

# Na Le Dang

## List of Publications by Year in descending order

Source: <https://exaly.com/author-pdf/1176806/publications.pdf>

Version: 2024-02-01

16  
papers

553  
citations

933264

10  
h-index

1058333

14  
g-index

17  
all docs

17  
docs citations

17  
times ranked

1065  
citing authors

#	ARTICLE	IF	CITATIONS
1	Predictors of humoral response to SARS-CoV-2 mRNA vaccine BNT162b2 in patients receiving maintenance dialysis. <i>Antimicrobial Stewardship &amp; Healthcare Epidemiology</i> , 2022, 2, .	0.2	2
2	High Energy Channeling and Malleable Transition States: Molecular Dynamics Simulations and Free Energy Landscapes for the Thermal Unfolding of Protein U1A and 13 Mutants. <i>Biomolecules</i> , 2022, 12, 940.	1.8	0
3	“Black Box” to “Conversational” Machine Learning: Ondansetron Reduces Risk of Hospital-Acquired Venous Thromboembolism. <i>IEEE Journal of Biomedical and Health Informatics</i> , 2021, 25, 2204-2214.	3.9	24
4	Modeling the Bioactivation and Subsequent Reactivity of Drugs. <i>Chemical Research in Toxicology</i> , 2021, 34, 584-600.	1.7	11
5	SARS-CoV-2 Infection Risk Factors among Maintenance Hemodialysis Patients and Health Care Personnel In Outpatient Hemodialysis Centers. <i>Kidney360</i> , 2021, 2, 996-1001.	0.9	1
6	Deep Learning Coordinate-Free Quantum Chemistry. <i>Journal of Physical Chemistry A</i> , 2021, 125, 8978-8986.	1.1	5
7	Metabolic Forest: Predicting the Diverse Structures of Drug Metabolites. <i>Journal of Chemical Information and Modeling</i> , 2020, 60, 4702-4716.	2.5	11
8	XenoNet: Inference and Likelihood of Intermediate Metabolite Formation. <i>Journal of Chemical Information and Modeling</i> , 2020, 60, 3431-3449.	2.5	14
9	The Metabolic Rainbow: Deep Learning Phase I Metabolism in Five Colors. <i>Journal of Chemical Information and Modeling</i> , 2020, 60, 1146-1164.	2.5	26
10	Deep learning long-range information in undirected graphs with wave networks. , 2019, , .		8
11	Computationally Assessing the Bioactivation of Drugs by N-Dealkylation. <i>Chemical Research in Toxicology</i> , 2018, 31, 68-80.	1.7	30
12	Learning a Local-Variable Model of Aromatic and Conjugated Systems. <i>ACS Central Science</i> , 2018, 4, 52-62.	5.3	18
13	Computational Approach to Structural Alerts: Furans, Phenols, Nitroaromatics, and Thiophenes. <i>Chemical Research in Toxicology</i> , 2017, 30, 1046-1059.	1.7	32
14	Open Source Drug Discovery with the Malaria Box Compound Collection for Neglected Diseases and Beyond. <i>PLoS Pathogens</i> , 2016, 12, e1005763.	2.1	244
15	Modeling Reactivity to Biological Macromolecules with a Deep Multitask Network. <i>ACS Central Science</i> , 2016, 2, 529-537.	5.3	76
16	A simple model predicts UGT-mediated metabolism. <i>Bioinformatics</i> , 2016, 32, 3183-3189.	1.8	51