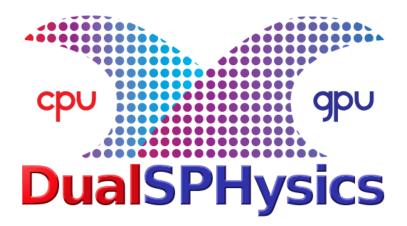
XML GUIDE FOR DUALSPHYSICS

Create your own case using the XML file



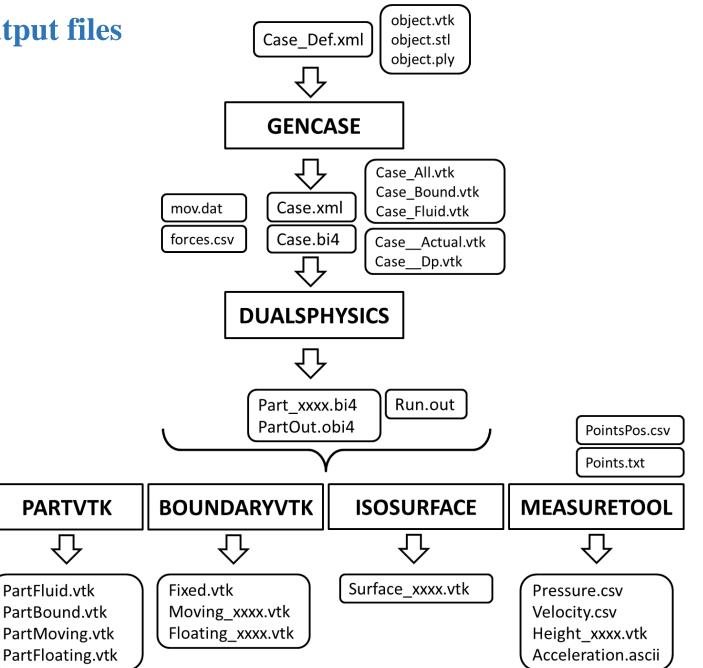
April 2016

DualSPHysics team

Input & output files

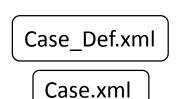
Pre-Processing





2

Input & output files: Format files



XML File

- The eXtensible Markup Language is textual data format compatible with any hardware and software.
- Information is structured and organised by using labels.
- They can be easily edited using any text editor.

Case.bi4

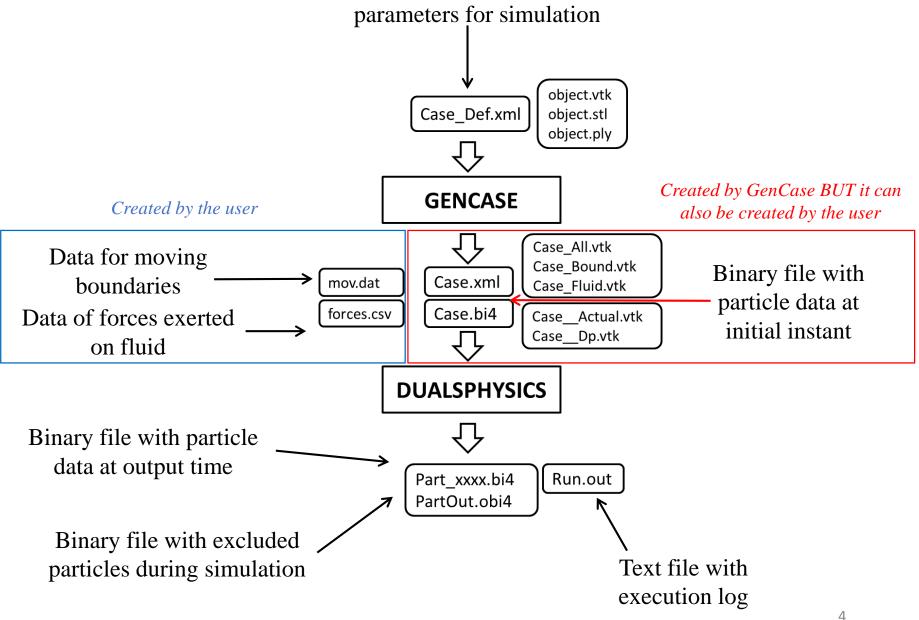
Part_xxxx.bi4 PartOut.obi4

BINARY File

- Binary format consumes six times less memory than text format.
- Reading or writing is several times faster using a binary format.
- A special code is required to read the data (JPartDataBi4.cpp/.h).
- ".bi4" is the new binary format that also includes double precision.
- The user can also define new arrays that post-processing tools can automatic manage.

Input & output files

Constants and configuration









co	case>
	<casedct> <constantsdef></constantsdef></casedct>
<pre>ccases> ccasedef></pre>	<pre>constantsantsants</pre>
<pre><coseuer> </coseuer></pre>	<pre><gravity comment="Gravitational acceleration" units_comment="m/s^2" x="0" y="0" z="-9.81"></gravity></pre>
<lattice bound="1" fluid="1"></lattice>	<cflnumber comment="Coefficient to multiply Dt" value="0.2"></cflnumber>
<pre><gravity comment="Gravitational acceleration" units_comment="m/s^2" x="0" y="0" z="-9.81"></gravity></pre>	<pre>chsvl value="0" auto="true" comment="Maximum still water level to calculate speedofsound using coefsound" units_comment="metres (m)" /></pre>
<pre><cfinumber comment="Coefficient to multiply Dt" value="0.2"></cfinumber></pre>	<pre>capeddystem value="0" att=""true" comment="Maximum system speed (by default the dam-break propagation is used)" /> foreframed maintering of the system is used in the system speed (by default the dam-break propagation is used)" /> </pre>
<pre>chsul value*"0" auto*"true" comment="Maximum still water level to calculate speedofsound using coeffound" units_comment="Maximum system speed (by default the dam-break propagation is used) ">></pre>	<pre><coefsound comment="Coefficient to multiply speedsystem" value="20"></coefsound> <pre><aprecisual auto="true" comment="Speed of sound to use in the simulation (by default speedspsud-coefsound*speedsystem)" value="0"></aprecisual></pre></pre>
<pre>cspecaystem value='o auto='true' comment='marinum system specu (by certaint the dam-break propagation is used)- /> ccoefsound value='2' coefficient to multiply specdystem' /> </pre>	<pre>csptetiound value = 0 also = true Comment= speed of sound to use in the simulation (p desail speed) sound-speed speed sound to use in the simulation (p desail speed) sound-speed speed speed sound to use in the simulation (p desail speed) sound-speed speed speed sound to use in the simulation (p desail speed) sound-speed speed sound to use in the simulation (p desail speed) sound-speed speed speed speed speed sound-speed speed sound-speed speed speed speed speed speed speed speed sound-speed speed speed spe</pre>
<pre><pre><aprecision: auto="true" comment="Speed of sound to use in the simulation (by default speedofsound=coefsound=speedsystem)" value="0"></aprecision:></pre></pre>	<pre><gamma comment="Politropic constant for water used in the state equation" value="7"></gamma></pre>
<pre><coefh comment="Coefficient to calculate the smoothing length (H=coefficient*sqrt(3*dp^2) in 3D)" value="1.0"></coefh></pre>	<rhop0 comment="Reference density of the fluid" units_comment="kg/m3" value="1000"></rhop0>
<gamma comment="Politropic constant for water used in the state equation" value*"7"=""></gamma>	
<pre>crhop value="1000" comment="Reference density of the fluid" units_comment="kg/m3" /> </pre>	<pre><mkconfig boundcount="240" fluidcount="10"></mkconfig></pre>
 <pre></pre>	<geometry></geometry>
<pre><geometry></geometry></pre>	<pre>cdefinition dp="0.01" units_comments"kettes (m)"></pre>
<definition dp="0.01" units_comment="metres (m)"></definition>	<pre><pre>cpointmain x=-1: y=0: z=-1: /> <pre><pre><pre><pre><pre><pre><pre><pre></pre></pre></pre></pre></pre></pre></pre></pre></pre></pre>
<pre><pointmin x="-1" y="0" z="-1"></pointmin></pre>	<pre></pre>
<pre><pre>cpcintmax x="4.5" y="0" z="3.5" /></pre></pre>	
	<pre><mainlist></mainlist></pre>
<pre><commands> </commands></pre>	<setdrawmode mode="full"></setdrawmode>
<pre><mainist> < setdrawmode mode**full* /></mainist></pre>	<setmkfluid mkm"0"=""></setmkfluid>
<pre><stcrawmode mode="rult=/"> <stcrawmode mode="rult=/"> <stcrawmode mode="rult=/"> </stcrawmode></stcrawmode></stcrawmode></pre>	<drawbox></drawbox>
	<boxfill>solid</boxfill>
<boxfill>molid</boxfill>	<pre><pre>cpoint x="0" y="-1" z="0" /></pre></pre>
<pre><pre>cpoint x="0" y="-1" z="0" /></pre></pre>	<pre><size x="1" y="2" z="2"></size></pre>
<pre><size x="1" y="2" z="2"></size></pre>	
	<pre><setmkbound mk="0"></setmkbound></pre>
<setakbound mk="0"></setakbound>	<pre><drawbox></drawbox></pre>
	<pre> <</br></br></pre>
<pre></pre>	<pre>cpoint x="0" y="-1" z="0" /> <cate x="4" y="2" z="3"></cate></pre>
<pre><pre><pre><pre><pre><pre><pre><pre></pre></pre></pre></pre></pre></pre></pre></pre>	<pre><dicale #~="" ''="" 4~="" y="_ 2="></dicale> </pre>
	<execution></execution>
<execution></execution>	<pre><pre>cparameters></pre></pre>
<pre><pre>cparameters></pre></pre>	<pre><parameter comment="Step Algorithm 1:Verlet, 2:Symplectic (default=1)" key="StepAlgorithm" value="1"></parameter></pre>
<pre>cprrameter key="StepAlgorithm" value="1" comment="Step Algorithm liverlat, 2;Symplectic (default=1)" /> cprrameter key="WeiteSteps" value="d" of comment="Weite only: Nuber of steps to apply Ruler timestepping (default=40)" /></pre>	<pre><parameter comment="Verlet only: Number of steps to apply Euler timestepping (default=40)" key="VerletSteps" value="40"></parameter></pre>
<pre>cprimeter key= force/comparise to comment="Interaction Kernel 1:Cubic Spline, 2:Wendland (default=2)" /></pre>	<pre><pre>state: key="Kernel" value="2" comment="Interaction Kernel 1:Cubic Spline, 2:Wendland (default=2)" /></pre></pre>
<pre><pre>cparameter key="ViscoTreatment" value="1" comment="Viscosity formulation 1:Artificial, 2:Laminar+SPS (default=1)" /></pre></pre>	<pre>cyparameter key="ViscoTreatment" value="1" comment="Viscosity formulation 1:Artificial, 2:Laminar+SPS (default=1)" /> </pre>
<pre><pre>cparameter key="Visco" value="0.02" comment="Viscosity value" /></pre></pre>	<pre>cparameter key=""visco" value="0.02" comment="Viscosity value" /> for a state of the state</pre>
<pre><parameter comment="Multiply viscosity value with boundary (default=1)" key="ViscoBoundFactor" value="1"></parameter></pre>	<pre>cparameter key="ViscoBoundFactor" value="1" comment="Multiply viscosity value with boundary (default=1)" /></pre>
<pre><pre>cparameter key="DeltaSPH" value="0" comment="DeltaSPH value, 0.1 is the typical value, with 0 disabled (default=0)" /></pre></pre>	<pre><pre>cparameter key" beatcaser "value" or comment="backaser value, of is the typical value, with o disabled (default=0)" /> <pre><pre></pre></pre></pre></pre>
<pre>cparameter key=#\$Shifting value="0" comment="Shifting mode 0:None, 1:Ignore bound, 2:Ignore fixed, 3:Full (default=0)" /> cparameter key=#\$Khifting " value="0" comment="Shifting mode 0:None, 1:Ignore bound, 2:Ignore fixed, 3:Full (default=0)" /></pre>	<pre>cparameter key="#ShiftCoof" value="-2" comment="Coofficient for shifting computation (default=2)" /></pre>
<pre>cparameter key="#shiftCoaf" values"-2" comment="Coafficient for shifting computation (default=2)" /> cparameter key="#shiftToaf" values"-1.5" comment="Threshold to defact free surface. Typically 1.5 for 2D and 2.75 for 3D (default=0)" /></pre>	<pre>cparameter key="#ShiftTFS" value="1.5" comment="Threshold to detect free surface. Typically 1.5 for 2D and 2.75 for 3D (default=0)" /></pre>
<pre>cparameter key="shifting value="in" comment="intended to detect the surface. typically in for 2D and 2.75 for aD (default=0)" /> cparameter key="shifting value="in" comment="shifting d Algorithm 1:5PR, 2:DBM (default=1)" /></pre>	<pre>cparameter key="RigidAlgorithm" value="1" comment="Rigid Algorithm 1:SPH, 2:DEM (default=1)" /></pre>
<pre>cyprimetry key="fifthuse" value" 1 Community fifth copyright copyright for (control of the simulation start (warmup) (default=0)" units_comment="seconds" /></pre>	<pre><pre><pre><pre>cparameter key="FtPause" value="0.0" comment="Time to freeze the floatings at simulation start (warmup) (default=0)" units_comment="seconds" /></pre></pre></pre></pre>
<pre><pre>cparameter key="CoefDtWin" value="0.05" comment="Coefficient to calculate minimum time step dtmin=coefdtmin=h/speedsound (default=0.05)" /></pre></pre>	<pre><parameter comment="Coefficient to calculate minimum time step dtmin=coefdtemin*h/speedsound (default=0.05)" key="CoefDtMin" value="0.05"></parameter></pre>
<pre><pre>cyarameter key="#DtIni" value="0.0001" comment="Initial time step (default=h/speedsound" units_comment="seconds" /></pre></pre>	<pre><pre>cparameter key="#DtIni" value="0.0001" comment="Initial time step (default=h/speedsound" units_comment="seconds" /></pre></pre>
<pre><pre><pre><pre>cparameter key="#DtMin" value="0.00001" comment="Minimum time step (default=coefdtmin*h/speedsound)" units_comment="seconds" /></pre></pre></pre></pre>	<pre><pre>cparameter key="#DtMin" value="0.00001" comment="Minimum time step (default=coefdtmin*h/speedsound)" units_comment="seconds" /></pre></pre>
<pre><pre><pre><cpre>parameter key="#DtFixed" value="DtFixed.dat" comment="Dt values are loaded from file (default=disabled)" /></cpre></pre></pre></pre>	<pre><parameter comment="Dt values are loaded from file (default=disabled)" key="#DtFixed" value="DtFixed.dat"></parameter></pre>
<pre><pre>cparameter key="DtAllParticles" value="0" comment="Velocity of particles used to calculate DT. 1:All, 0:Only fluid/floating (default=0)" /></pre></pre>	<pre><parameter comment="Velocity of particles used to calculate DT. 1:All, 0:Only fluid/floating (default=0)" key="DtAllParticles" value="0"></parameter></pre>
<pre>cprrameter key="TimeMax" value="0.12" comment="Time of simulation" units_comment="seconds" /> cprrameter key="timeMax" value="cprrameter" (cprrameter key="seconds" /> cprrameter key="timeMax" value="cprrameter key="seconds" key="seco</pre>	<pre>cparameter key="TimeMax" value="0.72" comment="Time of simulation" units_comment="seconds" /> </pre>
<pre>cparameter key="TimeOut" value="1.01" comment="Time out data" units_comment="seconds" /> cparameter key=TimeOut" units="TimeOut" forcesso of Z+" units_comment="decimal" /></pre>	<pre><pre>cparameter key="TimeOut" value="0.01" comment="Time out data" units_comment="seconds" /></pre></pre>
<pre>cparameter key="nincs" value="1" comment="infrease of i+" units_comment="accimant"/> cparameter key="PartisotUkax" value="1" comment="Allowed 4/100 of fluid particles out the domain (default=1)" units_comment="decimant" /></pre>	<pre>cparameter key="lno2" value="1" comment="increase of 5+" units comment="decimal" /> comment="Recordship" advantage of 5+" units comment="decimal" />> commen</pre>
<pre>cparameter key="knetsoutskk" value="10" comment="klowed #/loo riuta particules out the domain (derault=1)" units_comment="coefina" /> cparameter key="knetsoutskk" value="10" comment="klowed #/loo riuta (default=700)" units_comment="klowed" /> cparameter key="knetsoutskk" value="10" comment="klowed #/loo riuta (default=700)" units_comment="klowed" /> cparameter key="knetsoutskk" value="10" comment="klowed #/loo riuta particules out the domain (derault=1)" units_comment="klowed" /> cparameter key="klowed" /> comment="klowed #/loo riuta particules out the domain (derault=1)" units_comment="klowed" /> cparameter key="klowed" /> comment="klowed #/loo riuta particules out the domain (derault=1)" /> cparameter key="klowed" /> comment="klowed" /> comment="klo</pre>	<pre>cparameter key="PartsOutMax" value="1" comment="Allowed \$/100 of fluid particles out the domain (default=1)" units_comment="decimal" /> cparameter key="PartsOutMax" value="1" comment="Ministry comment="Min</pre>
<pre>cparameter key="hoppouthar" value="100" comment="Maximum thop valid (default=1300)" units_comment="key" //> </pre>	<pre>cparameter key="Rhopotkin" value="700" comment="Minimus rhop valid (default=700)" units_comment="kdyn3" /> cparameter key="Rhopotkin" value="100" comment="kdyna" valid (default=1300" units_comment="kdyn3" /> </pre>
	<pre>character Key="KnoboutMax" value= 100" comment= maximum thop value (dermate=100) units_comment= withs ''</pre>
executions</td <td><pre><pre>cparticles np="21001" nbf="1001" mkboundfirst="11" mkfluidfirst="1"></pre></pre></td>	<pre><pre>cparticles np="21001" nbf="1001" mkboundfirst="11" mkfluidfirst="1"></pre></pre>
	first mkbound*"O" mk*"11 begin*"O" count*"1001"/>
	<fluid begin="1001" count="20000" mk="1" mkfluid="0"></fluid>
	<constants></constants>
	<pre><gravity units_comment="m/s^2" x="0" y="0" z="-9.81"></gravity></pre>
	<cflnumber value="0.2"></cflnumber>
	<gamma value="7"></gamma>
	<pre><rhop0 units_comment="kg/m3" value="1000"></rhop0></pre>
	<pre><dp units_comment="metres (m)" value="0.01"></dp></pre>
	<h units_comment="metres (m)" value="1.41421356248-002"></h>
	<pre><b units_comment="metres (m)" value="1.11553714298+006"></pre>
	<pre>cmassbound value*1.0000000008-001* units_comment="kg" /> cmassfluid value*1.000000008-001* units_comment="kg" /></pre>
	<pre></pre> <pre><</pre>
· · · · · · · · · · · · · · · · · · ·	(motion />

</case>

STRUCTURE OF THE XML FILE

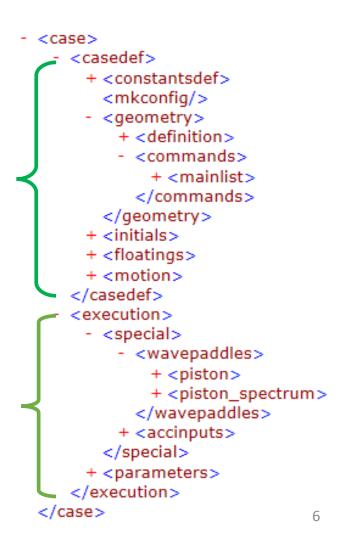
Divided in two sections:

"casedef"

Definition of the case with initial geometry and configuration. Created by the user and used by GenCase

"execution"

Information required to execute the case. Created by the user, modified by GenCase and only used by DualSPHysics



STRUCTURE OF THE XML FILE

- "casedef" :
- constantsdef constants needed in SPH
- mkconfig label configuration
- geometry system geometry (boundaries and fluid)
 - definition
 - commands (list & mainlist)
- initials special features for fluid particles
- floatings description of floating objects
- motion description of boundary movement
- "execution"
- **special** automatic wave generation and external forces
 - wavepaddles (piston & piston_spectrum)
 - accinputs
- parameters execution parameters in DualSPHysics

