

Uzi Landman

List of Publications by Year in descending order

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364
times ranked

16971
citing authors

#	ARTICLE	IF	CITATIONS
1	Nanotribology: friction, wear and lubrication at the atomic scale. Nature, 1995, 374, 607-616.	27.8	1,514
2	When Gold Is Not Noble: Nanoscale Gold Catalysts. Journal of Physical Chemistry A, 1999, 103, 9573-9578.	2.5	1,375
3	Charging Effects on Bonding and Catalyzed Oxidation of CO on Au ₈ Clusters on MgO. Science, 2005, 307, 403-407.	12.6	1,358
4	Nanocrystal gold molecules. Advanced Materials, 1996, 8, 428-433.	21.0	1,179
5	Ultrastable silver nanoparticles. Nature, 2013, 501, 399-402.	27.8	1,023
6	Bonding in Cu, Ag, and Au Clusters: Relativistic Effects, Trends, and Surprises. Physical Review Letters, 2002, 89, 033401.	7.8	611
7	On the Electronic and Atomic Structures of Small Au _N - (N = 4-14) Clusters: A Photoelectron Spectroscopy and Density-Functional Study. Journal of Physical Chemistry A, 2003, 107, 6168-6175.	2.5	598
8	Born-Oppenheimer molecular-dynamics simulations of finite systems: Structure and dynamics of (H ₂ O) ₂ . Physical Review B, 1993, 48, 2081-2097.	3.2	551
9	Structural, Electronic, and Impurity-Doping Effects in Nanoscale Chemistry: Supported Gold Nanoclusters. Angewandte Chemie - International Edition, 2003, 42, 1297-1300.	13.8	547
10	Total Structure and Electronic Properties of the Gold Nanocrystal Au ₃₆ (SR) ₂₄ . Angewandte Chemie - International Edition, 2012, 51, 13114-13118.	13.8	519
11	Structural Evolution of Smaller Gold Nanocrystals: The Truncated Decahedral Motif. Physical Review Letters, 1997, 79, 1873-1876.	7.8	460
12	Frictional Forces and Amontons' Law: From the Molecular to the Macroscopic Scale. Journal of Physical Chemistry B, 2004, 108, 3410-3425.	2.6	455
13	Gold clusters(Au _N , 2 ≤ N ≤ 10) and their anions. Physical Review B, 2000, 62, R2287-R2290.	3.2	454
14	Structure, Dynamics, and Thermodynamics of Passivated Gold Nanocrystallites and Their Assemblies. The Journal of Physical Chemistry, 1996, 100, 13323-13329.	2.9	410
15	Catalytic CO Oxidation by Free Au ₂ : Experiment and Theory. Journal of the American Chemical Society, 2003, 125, 10437-10445.	13.7	386
16	Charge Migration in DNA: Ion-Gated Transport. Science, 2001, 294, 567-571.	12.6	373
17	The energetics and structure of nickel clusters: Size dependence. Journal of Chemical Physics, 1991, 94, 7376-7396.	3.0	359
18	Interaction of O ₂ with Gold Clusters: Molecular and Dissociative Adsorption. Journal of Physical Chemistry A, 2003, 107, 4066-4071.	2.5	349

