Guy TrÃ©glia

List of Publications by Year
in descending order

162 papers	5,038 citations	87723 38 h-index	110170 64 g-index
162 all docs	162 docs citations	162 times ranked	2087 citing authors

#	Article	IF	Citations
1	Tight-binding modeling of interstitial ordering processes in metals: Application to zirconium hydrides. Physical Review B, 2020, 101, .	1.1	2
2	Bidimensional phases in Coâ€“Pt surface alloys: A theoretical study of ordering and surface segregation. Surface Science, 2019, 679, 128-138.	0.8	15
3	Theoretical study of xenon adsorption on UO₂surfaces. Journal of Physics Communications, 2018, 2, 035041.	0.5	1
4	Tight-binding modelling of ferromagnetic metals and alloys. Modelling and Simulation in Materials Science and Engineering, 2017, 25, 084004.	0.8	2
5	Effect of magnetism on surface segregation in FeNi alloys. Journal of Physics Condensed Matter, 2016, 28, 064003.	0.7	2
6	Stress influence on substitutional impurity segregation on dislocation loops in IVâ€“IV semiconductors. Computational Materials Science, 2016, 114, 23-32.	1.4	3
7	Atomistic modelling of residual stress at UO2surfaces. Journal of Physics Condensed Matter, 2016, 28, 015006.	0.7	2
8	Environment dependence of magnetic moment and atomic level shifts within tight-binding approximation: An illustration in the case of cobalt. Surface Science, 2016, 646, 261-268.	0.8	7
9	Tight-Bindingn-momentpotential for zirconium hydride atomistic modeling. Metallurgical Research and Technology, 2015, 112, 102.	0.4	3
10	Electronic structure of CoPt based systems: from bulk to nanoalloys. Journal of Physics Condensed Matter, 2015, 27, 455503.	0.7	5
11	Ordering and surface segregation in<mml:math xmlns:mml="http://www.w3.org/1998/Math/MathML"><mml:mrow><mml:msub><mml:mi>Co</mml:mi><mml:mrow><mml:mn>1</mml:mn> A theoretical study from surface alloys to nanoalloys. Physical Review B, 2015, 91, .	1.1	31
12	How to derive tight-bindingspdpotentials? Application to zirconium. Journal of Physics Condensed Matter, 2015, 27, 336301.	0.7	5
13	Theoretical study of xenon adsorption in UO₂nanoporous matrices. Journal of Physics Condensed Matter, 2014, 26, 485015.	0.7	7
14	Atomistic study of porosity impact on phonon driven thermal conductivity: Application to uranium dioxide. Journal of Applied Physics, 2014, 115, .	1.1	8
15	Theoretical investigation of Cottrell atmosphere in silicon. Acta Materialia, 2014, 65, 1-9.	3.8	11
16	Surface segregation trends in transition metal alloys. Physical Review B, 2013, 88, .	1.1	7
17	Disentangling coordination and alloy effects in transition-metal nanoalloys from their electronic structure. Physical Review B, 2013, 88, .	1.1	5
18	Surface Segregation in Transition Metal Alloys: From Electronic Structure to Phase Portraits. Progress of Theoretical Physics Supplement, 2013, 101, 159-180.	0.2	2

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19	Theoretical investigation of the influence of reaction and diffusion kinetics upon thin-film reactive diffusion. Physical Review B, 2012, 85, .	1.1	11
20	Electronic Structure of Nanoalloys: A Guide of Useful Concepts and Tools. Engineering Materials, 2012, , 159-195.	0.3	0
21	Structure and properties of nanoscale materials: theory and atomistic computer simulation. International Journal of Nanotechnology, 2012, 9, 576.	0.1	1
22	An atomistic modelling of the porosity impact on UO2 matrix macroscopic properties. Journal of Nuclear Materials, 2011, 415, 210-216.	1.3	23
