# A combined level set/ghost cell immersed boundary representation for floating body simulations 

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## SUMMARY

A six degrees of freedom (6DOF) algorithm is implemented in the open-source CFD code REEF3D. The model solves the incompressible Navier-Stokes equations. Complex free surface dynamics are modeled with the level set method based on a two-phase flow approach. The convection terms of the velocities and the level set method are treated with a high-order WENO discretization scheme. Together with the level set method for the free surface capturing, this algorithm can model the movement of rigid floating bodies and their interaction with the fluid. The 6DOF algorithm is implemented on a fixed grid. The solid-fluid interface is represented with a combination of the level set method and ghost cell immersed boundary method. As a result, re-meshing or overset grids are not necessary. The capability, accuracy and numerical stability of the new algorithm is shown through benchmark applications for the fluid-body interaction problem. Copyright (c) 2015 John Wiley \& Sons, Ltd.

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## 1. INTRODUCTION

Coupled fluid-structure interaction plays a major role in many engineering disciplines. In the fields of coastal, ocean and arctic engineering, fluid-structure interaction problems occur in the presence of a free water surface and result in complex floating body dynamics. Some examples are floating piers, floating oil platforms, ship motion prediction or floating ice floes. Advanced solutions for the hydrodynamics, which include more complex free surface phenomena such as wave breaking or water jets and viscous effects from turbulence, require the solution of the Navier-Stokes equations. In earlier studies, fluid-structure interaction problems based on the Navier-Stokes equations have been calculated with Arbitrary Lagrangian-Eulerian (ALE) methods ([25], [31]). Then the interface location between the solid and the fluid is tied to the numerical mesh. When the location of this interface changes, the numerical mesh needs to be adjusted in order to account for this. The remeshing procedure can have a detrimental effect on the numerical accuracy and stability, especially for more arbitrary solid body movements. Also, for free surface flows, this method can become prohibitively complex.

A way to avoid constant re-meshing is the usage of dynamic overset grids. The method consists of a base mesh, which covers the full flow domain. The overset mesh is placed in the vicinity of the solid structures and overlaps with the base mesh. Then the overset mesh follows the movement of the solid. Carrica et al. [9] presented a complete framework for the modeling of ship motion

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using dynamic overset grids. The free surface was treated with a single-phase level set method. The dynamic overset mesh approach has numerical stability advantages over ALE, as constant remeshing of the solution domain is not necessary. The challenge of the method lies in an accurate, yet stable scheme for establishing the connections between the overset mesh points and the underlying grid points in the overlapping region. Later, Yang and Stern [33] showed an extension of the dynamic overset mesh approach, where the solid-fluid interface was treated with a sharp interface method instead of overset grids. The free surface was treated with a two-phase level set method. A further iteration of the dynamic overset mesh approach featured a direct forcing immersed boundary method for the fluid-solid interface [34]. Special attention was given to the field extension method, which was earlier presented for one-phase fluid-structure interaction [32]. When the solid moves through the flow field, solid cells becomes fluid cells and vice versa. With the field extension, unphysical values for the pressure and the velocities are avoided, through interpolation and keeping the physical pressure gradients intact. More recently, Calderer et al. [8] presented a level set based two-phase flow solver for the simulation of floating structures. A curvilinear immersed boundary method was used for the fluid-solid interface [6]. Chen et al. [10] presented a strong, two-way, fluid-solid coupling algorithm, based on a variational-type cut cell methodology, implemented within a hybrid Eulerian-Lagrangian framework.

Smoothed Particle Hydrodynamics (SPH) methods have also been used to calculate fluid-floating structure interaction by Bouscasse et al. [7] using a ghost fluid method to enforce the solid-fluid boundary for two-dimensional problems. Omidvar et al. [22] used an SPH implementation for 3D floating bodies using variable mass distribution to make the model computationally efficient. A large number of particles are required to accurately simulate the hydrodynamics of floating bodies using these methods. Glowinski et al. [15] presented a Lagrange multiplier based fictitious domain decomposition method for the simulation of an airfoil and sedimentation of circular particles. Sueyoshi et al. [27] simulated wave induced nonlinear motions of a two-dimensional box-shaped floating body using a moving particle semi-implicit (MPS) method but had difficulties with shorter period waves in sway and heave motions and a phase shift in the wave elevation.

