

Charles T Campbell

List of Publications by Year in descending order

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14851
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#	ARTICLE	IF	CITATIONS
1	Influence of Adhesion on the Chemical Potential of Supported Nanoparticles as Modeled with Spherical Caps. ACS Catalysis, 2022, 12, 2302-2308.	11.2	9
2	Size-Dependent Adsorption and Adhesion Energetics of Ag Nanoparticles on Graphene Films on Ni(111) by Calorimetry. ACS Catalysis, 2022, 12, 2888-2897.	11.2	7
3	Acetonitrile Adsorption and Adhesion Energies onto the Pt(111) Surface by Calorimetry. ACS Catalysis, 2022, 12, 156-163.	11.2	2
4	Energetics of Ag Adsorption on and Adhesion to Rutile TiO ₂ (100) Studied by Microcalorimetry. Journal of Physical Chemistry C, 2021, 125, 3036-3046.	3.1	8
5	Predicting a Key Catalyst-Performance Descriptor for Supported Metal Nanoparticles: Metal Chemical Potential. ACS Catalysis, 2021, 11, 8284-8291.	11.2	25
6	Nature of the Active Sites on Ni/CeO ₂ Catalysts for Methane Conversions. ACS Catalysis, 2021, 11, 10604-10613.	11.2	37
7	Introduction: Advanced Materials and Methods for Catalysis and Electrocatalysis by Transition Metals. Chemical Reviews, 2021, 121, 563-566.	47.7	33
8	Analysis and prediction of reaction kinetics using the degree of rate control. Journal of Catalysis, 2021, 404, 647-660.	6.2	9
9	Effects of Solvents on Adsorption Energies: A General Bond-Additivity Model. Journal of Physical Chemistry C, 2021, 125, 24371-24380.	3.1	14
10	Adhesion Energies of Liquid Hydrocarbon Solvents onto Pt(111), MgO(100), Graphene, and TiO ₂ (110) from Temperature-Programmed Desorption Energies. Journal of Physical Chemistry C, 2021, 125, 27931-27937.	3.1	3
11	The degree of rate control of catalyst-bound intermediates in catalytic reaction mechanisms: Relationship to site coverage. Journal of Catalysis, 2020, 381, 53-62.	6.2	25
12	Silver Adsorption on Calcium Niobate(001) Nanosheets: Calorimetric Energies That Explain Sinter-Resistant Support. Journal of the American Chemical Society, 2020, 142, 15751-15763.	13.7	4
13	Calorimetric metal vapor adsorption energies for characterizing industrial catalyst support materials. Journal of Catalysis, 2020, 392, 209-216.	6.2	2
14	Ni Nanoparticles on CeO ₂ (111): Energetics, Electron Transfer, and Structure by Ni Adsorption Calorimetry, Spectroscopies, and Density Functional Theory. ACS Catalysis, 2020, 10, 5101-5114.	11.2	42
15	Energetics and Structure of Nickel Atoms and Nanoparticles on MgO(100). Journal of Physical Chemistry C, 2020, 124, 14685-14695.	3.1	10
16	Kinetic Isotope Effects: Interpretation and Prediction Using Degrees of Rate Control. ACS Catalysis, 2020, 10, 4181-4192.	11.2	24
17	Enhanced Bonding of Pentagon-Heptagon Defects in Graphene to Metal Surfaces: Insights from the Adsorption of Azulene and Naphthalene to Pt(111). Chemistry of Materials, 2020, 32, 1041-1053.	6.7	20
18	Aqueous phase catalytic and electrocatalytic hydrogenation of phenol and benzaldehyde over platinum group metals. Journal of Catalysis, 2020, 382, 372-384.	6.2	68

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19	Catalytic properties of model supported nanoparticles. <i>Journal of Chemical Physics</i> , 2020, 152, 140401.	3.0	3
20	Energetics of Adsorbed Phenol on Ni(111) and Pt(111) by Calorimetry. <i>Journal of Physical Chemistry C</i> , 2019, 123, 7627-7632.	3.1	32
