



Pre-processing and User Friendly Interface

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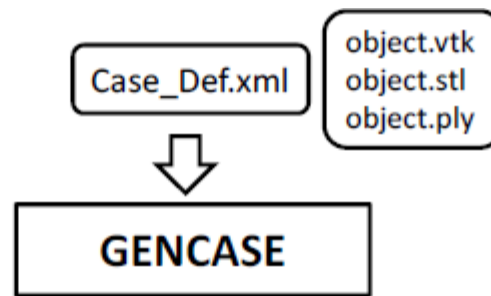
Universidade de Vigo
SPAIN

Outline of Presentation

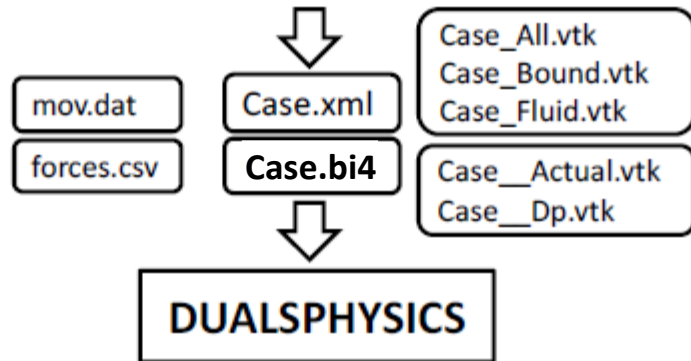
- Pre-processing tool: GenCase
- XML input file
- DesignSPHysics
- Future developments

Pre-processing: GenCase

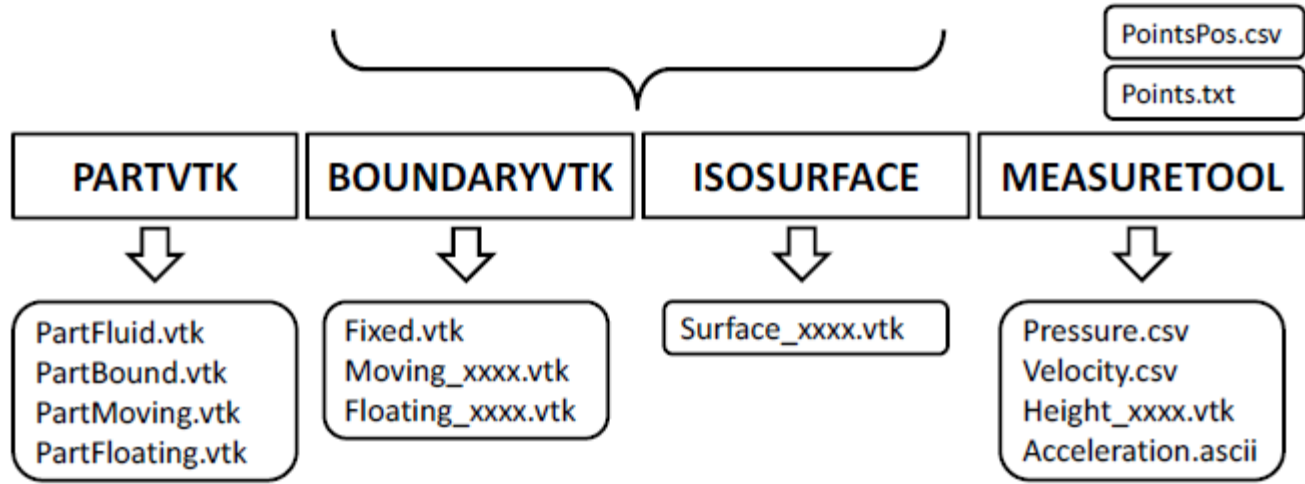
PRE-PROCESSING



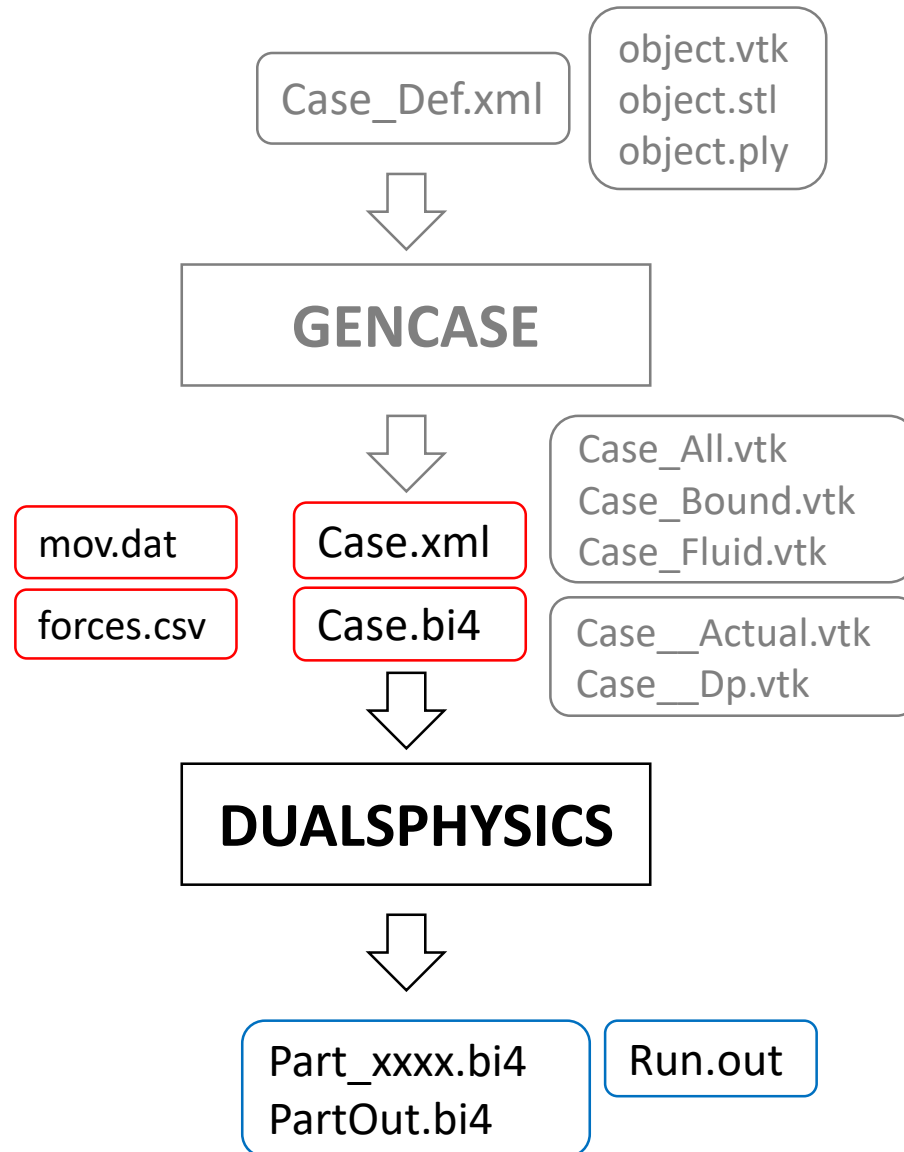
SPH SOLVER



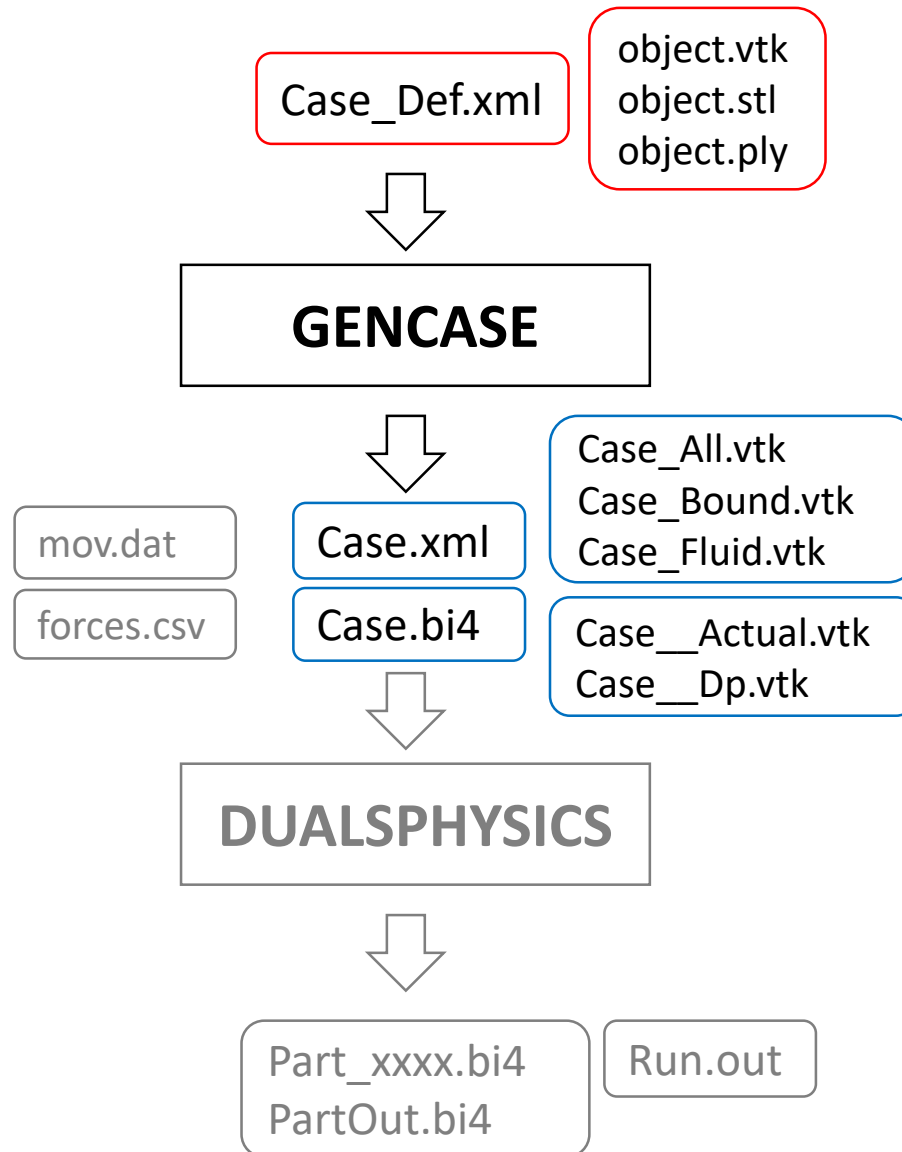
POST-PROCESSING



Pre-processing: GenCase



Pre-processing: GenCase



Pre-processing: GenCase

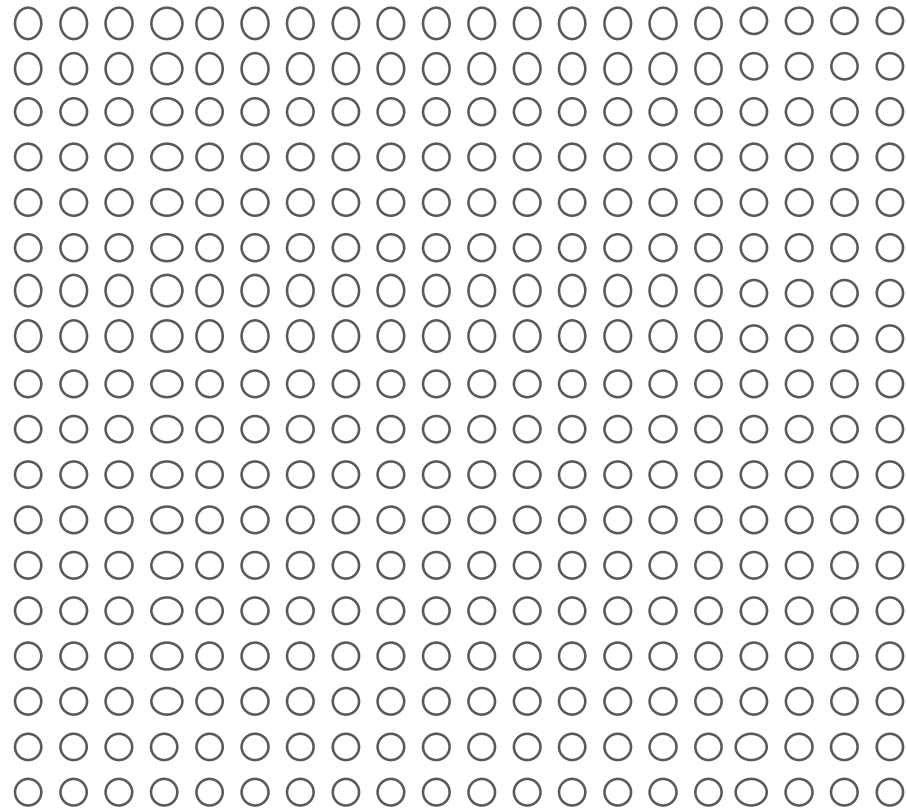
Draw points

- **GenCase** is a drawing application that creates **points** that will be converted **into particles**.
- It employs a **3D lattice to locate points** which represents possible positions of the particles.

