

Guillaume Maurin

List of Publications by Year in descending order

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355
papers

31,176
citations

4831

87
h-index

6024

165
g-index

375
all docs

375
docs citations

375
times ranked

23153
citing authors

#	ARTICLE	IF	CITATIONS
1	Metal-Organic Frameworks in Biomedicine. <i>Chemical Reviews</i> , 2012, 112, 1232-1268.	23.0	3,593
2	Flexible Porous Metal-Organic Frameworks for a Controlled Drug Delivery. <i>Journal of the American Chemical Society</i> , 2008, 130, 6774-6780.	6.6	1,564
3	A New Photoactive Crystalline Highly Porous Titanium(IV) Dicarboxylate. <i>Journal of the American Chemical Society</i> , 2009, 131, 10857-10859.	6.6	1,127
4	Gas/vapour separation using ultra-microporous metal-organic frameworks: insights into the structure/separation relationship. <i>Chemical Society Reviews</i> , 2017, 46, 3402-3430.	18.7	1,033
5	The new age of MOFs and of their porous-related solids. <i>Chemical Society Reviews</i> , 2017, 46, 3104-3107.	18.7	623
6	Why hybrid porous solids capture greenhouse gases?. <i>Chemical Society Reviews</i> , 2011, 40, 550-562.	18.7	603
7	An Explanation for the Very Large Breathing Effect of a Metal-Organic Framework during CO ₂ Adsorption. <i>Advanced Materials</i> , 2007, 19, 2246-2251.	11.1	501
8	A pressure-amplifying framework material with negative gas adsorption transitions. <i>Nature</i> , 2016, 532, 348-352.	13.7	490
9	Functionalization in Flexible Porous Solids: Effects on the Pore Opening and the Host-Guest Interactions. <i>Journal of the American Chemical Society</i> , 2010, 132, 1127-1136.	6.6	445
10	Rationale of Drug Encapsulation and Release from Biocompatible Porous Metal-Organic Frameworks. <i>Chemistry of Materials</i> , 2013, 25, 2767-2776.	3.2	412
11	A Series of Isorecticular, Highly Stable, Porous Zirconium Oxide Based Metal-Organic Frameworks. <i>Angewandte Chemie - International Edition</i> , 2012, 51, 9267-9271.	7.2	407
12	Co-adsorption and Separation of CO ₂ and CH ₄ Mixtures in the Highly Flexible MIL-53(Cr) MOF. <i>Journal of the American Chemical Society</i> , 2009, 131, 17490-17499.	6.6	398
13	How Linker's Modification Controls Swelling Properties of Highly Flexible Iron(III) Dicarboxylates MIL-88. <i>Journal of the American Chemical Society</i> , 2011, 133, 17839-17847.	6.6	383
14	Functionalizing porous zirconium terephthalate UiO-66(Zr) for natural gas upgrading: a computational exploration. <i>Chemical Communications</i> , 2011, 47, 9603.	2.2	345
15	A Water Stable Metal-Organic Framework with Optimal Features for CO ₂ Capture. <i>Angewandte Chemie - International Edition</i> , 2013, 52, 10316-10320.	7.2	303
16	Molecular Dynamics Simulations of Breathing MOFs: Structural Transformations of MIL-53(Cr) upon Thermal Activation and CO ₂ Adsorption. <i>Angewandte Chemie - International Edition</i> , 2008, 47, 8487-8491.	7.2	302
17	Fluorinated MOF platform for selective removal and sensing of SO ₂ from flue gas and air. <i>Nature Communications</i> , 2019, 10, 1328.	5.8	292
18	Influence of framework metal ions on the dye capture behavior of MIL-100 (Fe, Cr) MOF type solids. <i>Journal of Materials Chemistry A</i> , 2013, 1, 8534.	5.2	291

