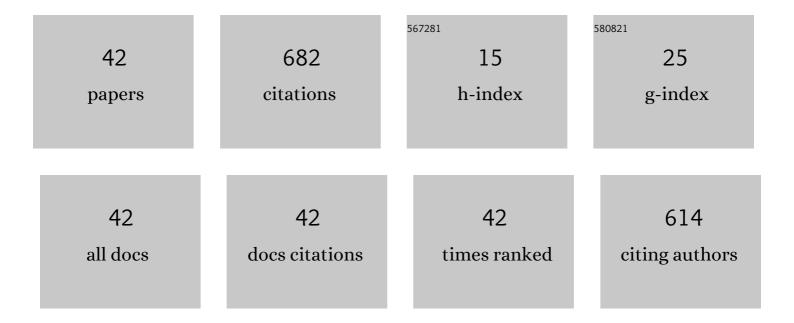
Carla Isabel Costa Pinheiro

List of Publications by Year in descending order

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#	Article	IF	CITATIONS
1	Solid–gas reactors driven by concentrated solar energy with potential application to calcium looping: A comparative review. Renewable and Sustainable Energy Reviews, 2022, 158, 112048.	16.4	31
2	Improved performance of modified CaO-Al2O3 based pellets for CO2 capture under realistic Ca-looping conditions. Journal of CO2 Utilization, 2022, 61, 102007.	6.8	13
3	Hydrogen Production with In Situ CO2 Capture at High and Medium Temperatures Using Solid Sorbents. Energies, 2022, 15, 4039.	3.1	15
4	Blending Wastes of Marble Powder and Dolomite Sorbents for Calcium-Looping CO2 Capture under Realistic Industrial Calcination Conditions. Materials, 2021, 14, 4379.	2.9	13
5	Enhancement of sintering resistance of CaO-based sorbents using industrial waste resources for Ca-looping in the cement industry. Separation and Purification Technology, 2020, 235, 116190.	7.9	23
6	Modeling the deactivation of CaO-based sorbents during multiple Ca-looping cycles for CO2 post-combustion capture. Computers and Chemical Engineering, 2020, 134, 106679.	3.8	16
7	Modelling Full Cycles of Carbonation-Calcination for Calcium Looping Process Simulation. Computer Aided Chemical Engineering, 2019, , 1009-1014.	0.5	1
8	Tailoring Synthetic Sol–Gel CaO Sorbents with High Reactivity or High Stability for Ca-Looping CO ₂ Capture. Industrial & Engineering Chemistry Research, 2019, 58, 8484-8494.	3.7	24
9	Multi-scale modelling and simulation of Ca-looping cycle process for CO2 post-combustion capture. Computer Aided Chemical Engineering, 2018, , 291-292.	0.5	0
10	Economic comparison of a reactive distillation-based process with the conventional process for the production of ethyl tert-butyl ether (ETBE). Computers and Chemical Engineering, 2017, 100, 9-26.	3.8	16
11	Development of Soft Sensors Based on Analytical and Spectral Data on a Real Small Size Wastewater Treatment Plant. Lecture Notes in Electrical Engineering, 2017, , 323-333.	0.4	1
12	CAPE in the Chemical Engineering Master's Integrated Programme at IST-ULisboa. Computer Aided Chemical Engineering, 2017, , 2959-2964.	0.5	0
13	Computer Aided Control Projects as Main Assessment Component of Master's Advanced Control Courses. Computer Aided Chemical Engineering, 2017, 40, 2953-2958.	0.5	0
14	A Multivariable Model Predictive Control Project in a Computer Aided Master's Degree Course. Computer Aided Chemical Engineering, 2016, 38, 871-876.	0.5	1
15	Waste Marble Powders as Promising Inexpensive Natural CaO-Based Sorbents for Post-Combustion CO ₂ Capture. Industrial & amp; Engineering Chemistry Research, 2016, 55, 7860-7872.	3.7	37
16	Model Based Fault Diagnosis for Performance Control of a Decentralized Wastewater Treatment Plant. Computer Aided Chemical Engineering, 2014, 33, 691-696.	0.5	2
17	Optimal design of reactive distillation systems: Application to the production of ethyl tert-butyl ether (ETBE). Computers and Chemical Engineering, 2014, 64, 81-94.	3.8	22
18	Fluid Catalytic Cracking (FCC) Process Modeling, Simulation, and Control. Industrial & Engineering Chemistry Research, 2012, 51, 1-29.	3.7	123

