

Ezequiel Pedro Marcos Leiva

List of Publications by Year in Descending Order

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The third column is the impact factor (IF) of the journal, and the fourth column is the number of citations of the article.

140
papers

2,236
citations

24
h-index

36
g-index

145
ext. papers

2,416
ext. citations

4.3
avg, IF

5.03
L-index

#	Paper	IF	Citations
140	On how interactions influence kinetic limitations in alkali-ion batteries. Application to Li-ion intercalation into graphite through voltammetric experiments. <i>Journal of Solid State Electrochemistry</i> , 2021 , 25, 2793	2.6	0
139	Role of the solvent in the activation of LiS as cathode material: a DFT study. <i>Journal of Physics Condensed Matter</i> , 2021 , 33,	1.8	1
138	Characterization of amorphous LiSi structures from ReaxFF accelerated exploration of local minima. <i>Physical Chemistry Chemical Physics</i> , 2021 , 23, 16776-16784	3.6	2
137	Kinetic Monte Carlo simulations applied to Li-ion and post Li-ion batteries: a key link in the multi-scale chain. <i>Progress in Energy</i> , 2021 , 3, 042001	7.7	12
136	Modeling of lithium-ion batteries is becoming viral: where to go?. <i>Journal of Solid State Electrochemistry</i> , 2020 , 24, 2117-2120	2.6	6
135	Fractional and integer stages of lithium ion-graphite systems: the role of electrostatic and elastic contributions. <i>Physical Chemistry Chemical Physics</i> , 2020 , 22, 16174-16183	3.6	3
134	Numerical simulations of cyclic voltammetry for lithium-ion intercalation in nanosized systems: finiteness of diffusion versus electrode kinetics. <i>Journal of Solid State Electrochemistry</i> , 2020 , 24, 3279-3287	2.6	5
133	Effect of Temperature on The Kinetics of Electrochemical Insertion of Li-Ions into a Graphite Electrode Studied by Kinetic Monte Carlo. <i>Journal of the Electrochemical Society</i> , 2020 , 167, 013533	3.9	12
132	Kinetic Monte Carlo applied to the electrochemical study of the Li-ion graphite system. <i>Electrochimica Acta</i> , 2020 , 331, 135439	6.7	13
131	Improving the polysulfide barrier by efficient carbon nanofibers coating on separator/cathode for Li-S batteries. <i>Journal of Solid State Electrochemistry</i> , 2020 , 24, 2341-2351	2.6	2
130	Thermodynamics of Nanoparticle Coalescence at Different Temperatures via Well-Tempered Metadynamics. <i>Journal of Physical Chemistry C</i> , 2020 , 124, 24009-24016	3.8	3
129	On the effect of the carbonaceous substrate in the nucleation of Sn nanoparticles for Li-ion anodes: experiments and first principles calculations. <i>Journal of Solid State Electrochemistry</i> , 2018 , 22, 1721-1733	2.6	2
128	Moře Hařsinsky: The Discoverer of Underpotential Deposition. <i>ChemElectroChem</i> , 2018 , 5, 849-854	4.3	6
127	Grand Canonical Monte Carlo Study of Li Intercalation into Graphite. <i>Journal of the Electrochemical Society</i> , 2018 , 165, A2019-A2025	3.9	11
126	A silica/carbon composite as anode for lithium-ion batteries with a large rate capability: Experiment and theoretical considerations. <i>Electrochimica Acta</i> , 2018 , 279, 289-300	6.7	22
125	Energetics of silica lithiation and its applications to lithium ion batteries. <i>Electrochimica Acta</i> , 2018 , 259, 1053-1058	6.7	24
124	The kinetic origin of the Daumas-Hěold model for the Li-ion/graphite intercalation system. <i>Electrochemistry Communications</i> , 2018 , 93, 133-137	5.1	27

