

COMPUTATIONAL ANALYSIS OF THE MIXING ZONE IN THE COMBUSTION CHAMBER OF RAMJET

Adil A. Al-Hemiri *and Sa'ad A. Fa'ek

*Chemical Engineering Department - College of Engineering - University of Baghdad – Iraq

Abstract

A theoretical analysis of mixing in the secondary combustion chamber of ramjet is presented. Theoretical investigations were initiated to insight into the flow field of the mixing zone of the ramjet combustor and a computer program to calculate axisymmetric, reacting and inert flow was developed. The mathematical model of the mixing zone of ramjet comprises differential equations for: continuity, momentum, stagnation enthalpy, concentration, turbulence energy and its dissipation rate. The simultaneous solution of these equations by means of a finite-difference solution algorithm yields the values of the variable at all internal grid nodes.

The results showed that increasing air mass flow (0.32 to 0.64 kg/s) increases the development of velocity profile due to the high turbulence generated resulting in very fast mixing and homogenous flow. And the occurrence of chemical reaction causes higher local temperature and composition resulting in faster development of the velocity profile

Keywords: oil Non-premixed chemically reacting ducted flows, Mixing of two streams, Ramjet modeling.

Introduction

Increased interest in ramjet propulsion system with higher performance requirements and tighter constraints on system size and weight has led to the need for improved techniques for analyzing and designing such system. A critical requirement for achieving high system performance within specified geometric limits is an accurate description of the secondary combustor flow field, including the effect of the mixing zone and chemical reaction.

In a conventional application (fig.1), the mixing zone is necessary for more complete combustion before the gases exit through the nozzle. It is important to be able to determine the combustion behavior within this region as a function of upstream conditions.

A recent study (Faek, 2004) gave a comprehensive survey on the subject and considered the effect of various parameters on the velocity profile. These parameters

were: air mass flow rate, air to fuel ratio, working pressure, air and fuel temperatures. He showed that:

Increasing air mass flow increases the development of velocity profile due to the high turbulence generated resulting in very fast mixing and homogenous flow.

Increasing air and fuel temperature reduces the cooling effect of air and gives higher mixing temperature because it increases the thermal energy exchange between the air and fuel streams.

Increasing upstream combustor pressure increases the development of velocity profile due to the development of highly recirculating region.

The fuel to air ratio has little effect on the mixing zone characteristics, since it only effect the inlet velocities of air and fuel thus it give the same profile and the mixing temperature.

The best air mass flow was reported to be 0.64kg/s, upstream combustor pressure of 8 bar, air inlet temperature of 500K, fuel inlet temperature of 1100K and fuel to air ratio of 0.125 (Faek,2004).

The purpose of this paper is to study the effect of air flow rate on the mixing zone velocity profile (for the two cases of reacting and non-reacting systems) by using a governing equations and solving them by a computer program developed for this goal.

MODEL OVERVIEW

The flow was assumed to be steady, two-dimensional, and subsonic. For simplicity, the value of specific heat (CP) was assumed to be constant although its dependence on temperature and/or composition could be easily included.

A modified Jones-Launder,(Jones and Launder,1972; Launder and Spalding,1974) two- parameter turbulence model was incorporated to calculate the effective viscosity. It uses five empirical constants (Table 1) and requires that two additional variables [turbulence kinetic energy (k) and turbulence dissipation rate (ε)] be evaluated. Effective viscosity was calculated using the formulas

$$\mu_{eff} = \mu_t + \mu \tag{1}$$

Where,

$$\mu_t \equiv \rho C_\mu k^2 / \epsilon \tag{2}$$

For reacting flows, four species were considered: oxygen, nitrogen, fuel, and products. Simple, one step, infinitely fast kinetics was assumed in which 1kg of fuel combines with i kg of oxidant to form (1+ i) kg of product without intermediaries (Moult and Srivatsa, 1977). Fuel and oxygen, therefore, could not exist simultaneously, and the combustion process was mixing limited. The turbulent Prandtl and Schmidt numbers were taken equal to unity, and therefore, the turbulent Lewis number was unity; the laminar Prandtl number was also taken to be unity.

The conservation equations for axisymmetrical flows with no tangential variations can be put into general form (Khalil, E.E., Spalding and Whitelaw,1975).

$$\frac{\partial}{\partial x}(\rho U \psi) + \frac{1}{r} \frac{\partial}{\partial r}(r \rho V \psi) = \frac{\partial}{\partial x} \left(\Gamma_\psi \frac{\partial \psi}{\partial x} \right) + \frac{1}{r} \frac{\partial}{\partial r} \left(r \Gamma_\psi \frac{\partial \psi}{\partial r} \right) + S_\psi \tag{3}$$

where ψ stands for the dependent variables (U,V,k,ε,h, etc....) being considered (ψ=1 for continuity equation), Γψ is the appropriate effective exchange coefficient for turbulent flow, and Sψ as the "source term" (Table 2). The energy equation in terms of stagnation enthalpy has

no source terms since the turbulent Prandtl and Schmidt numbers were chosen equal to unity and radiative transport was neglected. Thus, the stagnation enthalpy is given by

$$h = h' + (U^2 + V^2) / 2 \tag{4}$$

where for non-reacting flows

$$h' \equiv C_p T \tag{5}$$

and for reacting flows

$$h' \equiv m_{ox} \Delta H / i + C_p (T - T_{ref}) \tag{6}$$

Table (1) Turbulence model constants (Launder and Spalding, 1974)