23	Ordering trends in transition metal alloys from tight-binding electronic structure calculations. Physical Review B, 2011, 84, .	1.1	7
24	Unified picture of<mml:math xmlns:mml="http://www.w3.org/1998/Math/MathML" display="inline"><mml:mrow><mml:mi>d</mml:mi></mml:mrow></mml:math>-band and core-level shifts in transition metal alloys. Physical Review B, 2011, 83, .	1.1	15
25	Molecular dynamics simulation of silicon oxidization. Thin Solid Films, 2010, 518, 2422-2426.	0.8	5
26	Atomistic modeling of strain and diffusion at the<mml:math xmlns:mml="http://www.w3.org/1998/Math/MathML" display="inline"><mml:mrow><mml:mtext>Si</mml:mtext><mml:mo>/</mml:mo><mml:msub><mml:mrow><mml:mtext>SiO</mml:mtext></mml:mrow></mml:msub></mml:mrow></mml:math> Physical Review B, 2010, 81, .	1.1	20
27	Rules for tight-binding calculations in bi-metallic compounds based on density functional theory: the case of Coâ€“Au. Journal of Physics Condensed Matter, 2010, 22, 505503.	0.7	9
28	Physical origin of thickness-controlled sequential phase formation during reactive diffusion: Atomistic modeling. Physical Review B, 2010, 82, .	1.1	21
29	Subnanometric Si film reactive diffusion on Ni. Applied Physics Letters, 2009, 95, 023111.	1.5	16
30	Strain effect on self-diffusion in silicon: Numerical study. Physical Review B, 2009, 79, .	1.1	11
31	Self-organization of Ge tetramers on Ag(001) surface: A 2D realization of unusual substrate mediated interactions. Surface Science, 2008, 602, 506-510.	0.8	11
32	Role of spâ€“d hybridization in the formation of stacking defects at metal surfaces. Surface Science, 2008, 602, 2681-2688.	0.8	4
33	Model of surface segregation driving forces and their coupling. Physical Review B, 2008, 78, .	1.1	29
34	Coverage dependence of<mml:math xmlns:mml="http://www.w3.org/1998/Math/MathML" display="inline"><mml:mrow><mml:mi mathvariant="normal">Sb</mml:mi><mml:mo>â^•</mml:mo><mml:mi mathvariant="normal">Si</mml:mi><mml:mrow><mml:mo>(</mml:mo><mml:mn>111</mml:mn><mml:mo>)</mml:mo></mml:mrow></mml:mrow></mml:math> and desorption modes: Interplay between chemical interactions and site transitions. Physical Review B, 2008, 77, .	1.1	6
35	Formation of an unexpected ordered two-dimensional Ag2Pb surfacealloy on Ag(111): A SXRD and STM study. Journal of Physics and Chemistry of Solids, 2006, 67, 601-604.	1.9	5
36	Photoelectron spectroscopy study of Pb/Ag(111) in the submonolayer range. Surface Science, 2006, 600, 1227-1230.	0.8	20

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37	Site segregation in size-mismatched nanoalloys: Application to Cuâ€“Ag. Surface Science, 2006, 600, 5011-5020.	0.8	38
38	Influence of surface stress in the missing row reconstruction of fcc transition metals. Surface Science, 2006, 600, 5131-5135.	0.8	34
39	Combining Solid State Physics Concepts and X-Ray Absorption Spectroscopy to Understand DeNOx Catalysis. Oil and Gas Science and Technology, 2006, 61, 677-689.	1.4	3
40	Sb/Si(111)Adsorption: Hidden Phase Transitions Behind Langmuir-Like Isotherms. Physical Review Letters, 2005, 94, 076101.	2.9	6
41	Ordered surface alloy formation of immiscible metals: The case of Pb deposited on Ag(111). Physical Review B, 2005, 72, .	1.1	32
42	Linear time dependence of the surfactant effect: A local equilibrium under flux. Physical Review B, 2004, 69, .	1.1	11
43	Modeling free and supported metallic nanoclusters: structure and dynamics. Phase Transitions, 2004, 77, 101-113.	0.6	83
44	An unusual composition profile: a LEEDâ€“TBIM study of Pt25Cu75(). Surface Science, 2003, 527, 71-79.	0.8	15