#	Article	IF	CITATIONS
19	Estimation of catalyst deactivation parameters of ethyl tert-butyl ether (ETBE) reactors based on industrial plant data. Computer Aided Chemical Engineering, 2012, , 1002-1006.	0.5	2
20	Model Development and Validation of Ethyl <i>tert</i> Butyl Ether Production Reactors Using Industrial Plant Data. Industrial & Engineering Chemistry Research, 2012, 51, 15018-15031.	3.7	12
21	Investigation of a stable synthetic sol–gel CaO sorbent for CO2 capture. Fuel, 2012, 94, 624-628.	6.4	94
22	Influence of different catalyst deactivation models in a validated simulator of an industrial UOP FCC unit with high-efficiency regenerator. Fuel, 2012, 97, 97-108.	6.4	23
23	Methodologies for input-output data exchange between LabVIEW® and MATLAB®/Simulink®software for Real Time Control of a Pilot Scale Distillation Process. Computer Aided Chemical Engineering, 2011, 29, 708-712.	0.5	4
24	Model Development and Validation of an Industrial UOP Fluid Catalytic Cracking Unit with a High-Efficiency Regenerator. Industrial & Engineering Chemistry Research, 2008, 47, 850-866.	3.7	29
25	Closed-loop dynamic behavior of an industrial high-efficiency FCC unit. , 2008, , .		2
26	Dynamic modelling of an industrial R2R FCC unit. Chemical Engineering Science, 2007, 62, 1184-1198.	3.8	45
27	Steady state multiplicity in an UOP FCC unit with high-efficiency regenerator. Chemical Engineering Science, 2007, 62, 6308-6322.	3.8	28
28	Multiplicity of steady states in an UOP FCC unit with high efficiency regenerator. Computer Aided Chemical Engineering, 2006, , 1575-1580.	0.5	3
29	Determination of state-space model uncertainty using bootstrap techniques. Journal of Process Control, 2006, 16, 685-692.	3.3	4
30	Mechanistic dynamic modelling of an industrial FCC Unit. Computer Aided Chemical Engineering, 2005, , 589-594.	0.5	5
31	Transient microkinetic modelling of n-heptane catalytic cracking over H-USY zeolite. Chemical Engineering Science, 2004, 59, 1221-1232.	3.8	15
32	Subspace identification methods for a fast dynamic model structure screening. Computer Aided Chemical Engineering, 2004, 18, 697-702.	0.5	0
33	Title is missing!. Reaction Kinetics and Catalysis Letters, 2000, 69, 39-46.	0.6	1
34	Model predictive control of reactor temperature in a CSTR pilot plant operating at an unstable steady-state. Computers and Chemical Engineering, 1999, 23, S859-S862.	3.8	18
35	Dynamic modelling and network simulation of n-heptane catalytic cracking: influence of kinetic parameters. Chemical Engineering Science, 1999, 54, 1735-1750.	3.8	10
36	Network simulation of catalytic cracking reactions. Studies in Surface Science and Catalysis, 1997, , 529-534.	1.5	2

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37	Reactions of mixtures of light alkenes and n-heptane over USHY zeolite. Applied Catalysis A: General, 1994, 108, 107-114.	4.3	0
38	Temperature Programmed Desorption of Oxygen on Bismuth Molybdates and Reactivity For Olefin Oxidation. Studies in Surface Science and Catalysis, 1992, 72, 325-334.	1.5	1
39	Syntheses, electrochemistry, and bonding of bis(cyclopentadienyl)molybdenum alkyl complexes. Molecular structure of Mo(.eta.5-C5H5)2(C4H9)2. Thermochemistry of Mo(.eta.5-C5H5)2R2 and Mo(.eta.5-C5H5)2L (R = CH3, C2H5, C4H9; L = ethylene, diphenylacetylene). Organometallics, 1991, 10, 483-494.	2.3	23
40	A Comparison Between Low and High Temperature Bi2O3MoO3 Phases FOR 1-Butene Reactions. Studies in Surface Science and Catalysis, 1991, 67, 77-85.	1.5	5
41	New molybdenocene dihydrocarbyls. Journal of Organometallic Chemistry, 1987, 327, C59-C62.	1.8	4

42 Pelasgian fountains: learning Greek in the early Middle Ages. , 0, , 65-82.