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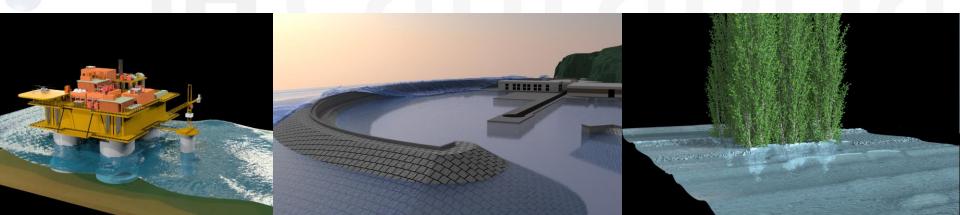
R+D+i for a Sustainiable Development

-



(Beginners Course) **IHFOAM** applied to Coastal Engineering Breaking solitary waves on a mild slope (2D)

Gabriel Barajas, Javier L. Lara, María Maza





- Now we are going to create a 3D case from a 2D one:
 - Copy a 2D case and change the folder name:
 - \$ cp -r ~/OpenFOAM-v1812/tutorials/multiphase/interFoam/laminar/ waveExampleSolitary ~/IHFoamCourse/.
- Rename the case:
 - Rename the case:
 - \$ mv ~/IHFoamCourse/waveExampleStokesV ~/IHFoamCourse/synolakis
 - Set OpenFOAM environment:
 - \$ source ~/OpenFOAM/OpenFOAM-v1812/etc/bashrc
 - Ensure everything you don't need is deleted
 - \$ cd ~/IHFoamCourse/synolakis
 - \$./Allclean



OpenFOAM workflow

Breaking solitary wave on a mild slope (2D)

Numeric Domain

Mesh generation

Boundary conditions, Initial conditions, Numerical parameters, Pre-processing

Run the Simulation

Numerical Setup

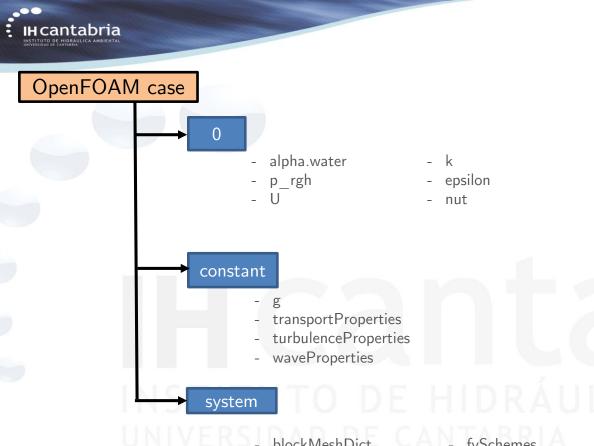
Geometric Domain

Execute serial or parallel

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Post-processing

Wave gauges, run-up, etc.



blockMeshDict

setFieldsDict _

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- snappyHexMeshDict
- extrudeMeshDict _

fvSchemes

- fvSolution _
- decomposeParDict -
- controlDict _

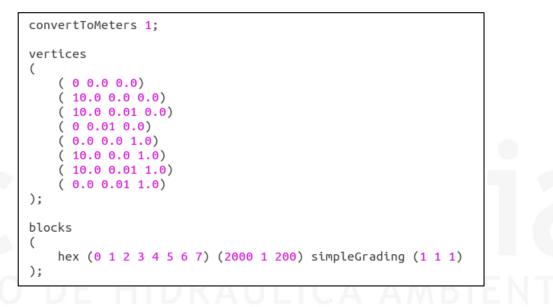
• Update *system/blockMeshDict* to fit the laboratory set-up (synolakis 1986):



\$ blockMesh

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Check the base mesh quality:
 \$ checkMesh





