Angel Terron

List of Publications by Year in descending order

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| 50 | 959 | 19 | 29 |
|----------|----------------|--------------|---------------------|
| papers | citations | h-index | g-index |
| 50 | 50 | 50 | 1040 citing authors |
| all docs | docs citations | times ranked | |

| # | Article | IF | CITATIONS |
|----|---|------------------|-------------------|
| 1 | Biological recognition patterns implicated by the formation and stability of ternary metal ion complexes of low-molecular-weight formed with amino acid/peptides and nucleobases/nucleosides. Coordination Chemistry Reviews, 2007, 251, 1973-1986. | 18.8 | 83 |
| 2 | Synthesis and mass spectroscopy kinetics of a novel ternary copper(II) complex with cytotoxic activity against cancer cells. Journal of Inorganic Biochemistry, 2007, 101, 649-659. | 3.5 | 69 |
| 3 | Experimental and theoretical study of uracil derivatives: the crucial role of weak fluorine–fluorine noncovalent interactions. CrystEngComm, 2010, 12, 3758. | 2.6 | 60 |
| 4 | Synthesis and characterization of nickel(II) complexes of purine and pyrimidine bases. Crystal and molecular structure of trans-bis(cytosine-O2)bis(ethylenediamine)nickel(II) bis(tetraphenylborate). An unusual metal binding mode of cytosine. Inorganic Chemistry, 1990, 29, 5168-5173. | 4.0 | 52 |
| 5 | Structural characterization, recognition patterns and theoretical calculations of long-chain N-alkyl substituted purine and pyrimidine bases as ligands: On the importance of anion–l€ interactions. Coordination Chemistry Reviews, 2013, 257, 2705-2715. | 18.8 | 42 |
| 6 | Xâ€ray Crystal Structure of a Metalled Doubleâ€Helix Generated by Infinite and Consecutive C*â€Ag ^I â€C* (C*:N ¹ â€Hexylcytosine) Base Pairs through Argentophilic and Hydrogen Bond Interactions. Chemistry - A European Journal, 2017, 23, 2103-2108. | 3. 3 | 41 |
| 7 | Lone pairâ€"Ï€ vs Ï€â€"Ï€ interactions in 5-fluoro-1-hexyluracil and 1-hexyluracil: a combined crystallographic and computational study. CrystEngComm, 2010, 12, 362-365. | 2.6 | 39 |
| 8 | Interactions of d10 metal ions with hippuric acid and cytosine. X-ray structure of the first cadmium (II)–amino acid derivative–nucleobase ternary compound. Journal of Inorganic Biochemistry, 2001, 85, 173-178. | 3.5 | 37 |
| 9 | Synthesis and structural characteristics of metal–acyclovir (ACV) complexes: [Ni(or) Tj ETQq1 1 0.784314 rgBT acyclovir by Ni–ACV. Journal of the Chemical Society Dalton Transactions, 1999, , 167-174. | /Overlock 1.1 | 10 Tf 50 42 32 |
| 10 | X-Ray Structural Studies of Metal-Nucleoside and Metal-Nucleoside Monophosphate Complexes: New Perspectives. Comments on Inorganic Chemistry, 1993, 14, 63-88. | 5.2 | 28 |
| 11 | Ternary complexes metal [Co(II), Ni(II), Cu(II) and Zn(II)] $\hat{a} \in \text{``ortho-iodohippurate (I-hip)} \hat{a} \in ``acyclovir. X-ray characterization of isostructural [(Co, Ni or Zn)(I-hip)2(ACV)(H2O)3] with stacking as a recognition factor. Journal of Inorganic Biochemistry, 2004, 98, 1703-1711.$ | 3.5 | 28 |
| 12 | Synthesis, X-ray characterization and regium bonding interactions of a trichlorido(1-hexylcytosine)gold(<scp>iii</scp>) complex. Chemical Communications, 2020, 56, 3524-3527. | 4.1 | 28 |
| 13 | Ruthenium(III) and iridium(III) complexes with nicotine. Polyhedron, 2010, 29, 34-41. | 2.2 | 27 |
| 14 | On the importance of antiparallel C Oâ√C–F interactions in N1-(3-hydroxypropyl)-5-fluorouracilate–Hg(II) complex: A combined X-ray and DFT study. Inorganica Chimica Acta, 2016, 452, 244-250. | 2.4 | 27 |
| 15 | Some new derivatives of Co(III) with uracil, uridine and pyrimidine nucleotides. Inorganica Chimica Acta, 1987, 135, 197-202. | 2.4 | 25 |
| 16 | Synthesis and characterization of a novel copper(II)-cytosine complex: tetrakis(cytosine)copper(II) chloride bis(dimethylacetamide) solvate. Polyhedron, 1994, 13, 2513-2518. | 2.2 | 22 |
| 17 | Molecular architecture by means of interactions between Ag(I) and glycine derivatives. Polyhedron, 2006, 25, 71-80. | 2.2 | 22 |
| 18 | Uracilato and 5-halouracilato complexes of Cu(II), Zn(II) and Ni(II). X-ray structures of [Cu(uracilato-N1)2(NH3)2]·2(H2O), [Cu(5-chlorouracilato-N1)2(NH3)2](H2O)2, [Ni(5-chlorouracilato-N1)2(en)2]·2H2O and [Zn(5-chlorouracilato-N1)(NH3)3]·(5-chlorouracilato-N1)·(H2O). Journal of Inorganic Biochemistry, 2004, 98, 632-638. | 3.5 | 21 |

