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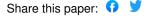
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# Simulation of Gravity Wave Propagation in Free Surface Flows by an Incompressible SPH Algorithm

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#### ABSTRACT

This paper presents an incompressible smoothed particle hydrodynamics (SPH) model to simulate wave propagation in a free surface flow. The Navier-Stokes equations are solved in a Lagrangian framework using a three-step fractional method. In the first step, a temporary velocity field is provided according to the relevant body forces. This velocity field is renewed in the second step to include the viscosity effects. A Poisson equation is employed in the third step as an alternative for the equation of state in order to evaluate pressure. This Poisson equation considers a trade-off between density and pressure which is utilized in the third step to impose the incompressibility effect. The computations are compared with the experimental as well as numerical data and a good agreement is observed. In order to validate proposed algorithm, a dam-break problem is solved as a benchmark solution and the computational results are compared with the previous numerical ones.

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NOMENCLATURE				
V	particle velocity	q=r/h	non-dimensional distance between particles	
$V^*$	particle first temporary velocity	A	a field variable	
$V^{**}$	particle second temporary velocity	B	a vector	
$\hat{V}$	particle corrected velocity without using XSPH	D	shear strain rate	
$\overline{V}$	particle velocity by using XSPH correction	T	period of signals	
P	particle pressure	Greek let	Greek letters	
t	time	ho	density of water	
g	gravitational acceleration	$\boldsymbol{\rho}^*$	a temporary density	
r	position vector	$ ho_0$	initial density	
X	horizontal coordinate	au	shear stress tensor	
у	vertical coordinate	$\phi$	a scalar or vector quantity	
W	interpolation kernel	$\eta$	small number to avoid singularity	
h	smoothing distance	$\mu$	constant viscosity of water	
m	particle mass	${oldsymbol{arepsilon}}$	constant XSPH correction	

#### 1. INTRODUCTION

The study of wave propagation in free surface flows has many benefits and is being given great attention in

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practice. Surface water waves have a high potential to cause damages and losses of lives in coastal areas. Predicting the damage of these waves is of importance when assessing risk and magnitude of flooding in these areas.

Previous investigations have been focused on the studies of wave propagation and overtopping through

using the experimental, analytical and numerical approaches. For example, Cox and Ortega [1] carried out an experimental study on the wave overtopping over a fixed deck using irregular and transient waves.

Unfortunately, due to the complexity of this phenomenon, number of reliable laboratory and field data is limited. Therefore, numerical methods are widely used for simulating of wave propagation.

Most of the numerical works on the recent wave problems have been carried out using the Eulerian gridbased methods with quite satisfactory results, such as finite element, finite volume and finite difference methods. These methods have provided many useful and satisfactory results. However, their successes largely rely on good quality meshes. The construction of such high quality mesh configurations is usually a difficult and time-consuming task, since it must be ensured that the aspect ratios of all elements are neither very large nor very small and thus the connectivity between nodes and elements must be carefully and accurately found and recorded. In addition, elements can frequently become over-distorted during the simulation of water wave evolutions. The over-distorted elements may be amended by remeshing. However, remeshing can be as expensive as the generation of original meshes and may take a major proportion of computational costs considering when this has to be done almost every time step.

A new class of methods has recently been developed, which can overcome the problems associated with mesh-based methods. These methods do not need using any mesh to discrete computational domains and they are only based on randomly-ordered and –distributed nodes. Many meshless (or particle) methods have been reported in the literature, such as meshless local Petrove–Galerkin (MLPG) method [2], moving particle semi-implicit (MPS) method [3], smoothed particle hydrodynamics method (SPH) [4] and so on. Among them, the SPH method has been used widely to simulate water wave problems [4-6].