CASEDEF-CONSTANTSDEF

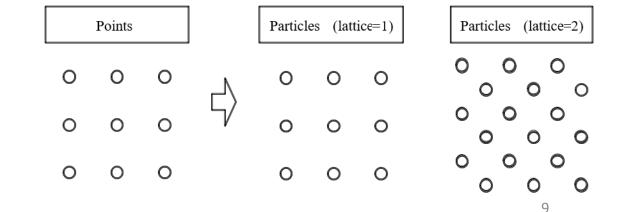
```
<constantsdef>
    <lattice bound="1" fluid="1" />
    <gravity x="0" y="0" z="-9.81" comment="Gravitational acceleration"</pre>
                                units comment="m/s^2" />
    <rhop0 value="1000" comment="Reference density of the fluid"
                                 units comment="kg/m^3" />
    <hswl value="0" auto="true" comment="Maximum still water level to calculate speedofsound"</pre>
                                 units comment="metres (m)" />
    <gamma value="7" comment="Polytropic constant for water used in the state equation" />
    <speedsystem value="0" auto="true" comment="Maximum system speed</pre>
                                 (by default the dam-break propagation is used) " />
    <coefsound value="20" comment="Coefficient to multiply speedsystem" />
    <speedsound value="0" auto="true" comment="Speed of sound to use in the simulation</pre>
                                 (by default speedofsound=coefsound*speedsystem) " />
    <coefh value="0.866025" comment="Coefficient to calculate the smoothing length
                                 (h=coefh*sqrt(3*dp^2) in 3D) " />
    <cflnumber value="0.2" comment="Coefficient to multiply dt" />
</constantsdef>
```

CASEDEF-CONSTANTSDEF

Lattice <constantsdef> <lattice bound="1" fluid="1" /> <gravity x="0" y="0" z="-9.81" comment="Gravitational acceleration"</pre> units comment="m/s^2" /> <rhop0 value="1000" comment="Reference density of the fluid" units comment="kg/m^3" /> <hswl value="0" auto="true" comment="Maximum still water level to calculate speedofsound"</pre> units comment="metres (m)" /> <gamma value="7" comment="Polytropic constant for water used in the state equation" /> <speedsystem value="0" auto="true" comment="Maximum system speed</pre> (by default the dam-break propagation is used) " /> <coefsound value="20" comment="Coefficient to multiply speedsystem" /> <speedsound value="0" auto="true" comment="Speed of sound to use in the simulation</pre> (by default speedofsound=coefsound*speedsystem) " /> <coefh value="0.866025" comment="Coefficient to calculate the smoothing length $(h=coefh*sqrt(3*dp^2) in 3D) " />$ <cflnumber value="0.2" comment="Coefficient to multiply dt" /> </constantsdef>

lattice: indicates the type of mesh

- to create particles:
- 1: one particle per point
- 2: two particles per point



CASEDEF-CONSTANTSDEF

Gravity

<constantsdef>

<lattice bound="1" fluid="1" /> <gravity x="0" y="0" z="-9.81" comment="Gravitational acceleration"</pre> units comment="m/s^2" /> <rhop0 value="1000" comment="Reference density of the fluid" units comment="kg/m^3" /> <hswl value="0" auto="true" comment="Maximum still water level to calculate speedofsound"</pre> units comment="metres (m)" /> <gamma value="7" comment="Polytropic constant for water used in the state equation" /> <speedsystem value="0" auto="true" comment="Maximum system speed</pre> (by default the dam-break propagation is used) " /> <coefsound value="20" comment="Coefficient to multiply speedsystem" /> <speedsound value="0" auto="true" comment="Speed of sound to use in the simulation</pre> (by default speedofsound=coefsound*speedsystem) " /> <coefh value="0.866025" comment="Coefficient to calculate the smoothing length $(h=coefh*sqrt(3*dp^2) in 3D) " />$ <cflnumber value="0.2" comment="Coefficient to multiply dt" /> </constantsdef>

$$\frac{d\boldsymbol{v}_a}{dt} = -\sum_b m_b \left(\frac{P_b + P_a}{\rho_b \cdot \rho_a} + \Pi_{ab} \right) \nabla_a W_{ab} + \boldsymbol{g}$$

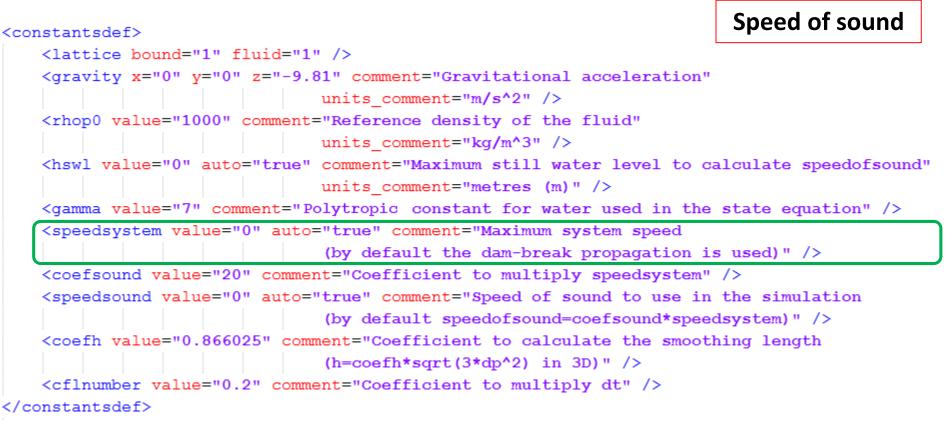
CASEDEF-CONSTANTSDEF

stantsdef>		Speed of sound
<pre><lattice <="" bound="1" fluid="1" pre=""></lattice></pre>	>	
	<pre>comment="Gravitational acceleration its comment="m/s^2" /></pre>	ב"
<rhop0 <br="" comment="R</td><th>eference density of the fluid" value="1000">its comment="kg/m^3" /><th></th></rhop0>		
<hswl auto="true" co<="" td="" value="0"><th>mment="Maximum still water level to</th><th>calculate speedofsound"</th></hswl>	mment="Maximum still water level to	calculate speedofsound"
	its comment="metres (m)" /> tropic constant for water used in t	ne state equation" />
<speedsystem auto="t</td><th>rue" comment="Maximum system speed</th><th></th></tr><tr><td>(b</td><th>y default the dam-break propagation</th><th>is used)" value="0"></speedsystem>		
<coefsound comment="</td" value="20"><th>"Coefficient to multiply speedsystem</th><th>n" /></th></coefsound>	"Coefficient to multiply speedsystem	n" />
<speedsound auto="tr</td><th>ue" comment="Speed of sound to use :</th><th>in the simulation</th></tr><tr><td>(b</td><th>y default speedofsound=coefsound*spe</th><th>eedsystem)" value="0"></speedsound>		
<coefh commen<="" td="" value="0.866025"><th>t="Coefficient to calculate the smoo</th><th>othing length</th></coefh>	t="Coefficient to calculate the smoo	othing length
(h	=coefh*sqrt(3*dp^2) in 3D)" />	
<cflnumber comment<="" td="" value="0.2"><th>="Coefficient to multiply dt" /></th><th></th></cflnumber>	="Coefficient to multiply dt" />	
onstantsdef>		

$$P = \frac{c_s^2 \rho_0}{\gamma} \left(\left(\frac{\rho}{\rho_0} \right)^{\gamma} - 1 \right)$$

$$B = \frac{c_s^2 \cdot \rho_0}{\gamma} = \frac{coef_{sound}^2 \cdot g}{\gamma} \frac{h_{swl} \cdot \rho_0}{\gamma}$$

CASEDEF-CONSTANTSDEF



$$P = \frac{c_s^2 \rho_0}{\gamma} \left(\left(\frac{\rho}{\rho_0} \right)^{\gamma} - 1 \right) \qquad \qquad c_s = coef_{sound} \cdot \sqrt{g \cdot h_{swl}} \\ B = \frac{c_s^2 \cdot \rho_0}{\gamma} = \frac{coef_{sound}^2 \cdot g \cdot h_{swl} \cdot \rho_0}{\gamma}$$

CASEDEF-CONSTANTSDEF

<pre>constantsdef> <lattice bound="1" fluid="1"></lattice> <gravity <="" comment="Gravitational acceleration" th="" x="0" y="0" z="-9.81"><th>ulate speedofsound</th></gravity></pre>	ulate speedofsound
<pre>units_comment="m/s^2" /> <rhop0 <="" comment="Reference density of the fluid" td="" value="1000"><td>ulate speedofsound</td></rhop0></pre>	ulate speedofsound
<pre><rhop <="" comment="Reference density of the fluid" td="" value="1000"><td>ulate speedofsound</td></rhop></pre>	ulate speedofsound
units_comment="kg/m^3" /> <hswl (m)"="" auto="true" comment="Maximum still water level to calc
units_comment=" metres="" value="0"></hswl> <gamma 0"="" 20"="" :<="" auto="true" comment="Coefficient to multiply speedsystem" td="" value="7"><td>culate speedofsound</td></gamma>	culate speedofsound
<pre><hswl (m)"="" auto="true" comment="Maximum still water level to cald</td><td>culate speedofsound</td></tr><tr><td>units_comment=" metres="" value="0"></hswl> <gamma 0"="" 20"="" :<="" auto="true" comment="Coefficient to multiply speedsystem" td="" value="7"><td>culate speedofsound</td></gamma></pre>	culate speedofsound
<pre><gamma 0"="" 20"="" auto="true" comment="Coefficient to multiply speedsystem" value="7"></gamma></pre>	ate equation" />
<pre><coefsound :<="" comment="Coefficient to multiply speedsystem" pre="" value="20"></coefsound></pre>	
	1sed) " />
	>
<pre><speedsound 0"="" auto="true" comment="Speed of sound to use in the
speedsound value=" i<="" in="" of="" sound="" speed="" speedsound="" td="" the="" to="" true"="" use="" value="0"><td>ne simulation</td></speedsound></pre>	ne simulation
(by default speedofsound=coefsound*speeds)	/stem)" />
<pre><coefh <="" comment="Coefficient to calculate the smoothin" pre="" value="0.866025"></coefh></pre>	ng length
(h=coefh*sqrt(3*dp^2) in 3D)" />	
<cflnumber comment="Coefficient to multiply dt" value="0.2"></cflnumber>	

</constantsdef>

$$P = \frac{c_s^2 \rho_0}{\gamma} \left(\left(\frac{\rho}{\rho_0} \right)^{\gamma} - 1 \right)$$

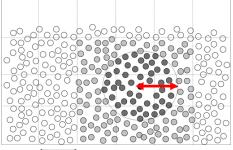
$$B = \frac{c_s^2 \rho_0}{\gamma} \left(\frac{\rho}{\rho_0} \right)^{\gamma} - 1 = \frac{c_s - \rho_0}{\gamma} \left(\frac{\rho}{\rho_0} \right)^{\gamma} - 1 = \frac{c_s - \rho_0}{\gamma} \left(\frac{\rho}{\rho_0} \right)^{\gamma} - 1 = \frac{c_s - \rho_0}{\gamma} \left(\frac{\rho}{\rho_0} \right)^{\gamma} - 1 = \frac{c_s - \rho_0}{\gamma} \left(\frac{\rho}{\rho_0} \right)^{\gamma} - 1 = \frac{c_s - \rho_0}{\gamma} \left(\frac{\rho}{\rho_0} \right)^{\gamma} - 1 = \frac{c_s - \rho_0}{\gamma} \left(\frac{\rho}{\rho_0} \right)^{\gamma} - 1 = \frac{c_s - \rho_0}{\gamma} \left(\frac{\rho}{\rho_0} \right)^{\gamma} - 1 = \frac{c_s - \rho_0}{\gamma} \left(\frac{\rho}{\rho_0} \right)^{\gamma} - 1 = \frac{c_s - \rho_0}{\gamma} \left(\frac{\rho}{\rho_0} \right)^{\gamma} - 1 = \frac{c_s - \rho_0}{\gamma} \left(\frac{\rho}{\rho_0} \right)^{\gamma} - 1 = \frac{c_s - \rho_0}{\gamma} \left(\frac{\rho}{\rho_0} \right)^{\gamma} - 1 = \frac{c_s - \rho_0}{\gamma} \left(\frac{\rho}{\rho_0} \right)^{\gamma} - 1 = \frac{c_s - \rho_0}{\gamma} \left(\frac{\rho}{\rho_0} \right)^{\gamma} - 1 = \frac{c_s - \rho_0}{\gamma} \left(\frac{\rho}{\rho_0} \right)^{\gamma} - 1 = \frac{c_s - \rho_0}{\gamma} \left(\frac{\rho}{\rho_0} \right)^{\gamma} - 1 = \frac{c_s - \rho_0}{\gamma} \left(\frac{\rho}{\rho_0} \right)^{\gamma} - 1 = \frac{c_s - \rho_0}{\gamma} \left(\frac{\rho}{\rho_0} \right)^{\gamma} - 1 = \frac{c_s - \rho_0}{\gamma} \left(\frac{\rho}{\rho_0} \right)^{\gamma} - 1 = \frac{c_s - \rho_0}{\gamma} \left(\frac{\rho}{\rho_0} \right)^{\gamma} - 1 = \frac{c_s - \rho_0}{\gamma} \left(\frac{\rho}{\rho_0} \right)^{\gamma} - 1 = \frac{c_s - \rho_0}{\gamma} \left(\frac{\rho}{\rho_0} \right)^{\gamma} - 1 = \frac{c_s - \rho_0}{\gamma} \left(\frac{\rho}{\rho_0} \right)^{\gamma} - 1 = \frac{c_s - \rho_0}{\gamma} \left(\frac{\rho}{\rho_0} \right)^{\gamma} - 1 = \frac{c_s - \rho_0}{\gamma} \left(\frac{\rho}{\rho_0} \right)^{\gamma} - 1 = \frac{c_s - \rho_0}{\gamma} \left(\frac{\rho}{\rho_0} \right)^{\gamma} - 1 = \frac{c_s - \rho_0}{\gamma} \left(\frac{\rho}{\rho_0} \right)^{\gamma} - 1 = \frac{c_s - \rho_0}{\gamma} \left(\frac{\rho}{\rho_0} \right)^{\gamma} - 1 = \frac{c_s - \rho_0}{\gamma} \left(\frac{\rho}{\rho_0} \right)^{\gamma} - 1 = \frac{c_s - \rho_0}{\gamma} \left(\frac{\rho}{\rho_0} \right)^{\gamma} - 1 = \frac{c_s - \rho_0}{\gamma} \left(\frac{\rho}{\rho_0} \right)^{\gamma} - 1 = \frac{c_s - \rho_0}{\gamma} \left(\frac{\rho}{\rho_0} \right)^{\gamma} - 1 = \frac{c_s - \rho_0}{\gamma} \left(\frac{\rho}{\rho_0} \right)^{\gamma} - 1 = \frac{c_s - \rho_0}{\gamma} \left(\frac{\rho}{\rho_0} \right)^{\gamma} - 1 = \frac{c_s - \rho_0}{\gamma} \left(\frac{\rho}{\rho_0} \right)^{\gamma} - 1 = \frac{c_s - \rho_0}{\gamma} \left(\frac{\rho}{\rho_0} \right)^{\gamma} - 1 = \frac{c_s - \rho_0}{\gamma} \left(\frac{\rho}{\rho_0} \right)^{\gamma} - 1 = \frac{c_s - \rho_0}{\gamma} \left(\frac{\rho}{\rho_0} \right)^{\gamma} - 1 = \frac{c_s - \rho_0}{\gamma} \left(\frac{\rho}{\rho_0} \right)^{\gamma} - 1 = \frac{c_s - \rho_0}{\gamma} \left(\frac{\rho}{\rho_0} \right)^{\gamma} - 1 = \frac{c_s - \rho_0}{\gamma} \left(\frac{\rho}{\rho_0} \right)^{\gamma} - 1 = \frac{c_s - \rho_0}{\gamma} \left(\frac{\rho}{\rho_0} \right)^{\gamma} - 1 = \frac{c_s - \rho_0}{\gamma} \left(\frac{\rho}{\rho_0} \right)^{\gamma} - 1 = \frac{c_s - \rho_0}{\gamma} \left(\frac{\rho}{\rho_0} \right)^{\gamma} - 1 = \frac{c_s - \rho_0}{\gamma} \left(\frac{\rho}{\rho_0} \right)^{\gamma} - 1 = \frac{c_s - \rho_0}{\gamma} \left(\frac{\rho}{\rho_0} \right)^{\gamma} - 1 = \frac{c_s - \rho_0}{\gamma} \left(\frac{\rho}{\rho_0} \right)^{\gamma} - 1 = \frac{c_s - \rho_0}{\gamma$$

CASEDEF-CONSTANTSDEF

stantsdef>	Kernel size
<lattice bound="1" fluid="1"></lattice>	
<pre><gravity)<="" comment="Gravitational acceleration" pre="" x="0" y="0" z="-9.81"></gravity></pre>	on"
units_comment="m/s^2" />	
<rhop0 <="" comment="Reference density of the fluid" td="" value="1000"><td></td></rhop0>	
units_comment="kg/m^3" />	
<hswl (m)"="" auto="true" comment="Maximum still water level to</td><td>calculate speedofsound</td></tr><tr><td>units_comment=" metres="" value="0"></hswl>	
<gamma comment="Polytropic constant for water used in t</td><td>the state equation" value="7"></gamma>	
<pre><speedsystem auto="true" comment="Maximum system speed</pre></td><td></td></tr><tr><td>(by default the dam-break propagation</td><td>is used)" value="0"></speedsystem></pre>	
<coefsound comment="Coefficient to multiply speedsyste</td><td>em" value="20"></coefsound>	
<pre><speedsound auto="true" comment="Speed of sound to use</pre></td><td>in the simulation</td></tr><tr><td>(by default speedofsound=coefsound*sp</td><td>eedsystem)" value="0"></speedsound></pre>	
<coefh comment="Coefficient to calculate the smo</td><td>othing length</td></tr><tr><td>(h=coefh*sqrt(3*dp^2) in 3D)" value="0.866025"></coefh>	
<cflnumber comment="Coefficient to multiply dt" value="0.2"></cflnumber>	

</constantsdef>

		$h = coefh \cdot \sqrt{dx^2 + dy^2 + dz^2}$	
coefh=1	typical value		00
coofb-1 2 1 5	better for wave propagation	$h = coefh \cdot \sqrt{3 \cdot dp^2}$	0
coem=1.2, 1.5	better for wave propagation	$h = coefh \cdot \sqrt{3} \cdot dp$	0
) (



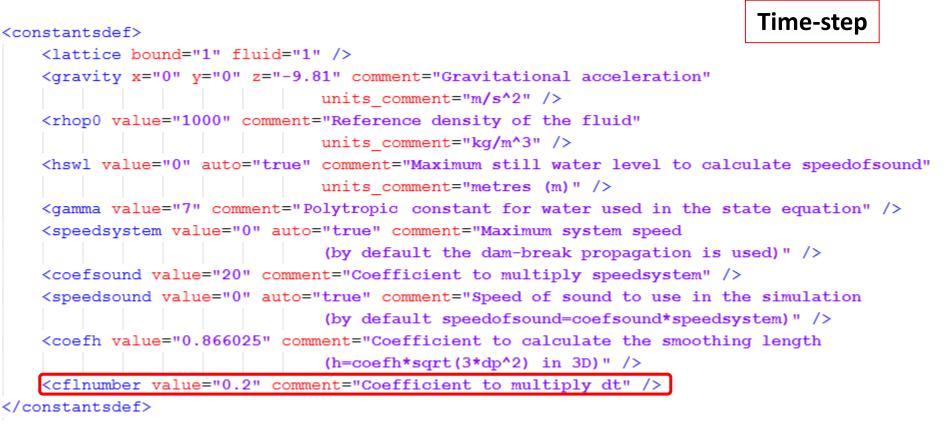
-

Other option is to define:

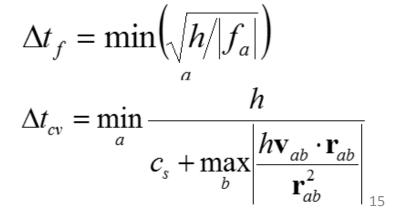
<hdp value="1.5" comment="Coefficient to calculate the smoothing length (hdp=h/dp)" />

$$hdp = h/dp$$

CASEDEF-CONSTANTSDEF



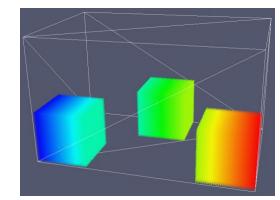
$$\Delta t = \mathbf{CFL} \cdot \min\left(\Delta t_f, \Delta t_{cv}\right)$$



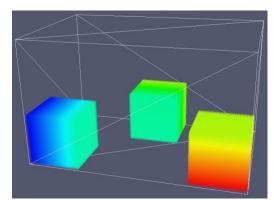
CASEDEF-MKCONFIG

- <mkconfig boundcount="240" fluidcount="10"> <mkorientbound mk="0" orient="YxZ"/> <mkorientfluid mk="1" orient="yzX"/> <mkorientfluid mk="2" orient="ZYx"/> </mkconfig>

mkorientfluid = "xyz"



mkorientfluid = "xyz" mkorientfluid = "yzX" mkorientfluid = "ZYx"



mk: label used to

- defines the order objects are created
- applies specific features to the different set of points such as movement, rigid motion...