21	A Simple Bond-Additivity Model Explains Large Decreases in Heats of Adsorption in Solvents Versus Gas Phase: A Case Study with Phenol on Pt(111) in Water. <i>ACS Catalysis</i> , 2019, 9, 8116-8127.	11.2	52
22	The kinetics of elementary thermal reactions in heterogeneous catalysis. <i>Nature Reviews Chemistry</i> , 2019, 3, 723-732.	30.2	31
23	Adhesion Energies of Solvent Films to Pt(111) and Ni(111) Surfaces by Adsorption Calorimetry. <i>ACS Catalysis</i> , 2019, 9, 11819-11825.	11.2	14
24	Apparent Activation Energies in Complex Reaction Mechanisms: A Simple Relationship via Degrees of Rate Control. <i>ACS Catalysis</i> , 2019, 9, 9465-9473.	11.2	71
25	Quantifying Adsorption of Organic Molecules on Platinum in Aqueous Phase by Hydrogen Site Blocking and in Situ X-ray Absorption Spectroscopy. <i>ACS Catalysis</i> , 2019, 9, 6869-6881.	11.2	40
26	Energies of Adsorbed Catalytic Intermediates on Transition Metal Surfaces: Calorimetric Measurements and Benchmarks for Theory. <i>Accounts of Chemical Research</i> , 2019, 52, 984-993.	15.6	38
27	Origin of Thermal and Hyperthermal CO ₂ from CO Oxidation on Pt Surfaces: The Role of Post-Transition State Dynamics, Active Sites, and Chemisorbed CO ₂ . <i>Angewandte Chemie</i> , 2019, 131, 6990-6994.	2.0	7
28	Origin of Thermal and Hyperthermal CO ₂ from CO Oxidation on Pt Surfaces: The Role of Post-Transition State Dynamics, Active Sites, and Chemisorbed CO ₂ . <i>Angewandte Chemie - International Edition</i> , 2019, 58, 6916-6920.	13.8	31
29	Energetics of Au Adsorption and Film Growth on Pt(111) by Single-Crystal Adsorption Calorimetry. <i>Journal of Physical Chemistry C</i> , 2019, 123, 5557-5561.	3.1	4
30	Heats of Adsorption of N ₂ , CO, Ar, and CH ₄ versus Coverage on the Zr-Based MOF NU-1000: Measurements and DFT Calculations. <i>Journal of Physical Chemistry C</i> , 2019, 123, 6586-6591.	3.1	8
31	Impact of pH on Aqueous-Phase Phenol Hydrogenation Catalyzed by Carbon-Supported Pt and Rh. <i>ACS Catalysis</i> , 2019, 9, 1120-1128.	11.2	55
32	Energetics of adsorbed benzene on Ni(111) and Pt(111) by calorimetry. <i>Surface Science</i> , 2018, 676, 9-16.	1.9	26
33	Adsorbed Hydroxyl and Water on Ni(111): Heats of Formation by Calorimetry. <i>ACS Catalysis</i> , 2018, 8, 1485-1489.	11.2	29
34	Energetics of van der Waals Adsorption on the Metal-Organic Framework NU-1000 with Zr ₆ -oxo, Hydroxo, and Aqua Nodes. <i>Journal of the American Chemical Society</i> , 2018, 140, 328-338.	13.7	11
35	Bond Energies of Adsorbed Intermediates to Metal Surfaces: Correlation with Hydrogen-Ligand and Hydrogen-Surface Bond Energies and Electronegativities. <i>Angewandte Chemie</i> , 2018, 130, 17119-17123.	2.0	2
36	Carbon-supported Pt during aqueous phenol hydrogenation with and without applied electrical potential: X-ray absorption and theoretical studies of structure and adsorbates. <i>Journal of Catalysis</i> , 2018, 368, 8-19.	6.2	49

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37	Bond Energies of Adsorbed Intermediates to Metal Surfaces: Correlation with Hydrogenâ€“Ligand and Hydrogenâ€“Surface Bond Energies and Electronegativities. <i>Angewandte Chemie - International Edition</i> , 2018, 57, 16877-16881.	13.8	7
38	Energetics of Adsorbed Methanol and Methoxy on Ni(111): Comparisons to Pt(111). <i>ACS Catalysis</i> , 2018, 8, 10089-10095.	11.2	7
39	The physical chemistry and materials science behind sinter-resistant catalysts. <i>Chemical Society Reviews</i> , 2018, 47, 4314-4331.	38.1	236
