

Daniel Reta

List of Publications by Year in descending order

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3,416
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304602

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docs citations

44
times ranked

2551
citing authors

#	ARTICLE	IF	CITATIONS
1	Ultrahard magnetism from mixed-valence dilanthanide complexes with metal-metal bonding. <i>Science</i> , 2022, 375, 198-202.	6.0	246
2	Studies of the Temperature Dependence of the Structure and Magnetism of a Hexagonal-Bipyramidal Dysprosium(III) Single-Molecule Magnet. <i>Inorganic Chemistry</i> , 2022, 61, 227-235.	1.9	13
3	Opening Magnetic Hysteresis by Axial Ferromagnetic Coupling: From Mono-Decker to Double-Decker Metallocrown. <i>Angewandte Chemie - International Edition</i> , 2021, 60, 5299-5306.	7.2	62
4	Opening Magnetic Hysteresis by Axial Ferromagnetic Coupling: From Mono-Decker to Double-Decker Metallocrown. <i>Angewandte Chemie</i> , 2021, 133, 5359-5366.	1.6	8
5	Ab Initio Prediction of High-Temperature Magnetic Relaxation Rates in Single-Molecule Magnets. <i>Journal of the American Chemical Society</i> , 2021, 143, 5943-5950.	6.6	110
6	A Cost-Effective Semi-Ab Initio Approach to Model Relaxation in Rare-Earth Single-Molecule Magnets. <i>Journal of Physical Chemistry Letters</i> , 2021, 12, 8826-8832.	2.1	35
7	Eight coordinated mononuclear dysprosium complexes of heptadentate aminophenol ligands: the influence of the phenol substituents and the ancillary donors on the magnetic relaxation. <i>Dalton Transactions</i> , 2021, 50, 15878-15887.	1.6	3
8	Strangely attractive: Collaboration and feedback in the field of molecular magnetism. <i>International Journal of Quantum Chemistry</i> , 2020, 120, e26248.	1.0	6
9	Understanding magnetic relaxation in single-ion magnets with high blocking temperature. <i>Physical Review B</i> , 2020, 101, .	1.1	94
10	Enhancing Magnetic Hysteresis in Single-Molecule Magnets by Ligand Functionalization. <i>CheM</i> , 2020, 6, 1777-1793.	5.8	103
11	A Study of Magnetic Relaxation in Dysprosium(III) Single-Molecule Magnets. <i>Chemistry - A European Journal</i> , 2020, 26, 5893-5902.	1.7	108
12	A double-dysprosocenium single-molecule magnet bound together with neutral ligands. <i>Chemical Communications</i> , 2020, 56, 5677-5680.	2.2	26
13	Reversible uptake of sulfur-containing gases by single crystals of a Cr ₈ metallocrown. <i>Dalton Transactions</i> , 2019, 48, 13184-13189.	1.6	3
14	Slow magnetic relaxation in a {EuCu ₅ } metallocrown. <i>Dalton Transactions</i> , 2019, 48, 1686-1692.	1.6	24
15	Studies of hysteresis and quantum tunnelling of the magnetisation in dysprosium(III) single molecule magnets. <i>Dalton Transactions</i> , 2019, 48, 8541-8545.	1.6	71
16	Light Lanthanide Metallocenium Cations Exhibiting Weak Equatorial Anion Interactions. <i>Chemistry - A European Journal</i> , 2019, 25, 7749-7758.	1.7	29
17	Uncertainty estimates for magnetic relaxation times and magnetic relaxation parameters. <i>Physical Chemistry Chemical Physics</i> , 2019, 21, 23567-23575.	1.3	200
18	Bis-Monophospholyl Dysprosium Cation Showing Magnetic Hysteresis at 48 K. <i>Journal of the American Chemical Society</i> , 2019, 141, 19935-19940.	6.6	123

