



POLITECNICO DI TORINO
Repository ISTITUZIONALE

Contribution of different turbulent scales in turbulent reacting flows: science or just a computational exercise?

Original

Contribution of different turbulent scales in turbulent reacting flows: science or just a computational exercise? / Vanni M.; Valerio S.; Pipino, M.; Barresi, A.A.; Baldi, G.. - STAMPA. - (1994), p. 450 [#185i]. ((Intervento presentato al convegno 1994 Annual AIChE Meeting tenutosi a San Francisco (CA, USA) nel 13-18 November 1994.

Availability:

This version is available at: 11583/2647805 since: 2016-09-09T20:55:53Z

Publisher:

American Institution of Chemical Engineers

Published

DOI:

Terms of use:

openAccess

This article is made available under terms and conditions as specified in the corresponding bibliographic description in the repository

Publisher copyright

(Article begins on next page)

EXTENDED ABSTRACTS

AMERICAN INSTITUTE OF CHEMICAL ENGINEERS 1994 ANNUAL MEETING

SAN FRANCISCO HILTON AND TOWERS HOTEL
SAN FRANCISCO, CALIFORNIA

NOVEMBER 13-18, 1994

**CONTRIBUTION OF DIFFERENT TURBULENT SCALES IN TURBULENT
REACTING FLOWS: SCIENCE OR JUST A COMPUTATIONAL EXERCISE?**

M. Vanni, S. Valerio, M. Pipino, A. A. Barresi, G. Baldi

450, paper #185i

COMPLETE TITLE OF PAPER

CONTRIBUTION OF DIFFERENT TURBULENT SCALES IN TURBULENT REACTING FLOWS: SCIENCE OR JUST A COMPUTATIONAL EXERCISE?

SPEAKER AND AUTHORS, WITH AFFILIATIONS AND COMPLETE MAILING ADDRESSES

M. VANNI, S. VALERIO, M. PIPINO, A. A. BARRESI, G. BALDI

Dipartimento di Scienza dei Materiali ed Ingegneria Chimica - Politecnico di Torino

Corso Duca degli Abruzzi 24, 10124 Torino, Italy

ABSTRACT

A turbulent reactive process is a function of two different simultaneous phenomena: the chemical reaction and the transport process; the rates of them can be evaluated by means of the characteristic times of the phenomena. The definition of a characteristic mixing time is difficult because turbulence is a phenomenon that occurs on a large range of scales from that of the characteristic length of the reactor to that of the dissipation range. In the literature a great number of models have been proposed to correlate the mixing time to a specific region of the spectrum. The models based on the Kolmogorov time scale assume that the main contribute to mixing is given by the high wavenumber region of the spectrum; others consider the contribution of the whole range of wavenumbers, thus assuming the scale of the energy containing eddies as the characteristic one for the mixing process.

To validate one of these models the authors are performing a simulation of the local concentration field generated by a fast chemical reaction in a tubular reactor with two coaxial flows in which unpremixed reactants are fed (non-homogeneous turbulence). The neutralization of NaOH with HCl is considered. The numerical results are compared with the experimental data relative to the axial concentration profiles of the key reactant, determined by spectrophotometrical analysis with a fiber-optic probe. By choosing properly the mixing constants it is possible to fit experimental data. A good agreement can be obtained not only taking into account the combined effects of both the previously defined scales, but also using a macro-mixing scale, related to the characteristic size of the reactor.

But the choice of the fitting constants, functions of reactor geometry, operative conditions, etc., to describe the experimental data, is really science or only empiricism? Is it a scientific work or only a numerical calculation exercise?