

Noriaki Takagi

List of Publications by Year in descending order

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116
papers

3,706
citations

172207

29
h-index

138251

58
g-index

116
all docs

116
docs citations

116
times ranked

3502
citing authors

#	ARTICLE	IF	CITATIONS
1	First-principles calculation of the graphene Dirac band on semi-infinite Ir(111). Physical Review B, 2020, 102, .	1.1	2
2	Scanning tunneling spectroscopy studies of topological materials. Journal of Physics Condensed Matter, 2020, 32, 243001.	0.7	7
3	Mechanically Tunable Spontaneous Vertical Charge Redistribution in Few-Layer WTe ₂ . Journal of Physical Chemistry C, 2020, 124, 2008-2012.	1.5	8
4	Unsubstituted and Fluorinated Copper Phthalocyanine Overlayers on Si(111)-(√7 × √3)-In Surface: Adsorption Geometry, Charge Polarization, and Effects on Superconductivity. Journal of Physical Chemistry C, 2019, 123, 8951-8958.	1.5	15
5	Quasiparticle scattering in type-II Weyl semimetal MoTe ₂ . Journal of Physics Condensed Matter, 2018, 30, 105703.	0.7	7
6	Silicene on Ag(111). , 2018, , 312-317.		0
7	Inelastic electron tunneling spectroscopy by STM of phonons at solid surfaces and interfaces. Progress in Surface Science, 2018, 93, 131-145.	3.8	8
8	Spin-orbit interaction in unoccupied surface states. Progress in Surface Science, 2018, 93, 177-188.	3.8	5
9	Surface structure of novel semimetal WTe ₂ . Applied Physics Express, 2017, 10, 045702.	1.1	9
10	Visualizing Type-II Weyl Points in Tungsten Ditelluride by Quasiparticle Interference. ACS Nano, 2017, 11, 11459-11465.	7.3	37
11	Structural evolution of Bi thin films on Au(111) revealed by scanning tunneling microscopy. Physical Review B, 2017, 96, .	1.1	20
12	Electron Transport Through a Single Molecule in Scanning Tunneling Microscopy Junction. Advances in Atom and Single Molecule Machines, 2017, , 355-379.	0.0	1
13	Single-molecule quantum dot as a Kondo simulator. Nature Communications, 2017, 8, 16012.	5.8	77
14	Transport characteristics of a silicene nanoribbon on Ag(110). Beilstein Journal of Nanotechnology, 2017, 8, 1699-1704.	1.5	10
15	Impact of reduced symmetry on magnetic anisotropy of a single iron phthalocyanine molecule on a Cu substrate. Journal of Chemical Physics, 2016, 144, 044701.	1.2	9
16	Comment on "Rashba Spin-Orbit Coupling in Image Potential States". Physical Review Letters, 2016, 117, 239701.	2.9	4
17	Atomic structure of "multilayer silicene" grown on Ag(111): Dynamical low energy electron diffraction analysis. Surface Science, 2016, 651, 70-75.	0.8	24
18	Linewidth analysis of image potential states on noble metal surfaces with high-energy resolved two-photon photoemission spectroscopy. Surface and Interface Analysis, 2016, 48, 1194-1198.	0.8	1