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19	Complex Adsorption of Short Linear Alkanes in the Flexible Metal-Organic-Framework MIL-53(Fe). <i>Journal of the American Chemical Society</i> , 2009, 131, 13002-13008.	6.6	281
20	Fast Variability of Tera-Electron Volt Å Rays from the Radio Galaxy M87. <i>Science</i> , 2006, 314, 1424-1427.	6.0	277
21	Hydrolytically stable fluorinated metal-organic frameworks for energy-efficient dehydration. <i>Science</i> , 2017, 356, 731-735.	6.0	275
22	Design of Hydrophilic Metal Organic Framework Water Adsorbents for Heat Reallocation. <i>Advanced Materials</i> , 2015, 27, 4775-4780.	11.1	253
23	Prediction of the Conditions for Breathing of Metal Organic Framework Materials Using a Combination of X-ray Powder Diffraction, Microcalorimetry, and Molecular Simulation. <i>Journal of the American Chemical Society</i> , 2008, 130, 12808-12814.	6.6	246
24	Multistep N ₂ Breathing in the Metal-Organic Framework Co(1,4-benzenedipyrazolate). <i>Journal of the American Chemical Society</i> , 2010, 132, 13782-13788.	6.6	220
25	A robust zirconium amino acid metal-organic framework for proton conduction. <i>Nature Communications</i> , 2018, 9, 4937.	5.8	218
26	A robust large-pore zirconium carboxylate metal-organic framework for energy-efficient water-sorption-driven refrigeration. <i>Nature Energy</i> , 2018, 3, 985-993.	19.8	217
27	Natural gas upgrading using a fluorinated MOF with tuned H ₂ S and CO ₂ adsorption selectivity. <i>Nature Energy</i> , 2018, 3, 1059-1066.	19.8	214
28	The Structure of the Aluminum Fumarate Metal-Organic Framework A520. <i>Angewandte Chemie - International Edition</i> , 2015, 54, 3664-3668.	7.2	206
29	Selective nitrogen capture by porous hybrid materials containing accessible transition metal ion sites. <i>Nature Materials</i> , 2017, 16, 526-531.	13.3	201
30	Structure and Dynamics of the Functionalized MOF Type UiO-66(Zr): NMR and Dielectric Relaxation Spectroscopies Coupled with DFT Calculations. <i>Chemistry of Materials</i> , 2012, 24, 2168-2177.	3.2	200
31	An Evaluation of UiO-66 for Gas-Based Applications. <i>Chemistry - an Asian Journal</i> , 2011, 6, 3270-3280.	1.7	192
32	Synthesis, Structure, Characterization, and Redox Properties of the Porous MIL-68(Fe) Solid. <i>European Journal of Inorganic Chemistry</i> , 2010, 2010, 3789-3794.	1.0	191
33	Adsorption Mechanism of Carbon Dioxide in Faujasites: Grand Canonical Monte Carlo Simulations and Microcalorimetry Measurements. <i>Journal of Physical Chemistry B</i> , 2005, 109, 16084-16091.	1.2	185
34	Explanation of the Adsorption of Polar Vapors in the Highly Flexible Metal Organic Framework MIL-53(Cr). <i>Journal of the American Chemical Society</i> , 2010, 132, 9488-9498.	6.6	185
35	CH ₄ storage and CO ₂ capture in highly porous zirconium oxide based metal-organic frameworks. <i>Chemical Communications</i> , 2012, 48, 9831.	2.2	180
36	Large breathing of the MOF MIL-47(VIV) under mechanical pressure: a joint experimental-modelling exploration. <i>Chemical Science</i> , 2012, 3, 1100.	3.7	176