45	Relation between surface stress and (1Ã—2) reconstruction for (1 1 0) fcc transition metal surfaces. Applied Surface Science, 2003, 212-213, 866-871.	3.1	26
46	Kinetics study of antimony adsorption on Si(1 1 1). Applied Surface Science, 2003, 212-213, 715-723.	3.1	7
47	Cu-Ag (111) Polymorphism Induced by Segregation and Advacancies. Physical Review Letters, 2003, 91, 176103.	2.9	22
48	Misfit dislocation loops or incommensurate structure at an interface: Vibrational and anharmonic effects. Physical Review B, 2002, 66, .	1.1	12
49	Theoretical investigation of chemical and morphological ordering inPdcCu1â^'cclusters. Physical Review B, 2002, 66, .	1.1	37
50	Atomistic simulations of relaxation and reconstruction phenomena in heteroepitaxy: Co/Au(1 1 1). Applied Surface Science, 2002, 188, 134-139.	3.1	7
51	Compressive strain versus tensile strain. Applied Surface Science, 2001, 177, 238-242.	3.1	11
52	Thermal dependence of surface polymorphism: the Ag/Cu (111) case. Applied Surface Science, 2001, 177, 252-257.	3.1	4
53	An atomistic approach for stress relaxation in materials. , 2001, , 119-150.		2
54	New opportunities to understand heterogeneous catalysis processes on nanoscale bimetallic particles through synchrotron radiation and theoretical studies. Applied Catalysis A: General, 2000, 200, 47-54.	2.2	52

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55	New trends in heterogeneous catalysis processes on metallic clusters from synchrotron radiation and theoretical studies. Applied Surface Science, 2000, 164, 140-146.	3.1	18
56	Molecular dynamics simulations for the Ag/Cu (111) system: from segregated to constitutive interfacial vacancies. Applied Surface Science, 2000, 162-163, 219-226.	3.1	26
57	New Structures and Atomistic Analysis of the Polymorphism for the â˜ = 5 (210) [001] Tilt Boundary. Journal of Materials Science, 2000, 8, 55-69.	1.2	17
58	Ge/Ag(111) semiconductor-on-metal growth: Formation of anAg2Gesurface alloy. Physical Review B, 2000, 62, 16653-16656.	1.1	86
59	Intergranular segregation and vibrational effects: A local analysis. Physical Review B, 2000, 61, 14470-14480.	1.1	22
60	Ge deposition on Ag surfaces: Dependence of the adsorption characteristics on the surface orientation. Physical Review B, 2000, 61, 8469-8474.	1.1	10
61	Theoretical determination of two critical sizes for strain relaxation during Co/Pt(111) heteroepitaxy. Surface Science, 2000, 446, 272-282.	0.8	22
62	Ag/Cu(111) structure revisited through an extended mechanism for stress relaxation. Physical Review B, 1999, 59, 10910-10917.	1.1	99
63	Pt/Co(0001) superstructures in the submonolayer range: A tight-binding quenched-molecular-dynamics study. Physical Review B, 1999, 60, 2781-2788.	1.1	58
64	Layer-by-layer versus surfactant dissolution modes in heteroepitaxy. Physical Review B, 1999, 60, 13890-13901.	1.1	22
65	Local-density approximation study of semiconductor/metal adsorption characteristics: Ge/Ag(100). Physical Review B, 1999, 59, 15337-15345.	1.1	11
66	How to compare superficial and intergranular segregation? A new analysis within the mixed SMAâ€“TBIM approach. Acta Materialia, 1999, 47, 2705-2715.	3.8	63
67	Theoretical prediction of new dissolution modes during metal heteroepitaxy. Journal of Crystal Growth, 1999, 198-199, 83-88.	0.7	2
68	Alloy surfaces: segregation, reconstruction and phase transitions. Computational Materials Science, 1999, 15, 196-235.	1.4	118