In the present manuscript the open-source CFD code REEF3D is used. The model has been used extensively for complex wave hydrodynamics problems in the field of coastal and ocean engineering such as breaking waves [2] [1], wave energy converter devices [21], non-breaking wave forces [20] and breaking wave forces [5]. A novel approach for the geometry description for the 6DOF (six degree of freedom) algorithm is proposed. With the combined use of triangular surface meshes and the level set method, the fluid force and momentum can be integrated in a straightforward manner. The solid body is immersed into the fluid and re-meshing or overset grids are avoided. This is achieved with the local directional immersed boundary [4] in this study. The presented results are all obtained with a weakly coupled scheme. In combination with an already robust two-phase flow solver, this results in a stable fluid-structure interaction model. Numerical results for a disc entry problem, a free falling wedge and roll decay and roll motion of a rectangular barge under the influence of waves are presented.

## 2. NUMERICAL MODEL

The governing equations of the numerical model are the continuity and the incompressible NavierStokes (NS) equations presented in compact tensor notation:

$$
\begin{gather*}
\frac{\partial u_{i}}{\partial x_{i}}=0  \tag{1}\\
\frac{\partial u_{i}}{\partial t}+u_{j} \frac{\partial u_{i}}{\partial x_{j}}=-\frac{1}{\rho} \frac{\partial p}{\partial x_{i}}+\frac{\partial}{\partial x_{j}}\left[\nu\left(\frac{\partial u_{i}}{\partial x_{j}}+\frac{\partial u_{j}}{\partial x_{i}}\right)\right]+g_{i} \tag{2}
\end{gather*}
$$

where $u$ is the velocity averaged over time $t, \rho$ is the fluid density, $p$ is the pressure, $\nu$ is the kinematic viscosity and $g$ the acceleration due to gravity.

A Cartesian grid is used to discretize the spatial domain, providing easy implementation of higher-order discretization schemes. The model employs a staggered numerical grid to ensure
better pressure-velocity coupling and avoid numerical instabilities. The convective terms of the NS equations are discretized with a fifth-order WENO (weighted essentially non-oscillatory) scheme by Jiang and Shu [18] in the conservative finite-difference framework. The conservative WENO scheme is used to treat the convective terms for the velocities $u_{i}$, while a Hamilton-Jacobi version [17] is used for the variables of the free surface algorithm. The Hamilton-Jacobi version approximates the spatial derivatives at integer grid points rather than at half-integer points in the conservative version and is more suitable for the approximation of the gradients in the level set function.

For the time treatment a second-order accurate TVD Runge-Kutta scheme is employed, solving the Poisson equation two times per full time-step consisting of two Euler steps [26]:

$$
\begin{align*}
\Phi^{(1)} & =\Phi^{n}+\Delta t L\left(\Phi^{n}\right) \\
\Phi^{(n+1)} & =\frac{1}{2} \Phi^{n}+\frac{1}{2} \Phi^{(1)}+\frac{1}{2} \Delta t L\left(\Phi^{(1)}\right) \tag{3}
\end{align*}
$$

Adaptive time stepping is used in order to control the time steps and ensure numerical stability by maintaining the required CFL number. Chorin's projection method for incompressible flow which is first-order accurate, is used for the treatment of the pressure [12]. During the solution of the NS equations at each Euler step of the Runge-Kutta time stepping procedure, the pressure gradient is excluded. Based on the resulting divergence of the flow, a Poisson equation for the pressure is formed with the right hand side containing the continuity defect $\frac{\partial u_{i}^{*}}{\partial x_{i}}$ in Eq. 4, where $u_{i}^{*}$ is the transient velocity. The Poisson equation is solved using the fully parallelized Jacobi-preconditioned BiCGStab algorithm [30]. The gradient of the new pressure field is then used to correct the velocity field, making it divergence free.

