## Alvin Yu

## List of Publications by Year in descending order

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840776 1058476 14 489 11 14 citations h-index g-index papers 18 18 18 572 citing authors all docs docs citations times ranked

#	Article	IF	CITATIONS
1	Cooperative multivalent receptor binding promotes exposure of the SARS-CoV-2 fusion machinery core. Nature Communications, 2022, 13, 1002.	12.8	30
2	Strain and rupture of HIV-1 capsids during uncoating. Proceedings of the National Academy of Sciences of the United States of America, 2022, 119, e2117781119.	7.1	21
3	A multiscale coarse-grained model of the SARS-CoV-2 virion. Biophysical Journal, 2021, 120, 1097-1104.	0.5	139
4	Integrin-based mechanosensing through conformational deformation. Biophysical Journal, 2021, 120, 4349-4359.	0.5	10
5	Stability and molecular pathways to the formation of spin defects in silicon carbide. Nature Communications, 2021, 12, 6325.	12.8	9
6	Temperature and Phase Transferable Bottom-up Coarse-Grained Models. Journal of Chemical Theory and Computation, 2020, 16, 6823-6842.	5.3	36
7	Atomic-scale characterization of mature HIV-1 capsid stabilization by inositol hexakisphosphate (IP) Tj ETQq1 1 (	).784314 10.3	rgBT/Overloo
8	TRIM5 $\hat{l}_{\pm}$ self-assembly and compartmentalization of the HIV-1 viral capsid. Nature Communications, 2020, 11, 1307.	12.8	51
9	Off-Pathway Assembly: A Broad-Spectrum Mechanism of Action for Drugs That Undermine Controlled HIV-1 Viral Capsid Formation. Journal of the American Chemical Society, 2019, 141, 10214-10224.	13.7	38
10	Neurotransmitter Funneling Optimizes Glutamate Receptor Kinetics. Neuron, 2018, 97, 139-149.e4.	8.1	25
11	Glutamate and Glycine Binding to the NMDA Receptor. Structure, 2018, 26, 1035-1043.e2.	3.3	42
12	Energetics of Glutamate Binding to an Ionotropic Glutamate Receptor. Journal of Physical Chemistry B, 2017, 121, 10436-10442.	2.6	18
13	Molecular lock regulates binding of glycine to a primitive NMDA receptor. Proceedings of the National Academy of Sciences of the United States of America, 2016, 113, E6786-E6795.	7.1	30
14	Computing Conformational Free Energies of iGluR Ligand-Binding Domains. Neuromethods, 2016, , 119-132.	0.3	4