

# Srinivasulu Aitipamula

## List of Publications by Year in descending order

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64  
papers

2,917  
citations

186265

28  
h-index

161849

54  
g-index

67  
all docs

67  
docs citations

67  
times ranked

2593  
citing authors

| #  | ARTICLE   | IF   | CITATIONS |
|----|---|------|-----------|
| 1  | Polymorphism and distinct physicochemical properties of the phloretin–nicotinamide cocrystal. <i>CrystEngComm</i> , 2022, 24, 560-570.  | 2.6  | 12        |
| 2  | Insights into the structure-property relationship of pharmaceutical co-crystals: Charge density and quantum chemical approaches. <i>Journal of Molecular Structure</i> , 2021, 1224, 129270.  | 3.6  | 1         |
| 3  | Cocrystal Formulations: Evaluation of the Impact of Excipients on Dissolution by Molecular Simulation and Experimental Approaches. <i>Crystal Growth and Design</i> , 2021, 21, 1006-1018.  | 3.0  | 10        |
| 4  | Directing Selectivity to Aldehydes, Alcohols, or Esters with Diphobane Ligands in Pd-Catalyzed Alkene Carbonylations. <i>Organometallics</i> , 2021, 40, 1914-1925.   | 2.3  | 7         |
| 5  | <i>gem</i> -Dialkyl Effect in Diphosphine Ligands: Synthesis, Coordination Behavior, and Application in Pd-Catalyzed Hydroformylation. <i>ACS Catalysis</i> , 2020, 10, 663-671.  | 11.2 | 9         |
| 6  | Cocrystal formulations: A case study of topical formulations consisting of ferulic acid cocrystals. <i>European Journal of Pharmaceutics and Biopharmaceutics</i> , 2020, 149, 95-104.  | 4.3  | 23        |
| 7  | Cocrystals of Leflunomide: Design, Structural, and Physicochemical Evaluation. <i>Crystal Growth and Design</i> , 2019, 19, 3923-3933.  | 3.0  | 19        |
| 8  | Agomelatine–hydroquinone (1:1) cocrystal: novel polymorphs and their thermodynamic relationship. <i>Acta Crystallographica Section B: Structural Science, Crystal Engineering and Materials</i> , 2019, 75, 969-977.  | 1.1  | 3         |
| 9  | Cocrystals of zonisamide: physicochemical characterization and sustained release solid forms. <i>CrystEngComm</i> , 2018, 20, 2923-2931.  | 2.6  | 24        |
| 10 | Reply to the “Comment on “Trimorphs of a pharmaceutical cocrystal involving two active pharmaceutical ingredients: potential relevance to combination drugs” by S. Aitipamula, P. S. Chow and R. B. H. Tan, <i>CrystEngComm</i> , 2009, 11, 1823–1823. <i>CrystEngComm</i> , 2018, 20, 373-374. | 2.6  | 0         |
| 11 | Evaluating Suspension Formulations of Theophylline Cocrystals With Artificial Sweeteners. <i>Journal of Pharmaceutical Sciences</i> , 2018, 107, 604-611.   | 3.3  | 21        |
| 12 | Novel solid forms of lonidamine: crystal structures and physicochemical properties. <i>CrystEngComm</i> , 2017, 19, 2925-2935.  | 2.6  | 11        |
| 13 | X-Ray Crystallography and its Role in Understanding the Physicochemical Properties of Pharmaceutical Cocrystals. <i>Journal of the Indian Institute of Science</i> , 2017, 97, 227-243.   | 1.9  | 42        |
| 14 | Design of Cocrystals for Molecules with Limited Hydrogen Bonding Functionalities: Propyphenazone as a Model System. <i>Crystal Growth and Design</i> , 2017, 17, 163-174.   | 3.0  | 35        |
| 15 | 1. Pharmaceutical co-crystals: crystal engineering and applications. , 2017, , 1-31.  |      | 1         |
| 16 | Synergistic enhancement of tableability and physicochemical properties through co-crystallization. <i>Acta Crystallographica Section A: Foundations and Advances</i> , 2017, 73, C59-C59.   | 0.1  | 0         |
| 17 | Using charge density to understand structure–property relationships in pharmaceutical co-crystals. <i>Acta Crystallographica Section A: Foundations and Advances</i> , 2017, 73, C799-C799.   | 0.1  | 0         |
| 18 | Novel solid forms of oxaprozin: cocrystals and an extended release drug–drug salt of salbutamol. <i>RSC Advances</i> , 2016, 6, 34110-34119.  | 3.6  | 28        |

