

# Francesco Musiani

## List of Publications by Year in descending order

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148  
papers

6,040  
citations

61984

43  
h-index

88630

70  
g-index

155  
all docs

155  
docs citations

155  
times ranked

5763  
citing authors

#	ARTICLE	IF	CITATIONS
1	Relevance of ARID1A Mutations in Endometrial Carcinomas. <i>Diagnostics</i> , 2022, 12, 592.	2.6	6
2	New Insights on Rotenone Resistance of Complex I Induced by the m.11778G>A/MT-ND4 Mutation Associated with Leber's Hereditary Optic Neuropathy. <i>Molecules</i> , 2022, 27, 1341.	3.8	3
3	Cloning the barley <i>nec3</i> disease lesion mimic mutant using complementation by sequencing. <i>Plant Genome</i> , 2022, , e20187.	2.8	0
4	Inhibition of Urease, a Ni-Enzyme: The Reactivity of a Key Thiol With Mono- and Di-Substituted Catechols Elucidated by Kinetic, Structural, and Theoretical Studies. <i>Angewandte Chemie - International Edition</i> , 2021, 60, 6029-6035.	13.8	12
5	Inhibition of Urease, a Ni-Enzyme: The Reactivity of a Key Thiol With Mono- and Di-Substituted Catechols Elucidated by Kinetic, Structural, and Theoretical Studies. <i>Angewandte Chemie</i> , 2021, 133, 6094-6100.	2.0	3
6	Combining Different Docking Engines and Consensus Strategies to Design and Validate Optimized Virtual Screening Protocols for the SARS-CoV-2 3CL Protease. <i>Molecules</i> , 2021, 26, 797.	3.8	14
7	The structural and functional characterization of <i>Malus domestica</i> double bond reductase MdDBR provides insights towards the identification of its substrates. <i>International Journal of Biological Macromolecules</i> , 2021, 171, 89-99.	7.5	6
8	A Blueprint for High Affinity SARS-CoV-2 Mpro Inhibitors from Activity-Based Compound Library Screening Guided by Analysis of Protein Dynamics. <i>ACS Pharmacology and Translational Science</i> , 2021, 4, 1079-1095.	4.9	44
9	Identification of Inhibitors of SARS-CoV-2 3CL-Pro Enzymatic Activity Using a Small Molecule in Vitro Repurposing Screen. <i>ACS Pharmacology and Translational Science</i> , 2021, 4, 1096-1110.	4.9	101
10	Nickel as a virulence factor in the Class I bacterial carcinogen, <i>Helicobacter pylori</i> . <i>Seminars in Cancer Biology</i> , 2021, 76, 143-155.	9.6	14
11	Kinetic and structural analysis of the inactivation of urease by mixed-ligand phosphine halide Ag(I) complexes. <i>Journal of Inorganic Biochemistry</i> , 2021, 218, 111375.	3.5	10
12	Definition of the Binding Architecture to a Target Promoter of HP1043, the Essential Master Regulator of <i>Helicobacter pylori</i> . <i>International Journal of Molecular Sciences</i> , 2021, 22, 7848.	4.1	8
13	Facilitating Nitrification Inhibition through Green, Mechanochemical Synthesis of a Novel Nitrapyrin Complex. <i>Crystal Growth and Design</i> , 2021, 21, 5792-5799.	3.0	10
14	Probing the transport of Ni(II) ions through the internal tunnels of the <i>Helicobacter pylori</i> UreDFG multimeric protein complex. <i>Journal of Inorganic Biochemistry</i> , 2021, 223, 111554.	3.5	6
15	Medicinal Au(III) compounds targeting urease as prospective antimicrobial agents: unveiling the structural basis for enzyme inhibition. <i>Dalton Transactions</i> , 2021, 50, 14444-14452.	3.3	10
16	Structure, dynamics, and function of SrnR, a transcription factor for nickel-dependent gene expression. <i>Metallomics</i> , 2021, 13, .	2.4	4
17	Nickel import and export in the human pathogen <i>Helicobacter pylori</i> , perspectives from molecular modelling. <i>Metallomics</i> , 2021, 13, .	2.4	6
18	Targeting the Protein Tunnels of the Urease Accessory Complex: A Theoretical Investigation. <i>Molecules</i> , 2020, 25, 2911.	3.8	13