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19	Formation, Stability, and Breakup of Nanojets. <i>Science</i> , 2000, 289, 1165-1169.	12.6	344
20	Layering Transitions and Dynamics of Confined Liquid Films. <i>Physical Review Letters</i> , 1997, 79, 705-708.	7.8	303
21	Oxidation of DNA: Damage to Nucleobases. <i>Accounts of Chemical Research</i> , 2010, 43, 280-287.	15.6	300
22	Photoelectron spectra of aluminum cluster anions: Temperature effects and ab initio simulations. <i>Physical Review B</i> , 1999, 60, R11297-R11300.	3.2	289
23	Bonding Trends and Dimensionality Crossover of Gold Nanoclusters on Metal-Supported MgO Thin Films. <i>Physical Review Letters</i> , 2006, 97, 036106.	7.8	268
24	Electron localization in water clusters. II. Surface and internal states. <i>Journal of Chemical Physics</i> , 1988, 88, 4429-4447.	3.0	251
25	Melting of gold clusters. <i>Physical Review B</i> , 1999, 60, 5065-5077.	3.2	242
26	Small is different: energetic, structural, thermal, and mechanical properties of passivated nanocluster assemblies. <i>Faraday Discussions</i> , 2004, 125, 1.	3.2	239
27	Structural evolution of Au nanoclusters: From planar to cage to tubular motifs. <i>Physical Review B</i> , 2006, 74, .	3.2	234
28	Structure and Thermodynamics of Self-Assembled Monolayers on Gold Nanocrystallites. <i>Journal of Physical Chemistry B</i> , 1998, 102, 6566-6572.	2.6	231
29	Gas-Phase Catalytic Oxidation of CO by Au ₂ ⁻ . <i>Journal of the American Chemical Society</i> , 2001, 123, 9704-9705.	13.7	230
30	Slip Diffusion and Lévy Flights of an Adsorbed Gold Nanocluster. <i>Physical Review Letters</i> , 1999, 82, 3835-3838.	7.8	229
31	Spontaneous Symmetry Breaking in Single and Molecular Quantum Dots. <i>Physical Review Letters</i> , 1999, 82, 5325-5328.	7.8	224
32	Reversible Manipulations of Room Temperature Mechanical and Quantum Transport Properties in Nanowire Junctions. <i>Physical Review Letters</i> , 1996, 77, 1362-1365.	7.8	219
33	Water-Enhanced Catalysis of CO Oxidation on Free and Supported Gold Nanoclusters. <i>Physical Review Letters</i> , 2005, 95, 106102.	7.8	211
34	Control and Manipulation of Gold Nanocatalysis: Effects of Metal Oxide Support Thickness and Composition. <i>Journal of the American Chemical Society</i> , 2009, 131, 538-548.	13.7	203
35	Structure and solvation forces in confined films: Linear and branched alkanes. <i>Journal of Chemical Physics</i> , 1997, 106, 4309-4318.	3.0	202
36	Melting of Gold Clusters: Icosahedral Precursors. <i>Physical Review Letters</i> , 1998, 81, 2036-2039.	7.8	199

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37	Structure and Magnetism of Neutral and Anionic Palladium Clusters. <i>Physical Review Letters</i> , 2001, 86, 2545-2548.	7.8	198
38	Symmetry breaking and quantum correlations in finite systems: studies of quantum dots and ultracold Bose gases and related nuclear and chemical methods. <i>Reports on Progress in Physics</i> , 2007, 70, 2067-2148.	20.1	198
39	Size-Dependent Structural Evolution and Chemical Reactivity of Gold Clusters. <i>ChemPhysChem</i> , 2007, 8, 157-161.	2.1	197
40	Preparation, structure, dynamics, and energetics of amorphous silicon: A molecular-dynamics study. <i>Physical Review B</i> , 1989, 40, 1164-1174.	3.2	192
41	Interfacial alkane films. <i>Physical Review Letters</i> , 1992, 69, 1967-1970.	7.8	191
42	Factors in gold nanocatalysis: oxidation of CO in the non-scalable size regime. <i>Topics in Catalysis</i> , 2007, 44, 145-158.	2.8	190
43	Metal-Semiconductor Nanocontacts: Silicon Nanowires. <i>Physical Review Letters</i> , 2000, 85, 1958-1961.	7.8	188
44	Preparation and melting of amorphous silicon by molecular-dynamics simulations. <i>Physical Review B</i> , 1988, 37, 4656-4663.	3.2	186
45	Predicted Oxidation of CO Catalyzed by Au Nanoclusters on a Thin Defect-Free MgO Film Supported on a Mo(100) Surface. <i>Journal of the American Chemical Society</i> , 2007, 129, 2228-2229.	13.7	167
46	Atomistic mechanisms of adhesive contact formation and interfacial processes. <i>Wear</i> , 1992, 153, 3-30.	3.1	164
47	Electron Localization in Alkali-Halide Clusters. <i>Physical Review Letters</i> , 1985, 54, 1860-1863.	7.8	162
48	Electron localization in water clusters. I. Electron-water pseudopotential. <i>Journal of Chemical Physics</i> , 1988, 88, 4421-4428.	3.0	158
49	Structural and dynamical consequences of interactions in interfacial systems. <i>Journal of Vacuum Science and Technology A: Vacuum, Surfaces and Films</i> , 1989, 7, 2829-2839.	2.1	154
50	Lattice relaxation at metal surfaces: An electrostatic model. <i>Physical Review B</i> , 1980, 21, 448-457.	3.2	148
51	Faceting at the silicon (100) crystal-melt interface: Theory and experiment. <i>Physical Review Letters</i> , 1986, 56, 155-158.	7.8	146
52	Atomic-Scale Issues in Tribology: Interfacial Junctions and Nano-elastohydrodynamics. <i>Langmuir</i> , 1996, 12, 4514-4528.	3.5	141
53	Au ₆₇ (SR) ₃₅ Nanomolecules: Characteristic Size-Specific Optical, Electrochemical, Structural Properties and First-Principles Theoretical Analysis. <i>Journal of Physical Chemistry A</i> , 2013, 117, 504-517.	2.5	140
54	Surface melting of Ni(110). <i>Physical Review B</i> , 1990, 41, 439-450.	3.2	137