40	Velocity-resolved kinetics of site-specific carbon monoxide oxidation on platinum surfaces. <i>Nature</i> , 2018, 558, 280-283.	27.8	92
41	Energetics of 2D and 3D Gold Nanoparticles on MgO(100): Influence of Particle Size and Defects on Gold Adsorption and Adhesion Energies. <i>ACS Catalysis</i> , 2017, 7, 2151-2163.	11.2	14
42	DFT-Based Method for More Accurate Adsorption Energies: An Adaptive Sum of Energies from RPBE and vdW Density Functionals. <i>Journal of Physical Chemistry C</i> , 2017, 121, 4937-4945.	3.1	80
43	The Degree of Rate Control: A Powerful Tool for Catalysis Research. <i>ACS Catalysis</i> , 2017, 7, 2770-2779.	11.2	327
44	Energetics of Adsorbed Methyl and Methyl Iodide on Ni(111) by Calorimetry: Comparison to Pt(111) and Implications for Catalysis. <i>ACS Catalysis</i> , 2017, 7, 1286-1294.	11.2	20
45	Trends in Adhesion Energies of Metal Nanoparticles on Oxide Surfaces: Understanding Support Effects in Catalysis and Nanotechnology. <i>ACS Nano</i> , 2017, 11, 1196-1203.	14.6	121
46	Direct Measurements of Half-Cycle Reaction Heats during Atomic Layer Deposition by Calorimetry. <i>Chemistry of Materials</i> , 2017, 29, 8566-8577.	6.7	33
47	Formic Acid Dissociative Adsorption on NiO(111): Energetics and Structure of Adsorbed Formate. <i>Journal of Physical Chemistry C</i> , 2017, 121, 28001-28006.	3.1	9
48	Chemical Potential of Metal Atoms in Supported Nanoparticles: Dependence upon Particle Size and Support. <i>ACS Catalysis</i> , 2017, 7, 8460-8466.	11.2	88
49	Energetics of adsorbed formate and formic acid on Ni(111) by calorimetry. <i>Journal of Catalysis</i> , 2017, 352, 300-304.	6.2	21
50	Calorimetric measurement of adsorption and adhesion energies of Cu on Pt(111). <i>Surface Science</i> , 2017, 657, 58-62.	1.9	5
51	Calcium Vapor Adsorption on the Metalâ€“Organic Framework NU-1000: Structure and Energetics. <i>Journal of Physical Chemistry C</i> , 2016, 120, 16850-16862.	3.1	16
52	Equilibrium Constants and Rate Constants for Adsorbates: Two-Dimensional (2D) Ideal Gas, 2D Ideal Lattice Gas, and Ideal Hindered Translator Models. <i>Journal of Physical Chemistry C</i> , 2016, 120, 10283-10297.	3.1	94
53	Electrocatalytic Hydrogenation of Phenol over Platinum and Rhodium: Unexpected Temperature Effects Resolved. <i>ACS Catalysis</i> , 2016, 6, 7466-7470.	11.2	86
54	Reply to â€œComment on â€“Equilibrium Constants and Rate Constants for Adsorbates: Two-Dimensional (2D) Ideal Gas, 2D Ideal Lattice Gas, and Ideal Hindered Translator Modelsâ€™â€–. <i>Journal of Physical Chemistry C</i> , 2016, 120, 20481-20482.	3.1	5

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55	Energies of Formation Reactions Measured for Adsorbates on Late Transition Metal Surfaces. Journal of Physical Chemistry C, 2016, 120, 25161-25172.	3.1	63
56	Water Dissociative Adsorption on NiO(111): Energetics and Structure of the Hydroxylated Surface. ACS Catalysis, 2016, 6, 7377-7384.	11.2	67
57	Catalysis: Quantifying charge transfer. Nature Energy, 2016, 1, .	39.5	5
58	Adsorption and Adhesion of Au on Reduced CeO ₂ (111) Surfaces at 300 and 100 K. Journal of Physical Chemistry C, 2016, 120, 12113-12124.	3.1	29
59	Hindered Translator and Hindered Rotor Models for Adsorbates: Partition Functions and Entropies. Journal of Physical Chemistry C, 2016, 120, 9719-9731.	3.1	113
60	Energetics of methanol and formic acid oxidation on Pt(111): Mechanistic insights from adsorption calorimetry. Surface Science, 2016, 650, 140-143.	1.9	17
