its Neutral. <i>Journal of Physical Chemistry A</i> , <b>2000</b> , 104, 5358-5365	2.8	6
58	Electronic instability of isolated SO42[and its solvation stabilization. <i>Journal of Chemical Physics</i> , <b>2000</b> , 113, 10837-10840	3.9	111
57	TEMPERATURE EFFFECTS IN ANION PHOTOELECTRON SPECTROSCOPY OF METAL CLUSTERS <b>2000</b> ,		17
56	Experimental Search for the Smallest Stable Multiply Charged Anions in the Gas Phase. <i>Physical Review Letters</i> , <b>1999</b> , 83, 3402-3405	7.4	99
55	Photodetachment photoelectron spectroscopy of multiply charged anions using electrospray ionization. <i>Review of Scientific Instruments</i> , <b>1999</b> , 70, 1957-1966	1.7	195
54	Photodetachment of free hexahalogenometallate doubly charged anions in the gas phase: [ML6]2[I(M=Re, Os, Ir, Pt; L=Cl and Br). <i>Journal of Chemical Physics</i> , <b>1999</b> , 111, 4497-4509	3.9	88
53	Combined photoelectron spectroscopy and ab initio study of the hypermetallic Al3C molecule. <i>Journal of Chemical Physics</i> , <b>1999</b> , 110, 8980-8985	3.9	36
52	First experimental photoelectron spectra of superhalogens and their theoretical interpretations. <i>Journal of Chemical Physics</i> , <b>1999</b> , 110, 4763-4771	3.9	243
51	Photoelectron spectra of aluminum cluster anions: Temperature effects and ab initio simulations. <i>Physical Review B</i> , <b>1999</b> , 60, R11297-R11300	3.3	247
50	Photodetachment photoelectron spectroscopy of doubly charged anions: S2O82\(\textit{Journal of Chemical Physics}\), 110, 3635-3638	3.9	31
49	Electronic structure and chemical bonding between the first row transition metals and C2: A photoelectron spectroscopy study of MC2[[M=Sc, V, Cr, Mn, Fe, and Co). <i>Journal of Chemical Physics</i> 1999, 111, 8389-8395	3.9	95

48	The electronic structure of MoC and WC by anion photoelectron spectroscopy. <i>Journal of Chemical Physics</i> , <b>1999</b> , 111, 2464-2469	3.9	33
47	High resolution photoelectron spectroscopy of C60□ <i>Journal of Chemical Physics</i> , <b>1999</b> , 110, 8217-8220	3.9	174
46	Observation of a spin-protected high-energy isomer of Al4NItluster. <i>Chemical Physics Letters</i> , <b>1999</b> , 301, 379-384	2.5	42
45	Electron tunneling through the repulsive Coulomb barrier in photodetachment of multiply charged anions. <i>Chemical Physics Letters</i> , <b>1999</b> , 307, 391-396	2.5	78
44	Vibrationally resolved photoelectron spectroscopy of POB and the electronic structure of PO3. <i>Chemical Physics Letters</i> , <b>1999</b> , 313, 179-183	2.5	12
43	Observation of negative electron-binding energy in a molecule. <i>Nature</i> , <b>1999</b> , 400, 245-248	50.4	190
42	Origin of the unusual stability of MnO4\(\textit{D}\)Chemical Physics Letters, <b>1999</b> , 312, 598-605	2.5	125
41	The electronic structure and chemical bonding of hypermetallic Al5C by ab initio calculations and anion photoelectron spectroscopy. <i>Journal of Chemical Physics</i> , <b>1999</b> , 111, 4993-4998	3.9	39
40	Tetracoordinated Planar Carbon in the Al4C- Anion. A Combined Photoelectron Spectroscopy and ab Initio Study. <i>Journal of the American Chemical Society</i> , <b>1999</b> , 121, 6033-6038	16.4	242
39	Investigation of Free Singly and Doubly Charged Alkali Metal Sulfate Ion Pairs: M+(SO42-) and [M+(SO42-)]2 (M = Na, K). <i>Journal of Physical Chemistry A</i> , <b>1999</b> , 103, 3423-3429	2.8	55
