

# TRUST Tutorial V1.8.1beta

CEA Saclay

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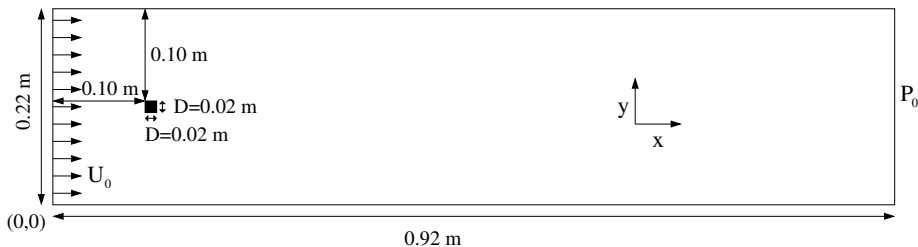
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# Initialization

- First, initialize the TRUST environment.
  - On CEA Saclay PCs and callisto cluster, TRUST versions are available with (e.g. X.Y.Z=1.7.6):  
**source /home/triou/env\_TRUST\_X.Y.Z.sh**
  - On your own computer, download and install the latest version of TRUST in your local folder \$MyPathToTRUSTversion (unless this was already performed), then write on the terminal:  
**source \$MyPathToTRUSTversion/env\_TRUST.sh**
- Second, several editors (**vim**, **emacs**, **nedit**, **gedit**) can be configured to highlight TRUST keywords in the data files. If you prefer using **nedit**, please do the following:
  - Run **nedit**, and select Preferences → Save Defaults.
  - Then run **trust -config nedit**, the message "nedit.rc updated" should appear.

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# Geometry



- Fluid:  $\mu = 3.7 \cdot 10^{-5} \text{ kg.m}^{-1}.\text{s}^{-1}$ ,  $\rho = 2 \text{ kg.m}^{-3}$  and  $Re = \frac{U_0 H_{inlet} \rho}{\mu} = \frac{1 \times 0.22 \times 2}{3.7 \cdot 10^{-5}} = 11891$
- Boundary conditions:
  - Inlet with uniform velocity:  $U_0 = 1 \text{ m.s}^{-1}$
  - Outlet with constant pressure:  $P_0 = 0$
  - Square cylinder: No-slip wall
  - Upper and Lower walls: Symmetry

# Create a study

- First, you must already have initialized TRUST environment.
- Open a terminal and run the commands to create a directory for your studies:  
**mkdir -p Formation\_TRUST/yourname**  
**cd Formation\_TRUST/yourname**
- Copy the test case from the TRUST database to your working directory with the command:  
**trust -copy Obstacle**  
**cd Obstacle**
- Ask for trust script options:  
**trust -help**
- Ask for help on the options of TRUST executable:  
**trust Obstacle -help\_trust**
- Run the test case with the command:  
**trust Obstacle**

# Probes and parameters

- Edit the data file `Obstacle.data` and set the time step to 0.004s:  
**nedit Obstacle.data &**
- Add to the post-processing block of `Obstacle.data` the following elements:
  - A pressure probes segment (22 probes between points (0.01, 0.12) and (0.91, 0.12)).
  - A velocity probes segment (22 probes between points (0.92, 0.00) and (0.92, 0.22)) to plot the velocity profile behind the square cylinder.
  - Add the vorticity to the fields being post-processed and change the saving period to 0.5s. To find the appropriate keyword for this field, you can open the Generic Guide with:  
**trust -doc &**
  - Replace the keyword "**format lml**" with "**format lata**" inside the block, just before the keyword **fields** in order to use the post-processing tool VisIt during and/or after the calculation.



# Visualization during the calculation

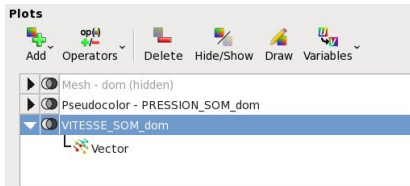
- You have access to useful resources in the `$TRUST_ROOT/index.html` file with your favorite browser (e.g.: firefox). Take few minutes to find test case examples containing a particular keyword using the Keywords link:  
**firefox \$TRUST\_ROOT/index.html & or trust -index &**
- Launch the "PLOT2D" tool with:  
**trust -evol Obstacle &**  
 This tool allows to launch calculation and visualize results
  - To run the calculation, click on the button "Start computation!" at the lower left corner of the window.
  - Visualization:
    - Select "PRESSION(X=0.13,Y=0.105)" in the left list and click on "Plot" to draw the evolution of the pressure at the probe location.
    - Check the velocity profile behind the square cylinder by plotting "VITESSE\_X(X=0.14,Y=0.115)" and "VITESSE\_Y(X=0.14,Y=0.115)".
    - Visualize the equation residuals on the same plot, select " $Ri = \max \left| \frac{dV}{dt} \right|$ " and " $residu = \max |Ri|$ " using the button "Plot on same" or select the two graphs with "Ctrl" button and "plot".

# VisIt

- To quit this tool, close the GUI.
- Once the calculation is finished, visualize the results with the graphical tool VisIt directly: **visit &**  
or using "PLOT2D" tool: **trust -evol Obstacle &** and click on "Visualisation" on the right menu.
  - First, we are going to configure VisIt: In the menu File → Open file, select Off instead of Smart for File grouping option. For the Filter, specify \*.lata to list only the lata files (results). Then save your choices, in the menu Options → Save Settings.
  - In the menu File → Open file, select the Obstacle.lata file.
  - Visualize the mesh in the "Plots" area with "Add → Mesh → dom" then click on the button "Draw". Zoom and move the mesh in the right window. You can un-zoom with right button (View → Reset view) or with a combination of "Ctrl" keypad and left button.
  - Visualize the pressure field (Plots area: "Add → Pseudocolor → PRESSION\_SOM\_dom" + Draw then select the last time on the Time slider)
  - Suppress or hide the mesh (Select Mesh then click on Delete or Hide/Show).

# VisIt

- Visualize the velocity field (Plots area: "Add → Vector → VITESSE\_SOM\_dom" + Draw). You can change each plot attributes:
  - ◇ click once onto the small arrow "►" then
  - ◇ double click on the item Vector (cf the figure below). For example, change the number of vectors being plotted (by default 400, set it to 40000 then click the button "Make default" and save definitively this modification with the menu Options → Save Settings). You need to click "Apply" to update. Then click "Dismiss" to close the window.



- Print your visualization (File → Save window): a PNG file is created into your working directory.

# VisIt

- Add a second screen with "Windows → Layouts → 1x2",
- Plot a pressure horizontal profile:
  - ◇ select the pressure field,
  - ◇ on the visualisation, use the right click and select "Mode → Line out",
  - ◇ then define your profile with left button,
  - ◇ click on the origin point, let the left button pushed, and release at the end point.
  - ◇ The profile is shown on the second window.
- You notice that it is necessary to update (button Draw) the right window after adding a new plot or changing an option. It is possible to automatically update by activating "Auto apply" on the top right of the VisIt's GUI.
- You can create new fields (expression) with "Controls → Expressions → New" by using existing variables and complex functions and visualize it.
- You can animate your visualization and/or create a movie (File → Save movie)
- You can operate calculations on variables with complex queries (Controls → Query),
- You can save a complex session (File → Save session) and reopen it during a next analyze with VisIt (File → Restore session),

# Outputs and resuming calculation

- During a 3D visualization, you will use one of the available Operators (In Plots, "Operators → Slicing → Slice") to create a 2D slice either in a 3D space, or projected to a 2D space.
- For more information on **VisIt**, you can refer to:
  - the **VisIt** website and its manuals:  
<https://wci.llnl.gov/simulation/computer-codes/visit/manuals>
  - the **VisIt** user community web site: <http://visitusers.org>
  - or send an email to the VisIt software users community at:  
[visit-users@elist.ornl.gov](mailto:visit-users@elist.ornl.gov)
- Edit the different output (\*.out) files to read the complete balances (mass, stress, energy, ...) on the whole domain or at the boundaries.
- Now we want to edit the data file in order to resume the calculation. So, open it using "PLOT2D" tool: **trust -evol Obstacle &**.

# Outputs and resuming calculation

- Find the last backup time of the previous calculation in the .err file (or in the bottom right file in the "PLOT2D" tool if it is still running).
  - Edit your data file with "Edit data" or "Edit with xdata", then modify **tinit**, **tmax** values in the object "mon\_schema".
  - Add in the problem description block just before the last "}":  
**reprise binaire Obstacle\_pb.sauv**  
(The file "Obstacle\_pb.sauv" must have been created during the first run.)
  - Save and close the window.
  - Resume the calculation again with "Start calculation!" button. You can see that values are added to the first probes during the new calculation.
- ⇒ *Remark:* to resume your calculation, you can also use the keyword **resume.last\_time** instead of **reprise** and only change the **tmax** value (cf Reference Manual).

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## Parallel calculation: 1st method

The goal of this exercise is to introduce parallelism in the data file of the previous exercise.

- Go to the previous study (should be done) and after you had suppressed the **reprise** keyword and set **tinit** to 0 again in the *Obstacle.data* file, create two new files:

```
cd Formation_TRUST/yourname/Obstacle
mkdir PARA1
cd PARA1
cp ../Obstacle.data DEC_Obstacle.data
cp ../Obstacle.data PAR_Obstacle.data
cp ../Obstacle.geo .
```

- Edit the first file (*DEC\_Obstacle.data*) to create the partition of the mesh.



# Parallel calculation: 1st method

- In this file, uncomment the block around the **Partition** keyword.
  - Here, the partitioning tool **Metis** is used. We cut in **nb\_parts** blocks, here in 2.
  - The overlapping width **Larg\_joint** between two parts of the partition should be defined according to the numerical scheme higher order, generally the convective scheme. Its value is generally 1 for a second-order scheme, and 2 for third- or fourth-order schemes such as Quick scheme.
  - In VEF, you should use **2** for **Larg\_joint** except when partitioning a domain where only the conduction equation will be solved.
  - At least, the keyword **zones\_name** is useful to define the name of the files containing the partitioned mesh and to write these files.
  - Notice the presence of the keyword **End** in the "Partition" block: the code will stop reading the data file at this line!
- Run the data file: **trust DEC\_Obstacle**
- Check that the partitioned mesh files DOM\_0000.Zones and DOM\_0001.Zones are generated inside your working directory: **ls \*.Zones**

## Parallel calculation: 1st method

- Now, edit the file `PAR_Obstacle.data` and comment the read of the mesh (using `#` tags of the 'BEGIN/END MESH' comments).
- Uncomment the **Scatter** keyword which will read the partitioned mesh.
- Visualize it with VisIt:  
**trust -mesh PAR\_Obstacle**
- Now, run a parallel calculation with TRUST:  
**trust PAR\_Obstacle 2**
- The post-processing step is identical in sequential or parallel mode. You have the probes into the `.son` files and the whole fields in the `.lata` files. To run VisIt with the command line:  
**visit -o PAR\_Obstacle.lata &**
- Select the last time step and visualize the blocks (with Plots: Add → Subset → blocks) which represent the parts of the domain partition, then the velocity fields. You can also visualize a field only on a selected part (block) with the menu Control → Subset.

## Parallel calculation: 2nd method

- To visualize probes after the end of the calculation, you can run the command line:  
**trust -evol PAR\_Obstacle &**
- The existing tool **trust -partition** is useful on the data files which have the marks MAILLAGE/MESH, DECOUPAGE/PARTITION and LECTURE/SCATTER. If you run the following commands:
- First we will create a new working directory with our data files:  
**cd Formation\_TRUST/yourname/Obstacle**  
**mkdir PARA2**  
**cd PARA2**  
**cp ../Obstacle.data exemple.data**  
**cp ../Obstacle.geo .**
- Then we run the command:  
**trust -partition exemple 3**  
**ls**

## Parallel calculation: 2nd method

- It creates:
  - a SEQ\_exemple.data file which is a copy of the sequential data file exemple.data,
  - a DEC\_exemple.data file which is the first data file to be run. It is immediately run by the command line **trust -partition** to create a partition (with 3 sub zones here), located in the **\*\*\*.Zones** files.

### ls \*.Zones

Note that the TRUST code stops reading this file at the keyword "End" just before the "# END PARTITION #" block.

- a PAR\_exemple.data file which is the data file for the parallel calculation. It uses the **\*\*\*.Zones** files to read the mesh through the line "Scatter DOM.Zones dom". Note that the meshing and cut of the mesh are commented here.
- Then you have to run the calculation by the usual command completed by the number of processors needed:  
**trust PAR\_exemple 3**

# Parallel calculation: 2nd method

## Useful information:

- Be careful when you want to modify your data file! You have two possibilities:
  - you want to modify your mesh,
  - you want to modify the calculation parameters.
- For the first one, you can modify:
  - the file `exemple.data` and run **trust -partition**. But it will erase the `DEC_exemple.data`, `SEQ_exemple.data` and `PAR_exemple.data` files and create new zones. Then it will run the new `DEC_exemple.data` file which gives your new `***.Zones` files or,
  - the meshing part of file `DEC_exemple.data` and run it with:  
**trust DEC\_exemple.data**

Then run the parallel calculation normally, on the new `***.Zones` files.  
**trust PAR\_exemple 3**

## Parallel calculation: 2nd method

- For the second possibility, you can modify:
  - the file `exemple.data` and run **trust -partition**. But it will erase the `DEC_exemple.data`, `SEQ_exemple.data` and `PAR_exemple.data` files and create new ones. Then it will run the new `DEC_exemple.data` file. Note that in that case, you don't need to re-create the mesh so you can use the second point below:
  - modify the `PAR_exemple.data` file without running **trust -partition**.

Then run the `PAR_exemple.data` file with:

**trust PAR\_exemple 3**

- Notice that if after a certain time, you want to reopen an old case and understand what you did in it without any doubts, you can create two files manually:
  - one "`BuildMeshes.data`" file only for the mesh and the cut of the mesh, and
  - one "`calculation.data`" file for the parallel calculation.

