

# Matic LozinÅek

## List of Publications by Year in descending order

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53  
papers

22,025  
citations

236612

25  
h-index

168136

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g-index

58  
all docs

58  
docs citations

58  
times ranked

23789  
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#	ARTICLE	IF	CITATIONS
1	How adsorbed H, O, OH, and Cl affect plain adsorption of imidazole on copper. <i>Corrosion Science</i> , 2022, 205, 110443.	3.0	13
2	Catalytic Stereoconvergent Synthesis of Homochiral $\hat{I}^2$ -CF <sub>3</sub> , $\hat{I}^2$ -SCF <sub>3</sub> , and $\hat{I}^2$ -OCF <sub>3</sub> Benzylic Alcohols. <i>ACS Organic &amp; Inorganic Au</i> , 2022, 2, 396-404.	1.9	12
3	Simplistic correlations between molecular electronic properties and inhibition efficiencies: Do they really exist?. <i>Corrosion Science</i> , 2021, 179, 108856.	3.0	86
4	Mixed Noble Gas Compounds of Krypton(II) and Xenon(VI); [F <sub>5</sub> Xe(FKrF)AsF <sub>6</sub> ] and [F <sub>5</sub> Xe(FKrF) <sub>2</sub> AsF <sub>6</sub> ]. <i>Angewandte Chemie - International Edition</i> , 2021, 60, 8149-8156.	7.2	20
5	Mixed Noble Gas Compounds of Krypton(II) and Xenon(VI); [F <sub>5</sub> Xe(FKrF)AsF <sub>6</sub> ] and [F <sub>5</sub> Xe(FKrF) <sub>2</sub> AsF <sub>6</sub> ]. <i>Angewandte Chemie</i> , 2021, 133, 8230-8237.	1.6	2
6	Electrochemical Performance and Mechanism of Calcium Metal Organic Battery. <i>Batteries and Supercaps</i> , 2021, 4, 214-220.	2.4	44
7	Study Of Mercaptobenzimidazoles As Inhibitors For Copper Corrosion: Down to the Molecular Scale. <i>Journal of the Electrochemical Society</i> , 2021, 168, 051504.	1.3	18
8	Expanding the boron peroxide chemistry on BODIPY scaffold. <i>Dyes and Pigments</i> , 2021, 190, 109290.	2.0	3
9	Molecular modeling of organic corrosion inhibitors: Calculations, pitfalls, and conceptualization of molecule surface bonding. <i>Corrosion Science</i> , 2021, 193, 109650.	3.0	70
10	Corrosion resistance of crystalline and amorphous CuZr alloys in NaCl aqueous environment and effect of corrosion inhibitors. <i>Journal of Alloys and Compounds</i> , 2021, 879, 160464.	2.8	12
11	How relevant are molecular electronic parameters for predicting corrosion inhibition efficiency: imidazoles as corrosion inhibitors of Cu/Zr materials in NaCl solution. <i>Corrosion Science</i> , 2021, 193, 109900.	3.0	16
12	Nitrosonium tetrafluoridoborate, NOBF <sub>4</sub> . <i>IUCrData</i> , 2021, 6, .	0.1	1
13	Towards dry and contaminant free Ca(BF <sub>4</sub> ) <sub>2</sub> -based electrolytes for Ca plating. <i>Journal of Power Sources Advances</i> , 2020, 6, 100032.	2.6	7
14	DFT study of aqueous-phase adsorption of cysteine and penicillamine on Fe(110): Role of bond-breaking upon adsorption. <i>Applied Surface Science</i> , 2020, 514, 145896.	3.1	34
15	On the importance of time-resolved electrochemical evaluation in corrosion inhibitor-screening studies. <i>Npj Materials Degradation</i> , 2020, 4, .	2.6	18
16	<i>trans</i> -Diastereoselective Ru(II)-Catalyzed Asymmetric Transfer Hydrogenation of $\hat{I}^{\pm}$ -Acetamido Benzocyclic Ketones via Dynamic Kinetic Resolution. <i>Organic Letters</i> , 2019, 21, 3644-3648.	2.4	34
17	New insights into adsorption bonding of imidazole: A viable C2-H bond cleavage on copper surfaces. <i>Applied Surface Science</i> , 2019, 479, 463-468.	3.1	13
18	Insight into the Bonding of Silanols to Oxidized Aluminum Surfaces. <i>Journal of Physical Chemistry C</i> , 2018, 122, 9417-9431.	1.5	25

