

Roda Bounaceur

List of Publications by Year in descending order

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docs citations

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times ranked

3277
citing authors

| # | ARTICLE | IF | CITATIONS |
|----|--|-----|-----------|
| 1 | Development of a Detailed Kinetic Model for the Oxidation of <i>n</i> -Butane in the Liquid Phase. Journal of Physical Chemistry B, 2021, 125, 6955-6967. | 2.6 | 8 |
| 2 | The Deep Composition of Uranus and Neptune from In Situ Exploration and Thermochemical Modeling. Space Science Reviews, 2020, 216, 1. | 8.1 | 16 |
| 3 | New chemical scheme for giant planet thermochemistry. Astronomy and Astrophysics, 2020, 634, A78. | 5.1 | 34 |
| 4 | Influence of H ₂ S on the thermal cracking of alkylbenzenes at high pressure (70â€‰MPa) and moderate temperature (583â€‰623â€‰K). Journal of Analytical and Applied Pyrolysis, 2019, 140, 423-433. | 5.5 | 6 |
| 5 | Establishing the Maximum Carbon Number for Reliable Quantitative Gas Chromatographic Analysis of Heavy Ends Hydrocarbons. Part 3. Coupled Pyrolysis-GC Modeling. Energy & Fuels, 2019, 33, 2045-2056. | 5.1 | 1 |
| 6 | Reduced chemical scheme for modelling warm to hot hydrogen-dominated atmospheres. Astronomy and Astrophysics, 2019, 624, A58. | 5.1 | 26 |
| 7 | Thermal cracking of <i>n</i> -butylbenzene at high pressure: Experimental study and kinetic modelling. Journal of Analytical and Applied Pyrolysis, 2018, 133, 234-245. | 5.5 | 13 |
| 8 | Kinetic parameters for the thermal cracking of simple hydrocarbons: From laboratory to geological time-temperature conditions. Journal of Analytical and Applied Pyrolysis, 2017, 125, 40-49. | 5.5 | 15 |
| 9 | Kinetic modeling of 1-methylnaphthalene pyrolysis at high pressure (100 bar). Journal of Analytical and Applied Pyrolysis, 2017, 124, 542-562. | 5.5 | 13 |
| 10 | Thermal cracking of 1- <i>n</i> -butyldecalin at high pressure (100 bar). Journal of Analytical and Applied Pyrolysis, 2017, 123, 204-215. | 5.5 | 3 |
| 11 | Evaluating the effects of CO ₂ capture benchmarks on efficiency and costs of membrane systems for post-combustion capture: A parametric simulation study. International Journal of Greenhouse Gas Control, 2017, 63, 449-461. | 4.6 | 23 |
| 12 | Rigorous variable permeability modelling and process simulation for the design of polymeric membrane gas separation units: MEMSIC simulation tool. Journal of Membrane Science, 2017, 523, 77-91. | 8.2 | 50 |
| 13 | Prediction of Auto-Ignition Temperatures and Delays for Gas Turbine Applications. Journal of Engineering for Gas Turbines and Power, 2016, 138, . | 1.1 | 17 |
| 14 | Development of a Model for Auto-Ignition Delays and its Use for the Prediction of Premix Combustion Reliability. , 2016, , . | | 0 |
| 15 | Interplay of inlet temperature and humidity on energy penalty for CO ₂ post-combustion capture: Rigorous analysis and simulation of a single stage gas permeation process. Energy, 2016, 116, 517-525. | 8.8 | 4 |
| 16 | Data supporting the validation of a simulation model for multi-component gas separation in polymeric membranes. Data in Brief, 2016, 9, 776-780. | 1.0 | 2 |
| 17 | Influence of pressure (100 Paâ€‰100 Mpa) on the pyrolysis of an alkane at moderate temperature (603) Tj ETQq1 1 0.784314 rgBT /Ove 442-451. | 5.5 | 8 |
| 18 | Thermal cracking of <i>n</i> -butylcyclohexane at high pressure (100 bar)â€”Part 2: Mechanistic modeling. Journal of Analytical and Applied Pyrolysis, 2016, 120, 174-185. | 5.5 | 16 |