69	Surface-induced ordering in phase separation systems: influence of concentration and orientation. Surface Science, 1999, 441, 225-239.	0.8	17
70	Step-driven molecular adsorption of Sb on Si(111). Surface Science, 1998, 395, 317-325.	0.8	18
71	Thermodesorption mass spectrometry study of the adsorption of Sb on misoriented Si(111). Surface Science, 1998, 417, 107-120.	0.8	6
72	Sb doping of Si molecular-beam epitaxial layers:â€ƒInfluence of the substrate misorientation. Physical Review B, 1997, 56, 7615-7622.	1.1	5

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73	Charge redistribution at Pd surfaces:Ab initiogrounds for tight-binding interatomic potentials. Physical Review B, 1997, 56, 12161-12166.	1.1	32
74	SURFACE ALLOY FORMATION IN THE Cuâ€“Pd(111) SYSTEM: A KTBIM APPROACH. Surface Review and Letters, 1997, 04, 1119-1122.	0.5	3
75	Microstructure of the surfactantlike effect in Ni/Ag(100) and (111). Physical Review B, 1997, 55, 10931-10937.	1.1	38
76	Step-descent mechanisms on Ag and Au(111). Surface Science, 1997, 377-379, 843-846.	0.8	9
77	On the â€œexoticâ€•behaviour of the Ptî—¸Sn system. Surface Science, 1997, 377-379, 1033-1037.	0.8	12
78	New magic numbers in metallic clusters: an unexpected metal dependence. Surface Science, 1997, 383, L719-L727.	0.8	102
79	Theoretical study of surface alloy formation through generation and annihilation of vacancies. Surface Science, 1996, 352-354, 552-556.	0.8	7
80	Flux dependence of the surfactant effect in : a theoretical study. Surface Science, 1996, 352-354, 562-566.	0.8	9
81	Competition or synergy between surface segregation and bulk ordering: the Cuî—¸Pd system. Surface Science, 1996, 352-354, 588-591.	0.8	12
82	Electronic structure of Pd clusters in the tight-binding approximation: influence of spd-hybridization. Surface Science, 1996, 352-354, 675-679.	0.8	32
83	Vacancy generation at steps and the kinetics of surface alloy formation. Surface Science, 1996, 364, 453-466.	0.8	25
84	High-Temperature Study of the Schwoebel Effect in Au(111). Physical Review Letters, 1996, 76, 2109-2112.	2.9	36
85	Intergranular Segregation in Cu(Ag) and Ag(Cu) Systems: Analysis of the Driving Force Using a Tight-Binding Scheme. Materials Science Forum, 1996, 207-209, 701-704.	0.3	4
86	Alliage de surface et Ã©quilibre local dans le systÃ¨me Pd/Cu (111). European Physical Journal Special Topics, 1996, 06, C7-155-C7-158.	0.2	3
87	CinÃ©tique de dissolution d'un dÃ©pÃ´t Fe/Cu. European Physical Journal Special Topics, 1996, 06, C7-151-C7-154.	0.2	0
88	Tight-binding molecular dynamics study of diffusion on Au and Ag(111). Surface Science, 1995, 331-333, 920-924.	0.8	42
89	Link between the surface wetting in Cu(Ag) and the layer-by-layer dissolution mode of a thick Ag deposit on a Cu substrate. Surface Science, 1995, 331-333, 805-810.	0.8	25
90	Anisotropy of diffusion along steps on the (111) faces of gold and silver. Physical Review B, 1994, 50, 12104-12117.	1.1	68

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91	Comparison of an experimental and a theoretical study of the surface alloy Pt/Cu(111). Journal of Electron Spectroscopy and Related Phenomena, 1994, 68, 357-362.	0.8	5
92	On the various terminations occurring in CuPt-ordered alloys: the TBIM approach. Surface Science, 1994, 307-309, 440-444.	0.8	16
93	On a â€œsurfactant-likeâ€•behaviour of deposit. Surface Science, 1994, 307-309, 531-537.	0.8	28
94	On the influence of topology on the energy profile in metallic Pd clusters. Surface Science, 1994, 307-309, 735-740.	0.8	17