123	Simulation of selective thermodynamic deposition in nanoholes. <i>Physical Chemistry Chemical Physics</i> , 2017 , 19, 1601-1609	3.6	2
122	Statistical mechanical modeling of the transition Stage II -&Stage I of Li-ion storage in graphite. A priori vs induced heterogeneity. <i>Electrochimica Acta</i> , 2017 , 245, 569-574	6.7	15
121	Lithium titanate as anode material for lithium ion batteries: Synthesis, post-treatment and its electrochemical response. <i>Journal of Electroanalytical Chemistry</i> , 2017 , 799, 142-155	4.1	38
120	Hydrogen Evolution Reaction on Nanostructures Electrodes - Scenario on Stepped Silver Surfaces. <i>Electrocatalysis</i> , 2017 , 8, 587-593	2.7	1
119	Computational study of nanostructured materials. <i>Current Opinion in Electrochemistry</i> , 2017 , 1, 1-6	7.2	
118	Criticality of the phase transition on stage two in a lattice-gas model of a graphite anode in a lithium-ion battery. <i>Physical Chemistry Chemical Physics</i> , 2017 , 19, 23138-23145	3.6	12
117	Coalescence of Nanoclusters Analyzed by Well-Tempered Metadynamics. Comparison with Straightforward Molecular Dynamics. <i>Journal of Chemical Theory and Computation</i> , 2017 , 13, 3874-3880	6.4	7
116	Shedding Light on the Entropy Change Found for the Transition Stage II-&Stage I of Li-Ion Storage in Graphite. <i>Journal of the Electrochemical Society</i> , 2017 , 164, A6154-A6157	3.9	17
115	Super-Nernstian Shifts of Interfacial Proton-Coupled Electron Transfers: Origin and Effect of Noncovalent Interactions. <i>Journal of Physical Chemistry C</i> , 2016 , 120, 15586-15592	3.8	11
114	What Is Coming Next?. <i>Monographs in Electrochemistry</i> , 2016 , 335-348	0.8	
113	Underpotential Deposition and Related Phenomena at the Nanoscale: Theory and Applications. <i>Monographs in Electrochemistry</i> , 2016 , 277-334	0.8	3
112	Modeling of Metal Electrodeposition at the Nanoscale 2016 , 971-1009		
111	New kinetic insight into the spontaneous oxidation process of lithium in air by EPMA. <i>Applied Surface Science</i> , 2016 , 383, 64-70	6.7	4
110	Monomolecular adsorption on nanoparticles with repulsive interactions: a Monte Carlo study. <i>Physical Chemistry Chemical Physics</i> , 2016 , 18, 14610-8	3.6	8
109	The Role of Tris(2-carboxyethyl)phosphine Reducing Agent in the Controlled Formation of Alkanedithiols Monolayers on Au(111) with Monocoordinated and Bicoordinated Configurations. <i>Langmuir</i> , 2016 , 32, 9428-36	4	3
108	Time Recovery for a Complex Process Using Accelerated Dynamics. <i>Journal of Chemical Theory and Computation</i> , 2015 , 11, 1725-34	6.4	7
107	Wind and solar hydrogen for the potential production of ammonia in the state of Cear�Brazil. <i>International Journal of Hydrogen Energy</i> , 2015 , 40, 9917-9923	6.7	24
106	Study of the natural resource and economic feasibility of the production and delivery of wind hydrogen in the province of C�doba, Argentina. <i>International Journal of Hydrogen Energy</i> , 2015 , 40, 4413-4425	6.7	12