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|----|---|-------|-----------|
| 19 | X-ray crystal structure of a ternary copper(II) peptide creatinine complex, (Aquo)(Creatinine)(Glycylglycinato) copper(II) sesquihydrate. Polyhedron, 1995, 14, 2537-2545. | 2.2 | 20 |
| 20 | The first X-ray structure of a silver–nucleotide complex: interaction of ion Ag(<scp>i</scp>) with cytidine-5′-monophosphate. CrystEngComm, 2017, 19, 5830-5834. | 2.6 | 18 |
| 21 | Experimental and theoretical study of thymine and cytosine derivatives: the crucial role of weak noncovalent interactions. CrystEngComm, 2012, 14, 5777. | 2.6 | 17 |
| 22 | Complexation in solution of magnesium(II) and cobalt(II) with purine $5\hat{a}\in^2$ -monophosphates and pyrimide $5\hat{a}\in^2$ -monophosphates: a potentiometric and calorimetric study. Polyhedron, 1998, 17, 3825-3833. | 2.2 | 16 |
| 23 | RNAs' uracil quartet model with a non-essential metal ion. Chemical Communications, 2011, 47, 4646. | 4.1 | 16 |
| 24 | Some new derivatives of Ni(II) with uracil, uridine and nucleotides. Inorganica Chimica Acta, 1986, 125, 159-166. | 2.4 | 15 |
| 25 | Ruthenium(III) complexes with modified nucleobases: N6-Substituted adenines. Polyhedron, 2008, 27, 2851-2858. | 2.2 | 13 |
| 26 | Crystal structures of N6-modified-amino acid related nucleobase analogs (II): hybrid adenine-β-alanine and adenine-GABA molecules. New Journal of Chemistry, 2019, 43, 9680-9688. | 2.8 | 13 |
| 27 | Complexation of Nickel(II) with Guanosine 5â€~-Monophosphate and Inosine 5â€~-Monophosphate: A Potentiometric and Calorimetric Study. Inorganic Chemistry, 1996, 35, 3786-3791. | 4.0 | 12 |
| 28 | Experimental and theoretical studies on the coordination chemistry of the N1-hexyl substituted pyrimidines (uracil, 5-fluorouracil and cytosine). Dalton Transactions, 2013, 42, 7631. | 3.3 | 12 |
| 29 | Synthesis, Xâ€ray characterization and density functional theory studies of N ⁶ â€benzylâ€N ⁶ â€methyladenineâ€"M(II) complexes (MÂ=ÂZn, Cd): The prominent ro Ï€â€"I€, Câ€"H···π and anionâ€"I€ interactions. Applied Organometallic Chemistry, 2019, 33, e4906. | ഠിഭത് | 12 |
| 30 | Synthesis and characterization of a new Ni(II) pyrimidine complex. Crystal and molecular structure of trans-bis(isocytosine-O4) bis(ethylenediamine) Ni(II) bis(tetraphenylborate). Inorganica Chimica Acta, 1997, 262, 85-89. | 2.4 | 10 |
| 31 | Ternary copper(II) complexes with hippurate derivatives and 1,10-phenanthroline: Synthesis and biological activity. Inorganica Chimica Acta, 2009, 362, 4744-4753. | 2.4 | 10 |
| 32 | Crystal structures of <i>N</i> ⁶ -modified-aminoacid/peptide nucleobase analogs: hybrid adenine–glycine and adenine–glycylglycine molecules. New Journal of Chemistry, 2018, 42, 14742-14750. | 2.8 | 9 |
| 33 | Experimental and theoretical study of N1-hexylcytosine and N1-hexylcytosinium nitrate: the crucial role of hydrophobic and anion–l€ interactions. Tetrahedron Letters, 2013, 54, 5355-5360. | 1.4 | 8 |
| 34 | A calorimetric study of the Ni(II)-5'AMP system. A base-stacking stabilization. FEBS Journal, 1991, 202, 401-404. | 0.2 | 7 |
| 35 | Intermolecular C—Hπ interactions in 1,5-diphenyl-3-(2-pyridyl)-2-pyrazoline. Acta Crystallographica Section C: Crystal Structure Communications, 2010, 66, o313-o316. | 0.4 | 7 |
| 36 | Iridium(III) coordination of N(6) modified adenine derivatives with aminoacid chains. Journal of Inorganic Biochemistry, 2020, 205, 111000. | 3.5 | 7 |