Constant	Value
C_1	1.44
C_2	1.92
C_μ	0.09
K	0.42
E	8.8

$$\sigma_k = \sigma_h = \sigma_f = 0.9$$

$$\sigma_\epsilon = K^2 / \left[(C_2 - C_1) C_\mu^{1/2} \right] = 1.22$$

Table (2) Governing equations parameters (Launder and Spalding, 1974)

Conse rvatio n of	ψ	Γ_ψ	S_ψ
Mass	1	0	0
Axial mome ntum	U	μ_{eff}	$\frac{\partial}{\partial x} \left(\mu_{eff} \frac{\partial U}{\partial x} \right) - \frac{\partial P}{\partial x} + \frac{1}{r} \frac{\partial}{\partial r} \left(r \mu_{eff} \frac{\partial U}{\partial r} \right)$
Radia l mome ntum	V	μ_{eff}	$\frac{1}{r} \frac{\partial}{\partial r} \left(r \mu_{eff} \frac{\partial V}{\partial r} \right) - \frac{\partial P}{\partial r} + \frac{\partial}{\partial x} \left(\mu_{eff} \frac{\partial U}{\partial r} \right) - 2 \mu_{eff} \frac{V}{r^2}$
Kineti c energ y	k	$\frac{\mu_{eff}}{\sigma_k}$	$G - \rho \epsilon$
Dissi pation rate	ϵ	$\frac{\mu_{eff}}{\sigma_\epsilon}$	$\frac{\epsilon}{k} (C_1 G - C_2 \rho \epsilon)$
Stagn ation enthal py	h	$\frac{\mu_{eff}}{\sigma_h}$	0
Mass fracti on of fuel	m	$\frac{\mu_{eff}}{\sigma_{fu}}$	0
Mixtu re fracti on	f	$\frac{\mu_{eff}}{\sigma_f}$	0

The particular ones employed for the equation of the three velocity components, k and ϵ , are:

$$U_P \frac{(C_\mu^{1/4} k_P^{1/2})}{(\tau/\rho)_w} = \frac{1}{\kappa} \ln(EC_\mu^{1/4} k_P^{1/2} y_P / \nu) \quad (7)$$

$$k_P = (\tau/\rho)_w C_\mu^{-1/2} \quad (8)$$

$$\epsilon_P = \frac{C_\mu^{3/4} k_P^{3/2}}{\kappa y_P} \quad (9)$$

The Solution Procedure

The set of differential equations are first reduced to finite-difference equations exhibiting "hybrid formulation of the coefficients", i.e., coefficient that contain combinations of the convective flux per unit mass F and the diffusive conductance. And then solved iteratively by "SIMPLE" procedure (Patankar and Spalding, 1972; Patankar, 1980; Versteeg and Malalasekera, 1995). The grid arrangement for discretisation is as shown in figure (2). And the logic diagram for SIMPLE algorithm is given by figure (3), where the procedure steps are as given below:

- Start
- Initial guess of p^* , u^* , v^* , Φ^* .
- STEP 1: Solving discretised momentum equations,

$$a_{i,j} U^*_{i,j} = \sum a_{nb} U^*_{nb} + (P^*_{I-1,J} - P^*_{I,J}) A_{i,J} + b_{i,J}$$

$$a_{i,j} V^*_{i,j} = \sum a_{nb} V^*_{nb} + (P^*_{I,J-1} - P^*_{I,J}) A_{i,j} + b_{i,j}$$
- STEP 2: Solving pressure correction equation,

$$a_{I,J} P'^*_{I,J} = a_{I+1,J} P'^*_{I+1,J} + a_{I-1,J} P'^*_{I-1,J} + a_{I,J+1} P'^*_{I,J+1} + a_{I,J-1} P'^*_{I,J-1} + b_{I,J}$$
- STEP 3: Correction of pressure and velocities,

$$P_{I,J} = P^*_{I,J} + P'_{I,J}$$

$$U_{I,j} = U^*_{I,j} + d_{i,j} (P'_{I-1,J} - P'_{I,J})$$

$$V_{I,j} = V^*_{I,j} + d_{I,j} (P'_{I-1,J} - P'_{I,J})$$
- STEP 4: Solving the discretised equations for ρ and T ,

Special Treatment near A Wall

To avoid the need for detailed calculations in the near-wall grid regions, algebraic relations are employed for the near wall grid nodes, which have to be spaced at such a distance from the neighboring walls that they lie within the so-called logarithmic layer. Such relations are termed wall functions which are derived so as to reproduce identically the implications of the logarithmic profiles with uniform shear stress prevailing up to the near wall grid node, with generation and dissipation of energy are in balance at this locality (Patankar and Spalding, 1970).