61	Toward Benchmarking in Catalysis Science: Best Practices, Challenges, and Opportunities. ACS Catalysis, 2016, 6, 2590-2602.	11.2	190
62	Using degrees of rate control to improve selective n-butane oxidation over model MOF-encapsulated catalysts: sterically-constrained Ag ₃ Pd(111). Faraday Discussions, 2016, 188, 21-38.	3.2	15
63	Ion scattering spectroscopy intensities for supported nanoparticles: The hemispherical cap model. Surface Science, 2015, 641, 166-169.	1.9	15
64	Surface-Bound Intermediates in Low-Temperature Methanol Synthesis on Copper: Participants and Spectators. ACS Catalysis, 2015, 5, 7328-7337.	11.2	77
65	Energetics of Cu Adsorption and Adhesion onto Reduced CeO ₂ (111) Surfaces by Calorimetry. Journal of Physical Chemistry C, 2015, 119, 17209-17217.	3.1	47
66	Degree of rate control approach to computational catalyst screening. Journal of Catalysis, 2015, 330, 197-207.	6.2	105
67	A benchmark database for adsorption bond energies to transition metal surfaces and comparison to selected DFT functionals. Surface Science, 2015, 640, 36-44.	1.9	396
68	Calcium Thin Film Growth on Phenyl-C ₆₁ -Butyric Acid Methyl Ester (PCBM): Interface Structure and Energetics. Journal of Physical Chemistry C, 2015, 119, 18444-18451.	3.1	13
69	Energy of Supported Metal Catalysts: From Single Atoms to Large Metal Nanoparticles. ACS Catalysis, 2015, 5, 5673-5678.	11.2	78
70	Quantitative modeling of electron spectroscopy intensities for supported nanoparticles: The hemispherical cap model for non-normal detection. Surface Science, 2015, 632, L5-L8.	1.9	6
71	Method for direct deconvolution of heat signals in transient adsorption calorimetry. Surface Science, 2015, 633, 17-23.	1.9	3
72	Energetics of Adsorbed CH ₂ and CH on Pt(111) by Calorimetry: The Dissociative Adsorption of Diiodomethane. Journal of Physical Chemistry C, 2014, 118, 29310-29321.	3.1	13

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73	Adsorption Energy of <i>tert</i> -Butyl on Pt(111) by Dissociation of <i>tert</i> -Butyl Iodide: Calorimetry and DFT. <i>Journal of Physical Chemistry C</i> , 2014, 118, 427-438.	3.1	22
74	Low-Temperature Growth Improves Metal/Polymer Interfaces: Vapor-Deposited Ca on PMMA. <i>Journal of Physical Chemistry C</i> , 2014, 118, 6352-6358.	3.1	3
75	Bond Energies of Molecular Fragments to Metal Surfaces Track Their Bond Energies to H Atoms. <i>Journal of the American Chemical Society</i> , 2014, 136, 4137-4140.	13.7	25
76	Energetics of Formic Acid Conversion to Adsorbed Formates on Pt(111) by Transient Calorimetry. <i>Journal of the American Chemical Society</i> , 2014, 136, 3964-3971.	13.7	44
77	Anchored metal nanoparticles: Effects of support and size on their energy, sintering resistance and reactivity. <i>Faraday Discussions</i> , 2013, 162, 9.	3.2	161
78	D. W. (â€œWayneâ€) Goodman: A Pioneer in Elucidating the Relationships Between Surface Structure of Catalysts and Their Performance, and in Using Model Catalysts for That Purpose. <i>Topics in Catalysis</i> , 2013, 56, 1273-1276.	2.8	4
79	Silver Nanoparticles on Fe ₃ O ₄ (111): Energetics by Ag Adsorption Calorimetry and Structure by Surface Spectroscopies. <i>Journal of Physical Chemistry C</i> , 2013, 117, 24932-24936.	3.1	23
80	Kinetic Prefactors of Reactions on Solid Surfaces. <i>Zeitschrift Fur Physikalische Chemie</i> , 2013, 227, .	2.8	81
81	A Sinter-Resistant Catalytic System Fabricated by Maneuvering the Selectivity of SiO ₂ Deposition onto the TiO ₂ Surface versus the Pt Nanoparticle Surface. <i>Nano Letters</i> , 2013, 13, 4957-4962.	9.1	101