The SPH method uses a purely Lagrangian approach and has been successfully employed in a wide range of applications, such as incompressible fluid flow, multiphase flow, turbulence, heat and mass transfer, elasticity and fracture, and etc [7]. Recently, SPH has also been employed to simulate fluid-structure interaction [8-10]. The meshless characteristic of SPH makes the method needless of using data connectivity as needed by the finite volume and finite element methods. This gives the method a very useful feature when dealing with complex flows exhibiting large deformations and/or free-surfaces.

The method was originally developed by Lucy [11] and Gingold and Monaghan [12] to solve compressible astrophysical problems, and then it was later extended to incompressible flows by Monaghan [4]. Several other

researchers have contributed to the method and solved various engineering problems including turbulent flows, interfacial flows and sloshing problems [13-15].

Dalrymple et al. [5] and Gomez-Gesteira et al. [16] employed this method to study the wave overtopping over coastal structures. In the previous simulations of incompressible flows using the SPH model, the incompressibility was achieved by employing an equation of state and the fluid was assumed to be slightly compressible. This approach is denoted as the weakly compressible SPH, since a large enough sound speed has to be introduced into the pressure equation to ensure the numerical stability. Based on the semiimplicit algorithm of the MPS method [3], a truly incompressible version of the SPH [17] has been proposed, in which the free surfaces were identified and tracked by particles without numerical diffusion. The difference between the original compressible SPH [11, 12] and the incompressible SPH [17] lies in that the former calculates the pressure using an equation of state, while the latter employs a strictly incompressible SPH formulation, in which the pressure is not an explicit thermodynamic variable but obtained implicitly through solving a pressure Poisson equation derived from the mass and momentum equations. In this sense, it is also very similar to the pressure projection method widely used in grid based methods and the projection SPH of Cummins and Rudman [18].

In this paper a new incompressible SPH algorithm based on Shao et al. [17] is presented to simulate gravity wave propagation in free surface flow problems. Before this simulation, in order to show the ability of the method for simulating of free surface problems, a dam-break problem is modeled as a benchmark and the results are compared with the results of standard SPH method.

The proposed algorithm is similar to three step explicit SPH algorithm proposed by Hosseini et al. [19] for simulation of incompressible fluid flows. In the first step of this algorithm, the momentum equation is solved in the presence of the body force neglecting all other forces. The calculated temporary velocities are renewed in the second step to include the viscosity effect. A Poisson equation is employed in the third step as an alternative of the equation of state in order to evaluate pressure of the particles. This Poisson equation considers a trade-off between density and pressure which is utilized in the third step to impose the incompressibility effect.

In fact there is no difference between the presented algorithm and the SPH algorithm proposed by Hosseini et al. [19], the purpose of this paper is to explore the potential of newly developed SPH method for modeling the propagation of nonlinear water waves. Wave propagation problems are of great importance to offshore engineering and any new numerical method,

even successfully applied to other fluid problems, is usually required to be carefully refined when it is applied to water waves. The SPH computations are compared with the experimental data of Cox and Ortega [1] and the numerical results of Gomez-Gesteira et al. [16].

#### 2. GOVERNING EQUATIONS

The governing equations for simulating free surface flow in 2-D dimensions are the mass and momentum conservation equations. With regard to fluid particles, they are written in Lagrangian form as follow:

$$\frac{1}{\rho} \frac{D \rho}{D t} + \nabla . V = 0 \tag{1}$$

$$\frac{DV}{Dt} = -\frac{1}{\rho} \nabla P + g + \frac{1}{\rho} \nabla \cdot \vec{\tau}$$
 (2)

Equation (1) is in the form of a compressible flow. The purpose is that the deviation of fluid densities at the particle can then be used to enforce incompressibility in the correction step of time integration.

#### 3. SPH FORMULATION

The SPH formulations as developed by Monaghan [20] are obtained by interpolating from a set of points that may be disordered. The interpolation is based on the theory of integral interpolants using kernels that approximate a delta function. The interpolants are analytic functions that can be differentiated without the use of grids. If the points are fixed in position, the equations reduce to finite difference equations, with different forms depending on the interpolation kernel. The SPH equations describe the motion of the interpolating points, which can be assumed as particles. Each particle carries a mass m, a velocity V, and other properties, depending on the problem.