You will run it like:

**trust BuildMeshes**

**trust calculation nb\_procs**

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## Parallel calculation on a cluster

NB: On CEA Clusters, TRUST is already installed and the procedure of launching calculation is described below. Out of CEA, your cluster administrator should install and configure TRUST. In addition, submission files and procedure depend on the cluster itself and could be different from those presented below.

- Log on to the CEA cluster callisto and initialize the TRUST environment:  
`ssh -X yourlogin@callisto-login1(.intra.cea.fr)`  
`source /panfs/ixion/home/triou/TRUST/TRUST-X.Y.Z/env_TRUST.sh`  
or `source /home/triou/env_TRUST_X.Y.Z.sh`
- Copy the study Obstacle:  
`cd /panfs/ixion/home/yourlogin`  
`mkdir -p Formation_TRUST/yourname`  
`cd Formation_TRUST/yourname`  
`trust -copy Obstacle`  
`cd Obstacle`
- Open Obstacle.data and set the **format** to **lata** in the post-traitement block.
- Create automatically a 2 partitioned mesh and a parallel data file with:  
`trust -partition Obstacle`



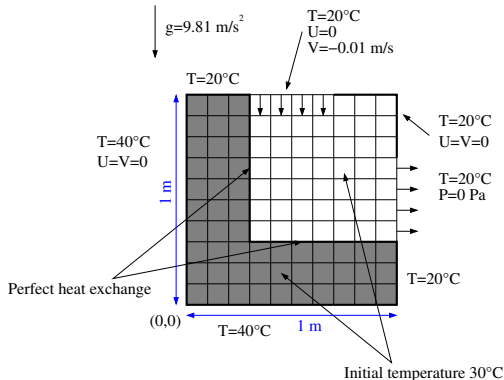
# Parallel calculation on a cluster

- For clusters, you have to create a submission file:  
**trust -create\_sub\_file PAR\_Obstacle 2**
- Open the file sub\_file and rename the job. Note that we will see only the first eight characters of the job name in the submitted jobs list.
- Submit the job with: **sbatch sub\_file**
- Check the state of the job with: **"squeue"** or **"squeue -u yourlogin"**
- You could run VisIt from the cluster but it is not recommended.
- So you can run VisIt on your PC and access/visualize the result file on callisto cluster.
  - Open your .bashrc file (or create it if needed).
  - Add the following lines:

```
if [[ $HOSTNAME:0:8 == "callisto" ]]
then
source /panfs/ixion/home/triou/TRUST/TRUST-X.Y.Z/env_TRUST.sh
fi
```
  - Open VisIt on your computer, then access to your files on callisto with:  
File → Open File and select callisto in the localhost menu.

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# Heat transfer (2D, VDF/VEF)



**Fluid:**

$$Pr = \frac{\mu C_p}{\lambda} = 1,$$

$$T_{ref} = 30^{\circ}\text{C},$$

$$\mu = 2 \cdot 10^{-3}\text{ kg}\cdot\text{m}^{-1}\cdot\text{s}^{-1},$$

$$\rho = 2\text{ kg}\cdot\text{m}^{-3},$$

$$\lambda = 1\text{ W}\cdot\text{m}^{-1}\cdot\text{K}^{-1}$$

$$C_p = 500\text{ J}\cdot\text{kg}^{-1}\cdot\text{K}^{-1},$$

$$\beta = 1\cdot 10^{-4}\text{ K}^{-1}$$

**Solid:**

$$\rho = 1000\text{ kg}\cdot\text{m}^{-3}$$

$$\lambda = 250\text{ W}\cdot\text{m}^{-1}\cdot\text{K}^{-1}$$

$$C_p = 100\text{ J}\cdot\text{kg}^{-1}\cdot\text{K}^{-1}$$

# Heat transfer (2D, VDF/VEF)

- Create a new study Coupling\_VDF by copying the docond study:  
**cd Formation\_TRUST/yourname**  
**trust -copy docond**  
**mv docond Coupling\_VDF**  
**cd Coupling\_VDF**
- Check the fluid and solid characteristics inside the docond.data file.
- This coupled problem is constituted by 2 domains of calculation with a mesh of 10x10 cells ( $\Delta x = \Delta y = 0.1m$ ) created with 3 blocks.
- Now open your data file with "Plot2D" tool:  
**trust -evol docond &**
- Click on "Edit data" (you can also use "Edit with xdata").
- We want to modify the data file to have the 2 domains on a mesh of 40x40 cells ( $\Delta x = \Delta y = 0.025m$ ).

# Heat transfer (2D, VDF/VEF)

- Change the number of nodes for each block like this:  
 First block (Cavite1): 4 11  $\rightarrow$  13 41  
 Second block (Cavite2): 8 4  $\rightarrow$  29 13  
 Third block (Cavite3): 8 8  $\rightarrow$  29 29
- Check your new mesh with:  
**trust -mesh docond**
- Change "**format lml**" to "**format lata**" into the two problems definition
- Click on "Save" and close the window.
- Run the calculation with "Start computation!" and check the evolution.
- Then post-process the temperature field with VisIt tool: "Visualization" button. A natural convection cell appears.
- Change the color tables for the temperature to have the same one on the 2 domains. Close VisIt.
- We are going to change the discretization of the test case: triangulate the domains with the keyword **Trianguler\_H** (refer to the Reference Manual).

# Heat transfer (2D, VDF/VEF)

- Then give an unstructured aspect to the 2 meshes using the following syntax:  
**Transformer name\_of\_domain  $x*(1-0.5*y*y)$   $y*(1+0.1*x*y)$**
- Substitute the discretization VDF (pressure nodes at the element center) to VEFPreP1B (pressure nodes at the element's center and nodes).
- Close the Plot2D tool.
- Check the meshes with:  
**trust -mesh docond**
- Run the calculation with:  
**trust docond**
- Open the IHM:  
**trust -evol docond**
- Select ' $Ri=\max\_pb1|dT/dt|$ ', ' $Ri=\max\_pb2|dT/dt|$ ', ' $Ri=\max\_pb2|dV/dt|$ ', ' $residu=\max|Ri|$ ' with "Ctrl" button and click on 'Plot on same'.
- To see when convergence is reached, select a probe (for example temperature) and click on 'Plot'.

# Heat transfer (2D, VDF/VEF)

- If the calculation is too long, open the docond.stop file, put a 1 instead the 0 and save. The calculation will stop after the current time step and make post-process.
- Post-process the results and compare the CPU performances with VDF discretization: the VEF calculation is running  $\approx 10$  times slower (because more pressure unknowns and shorter time steps). Check the docond.out file to see the time steps for each equation (click on "Edit .out" at the upper right corner of the GUI).
- Accelerate the calculation by impliciting the diffusive term of each equation with **diffusion\_implicite** option in the explicit Euler scheme (check again the Generic Guide: **trust -doc &**).
- Run the calculation without any option:  
**trust docond**
- Now, use a fully implicit scheme (suppress **diffusion\_implicite**), by substituting **Scheme\_Euler\_Explicit** by **Scheme\_Euler\_implicit** and adding the Implicit solver "**solveur implicite**".

# Heat transfer (2D, VDF/VEF)

- Have a look at the Reference Manual for the **gmres** options and define, according to the advice given on it, a value for **facsec**, **facsec\_max**.
- Your block will look like:  
**Solveur Implicite { solveur gmres { diag seuil 1e-30 nb\_it\_max 5 impr }  
 seuil\_convergence\_implicite 0.01 }**
- Run the calculation:  
**trust -evol docond &**



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  - 3D VEF mesh
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# Low Mach number flow (2D)

- Open a terminal and create a directory using Unix commands, and copy the study **TP\_Temp\_QC\_VEF** (it is a 2D simulation of helium gas flow from left to right between two heated walls):

```
mkdir -p Formation_TRUST/yourname
```

```
cd Formation_TRUST/yourname
```

```
trust -copy TP_Temp_QC_VEF
```

```
cd TP_Temp_QC_VEF
```

- Open the Generic Guide with (it will be useful to search for keywords in this exercise): **trust -doc &**
- Edit the data file with your favorite editor (**nedit** is recommended because it is configured to recognize the TRUST syntax):

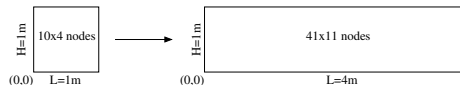
```
nedit TP_Temp_QC_VEF.data &
```

or

```
trust -evol TP_Temp_QC_VEF & and "Edit data" button.
```

# Low Mach number flow (2D)

- Edit the data file in order to:
  - Modify the geometry and the mesh:



- Add several probes (velocity, density, temperature) near the upper right corner of the geometry at location  $(x,y)=(4,1)$ .
  - Add a probe "**segment**" (with 9 points) between the locations  $(x,y)=(4,0.05)$  and  $(x,y)=(4,0.95)$  for the temperature field.
  - Write the results on the **lata** format and change the **dt\_post** period to 1s.
  - We are looking for the steady state, so suppress **tmax** keyword and change the **seuil\_statio**  $\varepsilon$  value to 10 ( $|dT/dt| < \varepsilon$  and  $dt \sim 0.001\text{s}$  so  $|dT| < 0.01$ ).
  - Add the keyword **impr** into the pressure solver to print its convergence.
  - If you use Plot2D tool, save and close the editor.
- Run the simulation with the TRUST command:  
**trust -evol TP\_Temp\_QC\_VEF &**

# Low Mach number flow (2D)

- Click on "Start computation!".
- Check mass flow rate (absolute and relative values) in the TP\_Temp\_QC\_VEF.out file:  
**nedit TP\_Temp\_QC\_VEF.out &**  
 or look at the upper small window on the right of the PLOT2D tool.
- Once the calculation finishes, visualize the results by running **VisIt**:  
**visit -o TP\_Temp\_QC\_VEF.lata &**  
 or  
 "Visualization" button on Plot2D tool.
  - Show the mesh (Plots: "Add → Mesh → dom → Draw").
  - Visualize the temperature field (Select the last Time with the slicer, then Plots: "Add → Pseudo Color → TEMPERATURE\_SOM\_dom → Draw").
  - Suppress or hide the mesh (Select "Mesh-dom" in the list of plots then "Delete" or "Hide/Show").
  - Visualize the velocity field (Add → Vector → VITESSE\_SOM\_dom → Draw).

# Low Mach number flow (2D)

- Select the Zoom mode with the right button of the mouse (Mode → Zoom) then zoom by selecting an area on the plot. To un-zoom push "Ctrl" button and select an area with the left button or with the right button select "View → Reset view".
- Print your visualization (File → Set Save options → File type → Select a type → Save): a file named visit\*\*\* is created into your working directory.
- Add a second screen with "Window → Layout → 1x2".
- Plot a horizontal profile of temperature (Select the temperature field and thanks to the right button, select "Mode → Lineout", and define your profile with left button): the profile is shown on the second window.
- Substitute the time scheme by an implicit time scheme (like **scheme\_euler\_implicit**).
- Use the **implicit** solver and specify **facsec** and **facsec\_max** parameters according to the advice given on the Reference Manual (search for the **scheme\_euler\_implicit** keyword). You can also see the instructions at the end of the Heat transfer VDF/VEF exercise on p.31.

# Low Mach number flow (2D)

- Run the calculation with this time scheme using the PLOT2D tool or:  
`trust TP_Temp_QC_VEF.data 1>TP_Temp_QC_VEF.out 2>TP_Temp_QC_VEF.err`
- Edit the file containing information about `dt` (used time step), `dt_stab` (stability time step), `facsec` ( $dt=dt\_stab*facsec$ ) and residuals evolution for each equation:  
`nedit TP_Temp_QC_VEF.dt_ev &`
- If everything is OK, try to enhance the convergence speed of the implicit solver with the value of **seuil\_convergence\_implicit** keyword (look at the `TP_Temp_QC_VEF.out` file, if the number of iterations for GMRES is comprised between 3 and 5 then it is enough to converge quickly).
- In order to resume a calculation, you will have to change the **tinit** value within the data file (pick up the last saved time in the `.err` file) and insert into the data file, in the problem definition block, the following keywords:  
`reprise binaire TP_Temp_QC_VEF_pb.sauv`

# Low Mach number flow (2D)

- Then run the calculation with:

```
trust TP_Temp_QC_VEF.data 1>TP_Temp_QC_VEF.out 2>TP_Temp_QC_VEF.err
```

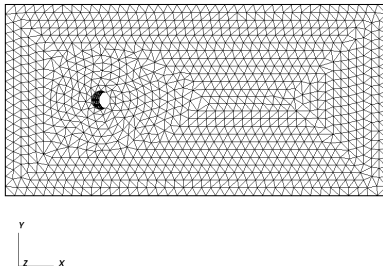
or

```
trust -evol TP_Temp_QC_VEF.data & which automatically creates the  
.out file
```

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# Periodic channel flow (3D)



**Fluid:**  $Re = 2000$ ,  $\rho = 2\text{kg.m}^{-3}$ ,  $\mu = 0.01\text{kg.m}^{-1}\text{s}^{-1}$ , initial velocity  $V0 = 1\text{m/s}$ , periodic boundary condition following the Z-direction.

- Copy the study named **P1toP1Bulle** as explained on page 7. It simulates a 3D incompressible laminar flow ( $Re = 2000$ ) with periodic boundary following the Z-direction only.
- Open the P1toP1Bulle.data file and use **RegroupeBord** keyword to merge Entree and Sortie boundaries into a single one named periox.

## Periodic channel flow (3D)

- Modify boundary conditions to apply a periodic boundary on the new boundary.
- Change the velocity initial condition to  $U_0 = (1, 0, 0)$ .
- Set the option **diffusion\_implicit** to 1 into the Euler scheme to implicit the diffusive term in the Navier-Stokes equations.
- You have now a 3D calculation with periodic boundary conditions on X- and Z-directions. Run the calculation for 30 time-steps (keyword **nb\_pas\_dt\_max**).
- Have a look at the P1toP1Bulle\_pb\_Debit.out file, check the flow rate on the periox boundary. Why does it decrease?
- Add the **Canal\_perio** source term in the Navier-Stokes equations of the data file and run again the calculation to check the flow rate evolution on 30 time steps.
- Look at pressure and viscous forces applied on the cylinder inside the .out files.

