Wilson Agerico Dio

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#	Paper	IF	Citations
200	First principles studies for the dissociative adsorption of H2 on graphene. <i>Journal of Applied Physics</i> , 2003 , 93, 3395-3400	2.5	129
199	Orientational effects in dissociative adsorption/associative desorption dynamics of H2(D2) on Cu and Pd. <i>Progress in Surface Science</i> , 2000 , 63, 63-134	6.6	96
198	H2 dissociative adsorption on Mg, Ti, Ni, Pd and La Surfaces. <i>Surface Science</i> , 2004 , 566-568, 703-707	1.8	61
197	Effective Pathway for Hydrogen Atom Adsorption on Graphene. <i>Journal of the Physical Society of Japan</i> , 2003 , 72, 995-997	1.5	61
196	Hydrogen pairing on graphene. <i>Carbon</i> , 2007 , 45, 218-220	10.4	57
195	Amino acid adsorption on single-walled carbon nanotubes. <i>European Physical Journal D</i> , 2006 , 38, 117-	1203	52
194	Potential energy of hydrogen atom motion on Pd(111) surface and in subsurface: A first principles calculation. <i>Journal of Applied Physics</i> , 2007 , 101, 123530	2.5	46
193	Dissociation and Sticking of H2on Mg(0001), Ti(0001) and La(0001) Surfaces. <i>Journal of the Physical Society of Japan</i> , 2004 , 73, 745-748	1.5	41
192	Photostimulated desorption and ortho-para conversion of H2 on Ag surfaces. <i>Physical Review Letters</i> , 2003 , 90, 096103	7.4	38
191	Effect of oxygen vacancy on the adsorption of O2 on anatase TiO2(001): A DFT-based study. <i>Surface Science</i> , 2015 , 633, 38-45	1.8	35
190	Role of Rotational Motion in the Dissociative Adsorption and Associative Desorption Dynamics of D2/Cu(111). <i>Physical Review Letters</i> , 1997 , 78, 286-289	7.4	34
189	Mg⊞ dissociation of magnesium hydride MgH2 catalyzed by 3d transition metals. <i>Thin Solid Films</i> , 2006 , 509, 157-159	2.2	31
188	Glycine adsorption on single-walled carbon nanotubes. <i>Thin Solid Films</i> , 2006 , 509, 218-222	2.2	31
187	H2 dissociative adsorption at the armchair edges of graphite. <i>Solid State Communications</i> , 2004 , 132, 713-718	1.6	31
186	Many body effects in elementary processes at metal surfaces. Surface Science Reports, 2001, 43, 1-43	12.9	31
185	Spatial and spectroscopic profiles of the Kondo resonance for magnetic atoms on metal surfaces. Journal of Applied Physics, 1999 , 86, 6970-6974	2.5	31
184	Potential Energy of H2Dissociation and Adsorption on Pt(111) Surface: First-Principles Calculation. Japanese Journal of Applied Physics, 2007 , 46, 4233-4237	1.4	28

183	Effects of Rotational State Excitations on the Dissociative Adsorption Dynamics of D2/Cu(111). Journal of the Physical Society of Japan, 1995 , 64, 2478-2487	1.5	27	
182	Orientation dependence of O2 dissociation from heme © 2 adduct. <i>Chemical Physics Letters</i> , 2005 , 402, 71-74	2.5	26	
181	Theoretical study of hydrazine adsorption on Pt(111): Anti or cis?. Surface Science, 2011, 605, 1347-1353	3 1.8	23	
180	Rotational and Vibrational Coupling Effects on the Dissociative Adsorption and Associative Desorption Dynamics of D2/Cu(111). <i>Journal of the Physical Society of Japan</i> , 1999 , 68, 887-892	1.5	23	
179	H2 Dissociative Adsorption at the Zigzag Edges of Graphite. <i>E-Journal of Surface Science and Nanotechnology</i> , 2004 , 2, 77-80	0.7	23	
178	Realizing a Carbon-Based Hydrogen Storage Material. <i>Japanese Journal of Applied Physics</i> , 2006 , 45, 176	6 51 76	722	
177	Dependence of oxygen dissociative adsorption on platinum surface structures. <i>Physical Review B</i> , 2005 , 72,	3.3	22	
176	Dynamical quantum filtering in hydrogenBurface reactions. <i>Surface Science</i> , 1998 , 418, L39-L44	1.8	21	