Pre-processing: GenCase

Draw points

Lattice



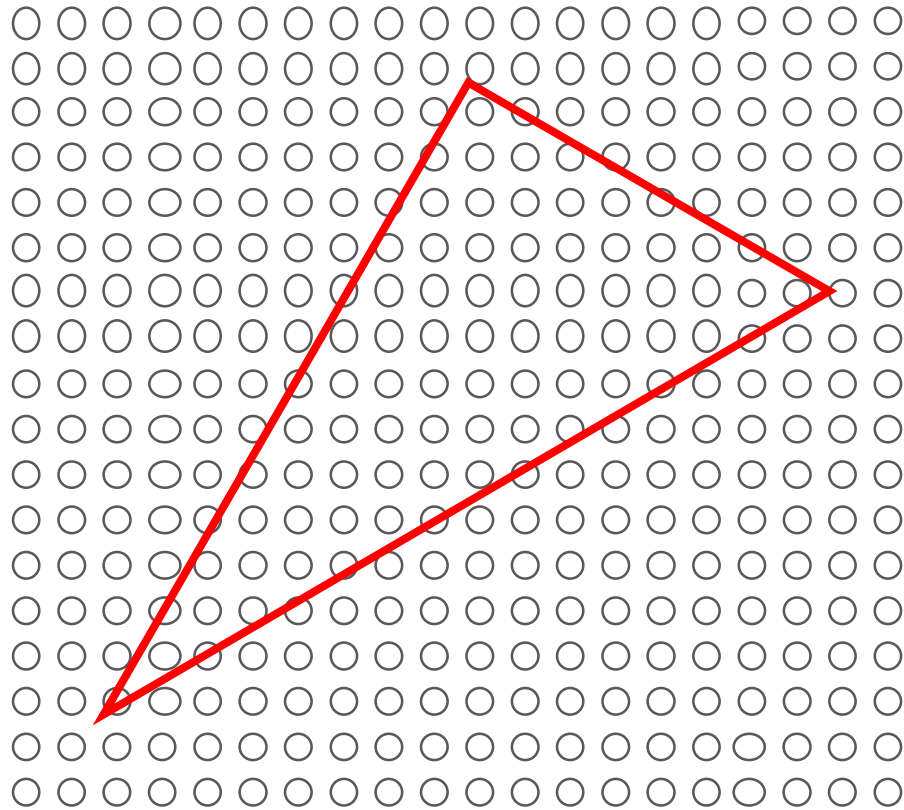
Pre-processing: GenCase

Draw points

Lattice



**Drawing
triangle**



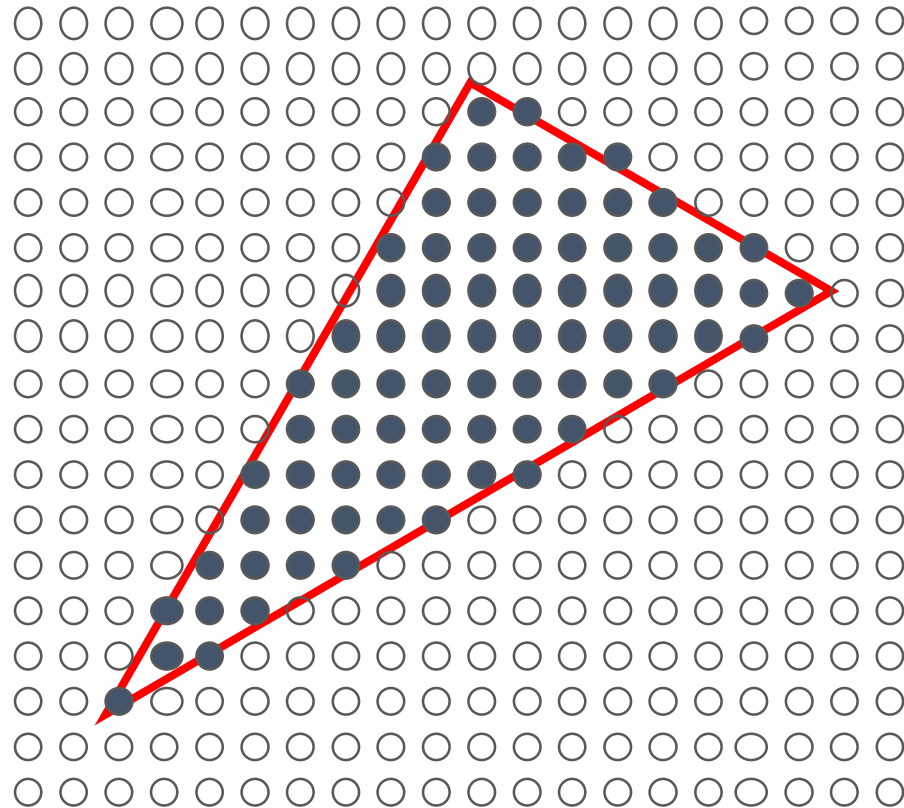
Pre-processing: GenCase

Draw points

Lattice



Drawing
triangle



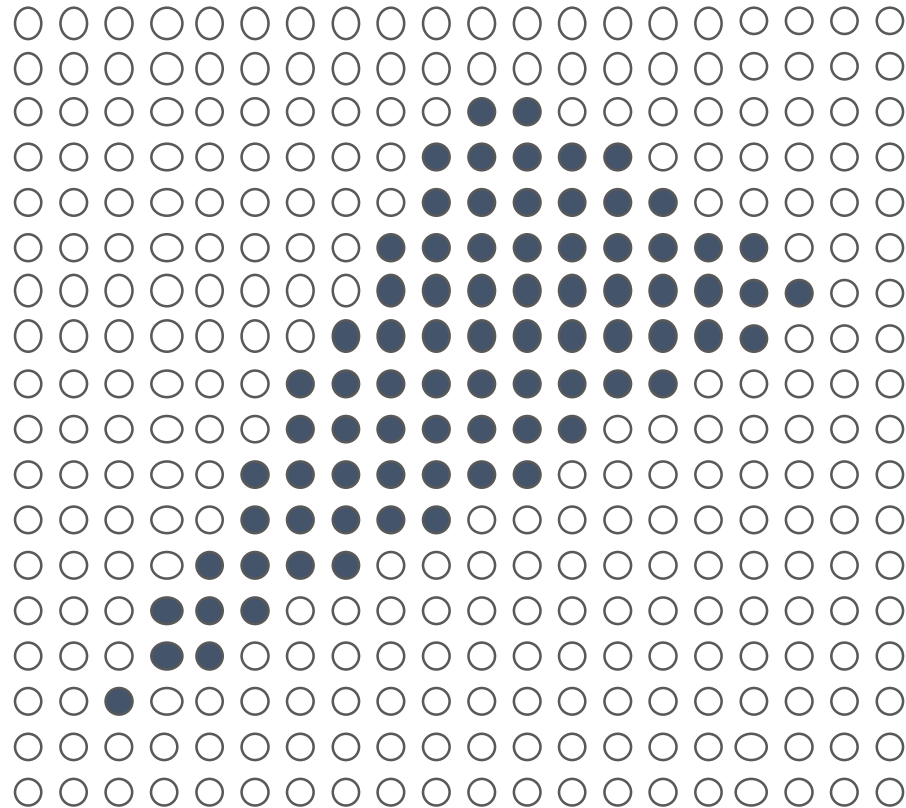
Pre-processing: GenCase

Draw points

Lattice



Triangle to
points



Pre-processing: GenCase

Draw points

Lattice



Triangle to
points

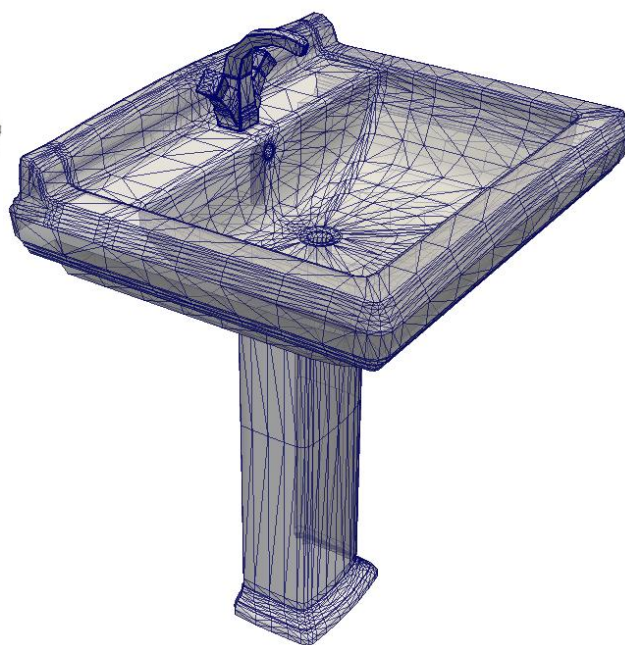


Pre-processing: GenCase

Draw points



3D model



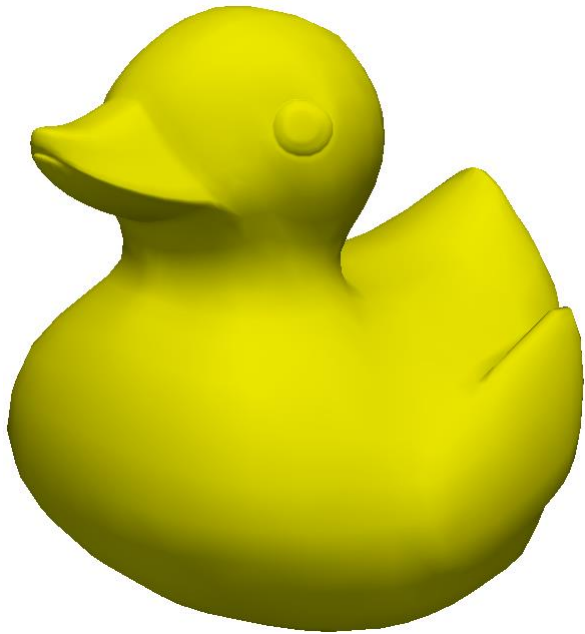
Triangles



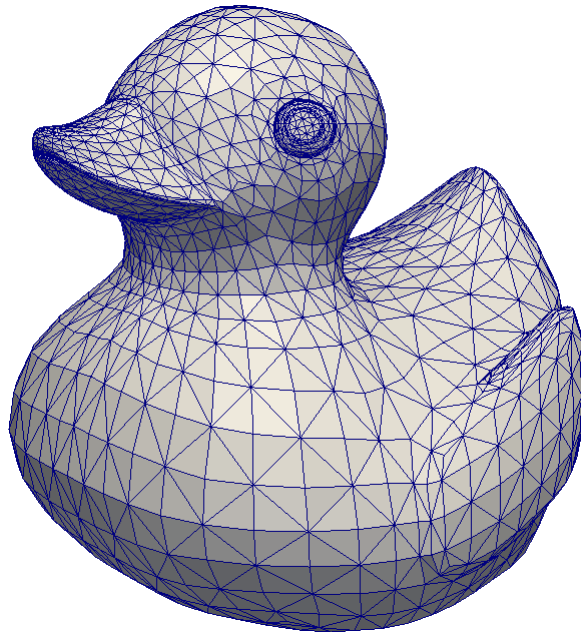
Points

Pre-processing: GenCase

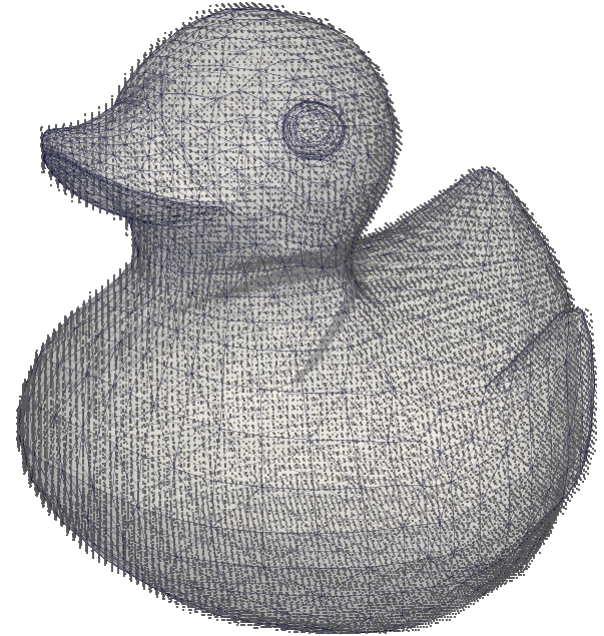
Draw points



3D model



Triangles

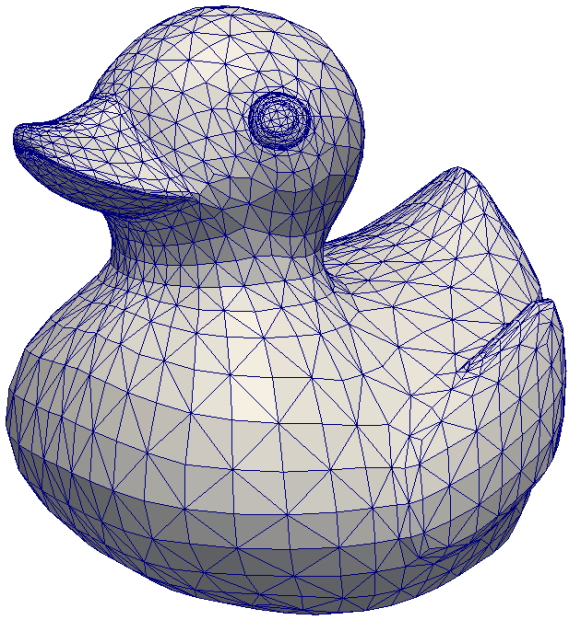


Points

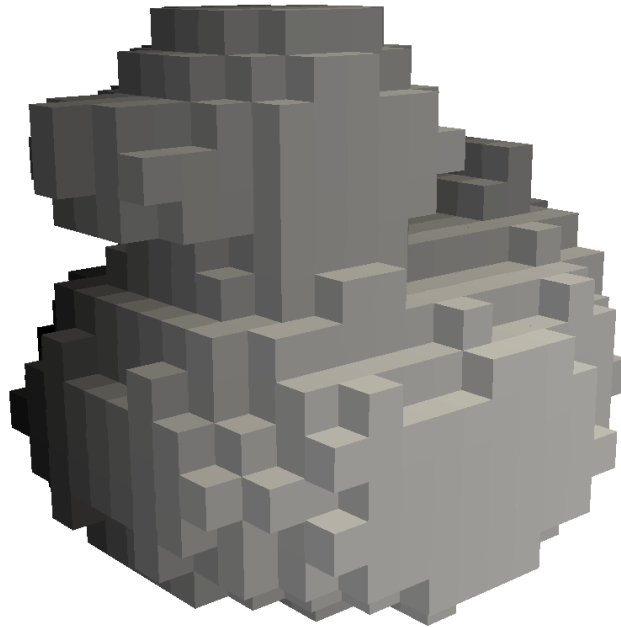
Pre-processing: GenCase

Draw points

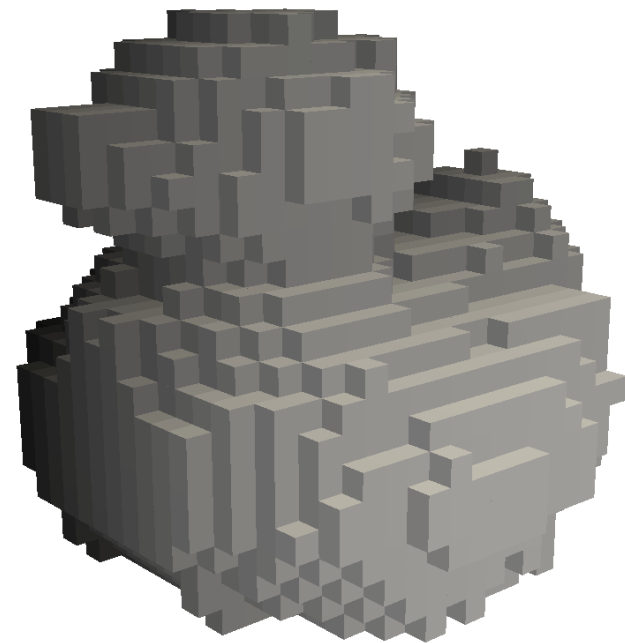
Particles are depicted as cubes



Triangles



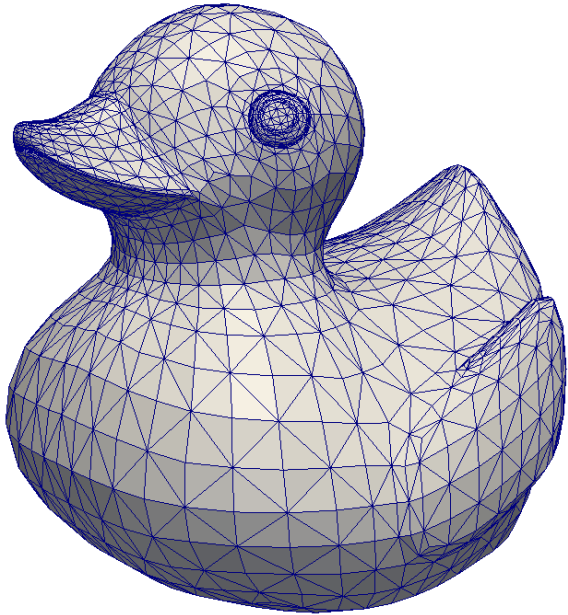
Low resolution
836 particles