## Periodic channel flow (3D)

- Now, the calculation domain is a rotating channel according to Z direction with a constant velocity  $\Omega = 1 \text{ rad/s}$ .
- Add the **Acceleration** source term in the Navier-Stokes equations. Suppress the **nb\_pas\_dt\_max** keyword and set **tmax** to 100s.
- Add, if you wish, velocity or statistic calculation to the post-processing instructions.
- Run the calculation.
- You can create a uniformly refined mesh using, for instance, the keyword **Raffiner\_Anisotrope**.
- Then improve the calculation speed on this mesh, you can use a coarse discretization **P1** (**Read dis { P1 }** keywords) with less pressure unknowns. On this latter, it runs 3 times faster than on P1Bulle discretization but it is less accurate: 8452 unknowns compared to 49221 unknowns.

## Periodic channel flow (3D)

- Then restart the calculation with **VEFPreP1B** discretization by reading the velocity field with **Champ\_fonc\_reprise** keyword in the initial conditions for the velocity:

**vitesse champ\_fonc\_reprise P1toP1Bulle\_pb.xyz pb vitesse last\_time**

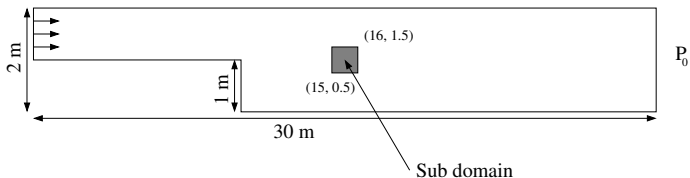
This will be useful to reach the quasi-stationary regime faster.

- You can also use implicit scheme (change the scheme to **Scheme\_Euler\_implicit** scheme and use an **Implicite** solver) only if you are looking for the stationary state.

You can also see the instructions at the end of the Heat transfer VDF/VEF exercise on p. 31.

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# Constituents and turbulent flow



**Fluid:**  $\mu = 3.7 \cdot 10^{-5} \text{ kg.m}^{-1}.\text{s}^{-1}$ ,  $\rho = 2 \text{ kg.m}^{-3}$ ,  $Re = \frac{U_0 H_{inlet} \rho}{\mu} = 54054$

## Boundary conditions:

Inlet with imposed velocity:  $U_0 = 1 \text{ m.s}^{-1}$  and constant values of  $k = 10^{-2}$  and  $\varepsilon = 10^{-3}$  (dimensionless values)

Outlet with constant pressure:  $P_0 = 0$  and constant values of  $k = 0$  and  $\varepsilon = 0$

Top and bottom walls: No-slip wall ( $U = 0$ ) and  $k$  standard flux,  $\varepsilon$  null.

- Copy the study named **Marche** as explained page 7. It simulates a 2D incompressible turbulent flow in the configuration described above using the  $k$ - $\varepsilon$  model.

# Constituents and turbulent flow

- We want to add a source of constituent's diffusion so copy the **Constituents** study which will give you examples for constituents keyword used for a 2D incompressible laminar flow.
- Edit your data file in the Marche directory. First, rename the problem in order to add concentration equations (look for the adequate keywords in the Reference Manual).
- Add 3 constituents of equal diffusivities ( $\alpha = 1m/s$ ) and associate the constituents to the problem.
- Define the concentration equation into the problem (remember that concentrations will be a vector of 3 components) with correct initial ( $C_1 = 0, C_2 = 0, C_3 = 0$ ) and boundary conditions.
- Use the Schmidt model to close the turbulence model in the concentration equation.
- Change the sources of the Navier-Stokes turbulence model to a **Source\_Transport\_K\_Eps\_aniso\_concen { C1\_eps 1.44 C2\_eps 1.92 C3\_eps 1. }** to fit with the new concentration equation.

# Constituents and turbulent flow

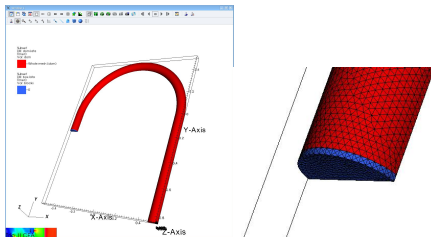
- Add into the fluid definition, the volume expansion coefficient for the concentration: **beta\_co** as a uniform field set to 0.
- You have also to add a gravity field which can be initialized to 0.
- Run the calculation to see if it is ok.
- Define a sub domain (in grey on the previous picture) with the keyword **Sous\_Zone** (like in **PCR** data file).
- Add a source term for the second constituent only ( $S_2 = 1m^{-3}$ ) applied on the sub domain thanks to the keyword **Champ\_Uniforme\_Morceaux**.
- Add **format lata** in the post-processing block.
- Add the keyword **concentration0**, **concentration1**, **concentration2** in the **fields** of the post-processing block to write the 3 concentrations into the .lata file.
- Run the calculation and check the results.



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# Turbulent flow in a curved pipe (3D)

## TrioCFD



### Goals:

- Use of a RANS or LES model.
- Use of a periodic box to initialize a fully developed turbulent flow.
- Use of the TrioCFD parallel capabilities.

# Turbulent flow in a curved pipe (3D)

## TrioCFD

- First initialize TrioCFD environment :  
`source /home/triou/env_TrioCFD_X.Y.Z.sh`  
`echo $exec`  
`echo $project_directory`
- Create a new directory and copy some data:  
`mkdir -p Formation_TRUST/yourname/PeriodicBox`  
`cd Formation_TRUST/yourname/PeriodicBox`  
`cp $project_directory/validation/share/Validation/Rapports_automatiques/Validant/pas_fini/PeriodicBox/src/* .`

Notice that this directory corresponds to an automated validation test case. If you want to run it and generate the pdf report, see the last exercise of this tutorial which is named "Validation form", but be careful this case needs more computational effort!

# Turbulent flow in a curved pipe (3D)

## TrioCFD

- There are several files :
  - *BuildMeshes.data*: To build the meshes
  - *PeriodicBoxRANS.data*: To run the flow in the box with RANS model
  - *DomainFlowRANS.data*: To run the flow in the domain with inlet steady conditions from the box domain
  - *PeriodicBoxLES.data*: To run the flow in the box with LES model
  - *DomainFlowLES.data*: To run the flow in the domain with inlet unsteady conditions from the box domain
- First, edit and read the BuildMeshes.data file.
- If you wish to run a RANS simulation, open the PeriodicBoxRANS.data and DomainFlowRANS.data files.
- Or if you wish to run a LES simulation, open the PeriodicBoxLES.data and DomainFlowLES.data files.

# Turbulent flow in a curved pipe (3D)

## TrioCFD

- Then build the meshes:

**./prepare**

**trust BuildMeshes**

Notice that we use "**trust**" command lines because this script will use the variable \$exec which is the path to the TrioCFD executable.

- You can visualize the partitioned meshes with (MODEL=RANS or LES):

**trust -mesh PeriodicBoxMODEL**

- Then, set the max number of time steps in the time scheme using **nb\_pas\_dt\_max** to 100 in the files PeriodicBoxRANS and PeriodicBoxLES and run a 2-cores parallel calculation, to initialize the turbulent flow in the box:

**trust PeriodicBoxRANS 2**

**trust PeriodicBoxLES 2**

(The full calculation takes approximately 1h in RANS and 10h in LES.)

# Turbulent flow in a curved pipe (3D)

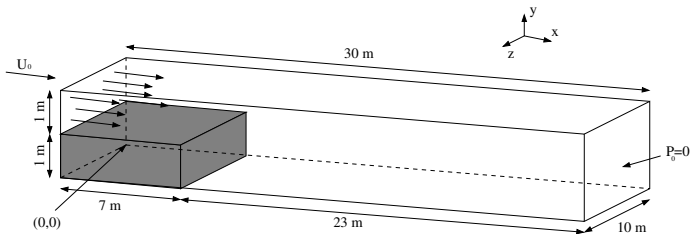
## TrioCFD

- Open the file PeriodicBoxRANS.dt.ev and PeriodicBoxLES.dt.ev to read the last time of the two calculations after 100 time steps.
- Once finished, open the file DomainFlowRANS.data and DomainFlowLES.data and change the maximal number of time steps to 10.
- You can see that these data files are constituted by 2 problems, one for the box and one for the domain.
  - We use the velocity and temperature fields of the last time step of the PeriodicBoxRANS (or LES) calculation as initial conditions for the pb\_box problem with the keyword "Champ\_fonc\_reprise".
  - In addition, the velocity and temperature fields of the pb\_box are used as boundary conditions for the pb\_dom through the keyword "champ\_front\_recyclage".
- Run a 6-cores parallel calculation over the domain (it will stop by default after 10 times steps):  
**trust DomainFlowRANS 6**  
**trust DomainFlowLES 6**

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# Turbulent flow over a backward-facing step

TrioCFD



**Meshing:**  $30 \times 10 \times 10$  ( $\Delta x = 1m, \Delta y = 0.2m, \Delta z = 1m$ )

**Fluid:**  $\mu = 5.10^{-5} kg.m^{-1}.s^{-1}, \rho = 2kg.m^{-3}$

**Boundary conditions:** with in entry  $Re = \frac{U_0 H_{inlet} \rho}{\mu} = \frac{1 \times 1 \times 2}{5.10^{-5}} = 40000$

Inlet:  $U_0 = 1m.s^{-1}$

Outlet:  $P_0 = 0$



# Turbulent flow over a backward-facing step

## TrioCFD

- First initialize TrioCFD environment :  
`source /home/triou/env_TrioCFD_X.Y.Z.sh`  
`echo $exec`  
`echo $project_directory`
- Copy the study named Marche3D: **trust -copy Marche3D**
- Edit the data file and:
  - Note that we use a "Pb\_Hydraulique\_Turbulent" problem with "Navier\_Stokes\_Turbulent" equations and a "modele\_turbulence" model.
  - Modify the fluid characteristics to perform a calculation at  $Re = 50000$ . For example, impose  $\rho = 1 \text{ kg.m}^{-3}$  and  $\mu = 2.10^{-5} \text{ kg.m}^{-1}.s^{-1}$ .
  - Select the sub-grid Smagorinsky turbulence model with standard wall law instead of the "sous\_maille" model (LES).
  - Select the Quick convection scheme.
  - Post-process of velocity, pressure, vorticity, turbulent viscosity at the nodes and elements.

# Turbulent flow over a backward-facing step

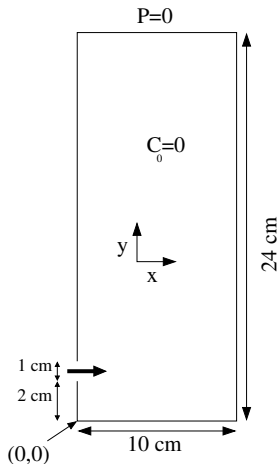
## TrioCFD

- Run the calculation and post-process the main calculated fields.  
**trust Marche3D**
- Notice that we use "**trust**" command lines because this script will use the variable \$exec which is the path to the TrioCFD executable.
- Replace the sub-grid model by the standard k\_eps model (RANS).
- Run the calculation and post-process of the velocity field to see the differences between the different turbulence models used.

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# Tank filling (2D, single-phase flow)

We want to simulate the following flow:



**Fluid:** Colored water  
 diffusion  $D = 10^{-9} m^2.s^{-1}$ ,  
 $\rho = 1000 kg.m^{-3}$ ,  $\mu = 10^{-3} kg.m^{-1}.s^{-1}$

## Boundary conditions:

*Inlet:* Velocity:  $(V_x, V_y) = (V(t), 0)$

$$\text{with } V(t) = \begin{cases} 1 - (y - 0.025/0.005)^2 & , t \leq 0.5s \\ 0 & , t > 0.5s \end{cases}$$

$$\text{Concentration: } C = \begin{cases} 1 & , t \leq 0.5s \\ 0 & , t > 0.5s \end{cases}$$

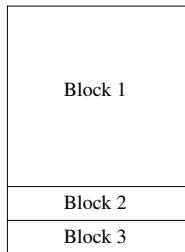
*Outlet:* Pressure  $P = 0$

*Wall:* Velocity  $V_x = 0$ ,  $V_y = 0$

**Initial conditions:** Concentration  $C_0 = 0$ ,  
 Velocity  $V = 0$

# Tank filling (2D, single-phase flow)

- Source the TRUST environment:  
`source /home/triou/env_TRUST_X.Y.Z.sh`
- Copy the study named **diagonale**. This test case deals with a 2D flow with Navier-Stokes and the equation for one constituent.
- Edit the data file and modify the fluid characteristics to the previous ones ( $\mu, \rho, D$ ).
- We want to modify the geometry of this problem to the previous picture. So we want to create 3 blocks like:



# Tank filling (2D, single-phase flow)

- Create the corresponding mesh with 3 blocks (start with  $dx = dy = 0.2cm$  which gives a total nodes number  $N_x = 51$  and  $N_y = 121$ ).
  - Create a first block "Block1" whose origin is (0, 0.03),  $N_x = 51$ ,  $N_y = 106$  (for  $dx = dy = 0.2cm$ ),  $L = 0.1m$ ,  $H = 0.21m$ . Name the wall boundaries Left1, Outlet(=Top1) and Right1. (Don't forget the comma between blocks definitions.)
  - Create the second block "Block2" whose origin is (0, 0.02),  $N_x = 51$ ,  $N_y = 6$  (for  $dx = dy = 0.2cm$ ),  $L = 0.1m$ ,  $H = 0.01m$ . Name the wall boundaries Inlet(=Left2) and Right2.
  - Create the third block "Block3" whose origin is (0, 0),  $N_x = 51$ ,  $N_y = 11$  (for  $dx = dy = 0.2cm$ ),  $L = 0.1m$ ,  $H = 0.02m$ . Name the wall boundaries Left3, Bottom3 and Right3.
- Define the boundary wall, using the keyword "**RegroupeBord**".
- You could also use `facteurs` and `symx`, `symy` keywords to define a refined mesh near the walls.
- Check the mesh with: **trust -mesh diagonale** and correct the mesh errors if necessary.