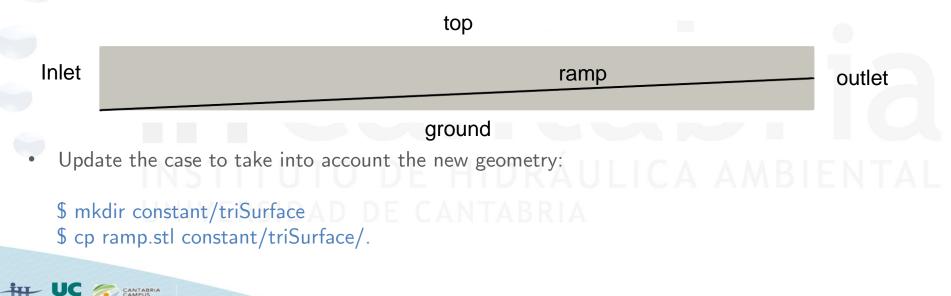
In the *0.org* folder, display the VoF (*alpha.water*), the velocity (*U*) and the pressure (*p_rgh*) files :

 alpha.water
 U
 p_rgh

		alphantatoi			0	Po.,
	dimensions internalField	[0 0 0 0 0 0 0]; uniform 0;	dimensions internalField	[0 1 -1 0 0 0 0]; uniform (0 0 0);	dimensions internalField	[1 -1 -2 0 0 0 0]; d uniform 0;
	boundaryField { { type value } outlet	waveAlpha; uniform 0;	boundaryField { inlet { type value } outlet	<pre>waveVelocity; uniform (0 0 0);</pre>	boundaryField { inlet { type value } outlet	fixedFluxPressure;
	{ type } ground	zeroGradient;	{ type value } sides	<pre>waveVelocity; uniform (0 0 0);</pre>	{ type valu } ground	fixedFluxPressure; e uniform 0;
	{ type } sides	zeroGradient;	{ type } ground	empty;	{ type value }	fixedFluxPressure; e uniform 0;
	type } top {	empty;	{ type value } top	fixedValue; uniform (0 0 0);	sides { type } top	empty;
+	type inletVa value } }	uniform 0;	{ type value } }	pressureInletOutletVeloc uniform (0 0 0);	{	totalPressure; uniform 0;
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- Define and create a mild slope (using Autocad, Rhino, etc.).
- Check the .stl file; open Paraview, load the .stl file and check that the geometry fits the base mesh:
 - \$ touch ih.foam && paraview

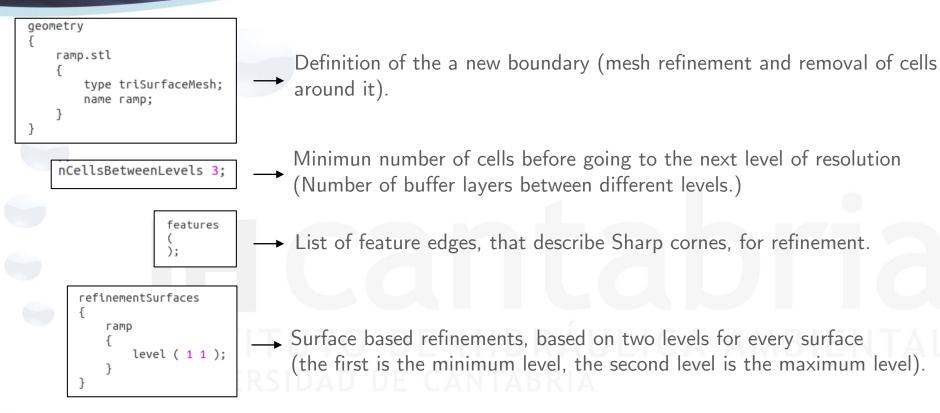




- Using *snappyHexMesh*, as mesh generator to take the existing base mesh and remesh it to fit the real geometry of the experiments.
- Copy snappyHexMeshDict from a tutorial:
 - \$ cp -r ~/OpenFOAM-v1806/tutorials/multiphase/interFoam/RAS/mixerVesselAMI/system/ snappyHexMeshDict ~/IHFoamCourse/overSetWaves/system/.
- This intermediate mesh, is created from the dictionary **system/snappyHexMeshDict**:
 - CastellatedMesh:
 - Mesh Refinement in prescribed regions.
 - Detection of the domain (surface and volume).
 - Removal of cells outside the domain.
 - Snap: NIVERSIDAD DE CANTABRIA
 - Mesh morphing to follow the provided geometry.
 - Layer addition could also be done.