$$
\begin{equation*}
-\frac{\partial}{\partial x_{i}}\left(\frac{1}{\rho\left(\Phi^{n+1}\right)} \frac{\partial p}{\partial x_{i}}\right)=-\frac{1}{\Delta t} \frac{\partial u_{i}^{*}}{\partial x_{i}} \tag{4}
\end{equation*}
$$

The location of the free water surface is represented implicitly by the zero level set of the smooth signed distance function $\Phi(\vec{x}, t)$ [23]. The level set function gives the closest distance to the interface and the two phases are distinguished by the change of the sign. This results in the following properties:

$$
\Phi(\vec{x}, t)\left\{\begin{array}{l}
>0 \text { if } \vec{x} \in \text { phase } 1  \tag{5}\\
=0 \text { if } \vec{x} \in \Gamma \\
<0 \text { if } \vec{x} \in \text { phase } 2
\end{array}\right.
$$

In addition, the Eikonal equation $|\nabla \phi|=1$ is valid. When the interface $\Gamma$ is moved under an externally generated velocity field $\vec{v}$, a convection equation for the level set function is obtained:

$$
\begin{equation*}
\frac{\partial \Phi}{\partial t}+u_{j} \frac{\partial \Phi}{\partial x_{j}}=0 \tag{6}
\end{equation*}
$$

When the interface evolves, the level set function loses its signed distance property. In order to maintain this property and to ensure mass conservation, the level set function is initialized after each time step. In the present paper a PDE based reinitialization equation is solved [28]:

$$
\begin{equation*}
\frac{\partial \Phi}{\partial t_{p}}+S(\Phi)(|\nabla \Phi|-1)=0 \tag{7}
\end{equation*}
$$

where $t_{p}$ is psuedo time and $S(\Phi)$ is the smoothed sign function [24]. The material properties of the two phases can be then be determined for the whole domain. Close to the interface, the density $\rho$ and the viscosity $\nu$ are smoothed out by calculating the density at the cell face with a regularized Heaviside function with an interface thickness of $\epsilon=2.1 d x$ :

$$
\begin{equation*}
\rho_{i+\frac{1}{2}}=\rho_{1} H\left(\Phi_{i+\frac{1}{2}}\right)+\rho_{2}\left(1-H\left(\Phi_{i+\frac{1}{2}}\right)\right) \tag{8}
\end{equation*}
$$

The level set function at the cell face is obtained using:

$$
\begin{equation*}
\Phi_{i+\frac{1}{2}}=\frac{1}{2}\left(\Phi_{i}+\Phi_{i+1}\right) \tag{9}
\end{equation*}
$$

## 3. 6 DOF ALGORITHM

For the calculation of fluid-structure interaction, the geometry of the solid body needs to be defined. In the current model, it is described by a primitive triangular surface mesh neglecting connectivity. By design, this resembles the structure of the STL format, a standard for surface meshes, available in most CAD and meshing softwares [13].


Figure 1. Representation of the solid body with a level set function.

The intersections of the surface mesh with the underlying Cartesian grid are determined with a ray-tracing algorithm [35]. Here, inside-outside information and the shortest distance to the closest triangle for the grid points along the coordinate axis can be calculated in a very efficient and reliable manner. After this step, the standard reinitialization algorithm [24] is used to get signed distance properties for the level set function in the vicinity of the solid body, see Fig. 1. In current literature, the level set method has been employed to represent moving solid bodies by [11] and [3]. It is important to mention, that the ray-tracing algorithm calculates the distances in an exact manner close to the solid boundary, which results in a sharp representation of the solid-fluid interface. The moving solid-fluid interface is treated with the ghost cell immersed boundary method [4]. The forces acting on the surface $\Omega$ of the floating body can be determined for each co-ordinate direction $i$ separately using the pressure $p$ and the viscous stress tensor $\tau$ in the following way:

$$
\begin{equation*}
F_{i, e}=\int_{\Omega}\left(-\mathbf{n}_{i} p+\mathbf{n}_{i} \cdot \tau\right) d \Omega \tag{10}
\end{equation*}
$$

