

# Albert Bruix

## List of Publications by Year in descending order

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Version: 2024-02-01

45

papers

3,668

citations

249298

26

h-index

252626

46

g-index

51

all docs

51

docs citations

51

times ranked

6702

citing authors

#	ARTICLE	IF	CITATIONS
1	Activating catalysts by adsorbate-induced reconstructions. <i>Nature Catalysis</i> , 2022, 5, 84-85.	16.1	1
2	Charting the Atomic C Interaction with Transition Metal Surfaces. <i>ACS Catalysis</i> , 2022, 12, 9256-9269.	5.5	6
3	Chemical ordering in Pt–Au, Pt–Ag and Pt–Cu nanoparticles from density functional calculations using a topological approach. <i>Materials Advances</i> , 2021, 2, 6589-6602.	2.6	12
4	Size-dependent phase stability in transition metal dichalcogenide nanoparticles controlled by metal substrates. <i>Nanoscale</i> , 2021, 13, 10167-10180.	2.8	4
5	CO oxidation activity of Pt/CeO <sub>2</sub> catalysts below 0 °C: platinum loading effects. <i>Applied Catalysis B: Environmental</i> , 2021, 286, 119931.	10.8	83
6	AgPd, AuPd, and AuPt Nanoalloys with Ag- or Au-Rich Compositions: Modeling Chemical Ordering and Optical Properties. <i>Journal of Physical Chemistry C</i> , 2021, 125, 17372-17384.	1.5	15
7	First-principles-based multiscale modelling of heterogeneous catalysis. <i>Nature Catalysis</i> , 2019, 2, 659-670.	16.1	197
8	Chemically-resolved determination of hydrogenated graphene–substrate interaction. <i>Physical Chemistry Chemical Physics</i> , 2019, 21, 13462-13466.	1.3	7
9	van der Waals exchange-correlation functionals over bulk and surface properties of transition metals. <i>Journal of Physics Condensed Matter</i> , 2019, 31, 315501.	0.7	10
10	Effects of Gas-Phase Conditions and Particle Size on the Properties of Cu(111)-Supported Zn <sub>x</sub> yO <sub>x</sub> Particles Revealed by Global Optimization and Ab Initio Thermodynamics. <i>Journal of Physical Chemistry C</i> , 2019, 123, 30903-30916.	1.5	17
11	Ab initio study of CO <sub>2</sub> hydrogenation mechanisms on inverse ZnO/Cu catalysts. <i>Journal of Catalysis</i> , 2018, 360, 168-174.	3.1	58
12	Step edge structures on the anatase TiO <sub>2</sub> (001) surface studied by atomic-resolution TEM and STM. <i>Faraday Discussions</i> , 2018, 208, 325-338.	1.6	13
13	Water Dissociation and Hydroxyl Ordering on Anatase TiO <sub>2</sub> (001) surface studied by atomic-resolution TEM and STM. Faraday Discussions, 2018, 208, 325-338. <i>Water Dissociation and Hydroxyl Ordering on Anatase TiO<sub>2</sub> (001) surface studied by atomic-resolution TEM and STM</i> xmlns:mml="http://www.w3.org/1998/Math/MathML" display="inline"><mml:mrow><mml:msub><mml:mrow><mml:mi>TiO</mml:mi></mml:mrow><mml:mrow><mml:mn>2</mml:mn></mml:mrow><mml:mrow><mml:mn>001</mml:mn></mml:mrow></mml:msub></mml:mrow><math>\text{Tj ETQq1 1 0.784314 rgBT /Overlock 10 Tf 50 252 Td (stretchy="false")</math>		