82	Energetics of Adsorbed CH ₃ on Pt(111) by Calorimetry. <i>Journal of the American Chemical Society</i> , 2013, 135, 5208-5211.	13.7	33
83	Surface kinetics and energetics from single crystal adsorption calorimetry lineshape analysis: Methyl from methyl iodide on Pt(111). <i>Journal of Catalysis</i> , 2013, 308, 114-121.	6.2	11
84	Enthalpies and Entropies of Adsorption on Well-Defined Oxide Surfaces: Experimental Measurements. <i>Chemical Reviews</i> , 2013, 113, 4106-4135.	47.7	211
85	The Energetics of Supported Metal Nanoparticles: Relationships to Sintering Rates and Catalytic Activity. <i>Accounts of Chemical Research</i> , 2013, 46, 1712-1719.	15.6	300
86	Energetics of Adsorbed CH ₃ and CH on Pt(111) by Calorimetry: Dissociative Adsorption of CH ₃ I. <i>Journal of Physical Chemistry C</i> , 2013, 117, 6325-6336.	3.1	24
87	Adsorption calorimetry during metal vapor deposition on single crystal surfaces: Increased flux, reduced optical radiation, and real-time flux and reflectivity measurements. <i>Review of Scientific Instruments</i> , 2013, 84, 123901.	1.3	21
88	Kinetic Prefactors of Reactions on Solid Surfaces. <i>Zeitschrift Fur Physikalische Chemie</i> , 2013, .	2.8	5
89	Energetics of Oxygen Adatoms, Hydroxyl Species and Water Dissociation on Pt(111). <i>Journal of Physical Chemistry C</i> , 2012, 116, 25772-25776.	3.1	62
90	The Entropies of Adsorbed Molecules. <i>Journal of the American Chemical Society</i> , 2012, 134, 18109-18115.	13.7	364

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91	Ca Carboxylate Formation at the Calcium/Poly(methyl methacrylate) Interface. <i>Journal of Physical Chemistry C</i> , 2012, 116, 20465-20471.	3.1	31
92	A Highly Reactive and Sinter-Resistant Catalytic System Based on Platinum Nanoparticles Embedded in the Inner Surfaces of CeO ₂ Hollow Fibers. <i>Angewandte Chemie - International Edition</i> , 2012, 51, 9543-9546.	13.8	121
93	Energetics of Adsorbed Methanol and Methoxy on Pt(111) by Microcalorimetry. <i>Journal of the American Chemical Society</i> , 2012, 134, 20388-20395.	13.7	70
94	Electronic perturbations. <i>Nature Chemistry</i> , 2012, 4, 597-598.	13.6	610
95	Built-In Potential in Conjugated Polymer Diodes with Changing Anode Work Function: Interfacial States and Deviation from the Schottky-Mott Limit. <i>Journal of Physical Chemistry Letters</i> , 2012, 3, 1202-1207.	4.6	50
96	Energy of Molecularly Adsorbed Water on Clean Pt(111) and Pt(111) with Coadsorbed Oxygen by Calorimetry. <i>Journal of Physical Chemistry C</i> , 2011, 115, 9164-9170.	3.1	61
97	The Energy of Adsorbed Hydroxyl on Pt(111) by Microcalorimetry. <i>Journal of Physical Chemistry C</i> , 2011, 115, 11586-11594.	3.1	47
98	Insights into catalysis by gold nanoparticles and their support effects through surface science studies of model catalysts. <i>Faraday Discussions</i> , 2011, 152, 227.	3.2	78
99	Growth, Structure, and Stability of Ag on CeO ₂ (111): Synchrotron Radiation Photoemission Studies. <i>Journal of Physical Chemistry C</i> , 2011, 115, 6715-6725.	3.1	78
100	The Energy of Hydroxyl Coadsorbed with Water on Pt(111). <i>Journal of Physical Chemistry C</i> , 2011, 115, 23008-23012.	3.1	45
101	Adsorption Microcalorimetry: Recent Advances in Instrumentation and Application. <i>Annual Review of Analytical Chemistry</i> , 2011, 4, 41-58.	5.4	26
102	An improved single crystal adsorption calorimeter for determining gas adsorption and reaction energies on complex model catalysts. <i>Review of Scientific Instruments</i> , 2011, 82, 024102.	1.3	58
