

Carles Pons

List of Publications by Year in descending order

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41
papers

3,871
citations

257101

24
h-index

301761

39
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45
all docs

45
docs citations

45
times ranked

5893
citing authors

#	ARTICLE	IF	CITATIONS
1	Analysing the yeast complexomeâ€”the Complex Portal rising to the challenge. <i>Nucleic Acids Research</i> , 2021, 49, 3156-3167.	6.5	5
2	Environmental robustness of the global yeast genetic interaction network. <i>Science</i> , 2021, 372, .	6.0	40
3	Natural variants suppress mutations in hundreds of essential genes. <i>Molecular Systems Biology</i> , 2021, 17, e10138.	3.2	13
4	Bioactivity Profile Similarities to Expand the Repertoire of COVID-19 Drugs. <i>Journal of Chemical Information and Modeling</i> , 2020, 60, 5730-5734.	2.5	3
5	A reference map of the human binary protein interactome. <i>Nature</i> , 2020, 580, 402-408.	13.7	724
6	Systematic genetics and singleâ€”cell imaging reveal widespread morphological pleiotropy and cellâ€”toâ€”cell variability. <i>Molecular Systems Biology</i> , 2020, 16, e9243.	3.2	37
7	Systematic analysis of bypass suppression of essential genes. <i>Molecular Systems Biology</i> , 2020, 16, e9828.	3.2	45
8	Systematic analysis of complex genetic interactions. <i>Science</i> , 2018, 360, .	6.0	201
9	A PanorOmic view of personal cancer genomes. <i>Nucleic Acids Research</i> , 2017, 45, W195-W200.	6.5	6
10	Features of the Chaperone Cellular Network Revealed through Systematic Interaction Mapping. <i>Cell Reports</i> , 2017, 20, 2735-2748.	2.9	47
11	A framework for exhaustively mapping functional missense variants. <i>Molecular Systems Biology</i> , 2017, 13, 957.	3.2	146
12	Mechanisms of suppression: The wiring of genetic resilience. <i>BioEssays</i> , 2017, 39, 1700042.	1.2	31
13	A geneâ€”centered <i>C.Âlegans</i> proteinâ€”DNA interaction network provides a framework for functional predictions. <i>Molecular Systems Biology</i> , 2016, 12, 884.	3.2	71
14	A global genetic interaction network maps a wiring diagram of cellular function. <i>Science</i> , 2016, 353, .	6.0	979
15	Exploring genetic suppression interactions on a global scale. <i>Science</i> , 2016, 354, .	6.0	157
16	Genetic Interactions Implicating Postreplicative Repair in Okazaki Fragment Processing. <i>PLoS Genetics</i> , 2015, 11, e1005659.	1.5	24
17	pyDockSAXS: proteinâ€”protein complex structure by SAXS and computational docking. <i>Nucleic Acids Research</i> , 2015, 43, W356-W361.	6.5	61
18	Transcription Factor Activity Mapping of a Tissue-Specific InÂVivo Gene Regulatory Network. <i>Cell Systems</i> , 2015, 1, 152-162.	2.9	64

#	ARTICLE	IF	CITATIONS
19	Linking Genetics to Structural Biology: Complex Heterozygosity Screening with Actin Alanine Scan Alleles Identifies Functionally Related Surfaces on Yeast Actin. <i>G3: Genes, Genomes, Genetics</i> , 2014, 4, 1491-1501.	0.8	1
20	Expanding the frontiers of protein-protein modeling: From docking and scoring to binding affinity predictions and other challenges. <i>Proteins: Structure, Function and Bioinformatics</i> , 2013, 81, 2192-2200.	1.5	20
21	pyDockWEB: a web server for rigid-body protein-protein docking using electrostatics and desolvation scoring. <i>Bioinformatics</i> , 2013, 29, 1698-1699.	1.8	214
22	Efficient Relaxation of Protein-Protein Interfaces by Discrete Molecular Dynamics Simulations. <i>Journal of Chemical Theory and Computation</i> , 2013, 9, 1222-1229.	2.3	13
23	Extensive Rewiring and Complex Evolutionary Dynamics in a <i>C.Âlegans</i> Multiparameter Transcription Factor Network. <i>Molecular Cell</i> , 2013, 51, 116-127.	4.5	83
24	Validated Conformational Ensembles Are Key for the Successful Prediction of Protein Complexes. <i>Journal of Chemical Theory and Computation</i> , 2013, 9, 1830-1837.	2.3	4
25	Cell-Dock: high-performance protein-protein docking. <i>Bioinformatics</i> , 2012, 28, 2394-2396.	1.8	14
26	H-bond network optimization in protein-protein complexes: Are all atom force field scores enough?. <i>Proteins: Structure, Function and Bioinformatics</i> , 2012, 80, 818-824.	1.5	12
27	Towards the prediction of protein interaction partners using physical docking. <i>Molecular Systems Biology</i> , 2011, 7, 469.	3.2	102
28	Scoring by Intermolecular Pairwise Propensities of Exposed Residues (SIPPER): A New Efficient Potential for Protein-Protein Docking. <i>Journal of Chemical Information and Modeling</i> , 2011, 51, 370-377.	2.5	70
29	Community-Wide Assessment of Protein-Interface Modeling Suggests Improvements to Design Methodology. <i>Journal of Molecular Biology</i> , 2011, 414, 289-302.	2.0	131
30	Prediction of protein-binding areas by small-world residue networks and application to docking. <i>BMC Bioinformatics</i> , 2011, 12, 378.	1.2	46
31	Present and future challenges and limitations in protein-protein docking. <i>Proteins: Structure, Function and Bioinformatics</i> , 2010, 78, 95-108.	1.5	76
32	Optimization of pyDock for the new CAPRI challenges: Docking of homology-based models, domain-domain assembly and protein-RNA binding. <i>Proteins: Structure, Function and Bioinformatics</i> , 2010, 78, 3182-3188.	1.5	25
33	Structural Characterization of Protein-Protein Complexes by Integrating Computational Docking with Small-angle Scattering Data. <i>Journal of Molecular Biology</i> , 2010, 403, 217-230.	2.0	64
34	STRUCTURAL PREDICTION OF PROTEIN-RNA INTERACTION BY COMPUTATIONAL DOCKING WITH PROPENSITY-BASED STATISTICAL POTENTIALS. , 2009, , 293-301.		33
35	Pushing Structural Information into the Yeast Interactome by High-Throughput Protein Docking Experiments. <i>PLoS Computational Biology</i> , 2009, 5, e1000490.	1.5	67
36	FRODOCK: a new approach for fast rotational protein-protein docking. <i>Bioinformatics</i> , 2009, 25, 2544-2551.	1.8	126

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37	On the Use of low-resolution Data to Improve Structure Prediction of Proteins and Protein Complexes. <i>Journal of Chemical Theory and Computation</i> , 2009, 5, 3129-3137.	2.3	7
38	Enhancing the prediction of protein pairings between interacting families using orthology information. <i>BMC Bioinformatics</i> , 2008, 9, 35.	1.2	21
39	Prediction and scoring of docking poses with pyDock. <i>Proteins: Structure, Function and Bioinformatics</i> , 2007, 69, 852-858.	1.5	40
40	TSEMA: interactive prediction of protein pairings between interacting families. <i>Nucleic Acids Research</i> , 2006, 34, W315-W319.	6.5	25
41	Computational paradigms for analyzing genetic interaction networks. , 0, , 12-35.		0