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19	Nickel and GTP Modulate <i>Helicobacter pylori</i> UreG Structural Flexibility. <i>Biomolecules</i> , 2020, 10, 1062.	4.0	9
20	Partitioning the structural features that underlie expansin-like and elicitor activities of cerato-platanin. <i>International Journal of Biological Macromolecules</i> , 2020, 165, 2845-2854.	7.5	5
21	The model structure of the copper-dependent ammonia monooxygenase. <i>Journal of Biological Inorganic Chemistry</i> , 2020, 25, 995-1007.	2.6	27
22	The structure-based reaction mechanism of urease, a nickel dependent enzyme: tale of a long debate. <i>Journal of Biological Inorganic Chemistry</i> , 2020, 25, 829-845.	2.6	92
23	DNMT1 mutations leading to neurodegeneration paradoxically reflect on mitochondrial metabolism. <i>Human Molecular Genetics</i> , 2020, 29, 1864-1881.	2.9	19
24	Intrinsic disorder in the nickel-dependent urease network. <i>Progress in Molecular Biology and Translational Science</i> , 2020, 174, 307-330.	1.7	6
25	Multifunctional Urea Cocrystal with Combined Ureolysis and Nitrification Inhibiting Capabilities for Enhanced Nitrogen Management. <i>ACS Sustainable Chemistry and Engineering</i> , 2019, 7, 13369-13378.	6.7	32
26	The Impact of pH on Catalytically Critical Protein Conformational Changes: The Case of the Urease, a Nickel Enzyme. <i>Chemistry - A European Journal</i> , 2019, 25, 12145-12158.	3.3	21
27	Molecular Modelling of the Ni(II)-Responsive <i>Synechocystis</i> PCC 6803 Transcriptional Regulator InrS in the Metal Bound Form. <i>Inorganics</i> , 2019, 7, 76.	2.7	7
28	Soyuretox, an Intrinsically Disordered Polypeptide Derived from Soybean ( <i>Glycine Max</i> ) Ubiquitous Urease with Potential Use as a Biopesticide. <i>International Journal of Molecular Sciences</i> , 2019, 20, 5401.	4.1	8
29	A Solvent-Exposed Cysteine Forms a Peculiar Ni II Binding Site in the Metallochaperone CooT from <i>Rhodospirillum rubrum</i> . <i>Chemistry - A European Journal</i> , 2019, 25, 15351-15360.	3.3	9
30	Urease Inhibitory Potential and Soil Ecotoxicity of Novel "Polyphenols" Deep Eutectic Solvents Formulations. <i>ACS Sustainable Chemistry and Engineering</i> , 2019, 7, 15558-15567.	6.7	23
31	The carbon monoxide dehydrogenase accessory protein CooJ is a histidine-rich multidomain dimer containing an unexpected Ni(II)-binding site. <i>Journal of Biological Chemistry</i> , 2019, 294, 7601-7614.	3.4	16
32	The Structure of the Elusive Urease-Urea Complex Unveils the Mechanism of a Paradigmatic Nickel-Dependent Enzyme. <i>Angewandte Chemie - International Edition</i> , 2019, 58, 7415-7419.	13.8	66
33	The Structure of the Elusive Urease-Urea Complex Unveils the Mechanism of a Paradigmatic Nickel-Dependent Enzyme. <i>Angewandte Chemie</i> , 2019, 131, 7493-7497.	2.0	7
34	Insights into Urease Inhibition by <i>N</i> -( <i>n</i> -Butyl) Phosphoric Triamide through an Integrated Structural and Kinetic Approach. <i>Journal of Agricultural and Food Chemistry</i> , 2019, 67, 2127-2138.	5.2	33
35	Bioinorganic Chemistry of Nickel. <i>Inorganics</i> , 2019, 7, 131.	2.7	5
36	The cytochrome b lysine 329 residue is critical for ubihydroquinone oxidation and proton release at the Q <sub>o</sub> site of bacterial cytochrome bc <sub>1</sub> . <i>Biochimica Et Biophysica Acta - Bioenergetics</i> , 2019, 1860, 167-179.	1.0	4