175	Dissociative adsorption of O2 on Pt and Au surfaces: Potential-energy surfaces and electronic states. <i>Physical Review B</i> , 2008 , 77,	3.3	21	
174	Ab InitioStudy of H2Desorption from Magnesium Hydride MgH2Cluster. <i>Journal of the Physical Society of Japan</i> , 2004 , 73, 2628-2630	1.5	20	
173	Hydrogen adsorption on Ag and Au monolayers grown on Pt(1 1 1). Surface Science, 2004, 566-568, 755	-71680	20	
172	Surface science-based reaction design: increasing the orthopara hydrogen conversion yield via molecular orientation, a case study. <i>Progress in Surface Science</i> , 2003 , 72, 53-86	6.6	20	
171	Effects of Correlation between Molecular Diffraction and Rotational Excitation on the Scattering Dynamics of H2from Cu(001). <i>Journal of the Physical Society of Japan</i> , 2000 , 69, 3878-3884	1.5	20	
170	Stable Hydrogen Configurations between Graphite Layers. <i>Journal of the Physical Society of Japan</i> , 2003 , 72, 1867-1870	1.5	19	
169	Vibrational properties of hydrogen atom adsorbed on Cu(111) and on Ir(111) surfaces. <i>Journal of Applied Physics</i> , 2004 , 96, 5020-5025	2.5	19	
168	Rotational Alignment in the Associative Desorption Dynamics of Hydrogen Molecules from Metal Surfaces. <i>Journal of the Physical Society of Japan</i> , 1998 , 67, 1517-1520	1.5	18	
167	Experimental and Theoretical Studies on Oxidation of Cu-Au Alloy Surfaces: Effect of Bulk Au Concentration. <i>Scientific Reports</i> , 2016 , 6, 31101	4.9	18	
166	Oxygen vacancy effects on electronic structure of Pt/NiO/Pt capacitor-like system. <i>Surface Science</i> , 2012 , 606, 239-246	1.8	17	

165	Scattering and dissociative adsorption of H2 on the armchair and zigzag edges of graphite. <i>Journal of Applied Physics</i> , 2004 , 96, 6331-6336	2.5	17	
164	Spectroscopic profiles of a magnetic dimer on a metal surface. <i>Solid State Communications</i> , 2009 , 149, 1241-1243	1.6	16	
163	Quantum states of hydrogen atom motion on the Pd(111) surface and in the subsurface. <i>Journal of Physics Condensed Matter</i> , 2007 , 19, 365214	1.8	16	
162	Quantum dynamics study on the interaction of H2 on a Pt(111) surface. <i>Thin Solid Films</i> , 2006 , 509, 227	-2 <u>29</u>	16	
161	Tuning of dissociative-adsorption processes on Cu{1 0 0} by controlling the kinetic energy of the impinging O2 molecule. <i>Chemical Physics</i> , 2004 , 301, 315-320	2.3	16	
160	Trapping hydrogen with a bimetallic interface. <i>Physical Review B</i> , 2005 , 71,	3.3	16	
159	Effect of antiferromagnetic RKKY interaction and magnetic field in a two-impurity Kondo system. <i>Physical Review B</i> , 2010 , 82,	3.3	15	
158	Magnetized/charged MgH2-based hydrogen storage materials. <i>Applied Physics Letters</i> , 2005 , 86, 21310	93.4	15	
157	Diameter Dependent Magnetic and Electronic Properties of Single-Walled Carbon Nanotubes with Fe Nanowires. <i>Japanese Journal of Applied Physics</i> , 2005 , 44, 882-888	1.4	15	
156	Combined effect of molecular rotational and surface vibrational excitations on the dissociative adsorption dynamics of. <i>Surface Science</i> , 1996 , 363, 52-61	1.8	15	
155	Hydrogen atom quantum migration on platinum. <i>E-Journal of Surface Science and Nanotechnology</i> , 2006 , 4, 619-623	0.7	15	
154	First Principles Studies on the Interaction of a Hydrogen Atom with a Single-Walled Carbon Nanotube. <i>Japanese Journal of Applied Physics</i> , 2003 , 42, 4626-4629	1.4	14	
153	Electric and Magnetic Properties of Co-filled Carbon Nanotube. <i>Journal of the Physical Society of Japan</i> , 2005 , 74, 742-745	1.5	14	