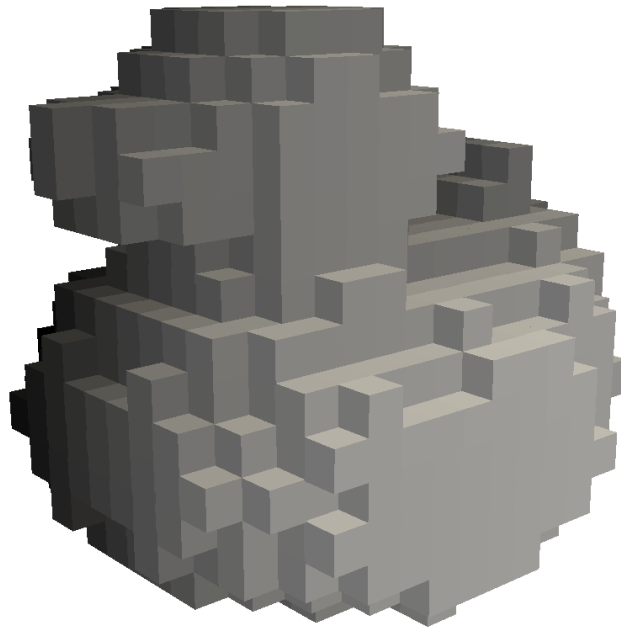
Higher resolution
1,836 particles

Pre-processing: GenCase

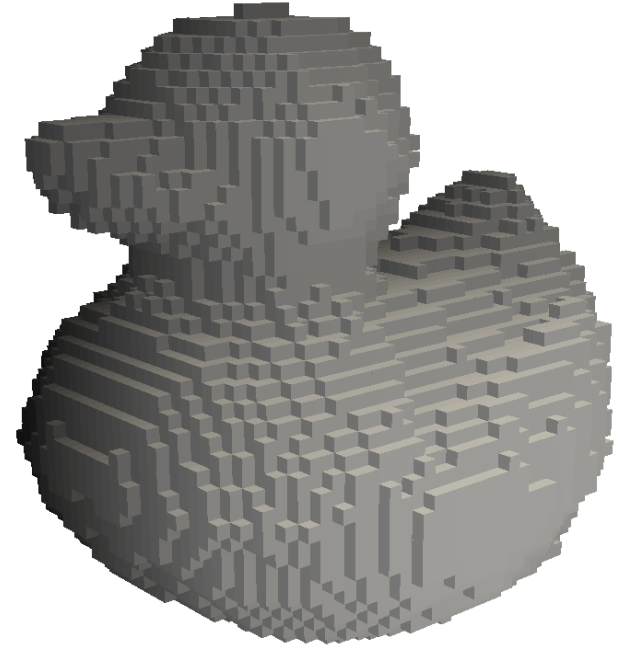
Draw points



Triangles



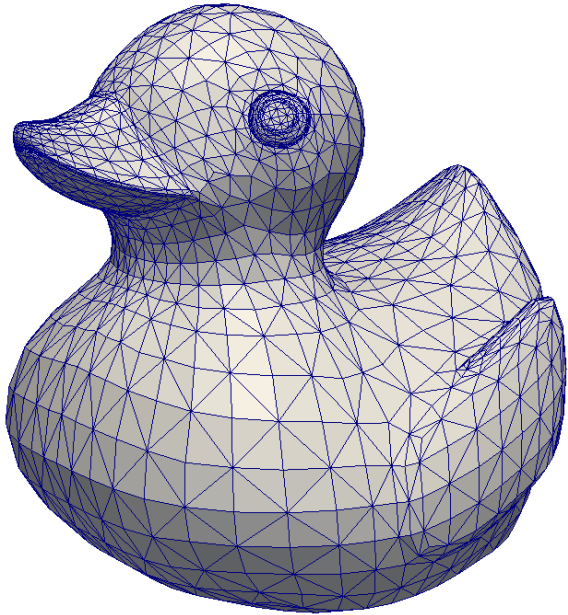
Low resolution
836 particles



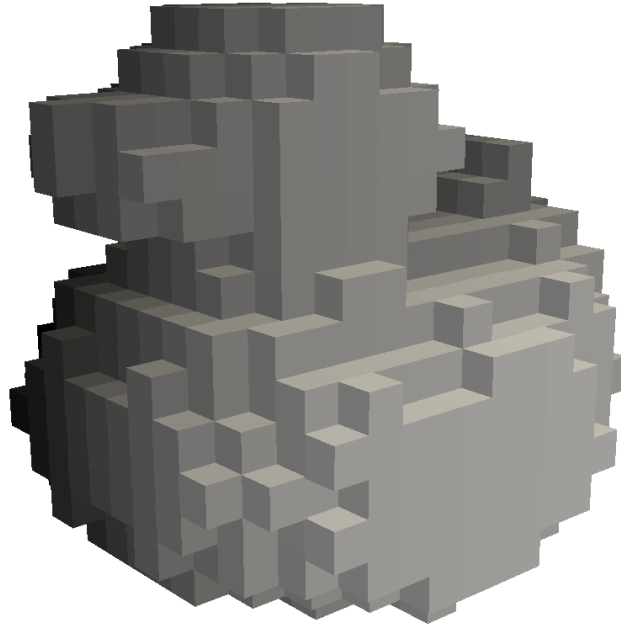
Higher resolution
6,934 particles

Pre-processing: GenCase

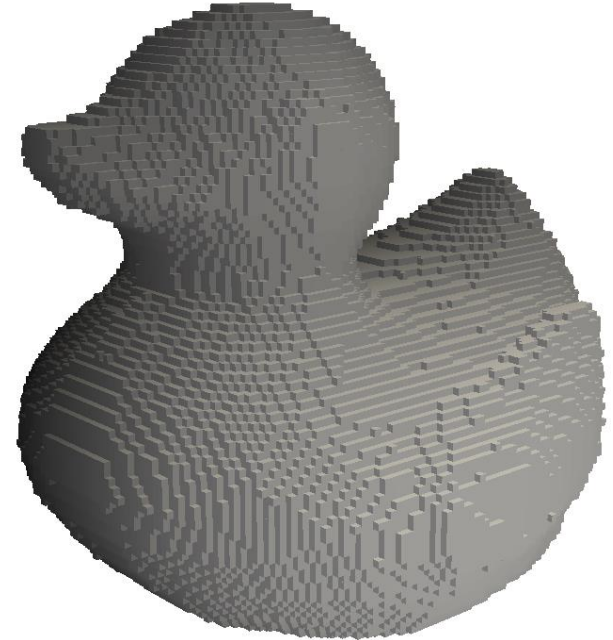
Draw points



Triangles



Low resolution
836 particles

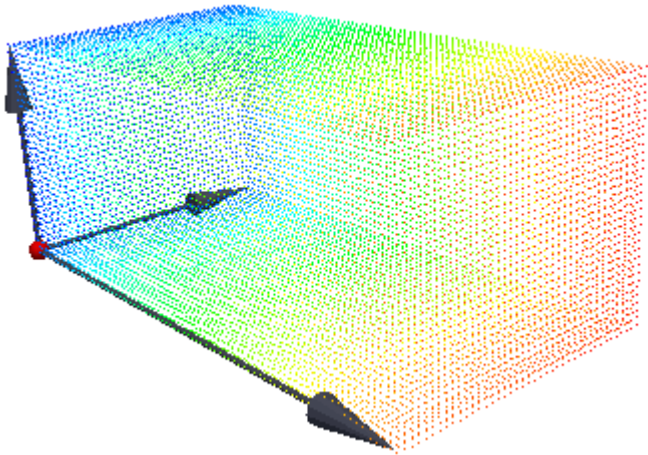


Higher resolution
27,474 particles

Pre-processing: GenCase

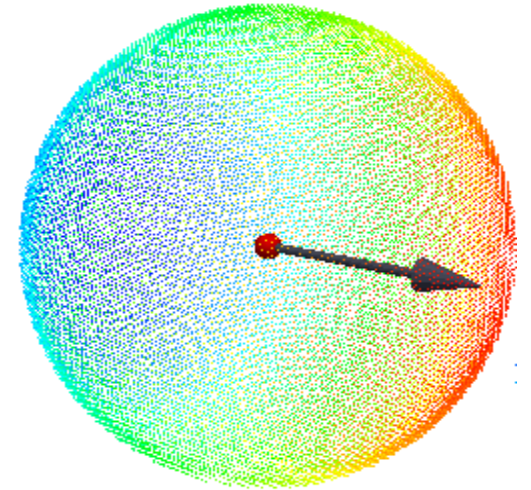
Predefined objects

Predefined shapes can be added to the simulation just by setting up some configuration parameters:



BOX: a **corner** and the **size** are required

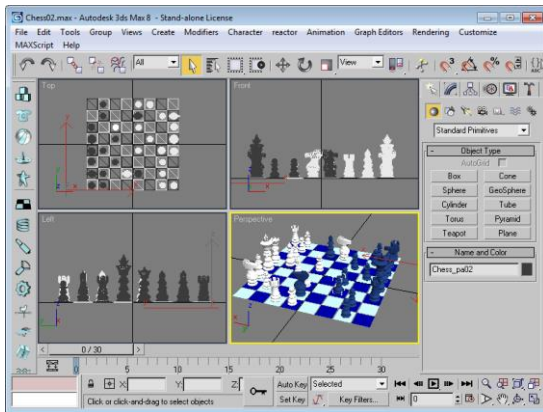
SPHERE: the **centre** and **radius** are needed



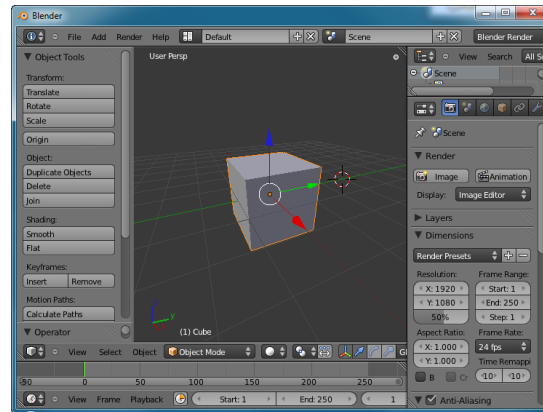
Pre-processing: GenCase

External objects

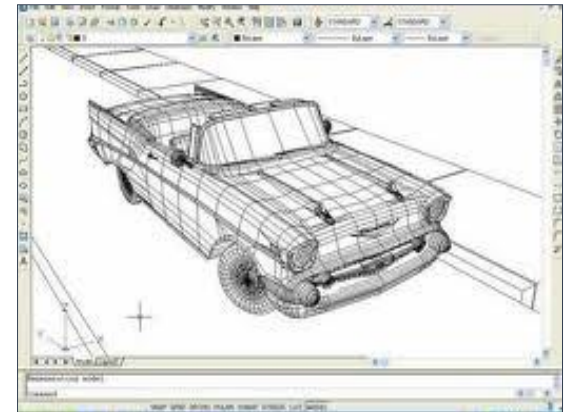
GenCase is able to load external object designed with different softwares.



