

# Penghua Li

## List of Publications by Year in descending order

Source: <https://exaly.com/author-pdf/30734/publications.pdf>

Version: 2024-02-01

10  
papers

185  
citations

1478505

6  
h-index

1372567

10  
g-index

10  
all docs

10  
docs citations

10  
times ranked

163  
citing authors

#	ARTICLE	IF	CITATIONS
1	Merging Reagent Modulation and Remote Anchimeric Assistance for Glycosylation: Highly Stereoselective Synthesis of $\alpha$ -Glycans up to a 30 $\mu$ mer. <i>Angewandte Chemie</i> , 2021, 133, 12705-12714.	2.0	6
2	Merging Reagent Modulation and Remote Anchimeric Assistance for Glycosylation: Highly Stereoselective Synthesis of $\beta$ -Glycans up to a 30 $\mu$ mer. <i>Angewandte Chemie - International Edition</i> , 2021, 60, 12597-12606.	13.8	47
3	An orthogonal and reactivity-based one-pot glycosylation strategy for both glycan and nucleoside synthesis: access to TMC-chitotriomycin, lipochitoooligosaccharides and capuramycin. <i>Chemical Science</i> , 2021, 12, 5143-5151.	7.4	32
4	Ortho-(1-phenylvinyl)benzyl glycosides: Ether-type glycosyl donors for the efficient synthesis of both O-glycosides and nucleosides. <i>Green Synthesis and Catalysis</i> , 2020, 1, 160-166.	6.8	11
5	Glycosyl ortho-(1-phenylvinyl)benzoates versatile glycosyl donors for highly efficient synthesis of both O-glycosides and nucleosides. <i>Nature Communications</i> , 2020, 11, 405.	12.8	57
6	Identification of potential AMPK activator by pharmacophore modeling, molecular docking and QSAR study. <i>Computational Biology and Chemistry</i> , 2019, 79, 165-176.	2.3	7
7	Discovery of novel indoleamine 2,3-dioxygenase 1 (IDO1) inhibitors by virtual screening. <i>Computational Biology and Chemistry</i> , 2019, 78, 306-316.	2.3	10
8	Discovery of FIXa inhibitors by combination of pharmacophore modeling, molecular docking, and 3D-QSAR modeling. <i>Journal of Receptor and Signal Transduction Research</i> , 2018, 38, 213-224.	2.5	6
9	Pharmacophore modeling, molecular docking and molecular dynamics simulations toward identifying lead compounds for Chk1. <i>Computational Biology and Chemistry</i> , 2018, 76, 53-60.	2.3	5
10	In silico Discovery of Novel FXa Inhibitors by Pharmacophore Modeling and Molecular Docking. <i>Natural Products and Bioprospecting</i> , 2017, 7, 249-256.	4.3	4