

Spray Combustion Modeling Including Detailed Chemistry

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- I. Why Detailed Chemisty?
- II. Structures of Spray Flames in the Counterflow Configuration
- III. Turbulent Spray-Flame Modeling
- IV. Summary and Conclusions

Modeling of Technical Spray Flames





Why Detailed Chemistry?

Detailed Chemical Reaction Mechanisms are Available for a Considerate Number of Relevant Combustion Systems (Alcanes, Alcohols, Hydrogen/Air, Hydrogen/Oxygen, ...)

- Combustion of liquid fuel sprays in air (e.g. internal engine combustion, industrial furnaces, gas turbine combustors)
- Liquid oxygen in (gaseous) hydrogen (liquid rocket propulsion)
- Liquid oxygen in gaseous hydrocarbons or alcohols (green propellants)

Advantages of Using Detailed Chemistry:

- Mechanism is independent of the experimental configuration, it depends only on pressure (not for hydrogen/air or hydrogen/oxygen)
- Mechanism is the base for development of reduced mechanisms (both manually or automatically developed systems)
- Prediction of pollutants and precursors of soot formation

Disadvantages of Using Detailed Chemistry:

- Stiffness of the conservation equations
- Consume a considerable amount of computer time

Applications:

Laminar Flames: Detailed mechanisms can be implemented directly for hydrogen and small hydrocarbons and alcohols Turbulent Flames: Detailed chemistry may be implemented through use of the flamelet model



Modeling of Laminar Spray Flames in the Counterflow Configuration



Motivation:

Investigation of laminar spray flame structures using detailed models for instance for chemical reactions

Flamelet modeling of turbulent spray diffusion flames

Properties:

- Planar or axisymmetric
- Two-dimensional
- Strained



n-Heptane/Air Spray Flame at Atmospheric Pressure



a = 500/s

Detailed Chemistry:

Solid Lines, Square

One-Step Chemistry:

Dashed Lines, Triangles

⇒ One-Step Chemistry is not Suitable to Correctly Predict Even the Outer Flame Structure

Gutheil, E., Sirignano, W. A.: *Counterflow Spray Combustion Modeling Including Detailed Transport and Detailed Chemistry*, Combustion and Flame: 113(2), 92-105 (1998).



Mathematical Model

Gas-phase with dilute spray

- Boundary layer approximation, low Mach number
- Dimensionless, steady equations
- ▷ Similarity transformation \Rightarrow 2D \rightarrow 1D equations
- ➤ Ideal gas law
- Detailed chemical reaction mechanisms.
 - > H₂/O₂ (8 species and 38 elementary reactions)
 - methanol/air (23 species and 170 elementary reactions)
- Detailed transport: molecular diffusion and thermo diffusion
- ➤ Gas-phase properties between 300 and 5000 K from NASA polynomials
- Physical properties of H₂ and O₂ in the range of 80 to 300 K and 1 to 200 bar from *JSME* tables



Mathematical Model

Liquid phase

- > Mono-, bi- and polydisperse sprays, single-component sprays
- Discrete droplet model
- Spherically symmetric droplets
- Convective droplet model for heating and vaporization (Abramzon-Sirignano model)
- Pressure and temperature dependent heat of vaporization
- > Assumption of thermodynamic equilibrium:
 - Ambrose's equation for the evaluation of the vapor pressure for methanol/air
 - Calculation of binary H₂/O₂ mixtures to obtain the gas mixture composition at the interface (replacement of Raoult's law)
- Droplet motion (drag)



Yang, V., Lin, N.N., Shuen, J.S., Combust. Sci. and Tech., 97: pp. 247-270, 1994

JSME Data Book, Thermophysical Properties of Fluids, 1983.

Prausnitz, J.M., Lichtenthaler, R.N., de Avezedo, E.G., *Molecular Thermo-dynamics of Fluid-Phase Equilibria*, Prentice-Hall, New Jersey, 1986. Litchford, R.R., Jeng, S.M., *AIAA Paper* 90-2191, 1990.

H₂/Air Spray Flame at Atmospheric Pressure

p = 1 bar, $T_{+} = T_{-} = 300$ K, a = 100/s

4000

3500

3000

2500

2000

1500 L____ 10²

10³

10⁴

10⁵

Strain Rate [1/s]

10⁶

10⁷

10⁸

Symbols: Experiment: T. M. Brown *et. al., Combust. Sci. and Tech.*, Vol. 129, pp. 71-88, 1997.



• Experiment: Sohn, C. H., Chung, S. H., Lee, S. R., Kim, J. S., *Combustion and Flame*, 115 (3): 299-312, 1998.



Schlotz, D., Gutheil, E.: *Modeling of Laminar Mono- and Bidisperse Liquid Oxygen/Hydrogen Spray Flames in the Counterflow Configuration*, Combustion Science and Technology, 158, 195-210 (2000).