castellatedMesh tr	eps to run ue;
	ue; lse;





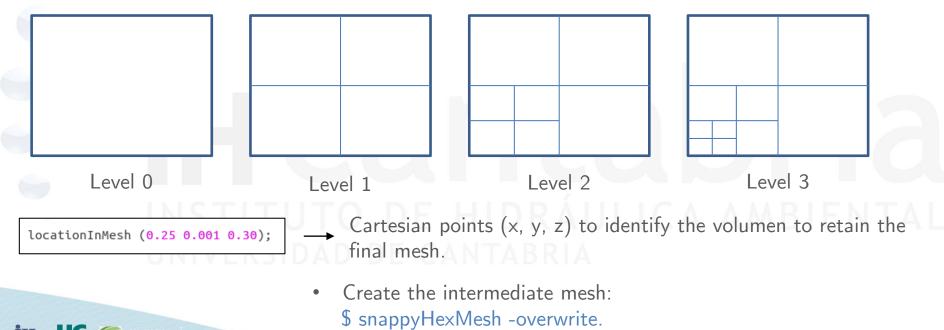




refinementRegions
{
}

Volume based refinements, based on two levels for every volume (the first is the minimum level, the second level is the maximum level).

• Refinement levels in OpenFOAM: increase in the refinement level reduces the cell size by half.





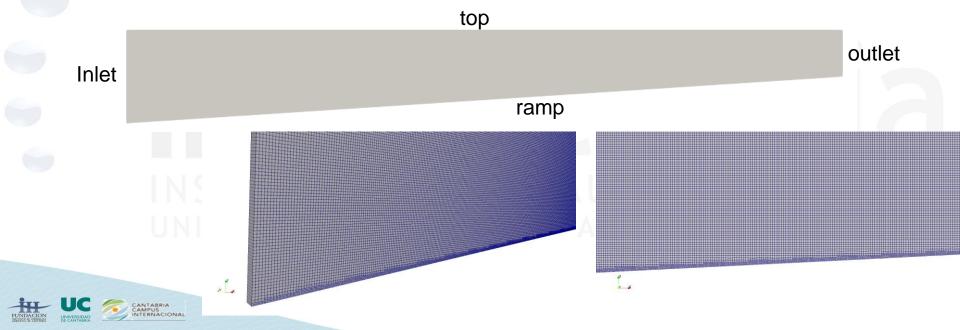
• Ajust the mesh: as it has been created by **snappyHexMesh** with several cells in the spanwise direction, it must be modified and extrude it to be purely two dimensional (2D).

```
constructFrom patch;
sourceCase ".";
sourcePatches (side1):
// If construct from patch: patch to use for back (can be same as sourcePatch)
exposedPatchName side2:
// Flip surface normals before usage.
flipNormals true;
//- Linear extrusion in point-normal direction
extrudeModel
                    linearNormal:
nLavers
                    1:
expansionRatio
                    1.0:
linearNormalCoeffs
    thickness
                    0.02;
// Do front and back need to be merged? Usually only makes sense for 360
// degree wedges.
mergeFaces false:
```

- **sourcePatches**: name of the patch to extrude.
- **exposedPatchName**: name of the patch opposed to the extruded one (sourcePatches)
- *nLayers*: number of divisions from sourcePatches to exposedPatchName
- thickness: length of the extrusion.
- Extrude the intermediate mesh:
 \$ extrudeMesh
- Check the final mesh quality:
 \$ checkMesh



- Check your final mesh with Paraview:
 \$ paraview
- Load the ih.foam file and press "Apply". (Remember to tick "Skip Zero Time", as the boundary conditions in the 0 folder have not been updated yet.)