The distance of the center of gravity from the origin of the body-fitted grid can be determined using the following equation 11 :

$$
\begin{equation*}
\mathbf{r}_{c g}=\frac{1}{m} \int_{V} \mathbf{r} \rho_{a} d V \tag{11}
\end{equation*}
$$

where $\mathbf{r}$ is the distance from each surface cell to the origin of the body-fitted co-ordinate system. Assuming that the origin of the body-fitted co-ordinate system is at the center of gravity of the floating body, $\mathbf{r}$ is the distance of each surface cell to the center of gravity and the moments are determined as follows:

$$
\begin{equation*}
L_{i, e}=\int_{\Omega} \mathbf{r} \times\left(-\mathbf{n}_{i} p+\mathbf{n}_{i} \cdot \tau\right) d \Omega \tag{12}
\end{equation*}
$$

Through the level set representation of the solid surface, the calculation of the discrete surface area in each grid cell can be accomplished with the help of a Dirac delta function [24]:

$$
\begin{equation*}
d \Omega=\int \delta(\Phi)|\nabla \Phi| d x \tag{13}
\end{equation*}
$$

The advantage of the level set description for the surface area of the solid body is, that the intersections of the surface mesh with the underlying grid do not need to be calculated explicitly. The implementation of the level set function was relatively easy, as all numerical routines are already available from the free surface algorithm and only small adjustments were required.

A free floating body has six degrees of freedom. The translation consists of the three linear velocities $u, v$, and $w$. The rotation has the three angular velocities $p, q$, and $r$ [14]. The location and the orientation of the floating body are given by the position vector and the Euler angles:

$$
\begin{equation*}
\boldsymbol{\eta}=\left(\boldsymbol{\eta}_{1}, \boldsymbol{\eta}_{2}\right)=\left(x_{c g}, y_{c g}, z_{c g}, \phi, \theta, \psi\right) \tag{14}
\end{equation*}
$$

The calculation of the six degrees of freedom for the solid body can be simplified with respect to the moments of inertia by introducing two separate coordinate systems. The fluid flow is calculated in the inertial coordinate system, and the floating body in the non-inertial coordinate system. Then the forces $X, Y$ and $Z$ and moments $K, M$ and $N$ acting on the body can be calculated in the inertial coordinate system. When the origin of the non-inertial coordinate system coincides with the center of gravity, the moments of inertia can be calculated by considering only the main diagonal of the moment of inertia tensor:

$$
\mathbf{I}=\left[\begin{array}{ccc}
I_{x} & 0 & 0  \tag{15}\\
0 & I_{y} & 0 \\
0 & 0 & I_{z}
\end{array}\right]=\left[\begin{array}{ccc}
m r_{x}^{2} & 0 & 0 \\
0 & m r_{y}^{2} & 0 \\
0 & 0 & m r_{z}^{2}
\end{array}\right]
$$

where $r_{x}, r_{y}$ and $r_{z}$ are the distances of a point from the center of gravity along the $x-, y-$ and $z-$ directions.

The calculated forces and moments from the inertial reference frame can be expressed in the noninertial coordinate system with a rotation matrix $\mathbf{J}_{1}^{-1}$, consisting of three elemental rotations around the axis of the coordinate system ( $s$ stands for $\sin$ and $c$ for $\cos$ ):

$$
\boldsymbol{a}_{f b}=\left[\begin{array}{ccc}
c \psi c \theta & s \psi c \theta & -s \theta  \tag{16}\\
-s \psi c \phi+s \phi s \theta c \psi & c \psi c \phi+s \phi s \theta s \psi & s \phi c \theta \\
s \theta s \psi+c \phi s \theta c \psi & -s \phi c \psi+c \phi s \theta s \psi & c \theta c \phi
\end{array}\right] \boldsymbol{a}_{e}=\mathbf{J}_{1}^{-1} \boldsymbol{a}_{e}
$$