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19	Rashba splitting in an image potential state investigated by circular dichroism two-photon photoemission spectroscopy. <i>Physical Review B</i> , 2016, 94, .	1.1	12
20	Surface phonon excitation on clean metal surfaces in scanning tunneling microscopy. <i>Physical Review B</i> , 2016, 93, .	1.1	13
21	Model Hamiltonian approach to the magnetic anisotropy of iron phthalocyanine at solid surfaces. <i>Physical Review B</i> , 2016, 94, .	1.1	5
22	Spectroscopic Identification of Ag-Terminated α -Multilayer Silicene Grown on Ag(111). <i>Journal of Physical Chemistry C</i> , 2016, 120, 6689-6693.	1.5	17
23	Silicene on Ag(111): Structure Evolution and Electronic Structure. <i>Springer Series in Materials Science</i> , 2016, , 143-165.	0.4	0
24	One-dimensional edge state of Bi thin film grown on Si(111). <i>Applied Physics Letters</i> , 2015, 107, .	1.5	35
25	Pragmatic Application of Abstract Algebra to Two-Dimensional Lattice Matching. <i>E-Journal of Surface Science and Nanotechnology</i> , 2015, 13, 361-365.	0.1	5
26	Comparison of electronic structure between monolayer silicenes on Ag (111). <i>Chinese Physics B</i> , 2015, 24, 087307.	0.7	8
27	Electronic structure of the 4 \times 4 silicene monolayer on semi-infinite Ag(111). <i>New Journal of Physics</i> , 2015, 17, 015013.	1.2	21
28	Silicene on Ag(111): Geometric and electronic structures of a new honeycomb material of Si. <i>Progress in Surface Science</i> , 2015, 90, 1-20.	3.8	58
29	Electronic and Geometric Structure of Silicene on Ag. <i>Journal of the Vacuum Society of Japan</i> , 2014, 57, 428-433.	0.3	0
30	Transport characteristics of a single C_{60} -molecule junction revealed by multiple Andreev reflections. <i>Physical Review B</i> , 2014, 90, .	1.1	10
31	Controlling orbital-selective Kondo effects in a single molecule through coordination chemistry. <i>Journal of Chemical Physics</i> , 2014, 141, 054702.	1.2	27
32	Determination of atomic positions in silicene on Ag(111) by low-energy electron diffraction. <i>Surface Science</i> , 2014, 623, 25-28.	0.8	97
33	Electronic decoupling by h-BN layer between silicene and Cu(111): A DFT-based analysis. <i>New Journal of Physics</i> , 2014, 16, 105019.	1.2	20
34	Adsorbed states of iron(II) phthalocyanine on Ag(111) studied by high-resolution electron energy loss spectroscopy. <i>Surface and Interface Analysis</i> , 2014, 46, 1253-1256.	0.8	5
35	Silicene grown on silver surface. <i>Journal of Surface Analysis (Online)</i> , 2014, 21, 63-70.	0.1	0
36	Comment on "Evidence for Dirac Fermions in a Honeycomb Lattice Based on Silicon". <i>Physical Review Letters</i> , 2013, 110, 229701.	2.9	56

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37	Substrate-Induced Symmetry Breaking in Silicene. <i>Physical Review Letters</i> , 2013, 110, 076801.	2.9	358
38	Structural transition of silicene on Ag(111). <i>Surface Science</i> , 2013, 608, 297-300.	0.8	169
39	Enhancement of Inelastic Electron Tunneling Conductance Caused by Electronic Decoupling in Iron Phthalocyanine Bilayer on Ag(111). <i>Journal of Physical Chemistry C</i> , 2013, 117, 21832-21837.	1.5	14
40	Mode-selective electron-phonon coupling in laser photoemission on Cu(110). <i>Physical Review B</i> , 2013, 88, .	1.1	6
41	High Energy- and Momentum-Resolved Two-Photon Photoemission Spectroscopy: Pure Dephasing Rate Measurement on Image States. <i>Hyomen Kagaku</i> , 2013, 34, 421-425.	0.0	1
42	Structure of Silicene Grown on Ag(111). <i>Applied Physics Express</i> , 2012, 5, 045802.	1.1	518
43	Symmetry-Driven Novel Kondo Effect in a Molecule. <i>Physical Review Letters</i> , 2012, 109, 086602.	2.9	138
44	Combined Scanning Tunneling Microscopy and High-Resolution Electron Energy Loss Spectroscopy Study on the Adsorption State of CO on Ag(001). <i>Langmuir</i> , 2012, 28, 13249-13252.	1.6	7
45	Density Functional Theory Calculation for Magnetism of Fe-Phthalocyanine Molecules on Au(111). <i>E-Journal of Surface Science and Nanotechnology</i> , 2012, 10, 38-44.	0.1	8
46	Nature of Electron Transport by Pyridine-Based Tripodal Anchors: Potential for Robust and Conductive Single-Molecule Junctions with Gold Electrodes. <i>Journal of the American Chemical Society</i> , 2011, 133, 3014-3022.	6.6	94
47	Evolution of Kondo Resonance from a Single Impurity Molecule to the Two-Dimensional Lattice. <i>Physical Review Letters</i> , 2011, 106, 187201.	2.9	138
48	Magnetism of Iron (II) Phthalocyanine at Surfaces -Spin, Magnetic Anisotropy and Kondo Effect-. <i>Hyomen Kagaku</i> , 2011, 32, 629-634.	0.0	0
49	Spin Excitation of a Single Iron (II) Phthalocyanine Molecule on the Cu(110) Surface. <i>Hyomen Kagaku</i> , 2009, 30, 433-438.	0.0	1
50	Mechanism of vibrational excitation in inelastic photoemission from solid surfaces. <i>Physical Review B</i> , 2009, 80, .	1.1	7
51	Adsorption-Induced Switching of Magnetic Anisotropy in a Single Iron(II) Phthalocyanine Molecule on an Oxidized Cu(110) Surface. <i>Physical Review Letters</i> , 2009, 102, 167203.	2.9	268
52	Synthesis of tripodal anchor units bearing selenium functional groups and their adsorption behaviour on gold. <i>Physical Chemistry Chemical Physics</i> , 2009, 11, 4949.	1.3	26
53	Structure and thermal fluctuation of one-dimensional AgO chains on Ag(110) surfaces studied with density functional theory and Monte Carlo simulations. <i>Journal of Chemical Physics</i> , 2008, 129, 154709.	1.2	4
54	Explosive evolution of hydrogen abstraction of water on oxidized Ag(110) surfaces studied by scanning tunnelling microscopy. <i>Physical Chemistry Chemical Physics</i> , 2007, 9, 5274.	1.3	3