152	Scattering Dynamics of Hydrogen Molecules on Metal Alloy Surfaces P robing Local Surface Reactivity with Hydrogen Molecules <i>Journal of the Physical Society of Japan</i> , 2001 , 70, 3491-3494	1.5	14	
151	High-uptake graphene hydrogenation: a computational perspective. <i>Journal of Physics Condensed Matter</i> , 2009 , 21, 474219	1.8	13	
150	Effect of RKKY Interaction on the System of Two Magnetic Atoms on a Metal Surface at Finite Temperatures. <i>Journal of the Physical Society of Japan</i> , 2010 , 79, 074702	1.5	13	
149	Rotational Effects in the Dissociative Adsorption Dynamics of Hydrogen on a Pd Surface. <i>Journal of the Physical Society of Japan</i> , 1997 , 66, 3344-3347	1.5	13	
148	Cyclohexane dehydrogenation catalyst design based on spin polarization effects. <i>Journal of Physics Condensed Matter</i> , 2004 , 16, S5721-S5724	1.8	13	

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147	Reactivity of gold thin films grown on iridium: Hydrogen dissociation. <i>Applied Catalysis A: General</i> , 2005 , 291, 55-61	5.1	13
146	Reactive gold thin films grown on iridium. <i>Applied Surface Science</i> , 2005 , 246, 68-71	6.7	13
145	First Principles Studies for the Interaction of Hydrogen with a Li(100) Surface. <i>Journal of the Physical Society of Japan</i> , 2005 , 74, 478-482	1.5	13
144	Initial stages of Cu3Au(111) oxidation: oxygen induced Cu segregation and the protective Au layer profile. <i>Physical Chemistry Chemical Physics</i> , 2014 , 16, 3815-22	3.6	12
143	Conformational effects on hydrazine and OH coadsorption on Ni(111): A first-principles investigation. <i>Surface Science</i> , 2017 , 664, 185-193	1.8	12
142	Magnetic domain patterns on strong perpendicular magnetization of Co/Ni multilayers as spintronics materials: II. Numerical simulations. <i>Journal of Physics Condensed Matter</i> , 2013 , 25, 395005	1.8	12
141	Magnetic domain patterns on strong perpendicular magnetization of Co/Ni multilayers as spintronics materials: I. Dynamic observations. <i>Journal of Physics Condensed Matter</i> , 2013 , 25, 406001	1.8	12
140	Quantum states of a hydrogen atom adsorbed on Cu(100) and (110) surfaces. <i>Physical Review B</i> , 2007 , 75,	3.3	12
139	Spatial and temperature dependence of the spectroscopic profile of a magnetic atom adsorbed on a metal surface (Io/Cu(111). <i>Journal of Applied Physics</i> , 2003 , 94, 334-341	2.5	12
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137	Vibrational and rotational coupling effects in the direct scattering of H2 from Cu(111). <i>Surface Science</i> , 1999 , 438, 254-260	1.8	12
136	CO-induced Pd segregation and the effect of subsurface Pd on CO adsorption on CuPd surfaces. Journal of Physics Condensed Matter, 2017 , 29, 025005	1.8	11
135	Stable Structures of Fe Nanowires on Cu(111). Japanese Journal of Applied Physics, 2003, 42, 4633-4635	1.4	11
134	Ab InitioStudy of Cyclohexane Dehydrogenation with a Transition Metal (Pt, Pd, Ni and Cu) Atom. <i>Journal of the Physical Society of Japan</i> , 2004 , 73, 1281-1284	1.5	11
133	Electronic structure of Fe, Co, Ni nanowires on Cu(111). Surface Science, 2004, 566-568, 1052-1056	1.8	11
132	Behavior of hydrogen atom at Nafion B t interface. <i>Solid State Communications</i> , 2005 , 134, 601-605	1.6	11
131	Spin Polarization Effects on O2Dissociation from Heme-O2Adduct. <i>Japanese Journal of Applied Physics</i> , 2005 , 44, L57-L59	1.4	11
130	Effects of surface corrugation on the molecular rotational dependence of H2 dissociative adsorption dynamics on Cu(100). <i>Applied Surface Science</i> , 2001 , 169-170, 30-35	6.7	11

