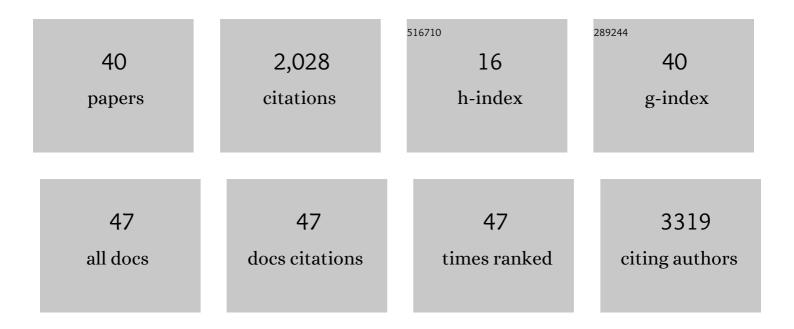
Vivek K Asati

List of Publications by Year in descending order

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#	Article	IF	CITATIONS
1	PI3K/Akt/mTOR and Ras/Raf/MEK/ERK signaling pathways inhibitors as anticancer agents: Structural and pharmacological perspectives. European Journal of Medicinal Chemistry, 2016, 109, 314-341.	5.5	452
2	Anti-cancer chalcones: Structural and molecular target perspectives. European Journal of Medicinal Chemistry, 2015, 98, 69-114.	5.5	382
3	Chalcones and their therapeutic targets for the management of diabetes: Structural and pharmacological perspectives. European Journal of Medicinal Chemistry, 2015, 92, 839-865.	5.5	188
4	Chalcone scaffolds as anti-infective agents: Structural and molecular target perspectives. European Journal of Medicinal Chemistry, 2015, 101, 496-524.	5.5	147
5	1,3,4-Oxadiazoles: An emerging scaffold to target growth factors, enzymes and kinases as anticancer agents. European Journal of Medicinal Chemistry, 2015, 97, 124-141.	5.5	142
6	Thiazolidine-2,4-diones as multi-targeted scaffold in medicinal chemistry: Potential anticancer agents. European Journal of Medicinal Chemistry, 2014, 87, 814-833.	5.5	110
7	Chalcone Derivatives: Anti-inflammatory Potential and Molecular Targets Perspectives. Current Topics in Medicinal Chemistry, 2017, 17, 3146-3169.	2.1	108
8	Perspectives of medicinally privileged chalcone based metal coordination compounds for biomedical applications. European Journal of Medicinal Chemistry, 2019, 174, 142-158.	5.5	71
9	PIM kinase inhibitors: Structural and pharmacological perspectives. European Journal of Medicinal Chemistry, 2019, 172, 95-108.	5.5	49
10	K-Ras and its inhibitors towards personalized cancer treatment: Pharmacological and structural perspectives. European Journal of Medicinal Chemistry, 2017, 125, 299-314.	5.5	39
11	An updated patent review of therapeutic applications of chalcone derivatives (2014-present). Expert Opinion on Therapeutic Patents, 2019, 29, 385-406.	5.0	38
12	MEK inhibitors in oncology: a patent review (2015-Present). Expert Opinion on Therapeutic Patents, 2017, 27, 887-906.	5.0	34
13	Design, synthesis and molecular modeling studies of novel thiazolidine-2,4-dione derivatives as potential anti-cancer agents. Journal of Molecular Structure, 2018, 1154, 406-417.	3.6	30
14	An exhaustive perspective on structural insights of SGLT2 inhibitors: A novel class of antidiabetic agent. European Journal of Medicinal Chemistry, 2020, 204, 112523.	5.5	25
15	Pyrazolopyrimidines as anticancer agents: A review on structural and target-based approaches. European Journal of Medicinal Chemistry, 2021, 225, 113781.	5.5	23
16	Virtual Screening, Molecular Docking, and DFT Studies of Some Thiazolidineâ€2,4â€diones as Potential PIMâ€1 Kinase Inhibitors. ChemistrySelect, 2018, 3, 127-135.	1.5	18
17	Integrated computational approach on sodium-glucose co-transporter 2 (SGLT2) Inhibitors for the development of novel antidiabetic agents. Journal of Molecular Structure, 2021, 1227, 129511.	3.6	18
18	Current Insights into the Chemistry and Antitubercular Potential of Benzimidazole and Imidazole Derivatives. Mini-Reviews in Medicinal Chemistry, 2021, 21, 643-657.	2.4	16