38	Eland ECoordinated Al in AlC2- and AlCSi A Combined Photoelectron Spectroscopy and ab Initio Study. <i>Journal of the American Chemical Society</i> , <b>1999</b> , 121, 10193-10197	16.4	30
37	New Magic Numbers in TixCy- Anion Clusters and Implications for the Growth Mechanisms of Titanium Carbide Clusters. <i>Journal of the American Chemical Society</i> , <b>1998</b> , 120, 6556-6562	16.4	45
36	Al3Oy (y=0B) clusters: Sequential oxidation, metal-to-oxide transformation, and photoisomerization. <i>Journal of Chemical Physics</i> , <b>1998</b> , 109, 449-458	3.9	88
35	Photoelectron Spectroscopy of Doubly Charged Anions: Intramolecular Coulomb Repulsion and Solvent Stabilization. <i>Journal of Physical Chemistry A</i> , <b>1998</b> , 102, 8633-8636	2.8	7 <sup>2</sup>
34	Photoelectron Spectroscopy and Electronic Structure of ScOn-(n= 1월) and YOn-(n= 1₿):□Strong Electron Correlation Effects in ScO-and YO <i>Journal of Physical Chemistry A</i> , <b>1998</b> , 102, 9129-9135	2.8	85
33	s- p Hybridization and Electron Shell Structures in Aluminum Clusters: A Photoelectron Spectroscopy Study. <i>Physical Review Letters</i> , <b>1998</b> , 81, 1909-1912	7.4	301
32	A photoelectron spectroscopic study of monovanadium oxide anions (VOxIx=1I). <i>Journal of Chemical Physics</i> , <b>1998</b> , 108, 5310-5318	3.9	98
31	The chemical bonding and electronic structure of RhC, RhN, and RhO by anion photoelectron spectroscopy. <i>Journal of Chemical Physics</i> , <b>1998</b> , 109, 5264-5268	3.9	37

30	Probing the Potential Barriers and Intramolecular Electrostatic Interactions in Free Doubly Charged Anions. <i>Physical Review Letters</i> , <b>1998</b> , 81, 2667-2670	7.4	171
29	Photodetachment Spectroscopy of a Doubly Charged Anion: Direct Observation of the Repulsive Coulomb Barrier. <i>Physical Review Letters</i> , <b>1998</b> , 81, 3351-3354	7.4	125
28	Photoelectron Spectroscopy of Transition Metal Clusters. <i>Zeitschrift Fur Physikalische Chemie</i> , <b>1998</b> , 203, 45-55	3.1	17
27	Photoelectron spectroscopy of small chromium clusters: Observation of even-odd alternations and theoretical interpretation. <i>Physical Review B</i> , <b>1997</b> , 55, 12884-12887	3.3	42
26	Si3Oy (y=1B) Clusters: Models for Oxidation of Silicon Surfaces and Defect Sites in Bulk Oxide Materials. <i>Physical Review Letters</i> , <b>1997</b> , 78, 4450-4453	7.4	105
25	A study of nickel monoxide (NiO), nickel dioxide (ONiO), and Ni(O2) complex by anion photoelectron spectroscopy. <i>Journal of Chemical Physics</i> , <b>1997</b> , 107, 16-21	3.9	106
24	Electronic structure and photoelectron spectroscopy of AlSi mixed dimer. <i>Journal of Chemical Physics</i> , <b>1997</b> , 107, 7667-7672	3.9	9
23	Growth Pathways of Metallocarbohedrenes: Cagelike or Cubic?. <i>Physical Review Letters</i> , <b>1997</b> , 78, 2983-	-2 <del>,</del> 9 <u>.</u> 86	36
22	Probing the Electronic Structure of Metallocarbohedrenes: M8C12 (M = Ti, V, Cr, Zr, Nb). <i>Journal of the American Chemical Society</i> , <b>1997</b> , 119, 7417-7422	16.4	53
21	Vibrationally Resolved Photoelectron Spectra of TiCx- (x = 25) Clusters. <i>Journal of Physical Chemistry A</i> , <b>1997</b> , 101, 7699-7701	2.8	51