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37	Novel Dual-Action Plant Fertilizer and Urease Inhibitor: Urea-Catechol Cocrystal. Characterization and Environmental Reactivity. <i>ACS Sustainable Chemistry and Engineering</i> , 2019, 7, 2852-2859.	6.7	42
38	Inhibition Mechanism of Urease by Au(III) Compounds Unveiled by X-ray Diffraction Analysis. <i>ACS Medicinal Chemistry Letters</i> , 2019, 10, 564-570.	2.8	30
39	Predicting ligand binding poses for low-resolution membrane protein models: Perspectives from multiscale simulations. <i>Biochemical and Biophysical Research Communications</i> , 2018, 498, 366-374.	2.1	32
40	<i>Pseudomonas pseudoalcaligenes</i> KF 707 grown with biphenyl expresses a cytochrome caa 3 oxidase that uses cytochrome c 4 as electron donor. <i>FEBS Letters</i> , 2018, 592, 901-915.	2.8	4
41	An Evaluation of Maleic-itaconic Copolymers as Urease Inhibitors. <i>Soil Science Society of America Journal</i> , 2018, 82, 994-1003.	2.2	9
42	Structure and dynamics of <i>Helicobacter pylori</i> nickel-chaperone HypA: an integrated approach using NMR spectroscopy, functional assays and computational tools. <i>Journal of Biological Inorganic Chemistry</i> , 2018, 23, 1309-1330.	2.6	20
43	Application of Molecular Dynamics to the Investigation of Metalloproteins Involved in Metal Homeostasis. <i>European Journal of Inorganic Chemistry</i> , 2018, 2018, 4661-4677.	2.0	12
44	The structure of urease inactivated by Ag( $\text{SCN}$ ): a new paradigm for enzyme inhibition by heavy metals. <i>Dalton Transactions</i> , 2018, 47, 8240-8247.	3.3	54
45	Targeting <i>Helicobacter pylori</i> urease activity and maturation: In-cell high-throughput approach for drug discovery. <i>Biochimica Et Biophysica Acta - General Subjects</i> , 2018, 1862, 2245-2253.	2.4	28
46	Deciphering OPA1 mutations pathogenicity by combined analysis of human, mouse and yeast cell models. <i>Biochimica Et Biophysica Acta - Molecular Basis of Disease</i> , 2018, 1864, 3496-3514.	3.8	36
47	Polymer-Mediated Delivery of siRNAs to Hepatocellular Carcinoma: Variables Affecting Specificity and Effectiveness. <i>Molecules</i> , 2018, 23, 777.	3.8	18
48	Smart urea ionic co-crystals with enhanced urease inhibition activity for improved nitrogen cycle management. <i>Chemical Communications</i> , 2018, 54, 7637-7640.	4.1	41
49	Protein Aggregation and Molecular Crowding. <i>International Review of Cell and Molecular Biology</i> , 2017, 329, 49-77.	3.2	22
50	Strategies to optimize siRNA delivery to hepatocellular carcinoma cells. <i>Expert Opinion on Drug Delivery</i> , 2017, 14, 797-810.	5.0	25
51	Protein Tunnels: The Case of Urease Accessory Proteins. <i>Journal of Chemical Theory and Computation</i> , 2017, 13, 2322-2331.	5.3	25
52	Glutamate Ligation in the Ni(II)- and Co(II)-Responsive <i>Escherichia coli</i> Transcriptional Regulator, RcnR. <i>Inorganic Chemistry</i> , 2017, 56, 6459-6476.	4.0	16
53	Development of a multisite model for Ni(II) ion in solution from thermodynamic and kinetic data. <i>Journal of Computational Chemistry</i> , 2017, 38, 1834-1843.	3.3	11
54	Glucose-1-phosphate uridylyltransferase from <i>Erwinia amylovora</i> : Activity, structure and substrate specificity. <i>Biochimica Et Biophysica Acta - Proteins and Proteomics</i> , 2017, 1865, 1348-1357.	2.3	13