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19	Fluorinated reduced graphene oxide as a protective layer on the metallic lithium for application in the high energy batteries. <i>Scientific Reports</i> , 2018, 8, 5819.	1.6	51
20	Manifestations of Weak O-H...F Hydrogen Bonding in M(H <sub>2</sub> O) <sub>n</sub> (B <sub>12</sub> F <sub>12</sub> ) Salt Hydrates: Unusually Sharp Fourier Transform Infrared 1/2(OH) Bands and Latent Porosity (M = Mg, Ba, Co, Ni, Zn). <i>Inorganic Chemistry</i> , 2018, 57, 14983-15000.	1.9	11
21	Reactivity of VOF <sub>3</sub> with N-Heterocyclic Carbene and Imidazolium Fluoride: Analysis of Ligand-VOF <sub>3</sub> Bonding with Evidence of a Minute $\pi$ Back-Donation of Fluoride. <i>Inorganic Chemistry</i> , 2018, 57, 13866-13879.	1.9	14
22	DFT Study of Azole Corrosion Inhibitors on Cu <sub>2</sub> O Model of Oxidized Copper Surfaces: I. Molecule-Surface and Cl-Surface Bonding. <i>Metals</i> , 2018, 8, 310.	1.0	30
23	DFT Study of Azole Corrosion Inhibitors on Cu <sub>2</sub> O Model of Oxidized Copper Surfaces: II. Lateral Interactions and Thermodynamic Stability. <i>Metals</i> , 2018, 8, 311.	1.0	14
24	Coordination of KrF <sub>2</sub> to a Naked Metal Cation, Mg <sup>2+</sup> . <i>Angewandte Chemie - International Edition</i> , 2017, 56, 6251-6254.	7.2	16
25	How relevant is the adsorption bonding of imidazoles and triazoles for their corrosion inhibition of copper?. <i>Corrosion Science</i> , 2017, 124, 25-34.	3.0	64
26	Latent Porosity in Alkali-Metal M <sub>2</sub> B <sub>12</sub> F <sub>12</sub> Salts: Structures and Rapid Room-Temperature Hydration/Dehydration Cycles. <i>Inorganic Chemistry</i> , 2017, 56, 12023-12041.	1.9	13
27	Coordination of KrF <sub>2</sub> to a Naked Metal Cation, Mg <sup>2+</sup> . <i>Angewandte Chemie</i> , 2017, 129, 6347-6350.	1.6	7
28	Origin of Surprising Attractive Interactions between Electronegative Oxygen Adatoms on Aluminum Surfaces. <i>Journal of Physical Chemistry C</i> , 2016, 120, 25915-25922.	1.5	9
29	The world of krypton revisited. <i>Nature Chemistry</i> , 2016, 8, 732-732.	6.6	5
30	Hydroxylammonium Tetrafluoridooxidovanadate(V) $\cdot$ (NH <sub>3</sub> OH)[VOF <sub>4</sub> ]. <i>Acta Chimica Slovenica</i> , 2015, 62, 378-384.	0.2	3
31	Ab initio modeling of the bonding of benzotriazole corrosion inhibitor to reduced and oxidized copper surfaces. <i>Faraday Discussions</i> , 2015, 180, 415-438.	1.6	106
32	The roles of mercapto, benzene, and methyl groups in the corrosion inhibition of imidazoles on copper: II. Inhibitor-copper bonding. <i>Corrosion Science</i> , 2015, 98, 457-470.	3.0	109
33	A DFT study of adsorption of imidazole, triazole, and tetrazole on oxidized copper surfaces: Cu <sub>2</sub> O(111) and Cu <sub>2</sub> O(111)-w/o-Cu <sub>2</sub> CUS. <i>Physical Chemistry Chemical Physics</i> , 2015, 17, 28602-28615.	1.3	45
34	Fluorinated Reduced Graphene Oxide as an Interlayer in Li-S Batteries. <i>Chemistry of Materials</i> , 2015, 27, 7070-7081.	3.2	124
35	The Effect of Surface Geometry of Copper on Adsorption of Benzotriazole and Cl. Part I. <i>Journal of Physical Chemistry C</i> , 2014, 118, 933-943.	1.5	42
36	Silver(I) tetrafluoridooxidovanadate(V) - Ag[VOF <sub>4</sub> ]. <i>Acta Chimica Slovenica</i> , 2014, 61, 542-7.	0.2	1