Here $\boldsymbol{a}_{f b}$ is a vector in the reference frame of the floating body, and $\boldsymbol{a}_{e}$ a vector in the inertial coordinate system. With the calculation of the forces, momentum and moments of inertia in place, the dynamic rigid body equations can be solved [9]:

$$
\left.\begin{array}{rl}
F_{i} & =\mathbf{J}_{1}^{-1} F_{i, e}  \tag{17}\\
L_{i} & =\mathbf{J}_{1}^{-1} L_{i, e}
\end{array}=[K, Y, Z], N\right]
$$

where

$$
\begin{align*}
{[m(\dot{u}-v r+w q)] } & =X \\
{[m(\dot{v}-w p+u r)] } & =Y \\
{[m(\dot{w}-u q+v p)] } & =Z \\
{\left[I_{x} \dot{p}+\left(I_{z}-I_{y}\right) q r\right] } & =K  \tag{18}\\
{\left[I_{y} \dot{q}+\left(I_{x}-I_{z}\right) r p\right] } & =M \\
{\left[I_{z} \dot{r}+\left(I_{y}-I_{x}\right) p q\right] } & =N
\end{align*}
$$

Here $u, v, w, p, q$ and $r$ are the values for the linear and angular velocities from the previous time step. Then $\dot{u}, \dot{v}, \dot{w}, \dot{p}, \dot{q}$ and $\dot{r}$ can be calculated in an explicit manner. Any of the linear and angular
velocities $\dot{\varphi}$ and any component of the position and orientation vector $\varphi$ of the floating body can be calculated with a second-order Adams-Bashforth scheme for the new time step:

$$
\begin{align*}
& \dot{\varphi}^{n+1}=\dot{\varphi}^{n}+\frac{\Delta t}{2}\left(3 \ddot{\varphi}^{n+1}-\ddot{\varphi}^{n}\right)  \tag{19}\\
& \varphi^{n+1}=\varphi^{n}+\frac{\Delta t}{2}\left(3 \dot{\varphi}^{n+1}-\dot{\varphi}^{n}\right)
\end{align*}
$$

As a result, the floating body dynamics are solved in a fully explicit way. Even though the weak coupling between the 6DOF algorithm and the flow solver has been reported to lead to numerical stability problems for complex cases (e.g. [9] or [8]), the current implementation shows good numerical stability throughout the range of applications.
The dynamic rigid body equations have been solved in the floating body reference frame. The translations and and orientations are also calculated there. They are transformed to the inertial reference frame with the matrix given in Eq. 16. The angular velocities are transformed using the following rotation matrix $\mathbf{J}_{2}$ [14] where $s, c$ and $t$ stand for $\sin$, cos and tan respectively:

$$
\dot{\boldsymbol{\eta}}_{2}=\left[\begin{array}{ccc}
1 & s \phi t \theta & c \phi t \theta  \tag{20}\\
0 & c \phi & -s \phi \\
0 & s \phi / c \theta & c \phi / c \theta
\end{array}\right] \boldsymbol{v}_{2}=\mathbf{J}_{2} \boldsymbol{v}_{2}
$$

Boundary conditions for the velocities on the solid-fluid interface result from the motion of the solid body with respect to its center of gravity.

$$
\begin{equation*}
u_{i}=\dot{\boldsymbol{\eta}}_{\mathbf{1}}+\dot{\boldsymbol{\eta}}_{\mathbf{2}} \times \mathbf{r} \tag{21}
\end{equation*}
$$

Pressure oscillations in the vicinity of the solid body can occur due to solid cells turning into fluid cells and vice versa. This is avoided by the implementing the field extension method [29][32] adapted to the ghost cell immersed boundary method. For non-moving boundaries, a zero-gradient boundary condition is used for the pressure. In order to maintain a physical pressure gradient near the floating body, the following boundary condition for the gradient of the pressure is given:

$$
\begin{equation*}
\frac{\partial p}{\partial x_{i}}=-\frac{1}{\rho} \frac{D u_{i}}{D t} \tag{22}
\end{equation*}
$$

In oder to determine $u_{i}$, one possibility is to evaluate the momentum equations. A simpler way is to differentiate Eq. 21 with respect to time and use this for the ghost cell values for the pressure:

$$
\begin{equation*}
\frac{D u_{i}}{D t}=\frac{d}{d t}\left(\dot{\boldsymbol{\eta}}_{\mathbf{1}}+\dot{\boldsymbol{\eta}}_{\mathbf{2}} \times \mathbf{r}\right) \tag{23}
\end{equation*}
$$

When cells are freshly cleared by the moving body and become fluid cells, unphysical values for the velocities and the pressure in those cells can cause numerical stability problems. This is avoided by assigning the velocities the values from Eq. 21 and the value for the pressure is found through interpolation from the fluid.