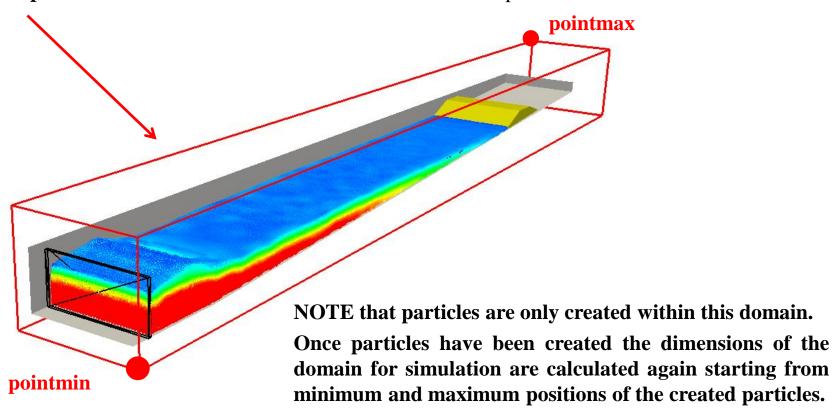
240 labels for boundary particles and 10 labels for fluid particles

mkorientation: determines the order of particles when creating one object (useful for visualization with the variable *idp*)

CASEDEF-GEOMETRY-DEFINITION

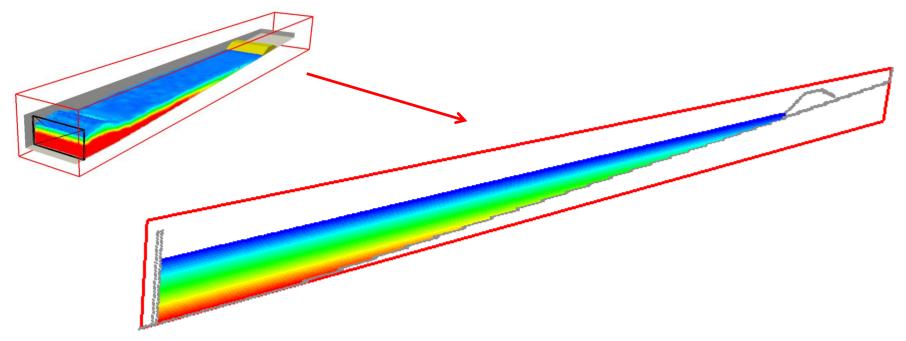
dp defines the distance between particles

WHEN CHANGING THIS PARAMETER, THE TOTAL NUMBER OF PARTICLES IS MODIFIED pointmin & pointmax defines the dimensions of the domain where particles can be created



CASEDEF-GEOMETRY-DEFINITION

A 2-D configuration can be generated by imposing the same values along Y-direction cpointmin> = <pointmax>



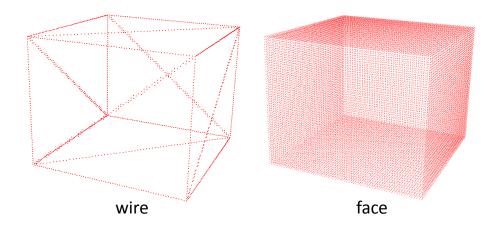
XML file CASEDEF-GEOMETRY-COMMANDS-MAINLIST

```
<commands>
    <mainlist>
        <setshapemode>dp | bound</setshapemode>
        <setdrawmode mode="full" />
        <!--CREATION OF FLUID PARTICLES (BOX OF WATER)-->
        <setmkfluid mk="0" />
                                                      Volume of fluid: setmkfluid mk=0,
        <drawbox>
                                                      solid to create particles within the specified volume
             <boxfill>solid</boxfill>
                                                      drawbox to plot a rectangular box defining a corner
             <point x="0" y="0" z="0" />
                                                      and its size in the 3 directions
             <size x="0.4" y="0.67" z="0.3" />
        </drawbox>
        <!--CREATION OF BOUNDARY PARTICLES (WALLS OF TANK) -->
        <setmkbound mk="0" />
        <drawbox>
             <boxfill>bottom | left | right | front | back</boxfill>
             <point x="0" v="0" z="0" />
                                                         Boundary Tank: setmkbound mk=0,
             <size x="1.6" v="0.67" z="0.4" />
                                                         specify box faces on which particles are
        </drawbox>
                                                         created (top is not used in this example)
        <shapeout file="" />
    </mainlist>
</commands>
```

This command indicates the mode to create points where particles will be generated

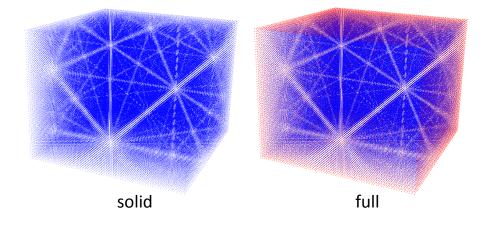
- <mainlist>

<setdrawmode mode="wire"/> <setdrawmode mode="face"/> <setdrawmode mode="solid"/> <setdrawmode mode="full"/> </mainlist>



<setdrawmode>:

- "wire": wire mode
- "face": draw faces
- "solid": draw inside
- "full": combines *face* and *solid*



```
- <mainlist>
    <setshapemode>dp | bound</setshapemode>
    <setshapemode>dp | bound | fluid</setshapemode>
    <setshapemode>real | void</setshapemode>
    </mainlist>
```

<setshapemode>: defines the draw operations to create VTK files (polygons)

- "real": using the real coordinates
- "**dp**": adjusting coordinates to *dp*
- "fluid": operations with *mk-fluid*
- "bound": operations with *mk-bound*
- "void": operations with *mk-void*

<setshapemode>: defines the draw operations to create a VTK files (polygons)

```
- <mainlist>
        <setshapemode>dp | bound</setshapemode>
        <setmkfluid mk="0"/>
        + <drawbox></drawbox>
        <setmkbound mk="0"/>
        + <drawbox></drawbox>
        <setmkvoid/>
        + <drawbox></drawbox>
        <setmkbound mk="1"/>
        + <drawbox></drawbox></drawbox>
        <setmkbound mk="1"/>
        + <drawbox></drawbox></drawbox></drawbox></drawbox></drawbox></drawbox></drawbox></drawbox></drawbox>
```

- <mainlist>
 <setshapemode>real | dp | bound</setshapemode>
 <setshapemode>real | dp | bound</setshapemode>
 <setdrawmode mode="full"/>
 <setmkfluid mk="0"/>
 + <drawprism mask="0"></drawprism>
 <setmkvoid/>
 + <drawbox></drawbox>
 <setdrawmode mode="face"/>
 <setdrawmode mode="face"/>
 <setmkbound mk="10"/>
 + <drawbox></drawbox>
 <setmkbound mk="0"/>
 + <drawbox></drawbox>
 <setmkbound mk="0"/>
 + <drawbox></drawbox>
 <setmkbound mk="0"/>
 + <drawprism mask="96"></drawprism>
 <setmkbound mk="10"/>
 + <drawprism mask="96"></drawprism>
 </setmkbound mk="10"/>
 </setmkbound mk="10"/>
 + <drawprism mask="96"></drawprism></setmkbound mk="10"/>
 </setmkbound mk="10"/>

shapeout: creates VTK files (polygons) of only some *bound* objects Case_Box_Dp.vtk Case_Building_Dp.vtk shapeout: creates VTK files (polygons) of all the *bound* objects Case__Real.vtk Case Dp.vtk

reset="true" objects created after this command will be saved on a different VTK file

These commands indicate the type of particles to be generated

- <commands>
- <commands>
- <mainlist>
 <setmkvoid/>
 <setmkfluid mk="0"/>
 <setmkbound mk="0"/>
 <setmknextfluid next="true"/>
 <setmknextfluid next="true"/>
 <setmknextbound next="false"/>
 <setmknextauto active="true"/>
 </mainlist>
 </commands>

<setmkvoid>, <setmkfluid>, <setmkbound>: defines the label *mk* to draw points of type: void (empty), fluid, bound

<setmknextfluid>, <setmknextbound>: increases (decreases) the value of *mk* with next=*true* (=*false*) <setmknextauto>: after each draw command *mk* is increased automatically

Transformation utilities

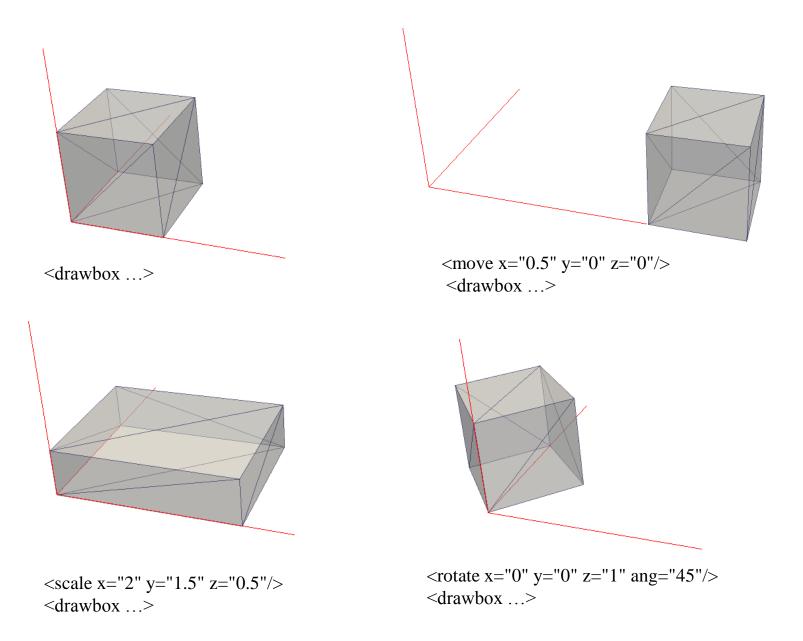
- <mainlist> <setshapemode>dp | bound</setshapemode> <setmkbound mk="0"/> <move x="0.5" v="0" z="0"/> + <drawbox></drawbox> <shapeout file="BoxMove" reset="true"/> <matrixreset/> <scale x="2" y="1.5" z="0.5"/> + <drawbox></drawbox> <shapeout file="BoxScale" reset="true"/> <matrixreset/> <rotate x="0" v="0" z="1" ang="45"/> + <drawbox></drawbox> <shapeout file="BoxRotate" reset="true"/> </mainlist>

<move>: a displacement is applied to the transformation matrix

<scale>: scaling is applied to matrix

<rotate>:a starting vector and angle are given for object rotation

<matrixreset>: the modified matrix is replaced by the original one (identity matrix)



```
- <mainlist>
    <setshapemode>dp | bound</setshapemode>
    <setmkbound mk="0"/>
  - < setlinebegin>
      <point x="0" y="0" z="0"/>
    </setlinebegin>
  - <drawlineto>
      <point x="0" y="1" z="0"/>
    </drawlineto>
    <setmknextbound next="true"/>
  - <drawline>
      <point x="0" y="1" z="0"/>
      <point x="1" y="1" z="0"/>
    </drawline>
    <setmknextbound next="true"/>
  - <drawline>
      <point x="1" y="1" z="0"/>
      <point x="1" y="0" z="0"/>
    </drawline>
    <setmknextbound next="true"/>
  - <drawlines>
      <point x="1" y="0" z="0"/>
      <point x="0" v="0" z="0.5"/>
      <point x="0" y="1" z="0.5"/>
      <point x="1" v="1" z="0.5"/>
      <point x="1" v="0" z="0.5"/>
    </drawlines>
    <shapeout file="Lines" reset="true"/>
  </mainlist>
```

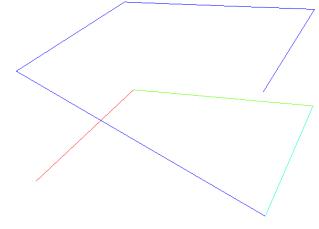
LINES

<setlinebegin>: sets the begining of the line with <drawlineto>

<drawlineto>: draws a line to a given point

<drawline>: draws a line between two points

<drawlines>: draws lines between several points

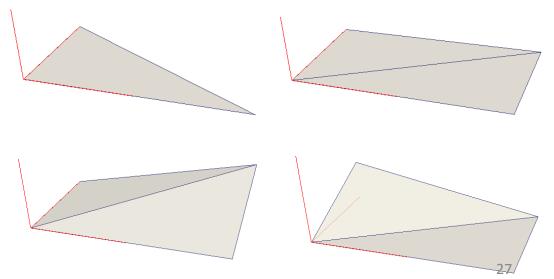


```
- <mainlist>
    <setshapemode>dp | bound</setshapemode>
    <setmkbound mk="0"/>
  - <drawtriangle>
      <point x="0" y="0" z="0"/>
      <point x="1" y="0" z="0"/>
      <point x="0" y="0.5" z="0"/>
    </drawtriangle>
    <shapeout file="Triangle" reset="true"/>
  - <drawquadri>
      > <point x="0" y="0" z="0"/>
      <point x="1" y="0" z="0"/>
      <point x="1" y="0.5" z="0"/>
      <point x="0" y="0.5" z="0"/>
    </drawquadri>
    <shapeout file="Quadri" reset="true"/>
  - <drawquadri>
      <point x="0" y="0" z="0"/>
      >point x="1" y="0" z="0"/>
      > <point x="1" y="0.5" z="0.2"/>
      </drawquadri>
    <shapeout file="Quadri2" reset="true"/>
  - <drawquadri>
      <point x="1" y="0" z="0"/>
      > point x="1" y="0.5" z="0"/>
      > point x="0" y="0.5" z="0.2"/>
    </drawquadri>
    <shapeout file="Quadri3" reset="true"/>
 </mainlist>
```

TRIANGLES

<drawtriangle>: draws a triangle with tree points (points must always go counterclockwise)

<drawquadri>: draws the quadrilateral described by four points (points may not be in the same plane)

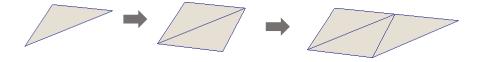


- <mainlist>

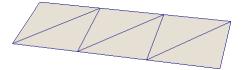
<setshapemode>dp | bound</setshapemode> <setmkbound mk="0"/> - <drawtrianglesstrip> <point x="0" y="1" z="0"/> <point x="0" y="0" z="0"/> <point x="1" y="1" z="0"/> <point x="1" y="0" z="0"/> <point x="2" y="1" z="0"/> <point x="2" y="0" z="0"/> <point x="3" y="1" z="0"/> >point x="3" y="0" z="0"/> <point x="4" y="1" z="0"/> </drawtrianglesstrip> <shapeout file="TrianglesStrip9" reset="true"/> </mainlist>

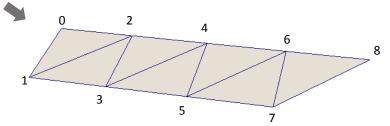
TRIANGLES

<drawtrianglesstrip>: draws a series of chained triangles





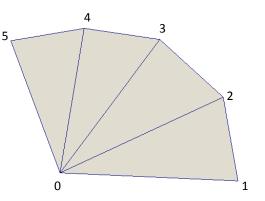


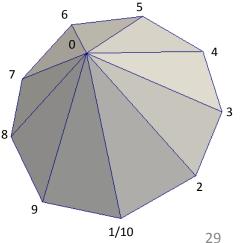


```
- <mainlist>
    <setshapemode>dp | bound</setshapemode>
    <setmkbound mk="0"/>
  - < drawtrianglesfan>
      <point x="0" y="0" z="0"/>
      <point x="1" y="0" z="0"/>
      > <point x="0.9" y="0.5" z="0"/>
      <point x="0.5" y="0.9" z="0"/>
      <point x="0" y="1" z="0"/>
      <point x="-0.5" y="0.9" z="0"/>
    </drawtrianglesfan>
    <shapeout file="TrianglesFan" reset="true"/>
    <setmkbound mk="0"/>
  - <drawtrianglesfan>
      <point x="0" y="0" z="1"/>
      <point x="1" y="0" z="0"/>
      > <point x="0.8" y="0.6" z="0"/>
      <point x="0.2" y="1" z="0"/>
      <point x="-0.5" y="0.9" z="0"/>
      >point x="-0.9" y="-0.3" z="0"/>
      >point x="-0.5" y="-0.9" z="0"/>
      <point x="0.2" v="-1" z="0"/>
      > <point x="0.8" v="-0.6" z="0"/>
      <point x="1" y="0" z="0"/>
    </drawtrianglesfan>
    <shapeout file="TrianglesFan2" reset="true"/>
  </mainlist>
```

TRIANGLES

<drawtrianglesfan>: draws a range of triangles

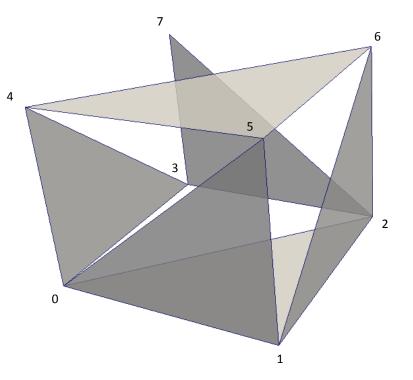




```
- <mainlist>
    <setshapemode>dp | bound</setshapemode>
    <setmkbound mk="0"/>
  - <drawtriangles>
    - <points>
         <point x="0" y="0" z="0"/>
         <point x="1" y="0" z="0"/>
         <point x="1" y="1" z="0"/>
         <point x="0" y="1" z="0"/>
         <point x="0" y="0" z="0.8"/>
         <point x="1" y="0" z="0.8"/>
         <point x="1" y="1" z="0.8"/>
         <point x="0" y="1" z="0.8"/>
       </points>
    - <triangles>
         <triangle x="0" y="1" z="5"/>
         <triangle x="1" y="2" z="6"/>
         <triangle x="2" y="3" z="7"/>
         <triangle x="3" y="0" z="4"/>
         <triangle x="0" y="2" z="1"/>
         <triangle x="4" y="5" z="6"/>
      </triangles>
    </drawtriangles>
    <shapeout file="Triangles" reset="true"/>
  </mainlist>
```

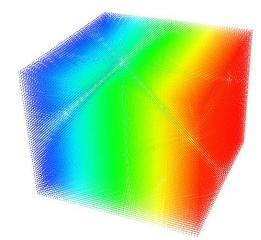
TRIANGLES

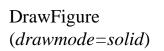
<drawtriangles>: draws a series of triangles defined by a set of points or a set of triangles

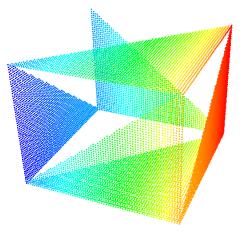


FIGURE

<drawfigure>: draws a solid figure consisting of all the interior points to the planes formed by the given triangles







DrawTriangles or DrawFigure (*drawmode=face*)

- <mainlist>

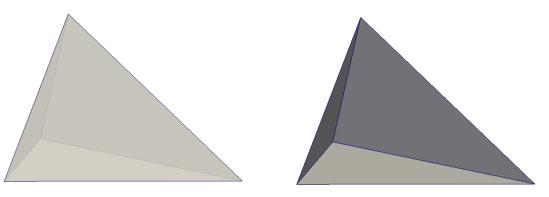
```
<setshapemode>dp | bound</setshapemode>
<setmkbound mk="0"/>
- <drawfigure>
- <points>
<point x="0" y="0" z="0"/>
```

```
<point x="1" y="0" z="0"/>
       >point x="1" y="1" z="0"/>
       <point x="0" y="1" z="0"/>
       > > > > > > 
       <point x="1" y="0" z="0.8"/>
       <point x="1" y="1" z="0.8"/>
       <point x="0" y="1" z="0.8"/>
    </points>
  - <triangles>
       <triangle x="0" y="1" z="5"/>
       <triangle x="1" y="2" z="6"/>
       <triangle x="2" y="3" z="7"/>
       <triangle x="3" y="0" z="4"/>
       <triangle x="0" y="2" z="1"/>
       <triangle x="4" y="5" z="6"/>
    </triangles>
  </drawfigure>
</mainlist>
```

```
- <mainlist>
    <setshapemode>dp | bound</setshapemode>
    <setdrawmode mode="full"/>
    <setmkbound mk="0"/>
  - <drawpyramid mask="0">
      > point x="0.25" y="0.25" z="0.7"/>
      >point x="0" y="0" z="0"/>
      <point x="1" y="0" z="0"/>
      <point x="0" y="1" z="0"/>
    </drawpyramid>
    <shapeout file="Pyramid1" reset="true"/>
  - <drawpyramid mask="2">
      > point x="0.25" y="0.25" z="0.7"/>
      <point x="0" y="0" z="0"/>
      <point x="1" y="0" z="0"/>
      <point x="0" y="1" z="0"/>
    </drawpyramid>
    <shapeout file="Pyramid2" reset="true"/>
  </mainlist>
```

<drawpyramid>: draws a pyramid with the top point and other points of the base (minimum 3)

mask indicates the faces to be hidden with bits the first bit always corresponds to the base and the rest to the faces following the order



Pyramid1 (*mask=0*)

Pyramid2 (*mask*=2=0010)