105	Counterion condensation on polyelectrolyte chains adsorbed on charged surfaces. <i>Colloids and Surfaces A: Physicochemical and Engineering Aspects</i> , 2015 , 487, 49-57	5.1	8
104	Underpotential deposition: From planar surfaces to nanoparticles. <i>Surface Science</i> , 2015 , 631, 23-34	1.8	33
103	Monte Carlo simulation of elongating metallic nanowires in the presence of surfactants. <i>Journal of Chemical Physics</i> , 2015 , 143, 244702	3.9	3
102	Mechanical Effects on the Electronic Properties of a Biphenyl-Based Molecular Switch. <i>Journal of Physical Chemistry C</i> , 2015 , 119, 5090-5097	3.8	3
101	Modeling of Metal Electrodeposition at the Nanoscale 2015 , 1-34		
100	First-principles studies concerning optimization of hydrogen storage in nanoporous reduced graphite oxide. <i>International Journal of Hydrogen Energy</i> , 2014 , 39, 4396-4403	6.7	10
99	A new model for the prediction of oxygen interference in hydrogen storage systems. <i>International Journal of Hydrogen Energy</i> , 2014 , 39, 5899-5905	6.7	4
98	Assessment of the potential for hydrogen production from renewable resources in Argentina. <i>International Journal of Hydrogen Energy</i> , 2014 , 39, 8204-8214	6.7	37
97	Non-covalent interactions at electrochemical interfaces: one model fits all?. <i>Physical Chemistry Chemical Physics</i> , 2014 , 16, 14281-6	3.6	7
96	An experimental and theoretical approach on the effect of presence of oxygen in milled graphite as lithium storage material. <i>Electrochimica Acta</i> , 2014 , 140, 160-167	6.7	13
95	Comment on Surface thermodynamics reconsidered. Derivation of the Gokhshtein relations from the Gibbs potential; and a new approach to surface stress by Stephen Fletcher. <i>Journal of Solid State Electrochemistry</i> , 2014 , 18, 2837-2839	2.6	1
94	Electrochemical, HR-XPS and SERS study of the self-assembly of biphenyl 4,4'-dithiol on Au(111) from solution phase. <i>Surface Science</i> , 2014 , 630, 101-108	1.8	10
93	First-principles studies of lithium storage in reduced graphite oxide. <i>Electrochimica Acta</i> , 2014 , 140, 232-237	6.7	17
92	A New Approach for the Calculation on the Entropic and Free Energy Contributions to Underpotential Deposition. <i>ECS Transactions</i> , 2014 , 58, 3-20	1	2
91	Comparative Study of Different Alkali (Na, Li) Titanate Substrates as Active Materials for Anodes of Lithium - Ion Batteries. <i>ECS Transactions</i> , 2014 , 63, 113-128	1	14
90	Curvature effect in the longitudinal unzipping carbon nanotubes. <i>Journal of Solid State Electrochemistry</i> , 2013 , 17, 1189-1200	2.6	7
89	A Reversible Molecular Switch Based on the Biphenyl Structure. <i>Journal of Physical Chemistry C</i> , 2013 , 117, 25724-25732	3.8	15
88	Anchoring sites to the STM tip can explain multiple peaks in single molecule conductance histograms. <i>Physical Chemistry Chemical Physics</i> , 2013 , 15, 1526-31	3.6	12

87	Stretching single atom contacts at multiple subatomic step-length. <i>Physical Chemistry Chemical Physics</i> , 2013 , 15, 12459-65	3.6	8
86	Computer simulation and detailed mean-field approximation applied to adsorption on nanoparticles. <i>Physical Review E</i> , 2013 , 88, 062407	2.4	4
85	Underpotential deposition on free nanoparticles: Its meaning and measurement. <i>Electrochemistry Communications</i> , 2012 , 16, 1-5	5.1	23
84	Thermodynamic stability of electrochemically decorated AuPd core@shell nanoparticles. <i>Electrochimica Acta</i> , 2012 , 76, 424-429	6.7	12
83	Configurational Behavior and Conductance of Alkanedithiol Molecular Wires from Accelerated Dynamics Simulations. <i>Journal of Chemical Theory and Computation</i> , 2012 , 8, 4539-45	6.4	15
82	A Straightforward Approach for the Determination of the Maximum Time Step for the Simulation of Nanometric Metallic Systems. <i>Journal of Chemical Theory and Computation</i> , 2012 , 8, 1744-9	6.4	6
81	The limits of underpotential deposition in the nanoscale. <i>Electrochemistry Communications</i> , 2012 , 21, 14-17	5.1	17
80	In silico and in vitro characterization of phospholipase A ₂ isoforms from soybean (Glycine max). <i>Biochimie</i> , 2012 , 94, 2608-19	4.6	14
79	Computer simulation of adsorption on nanoparticles: the case of attractive interactions. <i>Physical Review E</i> , 2012 , 86, 061602	2.4	8
78	On the selective decoration of facets in metallic nanoparticles. <i>Journal of Materials Research</i> , 2012 , 27, 1777-1786	2.5	15
77	Properties of rotating nanoalloys formed by cluster collision: a computer simulation study. <i>Journal of Chemical Physics</i> , 2011 , 134, 094701	3.9	14
76	Is hydrogen storage possible in metal-doped graphite 2D systems in conditions found on Earth?. <i>Physical Review Letters</i> , 2011 , 107, 158701	7.4	49
75	Quantitative study of non-covalent interactions at the electrode-electrolyte interface using cyanide-modified Pt(111) electrodes. <i>ChemPhysChem</i> , 2011 , 12, 2230-4	3.2	36
74	Interferents for hydrogen storage on a graphene sheet decorated with nickel: A DFT study. <i>International Journal of Hydrogen Energy</i> , 2011 , 36, 3537-3546	6.7	53
73	A model for underpotential deposition in the presence of anions. <i>Journal of Chemical Physics</i> , 2010 , 132, 184703	3.9	10
72	Role of metal contacts in the mechanical properties of molecular nanojunctions: Comparative ab initio study of Au/1,8-octanedithiol and Au/4,4-bipyridine. <i>Physical Review B</i> , 2010 , 81,	3.3	18
71	Kinetic model for the long term stability of contaminated monoatomic nanowires. <i>Physical Review B</i> , 2010 , 81,	3.3	9
70	On the occurrence of stable and supersaturated metastable states in metallic core-shell nanoparticles. <i>Physical Chemistry Chemical Physics</i> , 2010 , 12, 4580-9	3.6	16

