

Massimo Sammito

List of Publications by Year in descending order

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Version: 2024-02-01

47
papers

4,794
citations

516215

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454577

30
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docs citations

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times ranked

7957
citing authors

#	ARTICLE	IF	CITATIONS
1	Implications of <i>AlphaFold</i> ² for crystallographic phasing by molecular replacement. <i>Acta Crystallographica Section D: Structural Biology</i> , 2022, 78, 1-13.	1.1	65
2	Phasertng: directed acyclic graphs for crystallographic phasing. <i>Acta Crystallographica Section D: Structural Biology</i> , 2021, 77, 1-10.	1.1	10
3	Detection of translational noncrystallographic symmetry in Patterson functions. <i>Acta Crystallographica Section D: Structural Biology</i> , 2021, 77, 131-141.	1.1	5
4	Structure of the class XI myosin globular tail reveals evolutionary hallmarks for cargo recognition in plants. <i>Acta Crystallographica Section D: Structural Biology</i> , 2021, 77, 522-533.	1.1	0
5	Assessing the utility of <i>CASP14</i> models for molecular replacement. <i>Proteins: Structure, Function and Bioinformatics</i> , 2021, 89, 1752-1769.	1.5	47
6	The amphibian antimicrobial peptide uperin 3.5 is a cross- β /cross- β^2 chameleon functional amyloid. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2021, 118, .	3.3	41
7	<i>SEQUENCE SLIDER</i> : expanding polyalanine fragments for phasing with multiple side-chain hypotheses. <i>Acta Crystallographica Section D: Structural Biology</i> , 2020, 76, 221-237.	1.1	10
8	<i>ALEPH</i> : a network-oriented approach for the generation of fragment-based libraries and for structure interpretation. <i>Acta Crystallographica Section D: Structural Biology</i> , 2020, 76, 193-208.	1.1	16
9	Factors influencing estimates of coordinate error for molecular replacement. <i>Acta Crystallographica Section D: Structural Biology</i> , 2020, 76, 19-27.	1.1	6
10	Evaluation of template-based modeling in <i>CASP13</i> . <i>Proteins: Structure, Function and Bioinformatics</i> , 2019, 87, 1113-1127.	1.5	56
11	Evaluation of model refinement in <i>CASP13</i> . <i>Proteins: Structure, Function and Bioinformatics</i> , 2019, 87, 1249-1262.	1.5	28
12	Macromolecular structure determination using X-rays, neutrons and electrons: recent developments in <i>Phenix</i> . <i>Acta Crystallographica Section D: Structural Biology</i> , 2019, 75, 861-877.	1.1	4,060
13	Dealing with modulated macromolecular structures with translational non-crystallographic symmetry. <i>Acta Crystallographica Section A: Foundations and Advances</i> , 2019, 75, e197-e197.	0.0	0
14	<i>ALEPH</i> : a new software for structural analysis and generation of fragment libraries for molecular replacement. <i>Acta Crystallographica Section A: Foundations and Advances</i> , 2019, 75, e184-e184.	0.0	0
15	A new 3D reflection data viewer based on NGL. <i>Acta Crystallographica Section A: Foundations and Advances</i> , 2019, 75, e165-e165.	0.0	0
16	<i>Phaser.Voyager</i> : data-guided model generation and visualization. <i>Acta Crystallographica Section A: Foundations and Advances</i> , 2019, 75, e182-e182.	0.0	1
17	Combining the use of libraries from several distant homologs in <i>ARCIMBOLDO_SHREDDER</i> spheres. <i>Acta Crystallographica Section A: Foundations and Advances</i> , 2019, 75, e178-e178.	0.0	0
18	Exploiting distant homologues for phasing through the generation of compact fragments, local fold refinement and partial solution combination. <i>Acta Crystallographica Section D: Structural Biology</i> , 2018, 74, 290-304.	1.1	30