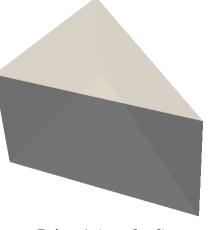
PRISM

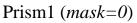
- <mainlist> <setshapemode>dp | bound</setshapemode> <setdrawmode mode="full"/> <setmkbound mk="0"/> - <drawprism mask="0"> <point x="0" y="0" z="0"/> <point x="1" y="0" z="0"/> <point x="0" y="1" z="0"/> <point x="0" y="0" z="0.5"/> <point x="1" y="0" z="0.5"/> > <point x="0" v="1" z="0.5"/> </drawprism> <shapeout file="Prism1" reset="true"/> - <drawprism mask="2"> <point x="0" y="0" z="0"/> <point x="1" y="0" z="0"/> <point x="0" y="1" z="0"/> >point x="0" v="0" z="0.5"/> >point x="1" y="0" z="0.5"/> <point x="0" v="1" z="0.5"/> </drawprism> <shapeout file="Prism2" reset="true"/> </mainlist>

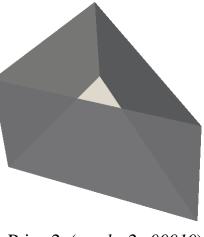
<drawprism>: draws a prism with a minimum of 6 points

The first half of points are the base and the second half the top (the number of points must be even)

mask indicates the faces to be hidden with bits The first bit corresponds to the base, the second to the top and the rest to the faces following the order







Prism2 (*mask*=2=00010)

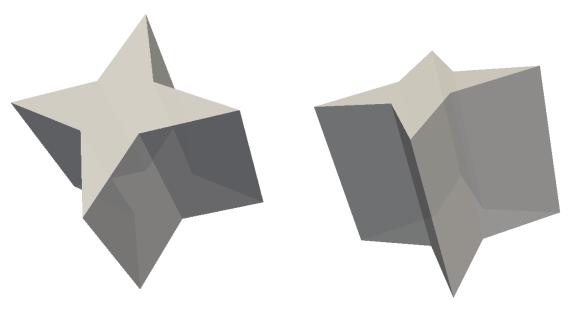
PRISM

- <mainlist> <setshapemode>dp | bound</setshapemode> <setdrawmode mode="full"/> <setmkbound mk="0"/> - <drawprism mask="0"> <point x="0" y="0" z="0"/> <point x="1" y="-3" z="0"/> <point x="2" y="0" z="0"/> <point x="5" y="1" z="0"/> <point x="2" v="2" z="0"/> <point x="1" y="5" z="0"/> <point x="0" y="2" z="0"/> <point x="-3" y="1" z="0"/> <point x="0" y="0" z="6"/> <point x="1" y="-3" z="6"/> <point x="2" y="0" z="6"/> <point x="5" v="1" z="6"/> <point x="2" y="2" z="6"/> <point x="1" y="5" z="6"/> <point x="0" v="2" z="6"/> <point x="-3" y="1" z="6"/> </drawprism> <shapeout file="Prism3" reset="true"/> </mainlist>

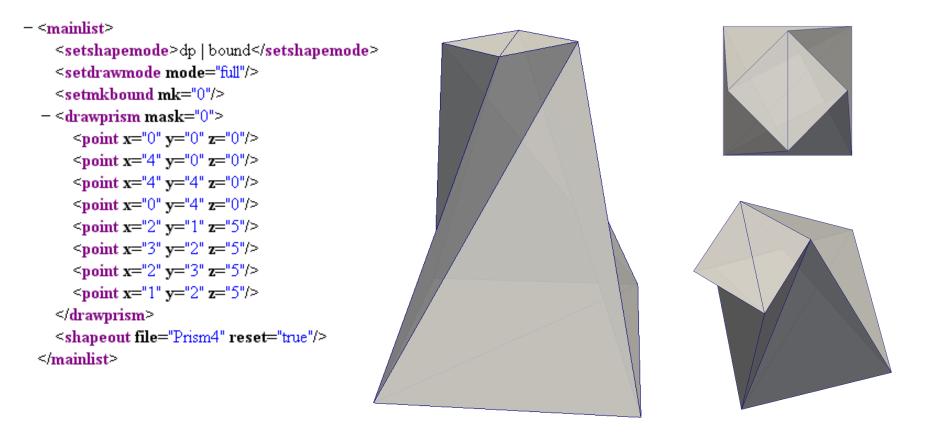
<drawprism>: draws a prism with a minimum of 6 points

The first half of points are the base and the second half the top (the number of points must be even)

mask indicates the faces to be hidden with bits The first bit corresponds to the base, the second to the top and the rest to the faces following the order



PRISM



MASK

mask indicates the faces to be hiddenInitially this is defined using BITS

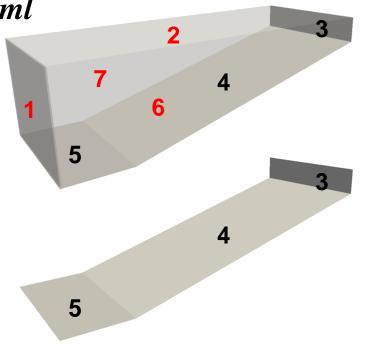
FOR EXAMPLE: OBJECT WITH 4 FACES:

mask="0"	decimal 0 is 0000 in binary	no faces are hidden
mask="1"	decimal 1 is 0001 in binary	first face is hidden
mask="2"	decimal 2 is 0010 in binary	second face is hidden
mask="4"	decimal 4 is 0100 in binary	third face is hidden
mask="8"	decimal 8 is 1000 in binary	fourth face is hidden
mask="12"	decimal 12 is 1100 in binary	third and fourth face are hidden

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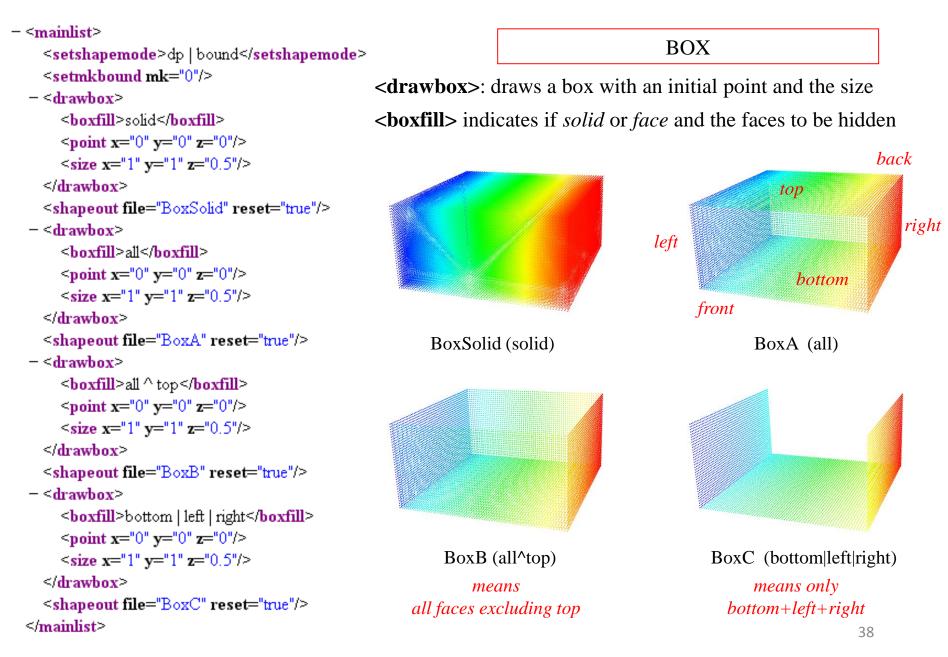
There is a second and easiest system to use mask mask can be also defined using the index of the faces instead of bits In this example, faces 1, 2, 6 and 7 are not created, only 3,4 and 5 It is important to use symbol "|" to detect this system!!!

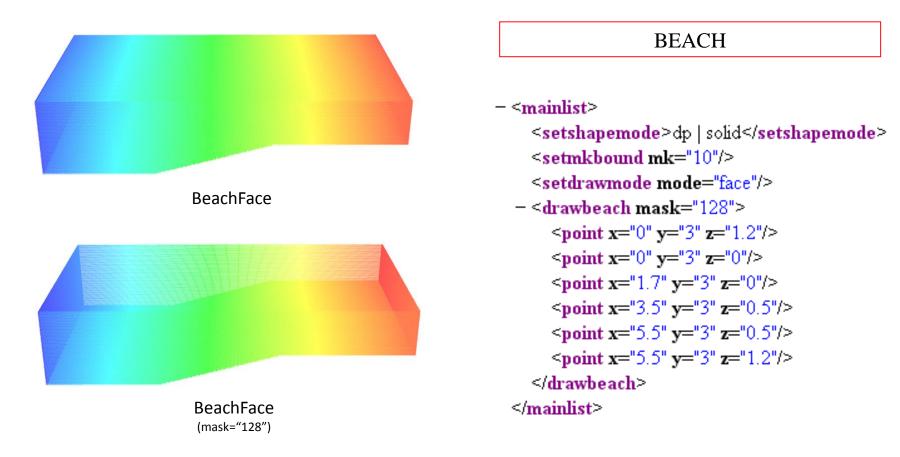
<setmkbound mk="0" /> <drawprism mask="1 | 2 | 6 | 7"> <point x="5" y="0" z="1.5" /> <point x="5" y="0" z="1.1" /> <point x="1" y="0" z="0" /> <point x="0" y="0" z="0" /> <point x="0" y="0" z="1.5" /> <point x="5" y="2" z="1.5" /> <point x="5" y="2" z="1.1" /> <point x="1" y="2" z="0" /> <point x="1" y="2" z="0" /> <point x="0" y="2" z="1.5" /> <point x="1" y="2" z="0" /> <point x="0" y="2" z="1.5" />



PRISM

EXAMPLE: CaseWavemaker_Def.xml





<drawbeach>: draws a beach with the lateral points that formed the profile of the beach mask indicates the faces to be hidden .

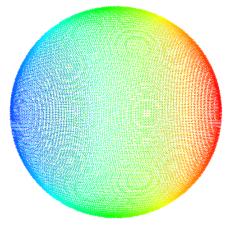
mask="128" decimal 128 is 10000000 in binary

eighth face is hidden

SPHERE

- <mainlist> <setmkbound mk="0"/> <setdrawmode mode="solid"/> <drawsphere radius="0.8"> <point x="1" y="1" z="1"/> </drawsphere> <setdrawmode mode="face"/> <drawsphere radius="0.8"> <point x="1" y="1" z="1"/> </drawsphere> </drawsphere radius="0.8"> </drawsphere radius="0.8"> </drawsphere radius="0.8"> </drawsphere radius="0.8"> </drawsphere> </drawsphere radius="0.8"> </drawsphere radius="0.8"> </drawsphere radius="0.8"> </drawsphere radius="0.8"> </drawsphere radius="0.8"> </drawsphere radius="0.8"> </drawsphere> </drawsphere> </drawsphere> </drawsphere> </drawsphere>

Sphere (*drawmode=solid*)



Sphere (*drawmode=face*)

when *face*:

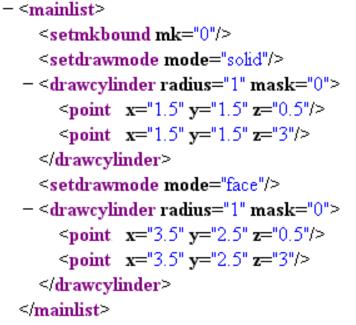
<setdpctes ctesphere="0.4"/> <setdpctes ctespherenumsides="40"/> ctesphere indicates the width of the sphere

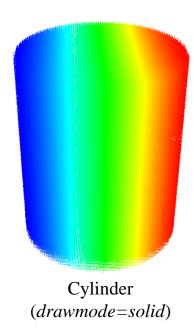
ctespherenumsides indicates the number of triangles used to create the VTK of polygons

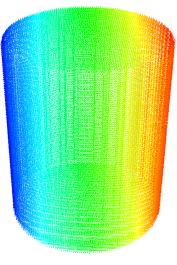
<drawsphere>: draws a sphere with the center point and the radius

CYLINDER

<drawcylinder>: draws a cylinder with two points and radius mask indicates the faces to be hide







Cylinder (*drawmode=face*)

when *face*:

<setdpctes ctecylindertube="0.6"/> <setdpctes ctecylindercover="0.7"/> <setdpctes ctecylindernumsides="40"/> ctecylindertube indicates the width of the tube ctecylindercover indicates the width of the covers ctespherenumsides indicates the number of triangles used to create the VTK of polygons

<setmkbound mk="0"/> <drawfilestl file="File.stl"/> <drawfileply file="File.ply"/> <drawfileply file="File.vtk"/> - <drawfilestl file="File.stl"> <drawmove x="0.5" v="0" z="0"/> <drawrotate angx="10" angy="15" angz="30"/> <drawscale x="1" y="1" z="0.8"/> </drawfilestl> - <drawfileply file="File.ply"> <drawmove x="0.5" y="0" z="0"/> </drawfileply> - <drawfileply file="File.ply"> <drawmove x="0.5" y="0" z="0"/> <drawrotate angx="10" angy="15" angz="30"/> </drawfileply> - <drawfileply file="File.ply"> <drawrotate angx="10" angy="15" angz="30"/> </drawfileply> - <drawfilevtk file="File.vtk"> <polyselec>points</polyselec> </drawfilevtk> - <drawfilevtk file="File.vtk"> <polyselec>points | lines</polyselec> </drawfilevtk> - <drawfilevtk file="File.vtk"> <polyselec>triangles</polyselec> </drawfilevtk> - <drawfilevtk file="File.vtk"> <polyselec>polygons</polyselec> </drawfilevtk> </mainlist>

- <mainlist>

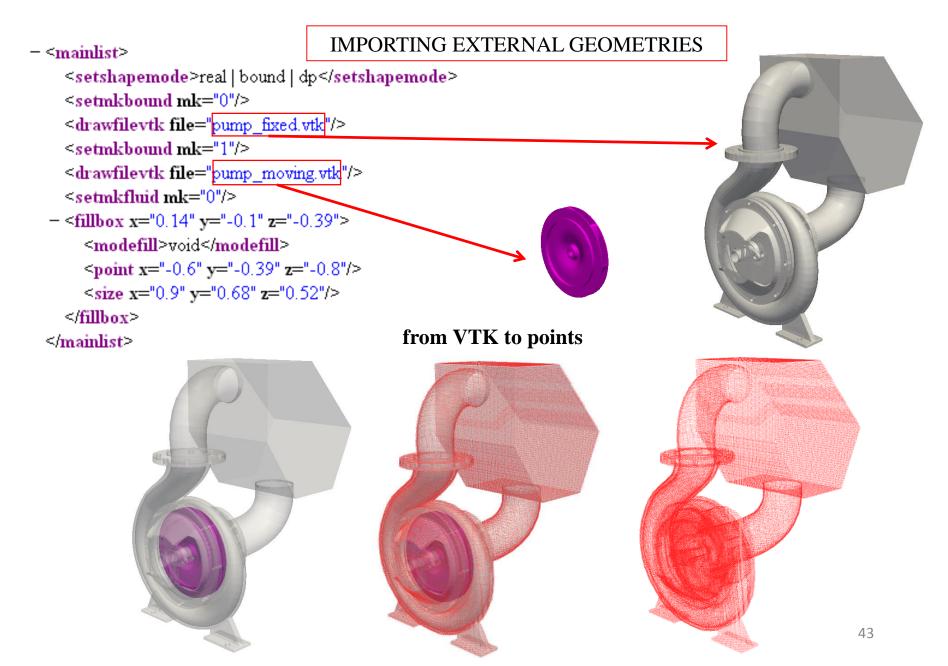
IMPORTING EXTERNAL GEOMETRIES

<drawfilevtk>: load a VTK file to be converted into points

<drawfileply>: load a PLY file to be converted into points

<drawfilestl>: load a STL file to be converted into points

Some modifications can be applied to the VTK, PLY or STL **drawmove** a displacement is applied to the external object **drawrotate** a rotation is applied to the external object **drawscale** scaling is applied to the external object



```
- <mainlist>
    <setmkfluid mk="0"/>
    <fillvoidpoint x="3" y="2" z="1"/>
  - <fillpoint x="3" y="2" z="1">
       <modefill>void</modefill>
    </fillpoint>
  - <fillpoint x="1" y="1" z="1" mkfluid="0">
       <modefill>fluid</modefill>
    </fillpoint>
  - <fillpoint x="1" y="1" z="1" mkbound="0">
       <modefill>bound</modefill>
    </fillpoint>
  - <fillpoint x="2" y="2" z="2" mkfluid="2" mkbound="8">
       <modefill>border | void | fluid | bound</modefill>
    </fillpoint>
  - < fillbox x = 0" v = 1" z = 0">
       <modefill>border</modefill>
      <point x="0.1" y="1" z="1.1"/>
      <size x="3" v="4" z="2"/>
    </fillbox>
  - <fillprism x="2" y="3" z="5">
       <point x="0" y="0" z="0"/>
      <point x="1" y="0" z="0"/>
      <point x="0" y="1" z="0"/>
      <point x="0" y="0" z="0.5"/>
      >point x="1" y="0" z="0.5"/>
       <point x="0" y="1" z="0.5"/>
       <modefill>void</modefill>
    </fillprism>
    <debugout/>
 </mainlist>
```

FILLING DOMAINS

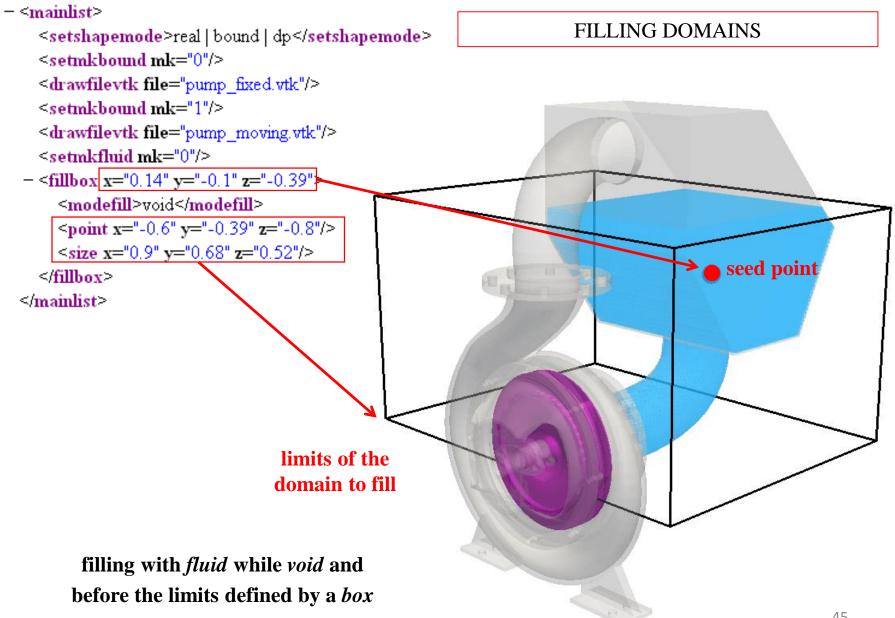
<fillpoint>: fills with points starting from the seed

<**fillbox>:** fills with points starting from the seed within the limits defined by a box

<**fillfigure>:** fills with points starting from the seed within the limits defined by a figure

<**fillprism>:** fills with points starting from the seed within the limits defined by a prism

<modefill> indicates what type of points can be filled with *void*, *fluid*, *bound*, it fills with that type of points inside the specified limits or the presence of a given type of point using *border*



