

Michel A Cuendet

List of Publications by Year in descending order

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49
papers

3,646
citations

377584

21
h-index

355658

38
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53
all docs

53
docs citations

53
times ranked

6196
citing authors

#	ARTICLE	IF	CITATIONS
1	Safety and Efficacy of Ipilimumab plus Nivolumab and Sequential Selective Internal Radiation Therapy in Hepatic and Extrahepatic Metastatic Uveal Melanoma. <i>Cancers</i> , 2022, 14, 1162.	1.7	9
2	A Graphic Encoding Method for Quantitative Classification of Protein Structure and Representation of Conformational Changes. <i>IEEE/ACM Transactions on Computational Biology and Bioinformatics</i> , 2021, 18, 1336-1349.	1.9	4
3	A lightweight method for evaluating in situ workflow efficiency. <i>Journal of Computational Science</i> , 2021, 48, 101259.	1.5	7
4	Digitalization is Fueling a Revolution in Precision Oncology. , 2021, , 342-346.		1
5	A Process Mining Approach to Statistical Analysis: Application to a Real-World Advanced Melanoma Dataset. <i>Lecture Notes in Business Information Processing</i> , 2021, , 291-304.	0.8	4
6	Probing the Conformational Dynamics of Affinity-Enhanced T Cell Receptor Variants upon Binding the Peptide-Bound Major Histocompatibility Complex by Hydrogen/Deuterium Exchange Mass Spectrometry. <i>Biochemistry</i> , 2021, 60, 859-872.	1.2	3
7	Unusual mode of dimerization of retinitis pigmentosa-associated F220C rhodopsin. <i>Scientific Reports</i> , 2021, 11, 10536.	1.6	7
8	Turning tumors from cold to inflamed to improve immunotherapy response. <i>Cancer Treatment Reviews</i> , 2021, 101, 102227.	3.4	42
9	Acquired resistance to anti-MAPK targeted therapy confers an immune-evasive tumor microenvironment and cross-resistance to immunotherapy in melanoma. <i>Nature Cancer</i> , 2021, 2, 693-708.	5.7	102
10	Truly privacy-preserving federated analytics for precision medicine with multiparty homomorphic encryption. <i>Nature Communications</i> , 2021, 12, 5910.	5.8	64
11	The combination of stereotactic radiosurgery with immune checkpoint inhibition or targeted therapy in melanoma patients with brain metastases: a retrospective study. <i>Journal of Neuro-Oncology</i> , 2020, 146, 181-193.	1.4	18
12	What Role Can Process Mining Play in Recurrent Clinical Guidelines Issues? A Position Paper. <i>International Journal of Environmental Research and Public Health</i> , 2020, 17, 6616.	1.2	12
13	Long-lasting, irreversible and late-onset immune-related adverse events (irAEs) from immune checkpoint inhibitors (ICIs): A real-world data analysis.. <i>Journal of Clinical Oncology</i> , 2020, 38, e15095-e15095.	0.8	3
14	A process mining approach to real-world advanced melanoma treatments.. <i>Journal of Clinical Oncology</i> , 2020, 38, e22040-e22040.	0.8	1
15	A Novel Metric to Evaluate In Situ Workflows. <i>Lecture Notes in Computer Science</i> , 2020, , 538-553.	1.0	2
16	CD8 Binding of MHC-Peptide Complexes in cis or trans Regulates CD8+ T-cell Responses. <i>Journal of Molecular Biology</i> , 2019, 431, 4941-4958.	2.0	7
17	Characterizing In Situ and In Transit Analytics of Molecular Dynamics Simulations for Next-Generation Supercomputers. , 2019, , .		4
18	Thermodynamic Coupling Function Analysis of Allosteric Mechanisms in the Human Dopamine Transporter. <i>Biophysical Journal</i> , 2018, 114, 10-14.	0.2	18