3D Studio Max



Blender



AutoCAD

3DS DXF DWG GIS H5PART CSV MAX
SHP CAD PLY STL VTK

Pre-processing: GenCase

External objects

3DS DXF DWG GIS H5PART CSV MAX
SHP CAD PLY STL VTK



PLY -> exportable using BLENDER
STL -> exportable using 3DSTUDIO
VTK -> PARAVIEW

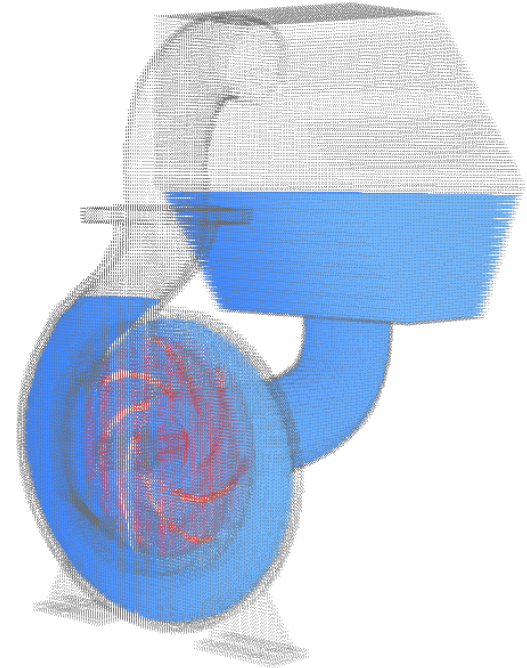
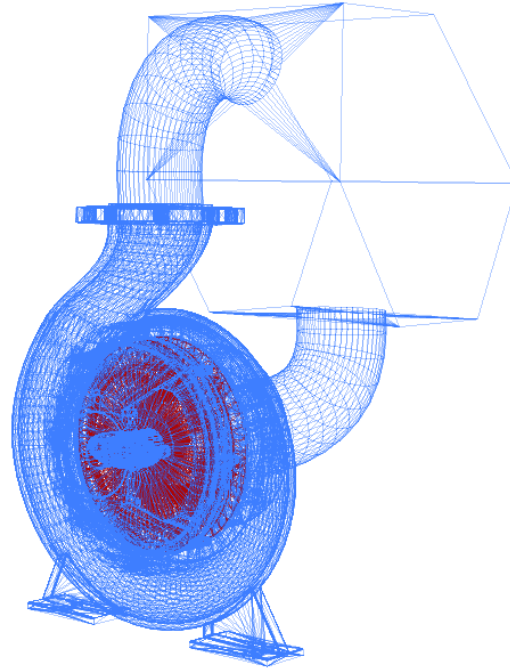
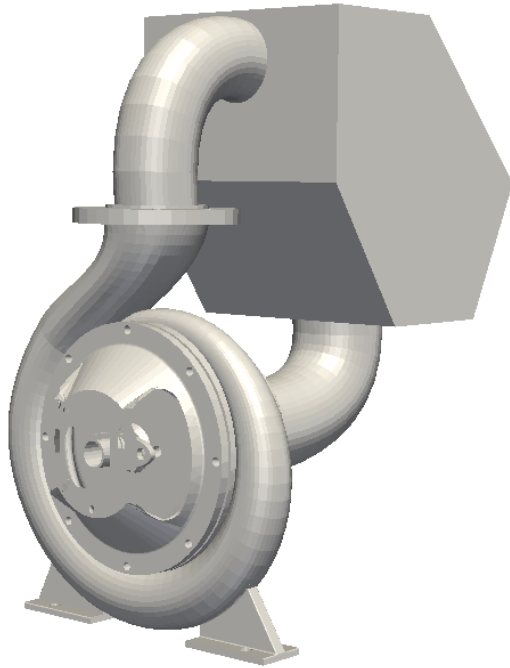


PLY, STL, VTK can be loaded by GenCase

Pre-processing: GenCase

External objects

CAD files

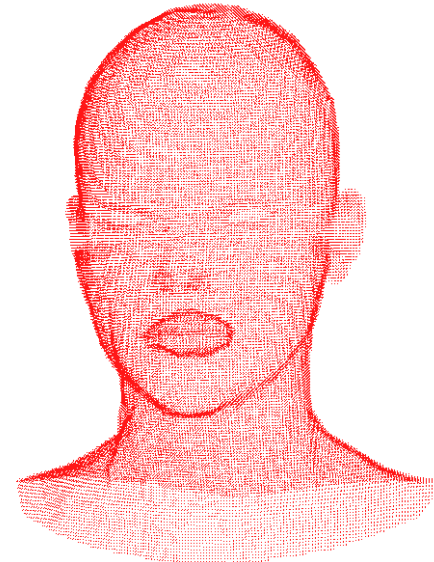
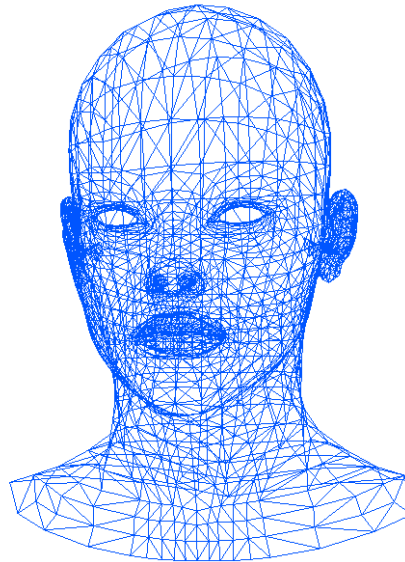


PLY, STL, VTK can be loaded by GenCase

Pre-processing: GenCase

External objects

3D Studio objects

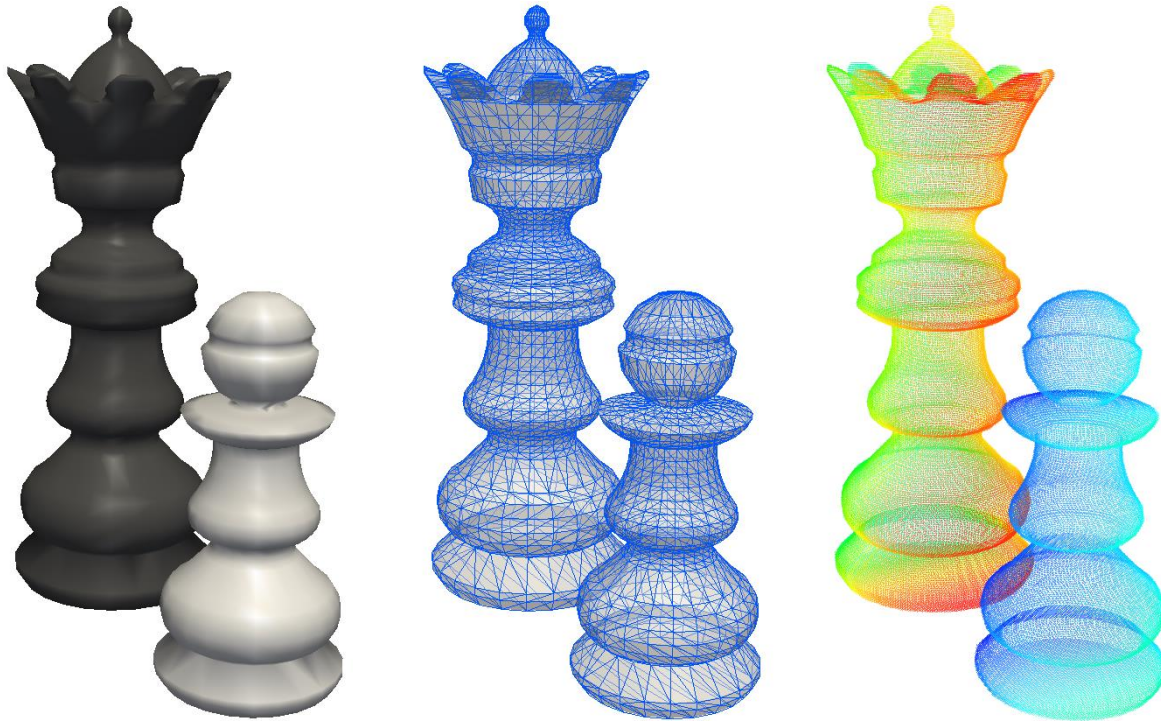


PLY, STL, VTK can be loaded by **GenCase**

Pre-processing: GenCase

External objects

3D Studio objects



PLY, STL, VTK can be loaded by GenCase

Outline of Presentation

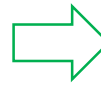
- Pre-processing tool: GenCase
- XML input file
- DesignSPHysics
- Future developments

XML file

Case_Def.xml



GENCASE



Case.xml

```
<case>
  <casedef>
    <constantsdef>
      <lattice bound="1" fluid="1" />
      <gravity x="0" y="0" z="-9.81" comment="Gravitational acceleration" units_comment="m/s^2" />
      <cofnumber value="0.2" comment="Coefficient to multiply Dt" />
      <chwl value="0" auto="true" comment="Maximum still water level to calculate speedofsound using coefsound" units_comment="metres (m)" />
      <speedsystem value="0" auto="true" comment="Maximum system speed (by default the dam-break propagation is used)" />
      <coefsound value="0" comment="Coefficient to multiply speedsystem" />
      <speedsound value="0" auto="true" comment="Speed of sound to use in the simulation (by default speedofsound=coefsound*speedsystem) />
      <coefh value="1.0" comment="Coefficient to calculate the smoothing length (B=coefficient*sqrt(3*dp/2) in 3D) />
      <gamma value="7" comment="Polytropic constant for water used in the state equation" />
      <rhop0 value="1000" comment="Reference density of the fluid" units_comment="kg/m3" />
    </constantsdef>
    <mkconfig boundcount="240" fluidcount="10" />
  </casedef>
  <geometry>
    <definition dp="0.01" units_comment="metres (m)">
      <pointmin x="-1" y="0" z="-1" />
      <pointmax x="4.5" y="0" z="3.5" />
    </definition>
    <commands>
      <mainlist>
        <setdrawmode mode="full" />
        <setmkfluid mk="0" />
        <drawbox>
          <boxfill>solid</boxfill>
          <point x="0" y="-1" z="0" />
          <size x="1" y="2" z="2" />
        </drawbox>
        <setmkbound mk="0" />
        <drawbox>
          <boxfill>bottom | left | right | front | back</boxfill>
          <point x="0" y="-1" z="0" />
          <size x="4" y="2" z="3" />
        </drawbox>
      </mainlist>
    </commands>
  </geometry>
</casedef>
  <execution>
    <parameters>
      <parameter key="StepAlgorithm" value="1" comment="Step Algorithm 1:Verlet, 2:Symplectic (default=1)" />
      <parameter key="VerletSteps" value="40" comment="Verlet only: Number of steps to apply Euler timestepping (default=40)" />
      <parameter key="Kernel" value="2" comment="Interaction Kernel 1:Cubic Spline, 2:Mesland (default=2)" />
      <parameter key="ViscoTreatment" value="1" comment="Viscosity formulation 1:Artificial, 2:Laminar+SPS (default=1)" />
      <parameter key="Visco" value="0.02" comment="Viscosity value" />
      <parameter key="ViscoBoundFactor" value="1" comment="Multiply viscosity value with boundary (default=1)" />
      <parameter key="DeltaSPH" value="0" comment="DeltaSPH value, 0.1 is the typical value, with 0 disabled (default=0)" />
      <parameter key="Shifting" value="0" comment="Shifting mode 0:None, 1:Ignore bound, 2:Ignore fixed, 3:Full (default=0)" />
      <parameter key="ShiftingCoef" value="-2" comment="Coefficient for shifting computation (default=-2)" />
      <parameter key="ShiftingTFS" value="1.5" comment="Threshold to detect free surface. Typically 1.5 for 2D and 2.75 for 3D (default=0)" />
      <parameter key="RigidAlgorithm" value="1" comment="Rigid Algorithms 1:SPH, 2:DBM (default=1)" />
      <parameter key="PFreeze" value="0.0" comment="Time to freeze the floatings at simulation start (warmup) (default=0) units_comment="seconds" />
      <parameter key="CoefDMin" value="0.05" comment="Coefficient to calculate minimum time step dtmin=coefdtmin*h/speedsound (default=0.05)" />
      <parameter key="DtIn" value="0.0001" comment="Initial time step (default=h/speedsound) units_comment="seconds" />
      <parameter key="DtMin" value="0.00001" comment="Minimum time step (default=coefdtmin*h/speedsound) units_comment="seconds" />
      <parameter key="DtFixed" value="DtFixed.dat" comment="Dt values are loaded from file (default=disabled)" />
      <parameter key="DtAllParticles" value="0" comment="Velocity of particles used to calculate DT. 1:All, 0:Only fluid/floating (default=0)" />
      <parameter key="TimeMax" value="0.72" comment="Time of simulation" units_comment="seconds" />
      <parameter key="TimeOut" value="0.01" comment="Time out data" units_comment="seconds" />
      <parameter key="InoZ" value="1" comment="Increase of I4" units_comment="decimal" />
      <parameter key="PartsOutMax" value="1" comment="Allowed k/100 of fluid particles out the domain (default=1) units_comment="decimal" />
      <parameter key="RhopOutMin" value="700" comment="Minimum rhop valid (default=700) units_comment="kg/m3" />
      <parameter key="RhopOutMax" value="1300" comment="Maximum rhop valid (default=1300) units_comment="kg/m3" />
    </parameters>
  </execution>
</case>
```

```
<case>
  <casedef>
    <constantsdef>
      <lattice bound="1" fluid="1" />
      <gravity x="0" y="0" z="-9.81" comment="Gravitational acceleration" units_comment="m/s^2" />
      <cofnumber value="0.2" comment="Coefficient to multiply Dt" />
      <chwl value="0" auto="true" comment="Maximum still water level to calculate speedofsound using coefsound" units_comment="metres (m)" />
      <speedsystem value="0" auto="true" comment="Maximum system speed (by default the dam-break propagation is used)" />
      <coefsound value="0" comment="Coefficient to multiply speedsystem" />
      <speedsound value="0" auto="true" comment="Speed of sound to use in the simulation (by default speedofsound=coefsound*speedsystem) />
      <coefh value="1.0" comment="Coefficient to calculate the smoothing length (B=coefficient*sqrt(3*dp/2) in 3D) />
      <gamma value="7" comment="Polytropic constant for water used in the state equation" />
      <rhop0 value="1000" comment="Reference density of the fluid" units_comment="kg/m3" />
    </constantsdef>
    <mkconfig boundcount="240" fluidcount="10" />
  </casedef>
  <geometry>
    <definition dp="0.01" units_comment="metres (m)">
      <pointmin x="-1" y="0" z="-1" />
      <pointmax x="4.5" y="0" z="3.5" />
    </definition>
    <commands>
      <mainlist>
        <setdrawmode mode="full" />
        <setmkfluid mk="0" />
        <drawbox>
          <boxfill>solid</boxfill>
          <point x="0" y="-1" z="0" />
          <size x="1" y="2" z="2" />
        </drawbox>
        <setmkbound mk="0" />
        <drawbox>
          <boxfill>bottom | left | right | front | back</boxfill>
          <point x="0" y="-1" z="0" />
          <size x="4" y="2" z="3" />
        </drawbox>
      </mainlist>
    </commands>
  </geometry>
</casedef>
  <execution>
    <parameters>
      <parameter key="StepAlgorithm" value="1" comment="Step Algorithm 1:Verlet, 2:Symplectic (default=1)" />
      <parameter key="VerletSteps" value="40" comment="Verlet only: Number of steps to apply Euler timestepping (default=40)" />
      <parameter key="Kernel" value="2" comment="Interaction Kernel 1:Cubic Spline, 2:Mesland (default=2)" />
      <parameter key="ViscoTreatment" value="1" comment="Viscosity formulation 1:Artificial, 2:Laminar+SPS (default=1)" />
      <parameter key="Visco" value="0.02" comment="Viscosity value" />
      <parameter key="ViscoBoundFactor" value="1" comment="Multiply viscosity value with boundary (default=1)" />
      <parameter key="DeltaSPH" value="0" comment="DeltaSPH value, 0.1 is the typical value, with 0 disabled (default=0)" />
      <parameter key="Shifting" value="0" comment="Shifting mode 0:None, 1:Ignore bound, 2:Ignore fixed, 3:Full (default=0)" />
      <parameter key="ShiftingCoef" value="-2" comment="Coefficient for shifting computation (default=-2)" />
      <parameter key="ShiftingTFS" value="1.5" comment="Threshold to detect free surface. Typically 1.5 for 2D and 2.75 for 3D (default=0)" />
      <parameter key="RigidAlgorithm" value="1" comment="Rigid Algorithms 1:SPH, 2:DBM (default=1)" />
      <parameter key="PFreeze" value="0.0" comment="Time to freeze the floatings at simulation start (warmup) (default=0) units_comment="seconds" />
      <parameter key="CoefDMin" value="0.05" comment="Coefficient to calculate minimum time step dtmin=coefdtmin*h/speedsound (default=0.05)" />
      <parameter key="DtIn" value="0.0001" comment="Initial time step (default=h/speedsound) units_comment="seconds" />
      <parameter key="DtMin" value="0.00001" comment="Minimum time step (default=coefdtmin*h/speedsound) units_comment="seconds" />
      <parameter key="DtFixed" value="DtFixed.dat" comment="Dt values are loaded from file (default=disabled)" />
      <parameter key="DtAllParticles" value="0" comment="Velocity of particles used to calculate DT. 1:All, 0:Only fluid/floating (default=0)" />
      <parameter key="TimeMax" value="0.72" comment="Time of simulation" units_comment="seconds" />
      <parameter key="TimeOut" value="0.01" comment="Time out data" units_comment="seconds" />
      <parameter key="InoZ" value="1" comment="Increase of I4" units_comment="decimal" />
      <parameter key="PartsOutMax" value="1" comment="Allowed k/100 of fluid particles out the domain (default=1) units_comment="decimal" />
      <parameter key="RhopOutMin" value="700" comment="Minimum rhop valid (default=700) units_comment="kg/m3" />
      <parameter key="RhopOutMax" value="1300" comment="Maximum rhop valid (default=1300) units_comment="kg/m3" />
    </parameters>
    <particles>
      <particles np="21001" nb="1001" nbf="1001" mkboundfirst="11" mkfluidfirst="1">
        <fixed mkbound="0" mk="11" begin="0" count="1001" />
        <fluid mkfluid="0" mk="1" begin="1001" count="20000" />
      </particles>
    </particles>
    <constants>
      <gravity x="0" y="0" z="-9.81" units_comment="m/s^2" />
      <cofnumber value="0.2" />
      <gamma value="7" />
      <rhop0 value="1000" units_comment="kg/m3" />
      <dp value="0.01" units_comment="metres (m)" />
      <b value="1.41421356237-00" units_comment="metres (m)" />
      <cb value="1.11553714298+006" units_comment="metres (m)" />
      <massfluid value="1.0000000000E-001" units_comment="kg" />
      <massfluid value="1.0000000000E-001" units_comment="kg" />
    </constants>
    <motion />
  </execution>
</case>
```