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|----|---|-----|-----------|
| 19 | Novel pharmaceutical cocrystals of triflusal: crystal engineering and physicochemical characterization. <i>CrystEngComm</i> , 2015, 17, 9323-9335.  | 2.6 | 14        |
| 20 | Palladium Complexes with Bulky Diphosphine Ligands as Highly Selective Catalysts for the Synthesis of (Bio-) Adipic Acid from Pentenoic Acid Mixtures.. <i>Organometallics</i> , 2015, 34, 4281-4292.                   | 2.3 | 33        |
| 21 | Polymorphism in Molecular Crystals and Cocrystals. , 2015, , 265-298.   |     | 2         |
| 22 | Charge density studies on polymorphic co-crystals. <i>Acta Crystallographica Section A: Foundations and Advances</i> , 2015, 71, s415-s416.   | 0.1 | 0         |
| 23 | Solvates of the antifungal drug griseofulvin: structural, thermochemical and conformational analysis. <i>Acta Crystallographica Section B: Structural Science, Crystal Engineering and Materials</i> , 2014, 70, 54-62. | 1.1 | 12        |
| 24 | Polymorphism in cocrystals: a review and assessment of its significance. <i>CrystEngComm</i> , 2014, 16, 3451.  | 2.6 | 242       |
| 25 | Cocrystallization with flufenamic acid: comparison of physicochemical properties of two pharmaceutical cocrystals. <i>CrystEngComm</i> , 2014, 16, 5793.  | 2.6 | 60        |
| 26 | Crystal Engineering of Tegafur Cocrystals: Structural Analysis and Physicochemical Properties. <i>Crystal Growth and Design</i> , 2014, 14, 6557-6569.  | 3.0 | 35        |
| 27 | Pharmaceutical Salts of Haloperidol with Some Carboxylic Acids and Artificial Sweeteners: Hydrate Formation, Polymorphism, and Physicochemical Properties. <i>Crystal Growth and Design</i> , 2014, 14, 2542-2556.      | 3.0 | 43        |
| 28 | Charge density studies on 1:1 co-crystals of ethenzamide and saccharin. <i>Acta Crystallographica Section A: Foundations and Advances</i> , 2014, 70, C964-C964.  | 0.1 | 0         |
| 29 | Novel solid forms of the anti-tuberculosis drug, Isoniazid: ternary and polymorphic cocrystals. <i>CrystEngComm</i> , 2013, 15, 5877.   | 2.6 | 97        |
| 30 | Polymorphism and phase transformations of a cocrystal of nicotinamide and pimelic acid. <i>CrystEngComm</i> , 2012, 14, 8193.   | 2.6 | 30        |
| 31 | Pharmaceutical cocrystals of ethenzamide: structural, solubility and dissolution studies. <i>CrystEngComm</i> , 2012, 14, 8515.   | 2.6 | 71        |
| 32 | The solvates of sulfamerazine: structural, thermochemical, and desolvation studies. <i>CrystEngComm</i> , 2012, 14, 691-699.  | 2.6 | 44        |
| 33 | Cocrystal Hydrate of an Antifungal Drug, Griseofulvin, with Promising Physicochemical Properties. <i>Crystal Growth and Design</i> , 2012, 12, 5858-5863.   | 3.0 | 61        |
| 34 | Polymorphs, Salts, and Cocrystals: Whatâ€™s in a Name?. <i>Crystal Growth and Design</i> , 2012, 12, 2147-2152.   | 3.0 | 767       |
| 35 | Co-crystals of caffeine and piracetam with 4-hydroxybenzoic acid: Unravelling the hidden hydrates of 1:1 co-crystals. <i>CrystEngComm</i> , 2012, 14, 2381.   | 2.6 | 36        |
| 36 | Correction for Polymorphs, Salts and Cocrystals: Whatâ€™s in a Name?. <i>Crystal Growth and Design</i> , 2012, 12, 4290-4291.   | 3.0 | 17        |