69	Effect of chain stiffness on the morphology of polyelectrolyte complexes. A Monte Carlo simulation study. <i>Polymer</i> , 2010 , 51, 3293-3302	3.9	39
68	Analysis of the potential for hydrogen production in the province of Córdoba, Argentina, from wind resources. <i>International Journal of Hydrogen Energy</i> , 2010 , 35, 5952-5956	6.7	35
67	Theoretical studies of preparation of core-shell nanoparticles by electrochemical metal deposition. <i>Electrochimica Acta</i> , 2010 , 55, 8244-8251	6.7	22
66	The origin of the catalysis of hydrogen peroxide reduction by functionalized graphene surfaces: A density functional theory study. <i>Electrochimica Acta</i> , 2010 , 56, 523-530	6.7	50
65	The behavior of single-molecule junctions predicted by atomistic simulations. <i>Electrochemistry Communications</i> , 2009 , 11, 987-989	5.1	4
64	Stochastic model for spontaneous formation of molecular wires. <i>Electrochimica Acta</i> , 2009 , 54, 2977-2982	7	9
63	Substituent Effect on the Mechanical Properties of Au Nanjunctions. <i>Journal of Physical Chemistry C</i> , 2009 , 113, 3850-3854	3.8	10
62	Atomistic computer simulations on the generation of bimetallic nanoparticles. <i>Faraday Discussions</i> , 2008 , 138, 89-104; discussion 119-35, 433-4	3.6	16
61	Polyelectrolyte Adsorption on a Charged Surface. Free Energy Calculation from Monte Carlo Simulations Using Jarzynski Equality. <i>Macromolecules</i> , 2008 , 41, 8267-8274	5.5	17
60	Thermodynamic considerations and computer simulations on the formation of core-shell nanoparticles under electrochemical conditions. <i>Physical Chemistry Chemical Physics</i> , 2008 , 10, 3561-8	3.6	28
59	Spontaneous nanoripple formation on metallic templates. <i>ACS Nano</i> , 2008 , 2, 2531-9	16.7	8
58	Diffusion mechanisms taking place at the early stages of cobalt deposition on Au(111). <i>Journal of Physics Condensed Matter</i> , 2008 , 20, 265010	1.8	3
57	Computer simulation of the effective double layer occurring on a catalyst surface under electro-chemical promotion conditions. <i>Journal of Applied Electrochemistry</i> , 2008 , 38, 1065-1073	2.6	14
56	When do nanowires break? A model for the theoretical study of the long-term stability of monoatomic nanowires. <i>Chemical Physics Letters</i> , 2008 , 460, 261-265	2.5	24
55	Polyelectrolyte Adsorption on a Charged Surface. A Study by Monte Carlo Simulations. <i>Macromolecules</i> , 2007 , 40, 7336-7342	5.5	22
54	Density functional theory study of a graphene sheet modified with titanium in contact with different adsorbates. <i>Physical Review B</i> , 2007 , 76,	3.3	85
53	Theory of electrochemical monoatomic nanowires. <i>Physical Review B</i> , 2006 , 74,	3.3	16
52	Externally applied electric fields on immiscible lipid monolayers: repulsion between condensed domains precludes domain migration. <i>Langmuir</i> , 2006 , 22, 9664-70	4	12