## 4. RESULTS

### 4.1. Disc Entry

At first the 6DOF algorithm is tested for the well-known disc entry problem. This case has been used by several authors to benchmark their models (e.g. [33] or [8]). The parameters for the geometry, the initial boundary conditions and flow parameters of this test are non-dimensional. The disc enters the water with the fixed vertical velocity of $V=-1$. The center of the disc is located at $H=1.25$ over the free surface at $T=V t / H=0$. The cylinder has a radius of $R=1$. The two-dimensional simulation domain has the size $30 R \times 22 R$. The uniform mesh size is $d x=0.025$, resulting in
$1200 \times 880$ grid cells. Acceleration due to gravity is $g=-1$, the density of the water is $\rho_{\text {water }}=1$, the density of the air is $\rho_{\text {air }}=0.001$, the viscosity of the water is $\nu_{\text {water }}=0.001$ and the viscosity of the air is $\nu_{\text {air }}=0.018$. The time step size is controlled with adaptive time stepping using a CFL number of 0.1 . No turbulence model is employed for this case.
The free surface location and contour for the velocity magnitude are shown in Fig. 2. The disc enters the liquid phase in Fig. 2a. In a symmetric fashion, breaking waves are generated from the solid impact at the sides of the disc and are traveling away from the disc. A fine enough mesh is necessary for capturing breaking waves, in general terms [2], but also specifically for the present case [8]. In Fig. 2b, the post-breaking waves are moving further away towards the side boundaries. At $T=3.0$, the displaced water has started to return to the location of the disc impact, and the solid body is now fully submerged in the liquid phase. At $T=4.0$, the returning water creates a vertical water jet. The free surface location compares well with other results reported in literature (e.g. [33] or [8]).


Figure 2. Disc entry problem, the contour shows the velocity magnitude.

### 4.2. Free Falling Wedge

In the previous section, the motion of the solid body was fixed. In this section, the free falling wedge has all degrees of freedom in the two-dimensional plane. The numerical setup is similar to
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Figure 3 . Free falling wedge, the contour shows the velocity magnitude.
Fig. 4a shows the vertical position of the center of gravity of the wedge, which is measured from the still water level. In the first 0.5 s of the free fall, the wedge moves downwards and hits the free surface, which can also be seen in Fig. 3d. Then the wedge moves into the body of water and starts to
oscillate around the still water level. Three different grid sizes are tested in 2D and one simulation is carried out in 3D. The result for the two finest grids are almost similar in 2D and comparable to the results in 3D. A marginal difference can be found for the grid with $d x=0.05 \mathrm{~m}$. Here, the wedge moves slightly deeper into the water after the impact than in the experiments and the fine grid results. The vertical velocity of the wedge is shown in Fig. 4b. In the free fall phase, the velocity is increasing linearly under acceleration due to gravity. After it reaches the peak velocity of $5.5 \mathrm{~m} / \mathrm{s}$, the wedge sharply decelerates due to the water impact. When the wedge starts to oscillate around the still water level, the vertical velocity becomes non-linear. The vertical velocity is modeled accurately on all three grids, with slight deviations on $d x=0.05 \mathrm{~m}$.

The result from the 3D simulation during water entry of the wedge at $t=0.8 \mathrm{~s}$ is presented in Fig. 5, showing the splash up of the water as the freely falling wedge impacts the free surface. This scenario involves large pressure gradients as the wedge impinges the fluid surface and typically requires strongly coupled schemes to solve the fluid structure interaction problem according to Calderer et al. [8]. In the current study, this complex fluid structure scenario is solved with a weakly coupled scheme with good results.