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55	Urease Inhibition in the Presence of <i>N</i> -( <i>n</i> -Butyl)thiophosphoric Triamide, a Suicide Substrate: Structure and Kinetics. <i>Biochemistry</i> , 2017, 56, 5391-5404.	2.5	53
56	Structural analysis of the interaction between Jaburetox, an intrinsically disordered protein, and membrane models. <i>Colloids and Surfaces B: Biointerfaces</i> , 2017, 159, 849-860.	5.0	10
57	The relationship between folding and activity in UreG, an intrinsically disordered enzyme. <i>Scientific Reports</i> , 2017, 7, 5977.	3.3	34
58	Inactivation of urease by catechol: Kinetics and structure. <i>Journal of Inorganic Biochemistry</i> , 2017, 166, 182-189.	3.5	57
59	Mathematical Modeling of Drug Release from Natural Polysaccharides Based Matrices. <i>Natural Product Communications</i> , 2017, 12, 1934578X1701200.	0.5	4
60	Characterization and 1.57 Å resolution structure of the key fire blight phosphatase AmsI from <i>Erwinia amylovora</i> . <i>Acta Crystallographica Section F, Structural Biology Communications</i> , 2016, 72, 903-910.	0.8	8
61	Isothermal Titration Calorimetry to Characterize Enzymatic Reactions. <i>Methods in Enzymology</i> , 2016, 567, 215-236.	1.0	20
62	Surface plasmon resonance and isothermal titration calorimetry to monitor the Ni(II)-dependent binding of <i>Helicobacter pylori</i> NikR to DNA. <i>Analytical and Bioanalytical Chemistry</i> , 2016, 408, 7971-7980.	3.7	14
63	Nickel impact on human health: An intrinsic disorder perspective. <i>Biochimica Et Biophysica Acta - Proteins and Proteomics</i> , 2016, 1864, 1714-1731.	2.3	151
64	On the role of a specific insert in acetate permeases (ActP) for tellurite uptake in bacteria: Functional and structural studies. <i>Journal of Inorganic Biochemistry</i> , 2016, 163, 103-109.	3.5	10
65	Inactivation of urease by 1,4-benzoquinone: chemistry at the protein surface. <i>Dalton Transactions</i> , 2016, 45, 5455-5459.	3.3	61
66	Kinetic and structural studies reveal a unique binding mode of sulfite to the nickel center in urease. <i>Journal of Inorganic Biochemistry</i> , 2016, 154, 42-49.	3.5	42
67	Conformational ensemble of human $\alpha$ -synuclein physiological form predicted by molecular simulations. <i>Physical Chemistry Chemical Physics</i> , 2016, 18, 5702-5706.	2.8	32
68	Transient Interactions of a Cytosolic Protein with Macromolecular and Vesicular Cosolutes: Unspecific and Specific Effects. <i>ChemBioChem</i> , 2015, 16, 2633-2645.	2.6	10
69	Evolution of Macromolecular Docking Techniques: The Case Study of Nickel and Iron Metabolism in Pathogenic Bacteria. <i>Molecules</i> , 2015, 20, 14265-14292.	3.8	3
70	Polysaccharides for the Delivery of Antitumor Drugs. <i>Materials</i> , 2015, 8, 2569-2615.	2.9	110
71	On the interaction of <i>Helicobacter pylori</i> NikR, a Ni(II)-responsive transcription factor, with the urease operator: in solution and in silico studies. <i>Journal of Biological Inorganic Chemistry</i> , 2015, 20, 1021-1037.	2.6	18
72	Intrinsic disorder and metal binding in UreG proteins from Archae hyperthermophiles: GTPase enzymes involved in the activation of Ni(II) dependent urease. <i>Journal of Biological Inorganic Chemistry</i> , 2015, 20, 739-755.	2.6	19

