

# Omprakash Tanwar

## List of Publications by Citations

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19  
papers

522  
citations

10  
h-index

19  
g-index

19  
ext. papers

607  
ext. citations

3.1  
avg, IF

3.17  
L-index

#	Paper	IF	Citations
19	Quinoline: A versatile heterocyclic. <i>Saudi Pharmaceutical Journal</i> , <b>2013</b> , 21, 1-12	4.4	248
18	Pyrazolines: a biological review. <i>Mini-Reviews in Medicinal Chemistry</i> , <b>2013</b> , 13, 921-31	3.2	100
17	Pyrazole-pyrazoline as promising novel antimalarial agents: A mechanistic study. <i>European Journal of Medicinal Chemistry</i> , <b>2018</b> , 149, 139-147	6.8	37
16	Recent updates on biological activities of oxadiazoles. <i>Mini-Reviews in Medicinal Chemistry</i> , <b>2013</b> , 13, 1027-46	3.2	19
15	Synthesis and biological evaluation of some new pyrazoline substituted benzenesulfonylurea/thiourea derivatives as anti-hyperglycaemic agents and aldose reductase inhibitors. <i>European Journal of Medicinal Chemistry</i> , <b>2014</b> , 80, 209-17	6.8	16
14	Structure based virtual screening of MDPI database: discovery of structurally diverse and novel DPP-IV inhibitors. <i>Bioorganic and Medicinal Chemistry Letters</i> , <b>2014</b> , 24, 3447-51	2.9	15
13	Structural comparison of Mtb-DHFR and h-DHFR for design, synthesis and evaluation of selective non-pteridine analogues as antitubercular agents. <i>Bioorganic Chemistry</i> , <b>2018</b> , 80, 319-333	5.1	12
12	Identification of ZINC02765569: a potent inhibitor of PTP1B by vHTS. <i>Medicinal Chemistry Research</i> , <b>2013</b> , 22, 28-34	2.2	12
11	Synthesis and antimalarial activity of quinoline-substituted furanone derivatives and their identification as selective falcipain-2 inhibitors. <i>Medicinal Chemistry Research</i> , <b>2015</b> , 24, 879-890	2.2	11
10	Synthesis, 3D-QSAR and docking studies of pyrimidine nitrile-pyrazoline: a novel class of hybrid antimalarial agents. <i>Medicinal Chemistry Research</i> , <b>2015</b> , 24, 1018-1037	2.2	11
9	Novel hydrazine derivatives as selective DPP-IV inhibitors: findings from virtual screening and validation through molecular dynamics simulations. <i>Journal of Molecular Modeling</i> , <b>2014</b> , 20, 2118	2	10
8	3D-QSAR of amino-substituted pyrido[3,2B]pyrazinones as PDE-5 inhibitors. <i>Medicinal Chemistry Research</i> , <b>2012</b> , 21, 202-211	2.2	8
7	Pharmacophore based virtual screening, synthesis and SAR of novel inhibitors of Mycobacterium sulfotransferase. <i>Bioorganic and Medicinal Chemistry Letters</i> , <b>2015</b> , 25, 701-7	2.9	6
6	Expansion of a novel lead targeting M. tuberculosis DHFR as antitubercular agents. <i>Bioorganic and Medicinal Chemistry</i> , <b>2019</b> , 27, 1421-1429	3.4	4
5	2-D QSAR studies of steroidal natural products oleanic acid and their semisynthetic derivatives as potent protein tyrosine phosphatase 1B inhibitors. <i>Medicinal Chemistry Research</i> , <b>2012</b> , 21, 351-361	2.2	4
4	Pharmacophore model generation and 3D-QSAR analysis of N-acyl and N-arylpiprazolines for enzymatic and cellular B-Raf kinase inhibition. <i>Medicinal Chemistry Research</i> , <b>2013</b> , 22, 2174-2187	2.2	4
3	Modeling VEGFR kinase inhibition of aminopyrazolopyridine urea derivatives using topological and physicochemical descriptors: a quantitative structure activity analysis study. <i>Medicinal Chemistry Research</i> , <b>2012</b> , 21, 3958-3964	2.2	3

- 2 3D quantitative structure–activity relationship for quinoline, benzimidazole and benzofuran-based analogs as phosphodiesterases IV (PDE-IV) inhibitors. *Medicinal Chemistry Research*, **2013**, 22, 5153-5166<sup>2,2</sup> 1
- 1 Ethyl Pyruvate as a Potential Defense Intervention against Cytokine Storm in COVID-19?. *ACS Omega*, **2021**, 6, 7754-7760 3-9 1