Figure 4. Motion of a wedge under free-fall


Figure 5.3D simulation of free falling wedge into water at $t=0.8 \mathrm{~s}$
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### 4.3. Roll motion of a rectangular barge

The case of roll motion of a rectangular barge in waves has been presented by Jung et al. [19]. They performed experiments in a wave flume with the barge fixed around the rotational direction axis in the center of gravity of the structure, resulting in a single degree of freedom. The barge is 0.3 m long, 0.1 m high and 0.9 m wide. It is made of solid acrylic glass with the density $\rho=1048 \mathrm{~kg} / \mathrm{m}^{3}$, resulting in a moment of inertia $I_{y}=0.236 \mathrm{~kg} \mathrm{~m}^{2}$ for the direction of the roll motion. The wave flume has a water depth of $h=0.9 \mathrm{~m}$. The center of gravity of the barge coincides with the still water free surface location.


Figure 6. Roll motion of a rectangular barge.

(a) $\mathrm{t} / \mathrm{T}=0.0$

(c) $\mathrm{t} / \mathrm{T}=0.5$

(b) $t / T=0.25$

(d) $t / T=0.75$

Figure 7. Roll motion of the rectangular barge, the contour shows the vertical velocity.

For the roll decay test, the barge is initially tilted to an angle of $15^{\circ}$ and then released. In the experiments, the roll angle is recorded with a rotary position sensor. In the numerical simulations the case is treated as two-dimensional, as it is symmetric along the width. The numerical domain for the roll decay test is 5 m long and 1.6 m high, the inflow and outflow boundaries are modeled as walls. A uniform mesh size of $d x=0.01 \mathrm{~m}$ is used, resulting in a moment of inertia of $I_{y}=0.0026 \mathrm{~kg} \mathrm{~m}^{2}$ for the 0.01 m wide barge. Following [8], the damping coefficient of $C=0.275$ for the threedimensional case is used in the rigid body equations and adjusted to the two-dimensional setup, in order to account for the friction of the experimental apparatus. No turbulence model is used. Fig. 6a shows the comparison between the experimental and numerical results for the roll decay test. For the acrylic glass barge, the calculated roll angle $\phi$ for $d x=0.01 \mathrm{~m}$ agrees well with measurements up to around 3 s . Small differences are observed for the roll decay on the coarser meshes. After that, the numerical results go out of phase.

The roll motion of the rectangular barge under the influence of regular periodic waves is shown in Fig. 6b. Waves with a length of $L=2.2 \mathrm{~m}$, a wave period of $T=1.2 \mathrm{~s}$ and a height $H=0.06 \mathrm{~m}$ are generated in the numerical model with the relaxation method [16][2]. The simulation domain is $L=12 \mathrm{~m}$ long, the wave generation zone is 2.2 m long and the numerical beach is 4.4 m long. Fig. 6 b shows the roll angle. When the first full wave has propagated to the barge from the generation zone, the roll angle shows two maximum peaks in the numerical results. After that, the roll angle maintains a constant amplitude and the measured and computed values from the finest grid show a good match. In Fig. 7, the rectangular barge is shown at different stages of a wave cycle. The contour shows the vertical velocity.

## 5. CONCLUSIONS

The implementation of a 6 DOF algorithm in the open-source CFD code REEF3D was shown. In a novel approach, the floating body is represented by the combination of a surface mesh, a level set function and the ghost cell immersed boundary method. This results in a method, that does not require re-meshing or overset grids. The level set method for the description of the surface area of the floating body has the advantage, that the forces and moments can be calculated without explicitly defining the intersections between the surface mesh and the grid of the flow domain. In addition, the numerical model uses the level set method for the capturing of the interface between water and air. The resulting model proved to be numerically stable and all simulations were performed with a weakly coupled fluid-structure interaction scheme. The three benchmark cases disc entry, free falling wedge and roll motion of a rectangular were calculated successfully and a high level of detail and accuracy was achieved. With the new 6DOF algorithm many interesting problems in coastal, ocean and arctic engineering can be investigated in the future.

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