CASEDEF-INITIALS

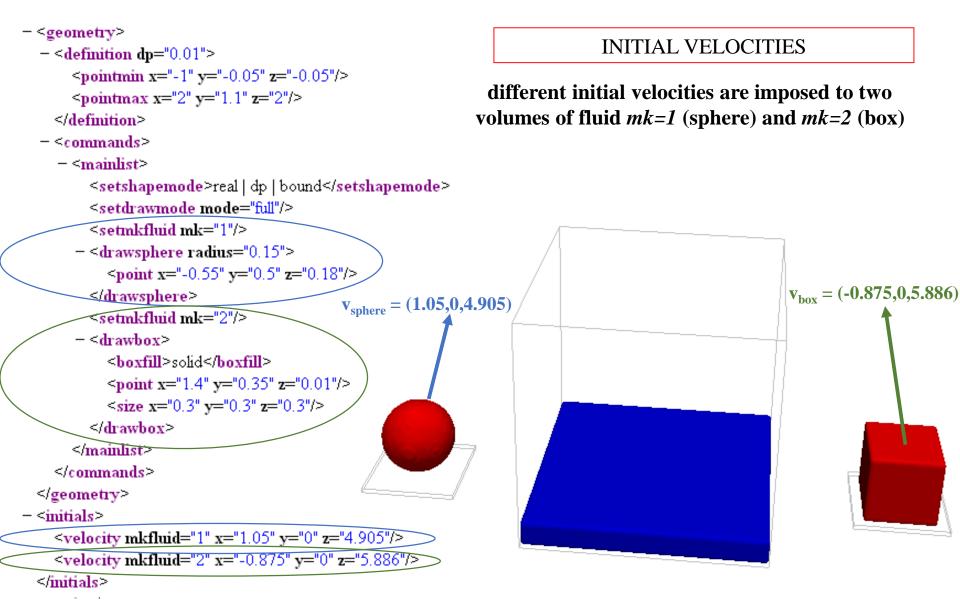
```
- <geometry>
  - <definition dp="0.01">
      >pointmin x="-1" y="-0.05" z="-0.05"/>
      <pointmax x="2" y="1.1" z="2"/>
    </definition>
  - <commands>
    - <mainlist>
         <setshapemode>real | dp | bound</setshapemode>
         <setdrawmode mode="full"/>
         <setmkfluid mk="1"/>
       - <drawsphere radius="0.15">
           > point x="-0.55" v="0.5" z="0.18"/>
         </drawsphere>
         <setmkfluid mk="2"/>
       - <drawbox>
           <boxfill>solid</boxfill>
           <point x="1.4" y="0.35" z="0.01"/>
           <size x="0.3" v="0.3" z="0.3"/>
        </drawbox>
      </mainlist>
    </commands>
 </geometry>
<initials>
    <velocity mkfluid="1" x="1.05" y="0" z="4.905"/>
    <velocity mkfluid="2" x="-0.875" y="0" z="5.886"/>
 </initials>
```

INITIAL VELOCITIES

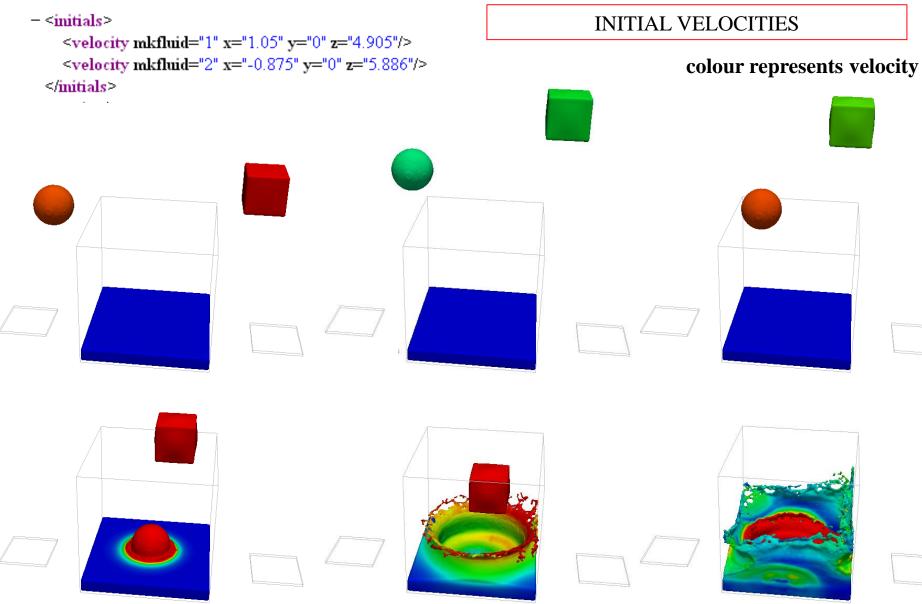
<initials>: special behaviours can be imposed to a set of fluid particles labeled with a *mk*, such as:

<velocity> initial velocity defined by a vector <velwave> a solitary wave defined by *depth* and *amplitude*

CASEDEF-INITIALS



CASEDEF-INITIALS



CASEDEF-FLOATINGS

```
- <floatings>
```

<floating mkbound="0" relativeweight="1.3"/>

- <floating mkbound="1" relativeweight="1.3">
<velini x="1" y="3" z="2"/>
<omegaini x="0.2" y="0.4" z="0.6"/>
<floating>

- <floating mkbound="2">
 <massbody value="1300"/>
 <center x="11" y="12" z="13"/>
 <inertia x="20" y="22" z="24"/>

</floating>

- <floating mkbound="3">
 <massbody value="1300"/>
 <center x="11" y="12" z="13"/>
 <inertia x="20" y="22" z="24"/>
 <velini x="1" y="3" z="2"/>
 <omegaini x="0.2" y="0.4" z="0.6"/>
 </floating>

```
- <floating mkbound="4">
        <massbody value="1300"/>
        <inertia x="20" y="22" z="24"/>
        </floating>
</floatings>
```

DEFINING FLOATINGS

<**floatings>:** indicates that a set of particles labelled with the same *mk* constitutes a floating object

Only one of these values can be defined: rhopbody density of the object relativeweight in relation to the reference density massbody total mass of the object

So that, the mass of a floating particles is: **masspart** = massbody / nfloat **or masspart** = relativeweight * rhop0 * dp^3 **or masspart** = rhopbody * dp^3

These variables are computed by GenCase or can be also specified in advance: **center** gravity center of the rigid object **inertia** momentum of inertia of the rigid object **velini** initial linear velocity of the object **omegaini** initial angular velocity of the object

CASEDEF-FLOATINGS

When the interaction of solids (boundaries or floatings) is computed using **Discrete Element Method (DEM)** some extra properties with parameters used in DEM are loaded from **"Floating_Materials.xml":**

```
<materials>
   <property name="steel">
        <Young_Modulus value="21000000000.0" comment="Young Modulus (N/m2)" />
        <PoissonRatio value="0.30" comment="Poisson Ratio (-)" />
        <Restitution_Coefficient value="0.80" comment="Restitution Coefficient (-)" />
        <Kfric value="0.45" comment="Kinetic friction coefficient" />
        </property>
</materials>
```

CASEDEF-PROPERTIES

DEFINING OTHER VARIABLES

Using section *properties*, users can define variables to be assigned to one or more *mk*.

```
<properties>
<links>
<link mkfluid="0" property="material_1" />
<link mkbound="3-6,1" property="material_2+data_x" />
</links>
<propertyfile file="run/ftdata_ext.xml" path="case.materials" />
<property name="material_1" weight="1.35" other="pepe"/>
<property name="material_2" begin="168" count="973">
<massbody value="4728.78" />
<center x="4.99" y="5" z="7.03" />
</property name="data_x" weight="1.35" />
</properties>
```

Each label of *property* has a name and can group several values that can be text (*other*) or a number (*weight*)

```
<property name="material 1" weight="1.35" other="pepe"/>
```

or other subvalues (*massbody* and *center*)

CASEDEF-PROPERTIES

These properties can be loaded from an external file using *propertyfile*. In this case, users have to indicate file name and path to access section with properties.

```
<propertyfile file="run/ftdata_ext.xml" path="case.materials" />
```

Example of file "ftdata_ext.xml":

```
<case>
<materials>
<property name="uno" value="1.35"/>
<property name="dos" value="168">
<property name="dos" value="168">
<massbody value="4728.78" />
</property>
</materials>
</case>
```

Section *links* assigns one or more *property* to one or several values of *mk*:

- Values of *material_1* are assigned to fluid particles with *mk*=0

```
<link mkfluid="0" property="material 1"/>
```

- Values of *material_2* and *data_x* are assigned to boundary particles with mk=1,3,4,5,6

```
<link mkbound="3-6,1" property="material_2+data_x"/>
```

It is also possible to indicate one *property* directly in the definition of the *floatings*:

CASEDEF-PROPERTIES

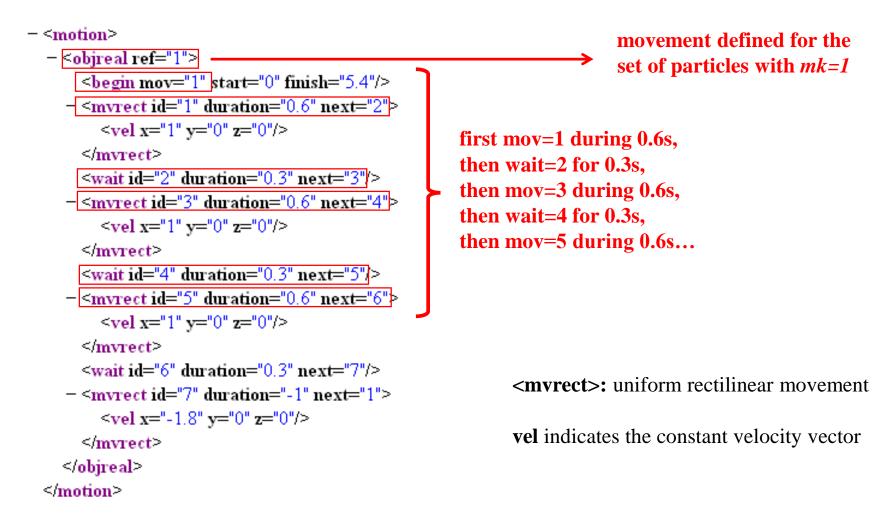
GenCase reads the information from case.casedef.properties and writes in case.execution.particles.properties.

```
<particles np="1494" nb="313" nbf="313" mkboundfirst="11" mkfluidfirst="1">
    <fixed mkbound="0" mk="11" begin="0" count="229" />
    <fixed mkbound="1" mk="12" begin="229" count="28" property="data x+material 2" />
    <fixed mkbound="2" mk="13" begin="257" count="28" />
    <fixed mkbound="4" mk="15" begin="285" count="28" property="data x+material 2" />
    <fluid mkfluid="0" mk="1" begin="313" count="1146" property="material 1" />
    <fluid mkfluid="1" mk="2" begin="1459" count="35" />
    <properties>
        <links>
            k mk="1" property="material 1" />
            k mk="12,15" property="data x+material 2" />
        </links>
        <property name="material 1" weight="1.35" other="pepe"/>
        <property name="material 2" begin="168" count="973"></property name="material 2" begin="168" count="973">
            <massbody value="4728.78" />
            <center x="4.99" y="5" z="7.03" />
        </property>
        <property name="data x" weight="1.35" />
    </properties>
</particles>
```

Thus, *DualSPHysics* can access to assigned values to each *mk*.

The object of type JSpaceParts is used to obtain the assigned properties to each block of particles

•*Motion01*: uniform rectilinear motion (**<mvrect** /**>**) that also includes pauses (**<wait** /**>**)



•*Motion01*: uniform rectilinear motion (**<mvrect** /**>**) that also includes pauses (**<wait** /**>**)



Time: 2.40 s

Time: 2.70 s

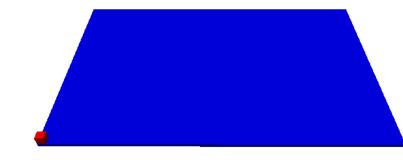
•*Motion02*: combination of two uniform rectilinear motions (**<mvrect** /**>**)

```
- <motion>
  - <objreal ref="1">
      <begin mov="1" start="0"/>
    - <mvrect id="1" duration="2" next="2">
         <vel x="1" v="0" z="0"/>
      </mvrect>
    - <mvrect id="2" duration="1" next="1">
        <vel x="-2" y="0" z="0"/>
      </mvrect>
      <br/>
<br/>
start="0.5"/>
    - <mvrect id="3" duration="1.3" next="4">
        <vel x="0" y="1" z="0"/>
      </mvrect>
    - <mvrect id="4" duration="1.3" next="3">
         <vel x="0" y="-1" z="0"/>
      </mvrect>
    </objreal>
 </motion>
```

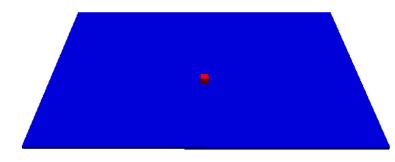
<mvrect>: uniform rectilinear movement

vel indicates the constant velocity vector

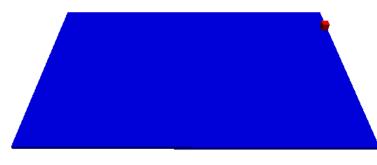
•*Motion02*: combination of two uniform rectilinear motions (**<mvrect** /**>**)



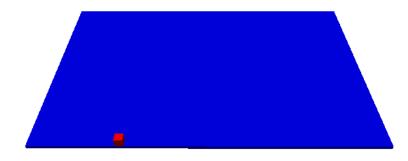
Time: 0.00 s



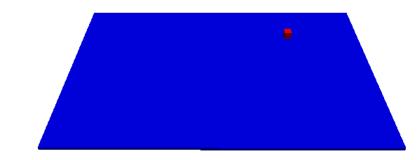
Time: 1.00 s



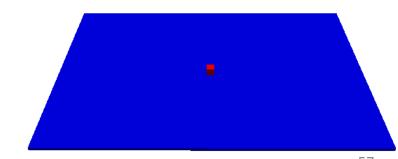
Time: 2.00 s



Time: 0.50 s



Time: 1.50 s



Time: 2.50 s

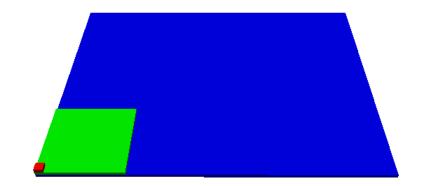
– <motion> - <objreal ref="1"> <begin mov="1" start="0"/> - <mvrect id="1" duration="1.5" next="2"> <vel x="1" y="0" z="0"/> </mvrect> - <mvrect id="2" duration="1.5" next="1"> <vel x="-1" y="0" z="0"/> </mvrect>

start="0.1"/> - <mvrect id="3" duration="1.1" next="4"> <vel x="0" y="1" z="0"/> </mvrect> - <mvrect id="4" duration="1.1" next="3"> <vel x="0" y="-1" z="0"/> </mvrect> - <objreal ref="2">
begin mov="1" start="0.2"/> - <mvrect id="1" duration="0.45" next="2"> <vel x="1" v="0" z="0"/> </mvrect> - <mvrect id="2" duration="0.45" next="3"> <vel x="0" y="1" z="0"/> </mvrect> - <mvrect id="3" duration="0.45" next="4"> <vel x="-1" y="0" z="0"/> </mvrect> - <mvrect id="4" duration="0.45" next="1"> <vel x="0" y="-1" z="0"/> </mvrect> </objreal> </objreal> </motion>

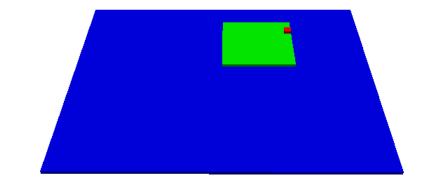
•*Motion03*: movement of an object depending on the movement of another (hierarchy of objects)

movement defined for the set of particles with mk=2 that also moves according to the movement defined for mk=1

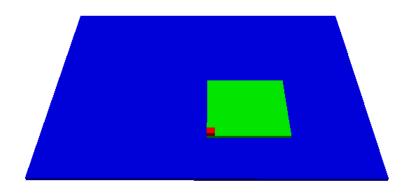
•Motion03: movement of an object depending on the movement of another (hierarchy of objects)



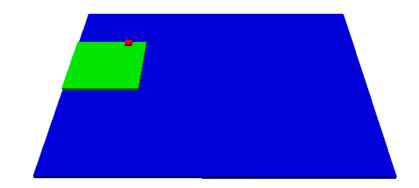
Time: 0.00 s



Time: 1.00 s



Time: 2.00 s



Time: 3.00 s

•*Motion04*: accelerated rectilinear motion (**<mvrectace** /**>**)

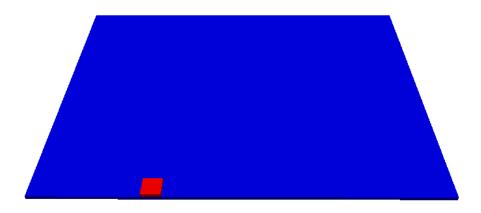
<mvrectace>: accelerated rectilinear movement

velini indicates the initial velocity vector **ace** indicates the acceleration vector

•*Motion04*: accelerated rectilinear motion (**<mvrectace** /**>**)



Time: 0.00 s



Time: 0.75 s



Time: 1.50 s



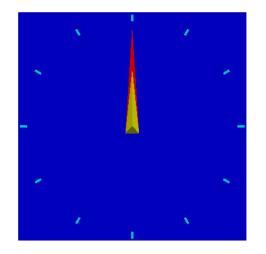
•*Motion*05: rotational motion (**<mvrot** />)

```
- <motion>
  - <objreal ref="3">
      <br/>
start="0"/>
    - <mvrot id="1" duration="1000">
         <vel ang="20"/>
         <axisp1 x="0.5" y="0.5" z="0"/>
         <axisp2 x="0.5" y="0.5" z="1"/>
      </mvrot>
    </objreal>
  - <objreal ref="4">
      <begin mov="1" start="0"/>
    - <mvrot id="1" duration="1000">
         <vel ang="240"/>
         <axisp1 x="0.5" y="0.5" z="0"/>
         <axisp2 x="0.5" y="0.5" z="1"/>
      </mvrot>
    </objreal>
 </motion>
```

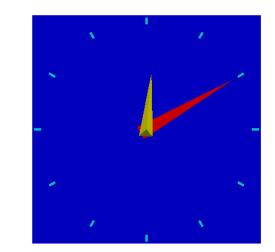
<mvrot>: rotational movement

vel indicates the angular velocityaxisp1 first point of the rotation axisaxisp2 second point of the rotation axis

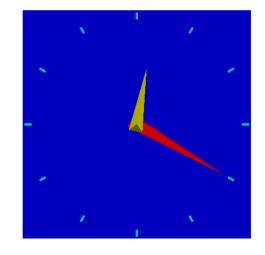
•*Motion*05: rotational motion (**<mvrot** />)



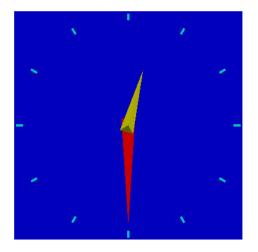
Time: 0.00 s

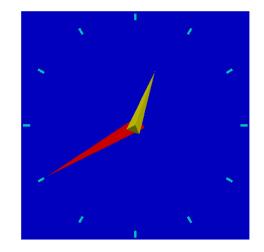


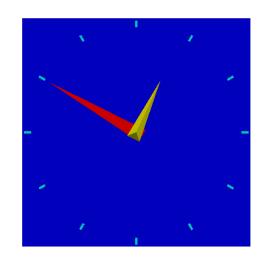
Time: 0.25 s



Time: 0.50 s







•*Motion06*: accelerated rotation motion (**<mvrotace** />) and accelerated circular motion (**<mvcirace** />).

```
- <motion>
  - <obj>
      <objreal ref="1"/>
      <objreal ref="3"/>
      <objreal ref="4"/>
      <begin mov="1" start="0"/>
    - <mvrotace id="1" duration="1000">
         <ace ang="9"/>
         <velini ang="-50"/>
         <axisp1 x="0" y="0" z="1.85"/>
         <axisp2 x="0" y="1" z="1.85"/>
      </mvrotace>
    </obj>
  - <objreal ref="5">
      <begin mov="1" start="0"/>
    - <mvcirace id="1" duration="1000">
         <ace ang="9"/>
         <velini ang="-50"/>
         <ref x="1.3" y="-0.7" z="1.85"/>
         <axisp1 x="0" y="0" z="1.85"/>
         <axisp2 x="0" y="1" z="1.85"/>
      </mvcirace>
    </objreal>
  </motion>
```

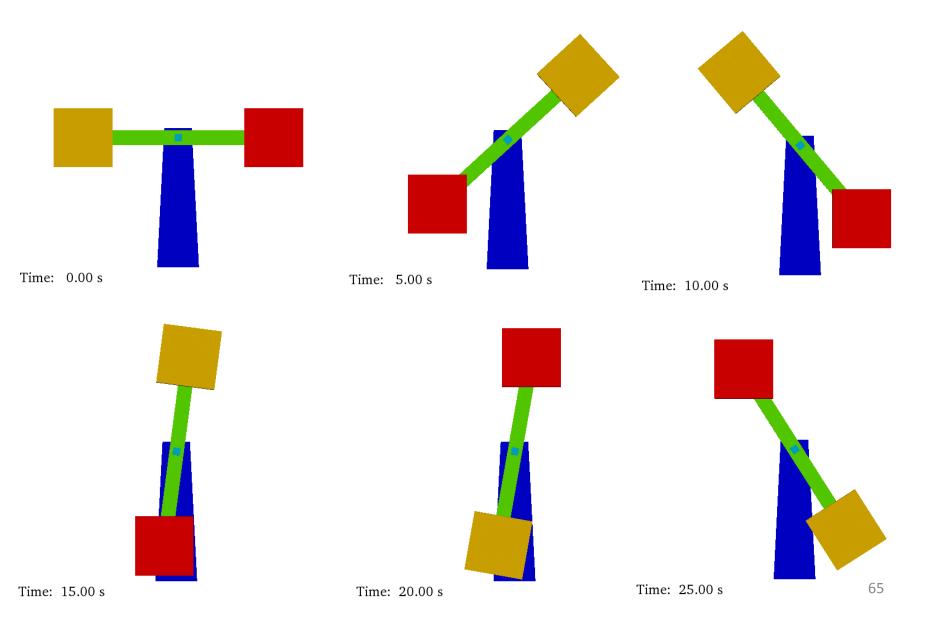
<mvrotace>: accelerated rotational movement

ace indicates the angular accelerationvelini indicates the initial angular velocityaxisp1 first point of the rotation axisaxisp2 second point of the rotation axis

<mvcirace>: accelerated circular movement

ace indicates the angular acceleration
ref indicates the point of the object that rotates
with the axis
velini indicates the initial angular velocity
axisp1 first point of the rotation axis
axisp2 second point of the rotation axis

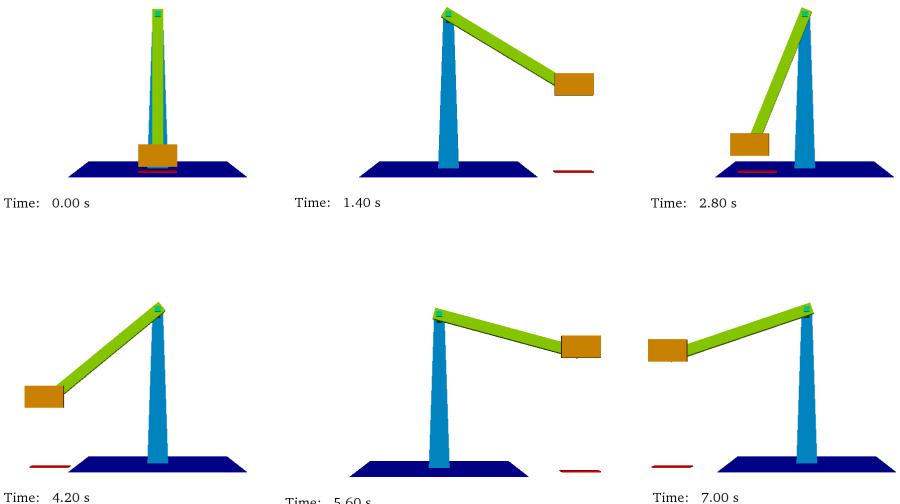
•*Motion06*: accelerated rotation motion (**<mvrotace** />) and accelerated circular motion (**<mvcirace** />).