 The final boundaries can be checked with Paraview (in the *Mesh Regions* dialog box) or they can be checked in the *constant/polyMesh/boundary* file.

 ✓ internalMesh inlet outlet ground top side1 side2 ramp 	
	R Á BRI

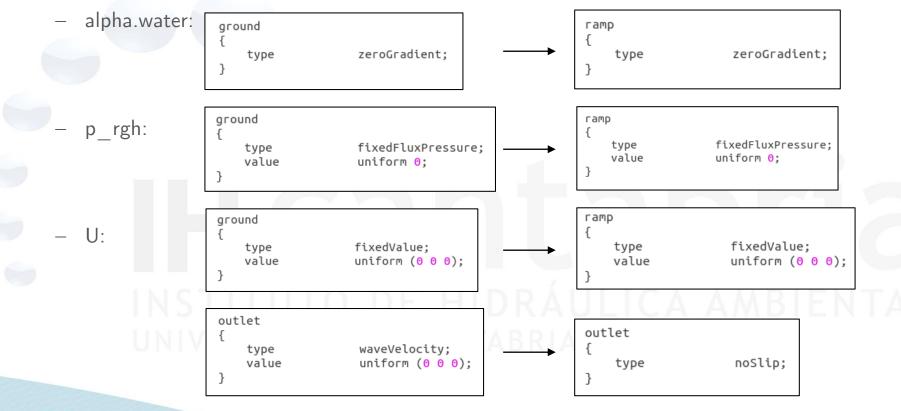


			_
inl	et		
{			
	type	patch;	
	nFaces	200;	
	startFace	708588;	
} .			
	let		
{	A		
	type	patch;	
	nFaces startFace	102;	
}	Startrace	708788;	
	und		
910 {	unu		
ι	type	wall;	
	inGroups	1(wall);	
	nFaces	1000;	
	startFace	708890;	
}		,	
top			
{ '			
	type	patch;	
	nFaces	2000;	
	startFace	709890;	
}			
sid	e1		
{			
	type	empty;	
	inGroups	<pre>1(empty);</pre>	
	nFaces	710788;	
	startFace	711890;	
}			
sid	ez		
{	+		
	type inGroups	empty;	
	nFaces	1(empty);	
	startFace	0; 1422678;	
}	startrace	1422070,	
л гам	n		
{	P		
L.	type	wall;	
	inGroups	1(wall);	
	nFaces	2000;	
	startFace	1422678;	
}		,	
-			



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• Once the final boundaries are known, update **0.org** folder:





 The case is defined as turbulent in: \$ more constant/turbulenceProperties

sim	ulationType	RAS;
RAS		
{	RASModel	kEpsilon;
	turbulence	on;
}	printCoeffs	on;

Therefore, the turbulent kinematic energy (k), the turbulent dissipation (epsilon) and the turbulent viscosity (nut) variables must be defined and added to the 0.org folder:

