

# Angela Rosa

## List of Publications by Year in descending order

Source: <https://exaly.com/author-pdf/9379945/publications.pdf>

Version: 2024-02-01

32  
papers

1,547  
citations

236925

25  
h-index

414414

32  
g-index

32  
all docs

32  
docs citations

32  
times ranked

1783  
citing authors

| #  | ARTICLE  | IF   | CITATIONS |
|----|--|------|-----------|
| 1  | Bis(CBT)palladium(II) Derivatives (CBT = <i>m</i> -carborane-1-thiolate): Synthesis, Molecular Structure, and Physicochemical Properties of <i>cis</i> -[(bipy)Pd(CBT) <sub>2</sub> ] and <i>trans</i> -[(py) <sub>2</sub> Pd(CBT) <sub>2</sub> ]. <i>Inorganic Chemistry</i> , 2021, 60, 10478-10491.         | 4.0  | 3         |
| 2  | Origin of the Enhanced Binding Capability toward Axial Nitrogen Bases of Ni(II) Porphyrins Bearing Electron-Withdrawing Substituents: An Electronic Structure and Bond Energy Analysis. <i>Inorganic Chemistry</i> , 2020, 59, 11528-11541.  | 4.0  | 4         |
| 3  | Nickel-macrocycle interaction in nickel(II) porphyrins and porphyrazines bearing alkylthio $\beta$ -substituents: A combined DFT and XPS study. <i>Journal of Porphyrins and Phthalocyanines</i> , 2017, 21, 371-380.  | 0.8  | 12        |
| 4  | Charge Effects on the Reactivity of Oxoiron(IV) Porphyrin Species: A DFT Analysis of Methane Hydroxylation by Polycationic Compound I and Compound II Mimics. <i>ACS Catalysis</i> , 2016, 6, 568-579.   | 11.2 | 27        |
| 5  | Phenyl Derivative of Iron 5,10,15-Tritolylcorrole. <i>Inorganic Chemistry</i> , 2014, 53, 4215-4227.   | 4.0  | 26        |
| 6  | Carboranyl-porphyrazines and derivatives for boron neutron capture therapy: From synthesis to in vitro tests. <i>Coordination Chemistry Reviews</i> , 2013, 257, 2213-2231.  | 18.8 | 37        |
| 7  | Reactivity of Compound II: Electronic Structure Analysis of Methane Hydroxylation by Oxoiron(IV) Porphyrin Complexes. <i>Inorganic Chemistry</i> , 2012, 51, 9833-9845.  | 4.0  | 29        |
| 8  | Near-Infrared-Emitting Phthalocyanines. A Combined Experimental and Density Functional Theory Study of the Structural, Optical, and Photophysical Properties of Pd(II) and Pt(II) $\beta$ -Butoxyphthalocyanines. <i>Inorganic Chemistry</i> , 2011, 50, 1135-1149.  | 4.0  | 46        |
| 9  | The role of the metal ion in the photophysical behavior of <i>Co(II)</i> , <i>Ni(II)</i> , and <i>Cu(II)</i> octabutoxynaphthalocyanines: insights from ultra-fast time-resolved spectroscopy and DFT/TDDFT calculations. <i>Journal of Porphyrins and Phthalocyanines</i> , 2010, 14, 689-700.                | 0.8  | 5         |
| 10 | Is [FeO] <sub>2</sub> the Active Center Also in Iron Containing Zeolites? A Density Functional Theory Study of Methane Hydroxylation Catalysis by Fe-ZSM-5 Zeolite. <i>Inorganic Chemistry</i> , 2010, 49, 3866-3880.  | 4.0  | 52        |
| 11 | On the photophysics of metallophthalocyanine-based photothermal sensitizers: Synergism between theory and experiment. <i>Journal of Inorganic Biochemistry</i> , 2008, 102, 406-413.   | 3.5  | 3         |
| 12 | Tetra-2,3-pyrazinoporphyrazines with Externally Appended Pyridine Rings. 6. Chemical and Redox Properties and Highly Effective Photosensitizing Activity for Singlet Oxygen Production of Penta- and Monopalladated Complexes in Dimethylformamide Solution. <i>Inorganic Chemistry</i> , 2008, 47, 8757-8766. | 4.0  | 34        |
| 13 | Photophysical Behavior of Open-Shell First-Row Transition-Metal Octabutoxynaphthalocyanines: CoNc(OBu) <sub>8</sub> and CuNc(OBu) <sub>8</sub> as Case Studies. <i>Inorganic Chemistry</i> , 2008, 47, 4275-4289.  | 4.0  | 15        |
| 14 | Synthesis and Liposome Insertion of a New Poly(carboranylalkylthio)porphyrazine to Improve Potentiality in Multiple-Approach Cancer Therapy. <i>Journal of the American Chemical Society</i> , 2007, 129, 2728-2729.   | 13.7 | 36        |
| 15 | Evidence for Tetraphenylporphyrin Monoacids. <i>Inorganic Chemistry</i> , 2007, 46, 5979-5988.   | 4.0  | 52        |
| 16 | Effects of Benzoannulation and $\beta$ -Octabutoxy Substitution on the Photophysical Behavior of Nickel Phthalocyanines: A Combined Experimental and DFT/TDDFT Study. <i>Inorganic Chemistry</i> , 2007, 46, 2080-2093.  | 4.0  | 43        |
| 17 | Synergism of Porphyrin-Core Saddling and Twisting of meso-Aryl Substituents. <i>Journal of Physical Chemistry A</i> , 2006, 110, 5180-5190.  | 2.5  | 79        |
| 18 | Structural, Optical, and Photophysical Properties of Nickel(II) Alkylthioporphyrins: Insights from Experimental and DFT/TDDFT Studies. <i>Inorganic Chemistry</i> , 2005, 44, 6609-6622.   | 4.0  | 48        |