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55	Methane Activation and Catalytic Ethylene Formation on Free Au ₂ ⁺ . <i>Angewandte Chemie - International Edition</i> , 2010, 49, 980-983.	13.8	137
56	Structures, solvation forces and shear of molecular films in a rough nano-confinement. <i>Tribology Letters</i> , 2000, 9, 3-13.	2.6	136
57	Gold Nanowires and Their Chemical Modifications. <i>Journal of Physical Chemistry B</i> , 1999, 103, 8814-8816.	2.6	135
58	Nanomechanics and dynamics of tip-substrate interactions. <i>Journal of Vacuum Science & Technology an Official Journal of the American Vacuum Society B, Microelectronics Processing and Phenomena</i> , 1991, 9, 414.	1.6	134
59	Hydration of sodium in water clusters. <i>Physical Review Letters</i> , 1993, 70, 1775-1778.	7.8	134
60	Cluster-derived structures and conductance fluctuations in nanowires. <i>Nature</i> , 1997, 387, 788-791.	27.8	131
61	Friction Control in Thin-Film Lubrication. <i>Journal of Physical Chemistry B</i> , 1998, 102, 5033-5037.	2.6	129
62	Hydrogen-bonded structure and mechanical chiral response of a silver nanoparticle superlattice. <i>Nature Materials</i> , 2014, 13, 807-811.	27.5	128
63	Patterns and barriers for fission of charged small metal clusters. <i>Physical Review Letters</i> , 1991, 67, 3058-3061.	7.8	126
64	Electronic shell effects in triaxially deformed metal clusters: A systematic interpretation of experimental observations. <i>Physical Review B</i> , 1995, 51, 1902-1917.	3.2	125
65	Origins of Solvation Forces in Confined Films. <i>Journal of Physical Chemistry B</i> , 1997, 101, 4013-4023.	2.6	125
66	Stabilized-jellium description of neutral and multiply charged fullerenes C _x ±60. <i>Chemical Physics Letters</i> , 1994, 217, 175-185.	2.6	123
67	Chemistry and Structure of Silver Molecular Nanoparticles. <i>Accounts of Chemical Research</i> , 2018, 51, 3104-3113.	15.6	123
68	Surface states of excess electrons on water clusters. <i>Physical Review Letters</i> , 1987, 59, 811-814.	7.8	122
69	STEM Electron Diffraction and High-Resolution Images Used in the Determination of the Crystal Structure of the Au ₁₄₄ (SR) ₆₀ Cluster. <i>Journal of Physical Chemistry Letters</i> , 2013, 4, 975-981.	4.6	122
70	CO Combustion on Supported Gold Clusters. <i>ChemPhysChem</i> , 2006, 7, 1871-1879.	2.1	121
71	The Superstable 25 kDa Monolayer Protected Silver Nanoparticle: Measurements and Interpretation as an Icosahedral Ag ₁₅₂ (SCH ₂ CH ₂ Ph) ₆₀ Cluster. <i>Nano Letters</i> , 2012, 12, 5861-5866.	9.1	121
72	Collective and Independent-Particle Motion in Two-Electron Artificial Atoms. <i>Physical Review Letters</i> , 2000, 85, 1726-1729.	7.8	119