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55	Vibrationally Induced Inelastic Structures in Laser Photoemission Spectra. Hyomen Kagaku, 2007, 28, 378-384.	0.0	0
56	Photochemistry of cyclohexane on Cu(111). Physical Chemistry Chemical Physics, 2006, 8, 179-185.	1.3	5
57	Quantum delocalization of hydrogen on metal surfaces. Surface Science Reports, 2005, 57, 113-156.	3.8	40
58	Excitation mechanism and ultrafast vibrational wavepacket dynamics of alkali-metal atoms on Pt(111). Surface Science, 2005, 593, 110-115.	0.8	7
59	Femtosecond wavepacket dynamics of potassium adsorbate on Pt(111). Springer Series in Chemical Physics, 2005, , 307-309.	0.2	1
60	Electronic structure and femtosecond electron transfer dynamics at noble metal/tris-(8-hydroxyquinoline) aluminum interfaces. Physical Review B, 2005, 71, .	1.1	24
61	Femtosecond wavepacket dynamics of Cs adsorbates on Pt(111): Coverage and temperature dependences. Physical Review B, 2005, 71, .	1.1	30
62	Mode-selective excitation of coherent surface phonons on alkali-covered metal surfaces. Physical Chemistry Chemical Physics, 2005, 7, 2697.	1.3	27
63	Electron Transfer Dynamics from Organic Adsorbate to a Semiconductor Surface: Zinc Phthalocyanine on TiO ₂ (110). Journal of Physical Chemistry B, 2005, 109, 18018-18024.	1.2	33
64	In Situ Observation of CO Oxidation on Ag(110)(2 \times 1)-O by Scanning Tunneling Microscopy: Structural Fluctuation and Catalytic Activity. Journal of Physical Chemistry B, 2005, 109, 14536-14543.	1.2	15
65	Direct Time-Domain Observation of Ultrafast Dephasing in Adsorbate-Substrate Vibration under the Influence of a Hot Electron Bath: Cs Adatoms on Pt(111). Physical Review Letters, 2004, 92, 057401.	2.9	48
66	Ultrafast excited state dynamics in 3,4,9,10-perylene tetracarboxylic dianhydride (PTCDA) thin films. Chemical Physics Letters, 2004, 383, 261-265.	1.2	10
67	Reactivity of molecular oxygen: conversion of methanol to formate at low temperatures on Pt(111). Chemical Physics Letters, 2004, 392, 334-339.	1.2	15
68	Title is missing!. Shinku/Journal of the Vacuum Society of Japan, 2004, 47, 412-417.	0.2	0
69	Structural changes of AgO chains on Ag(110) by photo- and CO-induced oxygen elimination. Surface Science, 2003, 528, 144-150.	0.8	5
70	Role of Structural Fluctuation in a Surface Reaction Studied by Scanning Tunneling Microscopy: The CO + O \rightarrow CO ₂ Clean-Off Reaction on Ag(110)-(2 \times 1)-O. Physical Review Letters, 2003, 90, 226105.	2.9	15
71	Reaction intermediates in the oxidation of methanol on a Pt(111)-(2 \times 2)O surface. Journal of Chemical Physics, 2003, 119, 4879-4886.	1.2	35
72	<title>Dynamic formation of reaction sites at nanostructured one-dimensional surface compounds</title>. , 2003, , .		0

