

# SÃ-ivia Osuna Oliveras

## List of Publications by Year in descending order

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97  
papers

4,370  
citations

93792

39  
h-index

134545

62  
g-index

107  
all docs

107  
docs citations

107  
times ranked

5201  
citing authors

#	ARTICLE	IF	CITATIONS
1	Discovery and In Vivo Proof of Concept of a Highly Potent Dual Inhibitor of Soluble Epoxide Hydrolase and Acetylcholinesterase for the Treatment of Alzheimer's Disease. <i>Journal of Medicinal Chemistry</i> , 2022, 65, 4909-4925.	2.9	22
2	Time Evolution of the Millisecond Allosteric Activation of Imidazole Glycerol Phosphate Synthase. <i>Journal of the American Chemical Society</i> , 2022, 144, 7146-7159.	6.6	24
3	The challenge of predicting distal active site mutations in computational enzyme design. <i>Wiley Interdisciplinary Reviews: Computational Molecular Science</i> , 2021, 11, e1502.	6.2	61
4	Insights into the molecular determinants of thermal stability in halohydrin dehalogenase HheD2. <i>FEBS Journal</i> , 2021, 288, 4683-4701.	2.2	5
5	Pervasive cooperative mutational effects on multiple catalytic enzyme traits emerge via long-range conformational dynamics. <i>Nature Communications</i> , 2021, 12, 1621.	5.8	72
6	From the Design to the <i>In Vivo</i> Evaluation of Benzohomoadamantane-Derived Soluble Epoxide Hydrolase Inhibitors for the Treatment of Acute Pancreatitis. <i>Journal of Medicinal Chemistry</i> , 2021, 64, 5429-5446.	2.9	12
7	Biocatalysis. <i>Nature Reviews Methods Primers</i> , 2021, 1, .	11.8	255
8	SBMOpenMM: A Builder of Structure-Based Models for OpenMM. <i>Journal of Chemical Information and Modeling</i> , 2021, 61, 3166-3171.	2.5	3
9	<i>In Silico</i> Identification and Experimental Validation of Distal Activity-Enhancing Mutations in Tryptophan Synthase. <i>ACS Catalysis</i> , 2021, 11, 13733-13743.	5.5	30
10	Protein-directed crystalline 2D fullerene assemblies. <i>Nanoscale</i> , 2020, 12, 3614-3622.	2.8	11
11	Computational NMR Spectra of <i>o</i> -Benzynes and Stable Guests and Their Hemicarceplexes. <i>Chemistry - A European Journal</i> , 2020, 26, 2626-2634.	1.7	4
12	2-Oxadadamant-1-yl Ureas as Soluble Epoxide Hydrolase Inhibitors: <i>In Vivo</i> Evaluation in a Murine Model of Acute Pancreatitis. <i>Journal of Medicinal Chemistry</i> , 2020, 63, 9237-9257.	2.9	14
13	Enzyme Conformation Influences the Performance of Lipase-powered Nanomotors. <i>Angewandte Chemie - International Edition</i> , 2020, 59, 21080-21087.	7.2	58
14	Mutational Analysis of Linalool Dehydratase Isomerase Suggests That Alcohol and Alkene Transformations Are Catalyzed Using Noncovalent Mechanisms. <i>ACS Catalysis</i> , 2020, 10, 11136-11146.	5.5	4
15	Conformational Landscapes of Halohydrin Dehalogenases and Their Accessible Active Site Tunnels. <i>Catalysts</i> , 2020, 10, 1403.	1.6	9
16	<i>In Vivo</i> Selection for Formate Dehydrogenases with High Efficiency and Specificity toward NADP <sup>+</sup> . <i>ACS Catalysis</i> , 2020, 10, 7512-7525.	5.5	51
17	Regio- and Stereoselective Steroid Hydroxylation at C7 by Cytochrome P450 Monooxygenase Mutants. <i>Angewandte Chemie - International Edition</i> , 2020, 59, 12499-12505.	7.2	83
18	Regioselective Synthesis and Characterization of Tris- and Tetra-Prato Adducts of M3N@C80 (M = Y, Tj) ETQq0 0 0 rgBT /Overlock 10 Tf	6.6	14