129	Behavior of a magnetic atom on a metal surface. <i>Journal of Electron Spectroscopy and Related Phenomena</i> , 2000 , 109, 63-70	1.7	11
128	A microscopic theory of STM-induced CO desorption from Cu(111). <i>Surface Science</i> , 1999 , 438, 283-288	1.8	11
127	Hemoglobin Components as Cathode Electrode Catalyst in Polymer Electrolyte Fuel Cells. <i>E-Journal of Surface Science and Nanotechnology</i> , 2004 , 2, 226-229	0.7	11
126	Enhanced molecular adsorption of ethylene on reduced anatase TiO2 (001): role of surface O-vacancies. <i>RSC Advances</i> , 2016 , 6, 92241-92251	3.7	11
125	Identifying Hydrogen Atoms on Graphite. Journal of the Physical Society of Japan, 2007, 76, 114703	1.5	10
124	H/Pd(111) Absorption and Desorption Dynamics -Hydrogen-Induced Lattice Relaxation Effects <i>Japanese Journal of Applied Physics</i> , 2003 , 42, 4630-4632	1.4	10
123	Spin Polarization of a Multiple-decked Sandwich Clusters: M(C6H6)2(M = Mn, Fe, Co). <i>Journal of the Physical Society of Japan</i> , 2004 , 73, 2292-2295	1.5	10
122	A theoretical analysis of quantum mirages on a Cu(111) surface. Surface Science, 2002, 514, 89-94	1.8	10
121	Photodesorption of hydrogen molecules physisorbed on Ag: Isotope dependence of translational-energy distribution. <i>Surface Science</i> , 2005 , 593, 229-234	1.8	10
120	Spatial and Spectroscopic Profiles of the Kondo Resonance for a Single Magnetic Atom on a Metal Surface under an External Magnetic Field. <i>Japanese Journal of Applied Physics</i> , 2000 , 39, 4359-4362	1.4	10
119	Molecular orientation dependence of o lp H 2 conversion on a 3d impurity sitting on a metal oxide surface. <i>Surface Science</i> , 2001 , 493, 285-291	1.8	10
118	Molecular Orientation Dependence ofoBConversion of H2Scattered from a 3dImpurity Sitting on a Metal Oxide Surface. <i>Journal of the Physical Society of Japan</i> , 2001 , 70, 3654-3659	1.5	10
117	Effect of RKKY Interaction on the Scanning Tunneling Spectra of a Classic Kondo System IIwo Magnetic Atoms Adsorbed on a Metal Surface IJournal of the Physical Society of Japan, 2009, 78, 084705	5 ^{1.5}	9
116	On the influence of incident angle in the scattering dynamics of D2 from NiAl(110). <i>Chemical Physics Letters</i> , 2002 , 359, 127-134	2.5	9
115	Reactive Ion Etching of NiFe Thin Films from First-Principles Study: A Case Study. <i>Japanese Journal of Applied Physics</i> , 2005 , 44, 893-894	1.4	9
114	Dissociative adsorption dynamics of H2 at the atop-Pt, atop-Cu, and Cu?Pt bridge sites of an ordered Cu3Pt(111) [brientational effects. <i>Applied Surface Science</i> , 2001 , 169-170, 36-41	6.7	9
113	Dynamical quantum filtering in the scattering dynamics of H2on Cu(001). <i>Journal of Physics Condensed Matter</i> , 2002 , 14, L479-L486	1.8	9
112	H 2 dissociation dynamics on an alloy surface Leontrolling the dynamics via orientation. <i>Surface Science</i> , 2001 , 493, 278-284	1.8	9

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111	Dissociative Adsorption Dynamics of H2(D2) on Metal Alloy Surfaces - A First Step Towards Surface Reaction Design/Control <i>Journal of the Physical Society of Japan</i> , 2000 , 69, 993-996	1.5	9
110	Cyclohexane Dehydrogenation Process Design Using Ni - Spin Polarization Effects <i>E-Journal of Surface Science and Nanotechnology</i> , 2004 , 2, 200-204	0.7	9
109	Effects of Adsorbates (CO, COH, and HCO) on the Arrangement of Pd Atoms in PdCu(111). <i>Journal of Physical Chemistry C</i> , 2017 , 121, 17818-17826	3.8	8