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|----|--|-----|-----------|
| 37 | Conformational Polymorphs of a Muscle Relaxant, Metaxalone. <i>Crystal Growth and Design</i> , 2011, 11, 4101-4109.  | 3.0 | 24        |
| 38 | Solvates and polymorphic phase transformations of 2-chloro-4-nitrobenzoic acid. <i>CrystEngComm</i> , 2011, 13, 1037-1045.   | 2.6 | 38        |
| 39 | Solvates and a monohydrate of N4-acetylsulfamerazine: Structural, thermochemical, and computational analysis. <i>Journal of Molecular Structure</i> , 2011, 1005, 134-140.                               | 3.6 | 7         |
| 40 | Structural, Spectroscopic and Thermal Analysis of Cocrystals of Carbamazepine and Piracetam with Hydroquinone. <i>Journal of Chemical Crystallography</i> , 2011, 41, 1604-1611.                         | 1.1 | 11        |
| 41 | Pyrimidin-2-amineâ€“1-phenylcyclopentane-1-carboxylic acid (1/1). <i>Acta Crystallographica Section E: Structure Reports Online</i> , 2011, 67, o552-o553.   | 0.2 | 2         |
| 42 | N,N-Dimethylpyridin-4-aminium 1-phenylcyclopentane-1-carboxylate monohydrate. <i>Acta Crystallographica Section E: Structure Reports Online</i> , 2011, 67, o1227-o1227.                                 | 0.2 | 0         |
| 43 | Conformational Polymorphism of Tolbutamide: A Structural, Spectroscopic, and Thermodynamic Characterization of Burgerâ€™s Forms Iâ€“IV. <i>Journal of Pharmaceutical Sciences</i> , 2010, 99, 2975-2990. | 3.3 | 62        |
| 44 | 2-Aminopyridinium 1-phenylcyclopropane-1-carboxylate. <i>Acta Crystallographica Section E: Structure Reports Online</i> , 2010, 66, o3339-o3340.   | 0.2 | 2         |
| 45 | Ethenzamideâ€“gentisic acidâ€“acetic acid (2/1/1). <i>Acta Crystallographica Section E: Structure Reports Online</i> , 2010, 66, o1045-o1046.  | 0.2 | 10        |
| 46 | Conformational and enantiotropic polymorphism of a 1â€“1 cocrystal involving ethenzamide and ethylmalonic acid. <i>CrystEngComm</i> , 2010, 12, 3691.  | 2.6 | 58        |
| 47 | Polymorphs and Solvates of a Cocrystal Involving an Analgesic Drug, Ethenzamide, and 3,5-Dinitrobenzoic Acid. <i>Crystal Growth and Design</i> , 2010, 10, 2229-2238.                                    | 3.0 | 109       |
| 48 | The amido-bridged zirconocene's reactivity and catalytic behavior for ethylenepolymerization. <i>Dalton Transactions</i> , 2010, 39, 807-814.  | 3.3 | 7         |
| 49 | Theophyllineâ€“gentisic acid (1/1). <i>Acta Crystallographica Section E: Structure Reports Online</i> , 2009, 65, o2126-o2127.   | 0.2 | 9         |
| 50 | Trimorphs of a pharmaceutical cocrystal involving two active pharmaceutical ingredients: potential relevance to combination drugs. <i>CrystEngComm</i> , 2009, 11, 1823-1827.                            | 2.6 | 134       |
| 51 | Dimorphs of a 1â€“1 cocrystal of ethenzamide and saccharin: solid-state grinding methods result in metastable polymorph. <i>CrystEngComm</i> , 2009, 11, 889.  | 2.6 | 73        |
| 52 | Microsphaerins Aâ€“D, four novel benzophenone dimers with activity against MRSA from the fungus <i>Microsphaeropsis</i> sp.. <i>Tetrahedron</i> , 2008, 64, 10181-10187.                                 | 1.9 | 24        |
| 53 | Polymorphs and Polymorphic Cocrystals of Temozolomide. <i>Chemistry - an Asian Journal</i> , 2008, 3, 1122-1133.   | 3.3 | 78        |
| 54 | Guest-Induced Supramolecular Isomerism in Inclusion Complexes of T-Shaped Host 4,4-Bis(4â€“hydroxyphenyl)cyclohexanone. <i>Chemistry - A European Journal</i> , 2005, 11, 6727-6742.                     | 3.3 | 90        |