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73	Relaxation dynamics following transition of solvated electrons. <i>Journal of Chemical Physics</i> , 1989, 90, 4413-4422.	3.0	117
74	Structure sensitivity in the non-scalable regime explored via catalysed ethylene hydrogenation on supported platinum nanoclusters. <i>Nature Communications</i> , 2016, 7, 10389.	12.8	115
75	Collective excitations of multishell carbon microstructures: Multishell fullerenes and coaxial nanotubes. <i>Physical Review B</i> , 1996, 53, 10225-10236.	3.2	113
76	Hydrogen-Promoted Oxygen Activation by Free Gold Cluster Cations. <i>Journal of the American Chemical Society</i> , 2009, 131, 8939-8951.	13.7	113
77	Aluminum cluster anions: Photoelectron spectroscopy and ab initio simulations. <i>Physical Review B</i> , 2000, 62, 13216-13228.	3.2	111
78	Controlled Deposition and Classification of Copper Nanoclusters. <i>The Journal of Physical Chemistry</i> , 1994, 98, 3527-3537.	2.9	110
79	Theory of Physisorption: He on Metals. <i>Physical Review B</i> , 1973, 8, 5484-5495.	3.2	109
80	Dynamics and spectra of a solvated electron in water clusters. <i>Journal of Chemical Physics</i> , 1988, 89, 2242-2256.	3.0	106
81	Optical spectra of localized excess electrons in alkali halide clusters. <i>Physical Review Letters</i> , 1990, 64, 2933-2936.	7.8	106
82	Surface premelting of Cu(110). <i>Physical Review B</i> , 1991, 44, 3226-3239.	3.2	106
83	Nanowire Gold Chains: Formation Mechanisms and Conductance. <i>Journal of Physical Chemistry B</i> , 2000, 104, 9063-9066.	2.6	106
84	Metallization of ionic clusters. <i>Physical Review Letters</i> , 1991, 67, 727-730.	7.8	104
85	Superheating, melting, and annealing of copper surfaces. <i>Physical Review Letters</i> , 1993, 71, 1023-1026.	7.8	103
86	Quantum dynamical simulations of nonadiabatic processes: Solvation dynamics of the hydrated electron. <i>Physical Review Letters</i> , 1991, 67, 1011-1014.	7.8	101
87	Structure and dynamics of n-alkanes confined by solid surfaces. I. Stationary crystalline boundaries. <i>Journal of Chemical Physics</i> , 1992, 97, 1937-1949.	3.0	99
88	Size dependence of the energetics of electron attachment to large water clusters. <i>Chemical Physics Letters</i> , 1988, 145, 382-386.	2.6	97
89	Multilayer lattice relaxation at metal surfaces: A total-energy minimization. <i>Physical Review B</i> , 1983, 28, 1685-1695.	3.2	96
90	Genetic Algorithms for Structural Cluster Optimization. <i>Journal of Physical Chemistry A</i> , 1998, 102, 6129-6137.	2.5	94