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		k	(epsilon	ľ	nut	
	dimensions internalField	[0 2 -2 0 0 0 0]; uniform 0.0084;		dimensions internalField	[0 2 -3 0 0 0 0]; uniform 0.0005094;		dimensions internalField	[0 2 -1 0 0 0 0]; uniform 0;
	boundaryField { inlet { type	zeroGradient;		boundaryField { inlet { type	zeroGradient;		boundaryField { inlet { type value	calculated; uniform 0;
	J outlet { type } side1	zeroGradient;		} outlet { type } side1	zeroGradient;		} outlet { type value }	calculated; uniform 0;
	{ type } side2 {	empty;		{ type } side2	empty;		side1 { type }	empty;
	type } ground {	empty;		type } ground	empty;		side2 { type }	empty;
	type value } ramp {	kqRWallFunction; uniform 0.0084;		type value } ramp	epsilonWallFunction; uniform 0.0005094;		ground { value }	nutkWallFur uniform 0;
	type value } top {	kqRWallFunction; uniform 0.0084;	А	t type value } top	epsilonWallFunction; uniform 0.0005094;		ramp { type value } top	nutkWallFur uniform 0;
LON	type inletVa value } }	inletOutlet; lue uniform 0.0084; uniform 0.0084;		type inletVa value } }	inletOutlet; alue uniform 0.0005094; uniform 0.0005094;		top { type value } }	calculated; uniform 0;

orm 0; calculated; uniform 0; calculated; uniform 0; empty; empty; nutkWallFunction; uniform 0; nutkWallFunction; uniform 0;



- Update the wave conditions in constant/waveProperties:
 - WaveModel: Boussinesq solitary wave
 - *nPaddle*: 1 single wave paddle (2d)
 - *waveHeight*: H = 0.07 m.
 - *activeAbsorption*: no (absorption yes/no at generation)

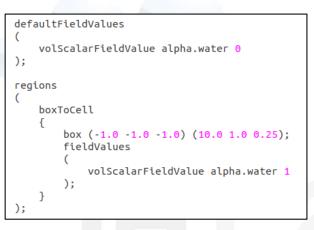
nlet					
	alpha	alpha.water;			
	waveModel	Boussinesq;			
	nPaddle	1;			
	waveHeight	0.07;			
	waveAngle	0.0;			
	rampTime	0.0;			
	activeAbsorption	n no;			
	wavePeriod	0.0;			

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• Update the initial set-up in *system/setFieldsDict*:



\$ cp -r 0.org 0\$ setFields

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 Open Paraview and plot the initial set-up to ensure eveything is correct:
 \$ paraview





Water and air properties are defined in:
 \$ more constant/transportProperties

phases (water air);	
water { transportModel nu rho }	Newtonian; [0 2 -1 0 0 0 0] 1e-06; [1 -3 0 0 0 0 0] 1000;
air { transportModel nu rho } sigma [1 (Newtonian; [0 2 -1 0 0 0 0] 1.48e-05; [1 -3 0 0 0 0 0] 1; 0 -2 0 0 0 0] 0.07;

Gravity is defined in:
 \$ more constant/g

dimensions	[0 1 -2 0 0 0 0];
value	(0 0 -9.81);

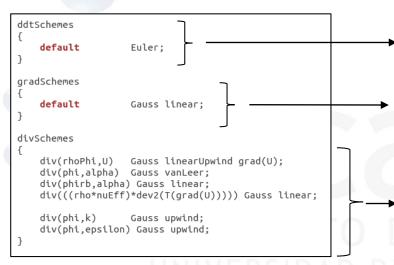
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Numerical Schemes:

• system/fvSchemes: is the file that sets the numerical scheme for the different terms.



Temporal discretization (*Euler* is a first order implicit discretisation scheme).

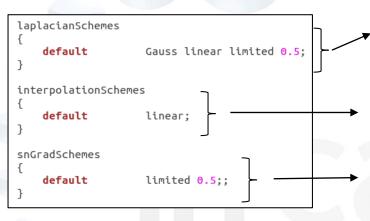
Gradient derivative terms, that is surface normal gradient terms (*Gauss linear* is second order discretisation scheme)

Divergence terms, such as advection terms and other terms that are often diffusive in nature (*Gauss linearUpwind grad(U)* is second order, *Gauss vanLeer* is second order and *Gauss upwind* is a first order numerical scheme)





Numerical Schemes:



Laplacian terms, such as the diffusion term in the momentum requation (Gauss linear limited is second order accuracy, values between 0 and 1 to handle a non-orthogonal mesh)