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19	Triggering PIK3CA Mutations in PI3K/Akt/mTOR Axis: Exploration of Newer Inhibitors and Rational Preventive Strategies. Current Pharmaceutical Design, 2016, 22, 6039-6054.	1.9	14
20	3D-QSAR and virtual screening studies of thiazolidine-2,4-dione analogs: Validation of experimental inhibitory potencies towards PIM-1 kinase. Journal of Molecular Structure, 2017, 1133, 278-293.	3.6	13
21	Mutant B-Raf Kinase Inhibitors as Anticancer Agents. Anti-Cancer Agents in Medicinal Chemistry, 2016, 16, 1558-1575.	1.7	13
22	Synthesis, characterization and antimicrobial evaluation of some 1,3-benzothiazole-2-yl-hydrazone derivatives. Arabian Journal of Chemistry, 2015, 8, 495-499.	4.9	10
23	The Recent Development of Piperazine and Piperidine Derivatives as Antipsychotic Agents. Mini-Reviews in Medicinal Chemistry, 2021, 21, 362-379.	2.4	10
24	Molecular modeling studies of some thiazolidine-2,4-dione derivatives as 15-PGDH inhibitors. Medicinal Chemistry Research, 2016, 25, 94-108.	2.4	9
25	In-silico studies for the development of novel RET inhibitors for cancer treatment. Journal of Molecular Structure, 2022, 1251, 132040.	3.6	8
26	Structural prediction of novel pyrazolo-pyrimidine derivatives against PIM-1 kinase: In-silico drug design studies. Journal of Molecular Structure, 2020, 1217, 128375.	3.6	7
27	A Computational approach to discover potential quinazoline derivatives against CDK4/6 kinase. Journal of Molecular Structure, 2021, 1245, 131079.	3.6	6
28	Nature Inspired Green Fabrication Technology for Silver Nanoparticles. Current Nanomedicine, 2017, 7, 5-24.	0.6	6
29	RECENT THERAPEUTIC PROGRESS OF CHALCONE SCAFFOLD BEARING COMPOUNDS AS PROSPECTIVE ANTI-GOUT CANDIDATES. Journal of Critical Reviews, 0, , 1-5.	0.1	5
30	Recent Advances in PI3 Kinase Inhibitors: Anticancer Activities and Structure-Activity Relationships. Mini-Reviews in Medicinal Chemistry, 2022, 22, 2146-2165.	2.4	5
31	Anti-Inflammatory activity of seeds extract of datura stramonium against carrageenan induced paw edema on albino wistar rats. Journal of Pharmaceutical and Biological Sciences, 2019, 7, 41-46.	0.4	4
32	Formulation and evaluation of sustain released matrix tablet of atenolol. Journal of Drug Delivery and Therapeutics, 2019, 9, 183-189.	0.5	4
33	SWFB and GA Strategies for Variable Selection in QSAR Studies for the Validation of Thiazolidine- 2,4-Dione Derivatives as Promising Antitumor Candidates. Indian Journal of Pharmaceutical Education and Research, 2017, 51, 436-451.	0.6	4
34	QSAR studies for some thiazolidine-2,4-dione derivatives as PIM-2 kinase inhibitors. Medicinal Chemistry Research, 2016, 25, 1329-1339.	2.4	3
35	Discovery of novel ALK2 inhibitors of pyrazolo-pyrimidines: A computational study. Journal of Biomolecular Structure and Dynamics, 2022, 40, 10422-10436.	3.5	3
36	Recent Perspectives Of Chalcone-Based Molecules As Protein Tyrosine Phosphatase 1b (Ptp1b)		3

Inhibitors. , 2019, , 235-251.

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37	Structural Aspects of mTOR Inhibitors: In Progress to Search Potential Compounds. Anti-Cancer Agents in Medicinal Chemistry, 2021, 21, .	1.7	2
38	3D-QSAR and Molecular Docking Studies on Oxadiazole Substituted Benzimidazole Derivatives: Validation of Experimental Inhibitory Potencies Towards COX-2. Current Computer-Aided Drug Design, 2019, 15, 277-293.	1.2	2
39	Discovery of Novel Antimalarial Drugs Based on Thiosemicarbazone Derivatives: An In Silico Approach. Current Signal Transduction Therapy, 2022, 17, 59-74.	0.5	0
40	An analytical review for the estimation of montelukast sodium. Separation Science Plus, 2022, 5, 120-137.	0.6	0