108	Scanning tunneling spectroscopic evidence of crossover transition in the two-impurity Kondo problem. <i>Surface Science</i> , 2010 , 604, 2139-2149	1.8	8
107	First-principles calculations-based model for the reactive ion etching of metal oxide surfaces. <i>Vacuum</i> , 2008 , 83, 599-601	3.7	8
106	Ab initio study of magnetic and electronic properties of Fe-filled single-walled carbon nanotubes. <i>Journal of Physics Condensed Matter</i> , 2004 , 16, S5755-S5758	1.8	8
105	PPS-metal adhesion: a density functional theory-based study. <i>Solid State Communications</i> , 2004 , 132, 405-408	1.6	8
104	Dynamics of STM-Induced Acetylene Dissociation on Cu(100). <i>Journal of the Physical Society of Japan</i> , 2002 , 71, 569-573	1.5	8
103	Hydrogenase-based nanomaterials as anode electrode catalyst in polymer electrolyte fuel cells. <i>Solid State Communications</i> , 2005 , 133, 589-591	1.6	8
102	Dynamics of STM-induced acetylene rotation on Cu(100). Surface Science, 2000, 454-456, 1052-1057	1.8	8
101	Dynamics of STM-Induced CO Desorption from Cu(111). <i>Journal of the Physical Society of Japan</i> , 1998 , 67, 4018-4021	1.5	8
100	Spin-Dependent O Binding to Hemoglobin. ACS Omega, 2018, 3, 9241-9245	3.9	8
99	Adsorption of Fe and Co Nanowires to (3,3) Single-Walled Carbon Nanotubes. <i>Japanese Journal of Applied Physics</i> , 2007 , 46, 1788-1791	1.4	7
98	Orthopara H2 conversion on multiple-decked sandwich clusters of M(C6H6)2 (M=Mn, Fe, Co) induced by an inhomogeneity of spin density distribution. <i>Thin Solid Films</i> , 2006 , 509, 223-226	2.2	7
97	First Principles Interpretation of Cyclohexane Dehydrogenation Process Using Pt. <i>Japanese Journal of Applied Physics</i> , 2005 , 44, 402-405	1.4	7
96	Quantum Dynamics of Abstraction in $H(g)+H(a)/Cu(111)$: Direct (Eley R ideal) and Indirect (Hot-Atom) Processes. <i>Journal of the Physical Society of Japan</i> , 2002 , 71, 222-227	1.5	7
95	Rotational state modification and fast ortho-para conversion of H2 trapped within the highly anisotropic potential of Pd(210). <i>Physical Review B</i> , 2018 , 97,	3.3	6
94	Analysis of the Changes in Electronic Structures and Work Function Variation in Alkali Metal [^] ndash;Metal Surface Systems. <i>Journal of the Vacuum Society of Japan</i> , 2014 , 57, 27-31		6

93	Tunneling Effect of O2 in Dissociative Adsorption on Pt(111) Surface. <i>E-Journal of Surface Science and Nanotechnology</i> , 2013 , 11, 72-75	0.7	6
92	First principles investigation on Fe-filled single-walled carbon nanotubes on Ni (111) and Cu (111). Journal of Magnetism and Magnetic Materials, 2007 , 310, e748-e750	2.8	6
91	Polybutylene terephthalate on metals: a density functional theory and cluster models investigation. <i>Journal of Physics Condensed Matter</i> , 2006 , 18, 1137-1142	1.8	6
90	Molecular orientation dependence of ortho-para conversion of a H2 interacting with a metal surface. <i>Journal of Applied Physics</i> , 2003 , 93, 644-648	2.5	6
89	Effects of H coverage on the dynamics of H abstraction from Cu(111). <i>Surface Science</i> , 2002 , 507-510, 838-844	1.8	6
88	Can we probe local surface reactivity with hydrogen molecules?. <i>Journal of Physics Condensed Matter</i> , 2002 , 14, 4379-4384	1.8	6
87	Examining Poly(Phenylene Sulfide) Adhesion using Cluster Models. <i>Shinku/Journal of the Vacuum Society of Japan</i> , 2005 , 48, 235-237		6
86	Effect of van der Waals Interaction on Ortho-Para Conversion of H2 on Ag(111) Surfaces. <i>Journal of the Vacuum Society of Japan</i> , 2012 , 55, 115-117		6
