

Atekeh Tarahhomi

List of Publications by Year in descending order

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38
papers

275
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1163117

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996975

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39
all docs

39
docs citations

39
times ranked

206
citing authors

| # | ARTICLE | IF | CITATIONS |
|----|--|-----|-----------|
| 1 | Different cyclic motifs in phosphoric triamides containing a C(O)NHP(O)(NH) ₂ skeleton and an <i>R</i> ₂ ² (10) graph set in three new compounds: a database analysis of hydrogen-bond strengths based on motifs. <i>Acta Crystallographica Section C: Crystal Structure Communications</i> , 2011, 67, o265-o272. | 0.4 | 34 |
| 2 | Hirshfeld surface analysis of new phosphoramidates. <i>Acta Crystallographica Section B: Structural Science, Crystal Engineering and Materials</i> , 2013, 69, 260-270. | 1.1 | 29 |
| 3 | Crystal structures, DFT calculations and Hirshfeld surface analyses of three new cobalt(III) Schiff base complexes derived from meso-1,2-diphenyl-1,2-ethylenediamine. <i>Journal of Molecular Structure</i> , 2016, 1122, 123-133. | 3.6 | 27 |
| 4 | Different orientations of C=O versus P=O in P(O)NHC(O) skeleton: the first study on an aliphatic diazaphosphorinane with a gauche orientation. <i>Structural Chemistry</i> , 2011, 22, 201-210. | 2.0 | 23 |
| 5 | Hirshfeld surface analysis of new organotin(IV)-phosphoramidate complexes. <i>Journal of Organometallic Chemistry</i> , 2014, 751, 508-518. | 1.8 | 21 |
| 6 | Two new XP(O)[NHC(CH ₃) ₃] ₂ phosphoramidates, with X= (CH ₃) ₂ N and [(CH ₃) ₃ CNH] ₂ P(O)(O). <i>Acta Crystallographica Section C: Crystal Structure Communications</i> , 2012, 68, o164-o169. | 0.4 | 14 |
| 7 | Crystal structures, DFT calculations, and Hirshfeld surface analyses of two new copper(II) and nickel(II) Schiff base complexes derived from meso-1,2-diphenyl-1,2-ethylenediamine. <i>Journal of Molecular Structure</i> , 2017, 1150, 214-226. | 3.6 | 11 |
| 8 | Syntheses and structures of four new mixed-amide phosphoric triamides. <i>Acta Crystallographica Section C, Structural Chemistry</i> , 2016, 72, 251-259. | 0.5 | 9 |
| 9 | A new six-coordinate organotin(IV) complex of OP[NC ₅ H ₁₀] ₃ : A comparison with an analogous five-coordinate complex by means of X-ray crystallography, Hirshfeld surface analysis and DFT calculations. <i>Journal of Coordination Chemistry</i> , 2018, 71, 1575-1592. | 2.2 | 8 |
| 10 | Synthesis and crystal structures of new phosphoric triamides: study of intermolecular interactions by semi-empirical calculations and Hirshfeld surface analysis. <i>Monatshefte für Chemie</i> , 2018, 149, 1759-1776. | 1.8 | 8 |
| 11 | A detailed theoretical and experimental study on the N H, P O and C O stretching frequencies in two new phosphoric triamides and a statistical comparison with analogous structures. <i>Polyhedron</i> , 2019, 158, 215-224. | 2.2 | 8 |
| 12 | Evaluation of Nâ€“Hâ€“O hydrogen bond interactions in two new phosphoric triamides with a P(O)[NHCH(CH ₃) ₃] ₂ segment by means of topological (AIM) calculations, Hirshfeld surface analysis and 3D energy framework approach. <i>Zeitschrift Fur Kristallographie - Crystalline Materials</i> , 2019, 234, 401-413. | 0.8 | 7 |
| 13 | Evaluation of structural, spectroscopic, bonding and electronic properties of some organotin(IV)-phosphoric triamide complexes by using help of DFT, QTAIM and Hirshfeld surface investigations. <i>Computational and Theoretical Chemistry</i> , 2021, 1195, 113098. | 2.5 | 7 |
| 14 | N,Nâ€“Di-tert-butyl-Nâ€“(2,6-difluorobenzoyl)phosphoric triamide. <i>Acta Crystallographica Section E: Structure Reports Online</i> , 2010, 66, o3159-o3159. | 0.2 | 6 |
| 15 | Structural studies on two new phosphoramidates with the same [OC ₄ H ₈ N]P(O) segment by means of single-crystal X-ray diffraction, Hirshfeld surface analysis and 3D energy framework approach. <i>Journal of Molecular Structure</i> , 2021, 1223, 128942. | 3.6 | 6 |
| 16 | The first coordination compounds of OP[NC ₄ H ₈ O] ₃ phosphoric triamide ligand: structural study and Hirshfeld surface analysis of Sn ^{IV} and Mn ^{II} complexes. <i>Journal of Coordination Chemistry</i> , 2017, 70, 1285-1302. | 2.2 | 5 |
| 17 | Energy frameworks and Hirshfeld surface analysis of supramolecular features in three new phosphoric triamides: tuning the intermolecular interactions via the substituent effect. <i>Journal of Molecular Structure</i> , 2022, 1259, 132742. | 3.6 | 5 |
| 18 | N-(2-Fluorobenzoyl)-Nâ€“,Nâ€“-bis(4-methylphenyl)phosphoric triamide. <i>Acta Crystallographica Section E: Structure Reports Online</i> , 2011, 67, o934-o934. | 0.2 | 4 |