# Tank filling (2D, single-phase flow)

- In the data file, change the values in the time scheme to stop the calculation at 1 second, and modify **dt\_min** and **dt\_max** values to let TRUST compute time step.
- Change values for the gravity to  $-9.81m.s^{-2}$  following y-axis.
- Note that the **beta\_co** keyword may be useful in order to have a Boussinesq coupling between momentum and concentration equations ( $\beta C_0 g (C - C_0)$  source term added to the Navier-Stokes equations).
- Change the initial and boundary conditions for Navier-Stokes equations:
  - for the Outlet boundary, you have to impose  $P = 0$ ,
  - for the Wall boundary, you have to impose  $V_x = V_y = 0$  with "**paroi\_fixe**" keyword.,
  - for the Inlet boundary, you have to impose  $(V_x, V_y) = (V(t), 0)$  with 
$$V(t) = \begin{cases} 1 - (y - 0.025/0.005)^2 & , t \leq 0.5s \\ 0 & , t > 0.5s \end{cases}$$
 . You will use the **Champ\_Front\_Fonc\_txyz** keyword for the velocity, to write something like: **Champ\_Front\_Fonc\_txyz 2 (1 - ((y - 0.025)/0.005)<sup>2</sup>) \* (t < 0.5) 0**.  
Note: Use  $(t[0.5])$  syntax if you prefer  $(t \leq 0.5)$

# Tank filling (2D, single-phase flow)

- Change the initial and boundary conditions for the constituent equation.
  - You will also use **Champ\_Front\_Fonc\_txyz** field for the Inlet boundary condition for concentration.
  - For the Outlet, use the following keywords to insure the external concentration is 0: **Frontiere\_ouverte C\_ext Champ\_front\_uniforme 1 0**.
  - For the Wall, the keyword for impermeable boundary condition for concentration is **paroi**.
- Check you have high-order schemes (i.e. "**Quick**" scheme) used in both equations to reduce numerical diffusion.
- Notice you could have suppressed diffusion term in concentration equation rather than using a small diffusion coefficient with:  
**Diffusion { negligeable }**
- Add a concentration probe near the inlet (e.g.: at (0,0.025)).
- Add a velocity segment probe (with 5 points between (0,0.021) and (0,0.029)) at the inlet boundary to see the time evolution of these two quantities (period 0.01s).



# Tank filling (2D, single-phase flow)

- Run the study and follow the time evolution with the probes:  
**trust -evol diagonale &**  
"Start computation!" button and "Plot" or "Plot on same" for probes.
- Check the flow rate in inlet boundary in the diagonale\_pb\_Debit.out file (plotted on the right of the PLOT2D window). You should find a value near  $6.8 \cdot 10^{-3} m^2.s^{-1}$ .
- Use VisIt to post-process the results at  $t=0.2$ ,  $t=0.4s$  and  $t=0.7s$ . VisIt has some interesting feature for this study. It can give concentration histogram to check the numerical diffusion in the concentration equation: Add → Histogram → CONCENTRATION\_ELEM\_dom. The volume of colored water (in  $m^3$ ) is given by  $Vol(t) = 6.66 \cdot 10^{-3} t$  before  $t = 0.5s$  and  $Vol(t) = 3.33 \cdot 10^{-3}$  after.

# Tank filling (2D, single-phase flow)

## → VEF

- Copy diagonale.data to diagonale\_VEF.data.
- Triangulate your mesh (**triangular** keyword).
- In this new file, change the discretization (**VEFPreP1B** instead of **VDF**).
- Use **muscl** instead of **quick** scheme.
- And you can switch **GCP** solver by **Cholesky** solver of the Petsc library (direct method which may need large amount of RAM memory) to increase the speed resolution of the pressure linear system:  
**GCP { precondition ssor { omega 1.5 } seuil 1.e-6 } → Petsc Cholesky { }**
- Run the calculation. You must have an error, and TRUST stop the calculation.

# Tank filling (2D, single-phase flow)

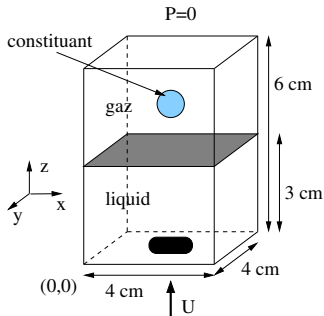
## → VEF

- As TRUST indicates, to avoid this problem, you can:
  - change the **trianguler** keyword to **trianguler\_h**,
  - or use the **VerifierCoin** keyword. For this, after this first error you must find a "diagonale\_VEF.decoupage\_som" file in your directory, so you can use it by adding:  
**VerifierCoin dom { read\_file diagonale\_VEF.decoupage\_som }**  
 just after "trianguler dom". This will subdivides inconsistent 2D/3D cells used with VEFPreP1B discretization (cf Reference Manual).
- Run the calculation and compare the results between **VDF/quick** and **VEFPreP1B/muscl** which must take much more time!

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# Tank filling (3D, two-phase flow)

TrioCFD



**Liquid:**  $\rho = 1000 \text{ kg.m}^{-3}$ ,  
 $\mu = 2,82.10^{-4} \text{ kg.m}^{-1}.\text{s}^{-1}$ ,  
 $\sigma = 0.05 \text{ N.m}^{-1}$ ,  $D = 10^{-6} \text{ m}^2.\text{s}^{-1}$

**Gas:**  $\rho = 100 \text{ kg.m}^{-3}$ ,  
 $\mu = 2,82.10^{-4} \text{ kg.m}^{-1}.\text{s}^{-1}$

**Boundary conditions:**

Up : Free outlet, Wall :  $V = 0$   
 Down:  $V(x, y, z) = (0, 0, 10^{-3} \text{ m.s}^{-1})$

**Initial conditions:**  $V = 0$ ,  
 $C = e^{(-((x-0.02)^2 + (y-0.02)^2 + (z-0.03)^2)/0.03^2)}$

N.B.: The interface between the air and the gas is a parabolic function.

# Tank filling (3D, two-phase flow)

## TrioCFD

- First initialize TrioCFD environment :  
`source /home/triou/env_TrioCFD_X.Y.Z.sh`  
`echo $exec`
- Copy the study named **FTD\_all\_VDF**: `trust -copy FTD_all_VDF`
- This test case deals with a 3D two-phase flow in a tank with one initial interface between liquid and gas, a droplet, and a rotating solid in the liquid. The Discontinuous Front Tracking method is used with a 3D structured mesh.
- Notice that:
  - 2D Discontinuous Front Tracking method has not been intensively tested yet.
  - the type of the problem: **Probleme\_FT\_disc\_gen** in the data file. This refers to the Discontinuous Front Tracking method.
  - the keyword **modele\_turbulence**. Navier-Stokes equations of the Discontinuous Front Tracking problem needs the read of this keyword even if the flow is laminar. In this case, use the **nul** keyword just after **modele\_turbulence**. Else, specify the turbulence model to use.

# Tank filling (3D, two-phase flow)

## TrioCFD

- Increase the height of the tank (from 0.06 to 0.12).
- Add a second drop above the first one, at  $z = 0.08$  (keywords **ajout\_phase0** could be useful to add other interfaces, cf Reference Manual for **ajout\_phase0/ajout\_phase1** keywords). Don't forget the comma between the two definition of the drops.
- Change the **dt\_post** period of the 3 post-processing blocks (0.05 to 0.01). The first one (add **format lata**) is the classical block for post-processing probes and fields. Here, we want to see the concentration field and the "**indicatrice\_interf**" field. Value of this field is 0 for liquid and 1 for gas, so the interface is located at "indicatrice" value 0.5.
- Change the interpolation location of **indicatrice\_interf** and the **concentration** fields in the first post-processing block, by adding the keyword **elem** just after the fields: the values in the post-processing tool will be plotted at the center of each element of the mesh.

# Tank filling (3D, two-phase flow)

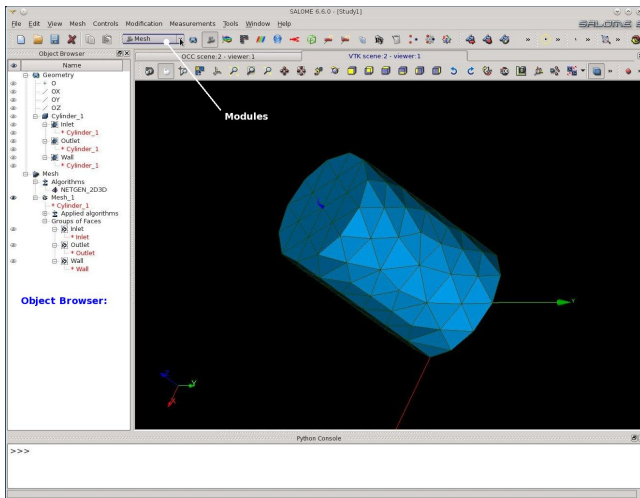
## TrioCFD

- The second post-processing block is the new syntax to post-process interfaces moving meshes. You can visualize it with **VisIt**.
- On each interface, you can plot several fields i.e.: curvature with **courbure** keyword and velocity interface with **vitesse** keyword, **pe** field is for debugging purpose, it is useless here, you can suppress it) on several locations (on nodes with **sommets** keyword, on cells with **elements** keyword).
- Run the calculation. Follow the time step evolution by having a look at the **dt\_ev** file. It contains on each line the physical time, the time step, security factor and residuals.
- Post-process to visualize the interface and the concentration field.
- You can increase the number of cells to have a finest simulation or also change to VEF discretization.



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## Salomé to create a 3D VEF mesh: Cylinder



# Cylinder

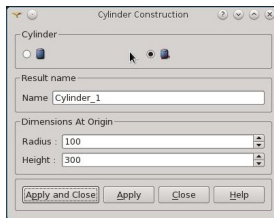
## Create a geometry

- Create a new folder:  
\$ mkdir -p Formation\_TRUST/yourname/salome/exo1  
\$ cd Formation\_TRUST/yourname/salome/exo1
- Launch Salomé (we suppose it is installed in \$PathToSalome):  
\$ \$PathToSalome/salome &
- Create a new study: File → New
- Select the Geometry module into the SALOME drop-down menu (contains all the modules).
- Save your study in hdf format (Salome format) frequently.
- Create a first geometry with: New Entity → Primitives → Cylinder

# Cylinder

## Create a geometry

- Specify Radius  $R = 100$  and Height  $H = 300$  for the cylinder (the default values). Then Apply and Close.

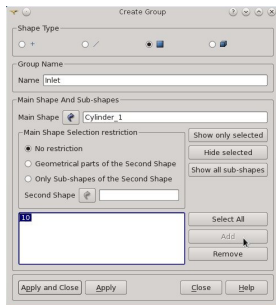


- Rotate, zoom, move the geometry by switching to "Interaction style switch": Mouse icon.
- Create groups for the geometry to define the top, the bottom and the lateral parts of the cylinder: New Entity → Group → Create Group
- Select the good Shape Type (→ □ surface).

# Cylinder

## Create a geometry

- Give a Group Name for the top: "Inlet".
- Click on the arrow button of the Main Shape field and select the "Cylinder\_1" in the "Object browser" or in the visualization window.
- Select the shape defining the top of the cylinder on the visualization window then Add → Apply.
- Select the shape defining the part on the window then Add → Apply.



# Cylinder

## Create a geometry

- Do the same for the two other parts:
  - For the lateral: "Wall"
  - For the bottom: "Outlet" (you can rotate the cylinder to click on the bottom).
- Close the window once the 3 groups has been created. Check that they appear in the Object Browser (by clicking on the "▷" in front of the "Cylinder\_1" object).

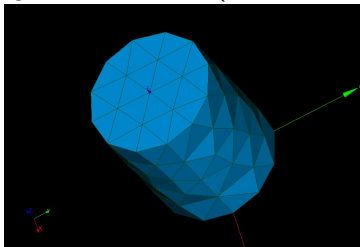
## Create a mesh

- Now, switch to the Mesh module in the SALOME drop-down menu.
- Select the Cylinder\_1 in the Object Browser and Right Click → 'Show' to visualize the geometry or click on the 'eye' next to the Cylindre\_1 object.
- Create a mesh with: Mesh → Create Mesh
- Select the Geometry used for the mesh if not selected by clicking on the Cylinder\_1 object in the Object Browser.

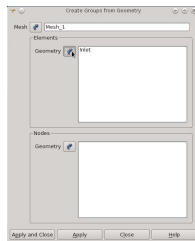
# Cylinder

## Create a mesh

- Choose Netgen 1D-2D-3D algorithm and click on "Apply and Close".
- Select the object Mesh\_1 in the Object Browser and Right Click → Compute (or Mesh → Compute).
- A tabular must appear with the number of triangles, quadrangles... Click on "Close".
- Hide the geometry by selecting the Cylinder\_1 in the Object Browser and Right Click → Hide (or click on the eye).



Mesh



Create Groups from Geometry

# Cylinder

## Export your mesh in MED format

- Now, we are going to create groups for the mesh, by using the geometry groups: Mesh → Create Groups from Geometry.
- Click on "Mesh\_1" object in the Object Browser to select it.
- Then push the arrow button of Elements/Geometry.
- Select together the 3 groups (Inlet, Wall and Outlet) with "Ctrl" +left click in the Cylinder\_1 object of the Object Browser, then press Apply and Close.
- Check that the 3 boundaries are in the "Group of Faces" of the Mesh\_1 object in the Object Browser.
- Export your mesh with the MED format :  
Select the Mesh\_1 object then Right Click → Export → MED file (or File → Export → MED file).



# Cylinder

## Read your mesh with TRUST

- Now build a data file named dom.data for TRUST:

```
dimension 3
domaine dom
Read_med family_names_from_group_names dom Mesh_1 Mesh_1.med
Postraiter_domaine { domaine dom fichier mesh format lata }
```

- Run the data file and post-process the mesh with VisIt:

```
source /home/triou/env_TRUST_X.Y.Z.sh
trust dom
visit -o mesh.lata
```

**Warning:** The more common error is to forget to define the boundaries with the groups for the mesh. The error in TRUST is printed and detected during the discretization where all the faces of the mesh (in particular the boundary faces) are built.