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37	Acid-functionalized UiO-66(Zr) MOFs and their evolution after intra-framework cross-linking: structural features and sorption properties. <i>Journal of Materials Chemistry A</i> , 2015, 3, 3294-3309.	5.2	174
38	Understanding the Thermodynamic and Kinetic Behavior of the CO ₂ /CH ₄ Gas Mixture within the Porous Zirconium Terephthalate UiO-66(Zr): A Joint Experimental and Modeling Approach. <i>Journal of Physical Chemistry C</i> , 2011, 115, 13768-13774.	1.5	166
39	Asymmetric pore windows in MOF membranes for natural gas valorization. <i>Nature</i> , 2022, 606, 706-712.	13.7	163
40	Rational design of mixed-matrix metal-organic framework membranes for molecular separations. <i>Science</i> , 2022, 376, 1080-1087.	6.0	160
41	Electrochemical intercalation of lithium into multiwall carbon nanotubes. <i>Chemical Physics Letters</i> , 1999, 312, 14-18.	1.2	159
42	Molecular Insight into the Adsorption of H ₂ S in the Flexible MIL-53(Cr) and Rigid MIL-47(V) MOFs: Infrared Spectroscopy Combined to Molecular Simulations. <i>Journal of Physical Chemistry C</i> , 2011, 115, 2047-2056.	1.5	157
43	Quasi-Elastic Neutron Scattering and Molecular Dynamics Study of Methane Diffusion in Metal Organic Frameworks MIL-47(V) and MIL-53(Cr). <i>Angewandte Chemie - International Edition</i> , 2008, 47, 6611-6615.	7.2	154
44	Understanding the origins of metal-organic framework/polymer compatibility. <i>Chemical Science</i> , 2018, 9, 315-324.	3.7	153
45	Proton Transport in a Highly Conductive Porous Zirconium-Based Metal-Organic Framework: Molecular Insight. <i>Angewandte Chemie - International Edition</i> , 2016, 55, 3919-3924.	7.2	152
46	A metal-organic framework for efficient water-based ultra-low-temperature-driven cooling. <i>Nature Communications</i> , 2019, 10, 3025.	5.8	145
47	On the breathing effect of a metal-organic framework upon CO ₂ adsorption: Monte Carlo compared to microcalorimetry experiments. <i>Chemical Communications</i> , 2007, , 3261.	2.2	137
48	Probing the Adsorption Sites for CO ₂ in Metal Organic Frameworks Materials MIL-53 (Al), <i>TJ ETQq0 0 0 rgBT /Overlock 10 T</i>	1.5	137
49	Probing the Dynamics of CO ₂ and CH ₄ within the Porous Zirconium Terephthalate UiO-66(Zr): A Synergic Combination of Neutron Scattering Measurements and Molecular Simulations. <i>Chemistry - A European Journal</i> , 2011, 17, 8882-8889.	1.7	137
50	A robust amino-functionalized titanium(IV) based MOF for improved separation of acid gases. <i>Chemical Communications</i> , 2013, 49, 10082.	2.2	135
51	Solvent-Induced Control over Breathing Behavior in Flexible Metal-Organic Frameworks for Natural Gas Delivery. <i>Angewandte Chemie - International Edition</i> , 2019, 58, 8073-8077.	7.2	132
52	Towards an Improved anti-HIV Activity of NRTI via Metal-Organic Frameworks Nanoparticles. <i>Advanced Healthcare Materials</i> , 2013, 2, 1630-1637.	3.9	130
53	Step-wise dealumination of natural clinoptilolite: Structural and physicochemical characterization. <i>Microporous and Mesoporous Materials</i> , 2010, 135, 187-196.	2.2	129
54	Microscopic Model of the Metal-Organic Framework/Polymer Interface: A First Step toward Understanding the Compatibility in Mixed Matrix Membranes. <i>ACS Applied Materials & Interfaces</i> , 2016, 8, 809-819.	4.0	129

