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Oxy-fuel combustion of pulverized fuels: Combustion fundamentals and modeling

Chungen Yin^{a,*}, Jinyue Yan^{b,c}

^a Department of Energy Technology, Aalborg University, 9220 Aalborg East, Denmark

^b School of Chemical Science and Engineering, Royal Institute of Technology, Sweden

^c School of Sustainable Development of Society and Technology, Mälardalen University, Sweden

HIGHLIGHTS

• The fundamentals underpinning oxy-fuel combustion development thoroughly reviewed.

• Oxy-fuel induced changes in combustion physics, chemistry and modeling explained.

• Generic modeling strategies for PF oxy-fuel combustion successfully proposed.

• Oxy-fuel based power generation and CCS systems and the key issues discussed.

• Research needs in oxy-fuel combustion fundamentals and their modeling identified.

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ABSTRACT

Oxy-fuel combustion of pulverized fuels (PF), as a promising technology for CO₂ capture from power plants, has gained a lot of concerns and also advanced considerable research, development and demonstration in the past years worldwide. The use of CO_2 or the mixture of CO_2 and H_2O vapor as the diluent in oxy-fuel combustion, instead of N₂ in conventional air-fuel combustion, induces significant changes to the combustion fundamentals, because of the great differences in the physical properties and chemical effects of the different diluents. Therefore, some fundamental issues and technological challenges need to be properly addressed to develop oxy-fuel combustion into an enabled technology. Computational Fluid Dynamics (CFD) modeling, which has been proven to be a very useful and cost-effective tool in research and development of conventional air-fuel combustion, is expected to play a similarly vital role in future development of oxy-fuel combustion technology. The paper presents a state-of-the-art review and an in-depth discussion of PF oxy-fuel combustion fundamentals and their modeling, which underpin the development of this promising technology. The focus is placed on the key issues in combustion physics (e.g., turbulent gas-solid flow, heat and mass transfer) and combustion chemistry (e.g., pyrolysis, gas phase combustion and char reactions), mainly on how they are affected in oxy-fuel conditions and how they are modeled and implemented into CFD simulations. The system performance of PF oxy-fuel combustion is also reviewed. Finally, the current status of PF oxy-fuel combustion fundamentals and modeling is concluded and the research needs in these regards are suggested.

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Review



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^{*} Corresponding author. Tel.: +45 30622577; fax: +45 98151411. *E-mail address:* chy@et.aau.dk (C. Yin).