| #  | ARTICLE   | IF   | CITATIONS |
|----|---|------|-----------|
| 19 | Photophysics of Octabutoxy Phthalocyaninato-Ni(II) in Toluene: Ultrafast Experiments and DFT/TDDFT Studies. <i>Journal of Physical Chemistry A</i> , 2005, 109, 2078-2089.  | 2.5  | 55        |
| 20 | Excitation Energies of Metal Complexes with Time-dependent Density Functional Theory. <i>Structure and Bonding</i> , 2004, , 49-116.  | 1.0  | 74        |
| 21 | Photodissociation of the Phosphine-Substituted Transition Metal Carbonyl Complexes Cr(CO) <sub>5</sub> L and Fe(CO) <sub>4</sub> L: A Theoretical Study. <i>Journal of the American Chemical Society</i> , 2003, 125, 3558-3567.  | 13.7 | 39        |
| 22 | Effects of Porphyrin Core Saddling, meso-Phenyl Twisting, and Counterions on the Optical Properties of meso-Tetraphenylporphyrin Diacids: The [H4TPP](X) <sub>2</sub> (X = F, Cl, Br, I) Series as a Case Study. <i>Journal of Physical Chemistry A</i> , 2003, 107, 11468-11482. | 2.5  | 131       |
| 23 | Electronic Structure, Chemical Bond, and Optical Spectra of Metal Bis(porphyrin) Complexes: A DFT/TDDFT Study of the Bis(porphyrin)M(IV) (M = Zr, Ce, Th) Series. <i>Journal of the American Chemical Society</i> , 2002, 124, 12319-12334.                                       | 13.7 | 54        |
| 24 | Electronic Spectra of M(CO) <sub>6</sub> (M = Cr, Mo, W) Revisited by a Relativistic TDDFT Approach. <i>Journal of the American Chemical Society</i> , 1999, 121, 10356-10365.  | 13.7 | 330       |
| 25 | Synthesis, Structure, and Physicochemical Properties of ((Ethylsulfanyl)porphyrinato)cobalt(II). Metal-Ligand Bonds in Co(OESPz) and in Related Cobalt(II) Tetrapyrroles: Insights from a Density Functional Study. <i>Inorganic Chemistry</i> , 1999, 38, 1422-1431.             | 4.0  | 37        |
| 26 | Metal-to-Ligand Charge Transfer Photochemistry: Homolysis of the Mn-Cl Bond in the mer-Mn(Cl)(CO) <sub>3</sub> (1,2-diimine) Complex and Its Absence in the fac-Isomer. <i>Inorganic Chemistry</i> , 1998, 37, 6244-6254.   | 4.0  | 20        |
| 27 | Structural Properties of M(dmit) <sub>2</sub> -Based (M = Ni, Pd, Pt; dmit <sub>2</sub> = 2-Thioxo-1,3-dithiole-4,5-dithiolato) Molecular Metals. Insights from Density Functional Calculations. <i>Inorganic Chemistry</i> , 1998, 37, 1368-1379.                                | 4.0  | 35        |
| 28 | Density Functional Study of the Primary Photoprocesses of Manganese Pentacarbonyl Chloride (MnCl(CO) <sub>5</sub> ). <i>Inorganic Chemistry</i> , 1997, 36, 1541-1551.  | 4.0  | 31        |
| 29 | Density Functional Study of the Photodissociation of Mn <sub>2</sub> (CO) <sub>10</sub> . <i>Inorganic Chemistry</i> , 1996, 35, 2886-2897.   | 4.0  | 64        |
| 30 | Metal-to-Ligand Charge Transfer (MLCT) Photochemistry of fac-Mn(Cl)(CO) <sub>3</sub> (H-DAB): A Density Functional Study. <i>The Journal of Physical Chemistry</i> , 1996, 100, 15346-15357.  | 2.9  | 39        |
| 31 | Density Functional Study of Ground and Excited States of Mn <sub>2</sub> (CO) <sub>10</sub> . <i>Inorganic Chemistry</i> , 1995, 34, 3425-3432.   | 4.0  | 54        |
| 32 | Porphyrin vs phthalocyanine metallomacrocycles based one-dimensional "molecular metals". Insights from density functional calculations. <i>Inorganic Chemistry</i> , 1993, 32, 5637-5639.   | 4.0  | 33        |