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73	Pliable natural biocide: Jaburetox is an intrinsically disordered insecticidal and fungicidal polypeptide derived from jack bean urease. <i>FEBS Journal</i> , 2015, 282, 1043-1064.	4.7	30
74	Nickel-responsive transcriptional regulators. <i>Metallomics</i> , 2015, 7, 1305-1318.	2.4	40
75	Evidence for a Transient Additional Ligand Binding Site in the TAS2R46 Bitter Taste Receptor. <i>Journal of Chemical Theory and Computation</i> , 2015, 11, 4439-4449.	5.3	70
76	Novel Lipid and Polymeric Materials as Delivery Systems for Nucleic Acid Based Drugs. <i>Current Drug Metabolism</i> , 2015, 16, 427-452.	1.2	26
77	Dynamic characterization and substrate binding of cis-2,3-dihydrobiphenyl-2,3-diol dehydrogenase an enzyme used in bioremediation. <i>Journal of Molecular Modeling</i> , 2014, 20, 2531.	1.8	6
78	FeON-FeOFF: the <i>Helicobacter pylori</i> Fur regulator commutates iron-responsive transcription by discriminative readout of opposed DNA grooves. <i>Nucleic Acids Research</i> , 2014, 42, 3138-3151.	14.5	38
79	The conformational response to Zn(II) and Ni(II) binding of <i>Sporosarcina pasteurii</i> UreG, an intrinsically disordered GTPase. <i>Journal of Biological Inorganic Chemistry</i> , 2014, 19, 1341-1354.	2.6	22
80	Nickel binding properties of <i>Helicobacter pylori</i> UreF, an accessory protein in the nickel-based activation of urease. <i>Journal of Biological Inorganic Chemistry</i> , 2014, 19, 319-334.	2.6	40
81	Molecular Dynamics Simulations Identify Time Scale of Conformational Changes Responsible for Conformational Selection in Molecular Recognition of HIV-1 Transactivation Responsive RNA. <i>Journal of the American Chemical Society</i> , 2014, 136, 15631-15637.	13.7	35
82	Fluoride inhibition of <i>Sporosarcina pasteurii</i> urease: structure and thermodynamics. <i>Journal of Biological Inorganic Chemistry</i> , 2014, 19, 1243-1261.	2.6	58
83	Topological characterization of a bacterial cellulose acrylic acid polymeric matrix. <i>European Journal of Pharmaceutical Sciences</i> , 2014, 62, 326-333.	4.0	15
84	Molecular landscape of the interaction between the urease accessory proteins UreE and UreG. <i>Biochimica Et Biophysica Acta - Proteins and Proteomics</i> , 2014, 1844, 1662-1674.	2.3	44
85	Nonredox Nickel Enzymes. <i>Chemical Reviews</i> , 2014, 114, 4206-4228.	47.7	235
86	Hot Biological Catalysis: Isothermal Titration Calorimetry to Characterize Enzymatic Reactions. <i>Journal of Visualized Experiments</i> , 2014, , .	0.3	7
87	Chemosensorial G-proteins-Coupled Receptors: A Perspective from Computational Methods. <i>Advances in Experimental Medicine and Biology</i> , 2014, 805, 441-457.	1.6	4
88	Structure of the Ure-Ure-UreE complex in <i>Helicobacter pylori</i> : a model study. <i>Journal of Biological Inorganic Chemistry</i> , 2013, 18, 571-577.	2.6	17
89	The crystal structure of <i>Sporosarcina pasteurii</i> urease in a complex with citrate provides new hints for inhibitor design. <i>Journal of Biological Inorganic Chemistry</i> , 2013, 18, 391-399.	2.6	49
90	Selectivity of Ni(II) and Zn(II) binding to <i>Sporosarcina pasteurii</i> UreE, a metallochaperone in the urease assembly: a calorimetric and crystallographic study. <i>Journal of Biological Inorganic Chemistry</i> , 2013, 18, 1005-1017.	2.6	21