95	Kinetics of segregation and dissolution in Cu1â^'cAgc and surface phase transition: comparison between mean field and Monte Carlo calculations. Surface Science, 1994, 307-309, 804-809.	0.8	17
96	Equilibrium and kinetics in the (111) surface of Cu-Ag alloys: Comparison between mean-field and Monte Carlo calculations. Physical Review B, 1994, 50, 1912-1921.	1.1	48
97	Experimental study of the growth and dissolution of Re/W(110) by core level photoemission spectroscopy: formation of a surface alloy. Surface Science, 1993, 286, 150-167.	0.8	15
98	Alloy surfaces and surface alloys: from equilibrium to kinetics. Surface Science, 1993, 287-288, 371-376.	0.8	30
99	Topology of a Ag monolayer on a Cu(111) substrate: a tight-binding quenched molecular dynamics study. Surface Science, 1993, 287-288, 476-481.	0.8	16
100	On a surprising anisotropy of surface segregation in CuPt alloys. Surface Science, 1993, 287-288, 851-856.	0.8	15
101	Structures of a Ag monolayer deposited on Cu(111), Cu(100), and Cu(110) substrates: An extended tight-binding quenched-molecular-dynamics study. Physical Review B, 1992, 46, 16018-16030.	1.1	129
102	Is the segregation-dissolution kinetics driven by a surface local equilibrium? An answer via the kinetic tight-binding Ising model. Surface Science, 1992, 274, 297-305.	0.8	69
103	Incomplete wetting of very dilute Cu(Ag) alloys by surface segregation. Surface Science, 1991, 251-252, 664-669.	0.8	16
104	Inversion of the core level shift between surface and subsurface atoms of the iridium (100)(1 Ã— 1) and (100)(5 Ã— 1) surfaces. Surface Science, 1991, 251-252, 717-721.	0.8	16
105	Surface segregation near the temperature of bulk phase separation: Incomplete wetting in Cu(Ag) alloys. Physical Review B, 1991, 44, 5842-5854.	1.1	77
106	Multilayer relaxation and reconstruction in bcc and fcc transition and noble metals. Vacuum, 1990, 41, 311-314.	1.6	29
107	On the temperature dependence of the surface sandwiches observed in PtNi and AgNi alloys. Vacuum, 1990, 41, 441-445.	1.6	2
108	Phase transitions in surface segregation ofPtcNi1â^'calloys from tight-binding Ising-model calculations. Physical Review B, 1990, 41, 4422-4434.	1.1	104

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109	Surface segregation in CuNi and AgNi alloys formulated as an area-preserving map. Surface Science, 1990, 225, 319-330.	0.8	44
110	Surface segregation in PtRh alloys revisited in the framework of the tight-binding Ising model. Surface Science, 1990, 236, 398-408.	0.8	50
111	Surface Segregation in Transition Metal Alloys. Progress of Theoretical Physics Supplement, 1990, 101, 159-180.	0.2	58
112	Comment on some â€˜â€˜newâ€™â€™ results on the multilayer reconstruction of W(100). Physical Review B, 1989, 40, 6440-6441.	1.1	9
113	Experimental and theoretical investigation of correlation effects in aggregates of Ni and Pd as a function of their size. Journal of Physics Condensed Matter, 1989, 1, 5875-5885.	0.7	5
114	The face dependence of the effective electron mean free path derived from spherical-wave corrections in photoelectron diffraction of W(110) and W(100) surfaces. Journal of Physics Condensed Matter, 1989, 1, 1879-1888.	0.7	8
115	Photoelectron diffraction study of the unreconstructed and the hydrogen reconstructed W(110) surfaces. Solid State Communications, 1989, 69, 1015-1018.	0.9	11
116	Segregation and Ordering at Surfaces of Transition Metal Alloys: The Tight-Binding Ising Model. Europhysics Letters, 1988, 7, 575-580.	0.7	165