Nomenclature

| Α | frequency factor in rate coefficient in Arrhenius form (s^{-1}) | U _{VM,C} | fraction of unburnt combustibles, $U_{VM,C} = \frac{\text{mass of volatiles and char in particle}}{\text{mass of volatiles of host in for each particle}} (-)$ |
|------------------------|---|------------------------|---|
| ABET | particle BET surface area (m ² /kg) | V | narticle velocity (m/s) |
| A_n | particle projected area (m ²) | V | cell volume (m^3) |
| A_{ni} | projected area of group <i>i</i> particles (m ²) | V | particle volume (m^3) |
| $A_{ns}^{P,}$ | particle surface area (m ²) | v p X | mole fraction (_) |
| beii | emissivity gas temperature polynomial coefficients in | л V | mole fraction of NO () |
| 0,1,1 | WSGGM (-) | л _{NO} V | more fraction of nitrogen in char () |
| Cc | concentration of PF particles (kg/m^3) | I N,char | mass machon of minogen in chai (-) |
| C _D | drag coefficient (–) | | |
| C_{n} | specific heat (I/(kg K)) | Greek let | ters 1 |
| d_n | particle size (m) | α | local gas absorption coefficient (m ⁻¹) |
| dm _p | conversion rate of particle in different sub-processes | $\alpha_1, \ \alpha_2$ | two yield factors (–) |
| at | (kg/s) | α_p | equivalent particle absorption coefficient (m ⁻¹) |
| D | pipe diameter (m) | ΔH | heat effects (J/kg) |
| D_{σ} | mass diffusivity (m^2/s) | 3 | total emissivity of local gas mixture (–) |
| E | activation energy in rate coefficient in Arrhenius form | Ер | particle emissivity (–) |
| 2 | (I/kmol) | $\varepsilon_{p,i}$ | emissivity of group <i>i</i> particles (–) |
| f.,; | scattering factor of group <i>i</i> particles $(-)$ | η | conversion factor (–) |
| f_{mo} | initial moisture fraction (–) | θ_R | radiation temperature (K) |
| J W,U g g | gravitational acceleration (m/s^2) | μ_{g} | air or gas dynamic viscosity (kg/(m s)) |
| в, в hм | convective mass transfer coefficients (m/s) | μ_t | turbulent viscosity (kg/(m s)) |
| h_{τ} | convective heat transfer coefficients (M/s) | $ ho_{	extsf{g}}$ | air or gas density (kg/m³) |
| $I(\vec{r} \ \hat{s})$ | radiative intensity at position \vec{r} in direction \hat{s} (W/ | $ ho_p$ | particle density (kg/m ³) |
| 1(1,0) | $(m^2 sr))$ | σ | Stefan–Boltzmann constant (5.67×10^{-6}) (W/(m ² K ⁴)) |
| k | kinetic rate (s^{-1}) | σ_p | equivalent particle scattering coefficient (m ⁻¹) |
| k _a | gas thermal conductivity (W/(mK)) | τ_v | particle momentum response time, $\tau_v = \rho_p d_p^2 / (18 \mu_g)$ |
| k; | absorption coefficient of <i>i</i> -th grav gas in WSGGM (1/ | | (m) |
| | (atm m)) | ϕ | phase function (–) |
| L | domain-based beam length (m) | Ω, Ω' | solid angle (sr) |
| Lc | characteristic length (m) | | |
| m_a | particle ash content (kg) | Abbreviat | tions |
| m_p | particle mass (kg) | CCS | carbon capture and storage |
| $m_{p,0}$ | initial particle mass at injection (kg) | CFD | Computational Fluid Dynamics |
| $m_{\nu}(t)$ | mass of volatile yield up to time t (kg) | CPD | Chemical Percolation Devolatilization |
| MW _N | molecular weight of N (kg/kmol) | CTF | combustion test facility |
| MW _{NO} | molecular weight of NO (kg/kmol) | DO | discrete ordinates (radiation model) |
| n _i | number density of group <i>i</i> particles $(1/m^3)$ | DTF | drop tube furnace |
| Nu | Nusselt number (–) | DTR | drop tube reactor |
| Р | sum of partial pressures of the participating gases (atm) | DTRM | discrete transfer radiation model |
| Patm | local gas pressure (atm) | EBU | eddy-breakup |
| $P_{\rm pa}$ | local gas pressure (Pa) | ED | Eddy Dissipation |
| Pr | Prandtl number (–) | EDC | Eddy Dissipation Concept |
| q_r | radiative flux (W/m ²) | EFR | entrained flow reactor |
| R | conversion rate (s ⁻¹) | EWBM | exponential wide band model |
| Re | Reynolds number (–) | FG-DVC | Functional Group – Depolymerisation Vaporisation |
| Rep | particle Reynolds number, $Re_p = \rho_\sigma \mathbf{u} - \mathbf{v} d_p / \mu_\sigma$ (-) | | Cross-linking |
| R_u | universal gas constant (8315) (J/(kmol K)) | FR/ED | Finite Rate/Eddy Dissipation |
| S | path length (m) | FSK | full spectrum k-distribution |
| Sc | char burnout rate (kg/s) | JL 4-step | Jones and Lindstedt 4-step |
| Sc | Schmidt number (–) | LES | large eddy simulation |
| Sh | Sherwood number (–) | PF | pulverized fuel |
| S _{NO} | NO source term (kg/(m ³ s)) | KANS | Reynolds-Averaged Navier-Stokes |
| t | time (s) | KFG | recycled flue gas |
| Т | temperature (K) | KIL | radiative transfer equation |
| T_g | local gas temperature (K) | IGA | thermogravimetric analysis |
| $T_{p,i}$ | temperature of group i particles (K) | UDF | user-defined function |
| u | rac volocity(m/s) | VIVI | volatile matters |
| | gas velocity (III/S) | | an Westhread and Driver 2 star |
| u_{\min} | the minimum (or saltation) velocity (m/s) | WD 2-ste | ep Westbrook and Dryer 2-step |

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