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91	Instability Driven Fragmentation of Nanoscale Fractal Islands. <i>Physical Review Letters</i> , 2002, 88, 196103.	7.8	94
92	Quantum electronic transport through three-dimensional microconstrictions with variable shapes. <i>Physical Review B</i> , 1996, 53, 4054-4064.	3.2	92
93	Molecular-orbital-self-consistent-field cluster model of H ₂ O adsorption on copper. <i>Physical Review B</i> , 1985, 32, 1430-1433.	3.2	91
94	Nanojets, Electrospray, and Ion Field Evaporation: Molecular Dynamics Simulations and Laboratory Experiments. <i>Journal of Physical Chemistry A</i> , 2008, 112, 9628-9649.	2.5	91
95	Infinite Conical Well: An Analytic Model for Quantum Mechanical Hindered Rotors. <i>Physical Review Letters</i> , 1982, 49, 426-430.	7.8	88
96	Dynamics and energy release in fission of small doubly charged clusters. <i>Physical Review Letters</i> , 1994, 72, 1636-1639.	7.8	87
97	Born-Oppenheimer dynamics using density-functional theory: Equilibrium and fragmentation of small sodium clusters. <i>Journal of Chemical Physics</i> , 1991, 94, 608-616.	3.0	85
98	Excited-state dynamics of rare-gas clusters. <i>Journal of Chemical Physics</i> , 1988, 88, 4273-4288.	3.0	84
99	All-quantum simulations: H ₃ O ⁺ and H ₅ O ₂ ⁺ . <i>Chemical Physics Letters</i> , 1995, 237, 161-170.	2.6	78
100	Molecular-dynamics simulations of epitaxial crystal growth from the melt. I. Si(100). <i>Physical Review B</i> , 1988, 37, 4637-4646.	3.2	77
101	Two-dimensional quantum dots in high magnetic fields: Rotating-electron-molecule versus composite-fermion approach. <i>Physical Review B</i> , 2003, 68, .	3.2	77
102	Size-Selected Monodisperse Nanoclusters on Supported Graphene: Bonding, Isomerism, and Mobility. <i>Nano Letters</i> , 2012, 12, 5907-5912.	9.1	76
103	Crystal-melt and melt-vapor interfaces of nickel. <i>Physical Review B</i> , 1989, 40, 924-932.	3.2	75
104	Multilayer lattice relaxation at metal surfaces. <i>Physical Review B</i> , 1983, 27, 6534-6537.	3.2	73
105	Edge states, Aharonov-Bohm oscillations, and thermodynamic and spectral properties in a two-dimensional electron gas with an antidot. <i>Physical Review B</i> , 1995, 52, 14067-14077.	3.2	73
106	Structure and Energetics of Ionized Water Clusters: $(\text{H}_2\text{O})_n^+$, $n = 2 \sim 5$. <i>Journal of Physical Chemistry A</i> , 1997, 101, 164-169.	2.5	73
107	Shell-correction method for calculating the binding energy of metal clusters: Application to multiply charged anions. <i>Physical Review B</i> , 1993, 48, 8376-8387.	3.2	72
108	Intercalation of Trioxatriangulenium Ion in DNA: Binding, Electron Transfer, X-ray Crystallography, and Electronic Structure. <i>Journal of the American Chemical Society</i> , 2003, 125, 2072-2083.	13.7	72

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109	Metal-on-metal thin-film growth: Au/Ni(001) and Ni/Au(001). <i>Physical Review B</i> , 1991, 44, 5970-5972.	3.2	71
110	Selective C-H Bond Cleavage in Methane by Small Gold Clusters. <i>Angewandte Chemie - International Edition</i> , 2017, 56, 13406-13410.	13.8	71
111	Surface and internal excess electron states in molecular clusters. <i>Accounts of Chemical Research</i> , 1989, 22, 350-357.	15.6	69
112	Molecular Dynamics in Shape Space and Femtosecond Vibrational Spectroscopy of Metal Clusters. <i>Journal of Physical Chemistry A</i> , 1998, 102, 2505-2508.	2.5	68
113	Cluster isomerization induced by electron attachment. <i>Journal of Chemical Physics</i> , 1987, 87, 2716-2723.	3.0	67
114	Energetics and structures of aluminum-lithium clusters. <i>Physical Review B</i> , 1993, 48, 1820-1824.	3.2	67
115	Trial wave functions with long-range Coulomb correlations for two-dimensional N-electron systems in high magnetic fields. <i>Physical Review B</i> , 2002, 66, .	3.2	67
116	Energetics and structures of neutral and charged S_n ($n \leq 10$) and sodium-doped S_nNa clusters. <i>Physical Review B</i> , 1997, 55, 7935-7944.	3.2	66
117	Hydrogen Welding and Hydrogen Switches in a Monatomic Gold Nanowire. <i>Nano Letters</i> , 2004, 4, 1845-1852.	9.1	66
118	Multiply Charged Metal Cluster Anions. <i>Physical Review Letters</i> , 2001, 86, 2996-2999.	7.8	64
119	Ethylene hydrogenation on supported Ni, Pd and Pt nanoparticles: Catalyst activity, deactivation and the d-band model. <i>Journal of Catalysis</i> , 2016, 333, 51-58.	6.2	62
120	Dynamics of excess electron migration, solvation, and spectra in polar molecular clusters. <i>Journal of Chemical Physics</i> , 1989, 91, 5567-5580.	3.0	60
121	Quantum simulations and ab initio electronic structure studies of $(H_2O)_n$. <i>Journal of Chemical Physics</i> , 1989, 91, 7797-7808.	3.0	59
122	Energetics, forces, and quantized conductance in jellium-modeled metallic nanowires. <i>Physical Review B</i> , 1998, 57, 4872-4882.	3.2	59
123	Effect of Base Sequence and Hydration on the Electronic and Hole Transport Properties of Duplex DNA: A Theory and Experiment. <i>Journal of Physical Chemistry A</i> , 2003, 107, 3525-3537.	2.5	58
124	Electric Field Control of Structure, Dimensionality, and Reactivity of Gold Nanoclusters on Metal-Supported MgO Films. <i>Physical Review Letters</i> , 2008, 100, 056102.	7.8	58
125	Temperature-Tunable Selective Methane Catalysis on Au_{20} : From Cryogenic Partial Oxidation Yielding Formaldehyde to Cold Ethylene Production. <i>Journal of Physical Chemistry C</i> , 2011, 115, 6788-6795.	3.1	57
126	Effect of Spatial Dispersion upon Physisorption Energies: He on Metals. <i>Physical Review Letters</i> , 1974, 33, 524-527.	7.8	56