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|----|--|------|-----------|
| 19 | Intramolecular effects on the kinetics of unimolecular reactions of \hat{I}^2 -HOROÖ™ and HOQÖ™OOH radicals. Physical Chemistry Chemical Physics, 2016, 18, 12231-12251. | 2.8 | 57 |
| 20 | Thermal cracking of n -butylcyclohexane at high pressure (100 bar)â€”Part 1: Experimental study. Journal of Analytical and Applied Pyrolysis, 2016, 117, 1-16. | 5.5 | 19 |
| 21 | Pyrolysis of n- octane at very low concentration and low temperature. Journal of Analytical and Applied Pyrolysis, 2016, 117, 282-289. | 5.5 | 5 |
| 22 | Prediction of Auto-Ignition Temperatures and Delays for Gas Turbine Applications. , 2015, , . | | 0 |
| 23 | New chemical scheme for studying carbon-rich exoplanet atmospheres. Astronomy and Astrophysics, 2015, 577, A33. | 5.1 | 64 |
| 24 | Pushing the limits of intensified CO2 post-combustion capture by gasâ€”liquid absorption through a membrane contactor. Chemical Engineering and Processing: Process Intensification, 2015, 91, 7-22. | 3.6 | 28 |
| 25 | Experimental and modeling study of burning velocities for alkyl aromatic components relevant to diesel fuels. Proceedings of the Combustion Institute, 2015, 35, 341-348. | 3.9 | 43 |
| 26 | Determining predictive uncertainties and global sensitivities for large parameter systems: A case study for n -butane oxidation. Proceedings of the Combustion Institute, 2015, 35, 607-616. | 3.9 | 31 |
| 27 | The oxidation of large alkylbenzenes: An experimental and modeling study. Proceedings of the Combustion Institute, 2015, 35, 349-356. | 3.9 | 34 |
| 28 | Measurements of Laminar Burning Velocities above Atmospheric Pressure Using the Heat Flux Methodâ€”Application to the Case of n -Pentane. Energy & Fuels, 2015, 29, 398-404. | 5.1 | 49 |
| 29 | Experimental and Modeling Investigation of the Low-Temperature Oxidation of Dimethyl Ether. Journal of Physical Chemistry A, 2015, 119, 7905-7923. | 2.5 | 85 |
| 30 | Kinetic modelling of char gasification by accounting for the evolution of the reactive surface area. Chemical Engineering Science, 2015, 138, 281-290. | 3.8 | 25 |
| 31 | Combustion and Oxidation Kinetics of Alternative Gas Turbines Fuels. , 2014, , . | | 6 |
| 32 | Laminar burning velocity of gasolines with addition of ethanol. Fuel, 2014, 115, 162-169. | 6.4 | 248 |
| 33 | Kinetic study of abatement of low concentration of dibenzofuran by oxidation â€” Effects of co-reactants. Chemical Engineering Journal, 2013, 218, 154-163. | 12.7 | 8 |
| 34 | Mechanistic modeling of the thermal cracking of methylcyclohexane near atmospheric pressure, from 523 to 1273K: Identification of aromatization pathways. Journal of Analytical and Applied Pyrolysis, 2013, 103, 240-254. | 5.5 | 27 |
| 35 | Establishing the Maximum Carbon Number for Reliable Quantitative Gas Chromatographic Analysis of Heavy Ends Hydrocarbons. Part 2. Migration and Separation Gas Chromatography Modeling. Energy & Fuels, 2013, 27, 2336-2350. | 5.1 | 6 |
| 36 | A chemical model for the atmosphere of hot Jupiters. Astronomy and Astrophysics, 2012, 546, A43. | 5.1 | 181 |