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55	Molecular Insight into the Adsorption and Diffusion of Water in the Versatile Hydrophilic/Hydrophobic Flexible MIL-53(Cr) MOF. <i>Journal of Physical Chemistry C</i> , 2011, 115, 10764-10776.	1.5	128
56	A phase transformable ultrastable titanium-carboxylate framework for photoconduction. <i>Nature Communications</i> , 2018, 9, 1660.	5.8	128
57	Enhanced gas separation performance of 6FDA-DAM based mixed matrix membranes by incorporating MOF UiO-66 and its derivatives. <i>Journal of Membrane Science</i> , 2018, 558, 64-77.	4.1	126
58	Effect of the organic functionalization of flexible MOFs on the adsorption of CO ₂ . <i>Journal of Materials Chemistry</i> , 2012, 22, 10266.	6.7	125
59	Gas adsorption microcalorimetry and modelling to characterise zeolites and related materials. <i>Comptes Rendus Chimie</i> , 2005, 8, 283-302.	0.2	124
60	Probing the adsorption performance of the hybrid porous MIL-68(Al): a synergic combination of experimental and modelling tools. <i>Journal of Materials Chemistry</i> , 2012, 22, 10210.	6.7	124
61	Adsorption of CO ₂ in metal organic frameworks of different metal centres: Grand Canonical Monte Carlo simulations compared to experiments. <i>Adsorption</i> , 2007, 13, 461-467.	1.4	123
62	Metal-organic frameworks as potential shock absorbers: the case of the highly flexible MIL-53(Al). <i>Chemical Communications</i> , 2014, 50, 9462-9464.	2.2	122
63	Rationalization of the entrapping of bioactive molecules into a series of functionalized porous zirconium terephthalate MOFs. <i>Journal of Materials Chemistry B</i> , 2013, 1, 1101.	2.9	118
64	Charge distribution in metal organic framework materials: transferability to a preliminary molecular simulation study of the CO ₂ adsorption in the MIL-53 (Al) system. <i>Physical Chemistry Chemical Physics</i> , 2007, 9, 1059-1063.	1.3	112
65	Hybrid Monte Carlo Simulations Combined with a Phase Mixture Model to Predict the Structural Transitions of a Porous Metal-Organic Framework Material upon Adsorption of Guest Molecules. <i>Journal of Physical Chemistry C</i> , 2010, 114, 6496-6502.	1.5	112
66	Revealing the Structure-Property Relationships of Metal-Organic Frameworks for CO ₂ Capture from Flue Gas. <i>Langmuir</i> , 2012, 28, 12094-12099.	1.6	110
67	Transport Diffusivity of CO ₂ in the Highly Flexible Metal-Organic Framework MIL-53(Cr). <i>Angewandte Chemie - International Edition</i> , 2009, 48, 8335-8339.	7.2	109
68	Self and Transport Diffusivity of CO ₂ in the Metal-Organic Framework MIL-47(V) Explored by Quasi-elastic Neutron Scattering Experiments and Molecular Dynamics Simulations. <i>ACS Nano</i> , 2010, 4, 143-152.	7.3	109
69	Thermodynamic insight into stimuli-responsive behaviour of soft porous crystals. <i>Nature Communications</i> , 2018, 9, 204.	5.8	104
70	Mechanical energy storage performance of an aluminum fumarate metal-organic framework. <i>Chemical Science</i> , 2016, 7, 446-450.	3.7	103
71	Achieving Superprotonic Conduction with a 2D Fluorinated Metal-Organic Framework. <i>Journal of the American Chemical Society</i> , 2018, 140, 13156-13160.	6.6	103
72	Electrically Induced Breathing of the MIL-53(Cr) Metal-Organic Framework. <i>ACS Central Science</i> , 2017, 3, 394-398.	5.3	102

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73	Polymer Infiltration into Metal-Organic Frameworks in Mixed-Matrix Membranes Detected in Situ by NMR. <i>Journal of the American Chemical Society</i> , 2019, 141, 7589-7595.	6.6	102
74	Thermo-Responsive MOF/Polymer Composites for Temperature-Mediated Water Capture and Release. <i>Angewandte Chemie - International Edition</i> , 2020, 59, 11003-11009.	7.2	101
75	Experimental Evidence Supported by Simulations of a Very High H_2 Diffusion in Metal Organic Framework Materials. <i>Physical Review Letters</i> , 2008, 100, 245901.	2.9	99
76	A biocompatible porous Mg-gallate metal-organic framework as an antioxidant carrier. <i>Chemical Communications</i> , 2015, 51, 5848-5851.	2.2	98
77	Mixed-Linker Hybrid Superpolyhedra for the Production of a Series of Large-Pore Iron(III) Carboxylate Metal-Organic Frameworks. <i>Angewandte Chemie - International Edition</i> , 2013, 52, 5056-5060.	7.2	97
78	Thermodynamic Insight in the High-Pressure Behavior of UiO-66: Effect of Linker Defects and Linker Expansion. <i>Chemistry of Materials</i> , 2016, 28, 5721-5732.	3.2	97
79	Porous metal organic framework nanoparticles to address the challenges related to busulfan encapsulation. <i>Nanomedicine</i> , 2011, 6, 1683-1695.	1.7	95
80	Metal-organic frameworks to satisfy gas upgrading demands: fine-tuning the MOF platform for the operative removal of H_2S . <i>Journal of Materials Chemistry A</i> , 2017, 5, 3293-3303.	5.2	94
81	Extended and functionalized porous iron(III) tri- or dicarboxylates with MIL-100/101 topologies. <i>Chemical Communications</i> , 2014, 50, 6872.	2.2	93
82	Toward an Understanding of the Microstructure and Interfacial Properties of PIMs/ZIF-8 Mixed Matrix Membranes. <i>ACS Applied Materials & Interfaces</i> , 2016, 8, 27311-27321.	4.0	93
83	The effect of crystallite size on pressure amplification in switchable porous solids. <i>Nature Communications</i> , 2018, 9, 1573.	5.8	92
84	Comparative Guest, Thermal, and Mechanical Breathing of the Porous Metal Organic Framework MIL-53(Cr): A Computational Exploration Supported by Experiments. <i>Journal of Physical Chemistry C</i> , 2012, 116, 13289-13295.	1.5	90
85	Adsorption and Diffusion of H_2 in the MOF Type Systems MIL-47(V) and MIL-53(Cr): A Combination of Microcalorimetry and QENS Experiments with Molecular Simulations. <i>Journal of Physical Chemistry C</i> , 2009, 113, 7802-7812.	1.5	89
86	Nano-mechanical cutting and opening of single wall carbon nanotubes. <i>Chemical Physics Letters</i> , 2000, 331, 125-131.	1.2	88
87	Tailoring the separation properties of flexible metal-organic frameworks using mechanical pressure. <i>Nature Communications</i> , 2020, 11, 1216.	5.8	88
88	Observations of the Sagittarius dwarf galaxy by the HESS experiment and search for a dark matter signal. <i>Astroparticle Physics</i> , 2008, 29, 55-62.	1.9	87
89	Adsorption of CO ₂ , CH ₄ and their binary mixture in Faujasite NaY: A combination of molecular simulations with gravimetry-manometry and microcalorimetry measurements. <i>Microporous and Mesoporous Materials</i> , 2009, 119, 117-128.	2.2	86
90	Adsorption and Diffusion of Light Hydrocarbons in UiO-66(Zr): A Combination of Experimental and Modeling Tools. <i>Journal of Physical Chemistry C</i> , 2014, 118, 27470-27482.	1.5	84