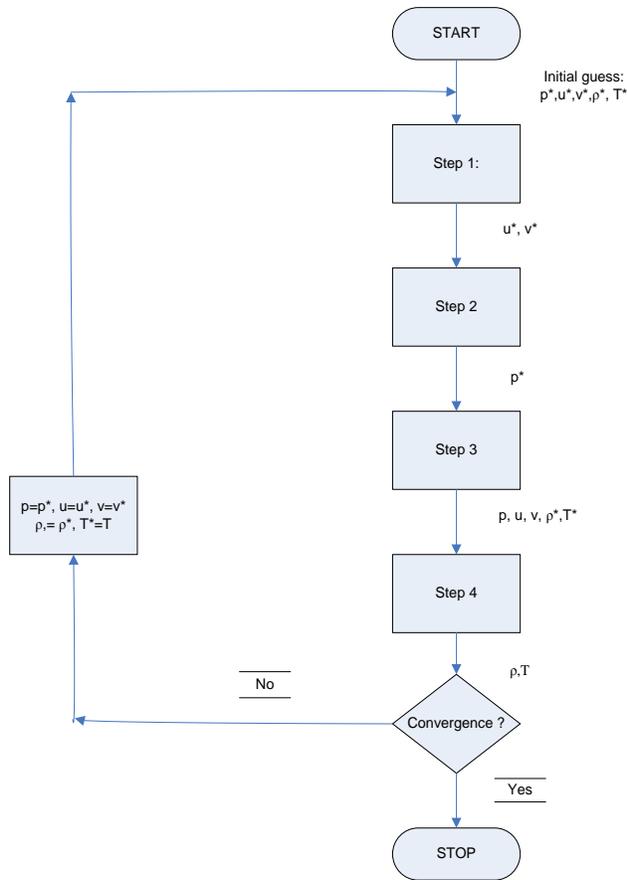


Figure (3); Logic Diagram For Program "SIMPLE"

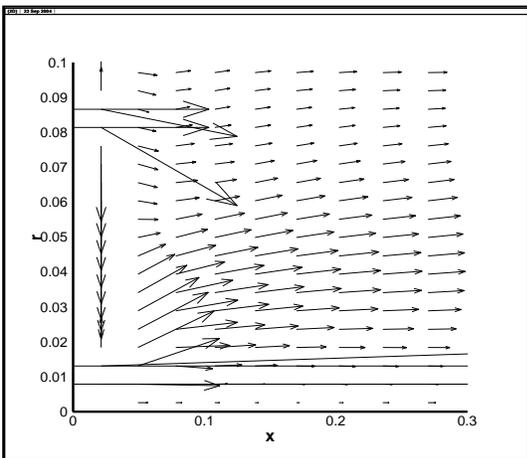


Fig.4: velocity vector plot at $\phi=0.125$, $m_a = 0.64$ kg/s, $p_c=8$ bar, $T_f=1100$ k and $T_a= 500$ k, and inert flow system.

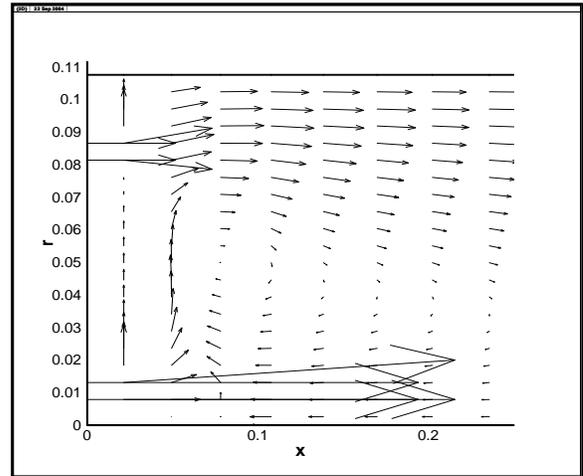


Fig.5: velocity vector plot at $\phi=0.125$, $m_a = 0.32$ kg/s, $p_c=8$ bar, $T_f=1100$ k and $T_a= 500$ k, and inert flow system.

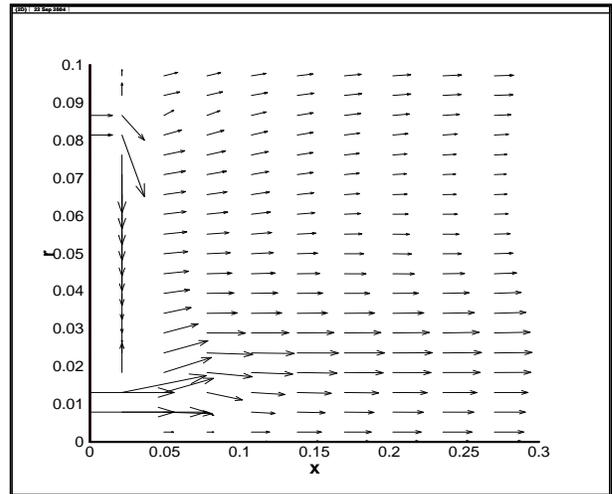


Fig.6: velocity vector plot at $\phi=0.125$, $m_a = 0.64$ kg/s, $p_c=8$ bar, $T_f=1100$ k, $T_a= 500$ k, and reacting flow system.