20	Chemical Bonding between Cu and OxygenCopper Oxides vs O2Complexes: A Study of CuOx(x= 0B) Species by Anion Photoelectron Spectroscopy. <i>Journal of Physical Chemistry A</i> , <b>1997</b> , 101, 2103-211	1 <sup>2.8</sup>	117
19	A study of the structure and bonding of small aluminum oxide clusters by photoelectron spectroscopy: AlxOy[[x=12], y=15]). <i>Journal of Chemical Physics</i> , <b>1997</b> , 106, 1309-1317	3.9	139
18	Small silicon oxide clusters: chains and rings. <i>Zeitschrift Fli Physik D-Atoms Molecules and Clusters</i> , <b>1997</b> , 40, 36-39		69
17	STUDY OF IRON©ARBON MIXED CLUSTERS FeCn (n=28): A POSSIBLE LINEAR-TO-CYCLIC TRANSITION FROM FeC3 TO FeC4. <i>Surface Review and Letters</i> , <b>1996</b> , 03, 423-427	1.1	20
16	Dimer Growth, Structural Transition, and Antiferromagnetic Ordering of Small Chromium Clusters. <i>Physical Review Letters</i> , <b>1996</b> , 77, 51-54	7.4	117
15	Observation and Photoelectron Spectroscopic Study of Novel Mono- and Diiron Oxide Molecules: FeOy- (y = $14$ ) and Fe2Oy- (y = $14$ ). <i>Journal of the American Chemical Society</i> , <b>1996</b> , 118, 5296-5301	16.4	111
14	Vibrationally resolved photoelectron spectroscopy of AlO[and AlO2[]International Journal of Mass Spectrometry and Ion Processes, <b>1996</b> , 159, 75-80		33
13	Electronic structure of small titanium clusters: Emergence and evolution of the 3d band. <i>Physical Review Letters</i> , <b>1996</b> , 76, 212-215	7.4	78

12	Electronic structure of small copper oxide clusters: From Cu2O to Cu2O4. <i>Physical Review B</i> , <b>1996</b> , 53, 8028-8031	3.3	97
11	Evolution of the Electronic Structure of Small Vanadium Clusters from Molecular to Bulklike. <i>Physical Review Letters</i> , <b>1996</b> , 77, 2436-2439	7.4	68
10	Sequential oxygen atom chemisorption on surfaces of small iron clusters. <i>Physical Review Letters</i> , <b>1996</b> , 76, 4853-4856	7.4	120
9	IRON CLUSTERS AND OXYGEN-CHEMISORBED IRON CLUSTERS. <i>Surface Review and Letters</i> , <b>1996</b> , 03, 695-699	1.1	23
8	Si3O4-: vibrationally resolved photoelectron spectrum and ab initio calculations. <i>Journal of the American Chemical Society</i> , <b>1995</b> , 117, 5417-5418	16.4	31
7	Probing the electronic structure of small iron clusters. <i>Chemical Physics Letters</i> , <b>1995</b> , 236, 57-63	2.5	12
6	Threshold photodetachment of cold CBO. Chemical Physics Letters, 1991, 182, 5-11	2.5	294
5	Fullerene triplet state production and decay: R2PI probes of C60 and C70 in a supersonic beam. <i>Chemical Physics Letters</i> , <b>1991</b> , 179, 449-454	2.5	186
4	High temperature and high resolution UV photoelectron spectroscopy using supersonic molecular beams. <i>Journal of Electron Spectroscopy and Related Phenomena</i> , <b>1990</b> , 51, 513-526	1.7	11
3	High resolution photoelectron spectroscopy of clusters of group V elements. <i>Physica Scripta</i> , <b>1990</b> , 41, 867-869	2.6	8
2	Vibrational spectra of Se2+ and Te2+ in their ground states. <i>Chemical Physics Letters</i> , <b>1989</b> , 158, 297-30	002.5	17
1	High resolution UV photoelectron spectroscopy of CO+2, COS+ and CS+2 using supersonic molecular beams. <i>Journal of Electron Spectroscopy and Related Phenomena</i> , <b>1988</b> , 47, 167-186	1.7	132