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91	Influence of Extra-Framework Cations on the Adsorption Properties of X-Faujasite Systems: \hat{A} Microcalorimetry and Molecular Simulations. <i>Journal of Physical Chemistry B</i> , 2005, 109, 125-129.	1.2	83
92	A Comparison of Barostats for the Mechanical Characterization of Metal-Organic Frameworks. <i>Journal of Chemical Theory and Computation</i> , 2015, 11, 5583-5597.	2.3	83
93	A Fine-Tuned MOF for Gas and Vapor Separation: A Multipurpose Adsorbent for Acid Gas Removal, Dehydration, and BTX Sieving. <i>CheM</i> , 2017, 3, 822-833.	5.8	83
94	Segmented and opened multi-walled carbon nanotubes. <i>Carbon</i> , 2001, 39, 1273-1278.	5.4	82
95	Adsorption of light hydrocarbons in the flexible MIL-53(Cr) and rigid MIL-47(V) metal-organic frameworks: a combination of molecular simulations and microcalorimetry/gravimetry measurements. <i>Physical Chemistry Chemical Physics</i> , 2010, 12, 6428.	1.3	82
96	MIL-91(Ti), a small pore metal-organic framework which fulfils several criteria: an upscaled green synthesis, excellent water stability, high CO ₂ selectivity and fast CO ₂ transport. <i>Journal of Materials Chemistry A</i> , 2016, 4, 1383-1389.	5.2	82
97	A Complete Separation of Hexane Isomers by a Functionalized Flexible Metal Organic Framework. <i>Advanced Functional Materials</i> , 2014, 24, 7666-7673.	7.8	81
98	Highly reversible sorption of H ₂ S and CO ₂ by an environmentally friendly Mg-based MOF. <i>Journal of Materials Chemistry A</i> , 2018, 6, 16900-16909.	5.2	81
99	Separation of CO ₂ -CH ₄ mixtures in the mesoporous MIL-100(Cr) MOF: experimental and modelling approaches. <i>Dalton Transactions</i> , 2012, 41, 4052.	1.6	78
100	Revisiting the Aluminum Trimesate-Based MOF (MIL-96): From Structure Determination to the Processing of Mixed Matrix Membranes for CO ₂ Capture. <i>Chemistry of Materials</i> , 2017, 29, 10326-10338.	3.2	78
101	A promising metal-organic framework (MOF), MIL-96(Al), for CO ₂ separation under humid conditions. <i>Journal of Materials Chemistry A</i> , 2018, 6, 2081-2090.	5.2	78
102	A quantitative structure activity relationship approach to probe the influence of the functionalization on the drug encapsulation of porous metal-organic frameworks. <i>Microporous and Mesoporous Materials</i> , 2012, 157, 124-130.	2.2	76
103	Conformation-Controlled Sorption Properties and Breathing of the Aliphatic Al-MOF [Al(OH)(CDC)]. <i>Inorganic Chemistry</i> , 2014, 53, 4610-4620.	1.9	74
104	Electrochemical lithium intercalation into multiwall carbon nanotubes: a micro-Raman study. <i>Solid State Ionics</i> , 2000, 136-137, 1295-1299.	1.3	73
105	Series of Porous 3-D Coordination Polymers Based on Iron(III) and Porphyrin Derivatives. <i>Chemistry of Materials</i> , 2011, 23, 4641-4651.	3.2	73
106	Towards general network architecture design criteria for negative gas adsorption transitions in ultraporous frameworks. <i>Nature Communications</i> , 2019, 10, 3632.	5.8	73
107	Caffeine Confinement into a Series of Functionalized Porous Zirconium MOFs: A Joint Experimental/Modeling Exploration. <i>Journal of Physical Chemistry C</i> , 2013, 117, 11694-11704.	1.5	70
108	High and energy-efficient reversible SO ₂ uptake by a robust Sc-based MOF. <i>Journal of Materials Chemistry A</i> , 2019, 7, 15580-15584.	5.2	70