85	CH3Cl/Cu(410): Interaction and Adsorption Geometry. <i>Journal of Physical Chemistry C</i> , 2018 , 122, 1182	25-3.1883	31 5
84	Kondo Effect in the Systems of Magnetic Trimers on a Metal Surface. <i>Journal of the Physical Society of Japan</i> , 2010 , 79, 113706	1.5	5
83	First Principles Based Investigation of Materials for Resistive RAM. <i>Journal of Computational and Theoretical Nanoscience</i> , 2008 , 5, 1976-1979	0.3	5
82	Applying computational nanomaterials design to the reactive ion etching of NiO thin films-a preliminary investigation. <i>Journal of Physics Condensed Matter</i> , 2007 , 19, 365210	1.8	5
81	Temperature and magnetic field dependence of the Yosidakondo resonance for a single magnetic atom adsorbed on a surface. <i>Thin Solid Films</i> , 2006 , 509, 168-172	2.2	5
80	Suppression of carrier spin polarization in diluted ferromagnetic semiconductors. <i>Journal of Physics Condensed Matter</i> , 2004 , 16, S5787-S5790	1.8	5
79	Change of magnetic properties of benzenes in multiple-decked sandwich clusters: Mnn(C6H6)n+1 (n = 1,2). <i>Journal of Physics Condensed Matter</i> , 2004 , 16, S5749-S5753	1.8	5
78	Adsorbate dynamics induced by STM. <i>Applied Surface Science</i> , 2001 , 169-170, 25-29	6.7	5
77	Isotope effects on the rotationally inelastic diffraction dynamics of hydrogen scattered from Cu(001). <i>Surface Science</i> , 2001 , 493, 298-304	1.8	5
76	Magnetic-Adsorbate-Induced Spin Polarization in Carbon Materials II wo Fe Atoms on Planar C10and C10H8 [Journal of the Physical Society of Japan, 2003, 72, 2413-2416	1.5	5

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75	C2H4 adsorption on Cu(210), revisited: bonding nature and coverage effects. <i>Physical Chemistry Chemical Physics</i> , 2016 , 18, 23621-7	3.6	5
74	Dynamics of Mu, H, D, and T Absorption into Pd(111): Isotope Effects. <i>Journal of the Physical Society of Japan</i> , 2014 , 83, 013601	1.5	4
73	Rotational Effects on the Dissociative Adsorption and Abstraction Dynamics of O2/Al(111). <i>Journal of the Physical Society of Japan</i> , 2013 , 82, 113602	1.5	4
72	Effect of Lithium Absorption at Tetrahedral Site and Isomorphic Substitution on Montmorillonite Properties: A Density Functional Theory Study. <i>Japanese Journal of Applied Physics</i> , 2011 , 50, 055701	1.4	4
71	Transport properties of a single-quantum dot Aharonov-Bohm interferometer. <i>European Physical Journal B</i> , 2007 , 57, 27-35	1.2	4
70	Molecular orientation dependence of ortho-para H2 conversion on Fe(OH)3 cluster induced by hyperfine contact interaction. <i>European Physical Journal D</i> , 2006 , 38, 99-101	1.3	4
69	Catalytic Reactivity of a Transition Metal (Pt, Pd, Ni and Cu) Atom on Cyclohexane Dehydrogenation. <i>Shinku/Journal of the Vacuum Society of Japan</i> , 2004 , 47, 155-158		4
68	Spin-dependent transport through Fe nanowires on Cu surfaces. <i>Journal of Magnetism and Magnetic Materials</i> , 2004 , 272-276, 1650-1651	2.8	4
67	Dynamical Cluster Approximation in Disordered Systems with Magnetic Impurities. <i>Journal of the Physical Society of Japan</i> , 2004 , 73, 3448-3452	1.5	4
66	Orientational effects on the dynamics of H2 dissociation at the atop-Cu, atop-Pt, and Cu P t bridge sites of an ordered Cu3Pt(1 1 1). <i>Surface Science</i> , 2001 , 482-485, 318-323	1.8	4
65	Band Structure and Surface Localized States of Fe Thin Film on Cu Surface. <i>Shinku/Journal of the Vacuum Society of Japan</i> , 2004 , 47, 232-234		4
64	Coadsorption of hydrazine (N2H4) and OH on NiZn surface: A DFT-based study. <i>Surface Science</i> , 2020 , 691, 121505	1.8	4