**3. 1. Interpolation** Using the above concepts, any quantity of particle i, whether scalar or vector, can be approximated by the direct summation of the relevant quantities of its neighboring particles j

$$\phi_i(r_i) = \sum_j m_j \frac{\phi_j(r_j)}{\rho_j(r_i)} W\left(r_i - r_j \mid h\right)$$
(3)

where i and j= reference particle and its neighbors.

**3. 2. Kernel** The use of different kernels is the SPH analogue of using different schemes in finite difference methods. While different equations can have different kernels, usually the same kernel is used throughout all formulations in one model. By balancing the

computational accuracy and efficiency, the following kernel based on the spline function and normalized in 2-D is adopted [20].

$$W(r,h) = \frac{10}{7\pi h^{2}} \times \begin{cases} \left(1 - \frac{3}{2}q^{2} + \frac{3}{4}q^{3}\right) & q < 1\\ \frac{1}{4}\left(2 - q^{3}\right) & 1 < q < 2\\ 0 & q > 2 \end{cases}$$
 (4)

The smoothing distance or kernel range h determines the degree which a particle interacts with neighboring particles.

This kernel has the advantage of possessing compact support, the second derivative being continuous and the dominant error term in the integral interpolant being the order of  $h^2$ . The continuity of the second derivative means that the kernel is not too sensitive to particle disorder and the errors in approximating the integral interpolants by summation interpolants are small provided the particle disorder is not too large.

**3. 3. Gradient, Divergence and Laplacian** The formulation of the gradient term in the Navier–Stokes equation has different forms depending on the derivation used [20]. The following symmetric forms are employed for gradient of a scalar *A* and divergence of a vector *B* since it conserves linear and angular momentum exactly

$$\frac{1}{\rho_i} \nabla_i A = \sum_i m_j \left( \frac{A_i}{\rho_i^2} + \frac{A_j}{\rho_i^2} \right) \nabla_i W_{ij} \tag{5}$$

$$\frac{1}{\rho_i} \nabla_i \cdot B_i = \sum_j m_j \left( \frac{B_i}{\rho_i^2} + \frac{B_j}{\rho_j^2} \right) \nabla_i W_{ij}$$
 (6)

where the summation is over all particles other than particle i and  $\nabla_{i}W_{ij}$  = gradient of the kernel taken with respect to the positions of particle i. In practice only near neighbors contribute, because the kernel has a finite range.

A simple way to formulate the Laplacian operator is to envisage it as dot product of the divergence and gradient operators. This approach proved to be problematic as the resulting second derivative of the kernel is very sensitive to particle disorder and when dealing with the Navier-Stokes equations can easily lead to pressure instability and decoupling in the computation due to the co-location of the velocity and pressure. In this paper, the following alternative approach is adopted [17]

$$\nabla \left(\frac{1}{\rho} \nabla A\right)_{i} = \sum_{j} m_{j} \frac{8}{\left(\rho_{i} + \rho_{j}\right)^{2}} \frac{A_{ij} r_{ij} \cdot \nabla_{i} W_{ij}}{\left|r_{ij}^{2}\right| + \eta^{2}}$$
(7)

where  $A_{ij} = A_i - A_j$ ,  $r_{ij} = r_i - r_j$  and  $\eta$  is a small number introduced to avoid a zero denominator during computations and is set to 0.1h.