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109	Impact of phosphorylation on the encapsulation of nucleoside analogues within porous iron(III) metal-organic framework MIL-100(Fe) nanoparticles. <i>Journal of Materials Chemistry B</i> , 2013, 1, 4231.	2.9	69
110	Computational exploration of a Zr-carboxylate based metal-organic framework as a membrane material for CO ₂ capture. <i>Journal of Materials Chemistry A</i> , 2014, 2, 1657-1661.	5.2	68
111	Rational design of a robust aluminum metal-organic framework for multi-purpose water-sorption-driven heat allocations. <i>Nature Communications</i> , 2020, 11, 5112.	5.8	68
112	A Water Stable Metal-Organic Framework with Optimal Features for CO ₂ Capture. <i>Angewandte Chemie</i> , 2013, 125, 10506-10510.	1.6	66
113	Modeling of Gas Transport through Polymer/MOF Interfaces: A Microsecond-Scale Concentration Gradient-Driven Molecular Dynamics Study. <i>Chemistry of Materials</i> , 2020, 32, 1288-1296.	3.2	64
114	Structural Origin of Unusual CO ₂ Adsorption Behavior of a Small-Pore Aluminum Bisphosphonate MOF. <i>Journal of Physical Chemistry C</i> , 2015, 119, 4208-4216.	1.5	63
115	Outstanding reversible H ₂ S capture by an Al(III)-based MOF. <i>Chemical Communications</i> , 2019, 55, 3049-3052.	2.2	63
116	CO ₂ adsorption in alkali cation exchanged Y faujasites: A quantum chemical study compared to experiments. <i>Chemical Physics Letters</i> , 2006, 426, 387-392.	1.2	62
117	Design of MoS ₂ /Graphene van der Waals Heterostructure as Highly Efficient and Stable Electrocatalyst for Hydrogen Evolution in Acidic and Alkaline Media. <i>ACS Applied Materials & Interfaces</i> , 2020, 12, 24777-24785.	4.0	62
118	Tuning Cellular Biological Functions Through the Controlled Release of NO from a Porous Ti-MOF. <i>Angewandte Chemie - International Edition</i> , 2020, 59, 5135-5143.	7.2	62
119	SPECTRAL ANALYSIS AND INTERPRETATION OF THE ¹³ C-RAY EMISSION FROM THE STARBURST GALAXY NGC 253. <i>Astrophysical Journal</i> , 2012, 757, 158.	1.6	61
120	Impact of the Metal Centre and Functionalization on the Mechanical Behaviour of MIL-53 Metal-Organic Frameworks. <i>European Journal of Inorganic Chemistry</i> , 2016, 2016, 4424-4429.	1.0	60
121	Covalent and Selective Grafting of Polyethylene Glycol Brushes at the Surface of ZIF-8 for the Processing of Membranes for Pervaporation. <i>ACS Sustainable Chemistry and Engineering</i> , 2019, 7, 6629-6639.	3.2	60
122	Highly Selective Removal of Perfluorinated Contaminants by Adsorption on Al-Silica Zeolite Beta. <i>Angewandte Chemie - International Edition</i> , 2020, 59, 14086-14090.	7.2	60
123	Modeling the Effect of Hydration in Zeolite Na-Mordenite. <i>Journal of Physical Chemistry B</i> , 2004, 108, 3739-3745.	1.2	59
124	Physics Behind the Guest-Assisted Structural Transitions of a Porous Metal-Organic Framework Material. <i>Journal of Physical Chemistry Letters</i> , 2010, 1, 2810-2815.	2.1	59
125	Multivariable Sieving and Hierarchical Recognition for Organic Toxics in Nonhomogeneous Channel of MOFs. <i>CheM</i> , 2019, 5, 1337-1350.	5.8	59
126	Molecular Dynamics Simulation of the Cation Motion upon Adsorption of CO ₂ in Faujasite Zeolite Systems. <i>Journal of Physical Chemistry B</i> , 2006, 110, 14372-14378.	1.2	58