STRUCTURE OF THE XML FILE

<http://dual.sphysics.org/index.php/downloads/>

- DUALSPHYSICS DOCUMENTATION:

- DualSPHysics_v4.0_GUIDE.pdf

- **XML_GUIDE_v4.0.pdf**

- ExternalModelsConversion_GUIDE.pdf

- PostprocessingCalculations.pdf

- DUALSPHYSICS PACKAGE:

- DualSPHysics_v4.0_Linux_x64.zip

- DualSPHysics_v4.0_Windows_x64.zip

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

```
- <case>
  - <casedef>
    + <constantsdef>
    <mkconfig/>
    - <geometry>
      + <definition>
      - <commands>
        + <mainlist>
        </commands>
      </geometry>
    + <initials>
    + <floatings>
    + <motion>
  </casedef>
  <execution>
    - <special>
      - <wavepaddles>
        + <piston>
        + <piston_spectrum>
      </wavepaddles>
    + <accinputs>
  </special>
  + <parameters>
</execution>
</case>
```

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

XML file

CASEDEF-CONSTANTSDEF

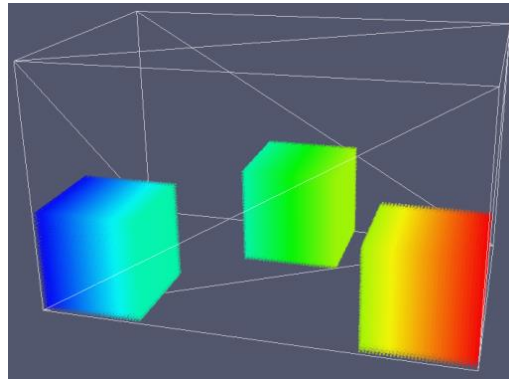
```
<constantsdef>
  <lattice bound="1" fluid="1" />
  <gravity x="0" y="0" z="-9.81" comment="Gravitational acceleration"
    units_comment="m/s^2" />
  <rhop0 value="1000" comment="Reference density of the fluid"
    units_comment="kg/m^3" />
  <hswl value="0" auto="true" comment="Maximum still water level to calculate speedofsound"
    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
    (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
    (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>
```

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

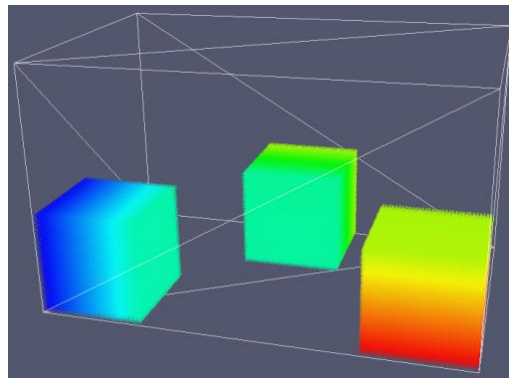
mk: label used to

- 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

mkorientfluid = "xyz"



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



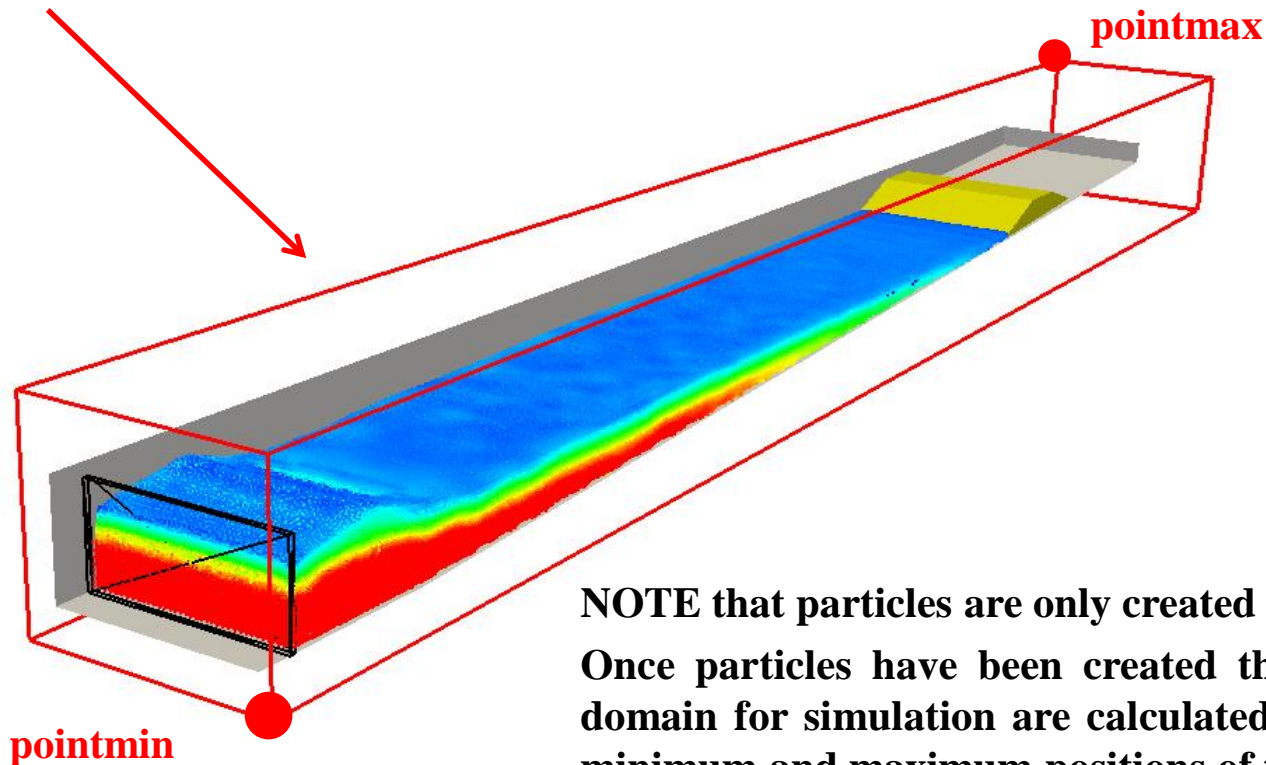
mkorientation: determines the order of particles when creating one object (useful for visualization with the variable *idp*)

```
<!--DEFINITION OF DOMAIN WHERE PARTICLES WILL BE CREATED -->  
<definition dp="0.005">  
  <pointmin x="-0.05" y="0.1" z="-0.05" />  
  <pointmax x=" 2.00" y="0.1" z=" 1.00" />  
</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



NOTE that particles are only created within this domain.
Once particles have been created the dimensions of the domain for simulation are calculated again starting from minimum and maximum positions of the created particles.

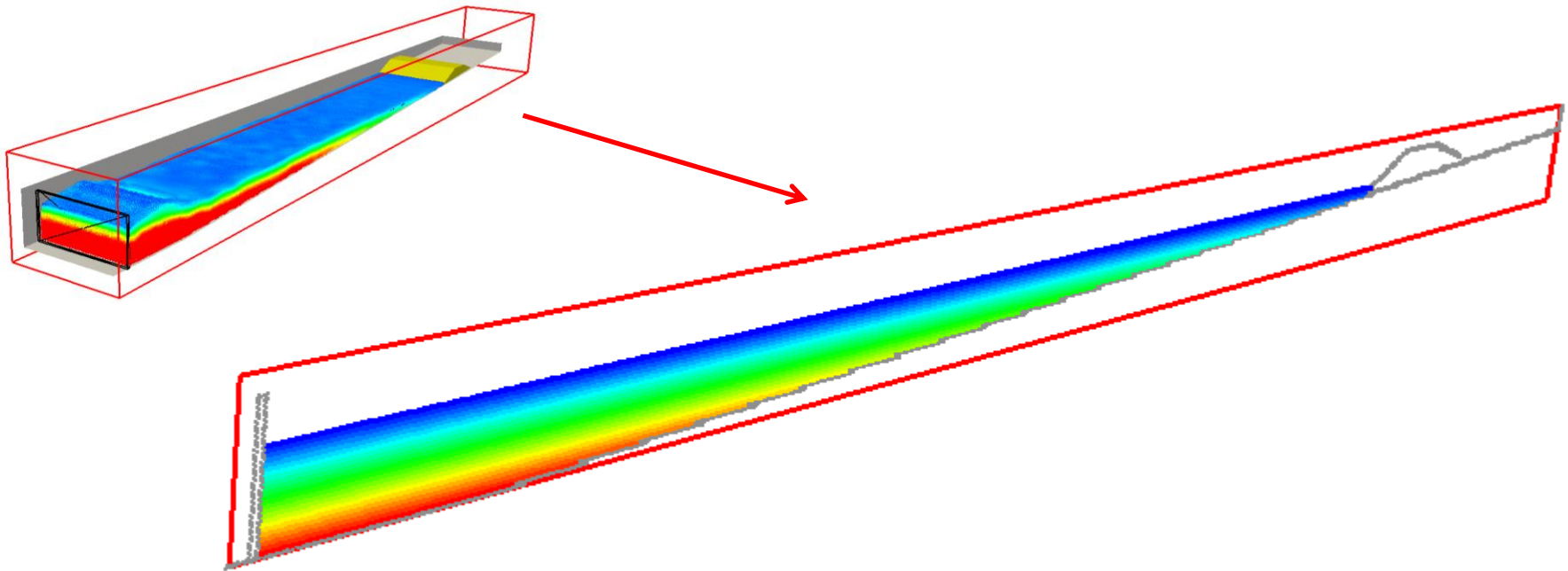
XML file

CASEDEF-GEOMETRY-DEFINITION

```
<!--DEFINITION OF DOMAIN WHERE PARTICLES WILL BE CREATED -->  
<definition dp="0.005">  
  <pointmin x="-0.05" y="0.1" z="-0.05" />  
  <pointmax x=" 2.00" y="0.1" z=" 1.00" />  
</definition>
```

A 2-D configuration can be generated by imposing the same values along Y-direction

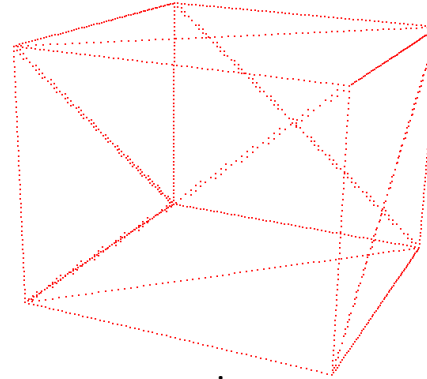
<pointmin> = <pointmax>



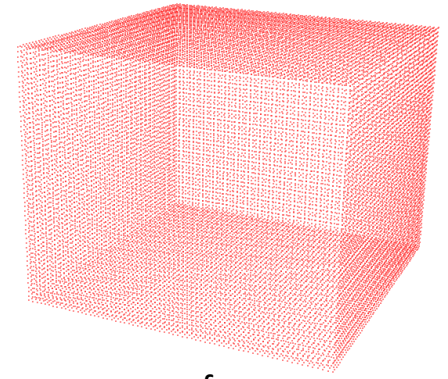
CASEDEF-GEOMETRY-COMMANDS-MAINLIST

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>
```



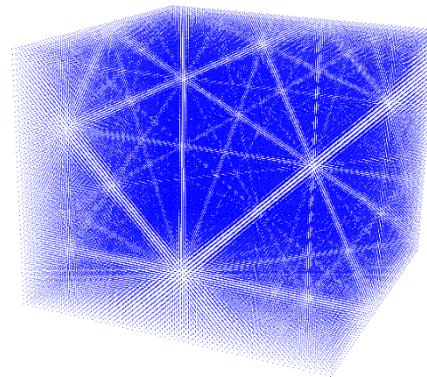
wire



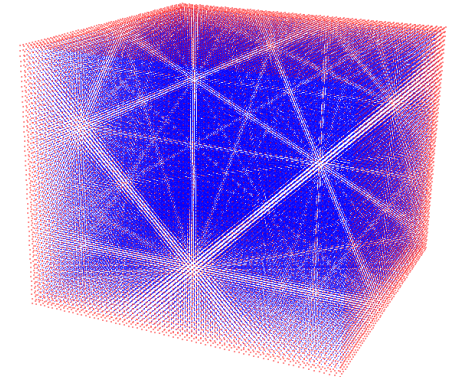
face

<setdrawmode>:

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



solid



full

CASEDEF-GEOMETRY-COMMANDS-MAINLIST

```
- <mainlist>
  <setmkbound mk="0"/>
  <drawfilestl file="File.stl"/>
  <drawfileply file="File.ply"/>
  <drawfilevtk file="File.vtk"/>
- <drawfilestl file="File.stl">
  <drawmove x="0.5" y="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>
```

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

CASEDEF-GEOMETRY-COMMANDS-MAINLIST

IMPORTING EXTERNAL GEOMETRIES

- `<mainlist>`

```
<setshapemode>real | bound | dp</setshapemode>
```

```
<setnkbound mk="0"/>
```

```
<drawfilevtk file="pump_fixed.vtk"/>
```

```
<setnkbound mk="1"/>
```

```
<drawfilevtk file="pump_moving.vtk"/>
```

```
<setnkfluid mk="0"/>
```

```
- <fillbox x="0.14" y="-0.1" z="-0.39">
```

```
  <modefill>void</modefill>
```

```
  <point x="-0.6" y="-0.39" z="-0.8"/>
```

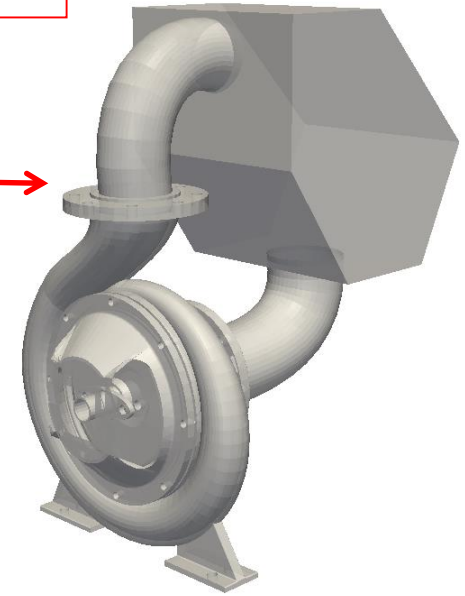
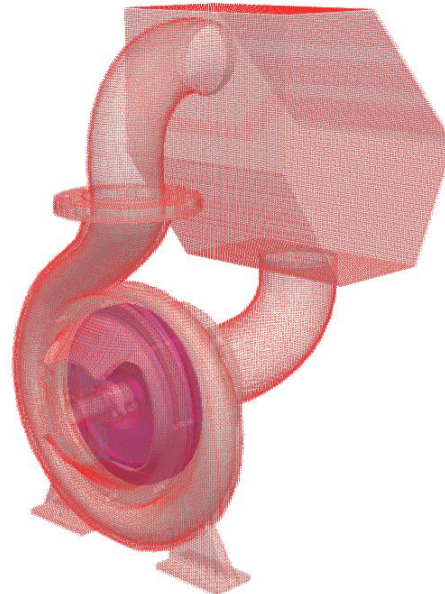
```
  <size x="0.9" y="0.68" z="0.52"/>
```

```
</fillbox>
```

```
</mainlist>
```



from VTK to points



CASEDEF-GEOMETRY-COMMANDS-MAINLIST

```
- <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" y="1" z="0">
  <modefill>border</modefill>
  <point x="0.1" y="1" z="1.1"/>
  <size x="3" y="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-GEOMETRY-COMMANDS-MAINLIST

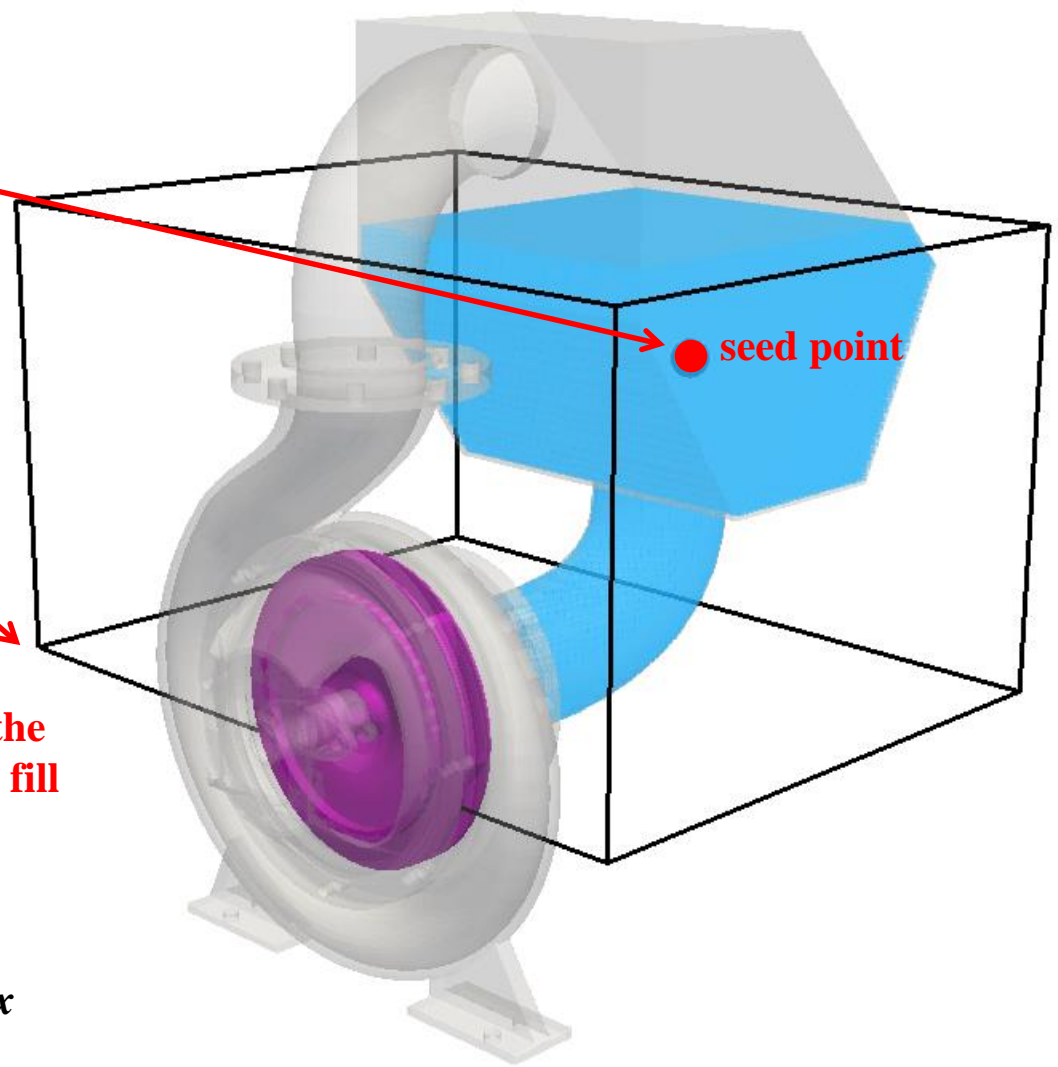
```
- <mainlist>  
  <setshapemode>real | bound | dp</setshapemode>  
  <setnkbound mk="0"/>  
  <drawfilevtk file="pump_fixed.vtk"/>  
  <setnkbound mk="1"/>  
  <drawfilevtk file="pump_moving.vtk"/>  
  <setnklfluid mk="0"/>  
- <fillbox x="0.14" y="-0.1" z="-0.39">  
  <modefill>void</modefill>  
  <point x="-0.6" y="-0.39" z="-0.8"/>  
  <size x="0.9" y="0.68" z="0.52"/>  
</fillbox>  
</mainlist>
```

FILLING DOMAINS

limits of the
domain to fill

seed point

filling with *fluid* while *void* and
before the limits defined by a *box*



```

- <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 * dp³ **or**

masspart = rhopbody * dp³

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

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

```

- <motion>
- <objreal ref="1">
  <begin mov="1" start="0" finish="5.4"/>
  - <mvrect id="1" duration="0.6" next="2">
    <vel x="1" y="0" z="0"/>
  </mvrect>
  <wait id="2" duration="0.3" next="3"/>
  - <mvrect id="3" duration="0.6" next="4">
    <vel x="1" y="0" z="0"/>
  </mvrect>
  <wait id="4" duration="0.3" next="5"/>
  - <mvrect id="5" duration="0.6" next="6">
    <vel x="1" y="0" z="0"/>
  </mvrect>
  <wait id="6" duration="0.3" next="7"/>
  - <mvrect id="7" duration="-1" next="1">
    <vel x="-1.8" y="0" z="0"/>
  </mvrect>
</objreal>
</motion>

```

movement defined for the set of particles with $mk=1$

first mov=1 during 0.6s,
then wait=2 for 0.3s,
then mov=3 during 0.6s,
then wait=4 for 0.3s,
then mov=5 during 0.6s...

<mvrect>: uniform rectilinear movement

vel indicates the constant velocity vector

XML file

CASEDEF-MOTION

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

```
- <motion>
- <objreal ref="4">
  <begin mov="1" start="0"/>
- <mvrotsinu id="1" duration="5" next="2">
  <axisp1 x="0" y="0" z="2.85"/>
  <axisp2 x="0" y="1" z="2.85"/>
  <freq v="0.2"/>
  <ampl v="60"/>
  <_phase v="0"/>
</mvrotsinu>
- <mvrotsinu id="2" duration="5" next="1">
  <axisp1 x="0" y="0" z="2.85"/>
  <axisp2 x="0" y="1" z="2.85"/>
  <freq v="0.4"/>
  <ampl v="75"/>
</mvrotsinu>
</objreal>
- <objreal ref="5">
  <begin mov="1" start="0"/>
- <mvcirsinu id="1" duration="5" next="2">
  <ref x="0" y="-0.7" z="0.2"/>
  <axisp1 x="0" y="0" z="2.85"/>
  <axisp2 x="0" y="1" z="2.85"/>
  <freq v="0.2"/>
  <ampl v="60"/>
  <phase v="0"/>
</mvcirsinu>
- <mvrectsinu id="2" duration="5" next="1">
  <freq x="0.4" y="0" z="0"/>
  <ampl x="2.55" y="0" z="0"/>
  <phase x="0" y="0" z="0"/>
</mvrectsinu>
</objreal>
</motion>
```

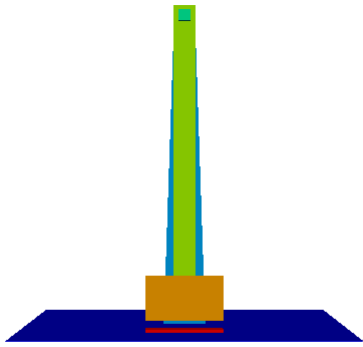
<mvrectsinu>: sinusoidal rectilinear movement

<mvrotsinu>: sinusoidal rotational movement

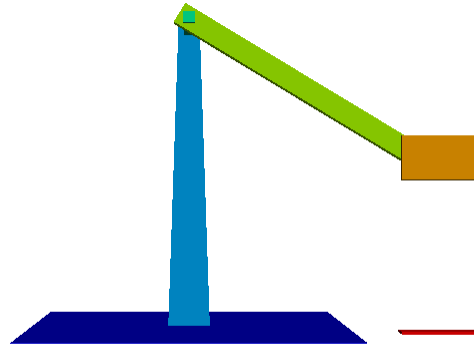
<mvcirsinu>: sinusoidal circular movement

axisp1 first point of the rotation axis
axisp2 second point of the axis
freq frequency
ampl amplitude
phase phase

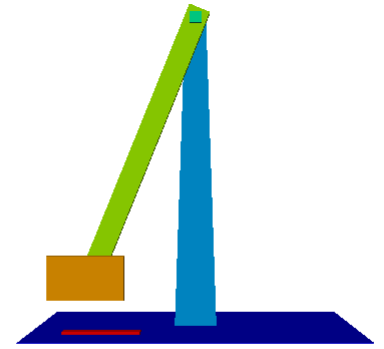
• *Motion07*: sinusoidal movement (<mvrectsinu />, <mvrotsinu />, <mvcircsinu />)



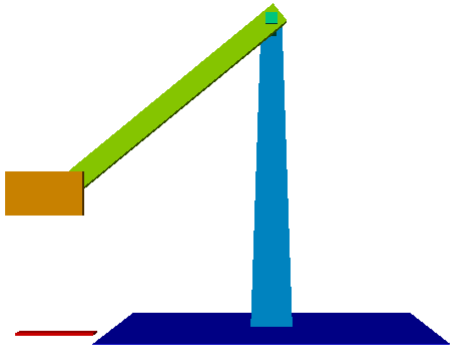
Time: 0.00 s



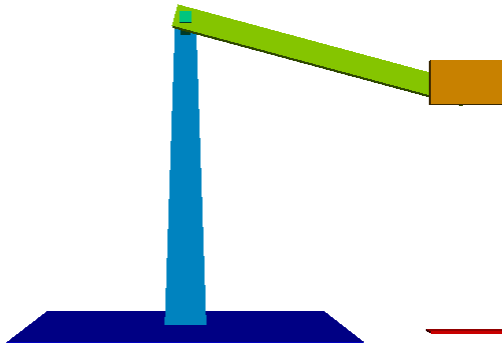
Time: 1.40 s



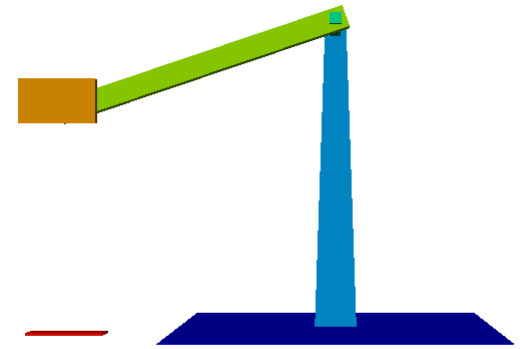
Time: 2.80 s



Time: 4.20 s



Time: 5.60 s



Time: 7.00 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>
- <objreal ref="150">
  <begin mov="1" start="0"/>
  - <mvpredef id="1" duration="8" next="2">
    <file name="motion08mov_f3.out" fields="4" fieldtime="0" fieldx="1" fieldy="2"/>
  </mvpredef>
  - <mvrect id="2" duration="-1">
    <vel x="0" y="0" z="-0.02"/>
  </mvrect>
</objreal>
- <objreal ref="151">
  <begin mov="1" start="0"/>
  - <mvpredef id="1" duration="10">
    <file name="motion08mov_f3.out" fields="4" fieldtime="0" fieldx="1" fieldz="3"/>
  </mvpredef>
</objreal>
- <objreal ref="152">
  <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>
```

<mvpredef /> or <mvfile />:
prescribed motion loaded from a file

name name of the file
fields number of columns of the file
fieldtime column with time
fieldx column with X-position
fieldy column with Y-position
fieldz column with Z-position

first field (or column) has reference "0"
second field (or column) has reference "1"

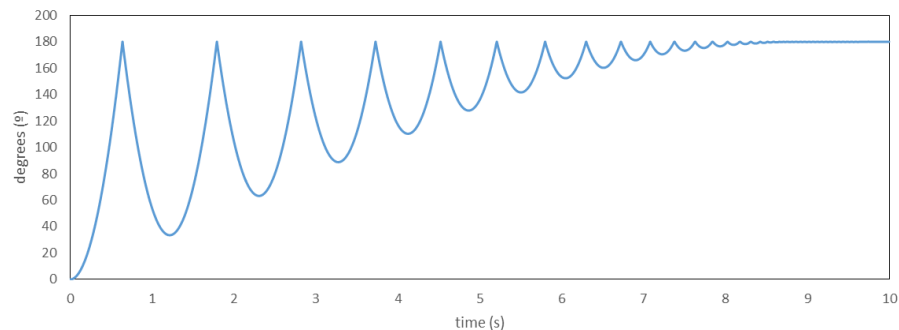
- *Motion09*: predefined movement with data from an external file (<mvrotfile />)

<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>
  <objreal ref="1">
    <begin 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>
    <mvrotfile id="2" duration="9" anglesunits="radians">
      <file name="Motion09mov_rad.csv" />
      <axisp1 x="1" y="-1" z="0.03" />
      <axisp2 x="1" y="1" z="0.03" />
    </mvrotfile>
  </objreal>
</motion>
```