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19	Regio- and Stereoselective Steroid Hydroxylation at C7 by Cytochrome P450 Monooxygenase Mutants. <i>Angewandte Chemie</i> , 2020, 132, 12599-12605.	1.6	19
20	Deciphering the Allosterically Driven Conformational Ensemble in Tryptophan Synthase Evolution. <i>Journal of the American Chemical Society</i> , 2019, 141, 13049-13056.	6.6	49
21	Intrinsic enzymatic properties modulate the self-propulsion of micromotors. <i>Nature Communications</i> , 2019, 10, 2826.	5.8	126
22	Structures of Gd <sub>3</sub> N@C <sub>80</sub> Prato Bis-Adducts: Crystal Structure, Thermal Isomerization, and Computational Study. <i>Journal of the American Chemical Society</i> , 2019, 141, 10988-10993.	6.6	16
23	Efficient reductive desymmetrization of bulky 1,3-cyclodiketones enabled by structure-guided directed evolution of a carbonyl reductase. <i>Nature Catalysis</i> , 2019, 2, 931-941.	16.1	68
24	Site-Selectivity of Prato Additions to C <sub>70</sub> : Experimental and Theoretical Studies of a New Thermodynamic Product at the [5,6]-Junction. <i>Organic Letters</i> , 2019, 21, 5162-5166.	2.4	13
25	Improved Electro- and Photocatalytic Water Reduction by Confined Cobalt Catalysts in Streptavidin. <i>ACS Catalysis</i> , 2019, 9, 5837-5846.	5.5	28
26	p38 $\beta$ is essential for cell cycle progression and liver tumorigenesis. <i>Nature</i> , 2019, 568, 557-560.	13.7	72
27	Molecular Dynamics Simulations on <i>Aspergillus niger</i> Monoamine Oxidase: Conformational Dynamics and Intermonomer Communication Essential for Its Efficient Catalysis. <i>Advanced Synthesis and Catalysis</i> , 2019, 361, 2718.	2.1	3
28	Hidden Conformations in <i>Aspergillus niger</i> Monoamine Oxidase are Key for Catalytic Efficiency. <i>Angewandte Chemie - International Edition</i> , 2019, 58, 3097-3101.	7.2	18
29	Hidden Conformations in <i>Aspergillus niger</i> Monoamine Oxidase are Key for Catalytic Efficiency. <i>Angewandte Chemie</i> , 2019, 131, 3129-3133.	1.6	0
30	Exploring the Conversion of a Sialic Acid Aldolase into a KDO Aldolase. <i>European Journal of Organic Chemistry</i> , 2018, 2018, 2603-2608.	1.2	4
31	Epoxide Hydrolase Conformational Heterogeneity for the Resolution of Bulky Pharmacologically Relevant Epoxide Substrates. <i>Chemistry - A European Journal</i> , 2018, 24, 12254-12258.	1.7	8
32	Toward Bioelectronic Nanomaterials: Photoconductivity in Protein-Porphyrin Hybrids Wrapped around SWCNT. <i>Advanced Functional Materials</i> , 2018, 28, 1704031.	7.8	25
33	Role of conformational dynamics in the evolution of novel enzyme function. <i>Chemical Communications</i> , 2018, 54, 6622-6634.	2.2	123
34	Exploring the reversal of enantioselectivity on a zinc-dependent alcohol dehydrogenase. <i>Organic and Biomolecular Chemistry</i> , 2017, 15, 4122-4129.	1.5	36
35	Rationalizing the relative abundances of trimetallic nitride template-based endohedral metallofullerenes from aromaticity measures. <i>Chemical Communications</i> , 2017, 53, 4140-4143.	2.2	5
36	Exploring the origins of selectivity in soluble epoxide hydrolase from <i>Bacillus megaterium</i> . <i>Organic and Biomolecular Chemistry</i> , 2017, 15, 8827-8835.	1.5	14