51	Energetic and entropic contributions to the underpotential/overpotential deposition shifts on single crystal surfaces from lattice dynamics. <i>Electrochimica Acta</i> , 2006 , 51, 3526-3536	6.7	18
50	Low-dimensional metallic nanostructures and their electrochemical relevance: Energetics and phenomenological approach. <i>Surface Science</i> , 2006 , 600, 4475-4483	1.8	6
49	Monte Carlo simulation of metal deposition on foreign substrates. <i>Surface Science</i> , 2006 , 600, 4741-4751	1.8	12
48	Aggregation of casein micelles by interactions with chitosans: a study by Monte Carlo simulations. <i>Journal of Agricultural and Food Chemistry</i> , 2005 , 53, 459-63	5.7	24
47	Reversible precipitation of casein micelles with a cationic hydroxyethylcellulose. <i>Journal of Agricultural and Food Chemistry</i> , 2005 , 53, 9031-8	5.7	5
46	Collision as a way of forming bimetallic nanoclusters of various structures and chemical compositions. <i>Journal of Chemical Physics</i> , 2005 , 123, 184505	3.9	81
45	On the generation of metal clusters with the electrochemical scanning tunneling microscope. <i>Surface Science</i> , 2005 , 597, 133-155	1.8	26
44	Off lattice Monte-Carlo simulations of low-dimensional surface defects and metal deposits on Pt(1 1 1). <i>Electrochemistry Communications</i> , 2005 , 7, 472-476	5.1	2
43	Kinetic Monte Carlo simulation of Pt discontinuous thin film formation adsorbed on Au. <i>Surface Science</i> , 2005 , 581, L109-L114	1.8	15
42	Effects of tip structure on the generation of metal clusters by an STM tip: a way to control the orientation of nanocrystallites?. <i>Nanotechnology</i> , 2005 , 16, 974-980	3.4	10
41	First principles calculations of mechanical properties of 4,4'-bipyridine attached to Au nanowires. <i>Physical Review Letters</i> , 2005 , 95, 045503	7.4	19
40	Monte Carlo simulation of cluster growth in surface defects induced by the tip of a scanning tunnelling microscope. <i>Surface Science</i> , 2004 , 571, L319-L324	1.8	6
39	Monte Carlo simulation of properties of monolayers and metal islands adsorbed on metallic (111) surfaces. <i>Langmuir</i> , 2004 , 20, 4279-88	4	5
38	NEMCA effect: why are the work function changes of the gas exposed catalyst-electrode surface one-to-one related to the changes in the catalyst working electrode potential?. <i>Journal of Solid State Electrochemistry</i> , 2003 , 7, 588-592	2.6	8
37	On the catalytic activity of palladium clusters generated with the electrochemical scanning tunnelling microscope. <i>Electrochemistry Communications</i> , 2003 , 5, 584-586	5.1	15
36	A combined experimental and theoretical study of the generation of palladium clusters on Au(111) with a scanning tunnelling microscope. <i>Electrochimica Acta</i> , 2003 , 48, 1287-1294	6.7	20
35	The basis for the formation of stable metal clusters on an electrode surface. <i>Nanotechnology</i> , 2003 , 14, 1009-1013	3.4	21
34	On the Reasons for Stepwise Changes in the Tunneling Current across Metallic Nanogaps. <i>Nano Letters</i> , 2003 , 3, 1633-1637	11.5	3

33	Comparative Monte Carlo Study of Monolayer Growth in a Heteroepitaxial System in the Presence of Surface Defects. <i>Langmuir</i> , 2003 , 19, 10538-10549	4	23
32	Model calculations for copper clusters on Au(111) surfaces. <i>Journal of Electroanalytical Chemistry</i> , 2002 , 518, 84-90	4.1	13
31	Theoretical Considerations of Electrochemical Phase Formation for an Ideal Frank-van der Merwe System. <i>Journal of the Electrochemical Society</i> , 2002 , 149, E109	3.9	26
30	On the Stability of Ag/Au(111) Expanded Structures. <i>Langmuir</i> , 2002 , 18, 6628-6632	4	11
29	Kinetic Monte Carlo Study of Electrochemical Growth in a Heteroepitaxial System. <i>Langmuir</i> , 2002 , 18, 9087-9094	4	23
28	2D-drop model applied to the calculation of step formation energies on a (111) substrate. <i>Surface Science</i> , 2002 , 499, L135-L140	1.8	9
27	Relevance of Heterometallic Binding Energy for Metal Underpotential Deposition. <i>Langmuir</i> , 2001 , 17, 2219-2227	4	23
26	Changes in surface stress caused by the adsorption of an epitaxial metal monolayer. <i>Chemical Physics Letters</i> , 2000 , 320, 393-397	2.5	18
25	Simulated annealing prediction of the crystal structure of ternary inorganic compounds using symmetry restrictions. <i>Dalton Transactions RSC</i> , 2000 , 4258-4262		7
24	Simulation Study of Pd Submonolayer Films on Au(hkl) and Pt(hkl) and Their Relationship to Underpotential Deposition. <i>Langmuir</i> , 2000 , 16, 9539-9546	4	22
23	Cu underpotential deposition on Au(111) and Au(100). Can this be explained in terms of the energetics of the Cu/Au system?. <i>Electrochimica Acta</i> , 1999 , 45, 691-697	6.7	30
22	Monte Carlo simulation for the formation and growth of low dimensionality phases during underpotential deposition of Ag on Au(100). <i>Electrochimica Acta</i> , 1999 , 45, 699-712	6.7	34
21	An embedded atom approach to underpotential deposition phenomena. <i>Surface Science</i> , 1999 , 421, 59-728		37
20	The oxygen and chlorine evolution reactions at titanium oxide electrodes modified with platinum. <i>Electrochimica Acta</i> , 1998 , 43, 1785-1794	6.7	36
19	Inclusion of symmetry for the enhanced determination of crystalline structures from powder diffraction data using simulated annealing. <i>Chemical Communications</i> , 1998 , 255-256	5.8	6
18	Theoretical Study about the Adsorption of Lead on (111), (100), (110) Monocrystalline Surfaces of Gold. <i>Zeitschrift Fur Physikalische Chemie</i> , 1994 , 185, 33-50	3.1	6
17	Thermodynamic derivation and model calculations of the metal underpotential dependence on electron work function differences. <i>Journal of Electroanalytical Chemistry</i> , 1993 , 350, 1-14	4.1	15
16	Semi-empirical calculations of the vibrational frequency of carbon monoxide adsorbed on noble metal single-crystal surfaces. <i>Journal of Electroanalytical Chemistry</i> , 1993 , 351, 65-79	4.1	7