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19	<i>ARCIMBOLDO</i> on coiled coils. Acta Crystallographica Section D: Structural Biology, 2018, 74, 194-204.	1.1	41
20	Gyreandgimble: a maximum-likelihood replacement for Patterson correlation refinement. Acta Crystallographica Section D: Structural Biology, 2018, 74, 279-289.	1.1	14
21	On the application of the expected log-likelihood gain to decision making in molecular replacement. Acta Crystallographica Section D: Structural Biology, 2018, 74, 245-255.	1.1	40
22	Overcoming phasing difficulties in coiled coils with ARCIMBOLDO_LITE: verifying solutions. Acta Crystallographica Section A: Foundations and Advances, 2018, 74, e229-e229.	0.0	0
23	The expected log-likelihood gain for decision making in molecular replacement. Acta Crystallographica Section A: Foundations and Advances, 2018, 74, e411-e411.	0.0	0
24	Expanding partial structures by assembling most probable side-chain composition. Acta Crystallographica Section A: Foundations and Advances, 2018, 74, e410-e410.	0.0	0
25	Revealing the properties of small local folds with ALEPH: from structure annotation to ab initio phasing. Acta Crystallographica Section A: Foundations and Advances, 2018, 74, e228-e229.	0.0	0
26	New in the <i>ARCIMBOLDO</i> toolbox for phasing with small fragments. Acta Crystallographica Section A: Foundations and Advances, 2017, 73, C881-C881.	0.0	0
27	Vaccinia Virus Immunomodulator A46: A Lipid and Protein-Binding Scaffold for Sequestering Host TIR-Domain Proteins. PLoS Pathogens, 2016, 12, e1006079.	2.1	19
28	PLA2s-like membrane perturbation mechanism: extracting the most of crystallography data. Acta Crystallographica Section A: Foundations and Advances, 2016, 72, s226-s226.	0.0	1
29	ARCIMBOLDO_SHREDDER's contribution to MR: phasing with fragments from distant homologs. Acta Crystallographica Section A: Foundations and Advances, 2016, 72, s198-s199.	0.0	0
30	BORGES_MATRIX: a tool to generate models for ab initio phasing and for structure interpretation. Acta Crystallographica Section A: Foundations and Advances, 2016, 72, s23-s23.	0.0	1
31	Macromolecular <i>ab initio</i> phasing enforcing secondary and tertiary structure. IUCr, 2015, 2, 95-105.	1.0	47
32	SEQUENCE SLIDER: a multi sequence evaluator and its application in venomics. Acta Crystallographica Section A: Foundations and Advances, 2015, 71, s167-s167.	0.0	1
33	BORGESlibraries: from phasing to structural bioinformatics. Acta Crystallographica Section A: Foundations and Advances, 2015, 71, s271-s272.	0.0	0
34	ARCIMBOLDO, an ab initio approach to MR phasing. Acta Crystallographica Section A: Foundations and Advances, 2015, 71, s20-s21.	0.0	0
35	<i>ARCIMBOLDO_LITE</i>: single-workstation implementation and use. Acta Crystallographica Section D: Biological Crystallography, 2015, 71, 1921-1930.	2.5	51
36	Combining phase information in reciprocal space for molecular replacement with partial models. Acta Crystallographica Section D: Biological Crystallography, 2015, 71, 1931-1945.	2.5	12

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37	Structure of a 13-fold superhelix (almost) determined from first principles. IUCr, 2015, 2, 177-187.	1.0	3
38	Use of clustering algorithms to combine partial solutions in reciprocal space. Acta Crystallographica Section A: Foundations and Advances, 2015, 71, s165-s165.	0.0	0
39	Structure solution with <scp>ARCIMBOLDO</scp> using fragments derived from distant homology models. FEBS Journal, 2014, 281, 4029-4045.	2.2	41
40	Structure solution of DNA-binding proteins and complexes with<i>ARCIMBOLDO</i>libraries. Acta Crystallographica Section D: Biological Crystallography, 2014, 70, 1743-1757.	2.5	19
41	Exploiting tertiary structure through local folds for crystallographic phasing. Nature Methods, 2013, 10, 1099-1101.	9.0	63
42	BORGES-ARCIMBOLDOexploiting tertiary folds as fragments libraries for phasing. Acta Crystallographica Section A: Foundations and Advances, 2013, 69, s670-s670.	0.3	1
43	Using<i>PHASER</i>for phasing in conjunction with wide-convergence refinement and model building techniques. Acta Crystallographica Section A: Foundations and Advances, 2013, 69, s40-s40.	0.3	2
44	Phasing Through Location of Small Fragments and Density Modification with ARCIMBOLDO. NATO Science for Peace and Security Series A: Chemistry and Biology, 2013, , 123-133.	0.5	1
45	BORGES-ARCIMBOLDOexploiting tertiary folds as fragments libraries for phasing. Acta Crystallographica Section A: Foundations and Advances, 2013, 69, s271-s271.	0.3	0
46	Reciprocal space clustering ofBORGES-ARCIMBOLDOpartial solutions: practical cases. Acta Crystallographica Section A: Foundations and Advances, 2013, 69, s294-s294.	0.3	0
47	Practical structure solution with <i><i>ARCIMBOLDO</i></i>. Acta Crystallographica Section D: Biological Crystallography, 2012, 68, 336-343.	2.5	50