#### 4. THREE-STEP INCOMPRESSIBLE SPH ALGORITHM

In this section, an algorithm is presented to show the sequence of computation of each term in the governing equations. In this paper, a fully explicit three-step algorithm is used. In the first step of this algorithm, the momentum equation is solved in the presence of the body forces neglecting all other forces. As a result, an intermediate velocity is computed as

$$V^* = V_t + g\Delta t \tag{8}$$

Our experience has shown that it is important to impose the body forces in the first step of the solution algorithm especially in highly viscous fluids. In the second step, the calculated intermediate velocities are employed to compute the divergence of the stress tensor. Generally speaking, viscosity of incompressible generalized Newtonian fluids depends only on the second principal invariant of the shear strain rate  $D = (\nabla V + \nabla V^T)/2$ , i.e.,

$$|D| = \sqrt{\sum_{i,j} D_{ij} D_{ij}} \tag{9}$$

A classical constitutive law for these generalized Newtonian fluids is given by

$$\tau = \mu(|D|)D \tag{10}$$

For Newtonian fluids the familiar form  $\tau = 2 \mu D$  is recovered.

Consequently, the stress tensor  $(\tau)$  can be calculated for any specified constitutive law. In this work, the divergence of the stress tensor in the momentum equation is obtained as:

$$\left(\frac{1}{\rho}\nabla \cdot \tau\right)_{i} = \sum_{j} m_{j} \left(\frac{\overrightarrow{\tau}_{i}}{\rho_{i}^{2}} + \frac{\overrightarrow{\tau}_{j}}{\rho_{j}^{2}}\right) \cdot \nabla_{i} W(r_{ij}, h)$$
(11)

At the end of second step, the velocity components of each particle is updated according to:

$$V^{**} = V^* + \left(\frac{1}{\rho}\nabla \cdot \tau\right) \Delta t \tag{12}$$

At this stage, each particle is moved according to its second intermediate velocity  $V^{**} = (u^{**}, v^{**})$ .

Thus far no constraint has been imposed to satisfy the incompressibility of the fluid and it is expected that the density of some particles change during this updating. In fact, with the help of the continuity equation one can calculate the temporal fluid density of each particle as:

$$\frac{D\rho_i}{Dt} = \sum_i m_j (V_i - V_j) . \nabla_i W(r_{ij}, h)$$
(13)

The velocity field  $(\hat{V} = (\hat{u}, \hat{v}))$  which is needed to restore the density of particles to their original values is now calculated. To do this, in the third step of the

algorithm, the momentum equation with the pressure gradient term as a source term is combined with the continuity Equation (1) as

$$\frac{1}{\rho_0} \frac{\rho_0 - \rho^*}{\Delta t} + \nabla \cdot \hat{V} = 0 \tag{14}$$

$$\widehat{V} = -\left(\frac{1}{\rho^*} \nabla P\right) \Delta t \tag{15}$$

to obtain the following pressure Poisson equation

$$\nabla \left( \frac{1}{\rho^*} \nabla P \right) = \frac{\rho_0 - \rho^*}{\rho_0 \Delta t^2} \tag{16}$$

Equation (16) can be discretized according to Equation (7) to obtain the pressure of each particle as:

$$P_{i} = \left(\frac{\rho_{0} - \rho_{i}^{*}}{\rho_{0} \Delta t^{2}} + \sum_{j} \frac{8m_{j}}{\left(\rho_{i}^{*} + \rho_{j}\right)^{2}} \frac{P_{j} r_{ij} \cdot \nabla_{i} W_{ij}}{\left|r_{ij}\right|^{2} + \eta^{2}}\right)$$

$$\left(\sum_{j} \frac{8m_{j}}{\left(\rho_{i}^{*} + \rho_{j}\right)^{2}} \frac{r_{ij} \cdot \nabla_{i} W_{ij}}{\left|r_{ij}\right|^{2} + \eta^{2}}\right)^{-1}$$
(17)

Using Equation (17) for the pressure of each particle, one can calculate  $\hat{V}$  according to Equations (15) and (5) as

$$\widehat{V_i} = -\Delta t \sum_{i} m_j \left( \frac{P_i}{\rho_i^{*2}} + \frac{P_j}{\rho_j^{2}} \right) \nabla_i W_{ij}$$
 (18)