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127	Investigation of CO ₂ adsorption in Faujasite systems: Grand Canonical Monte Carlo and molecular dynamics simulations based on a new derived Na ⁺ -CO ₂ force field. <i>Microporous and Mesoporous Materials</i> , 2007, 99, 70-78.	2.2	58
128	A new aluminium-based microporous metal-organic framework: Al(BTB) (BTB =) Tj ETQq0 0 0 rgBT /Overlock 10 Tf,50 702 Td (1,3,5-b	2.2	58
129	Rietveld Refinement of MIL-160 and Its Structural Flexibility Upon H ₂ O and N ₂ Adsorption. <i>European Journal of Inorganic Chemistry</i> , 2018, 2018, 3626-3632.	1.0	58
130	Toward a Rational Design of Titanium Metal-Organic Frameworks. <i>Matter</i> , 2020, 2, 440-450.	5.0	58
131	Enhanced Polymer Crystallinity in Mixed-Matrix Membranes Induced by Metal-Organic Framework Nanosheets for Efficient CO ₂ Capture. <i>ACS Applied Materials & Interfaces</i> , 2018, 10, 43095-43103.	4.0	55
132	Diffusion of Methanol in Zeolite NaY: A Molecular Dynamics Study. <i>Journal of Physical Chemistry B</i> , 2007, 111, 2836-2844.	1.2	54
133	Estimation of the breathing energy of flexible MOFs by combining TGA and DSC techniques. <i>Chemical Communications</i> , 2009, , 2733.	2.2	54
134	Structure and properties of Al-MIL-53-ADP, a breathing MOF based on the aliphatic linker molecule adipic acid. <i>Dalton Transactions</i> , 2016, 45, 4179-4186.	1.6	54
135	Engineering Structural Dynamics of Zirconium Metal-Organic Frameworks Based on Natural C ₄ Linkers. <i>Journal of the American Chemical Society</i> , 2019, 141, 17207-17216.	6.6	54
136	Exploration of the mechanical behavior of metal organic frameworks UiO-66(Zr) and MIL-125(Ti) and their NH ₂ -functionalized versions. <i>Dalton Transactions</i> , 2016, 45, 4283-4288.	1.6	53
137	Adsorption Contraction Mechanics: Understanding Breathing Energetics in Isoreticular Metal-Organic Frameworks. <i>Journal of Physical Chemistry C</i> , 2018, 122, 19171-19179.	1.5	52
138	Diffusion of Binary CO ₂ /CH ₄ Mixtures in the MIL-47(V) and MIL-53(Cr) Metal-Organic Framework Type Solids: A Combination of Neutron Scattering Measurements and Molecular Dynamics Simulations. <i>Journal of Physical Chemistry C</i> , 2013, 117, 11275-11284.	1.5	51
139	Gas Adsorption and Separation by the Al-Based Metal-Organic Framework MIL-160. <i>Journal of Physical Chemistry C</i> , 2017, 121, 26822-26832.	1.5	51
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