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37	The key role of aromaticity in the structure and reactivity of C <sub>60</sub> and endohedral metallofullerenes. <i>Inorganica Chimica Acta</i> , 2017, 468, 38-48.	1.2	8
38	Inducing high activity of a thermophilic enzyme at ambient temperatures by directed evolution. <i>Chemical Communications</i> , 2017, 53, 9454-9457.	2.2	41
39	Effect of incarcerated HF on the exohedral chemical reactivity of HF@C <sub>60</sub> . <i>Chemical Communications</i> , 2017, 53, 10993-10996.	2.2	26
40	Role of Conformational Dynamics in the Evolution of Retro-Aldolase Activity. <i>ACS Catalysis</i> , 2017, 7, 8524-8532.	5.5	103
41	Computational tools for the evaluation of laboratory-engineered biocatalysts. <i>Chemical Communications</i> , 2017, 53, 284-297.	2.2	84
42	Reaction Mechanism and Regioselectivity of the Bingel-Hirsch Addition of Dimethyl Bromomalonate to La@C <sub>2v</sub> -C <sub>82</sub> . <i>Chemistry - A European Journal</i> , 2016, 22, 5953-5962.	1.7	23
43	The Regioselectivity of Bingel-Hirsch Cycloadditions on Isolated Pentagon Rule Endohedral Metallofullerenes. <i>Angewandte Chemie</i> , 2016, 128, 2420-2423.	1.6	9
44	Reactivity of Single-Walled Carbon Nanotubes in the Diels-Alder Cycloaddition Reaction: Distortion-Interaction Analysis along the Reaction Pathway. <i>Chemistry - A European Journal</i> , 2016, 22, 12819-12824.	1.7	21
45	The Regioselectivity of Bingel-Hirsch Cycloadditions on Isolated Pentagon Rule Endohedral Metallofullerenes. <i>Angewandte Chemie - International Edition</i> , 2016, 55, 2374-2377.	7.2	37
46	(4 + 2) and (2 + 2) Cycloadditions of Benzyne to C <sub>60</sub> and Zig-Zag Single-Walled Carbon Nanotubes: The Effect of the Curvature. <i>Journal of Physical Chemistry C</i> , 2016, 120, 1716-1726.	1.5	34
47	Origins of stereoselectivity in evolved ketoreductases. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2015, 112, E7065-72.	3.3	104
48	Enantiospecific <i>cis</i> - <i>trans</i> Isomerization in Chiral Fulleropyrrolidines: Hydrogen-Bonding Assistance in the Carbanion Stabilization in H <sub>2</sub> O@C <sub>60</sub> . <i>Journal of the American Chemical Society</i> , 2015, 137, 1190-1197.	6.6	40
49	Extent of charge separation and exciton delocalization for electronically excited states in a triphenylamine-C <sub>60</sub> donor-acceptor conjugate: a combined molecular dynamics and TD-DFT study. <i>Theoretical Chemistry Accounts</i> , 2015, 134, 1.	0.5	13
50	Endohedral Metal-Induced Regioselective Formation of Bis-Prato Adduct of Y <sub>3</sub> N@Ih-C <sub>80</sub> and Gd <sub>3</sub> N@Ih-C <sub>80</sub> . <i>Journal of the American Chemical Society</i> , 2015, 137, 58-61.	6.6	33
51	Molecular Dynamics Explorations of Active Site Structure in Designed and Evolved Enzymes. <i>Accounts of Chemical Research</i> , 2015, 48, 1080-1089.	7.6	86
52	Bis-1,3-dipolar Cycloadditions on Endohedral Fullerenes M <sub>3</sub> N@Ih-C <sub>80</sub> (M = Sc, Lu): Remarkable Endohedral-Cluster Regiochemical Control. <i>Journal of the American Chemical Society</i> , 2015, 137, 11775-11782.	6.6	34
53	Interplay between R513 methylation and S516 phosphorylation of the cardiac voltage-gated sodium channel. <i>Amino Acids</i> , 2015, 47, 429-434.	1.2	23
54	Understanding the Exohedral Functionalization of Endohedral Metallofullerenes Metallofullerenes. <i>Carbon Materials</i> , 2015, , 67-99.	0.2	0