Cell to face interpolations of values (**linear** is second order discretisation scheme)

Component of gradient normal to a cell face (**limited** is second order accuracy, values between 0 and 1 to handle a non-orthogonal mesh)

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Algorithm control:

- In *system/fvSolutions* are defined the equations solvers, tolerances and algorithms.
- Controls for *MULES*, solver of the VoF equation:
 - *nAlphaCorr*: loops over VoF equation
 - *nAlphaSubcycles*: number of sub-cycles within the VoF equation
 - **cAlpha**: artificial compression velocity.

solvers				
{				
"alpha.water.*"				
{				
nAlphaCorr	1;			
nAlphaSubCycles	3;			
cAlpha	1;			
}				

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Algorithm control:

- For each variable solved in the particular equation, the type of solver and parameters that are used by the solver must be defined.
- Normally, the last iteration (variables are solved multiple times within a solution step) is solved with different parameters.

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```
рсогг
     solver
                      PCG;
     preconditioner
                      DIC:
     tolerance
                      1e-6:
     relTol
                      0.1;
pcorrFinal
     solver
                      PCG:
     preconditioner
                      DIC:
     tolerance
                      1e-7:
     relTol
                      0:
p_rgh
     solver
                      PCG;
     preconditioner
                      DIC:
     tolerance
                      1e-6:
     relTol
                      0.1:
p_rghFinal
     solver
                      PCG:
     preconditioner
                      DIC:
     tolerance
                      1e-7:
     relTol
                      Θ;
"(U|k|epsilon)"
                      PBiCG;
     solver
     preconditioner
                      DILU;
     tolerance
                      1e-6:
     relTol
                      0.1:
 "(U|k|epsilon)UFinal"
                      PBiCG:
     solver
     preconditioner
                      DILU:
     tolerance
                      1e-7:
     relTol
                      0:
```

Breaking solitary wave on a mild slope (2D)

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Algorithm control:

- **PIMPLE** algorithm solves the pressure-velocity coupling in the Navier-Stokes equations.
- **PIMPLE** algorithm combines PISO and SIMPLE.
 - *momentumPredictor*: switch the control for solving the momentum predictor.
 - *nCorrectors*: number of times the algorithm solves the pressure equation and momentum corrector in each step
 - **nOrthogonalCorrectors:** specifies repeated solutions of the pressure equation, used to update the explicit non-ortogonal correction.

PIMPLE				
{				
momentumPredictor no;				
nCorrectors 2;				
nNonOrthogonalCorrectors);			
1				

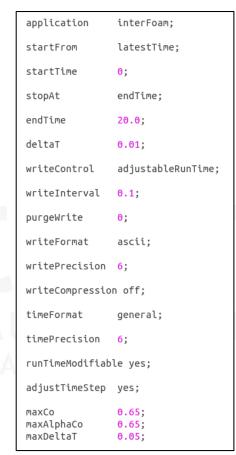




- Define simulation parameters in *system/controlDict*
 - Solver: *interFoam* (incompressible two phase flow)
 - startTime (start time for the simulation),
 endTime (end time for the simulation), deltaT (time step of the simulation).
 - writeInterval (controls the timing of write output),
 purgeWrite (integer representing a limit on the number of time directories that are stored).
 - maxCo (maximun Courant Number),
 maxAlphaCo (maximun Courant number for the pase fields), maxDeltaT (upper limit of the time step).

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- Update the runtime postprocessing sensors (*system/controlDict*)
 - to get the iso-Surface of the free surface elevation:





functions freeSurface type surfaces; functionObjectLibs "libsampling.so"): writeControl outputTime; outputInterval 1; surfaceFormat stl; interpolationScheme cellPoint; surfaces topFreeSurface type isoSurface; isoField alpha.water; isoValue 0.5; interpolate true;); fields alpha.water);



- Update the runtime postprocessing sensors (system/controlDict)
 - to get the free surface at some specific positions along the domain.