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127	Nanotribology and the Stability of Nanostructures. Japanese Journal of Applied Physics, 1993, 32, 1444-1462.	1.5	56
128	Theoretical studies of the spectroscopy of excess electrons in water clusters. Journal of Chemical Physics, 1990, 93, 6226-6238.	3.0	55
129	Solid and liquid junctions. Computational Materials Science, 1992, 1, 1-24.	3.0	55
130	On Mesoscopic Forces and Quantized Conductance in Model Metallic Nanowires. Journal of Physical Chemistry B, 1997, 101, 5780-5783.	2.6	55
131	Structure and dynamics of surface crystallization of liquidn-alkanes. Physical Review B, 1993, 48, 11313-11316.	3.2	54
132	Multilayer Relaxation of Interlayer Registry and Spacing at High-Index Metal Surfaces. Physical Review Letters, 1983, 51, 1359-1361.	7.8	53
133	Excess electrons in polar molecular clusters. Journal of Chemical Physics, 1988, 88, 6670-6671.	3.0	53
134	Materials by numbers: Computations as tools of discovery. Proceedings of the National Academy of Sciences of the United States of America, 2005, 102, 6671-6678.	7.1	53
135	Molecular-dynamics studies of the growth modes and structure of amorphous silicon films via atom deposition. Physical Review B, 1989, 40, 11733-11746.	3.2	52
136	Formation and control of electron molecules in artificial atoms: Impurity and magnetic-field effects. Physical Review B, 2000, 61, 15895-15904.	3.2	52
137	Einfluss der geometrischen und elektronischen Struktur sowie der elementaren Zusammensetzung von Clustern auf chemische Prozesse in der Nanometerskala. Angewandte Chemie, 2003, 115, 1335-1338.	2.0	52
138	Fundamental Insight into the Substrate-Dependent Ripening of Monodisperse Clusters. ChemCatChem, 2013, 5, 3330-3341.	3.7	52
139	Excess electrons in ammonia clusters. Chemical Physics Letters, 1988, 148, 249-252.	2.6	51
140	Molecular-dynamics simulations of epitaxial crystal growth from the melt. II. Si(111). Physical Review B, 1988, 37, 4647-4655.	3.2	51
141	Multiply charged anionic metal clusters. Chemical Physics Letters, 1993, 210, 437-442.	2.6	51
142	Electronic Entropy, Shell Structure, and Size-Evolutionary Patterns of Metal Clusters. Physical Review Letters, 1997, 78, 1424-1427.	7.8	49
143	Stochastic theory of multistate diffusion in perfect and defective systems. I. Mathematical formalism. Physical Review B, 1979, 19, 6207-6219.	3.2	48
144	Stability and Collapse of Metallic Structures on Surfaces. Physical Review Letters, 1994, 73, 569-572.	7.8	48