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91	Conformational Fluctuations of UreG, an Intrinsically Disordered Enzyme. <i>Biochemistry</i> , 2013, 52, 2949-2954.	2.5	33
92	Nickel and Human Health. <i>Metal Ions in Life Sciences</i> , 2013, 13, 321-357.	2.8	71
93	Urease. , 2013, , 2287-2292.		1
94	GOMoDo: A GPCRs Online Modeling and Docking Webserver. <i>PLoS ONE</i> , 2013, 8, e74092.	2.5	84
95	Structure/Function Relationships of Phospholipases C Beta. <i>Current Protein and Peptide Science</i> , 2013, 14, 650-657.	1.4	11
96	Crystallographic and X-ray absorption spectroscopic characterization of <i>Helicobacter pylori</i> UreE bound to Ni <sup>2+</sup> and Zn <sup>2+</sup> reveals a role for the disordered C-terminal arm in metal trafficking. <i>Biochemical Journal</i> , 2012, 441, 1017-1035.	3.7	52
97	Denaturant-Induced Conformational Transitions in Intrinsically Disordered Proteins. , 2012, 896, 197-213.		4
98	Insights in the (un)structural organization of <i>Bacillus pasteurii</i> UreG, an intrinsically disordered GTPase enzyme. <i>Molecular BioSystems</i> , 2012, 8, 220-228.	2.9	44
99	Engineered biosealant strains producing inorganic and organic biopolymers. <i>Journal of Biotechnology</i> , 2012, 161, 181-189.	3.8	19
100	Metal Ion-Mediated DNA-Protein Interactions. <i>Metal Ions in Life Sciences</i> , 2012, 10, 135-170.	2.8	21
101	Biochemical and structural studies on native and recombinant <i>Glycine max</i> UreG: a detailed characterization of a plant urease accessory protein. <i>Plant Molecular Biology</i> , 2012, 78, 461-475.	3.9	32
102	Unraveling the <i>Helicobacter pylori</i> UreG zinc binding site using X-ray absorption spectroscopy (XAS) and structural modeling. <i>Journal of Biological Inorganic Chemistry</i> , 2012, 17, 353-361.	2.6	32
103	Holo-Ni <sup>2+</sup> <i>Helicobacter pylori</i> NikR contains four square-planar nickel-binding sites at physiological pH. <i>Dalton Transactions</i> , 2011, 40, 7831.	3.3	28
104	Zinc Inhibition of Bacterial Cytochrome <i>bc<sub>1</sub></i> Reveals the Role of Cytochrome <i>b</i> E295 in Proton Release at the Q <sub>o</sub> Site. <i>Biochemistry</i> , 2011, 50, 4263-4272.	2.5	30
105	Chemistry of Ni <sup>2+</sup> in Urease: Sensing, Trafficking, and Catalysis. <i>Accounts of Chemical Research</i> , 2011, 44, 520-530.	15.6	224
106	Model Structures of <i>Helicobacter pylori</i> UreD(H) Domains: A Putative Molecular Recognition Platform. <i>Journal of Chemical Information and Modeling</i> , 2011, 51, 1513-1520.	5.4	14
107	Interaction of Selenoprotein W with 14-3-3 Proteins: A Computational Approach. <i>Journal of Proteome Research</i> , 2011, 10, 968-976.	3.7	29
108	Computational Study of the DNA-Binding Protein <i>Helicobacter pylori</i> NikR: The Role of Ni <sup>2+</sup> Francesco Musiani and Branimir BertoÅa contributed equally to the simulations presented here.. <i>Journal of Chemical Theory and Computation</i> , 2010, 6, 3503-3515.	5.3	32