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19	Water–Gas Shift over Metal-Free Nanocrystalline Ceria: An Experimental and Theoretical Study. ChemCatChem, 2017, 9, 1373-1377.	1.8	13
20	Oxide-based nanomaterials for fuel cell catalysis: the interplay between supported single Pt atoms and particles. Catalysis Science and Technology, 2017, 7, 4315-4345.	2.1	84
21	Metal-doped ceria nanoparticles: stability and redox processes. Physical Chemistry Chemical Physics, 2017, 19, 21729-21738.	1.3	30
22	Substrate-induced semiconductor-to-metal transition in monolayer $\text{WS}_2$ . Physical Review B, 2017, 96, .		
23	Modeling Ceria-Based Nanomaterials for Catalysis and Related Applications. Catalysis Letters, 2016, 146, 2053-2080.	1.4	63
24	High efficiency of $\text{Pt}^{2+}$ - $\text{CeO}_2$ novel thin film catalyst as anode for proton exchange membrane fuel cells. Applied Catalysis B: Environmental, 2016, 197, 262-270.	10.8	52
25	Crystalline and electronic structure of single-layer $\text{TaS}_2$ . Physical Review B, 2016, 94, .		
26	Band-gap engineering by Bi intercalation of graphene on Ir(111). Physical Review B, 2016, 93, .	1.1	30
27	Single-layer $\text{MoS}_2$ on Au(111): Band gap renormalization and substrate interaction. Physical Review B, 2016, 93, .		
28	Symmetry-Driven Band Gap Engineering in Hydrogen Functionalized Graphene. ACS Nano, 2016, 10, 10798-10807.	7.3	55
29	Designing new catalysts: synthesis of new active structures: general discussion. Faraday Discussions, 2016, 188, 131-159.	1.6	4
30	Catalyst design from theory to practice: general discussion. Faraday Discussions, 2016, 188, 279-307.	1.6	2
31	Bridging model and real catalysts: general discussion. Faraday Discussions, 2016, 188, 565-589.	1.6	3
32	Effects of particle size and edge structure on the electronic structure, spectroscopic features, and chemical properties of Au(111)-supported $\text{MoS}_2$ nanoparticles. Faraday Discussions, 2016, 188, 323-343.	1.6	22
33	Towards stable single-atom catalysts: strong binding of atomically dispersed transition metals on the surface of nanostructured ceria. Catalysis Science and Technology, 2016, 6, 6806-6813.	2.1	92
34	Growth and electronic structure of epitaxial single-layer $\text{WS}_2$ on Au(111). Physical Review B, 2015, 92, .		
35	Theoretical Study of the Stoichiometric and Reduced Ce-Doped $\text{TiO}_2$ Anatase (001) Surfaces. Journal of Physical Chemistry C, 2015, 119, 4805-4816.	1.5	24
36	<i>In Situ</i> Detection of Active Edge Sites in Single-Layer $\text{MoS}_2$ Catalysts. ACS Nano, 2015, 9, 9322-9330.	7.3	144

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37	DFT Study on Ce-Doped Anatase TiO <sub>2</sub> : Nature of Ce <sup>3+</sup> and Ti <sup>3+</sup> Centers Triggered by Oxygen Vacancy Formation. <i>Journal of Physical Chemistry C</i> , 2014, 118, 9677-9689.		1.5	51
38	The Unique Properties of the Oxide-Metal Interface: Reaction of Ethanol on an Inverse Model CeO <sub>x</sub> <sup>i</sup> Au(111) Catalyst. <i>Journal of Physical Chemistry C</i> , 2014, 118, 25057-25064.		1.5	22
39	Maximum Noble Metal Efficiency in Catalytic Materials: Atomically Dispersed Surface Platinum. <i>Angewandte Chemie - International Edition</i> , 2014, 53, 10525-10530.		7.2	384
40	A New Type of Strong Metal-Support Interaction and the Production of H <sub>2</sub> through the Transformation of Water on Pt/CeO <sub>2</sub> (111) and Pt/CeO <sub>x</sub> /TiO <sub>2</sub> (110) Catalysts. <i>Journal of the American Chemical Society</i> , 2012, 134, 8968-8974.		6.6	682
41	Density functional theory model study of size and structure effects on water dissociation by platinum nanoparticles. <i>Journal of Chemical Physics</i> , 2012, 137, 034701.		1.2	56
42	Effects of deposited Pt particles on the reducibility of CeO <sub>2</sub> (111). <i>Physical Chemistry Chemical Physics</i> , 2011, 13, 11384.		1.3	89
43	Support nanostructure boosts oxygen transfer to catalytically active platinum nanoparticles. <i>Nature Materials</i> , 2011, 10, 310-315.		13.3	748
44	On the adsorption and formation of Pt dimers on the CeO <sub>2</sub> (111) surface. <i>Journal of Chemical Physics</i> , 2011, 135, 244708.		1.2	14
45	Adsorption, Oxidation State, and Diffusion of Pt Atoms on the CeO <sub>2</sub> (111) Surface. <i>Journal of Physical Chemistry C</i> , 2010, 114, 14202-14207.		1.5	71