15	Electrochemical behaviour of passive zirconium alloys. <i>Electrochimica Acta</i> , 1992 , 37, 281-287	6.7	10
14	An extended Hückel/point dipole model for the calculation of dipole potentials for sp metals. <i>Journal of Electroanalytical Chemistry and Interfacial Electrochemistry</i> , 1991 , 303, 55-63		4
13	Contribution of the metal to the capacitance of the double layer: a self-consistent calculation including pseudopotentials. <i>Chemical Physics Letters</i> , 1991 , 187, 143-148	2.5	20
12	CO adsorbate on Pt(111) single crystal surfaces. <i>Electrochimica Acta</i> , 1991 , 36, 555-561	6.7	52
11	Potentiodynamic and AC Impedance Investigation of Anodic Zirconium Oxide Films. <i>Journal of the Electrochemical Society</i> , 1990 , 137, 524-530	3.9	59
10	A jellium/point dipoles model for water adsorption on Ag(110). <i>Surface Science</i> , 1990 , 227, L121-1124	1.8	8
9	Comparative study of CO adsorbates for different structures of platinum surfaces. <i>Journal of Electroanalytical Chemistry and Interfacial Electrochemistry</i> , 1987 , 227, 199-211		34
8	The effect of adsorbed carbon monoxide on hydrogen adsorption and hydrogen evolution on platinum. <i>Journal of Electroanalytical Chemistry and Interfacial Electrochemistry</i> , 1986 , 215, 357-367		48
7	Voltammetric Electro-Oxidation of Carbon Monoxide Previously Adsorbed on Electrochemically Modified Platinum Electrodes. <i>Journal of the Electrochemical Society</i> , 1986 , 133, 1660-1662	3.9	13
6	The influence of solution composition on the kinetics of reduced CO ₂ electrooxidation at polycrystalline platinum. <i>Journal of Electroanalytical Chemistry and Interfacial Electrochemistry</i> , 1985 , 189, 257-269		11
5	Electrodesorption spectra of residues formed on electrochemically modified polycrystalline platinum from carbon dioxide, formic acid, methanol and ethylene glycol adsorption. <i>Electrochimica Acta</i> , 1985 , 30, 1111-1114	6.7	20
4	Complex kinetic behaviour of reduced CO ₂ electro-oxidation at Pt electrodes. <i>Journal of Electroanalytical Chemistry and Interfacial Electrochemistry</i> , 1983 , 158, 103-114		7
3	Multiple current components for methanol adsorption and electro-oxidation at platinum in acidic solutions. <i>Journal of Electroanalytical Chemistry and Interfacial Electrochemistry</i> , 1983 , 158, 115-130		5
2	The Influence of Platinum Electrode Surface on the Electroadsorption and Electro-Oxidation of Methanol in Acid Solutions. <i>Journal of the Electrochemical Society</i> , 1983 , 130, 1305-1312	3.9	25
1	Computer Simulations of Electrochemical Low-Dimensional Metal Phase Formation		30-60 1