Numerical Simulation of Gravitonal Turbulent Mixing with the Acceleration Sign Changed

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The problem of turbulent mixing under constant gravity (constant acceleration) creating an unstable situation on a plane interface of two incompressible fluids is studied experimentally in a number of papers [1-4]. Some papers [2,5-8] also study the problem numerically through direct 3D numerical simulations with code TREK.

Refs. [9,10] perform experiments with alternating sign of acceleration resulting in an unstable situation at the interface. The relevant problem is numerically studied in [11] with a turbulence type phenomenological model [12], however the mixing zone decrease effect at the steady-state stage can not be described with the model.

This paper studies the problem of turbulent mixing at a plane interface of two incompressible fluids of density difference n=3 with alternating acceleration sign through direct numerical simulation with 3D code TREK. The computation is parallelized into several tens of processors.

The problem is set up much like in refs. [6-8]: at the initial time two semi-spaces separated with plane $z=z_z=0$ are filled with resting ideal gases of densities $\rho_1=1$ and $\rho_2=n$

(n=3, Atwood number $A = \frac{p_2 - p_1}{q_2 - q_1}$ $_{2}$ + μ_{1} $A \equiv \frac{\rho_2 - \rho_1}{\rho_2}$ $\equiv \frac{\rho_2 - \rho_1}{\rho_2 + \rho_1}$). If $t \le t^* = 3$, gravity $g_z = -g = -1$ is directed from the heavy

material to the light; if $t > t^*$, $g_z = g_{12} = 22/75$.

At the initial time the random-number generator sets random density perturbations at the interface (in a layer one cell in thickness): $\delta \rho = \pm \rho_1 \cdot \delta$, where $\delta = 0.1$.

The gas dynamics equations for ideal two-component medium (with zero molecular viscosity and heat conductivity) are solved.

The computational domain is a parallelepiped with the vertical side Λ =2 of side face. Its horizontal face is a square with side $L_x = L_y = 1$.

The initial pressure profile was given reasoning from the hydrostatic equilibrium condition.

The coordinate of the upper face is $z_2 = 0.85$, that of the lower $z_1 = -1.15$.

The initial pressure is such, that for the turbulent flow type under discussion the incompressibility condition was adequately met:

The equation of state is ideal gas with adiabatic constant $\gamma=1.4$.

The computational grid is uniform with $N_x=201$, $N_y=201$, $N_z=401$ cells. The cell size is $h_x = h_y = h_z = 0.005$.

The condition of "rigid wall" type was imposed on all computational domain boundaries.

The computed data is reported in the form of pixel patterns of light material concentrations for different times in different horizontal sections.

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The flow time history observed at the first (unstable) problem step is similar on the whole to the previous computations [2,5-8]: vortex enlargement with time and tapering-off to a selfsimilar regime at the end of the above step are observed.

The self-similar regime for that step manifests itself, in particular, in the tapering-off to the linear time function of turbulent mixing zone width $L_t(t)$:

$$
F \equiv \frac{1}{t_0} \sqrt{\frac{L_t}{Ag}},
$$

where $t_0 \equiv \sqrt{\frac{L_x}{g}}, \quad (\tau \equiv t / t_0).$ (1)

Here the slope angle $dF/d\tau$ determines coefficient 2 *a dF* $\alpha_a = \frac{d}{d}$ $=\left(\frac{dF}{d\tau}\right)^2$ in the relation for the

mixing zone width at the first (self-similar) step:

$$
L_{ia} = \alpha_a A g t^2. \tag{2}
$$

The calculated $\alpha_a \approx 0.05$ is slightly less than the data of our previous calculations [7,8] and that presented elsewhere $[2,5]$ – the difference is explained by the high magnitude of ε =0.1, the same as in expeiments [9,10].

At the second step, $F(\tau)$ linearly decreases on the whole (to an accuracy of small fluctuations relating to the wave passage) with the slope angle $\alpha^{(-)} = \frac{dF}{dt} \approx -0.084$ $\alpha^{\vee} = \frac{1}{d}$ σ^{-1} = $\frac{dr}{d\tau}$ ≈ -0.084 -

$$
Refs. [9,10] measure
$$

$$
f \equiv \sqrt{2A\alpha_s} \equiv \frac{d\sqrt{z_2 - z_c}}{d\sqrt{S'}}
$$
\n(3)

where

$$
S' \equiv g_{12} \frac{\left(t - t_c\right)^2}{2},
$$

 t_c is the time of F_2 peak.

From our calculation, the expression for f is

$$
f = \frac{dF_2}{d\tau} \cdot \sqrt{A \cdot \frac{2g}{g_{12}}} \approx -0.074,
$$

which is somewhat less (modulo) than the value of $f \approx -0.1 (\alpha_s \approx 0.01)$ suggested in [9,10] reasoning from the experiments.

For this problem the self-similar regime also manifests itself in tapering-off to the timedependent value

$$
E_m(t) \equiv \max(E)
$$

– the scaled turbulent energy maximal in the TMZ width:

$$
E \equiv \frac{k}{L_{i}g}
$$

With the change in the acceleration sign, the E_m first noticeably drops and then increases slowly, with severe fluctuations.

The paper treats squared density and mass flow fluctuations σ maximal in TMZ.

It is shown that at the second step the resultant flow on the whole reverses sign relative to the normal (steady) situation (to an accuracy of the wave passage related fluctuations).

The paper makes comparison to the results of the calculations by the $k - \varepsilon$ turbulence model performed by the adequate, 1D EGAK-T computation on Lagrangian computational grid.

The flow time history observed at the first (unstable) step of the problem is similar on the whole to the previous computations: the tapering-off to the self-similar regime with time is observed.

For that step it manifests itself, in particular, in the tapering-off to the linear time function F of turbulent mixing zone width L_{μ} (t).

However, at the second step (on the acceleration sign change), although the TMZ width growth rate slowing is noticeable, it can be seen that the $k - \varepsilon$ turbulence model can not decrease the width, its growth slows down and tapers off to a constant at late times, although an abrupt decrease in the turbulent energy is observed on the acceleration sign change.

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