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|----|--|--------------|------------|
| 37 | Interactions of nickel(II) with adenosine, uridine and cytidine monophosphates. A calorimetric study. Polyhedron, 1995, 14, 1771-1777. | 2.2 | 6 |
| 38 | A calorimetric study of 3d metal ions–acyclovir interactions. The 2-hydroxyethoxymethyl group of acyclovir mimics the role of ribose in deoxy-guanosine and guanosine promoting the coordination through N(7). Journal of Inorganic Biochemistry, 2001, 86, 677-680. | 3 . 5 | 6 |
| 39 | Crystal structures and DFT calculations of new chlorido-dimethylsulfoxide-MIII (M = Ir, Ru, Rh) complexes with the N-pyrazolyl pyrimidine donor ligand: kinetic vs. thermodynamic isomers. Dalton Transactions, 2014, 43, 6353. | 3.3 | 6 |
| 40 | Synthesis, reactivity, X-ray characterization and docking studies of N7/N9-(2-pyrimidyl)-adenine derivatives. Journal of Inorganic Biochemistry, 2020, 203, 110879. | 3. 5 | 6 |
| 41 | Models for thyroxine: Aromatic iodine-assisted self-assemblies. Polyhedron, 2007, 26, 1417-1426. | 2.2 | 5 |
| 42 | Crystal structures of <i>N</i> ⁶ -modified-amino acid nucleobase analogs(<scp>iii</scp>): adenineâ€"valeric acid, adenineâ€"hexanoic acid and adenineâ€"gabapentine. New Journal of Chemistry, 2020, 44, 12236-12246. | 2.8 | 5 |
| 43 | Di-ν-chlorido-bis{chlorido[(<i>R</i>)/(<i>S</i>)-1,5-diphenyl-3-(2-pyridyl-β <i>N</i>)-2-pyrazoline-β <i>N</i> <sup Acta Crystallographica Section E: Structure Reports Online, 2010, 66, m899-m900.</sup | >2<[sup>] | zinc(II)}. |
| 44 | Synthesis, spectroscopic and magnetic characterization of some iron(III)-nucleotide compounds. Transition Metal Chemistry, 1985, 10, 90-93. | 1.4 | 3 |
| 45 | Some new derivatives of Cr(III) with uracil, uridine and 5′-UMP. Polyhedron, 1986, 5, 1125-1130. | 2.2 | 3 |
| 46 | Interactions in solution of cobalt(II) and nickel(II) with nicotinamide adenine dinucleotide: a potentiometric and calorimetric study. Journal of Biological Inorganic Chemistry, 2002, 7, 313-317. | 2.6 | 3 |
| 47 | New chloride-dimethylsulfoxide-iridium(III) complex with histaminium. Polyhedron, 2015, 102, 735-740. | 2,2 | 2 |
| 48 | Cu(II)–N6-Alkyladenine Complexes: Synthesis, X-ray Characterization and Magnetic Properties. Magnetochemistry, 2018, 4, 24. | 2.4 | 2 |
| 49 | Metal removal from the secondary building unit of bio-MOF-1 by adenine N6-alkylation while retaining the overall 3D porous topology. CrystEngComm, 2020, 22, 4201-4205. | 2.6 | 2 |
| 50 | Modified-amino acid/peptide pyrimidine analogs: synthesis, structural characterization and DFT studies of N-(pyrimidyl)gabapentine and N-(pyrimidyl)baclofen. New Journal of Chemistry, 0, , . | 2.8 | 1 |