103	A Sinter-Resistant Catalytic System Based on Platinum Nanoparticles Supported on TiO ₂ Nanofibers and Covered by Porous Silica. <i>Angewandte Chemie - International Edition</i> , 2010, 49, 8165-8168.	13.8	125
104	Improved pyroelectric detectors for single crystal adsorption calorimetry from 100 to 350 K. <i>Review of Scientific Instruments</i> , 2010, 81, 024102.	1.3	54
105	Ag Adsorption on Reduced CeO ₂ (111) Thin Films. <i>Journal of Physical Chemistry C</i> , 2010, 114, 17166-17172.	3.1	67
106	Ceria Maintains Smaller Metal Catalyst Particles by Strong Metal-Support Bonding. <i>Science</i> , 2010, 329, 933-936.	12.6	763
107	Particle-size dependent heats of adsorption of CO on supported Pd nanoparticles as measured with a single-crystal microcalorimeter. <i>Physical Review B</i> , 2010, 81, .	3.2	77
108	Degree of Rate Control: How Much the Energies of Intermediates and Transition States Control Rates. <i>Journal of the American Chemical Society</i> , 2009, 131, 8077-8082.	13.7	461

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109	Catalytic reaction energetics by single crystal adsorption calorimetry: hydrocarbons on Pt(111). <i>Chemical Society Reviews</i> , 2008, 37, 2172.	38.1	46
110	Calcium Adsorption on MgO(100): Energetics, Structure, and Role of Defects. <i>Journal of the American Chemical Society</i> , 2008, 130, 2314-2322.	13.7	45
111	Energetics of Cyclohexene Adsorption and Reaction on Pt(111) by Low-Temperature Microcalorimetry. <i>Journal of the American Chemical Society</i> , 2008, 130, 10247-10257.	13.7	65
112	Kinetic model for sintering of supported metal particles with improved size-dependent energetics and applications to Au on TiO ₂ (110). <i>Physical Review B</i> , 2007, 75, .	3.2	136
113	SPR microscopy and its applications to high-throughput analyses of biomolecular binding events and their kinetics. <i>Biomaterials</i> , 2007, 28, 2380-2392.	11.4	367
114	Reactivity and sintering kinetics of Au/TiO ₂ (110) model catalysts: particle size effects. <i>Topics in Catalysis</i> , 2007, 44, 3-13.	2.8	74
115	Thermodynamics of Statherin Adsorption onto Hydroxyapatite. <i>Biochemistry</i> , 2006, 45, 5576-5586.	2.5	74
116	Heat of Adsorption of Naphthalene on Pt(111) Measured by Adsorption Calorimetry. <i>Journal of Physical Chemistry B</i> , 2006, 110, 17539-17545.	2.6	73
117	n-alkanes on Pt(111) and on C(0001)-Pt(111): Chain length dependence of kinetic desorption parameters. <i>Journal of Chemical Physics</i> , 2006, 125, 234308.	3.0	170
118	Transition Metal Oxides: Extra Thermodynamic Stability as Thin Films. <i>Physical Review Letters</i> , 2006, 96, 066106.	7.8	78
119	Pyroelectric heat detector for measuring adsorption energies on thicker single crystals. <i>Sensors and Actuators B: Chemical</i> , 2005, 107, 454-460.	7.8	19
120	Heats of adsorption of Pb on pristine and electron-irradiated poly(methyl methacrylate) by microcalorimetry. <i>Surface Science</i> , 2005, 598, 22-34.	1.9	20
121	n-alkanes on MgO(100). I. Coverage-dependent desorption kinetics of n-butane. <i>Journal of Chemical Physics</i> , 2005, 122, 164707.	3.0	120
122	n-alkanes on MgO(100). II. Chain length dependence of kinetic desorption parameters for small n-alkanes. <i>Journal of Chemical Physics</i> , 2005, 122, 164708.	3.0	156
123	Small Pd Clusters, up to the Tetramer At Least, Are Highly Mobile on the MgO(100) Surface. <i>Physical Review Letters</i> , 2005, 95, 146103.	7.8	87