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37	Synthesis and Crystal Structure of Triammine Pentafluorido Tantalum(V) [TaF <sub>5</sub> (NH <sub>3</sub> ) <sub>3</sub> ]. Zeitschrift Fur Anorganische Und Allgemeine Chemie, 2013, 639, 2586-2588.	0.6	8
38	Formation and structure of inhibitive molecular film of imidazole on iron surface. Corrosion Science, 2013, 68, 195-203.	3.0	87
39	The relation between adsorption bonding and corrosion inhibition of azole molecules on copper. Corrosion Science, 2013, 73, 7-17.	3.0	90
40	Lead Fluoridooxidovanadates(V), Pb(V <sub>2</sub> O <sub>8</sub> F <sub>8</sub> ), Pb(V <sub>5</sub> O <sub>18</sub> F <sub>18</sub> ), and Mixed Valent Fluoridooxidovanadate(IV, V), Pb <sub>3</sub> F(V <sub>4</sub> O <sub>3</sub> F <sub>18</sub> ). Zeitschrift Fur Anorganische Und Allgemeine Chemie, 2012, 638, 2123-2128.	0.6	3
41	On the HSAB based estimate of charge transfer between adsorbates and metal surfaces. Chemical Physics, 2012, 393, 1-12.	0.9	283
42	DFT Study of Interaction of Azoles with Cu(111) and Al(111) Surfaces: Role of Azole Nitrogen Atoms and Dipole-Dipole Interactions. Journal of Physical Chemistry C, 2011, 115, 24189-24197.	1.5	159
43	Electrostatic model for treating long-range lateral interactions between polar molecules adsorbed on metal surfaces. Physical Review B, 2011, 84, .	1.1	52
44	DFT study of gas-phase adsorption of benzotriazole on Cu(111), Cu(100), Cu(110), and low coordinated defects thereon. Physical Chemistry Chemical Physics, 2011, 13, 20408.	1.3	69
45	On the consistent use of electrophilicity index and HSAB-based electron transfer and its associated change of energy parameters. Chemical Physics Letters, 2011, 507, 181-184.	1.2	52
46	Syntheses, structures and Raman spectra of Cd(BF <sub>4</sub> )(AF <sub>6</sub> ) (A=Ta, Bi) compounds. Journal of Fluorine Chemistry, 2011, 132, 767-771.	0.9	14
47	Is the analysis of molecular electronic structure of corrosion inhibitors sufficient to predict the trend of their inhibition performance. Electrochimica Acta, 2010, 56, 745-755.	2.6	258
48	What Determines the Inhibition Effectiveness of ATA, BTAH, and BTAOH Corrosion Inhibitors on Copper?. Journal of the American Chemical Society, 2010, 132, 16657-16668.	6.6	278
49	HF molecules and poly(hydrogen fluoride) anions as ligands to metal centers. Journal of Fluorine Chemistry, 2009, 130, 1093-1098.	0.9	12
50	Syntheses, crystal structures and Raman spectra of Ba(BF <sub>4</sub> )(PF <sub>6</sub> ), Ba(BF <sub>4</sub> )(AsF <sub>6</sub> ) and Ba <sub>2</sub> (BF <sub>4</sub> ) <sub>2</sub> (AsF <sub>6</sub> )(H <sub>3</sub> F <sub>4</sub> ); the first examples of metal salts containing simultaneously tetrahedral BF <sub>4</sub> <sup>-</sup> and octahedral AF <sub>6</sub> <sup>-</sup> anions. Journal of Solid State Chemistry, 2009, 182, 2897-2903.	1.4	10
51	QUANTUM ESPRESSO: a modular and open-source software project for quantum simulations of materials. Journal of Physics Condensed Matter, 2009, 21, 395502.	0.7	18,183
52	Methane Dehydrogenation on Rh@Cu(111): A First-Principles Study of a Model Catalyst. Journal of the American Chemical Society, 2006, 128, 12448-12454.	6.6	60
53	XCrySDen—a new program for displaying crystalline structures and electron densities. Journal of Molecular Graphics and Modelling, 1999, 17, 176-179.	1.3	1,315