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19	Thermodynamic Coupling Function Analysis of Allosteric Coupling between Na ⁺ Release and Inward-Opening in the Human Dopamine Transporter. <i>Biophysical Journal</i> , 2018, 114, 421a.	0.2	0
20	Graphic Encoding of Macromolecules for Efficient High-Throughput Analysis. , 2018, , .		4
21	Substrate-modulated unwinding of transmembrane helices in the NSS transporter LeuT. <i>Science Advances</i> , 2018, 4, eaar6179.	4.7	47
22	Endpoint-restricted adiabatic free energy dynamics approach for the exploration of biomolecular conformational equilibria. <i>Journal of Chemical Physics</i> , 2018, 149, 072316.	1.2	11
23	Serum protein predictors of long term survival from combined ipilimumab and nivolumab therapy in metastatic melanoma patients.. <i>Journal of Clinical Oncology</i> , 2018, 36, e21513-e21513.	0.8	0
24	Rigorous Representation of Transporters as Allosteric Machines Enables a Quantitative Understanding of their Functional Mechanisms. <i>Biophysical Journal</i> , 2017, 112, 308a.	0.2	0
25	The Allosteric Landscape: Quantifying Thermodynamic Couplings in Biomolecular Systems. <i>Biophysical Journal</i> , 2017, 112, 354a.	0.2	0
26	Vectorial Cholesterol Transport by STARD4 is Mediated by Specific PIP 2 Membrane Composition. <i>Biophysical Journal</i> , 2017, 112, 87a.	0.2	0
27	Cholesterol Promotes Protein Binding by Affecting Membrane Electrostatics and Solvation Properties. <i>Biophysical Journal</i> , 2017, 113, 2004-2015.	0.2	38
28	The T-Cell Receptor Can Bind to the Peptide-Bound Major Histocompatibility Complex and Uncomplexed β_2 -Microglobulin through Distinct Binding Sites. <i>Biochemistry</i> , 2017, 56, 3945-3961.	1.2	8
29	Exploratory analysis of multiprotein serum predictors at baseline of progression-free survival of ipilimumab or ipilimumab and nivolumab in the Checkmate-069 study.. <i>Journal of Clinical Oncology</i> , 2017, 35, 9571-9571.	0.8	0
30	The Allosteric Landscape: Quantifying Thermodynamic Couplings in Biomolecular Systems. <i>Journal of Chemical Theory and Computation</i> , 2016, 12, 5758-5767.	2.3	33
31	Dissecting Functional Correlates of a Double Mutation Enhancing GltPh Transport Efficiency using Alchemical Free Energy Calculations. <i>Biophysical Journal</i> , 2016, 110, 627a.	0.2	0
32	From Physics to Phenotype: New Insights into Allosteric Transport Mechanisms in LeuT. <i>Biophysical Journal</i> , 2016, 110, 362a.	0.2	0
33	Allosteric Mechanisms of Molecular Machines at the Membrane: Transport by Sodium-Coupled Symporters. <i>Chemical Reviews</i> , 2016, 116, 6552-6587.	23.0	71
34	The SIB Swiss Institute of Bioinformaticsâ€™ resources: focus on curated databases. <i>Nucleic Acids Research</i> , 2016, 44, D27-D37.	6.5	64
35	Transport domain unlocking sets the uptake rate of an aspartate transporter. <i>Nature</i> , 2015, 518, 68-73.	13.7	144
36	Molecular Dynamics Simulation Study of a Mutant Construct of the Archaeal Glutamate Transporter GltPh with Transport Rates as Fast as its Human Counterpart. <i>Biophysical Journal</i> , 2015, 108, 198a.	0.2	0

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37	Free Energy Reconstruction from Metadynamics or Adiabatic Free Energy Dynamics Simulations. <i>Journal of Chemical Theory and Computation</i> , 2014, 10, 2975-2986.	2.3	53
38	How Accurately Do Current Force Fields Predict Experimental Peptide Conformations? An Adiabatic Free Energy Dynamics Study. <i>Journal of Physical Chemistry B</i> , 2014, 118, 6539-6552.	1.2	33
39	Heating and flooding: A unified approach for rapid generation of free energy surfaces. <i>Journal of Chemical Physics</i> , 2012, 137, 024102.	1.2	66
40	Alchemical Free Energy Differences in Flexible Molecules from Thermodynamic Integration or Free Energy Perturbation Combined with Driven Adiabatic Dynamics. <i>Journal of Chemical Theory and Computation</i> , 2012, 8, 3504-3512.	2.3	17
41	How T cell receptors interact with peptide-MHCs: A multiple steered molecular dynamics study. <i>Proteins: Structure, Function and Bioinformatics</i> , 2011, 79, 3007-3024.	1.5	43
42	SwissParam: A fast force field generation tool for small organic molecules. <i>Journal of Computational Chemistry</i> , 2011, 32, 2359-2368.	1.5	1,485
43	Use of the FACTS solvation model for protein-ligand docking calculations. Application to EADock. <i>Journal of Molecular Recognition</i> , 2010, 23, 457-461.	1.1	28
44	Implementation of the CHARMM Force Field in GROMACS: Analysis of Protein Stability Effects from Correction Maps, Virtual Interaction Sites, and Water Models. <i>Journal of Chemical Theory and Computation</i> , 2010, 6, 459-466.	2.3	866
45	Protein-Protein Interaction Investigated by Steered Molecular Dynamics: The TCR-pMHC Complex. <i>Biophysical Journal</i> , 2008, 95, 3575-3590.	0.2	117
46	Optical Spectra of Cu(II)-Azurin by Hybrid TDDFT-Molecular Dynamics Simulations. <i>Journal of Physical Chemistry B</i> , 2007, 111, 10248-10252.	1.2	38
47	On the calculation of velocity-dependent properties in molecular dynamics simulations using the leapfrog integration algorithm. <i>Journal of Chemical Physics</i> , 2007, 127, 184102.	1.2	95
48	The Jarzynski identity derived from general Hamiltonian or non-Hamiltonian dynamics reproducing NVT or NPT ensembles. <i>Journal of Chemical Physics</i> , 2006, 125, 144109.	1.2	41
49	Statistical Mechanical Derivation of Jarzynski's Identity for Thermostated Non-Hamiltonian Dynamics. <i>Physical Review Letters</i> , 2006, 96, 120602.	2.9	32