# Cylinder

## Refine your mesh and use viscous layers

Goal: Improve the mesh for TRUST near the wall by using viscous layers.

- Create a new mesh named "Refined\_mesh" with: Mesh → Create Mesh
- Select the Cylinder\_1 geometry in the Object Browser.
- Select the "Tetrahedron (Netgen)" or "MG-Tetra" 3D algorithm.
- Click on the wheel of "Add. Hypothesis" → "Viscous Layers" with:
  - Total thickness: 30
  - Number of layers: 3
  - Stretch factor: 1.1
  - Add to "Faces without layers" the 2 geometry groups "Inlet" and "Outlet" of Cylinder\_1 object (select or unselect the mouse icon).
  - Click OK
- Add a 2D algorithm: "Netgen 1D-2D" or "MG-CADSurf".

# Cylinder

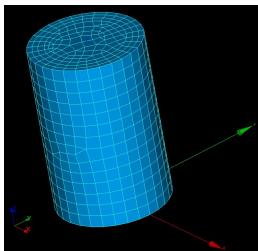
## Refine your mesh and use viscous layers

- Click on the wheel of "Hypothesis" → "Netgen 2D parameters" or "MG-CADSurf parameters":
  - For "Netgen 2D parameters":
    - Change "Fineness" from "Moderate" to "Very Fine".
    - Select "Allow Quadrangles".
    - Click OK.
  - For "MG-CADSurf parameters", change "User size" to 20. Click OK.
- "Apply and Close" the close mesh window.
- Select the Refined\_Mesh object in the Object Browser and Right click → Compute

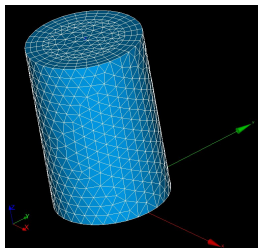
# Cylinder

## Refine your mesh and use viscous layers

- You should have a refined mesh with a mix of tetra, hexa, pyramid, prism elements for Netgen algorithms, and a mix of tetra and prisms for MG algorithms:



Netgen



MG

- As TRUST accepted only tetras elements, you can quickly tetraedrise:
  - Select the Refined\_Mesh in the Object Browser.
  - "Modification" → "Split Volumes" and select "Tetrahedron".
  - Don't change the parameters, and click "Apply and Close".

# Cylinder

## Refine your mesh and use viscous layers

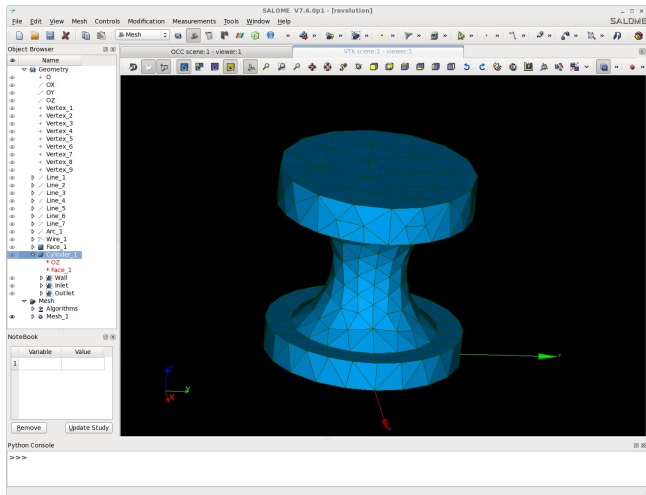
- As usual, define the mesh boundaries with Groups:
  - Select the Refined\_Mesh, Right click → Create Groups from Geometry.
  - Select "Inlet", "Outlet", "Wall" with "Ctrl"+click in the Cylinder\_1 object, then Apply and Close.
- Export the mesh:
  - Select the Refined\_mesh, Right click → Export → MED file.
  - Save into a Refined\_Mesh.med file.
- Save your work in hdf format ("File" → "Save/Save As..."), and in python format with "File" → "Dump Study..."

## Run with TRUST

- Edit your datafile or create a new one to read and visualize your refined mesh.
- **N.B.:** The solutions of the exercise (mesh.py file for the first mesh and prism.py file for the second mesh) are located here:  
\$TRUST\_ROOT/doc/TRUST/exercices/salome.

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# Salomé to create a 3D VEF mesh: Revolution



# Salomé to create a 3D VEF mesh: Revolution

## Create a geometry

- Create a directory and run Salomé (we suppose it is installed on \$PathToSalome):

```
$ mkdir -p Formation_TRUST/yourname/salome/exo2
```

```
$ cd Formation_TRUST/yourname/salome/exo2
```

```
$ $PathToSalome/salome &
```

- Create a new study: File → New.
- Select the Geometry module into the SALOME drop-down menu.

- Create points : New Entity → Basic → Point

Vertex_1 (0,0,0)	Vertex_2 (1,0,0)	Vertex_3 (1,0,0.3)
------------------	------------------	--------------------

Vertex_4 (0.75,0,0.3)	Vertex_5 (0.375,0,1)	Vertex_6 (0.75,0,1.6)
-----------------------	----------------------	-----------------------

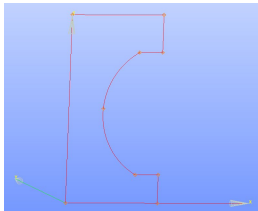
Vertex_7 (1,0,1.6)	Vertex_8 (1,0,2)	Vertex_9 (0,0,2)
--------------------	------------------	------------------

Then "Apply and Close"



# Revolution

## Create a geometry



- Create edges:
  - New Entity → Basic → Line
    - ◇ Line\_1 with Vertex\_1 and Vertex\_2
    - ◇ Line\_2 with Vertex\_2 and Vertex\_3
    - ◇ Line\_3 with Vertex\_3 and Vertex\_4
    - ◇ Line\_4 with **Vertex\_6** and **Vertex\_7**
    - ◇ Line\_5 with Vertex\_7 and Vertex\_8
    - ◇ Line\_6 with Vertex\_8 and Vertex\_9
    - ◇ Line\_7 with Vertex\_9 and Vertex\_1
    - ◇ Then Apply and Close.

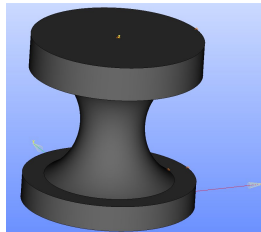
# Revolution

## Create a geometry

- Create edges:
  - New Entity → Basic → Arc
    - ◊ Arc\_1 with Vertex\_4, Vertex\_5 and Vertex\_6.
    - ◊ Then Apply and Close.
- Create a wire: New Entity → Build → Wire
  - Wire\_1 on edges with Line\_1,... , Line\_7 and Arc\_1 (with "Ctrl" button).
  - Then Apply and Close.
- Create a face: New Entity → Build → Face.
  - Face\_1 with Wire\_1 and "Apply and Close".
- Create a revolution cylinder: New Entity → Generation → Revolution.
  - named Cylinder\_1,
  - with Face\_1 in Objects,
  - click on the arrow button next "Axis" and select OZ in the Object Browser,
  - set the angle to  $360^\circ$  and "Apply and Close".

# Revolution

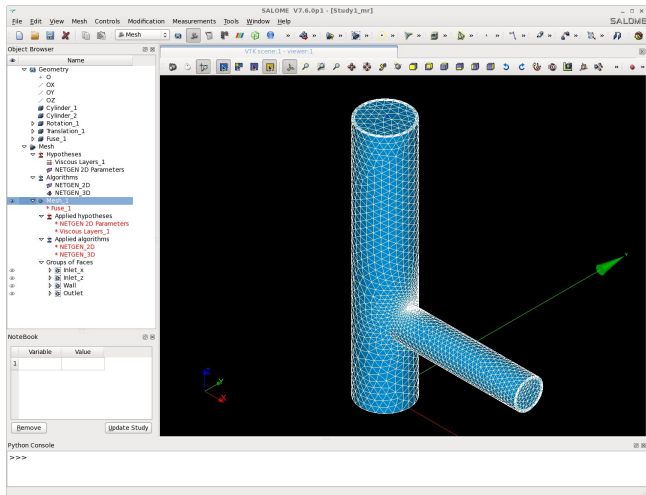
## Create a geometry



- Create groups for the geometry to define the top, the bottom and the lateral parts of the cylinder: New Entity → Group → Create Group.
- Save your study in hdf format ("File" → "Save/Save As..."), and in python format with "File" → "Dump Study..."
- Now you can create the mesh in the same way than page 99.
- **N.B.:** You can find the solutions of this exercise (revolution.py) in \$TRUST\_ROOT/doc/TRUST/exercices/salome.

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# Salomé to create a 3D VEF mesh: T-shape



# Salomé to create a 3D VEF mesh: T-shape

## Create a geometry

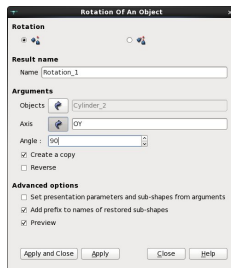
- Create a directory and run Salomé (we suppose it is installed on \$PathToSalome):  

```
$ mkdir -p Formation_TRUST/yourname/salome/exo3  
$ cd Formation_TRUST/yourname/salome/exo3  
$ $PathToSalome/salome &
```
- Create a new study: File → New.
- Select the Geometry module into the SALOME drop-down menu.
- Create two cylinders: New Entity → Primitives → Cylinders:  
Cylinder\_1: radius 0.5, height 5. Then "Apply".  
Cylinder\_2: radius 0.3, height 3. Then "Apply and Close".
- Save your study in hdf format (Salome format) frequently.

# T-shape

## Create a geometry

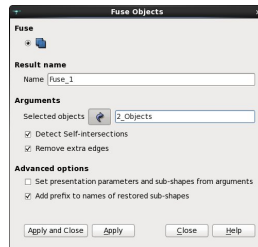
- Rotate Cylinder\_2: Operations → Transformation → Rotation  
Name: Rotation\_1, Object: Cylinder\_2, Axis: 'OY', Angle: 90°  
Then "Apply and Close".
- Translate Rotation\_1: Operations → Transformation → Translation  
Name: Translation\_1, Object: Rotation\_1, Dx=Dy=0, Dz=1.5  
Then "Apply and Close".



Rotation



Translation

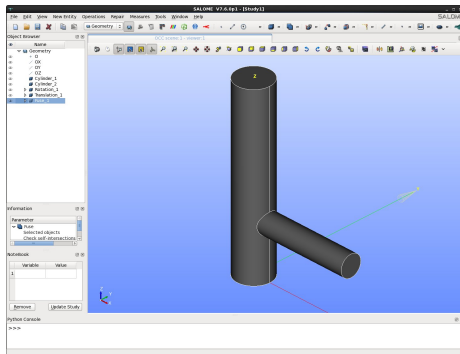


Fuse

# T-shape

## Create a geometry

- Fuse Cylinder\_1 and Translation\_1: Operations → Boolean → Fuse  
Name: Fuse\_1, Selected Objects: 2\_Objects (use "Ctrl" button to select Cylinder\_1 and Translation\_1 in the Object Browser).  
Then "Apply and Close".

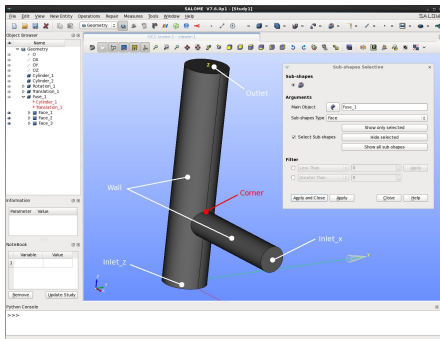




# T-shape

## Create a geometry

- We are now going to create the boundaries: New Entity → Explode
  - Main Object: Fuse\_1, Sub-shape type: Face, select "Select sub-shape" and click on the surface Outlet and "Apply".
  - This will create a face named "Face\_1" in the Fuse\_1 object (click on the "►"), rename it "Outlet" (by right-clicking and "rename").



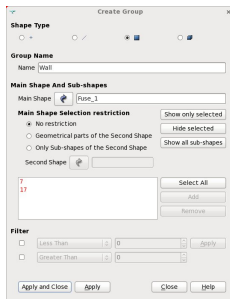
# T-shape

## Create a geometry

- Do the same for "Inlet\_x" and "Inlet\_z".
- We are now going to create the boundary Wall: New Entity → Group → Create group:

Shape Type: surface, Name: Wall, Main Shape: Fuse\_1.

Click on the surface of the Cylinder\_1 then "Add", click on the surface of Translation\_1 then "Add" and "Apply and Close".



# T-shape

## Create a geometry

- We are now going to create the point "Corner": New Entity → Explode
  - Main Object: Fuse\_1, Sub-shape type: Vertex, select "Select sub-shape" and click on the chosen point and "Apply and Close".
  - This will create a vertex name "Vertex\_1", rename it "Corner" (by right-clicking and "rename").

## Create a mesh

- Now, switch to the Mesh module in the SALOME drop-down menu.
- Select the Fuse\_1 in the Object Browser and Right Click → 'Show' to visualize the geometry or click on the 'eye' next to the Fuse\_1 object.
- Create a mesh with: Mesh → Create Mesh.
- Select the Geometry used for the mesh if not selected by clicking on the Fuse\_1 object in the Object Browser.