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55	Sponge-like molecular cage for purification of fullerenes. <i>Nature Communications</i> , 2014, 5, 5557.	5.8	162
56	The role of distant mutations and allosteric regulation on LovD active site dynamics. <i>Nature Chemical Biology</i> , 2014, 10, 431-436.	3.9	166
57	Essential Factors for Control of the Equilibrium in the Reversible Rearrangement of M <sub>3</sub> N@I <sub>h</sub> -C <sub>80</sub> Fulleropyrrolidines: Exohedral Functional Groups versus Endohedral Metal Clusters. <i>Chemistry - A European Journal</i> , 2014, 20, 14032-14039.	1.7	25
58	Why Bistetracenes Are Much Less Reactive Than Pentacenes in Diels-Alder Reactions with Fullerenes. <i>Journal of the American Chemical Society</i> , 2014, 136, 10743-10751.	6.6	52
59	Acceleration of an Aromatic Claisen Rearrangement via a Designed Spiroligozyme Catalyst that Mimics the Ketosteroid Isomerase Catalytic Dyad. <i>Journal of the American Chemical Society</i> , 2014, 136, 3817-3827.	6.6	27
60	The role of aromaticity in determining the molecular structure and reactivity of (endohedral) Tj ETQq0 0 0 rgBT /Overlock 10 Tf 50 542 T	18.7	97
61	Aromaticity as the driving force for the stability of non-IPR endohedral metallofullerene Bingel-Hirsch adducts. <i>Chemical Communications</i> , 2013, 49, 8767.	2.2	21
62	Maximum Aromaticity as a Guiding Principle for the Most Suitable Hosting Cages in Endohedral Metallofullerenes. <i>Angewandte Chemie - International Edition</i> , 2013, 52, 9275-9278.	7.2	55
63	Diels-Alder Reactions of Graphene: Computational Predictions of Products and Sites of Reaction. <i>Journal of the American Chemical Society</i> , 2013, 135, 17643-17649.	6.6	82
64	Confined organization of fullerene units along high polymer chains. <i>Journal of Materials Chemistry C</i> , 2013, 1, 5747.	2.7	16
65	A Complete Guide on the Influence of Metal Clusters in the Diels-Alder Regioselectivity of C <sub>80</sub> Endohedral Metallofullerenes. <i>Chemistry - A European Journal</i> , 2013, 19, 14931-14940.	1.7	37
66	Electrochemical control of the regioselectivity in the exohedral functionalization of C <sub>60</sub> : the role of aromaticity. <i>Chemical Communications</i> , 2013, 49, 1220.	2.2	44
67	Self-Assembled Tetragonal Prismatic Molecular Cage Highly Selective for Anionic I <sup>-</sup> Guests. <i>Chemistry - A European Journal</i> , 2013, 19, 1445-1456.	1.7	38
68	Scalable and Selective Dispersion of Semiconducting Arc-Discharged Carbon Nanotubes by Dithiafulvalene/Thiophene Copolymers for Thin Film Transistors. <i>ACS Nano</i> , 2013, 7, 2659-2668.	7.3	88
69	Covalently Patterned Graphene Surfaces by a Force-Accelerated Diels-Alder Reaction. <i>Journal of the American Chemical Society</i> , 2013, 135, 9240-9243.	6.6	121
70	Aromatic Claisen Rearrangements of O <sup>-</sup> -Prenylated Tyrosine and Model Prenyl Aryl Ethers: Computational Study of the Role of Water on Acceleration of Claisen Rearrangements. <i>European Journal of Organic Chemistry</i> , 2013, 2013, 2823-2831.	1.2	18
71	The Roles of Counterion and Water in a Stereoselective Cysteine-Catalyzed Rauhut-Currier Reaction: A Challenge for Computational Chemistry. <i>Chemistry - A European Journal</i> , 2013, 19, 14245-14253.	1.7	33
72	Editorial (Hot Topic: Nanoreactors and Molecular Prisons). <i>Current Organic Chemistry</i> , 2013, 17, 1469-1469.	0.9	0

