Carla Grijó Fonseca

List of Publications by Year in descending order

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Version: 2024-02-01

1163117 1372567 10 117 8 10 citations g-index h-index papers 10 10 10 129 docs citations times ranked citing authors all docs

#	Article	IF	CITATIONS
1	Structural and thermodynamic investigation of the hydration-dehydration process of Na+-Montmorillonite using DFTÂcalculations. Applied Clay Science, 2017, 143, 212-219.	5.2	24
2	Na+ as a probe to structural investigation of dehydrated smectites using NMR spectra calculated by DFT. Applied Clay Science, 2016, 126, 132-140.	5.2	16
3	Investigation of the initial stages of the montmorillonite acid-activation process using DFT calculations. Applied Clay Science, 2018, 165, 170-178.	5.2	16
4	Nb2O5 supported on mixed oxides catalyzed oxidative and photochemical conversion of anilines to azoxybenzenes. New Journal of Chemistry, 2019, 43, 5863-5871.	2.8	12
5	Experimental and theoretical study of deactivated HDT catalysts by Si species deposited on their surfaces: Models proposition, structural and thermodynamic analysis. Journal of Catalysis, 2020, 389, 578-591.	6.2	12
6	Structural analysis of dehydrated gibbsite-based layered double hydroxides Li–Al–X (X = Fâ^', Clâ^', Brâ^',) Tj The role of Zn 2+ dopants in the acid-basic catalysis on MgO(001) surface: Ab initio simulations of the	2.0	11
7	dissociative chemisorption of R-O-ka€² andAk-S-ka€² (k, <mmi:math) 0.784314="" 1="" 10="" etqq1="" ii<="" ij="" overlock="" rgbt="" td=""><td>1.9</td><td>d (xmins:mml</td></mmi:math)>	1.9	d (xmins:mml
8	DFT calculations for structural prediction and applications of intercalated lamellar compounds. Dalton Transactions, 2018, 47, 2852-2866.	3.3	10
9	DFT simulations of dissociative chemisorption of sulfur-, nitrogen- and oxygen-containing molecules on mixed oxides with Mg, Al and Zn. Applied Surface Science, 2019, 484, 524-533.	6.1	5
10	On the mixed oxides-supported niobium catalyst towards benzylamine oxidation. Catalysis Today, 2021, 381, 118-125.	4.4	1