Finally, the velocity of each particle at the end of time-step will be obtained as:

$$V_{t+\Delta t} = V^{**} + \widehat{V} \tag{19}$$

and the final position of particles are calculated using a central difference scheme in time.

$$r_{t+\Delta t} = r_t + \frac{\Delta t}{2} \left( V_{t+\Delta t} + V_t \right) \tag{20}$$

This completes the computations required for one time-step. The procedure should be repeated for every other time-step till a desired time is reached

## 5. TREATMENT OF FREE SURFACE AND WALL BOUNDARY CONDITION

No special treatment was applied on free surface particles in the computational domain. In fact in this SPH method free surface is modeled naturally and this is one of the main advantages of the method.

Generally, the solid walls are also simulated by particles, which balance the pressure of inner fluid particles and prevent them from penetrating the wall. Special provisions must be made at the edges of these wall particles in order for the motion of the adjacent fluid to be modeled correctly. The wall boundary conditions can be modeled either by fixed particles exerting a repulsive force on inner fluid particles through a force function [21], or by image particles that

mirror the physical properties of inner fluid particles, so that boundary conditions are strictly satisfied along the solid walls [18].

Here we follow the treatment used by Koshizuka et al. [22] to model the wall boundaries by fixed wall particles, which are spaced according to the initial configuration. The Poisson Equation (16) is solved on these wall particles to repulse the inner flow particles accumulating in the vicinity of the wall. In this sense, the wall particles are dependent on the inner flows. For example, the pressure of wall particles increases when the particle density in the vicinity of the wall increases and thus the inner fluid particles are repelled from the wall, and vice versa.

#### 6. MODEL VERIFICATION

In order to validate of the present model, a dam-break problem is solved as a benchmark problem and the computational results are compared with other numerical works. A schematic of the dam break model is presented in Figure 1. In this model, the tank is L=3.22m long and H=2.0m high. A column of water  $(L_w=1.2m$  and  $H_w=0.6m)$  is located in the left side of the tank.

Figure 2 presents some snapshots of the flow at different times. The computed results are compared with the numerical results in which the standard SPH method is used to simulate the global development of fluid flow in dam-break problem [14].

At time t = 0s (not shown in the figure) the water column is allowed to flow. A relatively high velocity and shallow water depth flow in the x direction quickly forms (e.g. t = 1.19s). As time progresses, the flow impacts on the vertical wall at the opposite side of the tank. An upward water jet is suddenly formed that rises until gravity overcomes the upward momentum (around t = 1.19s). At this moment, the jet becomes thicker and the flow starts to reverse. Due to the oncoming flow, an adverse momentum gradient is created that results in an overturning wave (around t = 1.415s). This wave formation continues until the wave tip reconnects with the incident shallow water flow that now has less forward momentum. (Before t = 1.526s).

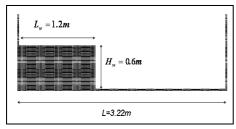
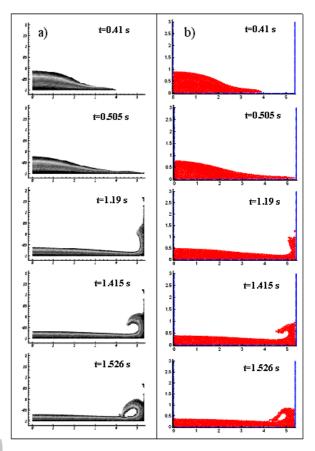


Figure 1. General layout of dam-break problem



**Figure 2.** Dam-break flow and impact against a vertical wall. a) Calculation by using standard SPH [14] b) Simulation by present incompressible SPH method

A good agreement in free surface shapes is observed by comparing between the results of standard SPH and present incompressible SPH method.