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|----|--|-----|-----------|
| 55 | Engineering the weak Nâ€“Hâ€“ hydrogen bond in 4-tritylbenzamide host and controlling the interaction through guest selection. <i>CrystEngComm</i> , 2005, 7, 44-52.   | 2.6 | 32        |
| 56 | Concomitant polymorphs of 2,2â€²,6,6â€²-tetramethyl-4,4â€²-terphenyldiol: the Î² <sup>2</sup> -quinol network reproduced in a metastable polymorph. <i>Chemical Communications</i> , 2005, , 3159.   | 4.1 | 29        |
| 57 | Thermochemical Analysis of Venlafaxine Hydrochloride Polymorphs1âˆ’5â€. <i>Crystal Growth and Design</i> , 2005, 5, 2268-2276.   | 3.0 | 43        |
| 58 | Hexagonal Host Framework of <i>sym</i>-Aryloxytriazines Stabilised by Weak Intermolecular Interactions. <i>Molecular Crystals and Liquid Crystals</i> , 2005, 440, 295-316.  | 0.9 | 17        |
| 59 | Ladder and Hexagonal Hydrogen-bond Networks from a Self-complementary H-shaped Tecton. <i>Supramolecular Chemistry</i> , 2005, 17, 17-25.  | 1.2 | 14        |
| 60 | Halogen Trimer-Mediated Hexagonal Host Framework of 2,4,6-Tris(4-halophenoxy)-1,3,5-triazine. Supramolecular Isomerism from Hexagonal Channel (X = Cl, Br) to Cage Structure (X = I). <i>Crystal Growth and Design</i> , 2005, 5, 887-899. | 3.0 | 98        |
| 61 | Hostâ€“guest and network structures of some tetraphenylmethane derivatives. <i>CrystEngComm</i> , 2004, 6, 120-125.  | 2.6 | 23        |
| 62 | Hydrogen-bond networks in tris(4-hydroxyphenyl)methane and its 1:1 molecular complex with 4,4â€²-bipyridine. <i>Acta Crystallographica Section C: Crystal Structure Communications</i> , 2003, 59, o481-o484.                              | 0.4 | 4         |
| 63 | Multiple molecules in the crystallographic asymmetric unit. Self hostâ€“guest and doubly interpenetrated hydrogen bond networks in a pair of keto-bisphenols. <i>CrystEngComm</i> , 2003, 5, 447.  | 2.6 | 44        |
| 64 | Topological Equivalences between Organic and Coordination Polymer Crystal Structures:â€“ An Organic Ladder Formed with Three-Connected Molecular and Supramolecular Synthons. <i>Organic Letters</i> , 2002, 4, 921-924.                   | 4.6 | 61        |