XML file EXECUTION-SPECIAL-WAVEPADDLER-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)" />
  <pistondir x="1" y="0" z="0" comment="Movement direction (def=(1,0,0))" />
  <waveheight value="0.1" comment="Wave height" />
  <waveperiod value="1.3" comment="Wave period" />
  <phase 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="2" zpos="-0.15"
    comment="Saves motion data. xpos and zpos are optional. zpos=-depth" />
</piston>
```

- *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)

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"
  | | |
  | | | 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"
  | | |
  | | | 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>
```

Parameters for execution in DualSPHysics

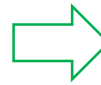
```
<parameters>
  <parameter key="PosDouble" value="1" comment="Precision in particle interaction 0:Simple, 1:Double, 2:Uses and saves double (default=0)" />
  <parameter key="StepAlgorithm" value="1" comment="Step Algorithm 1:Verlet, 2:Symplectic (default=1)" />
  <parameter key="VerletSteps" value="40"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


XML file

Case_Def.xml



GENCASE



Case.xml

```
<case>
<casedef>
<constantsdef>
<lattice bound="1" fluid="1" />
<gravity x="0" y="0" z="-9.81" comment="Gravitational acceleration" units_comment="m/s^2" />
<cofnumber value="0.2" comment="Coefficient to multiply Dt" />
<chwl value="0" auto="true" comment="Maximum still water level to calculate speedofsound using coefsound" units_comment="metres (m)" />
<speedsystem value="0" auto="true" comment="Maximum system speed (by default the dam-break propagation is used)" />
<coefsound value="0" comment="Coefficient to multiply speedsystem" />
<speedsound value="0" auto="true" comment="Speed of sound to use in the simulation (by default speedofsound=coefsound*speedsystem) />
<coefh value="1.0" comment="Coefficient to calculate the smoothing length (B=coefficient*sqrt(3*dp/2) in 3D) />
<gamma value="7" comment="Polytropic constant for water used in the state equation" />
<rhop0 value="1000" comment="Reference density of the fluid" units_comment="kg/m3" />
</constantsdef>
<mkconfig boundcount="240" fluidcount="10" />
<geometry>
<definition dp="0.01" units_comment="metres (m)">
<pointmin x="-1" y="0" z="-1" />
<pointmax x="4.5" y="0" z="3.5" />
</definition>
<commands>
<mainlist>
<setdrawmode mode="full" />
<setmkfluid mk="0" />
<drawbox>
<boxfill>solid</boxfill>
<point x="0" y="-1" z="0" />
<size x="1" y="2" z="2" />
</drawbox>
<setmkbound mk="0" />
<drawbox>
<boxfill>bottom | left | right | front | back</boxfill>
<point x="0" y="-1" z="0" />
<size x="4" y="2" z="3" />
</drawbox>
</mainlist>
</commands>
</geometry>
</casedef>
<execution>
<parameters>
<parameter key="StepAlgorithm" value="1" comment="Step Algorithm 1:Verlet, 2:Symplectic (default=1)" />
<parameter key="VerletSteps" value="40" comment="Verlet only: Number of steps to apply Euler timestepping (default=40)" />
<parameter key="Kernel" value="2" comment="Interaction Kernel 1:Cubic Spline, 2:Wendland (default=2)" />
<parameter key="ViscoTreatment" value="1" comment="Viscosity formulation 1:Artificial, 2:Laminar+SPS (default=1)" />
<parameter key="Visco" value="0.02" comment="Viscosity value" />
<parameter key="ViscoBoundFactor" value="1" comment="Multiply viscosity value with boundary (default=1)" />
<parameter key="DeltaSPH" value="0" comment="DeltaSPH value, 0.1 is the typical value, with 0 disabled (default=0)" />
<parameter key="Shifting" value="0" comment="Shifting mode 0:None, 1:Ignore bound, 2:Ignore fixed, 3:Full (default=0)" />
<parameter key="ShiftingCoef" value="-2" comment="Coefficient for shifting computation (default=-2)" />
<parameter key="ShiftingTFS" value="1.5" comment="Threshold to detect free surface. Typically 1.5 for 2D and 2.75 for 3D (default=0)" />
<parameter key="RigidAlgorithm" value="1" comment="Rigid Algorithms 1:SPH, 2:DBM (default=1)" />
<parameter key="PFreeze" value="0.0" comment="Time to freeze the floatings at simulation start (warmup) (default=0) units_comment="seconds" />
<parameter key="CoefDMin" value="0.05" comment="Coefficient to calculate minimum time step dtmin=coefdtmin*h/speedsound (default=0.05)" />
<parameter key="DtIn" value="0.0001" comment="Initial time step (default=h/speedsound) units_comment="seconds" />
<parameter key="DtMin" value="0.00001" comment="Minimum time step (default=coefdtmin*h/speedsound) units_comment="seconds" />
<parameter key="DtFixed" value="DtFixed.dat" comment="Dt values are loaded from file (default=disabled)" />
<parameter key="DtAllParticles" value="0" comment="Velocity of particles used to calculate DT. 1:All, 0:Only fluid/floating (default=0)" />
<parameter key="TimeMax" value="0.72" comment="Time of simulation" units_comment="seconds" />
<parameter key="TimeOut" value="0.01" comment="Time out data" units_comment="seconds" />
<parameter key="InoZ" value="1" comment="Increase of Z" units_comment="decimal" />
<parameter key="PartsOutMax" value="1" comment="Allowed k/100 of fluid particles out the domain (default=1)" units_comment="decimal" />
<parameter key="RhopOutMin" value="700" comment="Minimum rhop valid (default=700) units_comment="kg/m3" />
<parameter key="RhopOutMax" value="1300" comment="Maximum rhop valid (default=1300) units_comment="kg/m3" />
</parameters>
</execution>
</case>
```

```
<case>
<casedef>
<constantsdef>
<lattice bound="1" fluid="1" />
<gravity x="0" y="0" z="-9.81" comment="Gravitational acceleration" units_comment="m/s^2" />
<cofnumber value="0.2" comment="Coefficient to multiply Dt" />
<chwl value="0" auto="true" comment="Maximum still water level to calculate speedofsound using coefsound" units_comment="metres (m)" />
<speedsystem value="0" auto="true" comment="Maximum system speed (by default the dam-break propagation is used)" />
<coefsound value="0" comment="Coefficient to multiply speedsystem" />
<speedsound value="0" auto="true" comment="Speed of sound to use in the simulation (by default speedofsound=coefsound*speedsystem) />
<coefh value="1.0" comment="Coefficient to calculate the smoothing length (B=coefficient*sqrt(3*dp/2) in 3D) />
<gamma value="7" comment="Polytropic constant for water used in the state equation" />
<rhop0 value="1000" comment="Reference density of the fluid" units_comment="kg/m3" />
</constantsdef>
<mkconfig boundcount="240" fluidcount="10" />
<geometry>
<definition dp="0.01" units_comment="metres (m)">
<pointmin x="-1" y="0" z="-1" />
<pointmax x="4.5" y="0" z="3.5" />
</definition>
<commands>
<mainlist>
<setdrawmode mode="full" />
<setmkfluid mk="0" />
<drawbox>
<boxfill>solid</boxfill>
<point x="0" y="-1" z="0" />
<size x="1" y="2" z="2" />
</drawbox>
<setmkbound mk="0" />
<drawbox>
<boxfill>bottom | left | right | front | back</boxfill>
<point x="0" y="-1" z="0" />
<size x="4" y="2" z="3" />
</drawbox>
</mainlist>
</commands>
</geometry>
</casedef>
<execution>
<parameters>
<parameter key="StepAlgorithm" value="1" comment="Step Algorithm 1:Verlet, 2:Symplectic (default=1)" />
<parameter key="VerletSteps" value="40" comment="Verlet only: Number of steps to apply Euler timestepping (default=40)" />
<parameter key="Kernel" value="2" comment="Interaction Kernel 1:Cubic Spline, 2:Wendland (default=2)" />
<parameter key="ViscoTreatment" value="1" comment="Viscosity formulation 1:Artificial, 2:Laminar+SPS (default=1)" />
<parameter key="Visco" value="0.02" comment="Viscosity value" />
<parameter key="ViscoBoundFactor" value="1" comment="Multiply viscosity value with boundary (default=1)" />
<parameter key="DeltaSPH" value="0" comment="DeltaSPH value, 0.1 is the typical value, with 0 disabled (default=0)" />
<parameter key="Shifting" value="0" comment="Shifting mode 0:None, 1:Ignore bound, 2:Ignore fixed, 3:Full (default=0)" />
<parameter key="ShiftingCoef" value="-2" comment="Coefficient for shifting computation (default=-2)" />
<parameter key="ShiftingTFS" value="1.5" comment="Threshold to detect free surface. Typically 1.5 for 2D and 2.75 for 3D (default=0)" />
<parameter key="RigidAlgorithm" value="1" comment="Rigid Algorithms 1:SPH, 2:DBM (default=1)" />
<parameter key="PFreeze" value="0.0" comment="Time to freeze the floatings at simulation start (warmup) (default=0) units_comment="seconds" />
<parameter key="CoefDMin" value="0.05" comment="Coefficient to calculate minimum time step dtmin=coefdtmin*h/speedsound (default=0.05)" />
<parameter key="DtIn" value="0.0001" comment="Initial time step (default=h/speedsound) units_comment="seconds" />
<parameter key="DtMin" value="0.00001" comment="Minimum time step (default=coefdtmin*h/speedsound) units_comment="seconds" />
<parameter key="DtFixed" value="DtFixed.dat" comment="Dt values are loaded from file (default=disabled)" />
<parameter key="DtAllParticles" value="0" comment="Velocity of particles used to calculate DT. 1:All, 0:Only fluid/floating (default=0)" />
<parameter key="TimeMax" value="0.72" comment="Time of simulation" units_comment="seconds" />
<parameter key="TimeOut" value="0.01" comment="Time out data" units_comment="seconds" />
<parameter key="InoZ" value="1" comment="Increase of Z" units_comment="decimal" />
<parameter key="PartsOutMax" value="1" comment="Allowed k/100 of fluid particles out the domain (default=1)" units_comment="decimal" />
<parameter key="RhopOutMin" value="700" comment="Minimum rhop valid (default=700) units_comment="kg/m3" />
<parameter key="RhopOutMax" value="1300" comment="Maximum rhop valid (default=1300) units_comment="kg/m3" />
</parameters>
<particles>
<particles np="21001" nb="1001" nbf="1001" mkboundfirst="11" mkfluidfirst="1">
<fixed mkbound="0" mk="11" begin="0" count="1001" />
<fluid mkfluid="0" mk="1" begin="1001" count="2000" />
</particles>
<constants>
<gravity x="0" y="0" z="-9.81" units_comment="m/s^2" />
<cofnumber value="0.2" />
<gamma value="7" />
<rhop0 value="1000" units_comment="kg/m3" />
<dp value="0.01" units_comment="metres (m)" />
<ch value="1.41421356237-00" units_comment="metres (m)" />
<cb value="1.11553714298+006" units_comment="metres (m)" />
<massbound value="1.0000000000E-001" units_comment="kg" />
<massfluid value="1.0000000000E-001" units_comment="kg" />
</constants>
<motion />
</execution>
</case>
```

XML file

Summary of the number of created particles and computed constants