# T-shape

## Create a mesh

- Choose "Tetrahedron (Netgen)" for 3D algorithm.
- Click on the wheel of "Add. Hypothesis" → "Viscous Layers" and set:
  - Total thickness: 0.05
  - Number of layers: 3
  - Stretch factor: 1.1
  - Extrusion method: Node Offset
  - Add to "Faces with layers (Wall)" the geometry group "Wall" of Fuse\_1 object in the Object Browser (select or unselect the mouse icon). Click on "Add".
  - Click "OK".
- Choose "Netgen 1D-2D" for 2D algorithm.
- Click on the wheel of "Hypothesis" → "Netgen 2D parameters" and set for "Arguments" menu:
  - Max. Size: 0.6
  - Min. Size: 0
  - without "Second Order"
  - Finess: Custom

# T-shape

## Create a mesh

- Growth rate: 0.1
- Nb. segs per Edge: 2
- Nb. segs per Radius: 4
- Select "Limit size by Surface Curvature", "Optimize", "Fuse Coincident Nodes on Edges and Vertices".
- Unselect "Allow Quadrangles".
- For "Local Size" menu:
  - Select "Corner" object in the Object Browser and click on "On Vertex" in the "Hypothesis Construction" window.
  - Double-click on the value in the table and set it to "0.01".
  - Click "OK".
- Click on "Apply and Close".
- Select the Mesh\_1 object in the Object Browser and Right click → Compute.
- You should have a mesh with a mix of tetra and prism elements.

# T-shape

## Create a mesh

- As TRUST accepted only tetras elements, you can quickly tetraedrise:
  - Select Mesh\_1 in the Object Browser.
  - "Modification" → "Split Volumes" and select "Tetrahedron".
  - Don't change the parameters, and click "Apply and Close".
- Define the mesh boundaries with Groups:
  - Select the Mesh\_1, Right click → Create Groups from Geometry.
  - Select "Inlet\_x", "Inlet\_z", "Outlet" and "Wall" with "Ctrl"+click in the Fuse\_1 object.
  - Then Apply and Close.
- Export the mesh:
  - Select the Mesh\_1, Right click → Export → MED file.
  - Save into a Mesh\_1.med file.
- Save your study in hdf format ("File" → "Save/Save As..."), and in python format with "File" → "Dump Study..."
- **N.B.:** You can find the solutions of this exercise (T\_shape.py) in \$TRUST\_ROOT/doc/TRUST/exercices/salome.

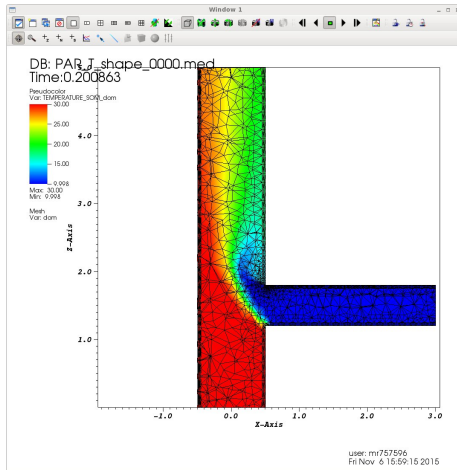
# T-shape

## Run with TRUST

- Copy the T\_shape.data file in your directory:  
**cp \$TRUST\_ROOT/doc/TRUST/exercices/salome/T\_shape.data .**
- Run it with TRUST:  
**trust T\_shape**  
or in parallel with:  
**trust -partition T\_shape**  
**trust PAR\_T\_shape 4**
- You can visualize the results with Visit or Salomé by opening the T\_shape\_0000.med file for sequential calculation or PAR\_T\_shape\_0000.med for parallel calculation.

# T-shape

## Visu with Visit





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# Salomé: Create domains for a TRUST coupled problem

Consider that we want to simulate a coupled problem with TRUST on a complex geometry. Suppose that this latter is drawn by means of Salomé.

The main difficulty arises from the fact that the mesh elements should be connected on the interface between the two domains in order to be correctly read by TRUST.

In this exercise, you will learn how to:

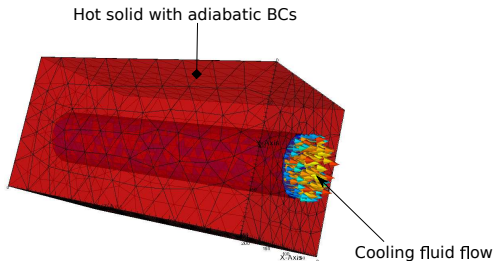
- Create two domains (domain 1 and domain 2) using Salomé
- Get a coherent mesh on the interface between the two domains. We recall that TRUST is able to treat only meshes with connected elements on the interface.
- Mesh both domains and export it into a single MED file.
- Read the MED file from TRUST datafile and simulate a coupled problem.

Note: for simple geometries, the internal TRUST mesher "Mailler" will be largely sufficient (see the exercise [3](#) for example).

# Mesh for coupled problem

## Description of the problem

Let us consider the cooling of a solid block by means of a fluid flowing inside circular cross-section channels. The channel is centered in the block of a square cross-section. The outer boundaries of the solid are adiabatic. Below is given a schematic description of the problem.



In order to build meshes using Salomé for such a simulation, we should create two domains: the first domain will represent the solid block and the second domain the fluid.

# Mesh for coupled problem

## In the Geometry module:

- Create a new folder for this exercise and launch Salomé:  

```
$ mkdir -p Formation_TRUST/yourname/salome/exo4  
$ cd Formation_TRUST/yourname/salome/exo4  
$ $PathToSalome/salome &
```
- Create a new study: File → New
- Select the Geometry module from drop-down menu of Salomé.
- Save your study in hdf format (Salomé format) frequently.
- Create the first geometry with: New Entity → Primitives → Box. Then, specify dimensions  $D_x = 200$ ,  $D_y = 200$  and  $D_z = 400$ . After that, Apply and Close.
- Create a vertex with: New Entity → Basic → Point. Specify the vertex coordinates  $X = 100$ ,  $Y = 100$  and  $Z = 0$ , then Apply and Close
- Create the second geometry with: New Entity → Primitives → Cylinder. Then, specify the Base Point: Vertex\_1 and Vector OZ, Radius  $R = 40$  and Height  $H = 400$  for the cylinder. After that, Apply and Close.

# Mesh for coupled problem

## In the Geometry module:

- Perform a cut with: Operations → Boolean → Cut. In the Main Object select: Box\_1 and in the Tool Objects select: Cylinder\_1 → Apply and Close.
- Create a partition with: Operations → Partition. In Objects select: Cylinder\_1 and Cut\_1. Then Apply and Close.
- Define 2 groups of volumes, one for each domain with: New entity → Group → Create Group.
  - 1 Shape Type: Volume. Name: Solid. Main Shape: Partition\_1. Select the hollow box then click on Add, after that on Apply.
  - 2 Shape Type: Volume. Name: Fluid. Main Shape: Partition\_1. Select the cylindrical channel then click on Add, after that on Apply and Close.

# Mesh for coupled problem

## In the Geometry module:

- Define the groups of faces for external boundaries and the interface with:  
New entity → Group → Create Group.
  - ① Shape Type: Surface. Name: Fluid\_inlet. Main Shape: Partition\_1. Then select the bottom of the cylinder and Add. Click on Apply.
  - ② Shape Type: Surface. Name: Fluid\_outlet. Main Shape: Partition\_1. Then select the top circular boundary of the cylinder then click on Add, then Apply.
  - ③ Shape Type: Surface. Name: Solid\_top. Main Shape: Partition\_1. Then select the top of the box then click on Add, then Apply.
  - ④ Shape Type: Surface. Name: Solid\_bottom. Main Shape: Partition\_1. Then select the bottom of the box then click on Add, then Apply.
  - ⑤ Shape Type: Surface. Name: Solid\_lateral\_walls. Main Shape: Partition\_1. Then select the remaining 4 lateral boundaries of the box then click on Add, then Apply.
  - ⑥ Shape Type: Surface. Name: Solid\_Fluid\_Interface. Main Shape: Partition\_1. Then select the top boundary of the box and click on Hide selected, then Click on a lateral boundary and click on Hide selected, then the lateral boundary of the cylinder will be visible. Select it and click on Add then Apply.

# Mesh for coupled problem

## In the Mesh module:

- Create a mesh based on the Partition\_1 with: Mesh → Create Mesh. Let the name be Mesh\_1 and in Geometry select Partition\_1. In the 3D algorithm, select NETGEN 1D-2D-3D. Click on the wheel of "Hypothesis" then on "NETGEN 3D Parameters". In Arguments, select the fineness "Fine" instead of "Moderate" then click on OK then Apply and Close.
- Right click on Mesh\_1, then Compute.
- Create groups with: Mesh → Create Groups from Geometry. In Mesh, select Mesh\_1 then in Elements select all the created groups in Partition\_1 (with keypad Ctrl) in Geometry then click on Apply And Close.
- Export the mesh in med format (if possible, choose MED 3.2).

# Mesh for coupled problem

## Launch the coupled problem datafile:

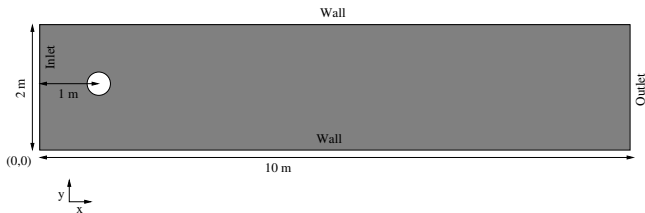
- Source TRUST environnement with:  
`source /home/triou/env_TRUST_X.Y.Z.sh`
- Copy the datafile:  
`$ cp $TRUST_ROOT/doc/TRUST/exercices/salome/Coupled_pb.data .`
- Run the test case using TRUST:  
`$ trust Coupled_pb.data`
- When the computation finishes, visualize results using VisIt:  
`$ visit -o Coupled_pb.lata`
- Draw the temperature profile on both domains and set the min and max on color bar to 300 and 400 respectively. When you visualize time evolution of temperature, you see that the solid is cooled and its temperature decreases. If we increase the time of the simulation, the temperature of the solid will be equal to that of the fluid at the steady state.



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# Gmsh to create a 2D VEF mesh

Geometry which will be created, based on a TrioCFD validation test case geometry



- Create a directory and copy an example:  

```
mkdir -p Formation_TRUST/yourname/gmsh
cd Formation_TRUST/yourname/gmsh
source /home/triou/env_TrioCFD_X.Y.Z.sh
echo $project_directory
dir=$project_directory/validation/share/Validation/Rapports_automatiques/Validant
cp $dir/pas_fini/Drag/src/shape.geo file.geo
nedit file.geo &
gmsh file.geo &
```

# Gmsh to create a 2D VEF mesh

- First configure gmsh to show points, lines, and surface numbers of the geometry. In menu Tools → Options → Geometry → Visibility, select Lines, Surfaces, Point labels, Line labels, Surface labels, Volume labels close this window.
- Save definitively your choices with File → Save Options As default.
- Now, look at the file.geo file, you can see the definition of parameters, points, lines... You can see the position of the points and lines with theirs numbers in gmsh.
- Modify the file to suppress the obstacle. We want to keep only 4 points and 4 lines, so you have to suppress the points 5,6,7,8 and the lines 1,3,4,5. For this, you have to:
  - comment (with "//") the definition of the points 5,6,7 and 8.
  - comment the definition of the lines 1,3,4,5. Note that the "Circle" is a line so "Circle(1)=Line(1)".
  - modify the line(2), it will now links points 1 and 2.
  - comment the "Physical Line" named "Shape" which use the line(1) (= Circle(1)) and the line(3).

# Gmsh to create a 2D VEF mesh

- suppress the numbers 4 and 5 in the "Physical Line" "Axis". It refers to the lines 4 and 5 which does not exist anymore.
- suppress the numbers 1,3,4 and 5 in the "Line Loop(1)".
- set H to 2 and L to 10. (You can comment D, E, param and X definitions.)
- Press "Reload" in the gmsh GUI → Geometry to update the geometry visualization.
- Now we will add the circle. Note that you can only create circle arcs with angle strictly smaller than  $\pi$ !
  - create the middle of the circle like "Point(10)={1,1,0,lc};" where the triplet "1,1,0" are the coordinates of the point and "lc" the thickness of the cells next this point.
  - create 4 points around this point, which will correspond to the 4 arc of the circle:  
Point(11)={1.25,1,0,lc2};  
Point(12)={1,1.25,0,lc2};  
Point(13)={0.75,1,0,lc2};  
Point(14)={1,0.75,0,lc2};

# Gmsh to create a 2D VEF mesh

- define the 4 arcs of the circle with this points:  
Circle(10)={11,10,12};  
Circle(11)={12,10,13};  
Circle(12)={13,10,14};  
Circle(13)={14,10,11}; "
- create a "Physical Line" for the circle:  
"Physical Line("Circle") = {10,11,12,13};"  
This name will be used in your TRUST data file as the name of your boundaries.
- create a line loop for the circle just after the first line loop:  
Line Loop(2) = {10,11,12,13};
- Add the number of this line loop in the "Plane Surface(1)":  
Plane Surface(1) = {1,2};
- Suppress the Physical lines "Axis" and "Top" and create a physical line named "Wall" which regroups the top and the bottom of this geometry (lines 2 and 7).

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# Gmsh to create a 3D VEF mesh

- Select "Mesh" in the drop-down menu of gmsh and mesh in 2D.
- Export it to a MED file: "File" → "Save As..." and name the file file.med. (Keep the default options.) You can verify your mesh by opening it with gmsh: **gmsh file.med &**.
- Build a TRUST data file with the **Postraiter\_domaine** keyword, to read the mesh (like in the Salome exercise page 81). Visualize the mesh with VisIt.
- Then we will try to create a 3D mesh, by using the Extrusions feature of Gmsh. See more about extrusions in <http://geuz.org/gmsh/doc/texinfo/gmsh.html>.
- Save your initial file in a new one named file3D.geo.
- Comment your "Physical lines", they will not be used here.
- Add the line "Extrude {0,0,1} { Surface{1} ; }" just before the definition of the physical surface.