117	Surface and bulk core-level lineshapes in tantalum. II. Experiments. Journal of Physics C: Solid State Physics, 1988, 21, 287-295.	1.5	10
118	Clean surface studies by photoelectron diffraction analysed within a single scattering theory. Journal De Physique, 1988, 49, 227-236.	1.8	25
119	Is ordering in PtNi alloys induced by spin-orbit interactions?. Journal of Physics F: Metal Physics, 1987, 17, 1935-1944.	1.6	29
120	Surface and bulk core-level lineshapes in tungsten and tantalum. I. Theory. Journal of Physics C: Solid State Physics, 1987, 20, 2647-2657.	1.5	11
121	Surface-sandwich segregation in Pt-Ni and Ag-Ni alloys: Two different physical origins for the same phenomenon. Physical Review B, 1987, 35, 4338-4344.	1.1	141
122	Experimental and Theoretical Evidence for a Strong Anisotropy of the Surface Debye-Waller Factor as Determined for a Monolayer of Cobalt on Copper (111) by Surface Extended X-Ray-Absorption Fine Structure. Physical Review Letters, 1986, 56, 1272-1275.	2.9	73
123	One- and two-hole excitation spectra of the 1d Hubbard model. Journal of Magnetism and Magnetic Materials, 1986, 54-57, 1029-1030.	1.0	1
124	A 'quenched molecular dynamics' approach to the atomic stability of the (100) face of BCC transition metals. Journal of Physics C: Solid State Physics, 1986, 19, 4463-4472.	1.5	49
125	One- and two-hole excitation spectra of the Hubbard model in the non-half-filled case. Physical Review B, 1986, 34, 5101-5110.	1.1	17
126	Theoretical study of the Debye-Waller factor in surface extended x-ray-absorption fine structure: Influence of an adsorbed monolayer. Physical Review B, 1986, 34, 6662-6668.	1.1	26

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127	Surface core level spectroscopy of transition metals: A new tool for the determination of their surface structure. Surface Science Reports, 1985, 5, 1-85.	3.8	260
128	Influence of spin-orbit coupling on the atomic stability of the (100) face of 5d BCC transition metals. Solid State Communications, 1985, 55, 961-965.	0.9	7
129	Electronic structure and pairwise interactions in substoichiometric transition metal carbides and nitrides. Journal De Physique, 1985, 46, 1001-1015.	1.8	43
130	One-hole excitation spectra of the one-dimensional Hubbard model. Physical Review B, 1985, 32, 2167-2177.	1.1	15
131	Core level spectroscopy of the low index faces of Tantalum. Surface Science, 1985, 162, 46-50.	0.8	27
132	Variation at the surface of mean correlated displacements in fcc transition and noble metals studied by a continued fraction technique. Surface Science, 1985, 162, 126-131.	0.8	8
133	Bulk and surface vibrational and thermodynamical properties of fcc transition and noble metals : a systematic study by the continued fraction technique. Journal De Physique, 1985, 46, 987-1000.	1.8	60
134	Interpretation of Cr (001) photoemission spectra: influence of correlations. Journal of Physics F: Metal Physics, 1984, 14, 1317-1324.	1.6	14
135	Core-level spectroscopy of clean and adsorbate-covered Ta(100). Physical Review B, 1984, 30, 5487-5493.	1.1	40
136	W(100) reconstruction studied by core level spectroscopy. Solid State Communications, 1984, 50, 393-396.	0.9	43
137	Surface core level spectroscopy of the stepped surface. Solid State Communications, 1984, 52, 635-639.	0.9	21
138	Computation of interatomic Green functions for transition metals using continued fraction techniques. Journal De Physique, 1984, 45, 283-290.	1.8	7