•*Motion07*: sinusoidal movement (**<mvrectsinu** /**>**, **<mvrotsinu** /**>**, **<mvcirsinu** /**>**)

```
- <motion>
  - <objreal ref="4">
      <br/>
<br/>
begin mov="1" start="0"/>
                                                                                           <mvrectsinu>: sinusoidal rectilinear
                                               - <mvcirsinu id="2" duration="5" next="1">
    = <mvrotsinu id="1" duration="5" next="2">
                                                   <ref x="0" y="-0.7" z="0.2"/>
                                                                                           movement
        <axisp1 x="0" y="0" z="2.85"/>
                                                   <axisp1 x="0" y="0" z="2.85"/>
        <axisp2 x="0" y="1" z="2.85"/>
                                                   <axisp2 x="0" y="1" z="2.85"/>
        <freq v="0.2"/>
                                                                                           <mvrotsinu>: sinusoidal rotational
                                                   <freq v="0.4"/>
        <ampl v="60"/>
                                                   <ampl v="75"/>
                                                                                           movement
        < phase v="0"/>
                                                   <phase v="0"/>
      </mvrotsinu>
                                                 </mvcirsinu>
    - <mvrotsinu id="2" duration="5" next="1">
                                                                                           <mvcirsinu>: sinusoidal circular
                                               </objreal>
        <axisp1 x="0" y="0" z="2.85"/>
                                              - <objreal ref="6">
                                                                                           movement
        <axisp2 x="0" y="1" z="2.85"/>
                                                  <begin mov="1" start="0"/>
        <freq v="0.4"/>
                                                - <mvrectsinu id="1" duration="5" next="2">
        <ampl v="75"/>
                                                    <freq x="0.2" y="0" z="0"/>
                                                                                            axisp1 first point of the rotation axis
      </mvrotsinu>
                                                    <ampl x="2.30" y="0" z="0"/>
                                                                                             axisp2 second point of the axis
   </obireal>
                                                    hase x="0" y="0" z="0"/>
                                                                                            freq frequency
  - <objreal ref="5">
                                                  </mvrectsinu>
      <br/>
start="0"/>
                                                - <mvrectsinu id="2" duration="5" next="1">
                                                                                             ampl amplitude
    - <mvcirsinu id="1" duration="5" next="2"</p>
                                                    <freq x="0.4" y="0" z="0"/>
                                                                                            phase phase
        <ref x="0" y="-0.7" z="0.2"/>
                                                    <ampl x="2.55" y="0" z="0"/>
        <axisp1 x="0" y="0" z="2.85"/>
                                                    hase x="0" y="0" z="0"/>
        <axisp2 x="0" y="1" z="2.85"/>
                                                  </mvrectsinu>
        <freq v="0.2"/>
                                                </objreal>
        <ampl v="60"/>
                                              </motion>
        <phase v="0"/>
      </mvcirsinu>
```

•*Motion07*: sinusoidal movement (**<mvrectsinu** /**>**, **<mvrotsinu** /**>**, **<mvcirsinu** /**>**)



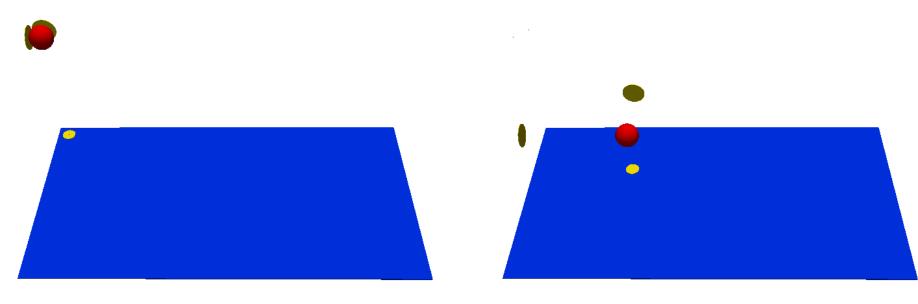
Time: 5.60 s

•*Motion08*: predefined movement with data from an external file (**<mvpredef** /**>** or **<mvfile** /**>**)

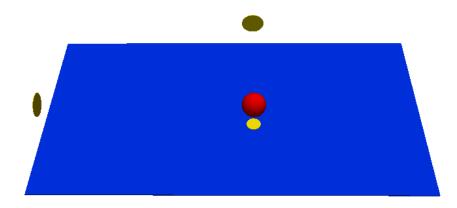
```
- <motion>
  - <objreal ref="200">
      <begin mov="1" start="0"/>
    - <mvpredef id="1" duration="10">
        <file name="motion08mov_f3.out" fields="4" fieldtime="0" fieldx="1" fieldy="2" fieldz="3"/>
      </mvpredef>
   </objreal>
                                                                                    <mvpredef /> or <mvfile />:
  - <objreal ref="150">
      <begin mov="1" start="0"/>
                                                                                    prescribed motion loaded from a file
    - <mvpredef id="1" duration="8" next="2">
        <file name="motion08mov_f3.out" fields="4" fieldtime="0" fieldx="1" fieldy="2"/>
                                                                                    name name of the file
      </mvpredef>
    - <mvrect id="2" duration="-1">
                                                                                    fields number of columns of the file
        <vel x="0" y="0" z="-0.02"/>
                                                                                    fieldtime column with time
      </mvrect>
                                                                                    fieldx column with X-position
   </objreal>
                                                                                    fieldy column with Y-position
  - <objreal ref="151">
      <begin mov="1" start="0"/>
                                                                                    filedz column with Z-position
    - <mvpredef id="1" duration="10">
        <file name="motion08mov_f3.out" fields="4" fieldtime="0" fieldx="1" fieldz="3"/>
      </mvpredef>
   </objreal>
                                                                             first field (or column) has reference "0"
  - <objreal ref="152">
                                                                             second field (or column) has reference "1"
      <begin mov="1" start="0"/>
    - <mvpredef id="1" duration="10">
        <file name="motion08mov_f3.out" fields="4" fieldtime="0" fieldy="2" fieldz="3"/>
      </mvpredef>
   </objreal>
```

</motion>

•*Motion08*: predefined movement with data from an external file (**<mvpredef** /**>** or **<mvfile** /**>**)



Time: 0.00 s



Time: 6.00 s

Time: 3.00 s



•*Motion09*: predefined movement with data from an external file (**<mvrotfile** /**>**)

<mvrotfile />: prescribed motion loaded from a file with degrees

name name of the file
axisp1 & axisp2 two points to define the axis of rotation

```
<motion>

<dojreal ref="1">

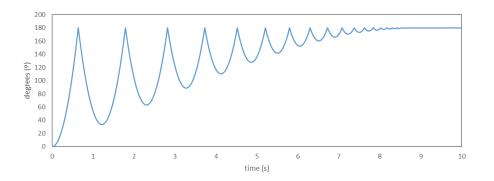
<degin mov="1" start="0" finish="100" />

<mvrotfile id="1" duration="9" next="2" anglesunits="degrees">

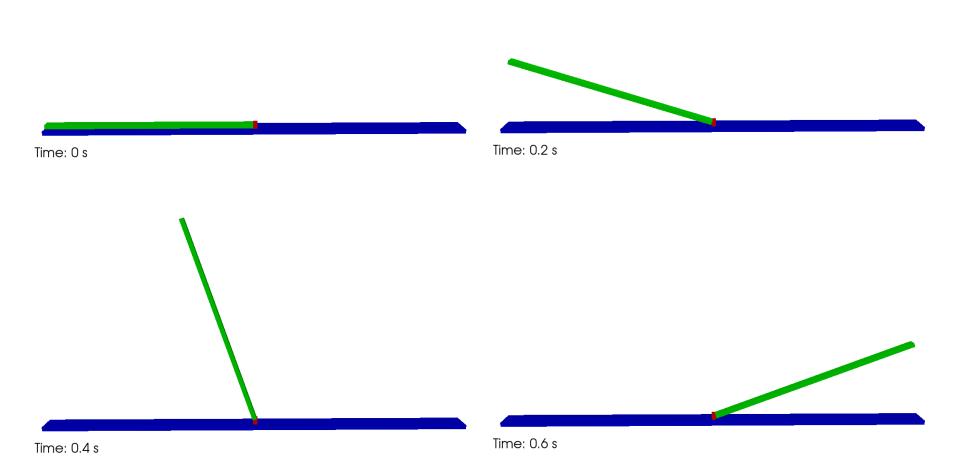
<file name="Motion09mov_deg.csv" />
<axisp1 x="1" y="1" z="0.03" />
<axisp2 x="1" y="-1" z="0.03" />
</mvrotfile>

</
```





•*Motion09*: predefined movement with data from an external file (**<mvrotfile** /**>**)



XML file EXECUTION-SPECIAL-WAVEPADDLES-PISTON

Generation of regular waves

<piston>

```
<mkbound value="10" comment="Mk-Bound of selected particles" />
<waveorder value="2" comment="Order wave generation 1:1st order, 2:2nd order (def=1)" />
<start value="0" comment="Start time (def=0)" />
<duration value="0" comment="Movement duration, Zero is the end of simulation (def=0)" />
<depth value="0.27" comment="Fluid depth (def=0)" />
<fixeddepth value="0" comment="Fluid depth without paddle (def=0)" />
<fixeddepth value="0" comment="Wave ment direction (def=(1,0,0))" />
<pistondir x="1" y="0" z="0" comment="Wave height" />
<waveheight value="0.1" comment="Wave period" />
square value="0" comment="Initial wave phase in function of PI (def=0)" />

<ramp value="0" comment="Periods of ramp (def=0)" />

<savemotion periods="24" periodsteps="20" xpos="-0.15"</p>
```

•*waveorder*: order of wave generation (1st order or 2nd order)

•*depth*: depth at front of the piston

•waveheight: wave height H

•waveperiod: wave period T

•ramp: number of periods to smooth the movement of the piston

•savemotion: saves theoretical results of elevation and orbital velocities at xpos and zpos

(being zpos=-depth of the measuring point)

XML file EXECUTION-SPECIAL-WAVEPADDLES-PISTON

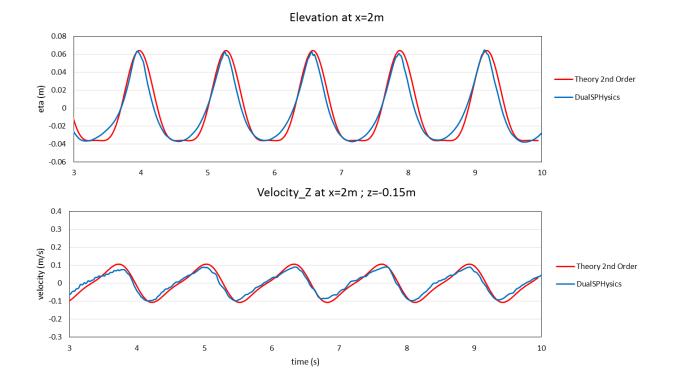
Generation of regular waves

- •*waveorder*: order of wave generation (1st order or 2nd order)
- •*depth*: depth at front of the piston
- •*waveheight*: wave height *H*
- •*waveperiod*: wave period *T*

```
•ramp: number of periods to smooth the movement of the piston (being zpos=-depth of the measuring point)
```

•savemotion: saves theoretical results of elevation and orbital velocities at xpos and zpos

<savemotion periods="24" periodsteps="20" xpos="2" zpos="-0.15"/>



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EXECUTION-SPECIAL-WAVEPADDLES-PISTON_SPECTRUM

Generation of irregular waves

```
<piston spectrum>
    <mkbound value="10" comment="Mk-Bound of selected particles" />
    <waveorder value="2" comment="Order wave generation 1:1st order, 2:2nd order (def=1)" />
    <start value="0" comment="Start time (def=0)" />
    <duration value="0" comment="Movement duration, Zero is the end of simulation (def=0)" />
    <depth value="0.27" comment="Fluid depth (def=0)" />
    <fixeddepth value="0" comment="Fluid depth without paddle (def=0)" />
    <pistondir x="1" y="0" z="0" comment="Movement direction (def=(1,0,0))" />
    <spectrum value="jonswap" comment="Spectrum type: jonswap,pierson-moskowitz" />
    <discretization value="stretched"</pre>
                comment="Spectrum discretization: regular, random, stretched, cosstretched (def=stretched)" />
    <waveheight value="0.1" comment="Wave height" />
    <waveperiod value="1.3" comment="Wave period" />
    <peakcoef value="3.3" comment="Peak enhancement coefficient (def=3.3)" />
    <waves value="128" comment="Number of waves to create irregular waves (def=50)" />
    <randomseed value="2" comment="Random seed to initialize a pseudorandom number generator" />
    <serieini value="2.8" comment="Initial time in irregular wave serie (def=0)" />
    <ramptime value="1" comment="Time of ramp (def=0)" />
    <savemotion time="50" timedt="0.05" xpos="2" zpos="-0.15"</pre>
                comment="Saves motion data. xpos and zpos are optional. zpos=-depth" />
    <saveserie timemin="0" timemax="1300" timedt="0.05" xpos="0" comment="Saves serie data (optional)" />
    <saveseriewaves timemin="0" timemax="1000" xpos="2" comment="Saves serie heights" />
</piston spectrum>
```

EXECUTION-SPECIAL-WAVEPADDLES-PISTON_SPECTRUM

Generation of irregular waves

•*waveorder*: order of wave generation (1st order or 2nd order)

•*spectrum*: type of spectrum (Jonswap or Pierson-Moskowitz)

•*waveheight*: significant wave height H_s

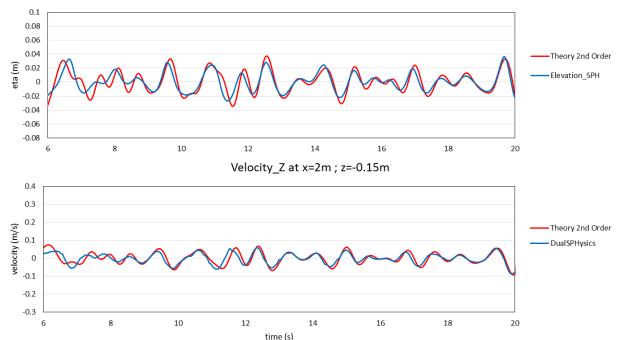
•*waveperiod*: peak wave period T_p

•serieini: initial series of the irregular train is chosen from "WavePaddle_mkb0010_Serie.csv"

•*ramptime*: time to slowly start a smoothed movement of the piston (being *zpos=-depth* of the measuring point)

•*savemotion*: saves theoretical results of elevation and orbital velocities at *xpos* and *zpos*

<savemotion time="50" timedt="0.05" xpos="2" zpos="-0.15" />



Elevation at x=2m

EXECUTION-PARAMETERS

Parameters for execution in DualSPHysics

<parameters>

<pre> <parameter< pre=""></parameter<></pre>	key="PosDouble" value="1" comment="Precision in particle interaction 0:Simple, 1:Double, 2:Uses and saves double (default=0)" />
<parameter< td=""><td>key="StepAlgorithm" value="1" comment="Step Algorithm 1:Verlet, 2:Symplectic (default=1)" /></td></parameter<>	key="StepAlgorithm" value="1" comment="Step Algorithm 1:Verlet, 2:Symplectic (default=1)" />
<parameter< td=""><td>key="VerletSteps" value="40" comment="Verlet only: Number of steps to apply Euler timestepping (default=40)" /></td></parameter<>	key="VerletSteps" value="40" comment="Verlet only: Number of steps to apply Euler timestepping (default=40)" />
<parameter< td=""><td>key="Kernel" value="1" comment="Interaction Kernel 1:Cubic Spline, 2:Wendland (default=2)" /></td></parameter<>	key="Kernel" value="1" comment="Interaction Kernel 1:Cubic Spline, 2:Wendland (default=2)" />
<pre><parameter< pre=""></parameter<></pre>	key="ViscoTreatment" value="1" comment="Viscosity formulation 1:Artificial, 2:Laminar+SPS (default=1)" />
<parameter< td=""><td>key="Visco" value="0.1" comment="Viscosity value" /></td></parameter<>	key="Visco" value="0.1" comment="Viscosity value" />
<parameter< td=""><td>key="ViscoBoundFactor" value="1" comment="Multiply viscosity value with boundary (default=1)" /></td></parameter<>	key="ViscoBoundFactor" value="1" comment="Multiply viscosity value with boundary (default=1)" />
<parameter< td=""><td>key="DeltaSPH" value="0" comment="DeltaSPH value, 0.1 is the typical value, with 0 disabled (default=0)" /></td></parameter<>	key="DeltaSPH" value="0" comment="DeltaSPH value, 0.1 is the typical value, with 0 disabled (default=0)" />
<parameter< td=""><td>key="#Shifting" value="0" comment="Shifting mode 0:None, 1:Ignore bound, 2:Ignore fixed, 3:Full (default=0)" /></td></parameter<>	key="#Shifting" value="0" comment="Shifting mode 0:None, 1:Ignore bound, 2:Ignore fixed, 3:Full (default=0)" />
<parameter< td=""><td><pre>key="#ShiftCoef" value="-2" comment="Coefficient for shifting computation (default=-2)" /></pre></td></parameter<>	<pre>key="#ShiftCoef" value="-2" comment="Coefficient for shifting computation (default=-2)" /></pre>
<parameter< td=""><td>key="#ShiftTFS" value="1.5" comment="Threshold to detect free surface. Typically 1.5 for 2D and 2.75 for 3D (default=0)" /></td></parameter<>	key="#ShiftTFS" value="1.5" comment="Threshold to detect free surface. Typically 1.5 for 2D and 2.75 for 3D (default=0)" />
<parameter< td=""><td><pre>key="RigidAlgorithm" value="1" comment="Rigid Algorithm 1:SPH, 2:DEM (default=1)" /></pre></td></parameter<>	<pre>key="RigidAlgorithm" value="1" comment="Rigid Algorithm 1:SPH, 2:DEM (default=1)" /></pre>
<parameter< td=""><td>key="FtPause" value="0.0" comment="Time to freeze the floatings at simulation start (warmup) (default=0)" units_comment="seconds" /></td></parameter<>	key="FtPause" value="0.0" comment="Time to freeze the floatings at simulation start (warmup) (default=0)" units_comment="seconds" />
<parameter< td=""><td>key="CoefDtMin" value="0.05" comment="Coefficient to calculate minimum time step dtmin=coefdtmin*h/speedsound (default=0.05)" /></td></parameter<>	key="CoefDtMin" value="0.05" comment="Coefficient to calculate minimum time step dtmin=coefdtmin*h/speedsound (default=0.05)" />
<parameter< td=""><td>key="#DtIni" value="0.0001" comment="Initial time step (default=h/speedsound)" units_comment="seconds" /></td></parameter<>	key="#DtIni" value="0.0001" comment="Initial time step (default=h/speedsound)" units_comment="seconds" />
<parameter< td=""><td>key="#DtMin" value="0.00001" comment="Minimum time step (default=coefdtmin*h/speedsound)" units_comment="seconds" /></td></parameter<>	key="#DtMin" value="0.00001" comment="Minimum time step (default=coefdtmin*h/speedsound)" units_comment="seconds" />
<parameter< td=""><td><pre>key="#DtFixed" value="DtFixed.dat" comment="Dt values are loaded from file (default=disabled)" /></pre></td></parameter<>	<pre>key="#DtFixed" value="DtFixed.dat" comment="Dt values are loaded from file (default=disabled)" /></pre>
<parameter< td=""><td>key="DtAllParticles" value="0" comment="Velocity of particles used to calculate DT. 1:All, 0:Only fluid/floating (default=0)" /></td></parameter<>	key="DtAllParticles" value="0" comment="Velocity of particles used to calculate DT. 1:All, 0:Only fluid/floating (default=0)" />
<parameter< td=""><td>key="TimeMax" value="1.5" comment="Time of simulation" units_comment="seconds" /></td></parameter<>	key="TimeMax" value="1.5" comment="Time of simulation" units_comment="seconds" />
<parameter< td=""><td>key="TimeOut" value="0.01" comment="Time out data" units_comment="seconds" /></td></parameter<>	key="TimeOut" value="0.01" comment="Time out data" units_comment="seconds" />
<parameter< td=""><td>key="IncZ" value="1" comment="Increase of Z+" units_comment="decimal" /></td></parameter<>	key="IncZ" value="1" comment="Increase of Z+" units_comment="decimal" />
<parameter< td=""><td>key="PartsOutMax" value="1" comment="Allowed %/100 of fluid particles out the domain (default=1)" units_comment="decimal" /></td></parameter<>	key="PartsOutMax" value="1" comment="Allowed %/100 of fluid particles out the domain (default=1)" units_comment="decimal" />
<parameter< td=""><td>key="RhopOutMin" value="700" comment="Minimum rhop valid (default=700)" units_comment="kg/m^3" /></td></parameter<>	key="RhopOutMin" value="700" comment="Minimum rhop valid (default=700)" units_comment="kg/m^3" />
<parameter< td=""><td>key="RhopOutMax" value="1300" comment="Maximum rhop valid (default=1300)" units_comment="kg/m^3" /></td></parameter<>	key="RhopOutMax" value="1300" comment="Maximum rhop valid (default=1300)" units_comment="kg/m^3" />