# Gmsh to create a 3D VEF mesh

- Comment the line "Physical Surface("domain") = {1};" in 3D we will have a "Physical Volume" which will be define at the end of the .geo file.
- Define the "Physical Surface" which will be the boundaries of your geometry with the number of the surfaces which can be read on the geometry plotted by gmsh:  
 $\text{Physical Surface("Inlet")} = \{38\};$   
 $\text{Physical Surface("Outlet")} = \{30\};$   
 $\text{Physical Surface("Wall")} = \{1,26,34,55\};$   
 $\text{Physical Surface("Obstacle")} = \{42,46,50,54\};$
- Define you physical volume, you can see its number in yellow in the window:  
 $\text{Physical Volume("dom")} = \{1\};$
- Select the "Mesh" tool in the drop-down menu pf gmsh and mesh in 3D your geometry. It takes a few minutes, to reduce this time, increase the size of your cells by changing the values of lc.



# Gmsh to create a 3D VEF mesh

- You can use the "Optimize 3D" algorithm to optimize your mesh.
- Export your mesh to a MED file. Run gmsh again on this exported MED file to check everything is defined:

**gmsh file.med &**

- Now, use your mesh in a TRUST calculation, for example:
  - copy the data file of the first exercise into a Obstacle\_VEF.data file,
  - read the MED file:  
**Lire\_med family\_names\_from\_group\_names dom file file.med**  
Notice that by default with Gmsh, the mesh name is the name of the file!
  - change the discretization type,
  - be careful to the choice of the convection scheme for your VEF calculation,
  - and run the simulation on the unstructured mesh.

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# TRUST Wizzard for 3D VEF mesh

- Open the TRUST wizzard:  
**trust -wiz**
- Create an hydraulic problem with:
  - in "Turbulent model" tab, set a laminar turbulent model.
  - in "Domain" tab, input the path to your 3D mesh file in med format. The wizzard will find the name of your mesh.
  - in "Boundary condition" tab, select a boundary condition of type:
    - "Input" for your Inlet boundary and set your velocity field to ( 1. 0. 0. ),
    - "Wall" with no velocity for your Obstacle boundary,
    - "Output" with pressure field set to 0. for your Outlet boundary,
    - "Symetry" for your Wall boundary.
  - in "Initial condition" tab, set your velocity field to ( 0. 0. 0. ).
  - in "Incompressible flow" tab, choose a predefined field like "Water 20 deg".
- Then click on "Validate" to see if your parameters are ok.
- You can see the message "All ok!" at the bottom left of the window.
- You can save your session at the .JSoN format by clicking on "Save ...".

# TRUST Wizzard for 3D VEF mesh

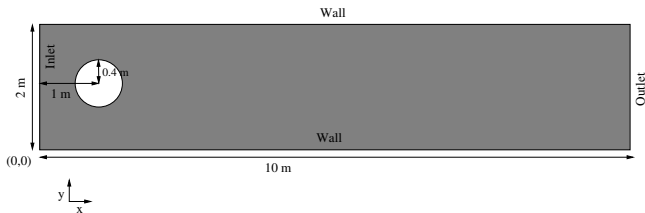
- Click on "Start ...", a window opens to save your data file. Name it for example "test\_wiz.data" and click on "Save".
- The TRUST PLOT2D tool opens, click on "Start computation!".
- An error occurs! Open your data file with "Edit data" button.
- Correct the expression of gravity field by adding a third component.
- Add a "**tmax**" value (for example 'tmax 0.1') and suppress the max number of iterations.
- Save it and run your simulation.
- It must be long because of the size of the cells. You can decrease it and create another mesh with gmsh to run faster.

**Warning:** TRUST Wizzard is still in development. Models are not yet all available.

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# Xprepro to create a 3D VDF simple mesh

## First exercise



In this exercise, you will learn how to create a 2D mesh with Xprepro. Note that in Xprepro, we dig a geometry in an initial block of matter.

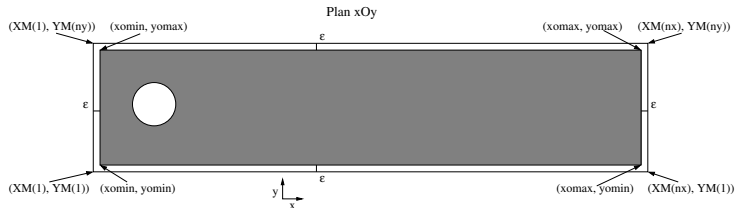
- Create a directory:

```
mkdir -p Formation_TRUST/yourname/Xprepro/exo1
cd Formation_TRUST/yourname/Xprepro/exo1
source /home/triou/env_TRUST_X.Y.Z.sh
Xprepro &
```

# Xprepro to create a 3D VDF simple mesh

- You can see two windows:
  - the command window: `xprepro.tcl` and
  - the list of existing objects: `viewlist` (empty for the moment).
- Click on "Default" button to begin from a cube. Read the pop-up window and click on "Ok". A new nedit window appears, opening a file named "maillagedefault" in which we will set the values of the parameters used in our geometry.
- We want to create a cube of length  $L=10\text{m}$ , height  $H=1\text{m}$  and width  $I=1\text{m}$  with a cylinder of radius  $R=0.4\text{m}$  at one meter from the left side of the cube. So the center of the cylinder is located at the point  $x = x_{\text{omin}} + L/10$ ,  $y = y_{\text{omin}} + H/2$  with  $(x_{\text{omin}}, y_{\text{omin}}, z_{\text{omin}})$  the origin of the frame.
- Set these values in the window "maillagedefault":
  - Add the declaration of the parameter "radius" in the first line of the "maillagedefault" file.
  - Set the values  $x_{\text{omin}} = 0$ ,  $x_{\text{omax}} = 10$ ,  $y_{\text{omin}} = 0$ ,  $y_{\text{omax}} = 2$ ,  $z_{\text{omin}} = 0$  and  $z_{\text{omax}} = 1$ .
  - Add the initialization of the radius:  $radius = 0.4$ .

# Xprepro to create a 3D VDF simple mesh



- Save your file.
- You can see the definition of  $XM(1/nx)$ ,  $YM(1/ny)$  and  $ZM(1/nz)$ , it represent the boundaries of the domain with a width of  $\varepsilon = 0.01$ .
- You can see on the top of the "viewlist" window the values of  $nx$ ,  $ny$  and  $nz$ . For the moment they are set to 4, it is the minimal number of nodes in an Xprepro mesh. Indeed the first point in the x-direction is on  $XM(1)$ , the second on  $xomin$ , the third on  $xomax$  and the fourth  $XM(nx)$ .
- Note that the "real" number of cells in the final domain is:  $Nx = nx - 3$ ,  $Ny = ny - 3$  and  $Nz = nz - 3$ .



# Xprepro to create a 3D VDF simple mesh

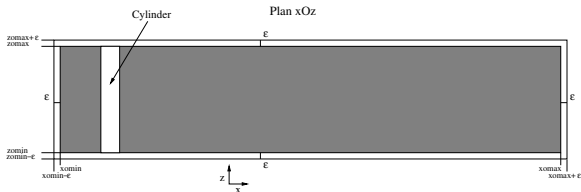
- Note that the "maillagefaut" file and your prepro file are saved in the directory Formation\_TRUST/yourname/Xprepro/model.
- Click on "Modify" on the top of the "viewlist" window to change the values of  $n_x$ ,  $n_y$  and  $n_z$  to have  $N_x = 100$ ,  $N_y = 20$  and  $N_z = 10$ . Then "Ok".
- Do not forget to save your geometry with "Save file prepro"!
- You can see in this window that there are some "?", we will complete them now.
- Note that lines ending by a "(Comm)" are commented lines.
- Choose the index number of the matter by double-clicking on the line 5 "filling up of the domain...", set the index of the matter to 1000. (For matters, we must use positive numbers. Matters with a negative index will be deleted.)
- Double-click on the line 7 "back boundary...", change the name of the boundary from "back boundary" to "Wall", and set the matter index to -1000. (Negative numbers for boundaries.)

# Xprepro to create a 3D VDF simple mesh

- Do the same thing with lines 8 to 12:
  - line 8, change the "front boundary" name to "Wall", INDMAT=-1000,
  - line 9, change the "left boundary" name to "Inlet", INDMAT=-2000,
  - line 10, change the "right boundary" name to "Outlet", INDMAT=-3000,
  - line 11, change the "bottom boundary" name to "Bottom", INDMAT=-1000,
  - line 12, change the "top boundary" name to "Top", INDMAT=-1000.
- Add the cylinder, click on "Add", then click on the "□" button and select "cylinder".
- Name it "Obstacle", set the values of (AC) and (BC) with C the center of the cylinder so (AC, BC) are the coordinates of C in the xOy plane, so  $AC = xomin + (xomax - xomin)/10$  and  $BC = yomin + (yomax - yomin)/2$ .
- Set the Radius to "radius".
- Set  $CMIN = zomin$  and  $CMAX = zomax$ ,  $CMAX - CMIN$  corresponds to the cylinder width.
- Set IDIR to 3, the cylinder axis and INDMAT to -9000 (to make a hole) then "Ok".

# Xprepro to create a 3D VDF simple mesh

- The new line defining the cylinder appears in the "viewlist" window (save your prepro file).
- Click on "Modeling run..." in the command window. A pop-up window appears with the nodes list, click "Ok".
- Click on "Pre-mesh visualisation", VisIt opens. You can:
  - visualize your mesh with Add → Mesh → dom\_IJK and
  - visualize the matter indexes with Add → Pseudocolor → INDMAT\_ELEM\_dom\_IJK.
- We have a box but no hole! In fact we cannot see the hole because it is inside the box and it doesn't pass through it.



# Xprepro to create a 3D VDF simple mesh

- So we must change the values of  $CMIN$  and  $CMAX$  in the cylinder parameters, to put  $CMIN = zomin - \varepsilon$  and  $CMAX = zomax + \varepsilon$
- Close VisIt, and click on "Modeling run..." and then "Pre-mesh visualization".
- Note that a cell is composed of a matter if its barycenter is in this matter's zone.
- Close VisIt, click on "Pre-processing run...", you can see the names of the boundaries in function of the indexes that we gave. For exemple, the boundary number 4 is made of the boundaries "Wall", "Wall", "Bottom" and "Top" and it's TRUST name will be "Wall\_Wall\_Bottom\_Top".
- You can change the names of these boundaries, to have the name you want.
- To give a name to a boundary is not mandatory for the boundaries with the same matter index. You have just to name at least one of it.
- Click on "Get geom", a nedit window opens with a TRUST data file named "default.mesh".

# Xprepro to create a 3D VDF simple mesh

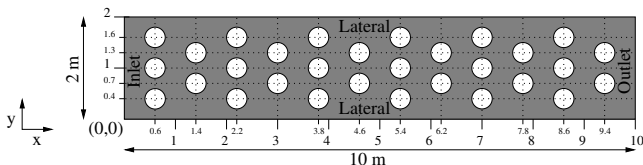
- Save this file with the name "channel.data", you can see that your geometry is in the file "default\_Pb1.geom".
- Quit Xprepro.
- Edit your channel.data file and add at the end of the file:  
**Postraiter\_domaine { domaine dom\_pb1 fichier dom\_pb1.lata format lata }**  
**discretiser\_domaine dom\_pb1**  
**End**
- Run the file: **trust channel**
- Visualize your mesh with: **visit -o dom\_pb1.lata &**

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# Xprepro to create a 2D VDF mesh

## Second exercise

In this exercise, you will learn how to create a 2D mesh with Xprepro.



The radius of the cylinders is 0.2m and the distances between cylinders are 0.2m in the y direction and 0.6m in the x direction.

- Create a directory: Run Xprepro in the TRUST environment:  
**mkdir -p Formation\_TRUST/yourname/Xprepro/exo2**  
**cd Formation\_TRUST/yourname/Xprepro/exo2**  
**source /home/triou/env\_TRUST\_X.Y.Z.sh**  
**Xprepro &**

# Xprepro to create a 2D VDF mesh

- First, have a look at the "Examples..." ("Picture" button and if you are interested by one "Copy and Read" button).
- Then, click on the button "Default" to define the initial block and edit the dimensions file.
- Note that the files "maillagedefault" and "default.prep" are saved in your directory (with "save" in nedit for "maillagedefault" and "Save file prepro" in Xprepro for default.prep).
- It is a 2D geometry, so we will define:
  - `xomin=0.`,
  - `xomax=10.`,
  - `yomin=0.`,
  - `yomax=2.`,
  - `zomin=0.`,
  - `zomax=0.01.`,
  - `eps=0.0001` (tolerance) and
  - `radius=0.2` (don't forget to declare it).
- Save the file.



# Xprepro to create a 2D VDF mesh

- Click "Modify" on the viewlist to modify the nodes number  $NX=203$ ,  $NY=43$ ,  $NZ=4$ .
- Set the matter index to 1000, line 5.
- Build the boundary blocks:
  - line 7, change the "back boundary" name to **lateral**,  $INDMAT=0$ ,
  - line 8, change the "front boundary" name to "",  $INDMAT=0$ ,
  - line 9, change the "left boundary" name to **inlet**,  $INDMAT=-1000$ ,
  - line 10, change the "right boundary" name to **outlet**,  $INDMAT=-2000$ ,
  - line 11, change the "bottom boundary" name to "",  $INDMAT=0$ ,
  - line 12, change the "top boundary" name to "",  $INDMAT=0$ .
- The coordinates of the center of the first cylinder, which after will be duplicated, are  $(x_0, y_0)=(0.6, 0.4)$ . So the distance between two cylinder center is  $dx = 0.8$  and  $dy = 0.3$ . Declare and initialize these parameters  $(x_0, y_0, dx, dy)$  in your "maillagedefault" file.
- Think about how to create 2 Fortran nested loops to copy the first cylinder in the two directions X and Y.

## Xprepro to create a 2D VDF mesh

- Create them with the button "Add", select "(fortran code)" instead of "☐" and write your loops in fortran.
- Check the other objects in the viewlist and save your work with the button "Save file" prepro then check the name of the boundaries with "Boundaries information".
- Run the model with "Modeling run...". A window is opened where you will check the nodes coordinates of the mesh.
- Click on "Pre-mesh visualization" to check your pre-mesh.
- Create a 2D cut slice in the XY plane: with "Add" button, choose an object of "Meshing creation 2D" type and name it coupe\_2D. Set  $POS = 0$ ,  $IDIR = 3$ ,  $INDMAT = -5000$ .
- Click "Modeling run..." then "Pre-mesh visualization".
- Warning, it is always a 3D model in Xprepro, even if you wish a 2D mesh. By default, you see the indexes between 1000 and -3000. If you don't see the cylinders, create a 2D slice in the XY plane in VisIt to see inside the 3D pre-mesh.