139	Electronic structure and relative stabilities of L12 and D022 ordered structures occuring in transition metal alloys. Solid State Communications, 1983, 45, 585-590.	0.9	60
140	Is the (110) face of tungsten reconstructed?. Solid State Communications, 1983, 47, 279-282.	0.9	13
141	Electronic structure and phase stability of A15 transition metals and alloys. Journal of Physics F: Metal Physics, 1983, 13, 2543-2567.	1.6	60
142	Hydrogen chemisorption on tungsten (110) studied by core level spectroscopy. Journal of Physics C: Solid State Physics, 1983, 16, 1555-1566.	1.5	8
143	A theoretical inquiry into the question of W and Ta (100) atomic structures. Journal of Physics C: Solid State Physics, 1983, 16, 2407-2419.	1.5	43
144	Band gaps and asymptotic behaviour of continued fraction coefficients. Journal of Physics C: Solid State Physics, 1982, 15, 2891-2924.	1.5	144

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145	W4f core level shift study on unreconstructed and hydrogen reconstructed W(100) faces. Journal of Physics C: Solid State Physics, 1982, 15, 4023-4032.	1.5	43
146	Electronic structure of NiZr2C16 compound. Journal of Physics F: Metal Physics, 1982, 12, 441-447.	1.6	17
147	Effect of coulomb correlations on energy bands in ferromagnetic transition metals : Ni, Co and Fe. Journal De Physique, 1982, 43, 341-353.	1.8	152
148	Photoemission from transition metals. Vacuum, 1981, 31, 453-454.	1.6	0
149	Coulomb correlations in nickel within the alloy analogy of the Hubbard model. Solid State Communications, 1981, 39, 1113-1116.	0.9	8
150	Electronic structure, pairwise interactions and ordering energies in binary f.c.c. transition metal alloys. Solid State Communications, 1981, 39, 149-153.	0.9	126
151	Study of the W (Ta) core level shifts induced by the adsorption of oxygen on tungsten (tantalum) (110). Journal of Physics C: Solid State Physics, 1981, 14, 3463-3473.	1.5	61
152	Correlation effects on Auger spectra in unfilled d band metals. Journal of Physics C: Solid State Physics, 1981, 14, 4347-4355.	1.5	92
153	Perturbation treatment of correlations in transition metals. Journal De Physique, 1980, 41, 281-289.	1.8	200
154	Influence of Coulomb correlations on the electronic excitations of transition metal surfaces: application to Mo(100) and Ni(100), (110) and (111). Journal of Physics C: Solid State Physics, 1980, 13, 6063-6072.	1.5	14
155	Thermodynamic derivation of the coherent potential approximation and ordering processes in transition alloys. Journal of Physics F: Metal Physics, 1980, 10, 2137-2146.	1.6	34
156	Comments on the electronic structure of nickel. Physical Review B, 1980, 21, 3729-3733.	1.1	135
157	Further comments on the electronic structure of nickel. Physical Review B, 1980, 22, 6472-6473.	1.1	23
158	Generalised perturbation theory in disordered transition metal alloys: application to the self-consistent calculation of ordering energies. Journal of Physics F: Metal Physics, 1978, 8, 1437-1456.	1.6	58
159	Theoretical and Experimental Evidences of Sequential Phase Formation during Sub-Nanometric-Thick Film Reactive Diffusion. Solid State Phenomena, 0, 172-174, 633-639.	0.3	0
160	Surface Segregation Maps Derived from Tight-Binding Ising Model. Solid State Phenomena, 0, 172-174, 1008-1015.	0.3	7
161	Nanometric-Size Effect upon Diffusion and Reaction in Semiconductors: Experimental and Theoretical Investigations. Defect and Diffusion Forum, 0, 323-325, 433-438.	0.4	0
162	Atomistic Model for Ge Condensation under SiGe Oxidation. Defect and Diffusion Forum, 0, 363, 210-216.	0.4	1