```
<particles np="21001" nb="1001" nbf="1001" mkboundfirst="11" mkfluidfirst="1">  
  <fixed mkbound="0" mk="11" begin="0" count="1001" />  
  <fluid mkfluid="0" mk="1" begin="1001" count="20000" />  
</particles>  
<constants>  
  <gravity x="0" y="0" z="-9.81" units_comment="m/s^2" />  
  <cflnumber value="0.2" />  
  <gamma value="7" />  
  <rhop0 value="1000" units_comment="kg/m^3" />  
  <dp value="0.01" units_comment="metres (m)" />  
  <h value="1.4142135624E-002" units_comment="metres (m)" />  
  <b value="1.1155371429E+006" units_comment="metres (m)" />  
  <massbound value="1.0000000000E-001" units_comment="kg" />  
  <massfluid value="1.0000000000E-001" units_comment="kg" />  
</constants>  
<motion />
```

np=total number of particles
nb=boundary particles
nbf=fixed boundary particles
and final **mk** of the objects

NOTE value of final "mk"
mk=mkbound+11
mk=mkfluid+1

$mass = \rho_{hop0} * dp * dp * dp$ in 3D
 $mass = \rho_{hop0} * dp * dp$ in 2D

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

Outline of Presentation

- Pre-processing tool: GenCase
- XML input file
- DesignSPHysics: **new GUI !!!**
- Future developments



DesignSPHysics

A simple user interface for DualSPHysics



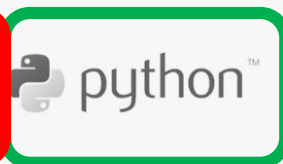
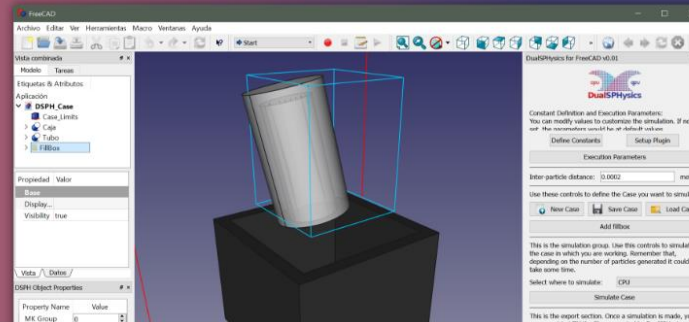
Andrés Vieira



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DesignSPHysics

A simple user interface for DualSPHysics





DesignSPHysics

A simple user interface for DualSPHysics

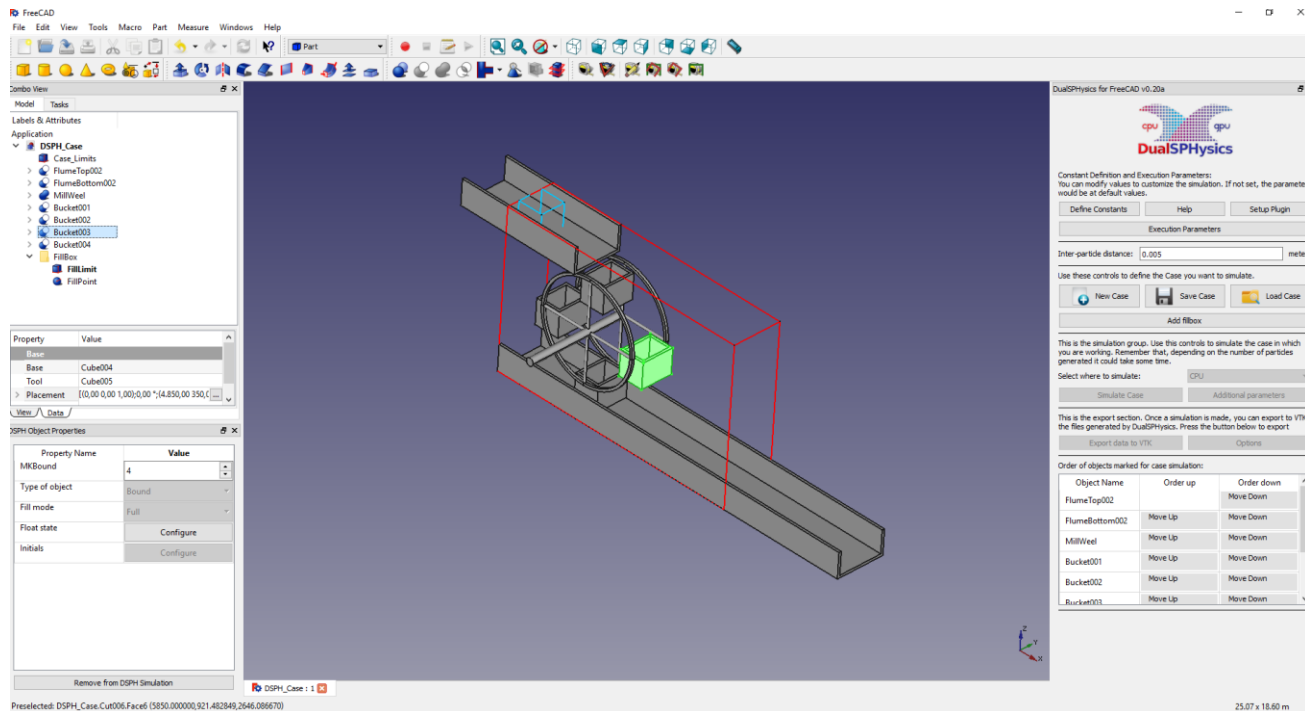


Modular: Integrates with FreeCAD, it is simple and easy to use.

Configurable: No need to edit external files or work with messy code.

Not only case design: It can run a simulation clicking on a button.

Coherent GUI: Easy to use, easy to learn.



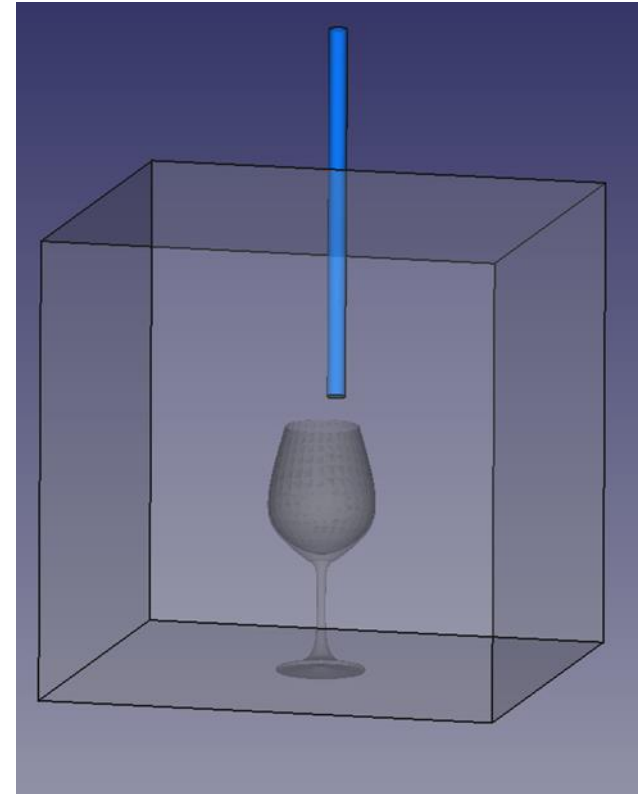
Let us try DesignSPHysics

The screenshot displays the DesignSPHysics v0.1 BETA SNAPSHOT.01 interface. The main 3D view shows a simulation setup within a wireframe cube. A blue cylinder is positioned vertically above a wine glass. The left sidebar shows the 'DSPH_Case' application with objects 'Case_Limits', 'Cube', 'Cylinder', and 'cup'. The 'Cylinder' object is selected, showing its properties: Radius: 10 mm, Height: 0.4 m, Angle: 360.00°. The 'DSPH Object Properties' panel shows 'MKFluid' set to 1, 'Type of object' as Fluid, and 'Fill mode' as Full. The right sidebar contains simulation controls, including 'Define Constants', 'Help', 'Setup Plugin', 'Execution Parameters', 'Inter-particle distance: 0.001 meters', 'New Case', 'Save Case', 'Load Case', 'Add filbox', 'Import XML', 'Simulate Case', 'Additional parameters', 'Export data to VTK', and 'Options'. The 'Order of objects marked for case simulation' table is as follows:

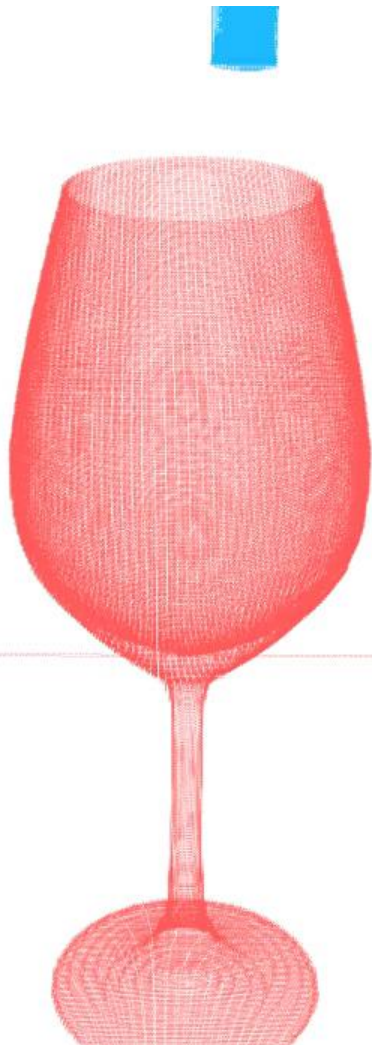
| Object Name | Order up | Order down |
|-------------|----------|------------|
| Cube | | Move Down |
| Cylinder | Move Up | Move Down |
| cup | Move Up | |

STEPS in DesignSPHysics

- Install FreeCAD version 0.16 or higher (<https://sourceforge.net/projects/free-cad/>)
- Download DesignSPHysics beta-version-> Installer.exe
- Open FreeCAD
- *Opens a dialog to let you execute a recorded macro: Use DSPH.py at %appdata%/FreeCAD/Macro*
- New Case: DSPH_Case (**CaseWine**)
 - Case_Limits
 - Base/Placement/Position: $x=-0.5\text{m}$, $y=-0.5\text{m}$, $z=-0.5\text{m}$
 - Box: Length=1m, Width=1m, Height=1.5m
 - Cube
 - Base/Placement/Position: $x=-0.25\text{m}$, $y=-0.25\text{m}$, $z=0\text{m}$
 - Box: Length=0.5m, Width=0.5m, Height=0.5m
 - ADD TO DSPH SIMULATION
 - MKBound=0, Type of object=Bound, Fill mode=Wire
 - Cylinder
 - Base/Placement/Position: $x=0.015\text{m}$, $y=0\text{m}$, $z=0.3\text{m}$
 - Cylinder: Radius=0.01m, Height=0.4m, Angle=360°
 - ADD TO DSPH SIMULATION
 - MKFluid=1, Type of object=Fluid, Fill mode=Full
 - Import: cup.stl (reads assuming units in mm)
 - ADD TO DSPH SIMULATION
 - MKBound=1, Type of object=Bound, Fill mode=Face
 - Inter-particle distance=0.001 meters
 - Define Constants
 - Execution Parameters
 - Time of simulation=1 seconds
 - Time out data=0.005 seconds
 - Setup Plugin: path of GenCase4, DualSPHysics4, PartVTK4
 - Save Case: Executes GenCase4 (creates .XML and initial .BI4)
 - Simulate Case: Executes DualSPHysics4 (CPU or GPU)
 - Export data to VTK: Executes PartVTK4 (VTK of all particles)



Visualisation in Paraview loading VTK files

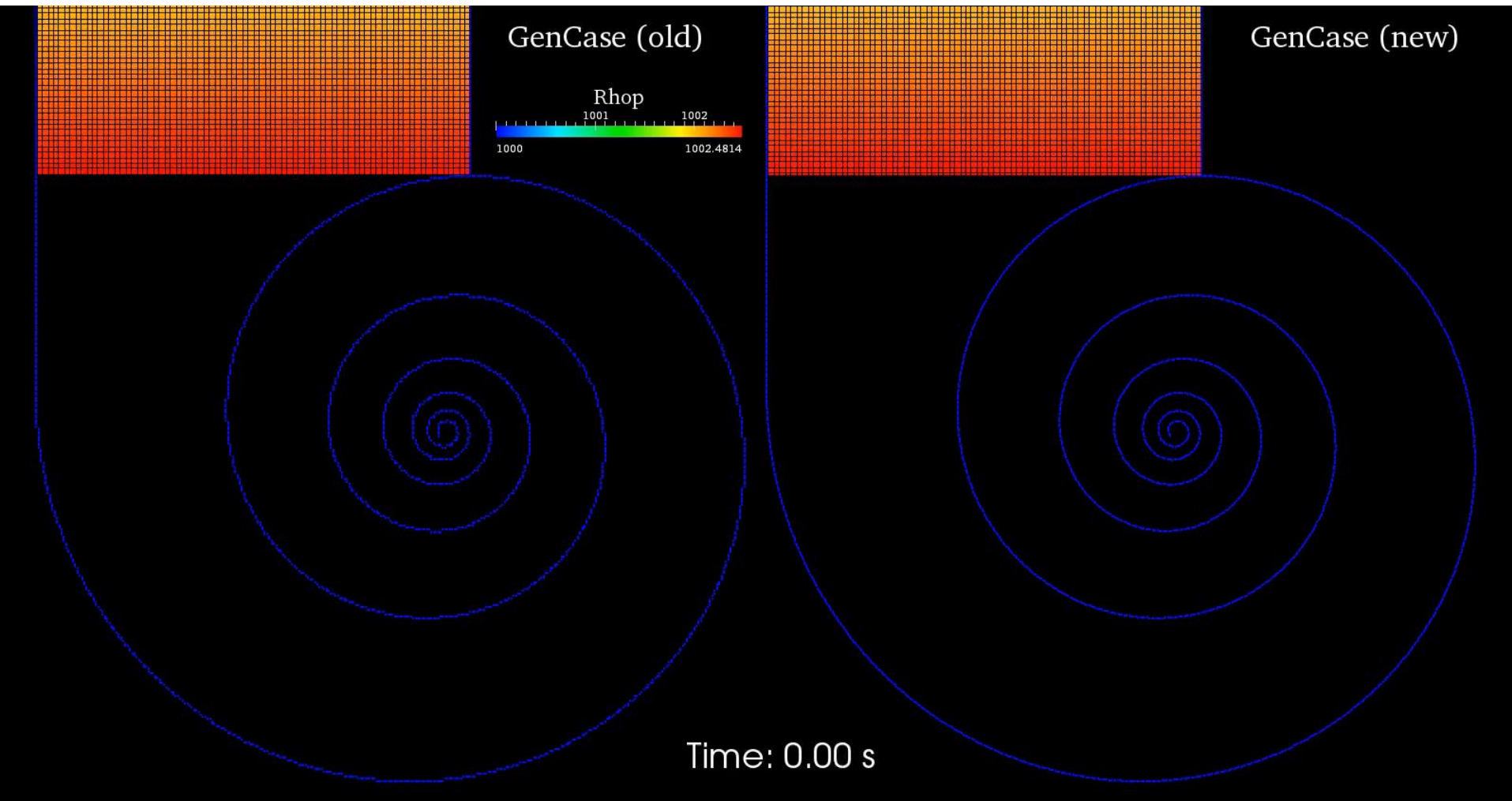


Outline of Presentation

- Pre-processing tool: GenCase
- XML input file
- DesignSPHysics: **new GUI !!!**
- **Future developments**

Future developments

- FreePoints



Future developments

- FreePoints
- Computation of normals (new BCs)
- Release of DesignSPHysics

Thank you

Acknowledgements

- DualSPHysics team: all developers and contributors
- Andrés Vieira

Website

Free open-source **DualSPHysics** code:

<http://www.dual.sphysics.org>