# Xprepro to create a 2D VDF mesh

- Zoom onto the boundaries to check that the boundary blocks are all defined.
- Click "Pre-processing run" to create the final mesh (check there is no error messages).
- Click on "Get geom" to generate the TRUST .geom file in your study. It opens a nedit window with a TRUST data file, save it in your repository with the name "dom\_1.data".
- Suppress the first lines in comments "#" and the "\n" ending a commentary line.
- Add at the end of the file:  
**Discretiser\_domaine dom\_1**  
**Postraiter\_domaine { domaine dom\_1 fichier mesh format lata }**
- Visualize your 2D mesh with VisIt.
- If you wish, build a data file to read your mesh and run the flow around the cylinders.

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# Validation form

## Example of Validation form:

2. COMPARISON BETWEEN FLOW RATE SPECIFIED BY DEBIT.IMPOSE OPTION AND COMPUTED FLOW RATE BY THE INITIAL CONDITION ON VELOCITY

Check the debit.impose option of canal\_perio keyword

### 1 Introduction

Validation made by : G.F.  
Report generated 04/04/2015.

#### 1.1 Description

#### 1.2 Parameters Trio.U

- Variable Trio.U :
- Variable Trio.U from out: /export/home/ser737086/gth/Bouvier\_171/Trio.U/conv/Trio.U\_inpt\_opt (1.7.1.beta)

#### 1.3 Test cases

- ./std.data :
- ./debit.data :
- ./debit2.data :
- ./debit3.data :

2. Comparison between flow rate specified by debit.impose option and computed flow rate by the initial condition on velocity

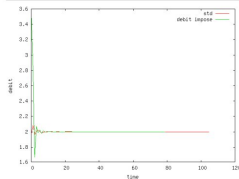
Data flow difference:

```
T5/T5
<
- vitesse champ.outflow 2 1. 0.
>
- vitesse champ.outflow 2 2 0.
T5/T2
<
- source { Canal.perio { bond.perio } }
>
- source { Canal.perio { bond.perio debit.impose 2. } }
```

In the first data file, the flow rate will be 2 m<sup>3</sup>/s (H=0.5 m/s and S=2m). In the second one, flow rate is 4 m<sup>3</sup>/s, and then will decrease to 2 m<sup>3</sup>/s.

Check the debit.impose option of canal\_perio keyword

3. INITIAL VELOCITY IS INCLINED INTO 2 DIRECTIONS, WITH A VERTICAL FLOW RATE WHICH SHOULD BE 0.

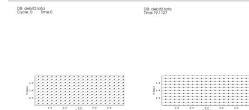


3. Initial velocity is inclined into 2 directions, with a vertical flow rate which should be 0.

When converged, the velocity profile shows horizontality.

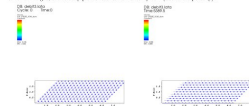
Check the debit.impose option of canal\_perio keyword

4. INITIAL VELOCITY IS INCLINED INTO 2 DIRECTIONS, WITH A VERTICAL FLOW RATE WHICH SHOULD BE 0.



4. Initial velocity is inclined into 2 directions, with a vertical flow rate which should be 0.

When converged, the velocity profile shows horizontality to 4 (due to the penalty) .



Check the debit.impose option of canal\_perio keyword

# Validation form

- First copy the validation form named Source\_canal\_perio:  
**mkdir -p Formation\_TRUST/yourname/validation**  
**cd Formation\_TRUST/yourname/validation**  
**VERIF=\$TRUST\_ROOT/Validation/Rapports\_automatiques/Verification**  
**cp -r \$VERIF/Verification\_codage/Source\_canal\_perio .**  
**cd Source\_canal\_perio**
- Ask for help about the Run\_fiche script:  
**Run\_fiche -help**
- Build the report:  
**Run\_fiche -xpdf**  
or  
**Run\_fiche**  
**evince build/rapport.pdf &**

# Validation form

- Now, we are going to change the validation form (Examples are given in pages 10 & 11 of the HowTo\_Validation.pdf note):

## **nedit src/Canal.prm &**

- Add the mesh plot in the report. For this, at end of .prm file, introduce a new block with "visu" keyword:

```
Chapter {
  Title "Additional information"
  Visu {
    Title "Mesh visualization"
    Mesh lata_file_name domain_name [color]
  }
}
```

- You can help with:
  - \$TRUST\_ROOT/Validation/Outils/Genere\_courbe/doc/manuel.xhtml#paramVisu
  - \$TRUST\_ROOT/Validation/Outils/Genere\_courbe/doc/exemples/visu.prm
- Save the .prm file and re-build the report without running the calculations:

**Run\_fiche -not\_run**

# Validation form

- Add the evolution of residuals in the report in log scale (see .dt\_ev file). For this, introduce a new block with "Figure" and "Curve" keywords:

- Complete the chapter "Additional information" with a new block "Figure":

```
Figure {
  Title "Residuals evolution"
  LabelX "Time (s)"
  LabelY "Residual"
  LogX
  LogY
  Include_description_Curves 0
  Curve {
    legend data_file_name.data
    file data_file_name.dt_ev
    columns $column_number_for_x $column_number_for_y
    style lines
  }
  ...
}
```

- If you need help, see:
  - ◇ \$TRUST\_ROOT/Validation/Outils/Genere\_courbe/doc/manuel.xhtml#paramFigure
  - ◇ \$TRUST\_ROOT/Validation/Outils/Genere\_courbe/doc/manuel.xhtml#paramCourbe
  - ◇ \$TRUST\_ROOT/Validation/Outils/Genere\_courbe/doc/exemples/impl.prm



# Validation form

- Visualize the pressure field at the last time: complete the chapter "Additional information" with a new block "Visu" using "PseudoColor" keywords. The name of the field and its localization must be uppercase letter:

```
Visu {
  Title "Visualization of pressure field at the last time"
  Description "Pressure field..."
  Pseudocolor data_file_name.lata domain_name FIELD POSITION
  Cycles cycle_numbers
  Width size_in_cm
}
```

NB: Cycle '-1' correspond to the last time.

NB: You can use the "magnitude" keyword to visualize the velocity field without using arrows (Syntax: Pseudocolor *data\_file\_name.lata dom\_magnitude VITESSE SOM*).

- Save and close the .prm file and build the report with the options: (Options are given in page 7 of the HowTo\_Validation.pdf note)

**Run\_fiche -xpdf -not\_run**

# Validation form

- Now, we are going to extract the number of cells and the last time from three .err files and write it in .dat files via a post\_run script.
  - Create a "post\_run" file in the src directory containing:
 

```
nb1=`grep "Total number of elements" std.err | awk '{print $NF}'`
nb2=`grep "Total number of elements" debit.err | awk '{print $NF}'`
nb3=`grep "Total number of elements" debit2.err | awk '{print $NF}'`
echo $nb1 $nb2 $nb3 > nbcells.dat
tp1=`grep "Backup of the field" std.err | awk '{print $NF}' | head -n 1`
tp2=`grep "Backup of the field" debit.err | awk '{print $NF}' | head -n 1`
tp3=`grep "Backup of the field" debit2.err | awk '{print $NF}' | head -n 1`
echo $tp1 $tp2 $tp3 > lasttime.dat
```
  - Add a table to display the results of .dat files: complete the chapter "Additional information" by introducing a new block with "table" and "line" keywords.

If you need help, see:

- ◇ [\\$TRUST\\_ROOT/Validation/Outils/Genere\\_courbe/doc/manuel.xhtml#paramTableau](#)
- ◇ [\\$TRUST\\_ROOT/Validation/Outils/Genere\\_courbe/doc/manuel.xhtml#paramLigne](#)
- ◇ [\\$TRUST\\_ROOT/Validation/Outils/Genere\\_courbe/doc/exemples/tableau.prm](#)

# Validation form

- The table block will look like:

```

Table {
  Title "Number of cells and last time results"
  Nb_columns number_of_colomns_without_the_first_one
  Label label_of_the_first_column | label_of_the_second_column ...
  Line {
    legend title_of_the_first_line
    file name_of_the_file.dat
  }
  ...
}

```

- Save the .prm file in src directory and build the report with a new option:  
**Run\_fiche -xpdf -post\_run**

# Validation form

- Now we are going to modify the "prepare" file in "src" directory in order to add a fourth test case: "debit4"
  - "debit4" correspond to "std" test case with zero initial velocity and imposed flow rate to  $2m^3/s$  on "periox" boundary.
  - Add a line in the "prepare" file using the "sed" command to create the new data file like the other ones:  
**nedit src/prepare &**
- We are going to update the validation form.
  - Modify the "Canal.prm" file in "src" directory to take into account the new test case debit4. Add in the **Parameters** block :  
**TestCase . Debit4.data**
  - Build the report by running the 4 test cases simultaneously (not sequentially):  
**Run\_fiche -parallel\_run -xpdf**
  - You can add the results of this test case to your "visu" and "table".

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# Unix Quick Reference

## File Commands

<b>ls</b>	Directory listing
<b>ls -al</b>	Formatted listing with hidden files
<b>ls -lt</b>	Sorting the Formatted listing by time modification
<b>cd dir</b>	Change directory to dir
<b>cd</b>	Change to home directory
<b>pwd</b>	Show current working directory
<b>mkdir dir</b>	Creating a directory dir
<b>cat &gt;file</b>	Places the standard input into the file
<b>more file</b>	Output the contents of the file
<b>head file</b>	Output the first 10 lines of the file
<b>tail file</b>	Output the last 10 lines of the file
<b>tail -f file</b>	Output the contents of file as it grows, starting with the last 10 lines
<b>touch file</b>	Create or update file
<b>rm file</b>	Deleting the file
<b>rm -r dir</b>	Deleting the directory

# Unix Quick Reference

## File Commands

<b>rm -f file</b>	Force to remove the file
<b>rm -rf dir</b>	Force to remove the directory dir
<b>cp file1 file2</b>	Copy the contents of file1 to file2
<b>cp -r dir1 dir2</b>	Copy dir1 to dir2;create dir2 if not present
<b>mv file1 file2</b>	Rename or move file1 to file2,if file2 is an existing directory
<b>ln -s file link</b>	Create symbolic link link to file

# Unix Quick Reference

## Process management

<b>ps</b>	To display the currently working processes
<b>top</b>	Display all running process
<b>kill pid</b>	Kill the process with given pid
<b>killall proc</b>	Kill all the process named proc
<b>pkill pattern</b>	Will kill all processes matching the pattern
<b>bg</b>	List stopped or background jobs, resume a stopped job in the background
<b>fg</b>	Brings the most recent job to foreground
<b>fg n</b>	Brings job n to the foreground

## File permission

<b>chmod octal file</b>	Change the permission of file to octal, which can be found separately for user, group, world by adding: 4-read(r) 2-write(w) 1-execute(x)
-------------------------	--



# Unix Quick Reference

## Searching

<b>grep pattern file</b>	Search for pattern in file
<b>grep -r pattern dir</b>	Search recursively for pattern in dir
<b>command   grep pattern</b>	Search pattern in the output of a command
<b>locate file</b>	Find all instances of file
<b>find . -name filename</b>	Searches in the current directory (represented by a period) and below it, for files and directories with names starting with filename
<b>pgrep pattern</b>	Searches for all the named processes , that matches with the pattern and, by default, returns their ID

# Unix Quick Reference

## System Info

<b>date</b>	Show the current date and time
<b>cal</b>	Show this month's calender
<b>uptime</b>	Show current uptime
<b>w</b>	Display who is on line
<b>whoami</b>	Who you are logged in as
<b>finger user</b>	Display information about user
<b>uname -a</b>	Show kernel information
<b>cat /proc/cpuinfo</b>	Cpu information
<b>cat proc/meminfo</b>	Memory information
<b>man command</b>	Show the manual for command
<b>df</b>	Show the disk usage
<b>du</b>	Show directory space usage
<b>free</b>	Show memory and swap usage
<b>whereis app</b>	Show possible locations of app
<b>which app</b>	Show which applications will be run by default

# Unix Quick Reference

## Compression

<b>tar cf file.tar file</b>	Create tar named file.tar containing file
<b>tar xf file.tar</b>	Extract the files from file.tar
<b>tar cf file.tar file</b>	Create tar named file.tar containing file
<b>tar xf file.tar</b>	Extract the files from file.tar
<b>tar czf file.tar.gz files</b>	Create a tar with Gzip compression
<b>tar xzf file.tar.gz</b>	Extract a tar using Gzip
<b>tar cjf file.tar.bz2</b>	Create tar with Bzip2 compression
<b>tar xjf file.tar.bz2</b>	Extract a tar using Bzip2
<b>gzip file</b>	Compresses file and renames it to file.gz
<b>gzip -d file.gz</b>	Decompresses file.gz back to file

# Unix Quick Reference

## Network

<b>ping host</b>	Ping host and output results
<b>whois domain</b>	Get whois information for domains
<b>dig domain</b>	Get DNS information for domain
<b>dig -x host</b>	Reverse lookup host
<b>wget file</b>	Download file
<b>wget -c file</b>	Continue a stopped download

# Unix Quick Reference

## Shortcuts

<b>"Ctrl" + c</b>	Halts the current command
<b>"Ctrl" + z</b>	Stops the current command, resume with fg in the foreground or bg in the background
<b>"Ctrl" + d</b>	Logout the current session, similar to exit
<b>"Ctrl" + w</b>	Erases one word in the current line
<b>"Ctrl" + u</b>	Erases the whole line
<b>"Ctrl" + r</b>	Type to bring up a recent command
<b>!!</b>	Repeats the last command
<b>exit</b>	Logout the current session

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## **Tutorial solutions:**

*`$TRUST_ROOT/doc/TRUST/exercices/Tutorial_solutions.pdf`*