EXECUTION-PARAMETERS

Double precision

<parameter key="PosDouble" value="2" comment="Precision in particle interaction 0:Simple, 1:Double, 2:Uses and saves double" />

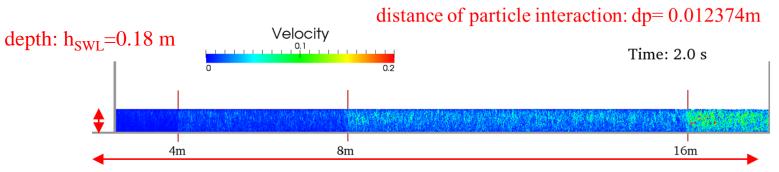
DualSPHysics v4.0 includes now implementation with double precision. Precision in particle interaction (the most time consuming part) can be:

0: particle interaction is performed using simple precision for variables of position

Necessary when "dp" is much smaller than size of the domain:

1: particle interaction is performed using double precision for variables of position but final position is stored using simple precision

2: particle interaction is performed using double precision for variables of position and final position is stored using double precision



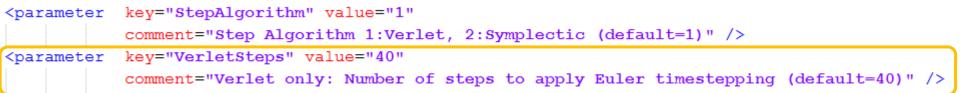
length=18 m

The problems of precision mainly appear when the domain is huge in comparison to the distance of interaction between particles

length>depth>>dp

EXECUTION-PARAMETERS

Time integrator scheme



Verlet algorithm

$$\boldsymbol{v}_{a}^{n+1} = \boldsymbol{v}_{a}^{n-1} + 2\Delta t \boldsymbol{F}_{a}^{n}$$
$$\boldsymbol{r}_{a}^{n+1} = \boldsymbol{r}_{a}^{n} + \Delta t \boldsymbol{V}_{a}^{n} + 0.5\Delta t^{2} \boldsymbol{F}_{a}^{n}$$
$$\rho_{a}^{n+1} = \rho_{a}^{n-1} + 2\Delta t D_{a}^{n}$$

once every M time steps

$$\boldsymbol{v}_{a}^{n+1} = \boldsymbol{v}_{a}^{n} + \Delta t \boldsymbol{F}_{a}^{n}$$
$$\boldsymbol{r}_{a}^{n+1} = \boldsymbol{r}_{a}^{n} + \Delta t \boldsymbol{V}_{a}^{n} + 0.5 \Delta t^{2} \boldsymbol{F}_{a}^{n}$$
$$\rho_{a}^{n+1} = \rho_{a}^{n} + \Delta t D_{a}^{n}$$

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<parameter key="StepAlgorithm" value="2"</pre> comment="Step Algorithm 1:Verlet, 2:Symplectic (default=1)" />

1

Symplectic algorithm

$$r_{a}^{n+\frac{1}{2}} = r_{a}^{n} + \frac{\Delta t}{2} v_{a}^{n} \qquad v_{a}^{n+1} = v_{a}^{n+\frac{1}{2}} + \frac{\Delta t}{2} F_{a}^{n+\frac{1}{2}}$$

$$\rho_{a}^{n+\frac{1}{2}} = \rho_{a}^{n} + \frac{\Delta t}{2} D_{a}^{n} \qquad r_{a}^{n+1} = r_{a}^{n+\frac{1}{2}} + \frac{\Delta t}{2} v_{a}^{n+1}$$
Predictor Corrector



Kernel function

$$\frac{d\boldsymbol{v}_a}{dt} = -\sum_b m_b \left(\frac{P_b + P_a}{\rho_b \cdot \rho_a} + \Pi_{ab}\right) \nabla_a W_{ab} + \boldsymbol{g}$$

<parameter key="Kernel" value="1"</pre>

comment="Interaction Kernel 1:Cubic Spline, 2:Wendland (default=2)" />

Cubic Spline

$$W(r,h) = \alpha_D \begin{cases} 1 - \frac{3}{2}q^2 + \frac{3}{4}q^3 & 0 \le q \le 1 \\ \frac{1}{4}(2-q)^3 & 1 \le q \le 2 \\ 0 & q \ge 2 \end{cases}$$

where α_D is equal to $10/7\pi h^2$ in 2-D and $1/\pi h^3$ in 3-D

Wendland

$$W(r,h) = \alpha_D \left(1 - \frac{q}{2}\right)^4 \left(2q + 1\right) \qquad 0 \le q \le 2$$

where α_D is equal to $7/4\pi h^2$ in 2-D and $21/16\pi h^3$ in 3-D



Viscosity treatment

<pre><parameter< pre=""></parameter<></pre>	key="ViscoTreatment" value="1"					
	<pre>comment="Viscosity formulation 1:Artificial, 2:Laminar+SPS (default=1)" /></pre>					
<parameter< td=""><td colspan="3"><pre>key="Visco" value="0.02" comment="Viscosity value" /></pre></td></parameter<>	<pre>key="Visco" value="0.02" comment="Viscosity value" /></pre>					
<parameter< td=""><td colspan="3">key="ViscoBoundFactor" value="1"</td></parameter<>	key="ViscoBoundFactor" value="1"					
	<pre>comment="Multiply viscosity value with boundary (default=1)" /></pre>					

$$\frac{d\boldsymbol{v}_{a}}{dt} = -\sum_{b} m_{b} \left(\frac{P_{b} + P_{a}}{\rho_{b} \cdot \rho_{a}} + \Pi_{ab} \right) \nabla_{a} W_{ab} + \boldsymbol{g}$$
$$\Pi_{ab} = \begin{cases} \frac{-\boldsymbol{\alpha} \overline{\boldsymbol{c}_{ab}} \mu_{ab}}{\overline{\rho_{ab}}} & \boldsymbol{v}_{ab} \cdot \boldsymbol{r}_{ab} < 0\\ 0 & \boldsymbol{v}_{ab} \cdot \boldsymbol{r}_{ab} > 0 \end{cases}$$

 α =0.01 for wave tanks higher values of α for dam-break (depends on dp)



Viscosity treatment

<parameter< th=""><th colspan="6">key="ViscoTreatment" value="1"</th></parameter<>	key="ViscoTreatment" value="1"					
	<pre>comment="Viscosity formulation 1:Artificial, 2:Laminar+SPS (default=1)" /></pre>					
<parameter< th=""><th colspan="4"><pre>key="Visco" value="0.02" comment="Viscosity value" /></pre></th></parameter<>	<pre>key="Visco" value="0.02" comment="Viscosity value" /></pre>					
<pre><parameter< pre=""></parameter<></pre>	key="ViscoBoundFactor" value="1"					
	<pre>comment="Multiply viscosity value with boundary (default=1)" /></pre>					

$$\frac{d\boldsymbol{v}_{a}}{dt} = -\sum_{b} m_{b} \left(\frac{P_{b} + P_{a}}{\rho_{b} \cdot \rho_{a}} + \Pi_{ab} \right) \nabla_{a} W_{ab} + \boldsymbol{g}$$
$$\Pi_{ab} = \begin{cases} \frac{-\alpha \, \overline{c_{ab}} \mu_{ab}}{\overline{\rho_{ab}}} & \boldsymbol{v}_{ab} \cdot \boldsymbol{r}_{ab} < 0\\ 0 & \boldsymbol{v}_{ab} \cdot \boldsymbol{r}_{ab} > 0 \end{cases}$$

 α_{FF} for interaction fluid-fluid α_{FB} for interaction fluid-boundary α_{FB} = ViscoBoundFactor $\cdot \alpha_{FF}$



Viscosity treatment

<p< th=""><th>parameter</th><th colspan="5">key="ViscoTreatment" value="2"</th></p<>	parameter	key="ViscoTreatment" value="2"				
		<pre>comment="Viscosity formulation 1:Artificial, 2:Laminar+SPS (default=1)" /></pre>				
<r< td=""><td>parameter</td><td>key="Visco" value="0.000001" comment="Viscosity value" units_comment="m^2/s"</td><td>" /></td></r<>	parameter	key="Visco" value="0.000001" comment="Viscosity value" units_comment="m^2/s"	" />			

$$\frac{d\boldsymbol{v}_{a}}{dt} = -\sum_{b} m_{b} \left(\frac{P_{b}}{\rho_{b}^{2}} + \frac{P_{a}}{\rho_{a}^{2}} \right) \nabla_{a} W_{ab} + \boldsymbol{g} + \sum_{b} m_{b} \left(\frac{4 \upsilon_{a} r_{ab} \cdot \nabla_{a} W_{ab}}{(\rho_{a} + \rho_{b})(r_{ab}^{2} + \eta^{2})} \right) \boldsymbol{v}_{ab} + \sum_{b} m_{b} \left(\frac{\vec{\tau}_{ij}^{b}}{\rho_{b}^{2}} + \frac{\vec{\tau}_{ij}^{a}}{\rho_{a}^{2}} \right) \nabla_{a} W_{ab}$$

 $\boldsymbol{v_o}$ is kinematic viscosity (typically 10⁻⁶ m²s for water

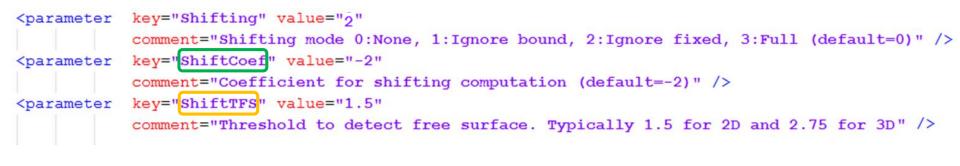


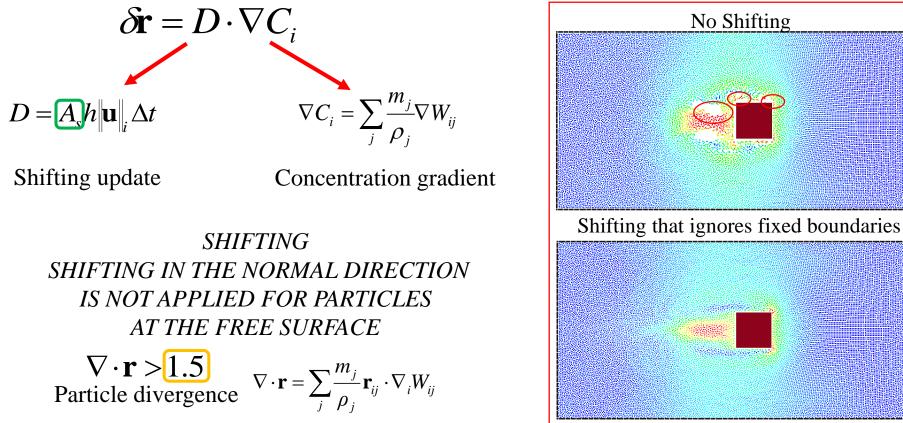
DeltaSPH formulation

$$\frac{d\rho_a}{dt} = \sum_b m_b \mathbf{v}_{ab} \cdot \nabla_a W_{ab} + 2 \partial h \sum_b m_b \overline{c_{ab}} \left(\frac{\rho_a}{\rho_b} - 1\right) \frac{\mathbf{r}_{ab}}{\mathbf{r}_{ab}^2 + \eta^2} \cdot \nabla_a W_{ab}$$



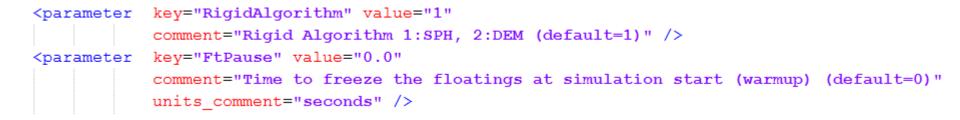
Shifting algorithm

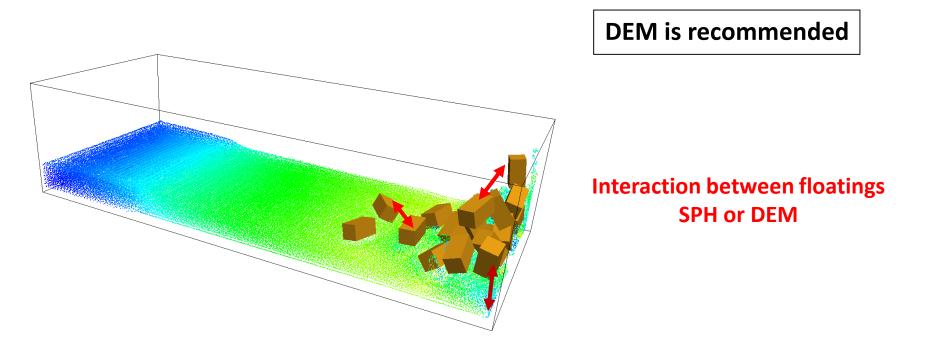






Interaction between solids





Interaction between floating and bottom SPH or DEM



Time step computation

<parameter< th=""><th>key="CoefDtMin" value="0.05"</th></parameter<>	key="CoefDtMin" value="0.05"
	comment="Coefficient to calculate minimum time step dtmin=coefdtmin*h/speedsound />
<parameter< td=""><td><pre>key="#DtIni" value="0.0001"</pre></td></parameter<>	<pre>key="#DtIni" value="0.0001"</pre>
	<pre>comment="Initial time step (default=h/speedsound" units_comment="seconds" /></pre>
<parameter< td=""><td>key="#DtMin" value="0.00001"</td></parameter<>	key="#DtMin" value="0.00001"
	<pre>comment="Minimum time step (default=coefdtmin*h/speedsound)" units_comment="seconds" /></pre>
<pre><parameter< pre=""></parameter<></pre>	<pre>key="#DtFixed" value="DtFixed.dat"</pre>
	<pre>comment="Dt values are loaded from file (default=disabled)" /></pre>

$$\Delta t_{minimum} = 0.05 \cdot \frac{h}{c_s}$$

$$\Delta t_{init} = \frac{h}{c_s}$$

<parameter key="DtAllParticles"</pre>

comment="Velocity of particles used to calculate DT. 1:All, 0:Only fluid/floating (default=0)" />

$$\Delta t = 0.3 \cdot \min(\Delta t_f, \Delta t_{cv})$$

$$\Delta t_f = \min\left(\sqrt{\frac{h}{|f_a|}}\right)$$

$$\Delta t_{cv} = \min_a \frac{h}{c_s + \max_b \left|\frac{h\mathbf{v}_{ab} \cdot \mathbf{r}_{ab}}{\mathbf{r}_{ab}^2}\right|}$$

$$0: b \in \text{fluid/floating}$$

$$OR$$

$$1: b \in \text{fluid/floating} + \text{boundaries}$$

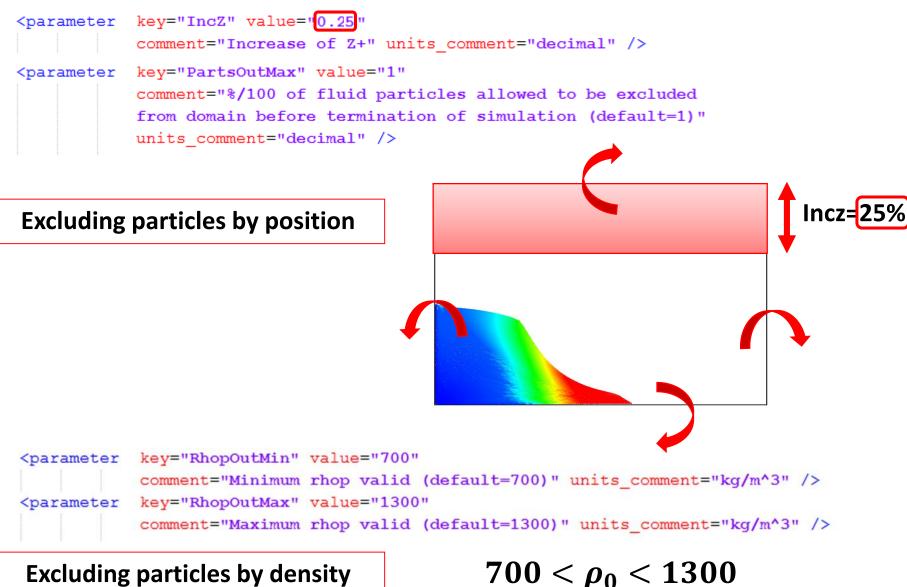


Physical time and frequency to store data

<parameter< th=""><th>key="TimeMax"</th><th>value="10"</th></parameter<>	key="TimeMax"	value="10"
	comment="Time	of simulation" units_comment="seconds" />
<parameter< td=""><td>key="TimeOut"</td><td>value="0.1"</td></parameter<>	key="TimeOut"	value="0.1"
	comment="Time	out data" units_comment="seconds" />