#### 7. THE WAVE PROPAGATION PROBLEM

In this section, the presented incompressible SPH model is employed for simulating waves generated by a wave-maker to reproduce the experimental data and previous numerical results.

**7.1. Introduction of Experiment** The detailed laboratory experiment can be found in Cox and Ortega [1]. In this experiment the wave flume was 36.0m long by 0.95m wide and 1.5 m high and equipped with a programmable wavemaker. The constant water depth in the flume was 65.0cm. The wavemaker signals consisted of two cycles of period t=1.0s sinusoidal waves followed by two and half cycles of period t=1.5s sinusoidal waves with a higher amplitude. The wave flume was long enough to prevent the reflections from

the far wall. The free surface positions were measured to cover the area of interest in the experiment.

**7. 2. Computational Parameters** Following the experimental setting of Cox and Ortega [1], the numerical tank is taken to be 17.3 m long and 1.2 m high. The length of numerical tank is shorter than the experimental one due to the constraint of the CPU time. The initial water depth is 0.65 m. The numerical wave tank used in this paper is shown in Figure 3.

The incident wave is generated by moving a numerical wave paddle located at the offshore boundary. The numerical wave signals are similar to those used by Gomez-Gesteira et al. [16].

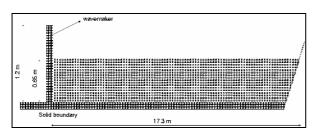
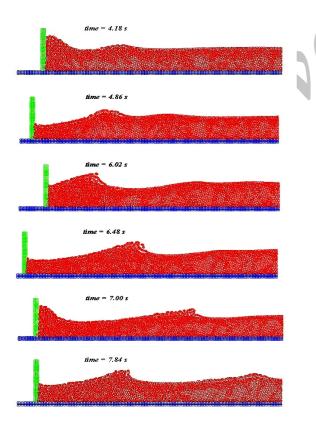
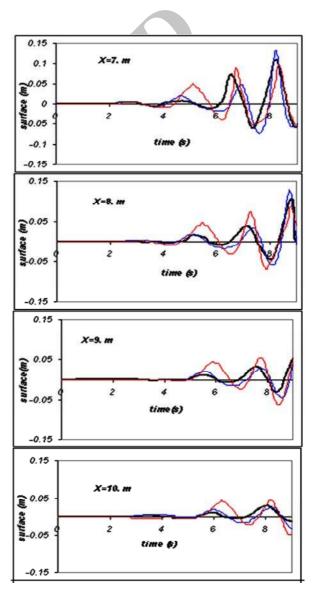


Figure 3. Numerical wave tank



**Figure 4.** Particle configurations during wave propagation computed by the present method at different times.

**7. 3. Discussion of Results** In this section, particle configuration due to the movement of the wavemaker is presented in Figure 4 at different times. In this computation, fluid particles are initially placed on a regular and equally spaced Cartesian grid with dx = dy = 0.04 m and thus approximately 9000 particles are used in the simulation. This number of particles ensures the accuracy of computations with respect to reduce the CPU time. A constant time step (dt = 0.00001s) and an initial particle velocity and pressure (V = P = 0) are used in the computations. The main computations using these quantities can be finished within 6 hours by a CPU 1.87GHz and RAM 1.0GB PC.



**Figure 5.** Free surface measurements. Comparison between presented incompressible SPH method (heavy black line), experimental results (blue line) and previous weakly compressible SPH method (red line).

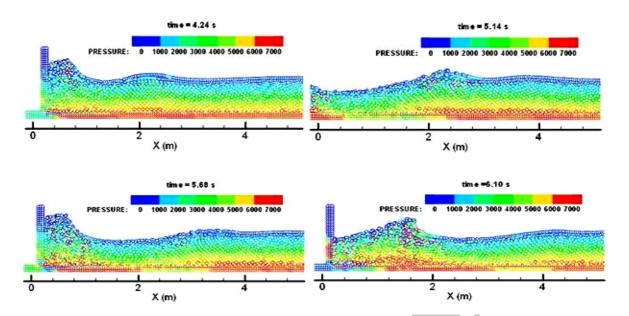


Figure 6. Pressure field of fluid during wave propagation at different times

The generated waves move toward the right wall corresponding to the wavemaker movement. The figure illustrates the stability of the present method to simulate the problems for long times.

