

Simulation of the QUEOS Experiment using Rigid Dynamic-Moving Particle Semi-Implicit (RD-MPS) Method

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1. Introduction

Particle dissipation and mixing in liquid are common multiphase phenomena not only in nature and but also in industrial processes. Also, It is one of important phenomena in nuclear safety analysis on severe accidents associated with the coolability of the corium debris bed, in which the characteristics of the porous corium debris bed such as local porosity distribution, debris bed configuration, debris size distribution, porous structure etc., determined by the corium jet break-up, precipitation, and mixing processes are considered to be of importance. Therefore, we developed a new computational tool, called ADDA (Analysis of Debris Dynamics and Agglomeration), based on an enhanced MPS (Moving Particle Semi-implicit) algorithm (Park, 2011) [1] to understand the complex debris dissipation and mixing phenomena and identify the roles of the debris characteristics in the process and verified against the QUEOS experiment (Meyer, 1996; 1997) [2, 3] performed at FzK in Germany.

2. Methods and Results

2.1. Computational algorithm

The description and formulations on the enhanced MPS can be found in Park's work [1]. In this section, the coupling way of these two methods is introduced.

At the stage of initializing particle configurations, initial velocities, positions, and pressures are specified. A particle's diameter for the rigid body dynamics calculation is set to $0.9l_0$. If the diameter is set to the same as l_0 , the fluid particles constantly conflict with each other and the contact point calculations become too complicated. If the diameter is much smaller than $0.9l_0$, the incompressibility is not well conserved and the overall calculation becomes identical to the result calculated only by the MPS method.

The entire calculation is divided into two stages for each time step. The first stage is the MPS calculation in which external forces induced by gravity, surface tension, and viscosity are calculated into temporal velocities and the pressure Poisson matrix is solved iteratively. The particle motion information gained from the first stage including \mathbf{u}_i^{**} , \mathbf{r}_i^{**} , and P_i^{**} is transferred to the rigid body dynamics calculation of the second step. In

the dynamics calculation, the velocity change generated by the MPS calculation is converted to the force by

$$\mathbf{F} = \frac{m_i(\mathbf{u}_i^{**} - \mathbf{u}_i^t)}{dt}. \quad (1)$$

Using the above forces acting on each fluid particle's center of mass, the colliding contacts and the resting contacts are calculated to obtain the velocities of the next time step.

2.2. Description on the QUEOS experiment

In this paper, the verification of the ADDA code was performed against the QUEOS experiments which was designed to establish the data base for testing the heat and momentum transfer models in the FCI (Fuel-Coolant Interaction) multi-fluid codes. For the code verification, we selected the Q21 test among the QUEOS tests conducted in a non-boiling condition (ZrO₂ particles at the ambient temperature and water temperature of 99 °C.) since the prime objective of the present verification is the hydrodynamic interactions between solid and liquid particles in the processes of jet breakup, mixing and precipitation. In the Q21 tests, a total of 18000 ZrO₂ particles in a shape of a jet with a diameter of 100 mm and a corresponding mass of 7 kg and a volume of 1830 cm³ was injected into a water pool for 55±5 ms, resulting the jet length of 27±3 cm. Each ZrO₂ particle has a diameter of 4.95 mm.

2.3. Particle precipitation and mixing simulation

To simulate the Q21 test of the QUEOS experiment, the initial configuration of the model is set as Figure 1. The width of the test section is 0.7 m and the water level is 1 m. The initial distance between particles, l_0 , is set to be 0.013 m, the radius of interaction, r_e , is $2.1 l_0$ and the kernel size for Laplacian is $4.0l_0$. The density of water is 1000 kg/m³, viscosity of water is 0.00109 Pa·s, the surface tension coefficient of water is 0.0728 N/m, and the friction coefficient of ZrO₂ particle is set to 1.0. To make a two-dimensional equivalent geometry, the mass of a particle is set to 0.097 kg. The initial downward velocity of the particles is 5.05 m/s.

Fig. 2 shows the side view of the experiment and the calculated results. Because the calculation was performed in two dimensions, the water level was overestimated comparing to the measured level. The calculated penetration velocity of the particles was lower

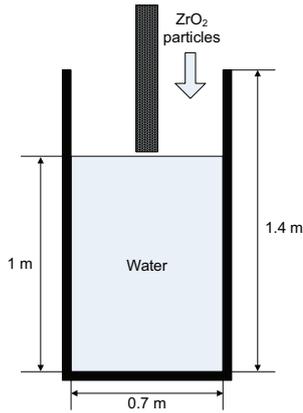


Figure 1: Schematic initial configuration of the Q21 test.

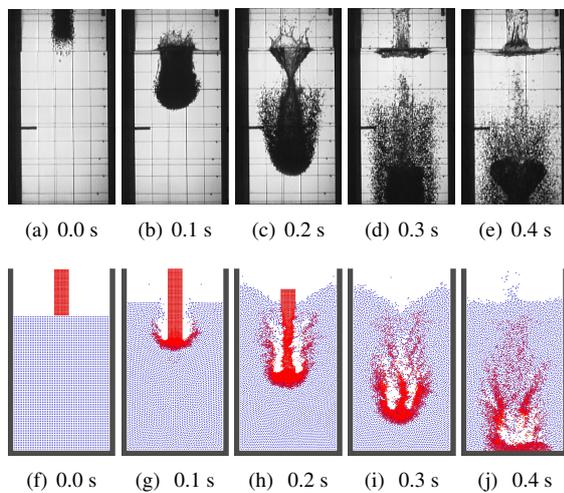


Figure 2: View from the side of the test section.

than the measured data. This mismatch was resulted from the fact that the calculation was performed in two-dimension. However, the simulation re-created the jet breakup and penetration behavior observed in the experiment, and revealed the structural details of the jet breakup process, showing the jet surface and the leading edge instability and boundary stripping phenomena.

Contrary to the previous MPS analysis [4], the ADDA simulation can simulate the complete process of the jet mixing phenomena from the jet penetration to the particle precipitation on the bottom of the test vessel due to its enhanced stability which was mentioned earlier. This capability makes it possible to simulate the final particle bed formation on the vessel bottom and to be compared with the experimental data, Q21.

Fig. 3 plots the normalized mass distribution of spheres on the bottom of the test vessel. The measured data at the center of the test vessel was compared to the calculation result. In the figure, the calculated result shows more flat leftover at the bottom because the simulation was two-dimensional.

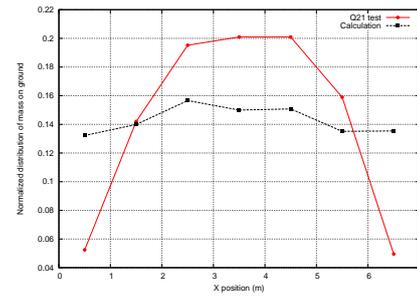


Figure 3: The normalized distribution of the particles on the bottom of the test vessel.

3. Conclusions

In this paper, the ADDA code with the RD-MPS algorithm[1] was used to simulate the non-boiling Q21 test of the QUEOS experiment in two dimensions. The code successfully created the complex particle jet mixing phenomena including particle penetration, mixing with surrounding coolant, dissipation and precipitation. Due to its limitation of three-dimensional calculation capability, the specific details of the phenomena such as the change of water level and penetration velocity were not accurately predicted. However, the code simulated the characteristic structure of particle jet breakup and mixing with surrounding liquid very well. Importantly, the enhanced stability of the code performance makes it possible to simulate the entire jet mixing process, predicting the particle bed formation on the bottom of the test vessel. The code is currently under heavy development to extend its capability to the three-dimensional model.

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