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73	Anomalous quenching of electronic states of nanographene on Pt(111) by deuterium edge termination. Physical Review B, 2002, 65, .	1.1	14
74	Coherent surface phonon at aGaAs(100) $\sqrt{2}\times\sqrt{2}$ surface. Physical Review B, 2002, 65, .	1.1	16
75	Thermal decomposition of acetylene on Pt(111) studied by scanning tunneling microscopy. Surface Science, 2002, 514, 414-419.	0.8	10
76	Impulsive excitation of a vibrational mode of Cs on Pt(111). Chemical Physics Letters, 2002, 366, 606-610.	1.2	58
77	Fermi Surface Nesting and Structural Transition on a Metal Surface: In/Cu(001). Physical Review Letters, 2001, 86, 854-857.	2.9	43
78	Growth mechanism of the Pd(100)-p(2 $\sqrt{2}\times\sqrt{2}$)-p4g-Al surface alloy. Surface Science, 2000, 460, 264-276.	0.8	10
79	Adsorbed states of K on the diamond (100)(2 $\sqrt{2}\times\sqrt{2}$) surface. Diamond and Related Materials, 2000, 9, 162-169.	1.8	11
80	Diels-Alder Reaction on the Clean Diamond (100) 2 $\sqrt{2}\times\sqrt{2}$ Surface. Japanese Journal of Applied Physics, 1999, 38, L1496-L1498.	0.8	29
81	Surface Phonons, Electronic Structure and Chemical Reactivity of Diamond (100)(2 $\sqrt{2}\times\sqrt{2}$) Surface. Japanese Journal of Applied Physics, 1999, 38, 6659-6666.	0.8	25
82	Surface phonons of theSi(001)(2 $\sqrt{2}\times\sqrt{2}$)surface. Physical Review B, 1999, 60, 10919-10925.	1.1	44
83	Structure and chemistry of Pd(100)-p(2 $\sqrt{2}\times\sqrt{2}$)-p4g-Al surface alloy. Surface Science, 1999, 427-428, 74-78.	0.8	7
84	A new reaction channel in H(g)+D(a)/Pd(100): absorption versus abstraction. Surface Science, 1999, 427-428, 277-281.	0.8	9
85	Chemisorbed states of atomic oxygen and its replacement by atomic hydrogen on the diamond (100)-(2 $\sqrt{2}\times\sqrt{2}$) surface. Surface Science, 1999, 436, 63-71.	0.8	50
86	Vibrational spectra of hydrogen on the Rh(111) surface. Surface Science, 1999, 441, 507-514.	0.8	33
87	Subsurface Hydrogen at Pd(100) Induced by Gas-Phase Atomic Hydrogen. Journal of Physical Chemistry B, 1999, 103, 7876-7881.	1.2	20
88	Mechanisms of the CO oxidation on the Pd(110)c(2 $\sqrt{2}\times\sqrt{2}$)-O surface. Surface Science, 1998, 397, 295-305.	0.8	5
89	Adsorbed states of CO on the Si(100)-K surface: electron energy-loss spectroscopy and thermal desorption studies. Surface Science, 1998, 395, L246-L251.	0.8	7
90	Path and mechanism of hydrogen absorption at Pd(100). Surface Science, 1998, 401, 344-354.	0.8	153