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145	Photoabsorption Spectra of Nan+ Clusters: Thermal Line-Broadening Mechanisms. <i>Physical Review Letters</i> , 2001, 87, 053401.	7.8	48
146	Group theoretical analysis of symmetry breaking in two-dimensional quantum dots. <i>Physical Review B</i> , 2003, 68, .	3.2	48
147	The Missing Link: Au ₁₉₁ (SPh-tBu) ₆₆ Janus Nanoparticle with Molecular and Bulk-Metal-like Properties. <i>Journal of the American Chemical Society</i> , 2020, 142, 15799-15814.	13.7	48
148	Low-Temperature CO Oxidation Catalyzed by Free Palladium Clusters: Similarities and Differences to Pd Surfaces and Supported Particles. <i>ACS Catalysis</i> , 2015, 5, 2275-2289.	11.2	47
149	Epitaxial Crystallization from a Melt: A Surface Molecular-Dynamics Study. <i>Physical Review Letters</i> , 1980, 45, 2032-2035.	7.8	46
150	Dimensionality crossovers of the ĩf plasmon in coaxial carbon nanotubes. <i>Physical Review B</i> , 1994, 50, 7977-7980.	3.2	45
151	Dynamical simulations of stress, strain, and finite deformations. <i>Physical Review B</i> , 1988, 38, 9522-9537.	3.2	44
152	Dynamics of Electron Localization, Solvation, and Migration in Polar Molecular Clusters. <i>Physical Review Letters</i> , 1989, 62, 106-109.	7.8	44
153	Crystalline Boson Phases in Harmonic Traps: Beyond the Gross-Pitaevskii Mean Field. <i>Physical Review Letters</i> , 2004, 93, 230405.	7.8	44
154	Adsorption on heterogeneous surfaces. I. Evaluation of the energy distribution function via the Wiener and Hopf method. <i>Journal of Chemical Physics</i> , 1976, 64, 1762-1767.	3.0	43
155	Further evaluation of the transform-deconvolution method for surface-structure determination by analysis of low-energy electron-diffraction intensities. <i>Physical Review B</i> , 1977, 15, 3775-3787.	3.2	43
156	Dielectrons in water clusters. <i>Journal of Chemical Physics</i> , 1992, 97, 1365-1377.	3.0	43
157	Oxidation State and Symmetry of Magnesia-Supported Pd ₁₃ O _x Nanocatalysts Influence Activation Barriers of CO Oxidation. <i>Journal of the American Chemical Society</i> , 2012, 134, 7690-7699.	13.7	43
158	Diffusion of Gold Clusters on Defective Graphite Surfaces. <i>Journal of Physical Chemistry B</i> , 2003, 107, 5882-5891.	2.6	42
159	Molecular Dynamics of a Laser-Annealing Experiment. <i>Physical Review Letters</i> , 1982, 49, 790-793.	7.8	41
160	Structure, collective hydrogen transfer, and formation of Si(OH) ₄ in SiO ₂ •(H ₂ O) _n clusters. <i>Journal of Chemical Physics</i> , 2002, 116, 9300-9304.	3.0	41
161	Microscopic theory of thermal desorption and dissociation processes catalyzed by a solid surface. <i>Physical Review B</i> , 1980, 21, 3256-3268.	3.2	40
162	Size-Dependent Binding Energies of Methane to Small Gold Clusters. <i>ChemPhysChem</i> , 2010, 11, 1570-1577.	2.1	40

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163	Oxidation of Magnesia-Supported Pd ₃₀ Nanoclusters and Catalyzed CO Combustion: Size-Selected Experiments and First-Principles Theory. <i>Journal of Physical Chemistry C</i> , 2012, 116, 9594-9607.	3.1	40
164	Atomically Precise Silver Clusters as New SERS Substrates. <i>Journal of Physical Chemistry Letters</i> , 2013, 4, 2769-2773.	4.6	40
165	Structural Variations and Multiple Charge Transfer Transitions between Chloranil and Carbazole Derivatives. <i>Macromolecules</i> , 1976, 9, 833-839.	4.8	39
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