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19	Post-B3LYP Functionals Do Not Improve the Description of Magnetic Coupling in Cu(II) Dinuclear Complexes. <i>Journal of Physical Chemistry A</i> , 2018, 122, 3423-3432.	1.1	12
20	The performance of density functional theory for the description of ground and excited state properties of inorganic and organometallic uranium compounds. <i>Journal of Organometallic Chemistry</i> , 2018, 857, 58-74.	0.8	30
21	Terbocenium: completing a heavy lanthanide metallocenium cation family with an alternative anion abstraction strategy. <i>Chemical Communications</i> , 2018, 54, 9182-9185.	2.2	30
22	Field- and temperature-dependent quantum tunnelling of the magnetisation in a large barrier single-molecule magnet. <i>Nature Communications</i> , 2018, 9, 3134.	5.8	170
23	Redox-Induced Gating of the Exchange Interactions in a Single Organic Diradical. <i>ACS Nano</i> , 2017, 11, 5879-5883.	7.3	50
24	Molecular magnetic hysteresis at 60 kelvin in dysprosocenium. <i>Nature</i> , 2017, 548, 439-442.	13.7	1,450
25	Calix[n]arene-based polyradicals: enhancing ferromagnetism by avoiding edge effects. <i>Physical Chemistry Chemical Physics</i> , 2017, 19, 24264-24270.	1.3	6
26	Synthesis and Electronic Structures of Heavy Lanthanide Metallocenium Cations. <i>Journal of the American Chemical Society</i> , 2017, 139, 18714-18724.	6.6	111
27	Design of multi-functional 2D open-shell organic networks with mechanically controllable properties. <i>Chemical Science</i> , 2017, 8, 1027-1039.	3.7	16
28	Helical Folding-Induced Stabilization of Ferromagnetic Polyradicals Based on Triarylmethyl Radical Derivatives. <i>Journal of the American Chemical Society</i> , 2016, 138, 5271-5275.	6.6	11
29	Magnetic Coupling Constants in Three Electrons Three Centers Problems from Effective Hamiltonian Theory and Validation of Broken Symmetry-Based Approaches. <i>Journal of Chemical Theory and Computation</i> , 2016, 12, 3228-3235.	2.3	14
30	Exchange Coupling Inversion in a High-Spin Organic Triradical Molecule. <i>Nano Letters</i> , 2016, 16, 2066-2071.	4.5	60
31	Discovery of the K_4 Structure Formed by a Triangular $\dot{\text{C}}$ Radical Anion. <i>Journal of the American Chemical Society</i> , 2015, 137, 7612-7615.	6.6	37
32	Spin Adapted versus Broken Symmetry Approaches in the Description of Magnetic Coupling in Heterodinuclear Complexes. <i>Journal of Chemical Theory and Computation</i> , 2015, 11, 1006-1019.	2.3	14
33	Triplet-singlet gap in structurally flexible organic diradicals. <i>Theoretical Chemistry Accounts</i> , 2015, 134, 1.	0.5	7
34	Handling Magnetic Coupling in Trinuclear Cu(II) Complexes. <i>Journal of Chemical Theory and Computation</i> , 2015, 11, 3650-3660.	2.3	13
35	Characterization of a Robust Co^{II} Fluorescent Complex Deposited Intact On HOPG. <i>Chemistry - A European Journal</i> , 2014, 20, 10439-10445.	1.7	9
36	Nanoparticles of Ni(II) and Co(II) metallo-organic molecular materials. <i>Journal of Nanoparticle Research</i> , 2014, 16, 1.	0.8	2

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37	The Triplet–Singlet Gap in the <i>m</i> -Xylylene Radical: A Not So Simple One. <i>Journal of Chemical Theory and Computation</i> , 2014, 10, 335-345.	2.3	56
38	Theoretical and computational investigation of meta-phenylene as ferromagnetic coupler in nitronyl nitroxide diradicals. <i>Theoretical Chemistry Accounts</i> , 2014, 133, 1.	0.5	12
39	Identifying atomic sites in N-doped pristine and defective graphene from ab initio core level binding energies. <i>Carbon</i> , 2014, 76, 155-164.	5.4	14
40	New nanostructured materials: Nanostructuring of a fluorescent magnet based on acridine yellow. <i>Polyhedron</i> , 2013, 66, 136-141.	1.0	2
41	Microwave assisted synthesis in coordination chemistry. <i>Polyhedron</i> , 2013, 52, 781-787.	1.0	17
42	Extraction of “hidden” relaxation times from AC susceptibility data. <i>Chemistry Squared</i> , 0, 4, 3.	0.0	3
43	Unparalleled selectivity and electronic structure of heterometallic [LnLn TM Ln] molecules as 3-qubit quantum gates. <i>Chemical Science</i> , 0, , .	3.7	6