63	Density functional study of methyl butanoate adsorption and its C-O bonds cleavage on MoS-based catalyst with various loads of Ni promoters. <i>Journal of Physics Condensed Matter</i> , 2019 , 31, 365001	1.8	3
62	Molecular Adsorption, Hindered Rotation, and Species Separation of H2/SrTiO3(001). <i>Journal of the Physical Society of Japan</i> , 2017 , 86, 073601	1.5	3
61	First principles study of orthopara H2 conversion on the O2(0.25[ML)/Ag(111) system. <i>Current Applied Physics</i> , 2012 , 12, S115-S118	2.6	3
60	First-Principles Calculations for Chemical Reaction between Sodium Diethyldithiocarbamate and Transition-Metal (Cr) atom to Produce Cr(DDC)3and Cr(DDC)2ODDC. <i>Japanese Journal of Applied Physics</i> , 2006 , 45, L1103-L1105	1.4	3
59	The effect of hydrogen adsorption on the magnetic properties of Fe adatoms on Si(001). <i>Journal of Physics Condensed Matter</i> , 2004 , 16, S5763-S5767	1.8	3
58	Steric effect on op conversion of a H2 interacting with a 3d impurity sitting on a metal oxide surface. <i>Surface Science</i> , 2002 , 514, 273-282	1.8	3

57	Magnetic properties in the high-temperature cuprate superconductors with nonmagnetic impurities. <i>Physical Review B</i> , 2003 , 68,	3.3	3
56	Isotope effects on direct and indirect processes of hydrogen abstraction from Cu(111). <i>Journal of Physics Condensed Matter</i> , 2002 , 14, 4345-4354	1.8	3
55	Rotational polarization and filtering of hydrogen molecules by metal surfaces lisotope effects. <i>Surface Science</i> , 1999 , 427-428, 358-363	1.8	3
54	Probing Local Surface Reactivity With Hydrogen Molecules-Realizing an Atom/Molecule Scanning Probe <i>Shinku/Journal of the Vacuum Society of Japan</i> , 2003 , 46, 391-396		3
53	Influence of reactive gas-phase species on the structure of an air/water interface. <i>Journal Physics D: Applied Physics</i> , 2017 , 50, 244002	3	2
52	Increasing the proton conductivity of sulfonated polyether ether ketone by incorporating graphene oxide: Morphology effect on proton dynamics. <i>Japanese Journal of Applied Physics</i> , 2018 , 57, 035201	1.4	2
51	First-principles study on oxygen ion conduction of La2GeO5based on the density functional theory. Journal of Physics: Conference Series, 2012 , 379, 012012	0.3	2
50	Influence of spin configuration on the transport properties of transition metal oxide nanostructures. <i>Solid State Communications</i> , 2007 , 142, 104-109	1.6	2
49	A nanoscale understanding of the adhesion of polybutylene terephthalate on aluminum. <i>Surface Science</i> , 2007 , 601, 5241-5245	1.8	2
48	Inducing spin polarization in carbon materials by introducing magnetic adsorbates E e2 on a planar C10. <i>Journal of Magnetism and Magnetic Materials</i> , 2004 , 272-276, 1602-1603	2.8	2
47	Transport properties via surface localized states of Ru, Rh and Pd thin films on Ag(111). <i>Solid State Communications</i> , 2005 , 135, 698-702	1.6	2
46	Temperature dependence of the spectral profile of the Yosidakondo resonance for a single magnetic atom adsorbed on a metal surface. <i>Surface Science</i> , 2005 , 593, 49-53	1.8	2
45	Dynamical phenomena including many body effects at metal surfaces. Surface Science, 2002, 500, 105-1	26 8	2
44	Some solutions for a system of interacting electrons confined in one dimension. <i>Physics Letters, Section A: General, Atomic and Solid State Physics</i> , 2000 , 277, 123-128	2.3	2
43	Orientational effects on the molecular diffraction dynamics of H2 scattered from Cu(0 0 1). <i>Surface Science</i> , 2001 , 482-485, 306-311	1.8	2
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