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91	Absorption of D in the H+D/Pd(100) reaction. Surface Science, 1998, 411, L849-L854.	0.8	10
92	Investigation on the Surface Electronic States of the Si(001) c(4 $\sqrt{2}$ -2) and c(8 $\sqrt{2}$ -8) Surfaces: An Electron Energy Loss Spectroscopy Study. Japanese Journal of Applied Physics, 1997, 36, L975-L978.	0.8	13
93	Adsorbed states of H on Ni(111) at 100 K: A vibrational study. Physical Review B, 1997, 56, 14952-14955.	1.1	17
94	Adsorption and Thermal Decomposition of Formic Acid on the Si(100)(2 $\sqrt{2}$ -1) $\sqrt{2}$ K Surface. Journal of Physical Chemistry B, 1997, 101, 7007-7011.	1.2	14
95	Adsorption and thermal decomposition of N ₂ O on Si(100): electron energy loss spectroscopy and thermal desorption studies. Surface Science, 1997, 382, 214-220.	0.8	13
96	Adsorbate-adsorbate interaction among NO and CO coadsorbed on Pd(100). Applied Surface Science, 1997, 121-122, 571-574.	3.1	4
97	Interaction of NO with CO on Pd(100): ordered coadsorption structures and explosive reaction. Surface Science, 1996, 350, 79-90.	0.8	47
98	Temporal and local reduction of adsorption potential energy under gas phase: CO on Ni(100) and Pt(111). Surface Science, 1996, 363, 85-90.	0.8	10
99	CO adsorption on the Pd(110)c(2 $\sqrt{2}$ -4)-O surface: formation of a p(2 $\sqrt{2}$ -4) structure. Surface Science, 1996, 365, 422-428.	0.8	2
100	Quantum delocalization of H on Pd(110): A vibrational study. Physical Review B, 1996, 53, 13767-13771.	1.1	25
101	REACTIONS OF GAS MOLECULES ON SILICON SURFACES STUDIED BY HIGH RESOLUTION ELECTRON ENERGY LOSS SPECTROSCOPY. , 1996, , 285-301.		0
102	Location of an O atom in the Pd(110)c(2 $\sqrt{2}$ -4)-O structure. An EELS study. Chemical Physics Letters, 1995, 232, 531-536.	1.2	32
103	Hydrogen desorption from Si(100)(2 $\sqrt{2}$ -1)-H induced by potassium adsorption. Surface Science, 1995, 325, 11-20.	0.8	11
104	Atomic-hydrogen-induced restructuring of the Si(100)(2 $\sqrt{2}$ -1)-K surface. Surface Science, 1995, 337, L783-L788.	0.8	3
105	Explosive production of CO ₂ from (NO + CO)/Pd(100). Surface Science, 1995, 341, L1096-L1100.	0.8	16
106	Adsorbed states of NH ₃ and C ₆ H ₆ on the Si(111)($\sqrt{3}\sqrt{3}$ -R30 $^\circ$ -B) surface: Thermal-desorption and electron-energy-loss-spectroscopy studies. Physical Review B, 1994, 50, 17440-17449.	1.1	8
107	Temporal and Local Reduction of a Potential Energy Well under Dynamic Equilibrium: CO on Ni(100). Physical Review Letters, 1994, 73, 292-295.	2.9	34
108	Stability of adsorbed states and site-conversion kinetics: CO on Ni(100). Physical Review B, 1994, 49, 16670-16677.	1.1	35

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109	Chemical reactivity of the Si(111) (1 × 1) R30°-B surface: An electron-energy-loss spectroscopy study. Applied Surface Science, 1994, 82-83, 434-436.	3.1	5
110	Site conversion of CO on Ni(100): binding-energy difference and role of low-energy hindered vibrations. Chemical Physics Letters, 1993, 211, 48-52.	1.2	28
111	Direct observation of isothermal adsorption and desorption processes of CO on the Ni(100) surface. Chemical Physics Letters, 1993, 215, 120-124.	1.2	14
112	The growth of ice clusters on the Si(100)(2 × 1)-H(D) surface: Electron energy loss spectroscopy and thermal desorption studies. Surface Science, 1993, 297, L43-L47.	0.8	10
113	Electron scattering from the K-exposed Si(100)(2 × 1)-H surface. Physical Review B, 1992, 45, 13524-13530.	1.1	6
114	Chemical reactivity of the Si(100)(2 × 1)-K surface: electron energy loss spectroscopy and thermal desorption studies. Surface Science, 1991, 242, 498-502.	0.8	17
115	Promoted oxidation of the K-modified Si(100) (2 × 1) surface: Electron-energy-loss-spectroscopy and thermal-desorption studies. Physical Review B, 1991, 44, 12945-12951.	1.1	13
116	Existence of two adsorbed states for K on the Si(100)(2 × 1) surface: A thermal desorption study. Physical Review B, 1990, 42, 1868-1871.	1.1	72