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109	The RNA Hydrolysis and the Cytokinin Binding Activities of PR-10 Proteins Are Differently Performed by Two Isoforms of the Pru p 1 Peach Major Allergen and Are Possibly Functionally Related. <i>Plant Physiology</i> , 2009, 150, 1235-1247.	4.8	66
110	<i>Helicobacter pylori</i> UreE, a urease accessory protein: specific Ni <sup>2+</sup> - and Zn <sup>2+</sup> -binding properties and interaction with its cognate UreG. <i>Biochemical Journal</i> , 2009, 422, 91-100.	3.7	83
111	Bifidobacterial enolase, a cell surface receptor for human plasminogen involved in the interaction with the host. <i>Microbiology (United Kingdom)</i> , 2009, 155, 3294-3303.	1.8	110
112	Zn <sup>2+</sup> -linked dimerization of UreG from <i>Helicobacter pylori</i> , a chaperone involved in nickel trafficking and urease activation. <i>Proteins: Structure, Function and Bioinformatics</i> , 2009, 74, 222-239.	2.6	73
113	Conformational Equilibria in Monomeric Î±-Synuclein at the Single-Molecule Level. <i>PLoS Biology</i> , 2008, 6, e6.	5.6	181
114	Structural Characterization of Binding of Cu(II) to Tau Protein. <i>Biochemistry</i> , 2008, 47, 10841-10851.	2.5	85
115	High-Affinity Ni <sup>2+</sup> Binding Selectively Promotes Binding of <i>Helicobacter pylori</i> NikR to Its Target Urease Promoter. <i>Journal of Molecular Biology</i> , 2008, 383, 1129-1143.	4.2	63
116	The Ni <sup>2+</sup> binding properties of <i>Helicobacter pylori</i> NikR. <i>Chemical Communications</i> , 2007, , 3649.	4.1	47
117	Biochemical Studies on <i>Mycobacterium tuberculosis</i> UreG and Comparative Modeling Reveal Structural and Functional Conservation among the Bacterial UreG Family. <i>Biochemistry</i> , 2007, 46, 3171-3182.	2.5	56
118	Urease: Recent Insights on the Role of Nickel. , 2007, , 241-277.		11
119	A model-based proposal for the role of UreF as a GTPase-activating protein in the urease active site biosynthesis. <i>Proteins: Structure, Function and Bioinformatics</i> , 2007, 68, 749-761.	2.6	36
120	The Nickel Site of <i>Bacillus pasteurii</i> UreE, a Urease Metallo-Chaperone, As Revealed by Metal-Binding Studies and X-ray Absorption Spectroscopy. <i>Biochemistry</i> , 2006, 45, 6495-6509.	2.5	49
121	Intrinsically Disordered Structure of <i>Bacillus pasteurii</i> UreG As Revealed by Steady-State and Time-Resolved Fluorescence Spectroscopy. <i>Biochemistry</i> , 2006, 45, 8918-8930.	2.5	47
122	An Italian contribution to structural genomics: Understanding metalloproteins. <i>Coordination Chemistry Reviews</i> , 2006, 250, 1419-1450.	18.8	14
123	Jack bean ( <i>Canavalia ensiformis</i> ) urease. Probing acid-base groups of the active site by pH variation. <i>Plant Physiology and Biochemistry</i> , 2005, 43, 651-658.	5.8	74
124	High potential iron-sulfur proteins and their role as soluble electron carriers in bacterial photosynthesis: tale of a discovery. <i>Photosynthesis Research</i> , 2005, 85, 115-131.	2.9	23
125	UreG, a Chaperone in the Urease Assembly Process, Is an Intrinsically Unstructured GTPase That Specifically Binds Zn <sup>2+</sup> . <i>Journal of Biological Chemistry</i> , 2005, 280, 4684-4695.	3.4	91
126	Structure of the Intermolecular Complex between Plastocyanin and Cytochrome f from Spinach*. <i>Journal of Biological Chemistry</i> , 2005, 280, 18833-18841.	3.4	20

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127	Nickel trafficking: insights into the fold and function of UreE, a urease metallochaperone. <i>Journal of Inorganic Biochemistry</i> , 2004, 98, 803-813.	3.5	43
128	The Asn 38 <sup>→</sup> Cys 84 H-Bond in Plastocyanin. <i>Journal of Physical Chemistry B</i> , 2004, 108, 7495-7499.	2.6	7
129	Molecular Details of Urease Inhibition by Boric Acid: Insights into the Catalytic Mechanism. <i>Journal of the American Chemical Society</i> , 2004, 126, 3714-3715.	13.7	142
130	Electron Transfer from HiPIP to the Photooxidized Tetraheme Cytochrome Subunit of <i>Allochrochromatium vinosum</i> Reaction Center: New Insights from Site-Directed Mutagenesis and Computational Studies. <i>Biochemistry</i> , 2004, 43, 437-445.	2.5	10
131	Molecular characterization of <i>Bacillus pasteurii</i> UreE, a metal-binding chaperone for the assembly of the urease active site. <i>Journal of Biological Inorganic Chemistry</i> , 2002, 7, 623-631.	2.6	39
132	NMR Solution Structure, Backbone Mobility, and Homology Modeling of c-Type Cytochromes from Gram-Positive Bacteria. <i>ChemBioChem</i> , 2002, 3, 299-310.	2.6	23
133	The First Solution Structure of a Paramagnetic Copper(II) Protein: The Case of Oxidized Plastocyanin from the Cyanobacterium <i>Synechocystis</i> PCC6803. <i>Journal of the American Chemical Society</i> , 2001, 123, 2405-2413.	13.7	65
134	Structure-based computational study of the catalytic and inhibition mechanisms of urease. <i>Journal of Biological Inorganic Chemistry</i> , 2001, 6, 300-314.	2.6	110
135	Backbone Dynamics of Plastocyanin in Both Oxidation States. <i>Journal of Biological Chemistry</i> , 2001, 276, 47217-47226.	3.4	50
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