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Multiple Structures of Spray Flames

Methanol/Air Spray Flame at Atmospheric Pressure

a = 100/s



Multiple Structures of Spray Flames

Methanol/Air Spray Flame at Atmospheric Pressure

a = 300/s





Methanol/Air Spray Flame at Atmospheric Pressure

a= 500/s



Multiple Structures of Spray Flames

Methanol/Air Spray Flame at Atmospheric Pressure

a = 300/s



Comparison: Gas-Sided Flame and Pure Gas Flames

Methanol/Air Spray Flame at Atmospheric Pressure

a = 300/s

Gas Side (Spray Flame)

Gas Flame





Methanol/Air Spray Flame at Atmospheric Pressure

Comparison of Spray and Gas Flame



Structures of Laminar Spray Flames in the Counterflow Configuration

- The LOX/H₂ Spray Flames are very stable and persist to strain rates of 25,000/s. The non-monotonicity of the gaseous oxygen profile on the spray side stems from the competition of vaporization and combustion.
- Multiple structures of methanol/air spray flames have been found for strain rates up to 400/s. The inner structure of the gas-sided flame is the same as a pure gas flamelet with appropriate initial conditions.
- At high strain, the gas-sided flame is extinguished and the spray-sided flame moves towards the gas-side of the counterflow configuration.
 - Question: How does the finding affect models such as the flamelet model for turbulent spray diffusion flames?



Flamelet-Model for Turbulent Diffusion Flames

Turbulent Flame



Library of laminar flame structures in the counterflow configuration



- Gas flames
 - Strain rate
- Spray flames
 - Strain rate
 - Droplet size
 - Droplet velocity
 - Equivalence ratio

Laminar Spray Flame Structures for Use in Flamelet Models for

Turbulent Spray Diffusion Flames (Methanol/Air)



Experiment: McDonell V.G., Samuelsen, G.S., *UCI-Laboratory Report* UCI-ARTR-90-17A-C (1990) Simulation: Hollmann, C., Gutheil. E., *Combust. Sci. and Tech.* **135** 1-6, 175 (1998).





Modeling of Turbulent LOX/H₂ Spray Flames Micro Combustion Chamber M3 (DLR Lampoldshausen)

OH-Emission, p = 5 bar, $T_0 = 100$ K



Sender, J., et al., Proceedings of the 13th Annual Conference on Liquid Atomization on Spray Systems, Florence, Italy, 145-154 (1997).



Modeling of Turbulent Spray Flames



Experiment: Sender, J., et al., Proceedings of the 13th Annual Conference on Liquid Atomization on Spray Systems, Florence, Italy, 145-154 (1997).
Simulation: Schlotz, D., Brunner, M., Gutheil, E.: Modeling of Turbulent LOX/H₂ Combustion under Cryogenic and Elevated Pressure Conditions, ILASS Europe Conference, Zürich, September 2-6 (2001).



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Mixing in Turbulent Sprays

- The β -function that is typically used to describe the mixing in turbulent diffusion flames does not perform well in regions where vaporization is present¹.
- Here: Modification of the description of the β -function through use of a transport equation for the probability density function of the mixture fraction, \tilde{f} , in turbulent sprays²:

$$\overline{\rho}_{g}\frac{\tilde{\mathcal{J}}}{\partial t}+\overline{\rho}_{g}U_{j}\frac{\tilde{\mathcal{J}}}{\partial x_{j}}+\frac{\partial(\overline{\rho}_{g}\overline{S}_{s}\widetilde{f})}{\partial\zeta_{c}}=-\frac{\partial}{\partial\zeta_{c}}\left[\overline{\rho}_{g}\left\langle\frac{\partial}{\partial x_{j}}\left(D_{M}\frac{\partial\xi_{c}}{\partial x_{j}}\right)\middle|\zeta_{c}\right\rangle\widetilde{f}\right].$$

¹Miller R.S. Bellan, J. On the Validity of the Assumed Probability Density Function Method for Modeling Binary Mixing/ Reaction of Evaporated Vapor in Gas/Liquid-Droplet Turbulent Shear Flow, *Proc. Combust. Inst.* **27**: 1065-1072, 1998.

²Ge, H.-W., Gutheil, E.: PDF Simulation of Turbulent Spray Flows, *Atomiz. Sprays*, 2004, submitted.

Mixing in Turbulent Methanol/Air Sprays

Methanol Vapor Fraction and PDF of the Mixture Fraction



Ge, H.-W., Gutheil, E.: PDF Simulation of Turbulent Spray Flows, Atomiz. Sprays, 2004, submitted.

Probability Density Functions at Various Positions



Ge, H.-W., Gutheil, E.: PDF Simulation of Turbulent Spray Flows, Atomiz. Sprays, 2004, submitted.

Comparison of Results with Presumed and Monte-Carlo PDF, and with Experiment

x = 25 mm





Simulation: Ge, H.-W., Gutheil, E.: PDF Simulation of Turbulent Spray Flows, Atomiz. Sprays, 2004, submitted.



Summary and Conclusions

- LOX/H₂ spray flames in the counter-flow configuration have been studied, and the gaseous oxygen profile shows a non-monotonic behavior because of the high reactivity of the system. The flames persist to strain rates up to 25,000/s, and extinction has not yet been found.
- Multiple structures of laminar methanol/air counter-flowing spray flames have been identified at low strain rates up to 400/s on the spray side of the configuration for the present conditions. The gas-sided spray flame shows the same inner structure as a pure gas flamelet with appropriate boundary conditions, and this simplifies the implementation of the flamelet model for turbulent spray diffusion flames.
- The assumed β -function for the turbulent mixing in spray flames is poor in regions where vaporization exists, and it has been replaced by a PDF transport equation for the mixture fraction. A modified β -function is suitable to predict the shape of the PDF of the mixture fraction.



Future Research

- Extension of the model to unsteady flamelets
- Application of the PDF method to turbulent spray flame simulations
- Extension to other liquids