In order to compare our results to the experimental ones [1] and numerical results in which the weakly compressible SPH is used to simulate propagation of waves in the tank [16], the numerical free surface position at different times is shown in Figure 5. The four frames correspond to the different positions from x=7.0 to 10.0 m, measured from the wavemaker. The numerical signal is observed to be in good agreement with the experimental one, both in phase and amplitude, although there are several slight discrepancies between the experimental and numerical profiles.

This is probably due to the fact that the numerical signal was filtered using a low pass filter, in such a way that sharp peaks are considerably smoothed out. Moreover, it is evident that the results provided by the current presented SPH algorithm shows more compatible results with experiments than those reported by weakly compressible SPH [16] especially in the case of amplitude.

It is noteworthy that in the previous simulations using weakly compressible SPH a tendency to form clump has been reported. However, investigation of Hosseini et al. [8] illustrates that the three step SPH algorithm shows no tendency to particle clustering. Although this numerical instability is insignificant in fluid simulations, it can result in an artificial surface tension. This characteristic prevents forming high

curvature regions because of large density discontinuities.

In Figure 6 the computed pressure field at different times is presented. These pressure contours show that the pressure of free surface particles is about zero and grows gradually with increase in water depth, hence the maximum pressure of fluid (about 7000 pa) belongs to the beneath particles.

#### 8. CONCLUSION

In this paper, the application of an incompressible SPH method for numerical simulation of free surface problems e.g. gravity wave propagation in free surface flows is investigated. One of the main differences between the utilized algorithm and the previous standard SPH lies in the pressure computation. By employing a new form of source term to the Poisson equation of pressure and using a three-step fractional algorithm, accuracy of this method in comparison with weakly standard SPH method is improved whereas, no special treatment is applied on free surface particles in the computational domain. In fact, by using the SPH form of the continuity equation (Equation (13)) instead of the density summation interpolant, which is common in SPH simulation, as well as using the Poisson equation, sensitivity of the method to density discontinuities is decreased. The computational results were in good agreement with the experimental data and previous numerical results.

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# Simulation of Gravity Wave Propagation in Free Surface Flows by an Incompressible SPH Algorithm

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Keywords: Smoothed Particle Hydrodynamics (SPH) Lagrangian Method Wave Propagation Free Surface این مقاله، یک مدل تراکمناپذیر هیدرودینامیک ذرات هموار (SPH) را برای شبیهسازی انتشار موج در جریان سطح آزاد ارائه می دهد. معادلات ناویر- استوکس، به صورت لاگرانژی با استفاده از یک الگوریتم سه مرحلهای حل می شوند. در گام اول، با استفاده از نیروهای حجمی، یک میدان سرعت میانی به دست می آید. این میدان سرعت در گام دوم، با اعمال اثرات لزجت، تجدید می شود. در گام سوم، یک معادله پواسون به عنوان جایگزین معادله حالت برای محاسبه فشار به کارگرفته می شود. این معادله پواسون، موازنه ای را بین چگالی و فشار ایجاد می کند که از آن در گام سوم برای اعمال تراکم ناپذیری استفاده می شود. نتایج محاسبات با نتایج عددی و آزمایشگاهی دیگر مقایسه شده و توافق خوبی مشاهده شده است. به منظور بررسی صحت روش ارائه شده، مسأله شکسته شدن سد، به عنوان یک مسأله پایه ای حل شده و نتایج به دست آمده با نتایج حاصل از روش های عددی دیگر مقایسه شده است.

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