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|----|--|------|-----------|
| 37 | Membrane contactors for intensified post combustion carbon dioxide capture by gas-liquid absorption in MEA: A parametric study. Chemical Engineering Research and Design, 2012, 90, 2325-2337. | 5.6 | 14 |
| 38 | Establishing the Maximum Carbon Number for Reliable Quantitative Gas Chromatographic Analysis of Heavy Ends Hydrocarbons. Part 1: Low-Conversion Thermal Cracking Modeling. Energy & Fuels, 2012, 26, 2600-2610. | 5.1 | 8 |
| 39 | JTHERGAS: Thermodynamic estimation from 2D graphical representations of molecules. Energy, 2012, 43, 161-171. | 8.8 | 12 |
| 40 | Experimental and modeling study of the oxidation of n-butylbenzene. Combustion and Flame, 2012, 159, 1399-1416. | 5.2 | 59 |
| 41 | Evaluating the intensification potential of membrane contactors for gas absorption in a chemical solvent: A generic one-dimensional methodology and its application to CO ₂ absorption in monoethanolamine. Journal of Membrane Science, 2012, 389, 1-16. | 8.2 | 25 |
| 42 | Towards cleaner combustion engines through groundbreaking detailed chemical kinetic models. Chemical Society Reviews, 2011, 40, 4762. | 38.1 | 111 |
| 43 | Measurements of Laminar Flame Velocity for Components of Natural Gas. Energy & Fuels, 2011, 25, 3875-3884. | 5.1 | 181 |
| 44 | Thermal evolution of n- and iso-alkanes in oils. Part 1: Pyrolysis model for a mixture of 78 alkanes (C ₁ -C ₃₂) including 13,206 free radical reactions. Organic Geochemistry, 2011, 42, 439-450. | 1.8 | 27 |
| 45 | DME as a Potential Alternative Fuel for Gas Turbines: A Numerical Approach to Combustion and Oxidation Kinetics. , 2011, , . | | 7 |
| 46 | Laminar Flame Velocity of Components of Natural Gas. , 2011, , . | | 2 |
| 47 | Modelling intra-particle phenomena of biomass pyrolysis. Chemical Engineering Research and Design, 2011, 89, 2136-2146. | 5.6 | 56 |
| 48 | Oxidation of small unsaturated methyl and ethyl esters. International Journal of Chemical Kinetics, 2011, 43, 204-218. | 1.6 | 32 |
| 49 | Thermal and Kinetic Impact of CO, CO ₂ , and H ₂ O on the Postoxidation of IC-Engine Exhaust Gases. Combustion Science and Technology, 2010, 182, 39-59. | 2.3 | 23 |
| 50 | Ethanol as an Alternative Fuel in Gas Turbines: Combustion and Oxidation Kinetics. , 2010, , . | | 5 |
| 51 | An extended mechanism including high pressure conditions (700bar) for toluene pyrolysis. Journal of Analytical and Applied Pyrolysis, 2010, 87, 236-247. | 5.5 | 21 |
| 52 | Adiabatic flame temperature from biofuels and fossil fuels and derived effect on NO _x emissions. Fuel Processing Technology, 2010, 91, 229-235. | 7.2 | 104 |
| 53 | An experimental and kinetic modeling study of the autoignition of 1-methylnaphthalene/air and 1-methylnaphthalene/n-decane/air mixtures at elevated pressures. Combustion and Flame, 2010, 157, 1976-1988. | 5.2 | 67 |
| 54 | Modelling of Weak Acid Conversion in an EDI Cell. Separation Science and Technology, 2010, 45, 1015-1024. | 2.5 | 4 |