Number of output files	= TimeMax / TimeOut
	= 10 / 0.1 = 100 files

EXECUTION-PARAMETERS XML file

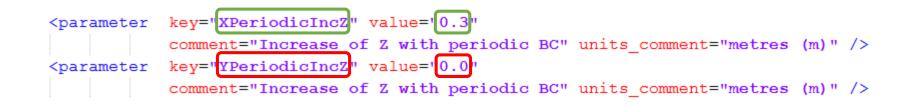


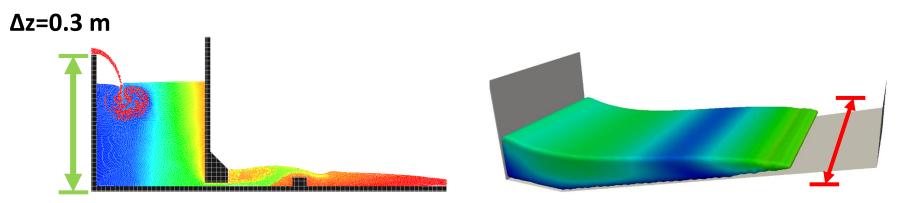
Excluding particles by density

88



Periodicity





Δz=0 m







ce	case> <casedef></casedef>
	<pre><caseder> </caseder></pre>
<pre>ccases> ccasedef></pre>	<pre>constantSantSantSantSantSantSantSantSantSant</pre>
<pre><coaseur> </coaseur></pre>	<pre><gravity comment="Gravitational acceleration" units_comment="m/a^2" x="0" y="0" z="-9.81"></gravity></pre>
<lattice bound="1" fluid="1"></lattice>	<cflnumber comment="Coefficient to multiply Dt" value="0.2"></cflnumber>
<pre><gravity comment="Gravitational acceleration" units_comment="m/s^2" x="0" y="0" z="-9.61"></gravity></pre>	<pre>chswl value="0" auto="true" comment="Maximum still water level to calculate speedofsound using coefsound" units_comment="metres (m)" /></pre>
<pre><cfinumber comment="Coefficient to multiply Dt" value="0.2"></cfinumber></pre>	<pre>cspeedsystem value="10" auto=""true" comment="Maximum system speed (by default the dam-break propagation is used)" /> foreframed "bluer100" auto="true" comment="Maximum system speed (by default the dam-break propagation is used)" /> </pre>
<pre>chsul value*"0" auto*"true" comment="Maximum still water level to calculate speedofsound using conformat" units_comment="matters (m)" /> capedayster value*"0" auto*"true" comment="Maximum systems speed (by default the dam-break propagation is used) /> </pre>	<pre><coefsound comment="Coefficient to multiply speedsystem" value="0"></coefsound> <pre><apredsound auto="true" comment="Speed of sound to use in the simulation (by default speedsfound+coefsound+speedsystem)" value="0"></apredsound></pre></pre>
<pre>cspecusystem value='0' auto='true' comment='maximum system specusystem '0' init the dam-break propagation is used)' /> ccoefsound value='0' comment='Coefficient to multiply specisystem' /> </pre>	<pre>cypetations value = 0 also = true comment= speed of sound to use in the simulation (p detail speedorsound-operative) // <coeff <="" comment="Coefficient to calculate the smoothing length (B-coefficient*sqrt(3*dp^2) in 30)" pre="" value="1.0°"></coeff></pre>
<pre><pre><aprecision auto="true" comment="Speed of sound to use in the simulation (by default speedofsound=coefsound=speedsystem)" value="0"></aprecision></pre></pre>	<pre><qamma comment="Politropic constant for water used in the state equation" value="7"></qamma></pre>
<pre><coefh comment="Coefficient to calculate the smoothing length (H=coefficient*sqrt(3*dp^2) in 3D)" value="1.0"></coefh></pre>	<rhop0 comment="Reference density of the fluid" units_comment="kg/m3" value="1000"></rhop0>
<gamma comment="Politropic constant for water used in the state equation" value="7"></gamma>	
<pre>crhop0 value="1000" comment="Reference density of the fluid" units_comment="kg/m3" /> </pre>	<mkconfig boundcount="240" fluidcount="10"></mkconfig>
 <pre>cmkconfig boundcount="240" fluidocunt="10" /></pre>	<geometry></geometry>
<pre><geometry></geometry></pre>	<pre><definition dp="0.01" units_comment="metres (m)"> <pointmin x="-1" y="0" z="-1"></pointmin></definition></pre>
<definition dp="0.01" units_comment="metres (m)"></definition>	<pre><pointmin s.5'="" x="-ir" y="0 z=-ir /> <pointmin x="></pointmin> <pointmin x="s.5" y="0" z="s.5"></pointmin></pre>
<pre><pointmin x="-1" y="0" z="-1"></pointmin></pre>	<pre></pre>
<pre><pre>cpcintmax x="4.5" y="0" z="3.5" /></pre></pre>	<commands></commands>
	<pre><mainlist></mainlist></pre>
<pre><commands> </commands></pre>	<setdrawmode mode="full"></setdrawmode>
<pre><mainist> < setdrawmode mode*"full" /></mainist></pre>	<setmkfluid mk="0"></setmkfluid>
<pre><setukfuld full="" meter=""></setukfuld> <setukfuld meter=""></setukfuld></pre>	<drawbox></drawbox>
	<boxfill>solid</boxfill>
<boxfill>molid</boxfill>	<pre><pre>cpoint x="0" y="-1" z="0" /></pre></pre>
<pre><pre>cpoint x="0" y="-1" z="0" /></pre></pre>	<pre><size x="1" y="2" z="2"></size></pre>
<size x="1" y="2" z="2"></size>	
	<pre><aetmkbound mk="0"></aetmkbound></pre>
<setakbound mk="0"></setakbound>	<pre><drawbox></drawbox></pre>
<pre><drawbox></drawbox></pre>	<pre>cboxfill>bottom left right front back cpoint x="0" y=-1" z=="0" /></pre>
<pre> clostillabottom loft right front backcopint x="0" y="1" z=z"" /></pre>	<pre>cpoint x="u" y="-i" z="u" /> <sit '="" =="" x="i" y="' z="></sit></pre>
<pre><pre><pre><pre><pre><pre><pre><pre></pre></pre></pre></pre></pre></pre></pre></pre>	<pre><dicale #-4"="" l="3" y="2"></dicale> </pre>
	<execution></execution>
<#Regulion>	<pre><pre>cparameters></pre></pre>
<pre><pre>cparameters></pre></pre>	<pre><parameter comment="Step Algorithm 1:Verlet, 2:Symplectic (default=1)" key="StepAlgorithm" value="1"></parameter></pre>
<pre>cparameter key="stepAlgorithm" value="!" comment="Step Algorithm inverter, 2:Symplectic (default=1)" /> cparameter key="valueSteps" value="d" comment="Value" value" value of steps to apply Euler timestepping (default=40)" /></pre>	<pre><parameter comment="Verlet only: Number of steps to apply Euler timestepping (default=40)" key="VerletSteps" value="40"></parameter></pre>
<pre><pre><pre><pre><pre><pre><pre><pre></pre></pre></pre></pre></pre></pre></pre></pre>	<pre><pre>cparameter key="Kernel" value="2" comment="Interaction Kernel 1:Cubic Spline, 2:Wendland (default=2)" /></pre></pre>
<pre><pre>cparameter key="ViscoTreatment" value="1" comment="ViscoSity formulation 1:Artificial, 2:LaminartSPS (default=1)" /></pre></pre>	<pre><parameter comment="Viscosity formulation 1:Artificial, 2:Laminar+SFS (default=1)" key="ViscoTreatment" value="1"></parameter></pre>
<pre><pre><pre><pre>cparameter key="Visco" value="0.02" comment="Viscosity value" /></pre></pre></pre></pre>	<pre>cparameter key="Visco" value="0.02" comment="Viscosity value"/> cparameter key="Viscosity value"/> cparameter key="V</pre>
<pre><pre>cparameter key="ViscoBoundFactor" value="1" comment="Multiply viscosity value with boundary (default=1)" /></pre></pre>	<pre>cparameter key="ViscoBoundFactor" value="1" comment="Multiply viscosity value with boundary (default=1)" /></pre>
<pre><pre><cpre>areter key="DeltaSPH" value="0" comment="DeltaSPH value, 0.1 is the typical value, with 0 disabled (default=0)" /></cpre></pre></pre>	<pre>cparameter kmy="bettash" value="0" comment="bettash" value, 0.1 is the typical value, vitor o insuled (default=0)" /> <pre><pre><pre><pre><pre><pre><pre><pre></pre></pre></pre></pre></pre></pre></pre></pre></pre>
<pre><pre><pre><pre><pre><pre><pre><pre></pre></pre></pre></pre></pre></pre></pre></pre>	<pre>cprimeter key="shifting wate" = Comment= shifting mode / index / index / index / index / index / // // // // // // // // // // // //</pre>
$c_{parameter} = k_{parameter} = k_{parameter} = c_{parameter} = c_{parameter} = k_{parameter} = k_{parameter$	<pre>cprameter key="#shiftTFS" value="1.5" comment="Threshold to detect free surface. Typically 1.5 for 2D and 2.75 for 3D (default=0)" /></pre>
<pre>cparameter key="#ighitTF5" value="1.5" comment="Threshold to detext free surface. Typically 1.5 for 2D and 2.75 for 3D (default=0)= /> cparameter key="kighidhoprithm 'value="l' comment="Righidhoprithm 1:5FH, 2:DEM (default=1)= /></pre>	<pre><pre>cyparameter key="RigidAlgorithm" value="1" comment="Rigid Algorithm 1:SPH, 2:DEM (default=1)" /></pre></pre>
<pre>cparameter key="filesses" value="1" comment- high Adjoints issn; files (usualion start (warmup) (default=0)" units_comment="seconds" /></pre>	<pre>cparameter key="FtFause" value="0.0" comment="Time to freeze the floatings at simulation start (warmup) (default=0)" units_comment="seconds" /></pre>
<pre>cparameter key="CoefDMMin" value="0.05" comment="Coefficient to calculate minimum time step dtmin=coefdtemin=h/speedsound (default=0.05)" /></pre>	<pre><pre>coefDtMin" value="0.05" comment="Coefficient to calculate minimum time step dtmin=coefdtemin*h/speedsound (default=0.05)" /></pre></pre>
<pre><pre>comment="seconds" /></pre></pre>	<pre><pre>cparameter key="#DtIni" value="0.0001" comment="Initial time step (default=h/speedsound" units_comment="seconds" /></pre></pre>
<pre><pre><pre>cparameter key="#DtMin" value="0.00001" comment="Minimum time step (default=coefdtmin*h/speedsound)" units_comment="seconds" /></pre></pre></pre>	<pre><pre><pre>cparameter key="#DtMin" value="0.00001" comment="Minimum time step (default=coefdtmin*h/speedsound)" units_comment="seconds" /></pre></pre></pre>
<pre><pre><pre>cparameter key="#DtFixed" value="DtFixed.dat" comment="Dt values are loaded from file (default=disabled)" /></pre></pre></pre>	<pre><parameter comment="Dt values are loaded from file (default=disabled)" key="#DtFixed" value="DtFixed.dat"></parameter></pre>
<pre><pre><pre><pre><pre><pre><pre><pre></pre></pre></pre></pre></pre></pre></pre></pre>	<pre><parameter comment="Velocity of particles used to calculate DT. 1:All, 0:Only fluid/floating (default=0)" key="DtAllParticles" value="0"></parameter></pre>
<pre>cparameter key="fineMax" value="0.72" comment="fine of simulation" units_comment="seconds" /></pre>	<pre><pre>cparameter key="TimeMax" value="0.72" comment="Time of simulation" units_comment="seconds" /></pre></pre>
<pre>cparameter key="TimeOut" value="1.01" comment="Time out data" units_comment="seconds" /> cparameter key=TimeOut" value="1" comment="Increase of E** units_comment="decimal/></pre>	<pre>cparameter key="TimeOut" value="0.01" comment="Time out data" units_comment="seconds" /></pre>
<pre>cparameter key="ning: value=1" comment="inforcesse of i+" units_comment="decimal" /> cparameter key="Farthoutkax" value=1" comment="Allowed 4/100 of fluid particles out the domain (default=1)" units_comment="decimal" /></pre>	<pre>cparameter key="lnot" value="1" comment="increase of 5+" units comment="decimal" /> comment="Basedouble" molecular and the second second</pre>
<pre>cparameter xey="ratesouteax" value="10" comment="Allowed #/loo i fuid particles out the domain (default=70)" (default=700)" units_comment="kg/main/s/> cparameter xey="Rhopothin" value="10" comment="kg/main/s/</pre>	<pre>cparameter key="PartsOutMax" value="1" comment="Allowed %/100 of fluid particles out the domain (default=1)" units_comment="decimal" /> comments="Multiplus" units_comment="Multiplus" the units_comment="Multiplus"</pre>
<pre><pre><pre><pre><pre><pre><pre><pre></pre></pre></pre></pre></pre></pre></pre></pre>	<pre>cparameter key="RhopOutKin" value="700" comment="Minismu rhop valid (default=700)" unitg_comment="Kg/M3" /> cparameter key="RhopOutKin" value="700" comment="Kg/M3" (default=1300)" unitg_comment="Kg/M3" /> </pre>
	<pre>cparameter key="KnopputKax" value= 1000" comment= maximum thop value (dereart=1000) units_comment= kd/mp //</pre>
<pre>c/executions</pre>	<pre><pre>cparticles np="21001" nbf="1001" mkboundfirst="11" mkfluidfirst="1"></pre></pre>
	<pre>cfixed mkbound="0" mk="11" begin="0" count="1001" /></pre>
	<fluid begin="1001" count="20000" mk="1" mkfluid="0"></fluid>
	<pre><constants></constants></pre>
	<pre><gravity units_comment="m/s^2" x="0" y="0" z="-9.81"></gravity></pre>
	<cflnumber value="0.2"></cflnumber>
	<ganma value="7"></ganma>
	<pre><rhop0 units_comment="kg/m3" value="1000"></rhop0></pre>
	<pre><dp units_comment="metres (m)" value="0.01"></dp></pre>
	<h units_comment="metres (m)" value="1.4142135624E-002"></h>
	<pre><b units_comment="metres (m)" value='1.11553714998+006"'></pre>
	<pre>cmassbound value*1.0000000008-001" units_comment="kg" /> cmassfluid value*1.000000008-001" units_comment="kg" /></pre>
	<pre>cmassium value-1.000000000-001 units_comment=rkg /> </pre>
	(notion />

</case>

Summary of the number of created particles and computed constants

<particles np="21001" nb="1001" nbf="1001" mkboundfirst="11" mkfluidfirst="1">



YOU SHOULD ALWAYS CHECK Case_All.vtk, Case_Bound.vtk, Case_Fluid.vtk

Run.out

Text file with execution log

	1	DualSPHysics v4 (10-11-2015)	ר	
es	2			
	з	[Select CUDA Device]		
	4	Device 0: "GeForce GTX 590"		
	5	Compute capability:	2.0	
	6	Multiprocessors:	16 (512 cores)	
	7	Memory global:	1536 MB	
	8	Clock rate:	1.23 GHz	
	9	Run time limit on kernels:	Yes	
	10	ECC support enabled:	No	
g	11	Device 1: "GeForce 8400 GS"		
D	12	Compute capability:	1.1	
	13	Multiprocessors:	1 (8 cores)	
	14	Memory global:	512 MB	
	15	Clock rate:	1.63 GHz	
	16	Run time limit on kernels:		
	17	ECC support enabled:	No	
	18			
		[GPU Hardware]		
		Device default: 0 "GeForce	GTX 590"	
		Compute capability: 2.0		
		Memory global: 1536 MB		
		Memory shared: 49152 Bytes		
		[Initialising JSphGpuSingle]	v0.70 24-11-201	5 15:40:361
		**Basic case configuration i		
		**Special case configuration		
		Loading initial state of par		
		Loaded particles: 5281		
			e-006.0.0999929.	-7.07107e-006)-(1.60001,0.100007,0.400007)
	2.2			7.07107e-006) - (1.60001,0.100007,0.600014)
	31	**Initial state of particles		
		**2D-Simulation parameters:	10 104464	
		CaseName="CaseDambreak2D"		
		RunName="CaseDambreak2D"		
		SvDouble=False		
		SvTimers=True		
		SvTimersStep=0.000000		
n		StepAlgorithm="Verlet"		
		VerletSteps=40		
n l		Kernel="Wendland"		
		Viscosity="Artificial"		
es l		Visco=0.300000		
	43	ViscoBoundFactor=0.000000		
	44	DeltaSph="None"		
		Shifting="None"		
		RenCorrection=0.000000		
	47	Splitting=False		
	48	FloatingFormulation="None"		92
	49	FloatingCount=0		
		- Total ago and o		

Domain dimensions computed starting from minimum and maximum positions of the particles created initially

Run.out

Text file with execution log

50	CaseNp=5281	
	CaseNbound=481	
2	CaseNfixed=481	
	CaseNmoving=0	
	CaseNfloat=0	
	CaseNfluid=4800	
	PeriodicActive=0	
	Dx=0.005	
	H=0.007071	
	CoefficientH=1	
	CteB=165368.578125	
	Gamma=7.000000	
	RhopZero=1000.000000	
	Eps=0	
	Cs0=34.0232	
	CFLnumber=0.200000	
	DtIni=0.000207831	
	DtMin=1.03915e-005	
	DtAllParticles=False	
	MassFluid=0.025000	
	MassBound=0.025000	
	Bwen (wendland)=-7877736.000000	
	TimeMax=2	
	TimePart=0.02	
	Gravity=(0.000000,0.000000,-9.810000)	
	NpMinimum=481	
	RhopOut=True	
	RhopOutMin=700.000000	
	RhopOutMax=1300.000000	
	**Requested gpu memory for 5281 particles: 0.6 MB.	
	CellOrder="XYZ"	
	CellMode="2H"	
	Hdiv=1	
	MapCells=(114,1,43)	
	DomCells=(114,1,43)	
	DomCellCode="13_8_11"	
	PtxasFile="//EXECS/DualSPHysics_win64_ptxasinfo"	
	Use code for compute capability 2.0 on hardware 2.0	
	BsForcesBound=128 (36 regs)	
	BsForcesFluid=128 (50 regs)	
	**CellDiv: Requested gpu memory for 5545 particles: 0.0 MB.	
	**CellDiv: Requested gpu memory for 1488 cells (CellMode=2H): 0.0 MB.	
	RunMode="Pos-Simple, Single-Gpu"	
5	Allocated memory in CPU: 475290 (0.45 MB)	
	Allocated memory in GPU: 745752 (0.71 MB)	-
	Part 0000 5281 particles successfully stored	9

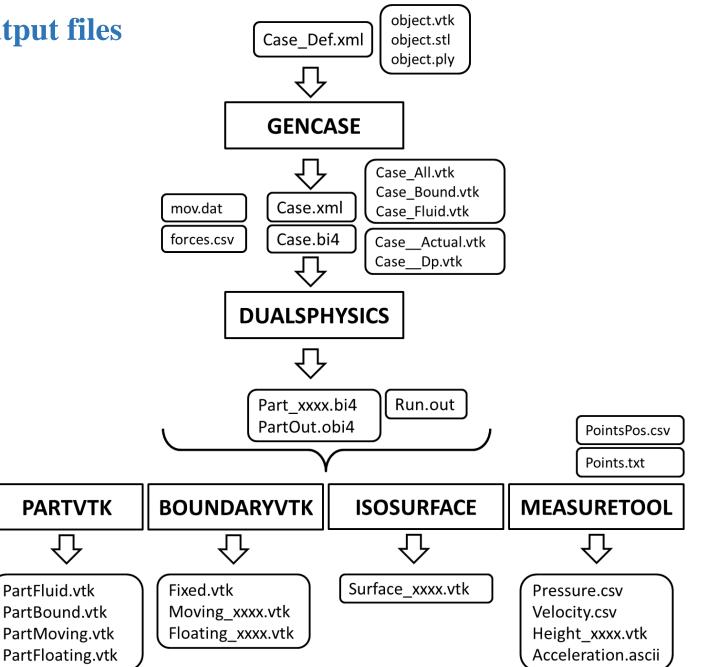
Run.out

Text file with execution log

92	**CellDiv:	Requested gpu	memory for 55	45 partic	les: 0.0 ME	3.	
93	**CellDiv:	Requested gpu	memory for 14	88 cells	(CellMode=2	2H): 0.0 MB.	
94	RunMode="Po	os-Simple, Sin	gle-Gpu"				
95	Allocated r	memory in CPU:	475290 (0.45	MB)			
96	Allocated r	memory in GPU:	745752 (0.71	MB)			
97	Part_0000	5281 pa	rticles succes	sfully st	ored		
98							
99	-	-	(m15001k0) 2		-		
.00	PART ========	PartTime	TotalSteps	Steps	Time/Seg		
.01 .02	Part 0001	0.020027	484	484		24-11-2015 1	5.42.10
.03	Part 0002	0.040010	970			24-11-2015 1	
.04	Part 0003	0.060019	1460	490		24-11-2015 1	
.05	Part 0004	0.080034	1953	493	47.73	24-11-2015 1	5:42:11
.06	Part 0005	0.100001	2447	494	48.86	24-11-2015 1	5:42:11
.07	–						
.08							
.09							
.10	Part_0100	2.000038	49060	484	67.06	24-11-2015 1	5:42:14
.11							
.12			-11-2015 15:42	-			
.13			(initial): 528	1			
.14 .15	-	ed to DtMin articles					
.15				489380 88	C.		
.17							
.18			ation: 49.				
.19	Steps per s	second	498	.439697			
.20	Steps of s:	imulation	490	60			
.21	PART files		101				
.22	Maximum nur	mber of partic	les: 528	1			
.23			285				
.24	-		475	-			
.25	GPU Memory	•••••	777	144 (0.74	MB)		
.26 .27	[GPU Timers	-1					
.27	-	-	0.0	36532 860			
.20							
.30							
.31			33.				
.32			4.9				
.33	NL-SortData	a	0.6	39390 sec			
.34	NL-OutChecl	k	0.1	14177 sec			
.35	CF-PreForce	28	: 3.7	01347 sec			
.36	CF-Forces.		11.	354674 se	с.		
.37	-	-	0.9				
.38		-	0.0				
.39	SU-Motion.			00000 sec			94

Pre-Processing





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