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|----|--|-----|-----------|
| 19 | Three new phosphoric triamides with a [C(O)NH]P(O)[N(C)(C)] ₂ skeleton: a database analysis of C-N-C and P-N-C bond angles. <i>Acta Crystallographica Section C, Structural Chemistry</i> , 2014, 70, 998-1002. | 0.5 | 4 |
| 20 | Hirshfeld surface analysis of the 1,1-(ethane-1,2-diyl)dipyridinium dication in two new salts: perchlorate and peroxodisulfate. <i>Acta Crystallographica Section C, Structural Chemistry</i> , 2014, 70, 230-235. | 0.5 | 4 |
| 21 | Association of Non-covalent Interactions X (X = O, F, Cl, I) and Cl with Hydrogen Bond Interactions N-H-O in Molecular Assembly of New Phosphoramides: A Combined X-Ray Crystallography and Topology (AIM and Hirshfeld) Analysis. <i>ChemistrySelect</i> , 2020, 5, 185-195. | 1.5 | 4 |
| 22 | Coordination versus hydrogen bonds in the structures of different tris(pyridin-2-yl)phosphoric triamide derivatives. <i>RSC Advances</i> , 2021, 11, 8178-8197. | 3.6 | 4 |
| 23 | N,N-Dibenzyl-N-(2,6-difluorobenzoyl)-N,N-dimethylphosphoric triamide. <i>Acta Crystallographica Section E: Structure Reports Online</i> , 2010, 66, o2524-o2524. | 0.2 | 4 |
| 24 | Experimental and theoretical study of novel amino-functionalized P(V) coordination compounds suggested as inhibitor of M ^{Pro} of SARS-CoV-2 by molecular docking study. <i>Applied Organometallic Chemistry</i> , 2022, 36, e6636. | 3.5 | 4 |
| 25 | O-Phenyl (tert-butylamido)(p-tolylamido)phosphinate. <i>Acta Crystallographica Section E: Structure Reports Online</i> , 2011, 67, o3405-o3406. | 0.2 | 3 |
| 26 | Synthesis, crystal structures and Hirshfeld surface analyses of two new Salen type nickel/sodium heteronuclear complexes. <i>Journal of Molecular Structure</i> , 2016, 1110, 119-127. | 3.6 | 3 |
| 27 | Three new amidophosphoric acid esters with a P(O)[OCH ₂ C(CH ₃) ₂ CH ₂ O] segment: X-ray diffraction, DFT, AIM and Hirshfeld surface investigations of bi- and tri-furcated (three and four-center) hydrogen bond interactions. <i>Zeitschrift Fur Kristallographie - Crystalline Materials</i> , 2020, 235, 69-84. | 0.8 | 3 |
| 28 | A second monoclinic polymorph of N-[bis(morpholin-4-yl)phosphinoyl]-4-fluorobenzamide with the P ₂ 1n space group. <i>Acta Crystallographica Section E: Structure Reports Online</i> , 2012, 68, o2444-o2444. | 0.2 | 2 |
| 29 | A novel amido-pyrophosphate MnII chelate complex with the synthetic ligand O{P(O)[NHC(CH ₃) ₃] ₂ }(L): [Mn(L) ₂ {OC(H)N(CH ₃) ₂ }]Cl ₂ ·2H ₂ O. <i>Acta Crystallographica Section C: Crystal Structure Communications</i> , 2013, 69, 225-228. | 0.4 | 2 |
| 30 | Synthesis and structural study of five new phosphoric triamides: interplay between classical and non-classical intermolecular interactions. <i>Zeitschrift Fur Kristallographie - Crystalline Materials</i> , 2021, 236, 301-312. | 0.8 | 2 |
| 31 | N-(2,6-Difluorobenzoyl)-P,P-bis(pyrrrolidin-1-yl)phosphinic amide. <i>Acta Crystallographica Section E: Structure Reports Online</i> , 2011, 67, o2444-o2444. | 0.2 | 1 |
| 32 | N,N,N,N-Tetraethyl-N-(2-fluorobenzoyl)phosphoric triamide. <i>Acta Crystallographica Section E: Structure Reports Online</i> , 2011, 67, o2643-o2643. | 0.2 | 1 |
| 33 | N,N-Dicyclohexyl-N-(2,6-difluorobenzoyl)-N,N-dimethylphosphoric triamide. <i>Acta Crystallographica Section E: Structure Reports Online</i> , 2011, 67, o3027-o3027. | 0.2 | 1 |
| 34 | A new C _{2v} amidophosphoester structure: Hirshfeld surface analysis of symmetry-independent molecules. <i>Zeitschrift Fur Kristallographie - Crystalline Materials</i> , 2017, 232, 453-462. | 0.8 | 1 |
| 35 | Cyclohexyl(methyl)ammonium {bis[cyclohexyl(methyl)amino]phosphoryl}(4-methylphenylsulfonyl)azanide. <i>Acta Crystallographica Section E: Structure Reports Online</i> , 2011, 67, o2795-o2795. | 0.2 | 0 |
| 36 | N,N,N,N-Tetrabenzyl-N-(2,6-difluorobenzoyl)phosphoric triamide. <i>Acta Crystallographica Section E: Structure Reports Online</i> , 2012, 68, o3186-o3186. | 0.2 | 0 |

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| 37 | N-(4-Fluorobenzoyl)-N,N'-diisopropylphosphoric triamide. Acta Crystallographica Section E: Structure Reports Online, 2012, 68, o3406-o3407. | 0.2 | 0 |
| 38 | Crystal structure of N,N',N'',N'''-tetrakis(4-methylphenyl)-diphosphoric tetraamide, C ₂₈ H ₃₂ N ₄ O ₃ P ₂ . Zeitschrift Fur Kristallographie - New Crystal Structures, 2012, 227, . | 0.3 | 0 |