## Atish T Paul

## List of Publications by Year in descending order

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Version: 2024-02-01

840585 526166 772 29 11 27 h-index citations g-index papers 29 29 29 620 docs citations citing authors all docs times ranked

#	Article	IF	CITATIONS
1	Synthesis, molecular modelling, <i>in vitro</i> and <i>in vivo</i> evaluation of conophylline inspired novel benzyloxy substituted indole glyoxylamides as potent pancreatic lipase inhibitors. Journal of Biomolecular Structure and Dynamics, 2022, 40, 9530-9542.	2.0	6
2	Mechanistically acting anti-obesity compositions/formulations of natural origin: a patent review (2010–2021). Expert Opinion on Therapeutic Patents, 2022, 32, 29-46.	2.4	3
3	Design and Synthesis of Echitamine Inspired Hybrid Analogues Containing Thiazolidinediones as Potential Pancreatic Lipase Inhibitors. Letters in Drug Design and Discovery, 2022, 19, .	0.4	O
4	Metagenomic analysis for taxonomic and functional potential of Polyaromatic hydrocarbons (PAHs) and Polychlorinated biphenyl (PCB) degrading bacterial communities in steel industrial soil. PLoS ONE, 2022, 17, e0266808.	1.1	18
5	Design, synthesis, biological evaluation and molecular modelling studies of oxoacetamide warhead containing indole-quinazolinone based novel hybrid analogues as potential pancreatic lipase inhibitors. New Journal of Chemistry, 2022, 46, 11648-11661.	1.4	12
6	U.S. FDA Approved Drugs from 2015–June 2020: A Perspective. Journal of Medicinal Chemistry, 2021, 64, 2339-2381.	2.9	314
7	Design, synthesis and biological evaluation of <i>N</i> à€substituted indoleâ€thiazolidinedione analogues as potential pancreatic lipase inhibitors. Chemical Biology and Drug Design, 2021, 98, 49-59.	1.5	12
8	Preparation and Evaluation of Quinapyramine Sulphate-Docusate Sodium Ionic Complex Loaded Lipidic Nanoparticles and Its Scale Up Using Geometric Similarity Principle. Journal of Pharmaceutical Sciences, 2021, 110, 2241-2249.	1.6	6
9	Design, synthesis, <i>in silico </i> molecular modelling studies and biological evaluation of novel indole-thiazolidinedione hybrid analogues as potential pancreatic lipase inhibitors. New Journal of Chemistry, 2021, 45, 1381-1394.	1.4	22
10	Pharmaceutical Application of Bio-actives from Alstonia Genus: Current Findings and Future Directions. Advanced Structured Materials, 2021, , 463-533.	0.3	O
11	Recent advances in the pharmacological diversification of quinazoline/quinazolinone hybrids. RSC Advances, 2020, 10, 41353-41392.	1.7	132
12	Investigation of synergistic potential of green tea polyphenols and orlistat combinations using pancreatic lipase assay-based synergy directed fractionation strategy. South African Journal of Botany, 2020, 135, 50-57.	1.2	10
13	Rapid and costâ€effective LC–MS/MS method for determination of hydroxycitric acid in plasma: Application in the determination of pharmacokinetics in commercial ⟨i⟩Garcinia⟨ i⟩ preparations. Biomedical Chromatography, 2020, 34, e4902.	0.8	5
14	Nanotechnological interventions for treatment of trypanosomiasis in humans and animals. Drug Delivery and Translational Research, 2020, 10, 945-961.	3.0	14
15	Design, synthesis, evaluation, and molecular modeling studies of indolyl oxoacetamides as potential pancreatic lipase inhibitors. Archiv Der Pharmazie, 2020, 353, e2000048.	2.1	7
16	Design, synthesis, biological evaluation and molecular modelling studies of conophylline inspired novel indolyl oxoacetamides as potent pancreatic lipase inhibitors. New Journal of Chemistry, 2020, 44, 12355-12369.	1.4	9
17	Optimisation of an extraction conditions for Rumex nepalensis anthraquinones and its correlation with pancreatic lipase inhibitory activity. Journal of Food Composition and Analysis, 2020, 92, 103575.	1.9	7
18	Impact of quercetin on pharmacokinetics of quetiapine: insights from <i>in-vivo</i> studies in wistar rats. Xenobiotica, 2020, 50, 1483-1489.	0.5	1

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19	Evaluation of biphenyl- and polychlorinated-biphenyl (PCB) degrading <i>Rhodococcus</i> sp. MAPN-1 on growth of <i>Morus alba</i> by pot study. International Journal of Phytoremediation, 2020, 22, 1487-1496.	1.7	10
20	Design, synthesis, biological evaluation, and molecular modeling studies of rhodanine derivatives as pancreatic lipase inhibitors. Archiv Der Pharmazie, 2019, 352, e1900029.	2.1	23
21	Development and validation of a new HPTLC-HRMS method for the quantification of a potent pancreatic lipase inhibitory lead Echitamine from Alstonia scholaris. Natural Product Research, 2019, 35, 1-5.	1.0	9
22	Design, synthesis, biological evaluation and molecular modelling studies of indole glyoxylamides as a new class of potential pancreatic lipase inhibitors. Bioorganic Chemistry, 2019, 85, 373-381.	2.0	22
23	Development and validation of a new HPTLC method for quantification of conophylline in Tabernaemontana divaricata samples obtained from different seasons and extraction techniques: Insights into variation of pancreatic lipase inhibitory activity. Industrial Crops and Products, 2018, 111, 462-470.	2.5	8
24	Unanticipated Cleavage of 2-Nitrophenyl-Substituted <i>N</i> Formyl Pyrazolines under Bechamp Conditions: Unveiling the Synthesis of 2-Aryl Quinolines and Their Mechanistic Exploration via DFT Studies. ACS Omega, 2018, 3, 18783-18790.	1.6	4
25	Bis-indole alkaloids from Tabernaemontana divaricata as potent pancreatic lipase inhibitors: molecular modelling studies and experimental validation. Medicinal Chemistry Research, 2017, 26, 1268-1278.	1.1	20
26	Design, synthesis, biological evaluation and molecular modelling studies of novel diaryl substituted pyrazolyl thiazolidinediones as potent pancreatic lipase inhibitors. Bioorganic and Medicinal Chemistry Letters, 2017, 27, 3749-3754.	1.0	37
27	Synthesis, evaluation and molecular modelling studies of 2-(carbazol-3-yl)-2-oxoacetamide analogues as a new class of potential pancreatic lipase inhibitors. Bioorganic and Medicinal Chemistry, 2017, 25, 609-620.	1.4	42
28	Evaluation of Apoptosis and Autophagy Inducing Potential of Berberis aristata, Azadirachta indica, and Their Synergistic Combinations in Parental and Resistant Human Osteosarcoma Cells. Frontiers in Oncology, 2017, 7, 296.	1.3	15
29	Lupane Analogue from Bark of Pithecellobium dulce and in vitro α-Glucosidase and α-Amylase Enzyme Inhibition Assay of Extract for Potential Antidiabetic Activity. Chemistry of Natural Compounds, 2016, 52, 359-362.	0.2	4