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|----|---|-----|-----------|
| 55 | Reassessment of the Kinetic Influence of Toluene on n -Alkane Pyrolysis. Energy & Fuels, 2010, 24, 3817-3830. | 5.1 | 19 |
| 56 | New insights into secondary gas generation from the thermal cracking of oil: Methylated monoaromatics. A kinetic approach using 1,2,4-trimethylbenzene. Part I: A mechanistic kinetic model. Organic Geochemistry, 2010, 41, 146-167. | 1.8 | 57 |
| 57 | Gas Turbines and Biodiesel: A Clarification of the Relative NOX Indices of FAME, Gasoil, and Natural Gas. , 2009, , . | | 2 |
| 58 | Influence of EGR compounds on the oxidation of an HCCI-diesel surrogate. Proceedings of the Combustion Institute, 2009, 32, 2851-2859. | 3.9 | 31 |
| 59 | Biogas, membranes and carbon dioxide capture. Journal of Membrane Science, 2009, 328, 11-14. | 8.2 | 106 |
| 60 | Influence of the position of the double bond on the autoignition of linear alkenes at low temperature. Proceedings of the Combustion Institute, 2009, 32, 387-394. | 3.9 | 66 |
| 61 | A hybrid process combining oxygen enriched air combustion and membrane separation for post-combustion carbon dioxide capture. Separation and Purification Technology, 2009, 68, 30-36. | 7.9 | 70 |
| 62 | Modeling of autoignition and NO sensitization for the oxidation of IC engine surrogate fuels. Combustion and Flame, 2009, 156, 505-521. | 5.2 | 61 |
| 63 | CO_2/N_2 Reverse Selective Gas Separation Membranes: Technological Opportunities and Scientific Challenges. Industrial & Engineering Chemistry Research, 2009, 48, 3700-3701. | 3.7 | 21 |
| 64 | Kinetic modelling of a surrogate diesel fuel applied to 3D auto-ignition in HCCI engines. International Journal of Vehicle Design, 2007, 44, 124. | 0.3 | 28 |
| 65 | CHEMICAL IMPACT OF CO AND H_2 ADDITION ON THE AUTO-IGNITION DELAY OF HOMOGENEOUS N-HEPTANE/AIR MIXTURES. Combustion Science and Technology, 2007, 179, 1937-1962. | 2.3 | 27 |
| 66 | Primary Mechanism of the Thermal Decomposition of Tricyclodecane. Journal of Physical Chemistry A, 2006, 110, 11298-11314. | 2.5 | 82 |
| 67 | Membrane processes for post-combustion carbon dioxide capture: A parametric study. Energy, 2006, 31, 2556-2570. | 8.8 | 260 |
| 68 | Experimental and modeling study of the oxidation of xylenes. International Journal of Chemical Kinetics, 2006, 38, 284-302. | 1.6 | 88 |
| 69 | Progress toward a unified detailed kinetic model for the autoignition of alkanes from C4 to C10 between 600 and 1200 K. Combustion and Flame, 2005, 142, 170-186. | 5.2 | 272 |
| 70 | Experimental and modeling study of the oxidation of toluene. International Journal of Chemical Kinetics, 2005, 37, 25-49. | 1.6 | 176 |
| 71 | Experimental Study and Modeling of the Role of Hydronaphthalenics on the Thermal Stability of Hydrocarbons under Laboratory and Geological Conditions. Industrial & Engineering Chemistry Research, 2005, 44, 8972-8987. | 3.7 | 17 |
| 72 | Experimental and modelling investigation of the thermal decomposition of n-dodecane. Journal of Analytical and Applied Pyrolysis, 2004, 71, 865-881. | 5.5 | 118 |

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|----|--|-----|-----------|
| 73 | Contribution to Scramjet Active Cooling Analysis Using N-dodecane Decomposition Model. , 2003, , . | | 16 |
| 74 | Inhibiting Effect of Tetralin on the Pyrolytic Decomposition of Hexadecane. Comparison with Toluene. Industrial & Engineering Chemistry Research, 2002, 41, 4689-4701. | 3.7 | 31 |
| 75 | Up to what temperature is petroleum stable? New insights from a 5200 free radical reactions model. Organic Geochemistry, 2002, 33, 1487-1499. | 1.8 | 54 |
| 76 | Modeling of hydrocarbons pyrolysis at low temperature. Automatic generation of free radicals mechanisms. Journal of Analytical and Applied Pyrolysis, 2002, 64, 103-122. | 5.5 | 51 |
| 77 | Mechanistic Modeling of the Thermal Cracking of Tetralin. Industrial & Engineering Chemistry Research, 2000, 39, 4152-4165. | 3.7 | 42 |
| 78 | Modeling the Laminar Flame Speed of Natural Gas and Gasoline Surrogates. , 0, , . | | 13 |
| 79 | Reactive Transport and Its Implications on Heavy Oil HTGC Analysis. A Coupled Thermo-Hydro-Chemical (THC) Multiphysics Modelling Approach. , 0, , . | | 0 |