

# Billy K Poon

## List of Publications by Year in descending order

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Version: 2024-02-01

25  
papers

7,844  
citations

623734  
14  
h-index

677142  
22  
g-index

26  
all docs

26  
docs citations

26  
times ranked

12283  
citing authors

#	ARTICLE	IF	CITATIONS
1	Macromolecular structure determination using X-rays, neutrons and electrons: recent developments in <i>&lt; i&gt;Phenix&lt;/i&gt;</i> . <i>Acta Crystallographica Section D: Structural Biology</i> , 2019, 75, 861-877.	2.3	4,060
2	Real-space refinement in <i>&lt; i&gt;PHENIX&lt;/i&gt;</i> for cryo-EM and crystallography. <i>Acta Crystallographica Section D: Structural Biology</i> , 2018, 74, 531-544.	2.3	2,065
3	New tools for the analysis and validation of cryo-EM maps and atomic models. <i>Acta Crystallographica Section D: Structural Biology</i> , 2018, 74, 814-840.	2.3	575
4	Polder maps: improving OMIT maps by excluding bulk solvent. <i>Acta Crystallographica Section D: Structural Biology</i> , 2017, 73, 148-157.	2.3	500
5	Shapes and vorticities of superfluid helium nanodroplets. <i>Science</i> , 2014, 345, 906-909.	12.6	197
6	A New Method for Coarse-Grained Elastic Normal-Mode Analysis. <i>Journal of Chemical Theory and Computation</i> , 2006, 2, 464-471.	5.3	81
7	Normal mode refinement of anisotropic thermal parameters for a supramolecular complex at 3.42-Å crystallographic resolution. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2007, 104, 7869-7874.	7.1	56
8	Announcing mandatory submission of PDBx/mmCIF format files for crystallographic depositions to the Protein Data Bank (PDB). <i>Acta Crystallographica Section D: Structural Biology</i> , 2019, 75, 451-454.	2.3	46
9	Three-dimensional single-particle imaging using angular correlations from X-ray laser data. <i>Acta Crystallographica Section A: Foundations and Advances</i> , 2013, 69, 365-373.	0.3	45
10	Cryo_fit: Democratization of flexible fitting for cryo-EM. <i>Journal of Structural Biology</i> , 2019, 208, 1-6.	2.8	30
11	Normal-Mode Refinement of Anisotropic Thermal Parameters for Potassium Channel KcsA at 3.2 Å... Crystallographic Resolution. <i>Structure</i> , 2007, 15, 955-962.	3.3	29
12	Autoindexing with outlier rejection and identification of superimposed lattices. <i>Journal of Applied Crystallography</i> , 2010, 43, 611-616.	4.5	24
13	Structure determination of Pt-coated Au dumbbells <i>&lt; i&gt;via&lt;/i&gt;</i> fluctuation X-ray scattering. <i>Journal of Synchrotron Radiation</i> , 2012, 19, 695-700.	2.4	23
14	Structural improvement of unliganded simian immunodeficiency virus gp120 core by normal-mode-based X-ray crystallographic refinement. <i>Acta Crystallographica Section D: Biological Crystallography</i> , 2009, 65, 339-347.	2.5	15
15	<i>&lt; i&gt;CERES&lt;/i&gt;</i> : a cryo-EM re-refinement system for continuous improvement of deposited models. <i>Acta Crystallographica Section D: Structural Biology</i> , 2021, 77, 48-61.	2.3	14
16	Detection and correction of underassigned rotational symmetry prior to structure deposition. <i>Acta Crystallographica Section D: Biological Crystallography</i> , 2010, 66, 503-513.	2.5	13
17	Computation of fluctuation scattering profiles <i>&lt; i&gt;via&lt;/i&gt;</i> three-dimensional Zernike polynomials. <i>Acta Crystallographica Section A: Foundations and Advances</i> , 2012, 68, 561-567.	0.3	13
18	Interactive comparison and remediation of collections of macromolecular structures. <i>Protein Science</i> , 2018, 27, 182-194.	7.6	13

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19	Application of normal-mode refinement to X-ray crystal structures at the lower resolution limit. <i>Acta Crystallographica Section D: Biological Crystallography</i> , 2009, 65, 633-643.		2.5	12
20	Integration of software tools for integrative modeling of biomolecular systems. <i>Journal of Structural Biology</i> , 2022, 214, 107841.		2.8	7
21	Free-electron laser data for multiple-particle fluctuation scattering analysis. <i>Scientific Data</i> , 2018, 5, 180201.		5.3	6
22	Sizes of pure and doped helium droplets from single shot x-ray imaging. <i>Journal of Chemical Physics</i> , 2022, 156, 041102.		3.0	3
23	Polder maps: improving OMIT maps for ligand building and validation. <i>Acta Crystallographica Section A: Foundations and Advances</i> , 2017, 73, C48-C48.		0.1	0
24	Video tutorials for the <i>Phenix</i> software suite. <i>Acta Crystallographica Section A: Foundations and Advances</i> , 2017, 73, C1134-C1134.		0.1	0
25	Polder maps: improving OMIT maps for ligand building and validation. <i>Acta Crystallographica Section A: Foundations and Advances</i> , 2018, 74, a308-a308.		0.1	0