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73	Product formation in the Prato reaction on Sc <sub>3</sub> N@D <sub>5</sub> h-C <sub>80</sub> : preference for [5,6]-bonds, and not pyracylenic bonds. <i>Chemical Communications</i> , 2012, 48, 2486.	2.2	26
74	The Frozen Cage Model: A Computationally Low-Cost Tool for Predicting the Exohedral Regioselectivity of Cycloaddition Reactions Involving Endohedral Metallofullerenes. <i>Journal of Chemical Theory and Computation</i> , 2012, 8, 1671-1683.	2.3	18
75	The Exohedral Diels-Alder Reactivity of the Titanium Carbide Endohedral Metallofullerene Ti <sub>2</sub> C <sub>2</sub> @D <sub>3</sub> h-C <sub>78</sub> : Comparison with M <sub>3</sub> N@D <sub>3</sub> h-C <sub>78</sub> and M <sub>3</sub> N@D <sub>3</sub> h-C <sub>78</sub> (M=Sc and Y) Reactivity. <i>Chemistry - A European Journal</i> , 2012, 18, 7141-7154.	1.7	54
76	Full Exploration of the Diels-Alder Cycloaddition on Metallofullerenes M <sub>3</sub> N@C <sub>80</sub> (M=Sc, Lu, Gd): The D <sub>5</sub> h versus I <sub>h</sub> Isomer and the Influence of the Metal Cluster. <i>Chemistry - A European Journal</i> , 2012, 18, 8944-8956.	1.7	49
77	The reactivity of endohedral fullerenes. What can be learnt from computational studies?. <i>Physical Chemistry Chemical Physics</i> , 2011, 13, 3585-3603.	1.3	128
78	Dispersion Corrections Essential for the Study of Chemical Reactivity in Fullerenes. <i>Journal of Physical Chemistry A</i> , 2011, 115, 3491-3496.	1.1	117
79	The Chemical Reactivity of Fullerenes and Endohedral Fullerenes: A Theoretical Perspective. <i>Carbon Materials</i> , 2011, , 57-78.	0.2	2
80	On the Mechanism of Action of Fullerene Derivatives in Superoxide Dismutation. <i>Chemistry - A European Journal</i> , 2010, 16, 3207-3214.	1.7	49
81	Reactivity and Regioselectivity of Noble Gas Endohedral Fullerenes Ng@C <sub>60</sub> and Ng <sub>2</sub> @C <sub>60</sub> (Ng=He-Xe). <i>Chemistry - A European Journal</i> , 2010, 16, 3878-3878.	1.7	6
82	Density Functional Study of the [2+2+2] Cyclootrimerization of Acetylene Catalyzed by Wilkinson's Catalyst, RhCl(PPh <sub>3</sub> ) <sub>3</sub> . <i>Organometallics</i> , 2010, 29, 562-569.	1.1	68
83	Reaction Mechanisms for Graphene and Carbon Nanotube Fluorination. <i>Journal of Physical Chemistry C</i> , 2010, 114, 3340-3345.	1.5	56
84	Rhodium(I)-Catalysed Intramolecular [2+2+2] Cyclootrimerisations of 15-, 20- and 25-Membered Azamacrocycles: Experimental and Theoretical Mechanistic Studies. <i>Chemistry - A European Journal</i> , 2009, 15, 5289-5300.	1.7	49
85	Reactivity and Regioselectivity of Noble Gas Endohedral Fullerenes Ng@C <sub>60</sub> and Ng <sub>2</sub> @C <sub>60</sub> (Ng=He-Xe). <i>Chemistry - A European Journal</i> , 2009, 15, 13111-13123.	1.7	45
86	Cycloaddition Reactions of Butadiene and 1,3-Dipoles to Curved Arenes, Fullerenes, and Nanotubes: Theoretical Evaluation of the Role of Distortion Energies on Activation Barriers. <i>Chemistry - A European Journal</i> , 2009, 15, 13219-13231.	1.7	92
87	Homolytic versus Heterolytic Dissociation of Alkali Metal Halides: The Effect of Microsolvation. <i>ChemPhysChem</i> , 2009, 10, 2955-2965.	1.0	14
88	On the Regioselective Intramolecular Nucleophilic Addition of Thiols to C <sub>60</sub> . <i>European Journal of Organic Chemistry</i> , 2009, 2009, 6231-6238.	1.2	16
89	Diels-Alder Reaction between Cyclopentadiene and C <sub>60</sub> : An Analysis of the Performance of the ONIOM Method for the Study of Chemical Reactivity in Fullerenes and Nanotubes. <i>Journal of Physical Chemistry A</i> , 2009, 113, 9721-9726.	1.1	63
90	The Diels-Alder Reaction on Endohedral Y <sub>3</sub> N@C <sub>78</sub> : The Importance of the Fullerene Strain Energy. <i>Journal of the American Chemical Society</i> , 2009, 131, 129-139.	6.6	76

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91	H-Bond-Assisted Regioselective ( <i>cis</i> -1) Intramolecular Nucleophilic Addition of the Hydroxyl Group to [60]Fullerene. <i>Journal of Organic Chemistry</i> , 2009, 74, 1480-1487.	1.7	37
92	Competitive Retro-Cycloaddition Reaction in Fullerene Dimers Connected through Pyrrolidinopyrazolino Rings. <i>Journal of Organic Chemistry</i> , 2009, 74, 8174-8180.	1.7	25
93	Regioselective Intramolecular Nucleophilic Addition of Alcohols to C <sub>60</sub> : One-Step Formation of a <i>cis</i> -1 Bicyclic-Fused Fullerene. <i>Journal of Organic Chemistry</i> , 2009, 74, 6253-6259.	1.7	33
94	Local Aromaticity of Pristine and Fluorinated Carbon Nanotubes. <i>Journal of Nanoscience and Nanotechnology</i> , 2009, 9, 6078-6083.	0.9	8
95	On the Mechanism of the Thermal Retrocycloaddition of Pyrrolidinofullerenes (Retro-Prato) <i>Tj ETQq1 1 0.784314,rgBT /Overlock 10</i>	1.7	56
96	Chemical Reactivity of D <sub>3h</sub> C <sub>78</sub> (Metallo)Fullerene: Regioselectivity Changes Induced by Sc <sub>3</sub> N Encapsulation. <i>Journal of the American Chemical Society</i> , 2008, 130, 6206-6214.	6.6	75
97	Are nucleus-independent (NICS) and <sup>1</sup> H NMR chemical shifts good indicators of aromaticity in $\pi$ -stacked polyfluorenes?. <i>Chemical Physics Letters</i> , 2006, 428, 191-195.	1.2	33