124	CHEMISTRY: Oxygen Vacancies and Catalysis on Ceria Surfaces. <i>Science</i> , 2005, 309, 713-714.	12.6	1,103
125	Calorimeter for adsorption energies of larger molecules on single crystal surfaces. <i>Review of Scientific Instruments</i> , 2004, 75, 4471-4480.	1.3	60
126	Calorimetric Measurement of the Heat of Adsorption of Benzene on Pt(111). <i>Journal of Physical Chemistry B</i> , 2004, 108, 14627-14633.	2.6	130

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127	PHYSICS: The Active Site in Nanoparticle Gold Catalysis. <i>Science</i> , 2004, 306, 234-235.	12.6	327
128	Cyclohexane Dehydrogenation and H ₂ Adsorption on Pt Particles on ZnO(0001)~O. <i>Journal of Physical Chemistry B</i> , 2003, 107, 1180-1188.	2.6	20
129	Benzene Adsorption and Dehydrogenation on Pt/ZnO(0001)~O Model Catalysts. <i>Journal of Physical Chemistry B</i> , 2003, 107, 1174-1179.	2.6	12
130	SURFACE SCIENCE: Enhanced: Waltzing with O ₂ . <i>Science</i> , 2003, 299, 357-357.	12.6	21
131	Metal Adsorption and Adhesion Energies on MgO(100). <i>Journal of the American Chemical Society</i> , 2002, 124, 9212-9218.	13.7	79
132	The Effect of Size-Dependent Nanoparticle Energetics on Catalyst Sintering. <i>Science</i> , 2002, 298, 811-814.	12.6	907
133	Methanol Decomposition on Pt/ZnO(0001)~Zn Model Catalysts. <i>Journal of Physical Chemistry B</i> , 2001, 105, 9273-9279.	2.6	26
134	Enthalpies of adsorption of metal atoms on single-crystalline surfaces by microcalorimetry. <i>Journal of Chemical Thermodynamics</i> , 2001, 33, 333-345.	2.0	18
135	Finding the Rate-Determining Step in a Mechanism. <i>Journal of Catalysis</i> , 2001, 204, 520-524.	6.2	255
136	Pyroelectric detector for single-crystal adsorption microcalorimetry: analysis of pulse shape and intensity. <i>Sensors and Actuators B: Chemical</i> , 2000, 62, 13-22.	7.8	33
137	Title is missing!. <i>Topics in Catalysis</i> , 2000, 14, 43-51.	2.8	67
138	The influence of chlorine on the dispersion of Cu particles on Cu/ZnO(0001) model catalysts. <i>Catalysis Letters</i> , 2000, 65, 159-168.	2.6	14
139	Quantification of Tight Binding to Surface-Immobilized Phospholipid Vesicles Using Surface Plasmon Resonance:~ Binding Constant of Phospholipase A ₂ . <i>Journal of the American Chemical Society</i> , 2000, 122, 4177-4184.	13.7	100
140	Surface Characterization of Hydroxyapatite and Related Calcium Phosphates by XPS and TOF-SIMS. <i>Analytical Chemistry</i> , 2000, 72, 2886-2894.	6.5	300
141	Sticking Probabilities in Adsorption of Alkanethiols from Liquid Ethanol Solution onto Gold. <i>Journal of Physical Chemistry B</i> , 2000, 104, 11168-11178.	2.6	107
142	The kinetics of CO oxidation by adsorbed oxygen on well~defined gold particles on TiO ₂ (110). <i>Catalysis Letters</i> , 1999, 63, 143-151.	2.6	203
143	A microcalorimetric study of the heat of adsorption of copper on well-defined oxide thin film surfaces: MgO(100), p(2~1) oxide on Mo(100) and disordered W oxide. <i>Faraday Discussions</i> , 1999, 114, 195-208.	3.2	51
144	Organofunctionalization of TiO ₂ (110):~ (3,3,3-Trifluoropropyl)trimethoxysilane Adsorption. <i>Journal of Physical Chemistry B</i> , 1998, 102, 4536-4543.	2.6	30

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146	Metal adsorption calorimetry and adhesion energies on clean single-crystal surfaces. Journal of Chemical Physics, 1997, 107, 5547-5553.	3.0	57
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