```
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```



```
line
    type
                     sets;
                     ("libsampling.so");
    libs
    enabled
                     true:
    writeControl
                    writeTime:
    writeInterval
                    1;
    interpolationScheme cellPoint;
    setFormat
                    raw;
    sets
        line1
            type
                    uniform:
                    distance;
            axis
            start
                    (1.0\ 1.0\ 0.0);
                    (1.0\ 1.0\ 0.7);
            end
            nPoints 1001;
        line2
            type
                    uniform;
            axis
                    distance;
                    ( 2.0 1.0 0.0 );
            start
                    (2.0\ 1.0\ 0.7);
            end
            nPoints 1001;
        line3
                    uniform:
            type
            axis
                     distance;
                     (3.0\ 1.0\ 0.0);
            start
            end
                     (3.0\ 1.0\ 0.7);
            nPoints 1001;
```

```
Breaking solitary wave on
a mild slope (2D)
```

```
line4
        type
                uniform:
        axis
                distance;
                (5.0\ 1.0\ 0.0);
        start
                (5.0 1.0 0.7);
        end
        nPoints 1001;
    line5
                uniform;
        type
                distance;
        axis
        start
                (7.0 1.0 0.0);
        end
                (7.0\ 1.0\ 0.7);
        nPoints 1001:
    line6
                uniform;
        type
        axis
                distance:
                (9.0\ 1.0\ 0.0);
        start
        end
                (9.0\ 1.0\ 0.7);
        nPoints 1001;
);
fixedLocations false;
fields
    U alpha.water
);
```



- Decompose case:
 - *system/decomposeParDict*: if we want to run our simulation in parallel we can decompose it using this file:

numberOffubdem

- numberOfSubdomains: set the number of parts in which we are going to split our domain.
- *n*: it should be equal to the number of subdomains

Run the command:

\$ decomposePar

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numberorsubdor	nains Z;
method	hierarchical;
hierarchicalCo { n delta order	0.001; xyz;
ſ	

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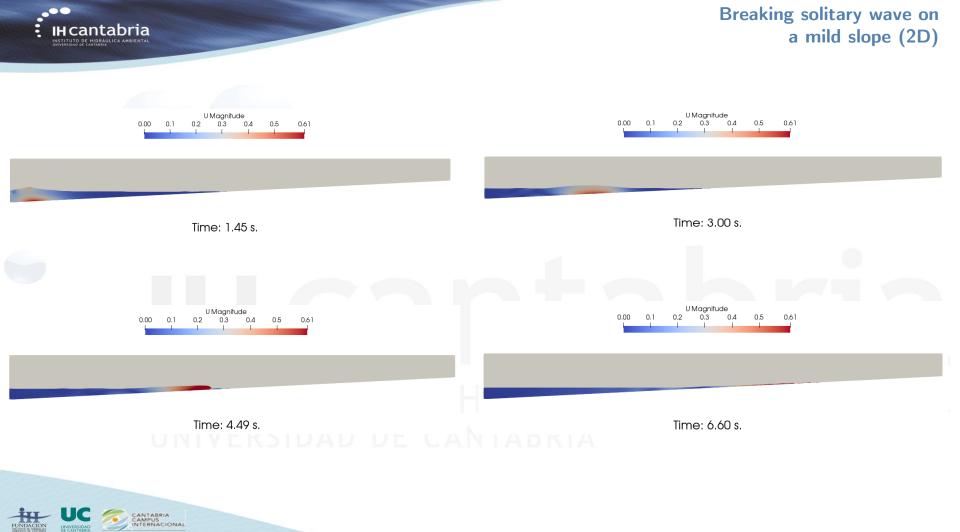
• Run the case!

\$ mpirun -np 2 interFoam -parallel > log.synolakis &
\$ tail -f log.synolakis
\$ kill PID number

Postprocessing with Paraview:
 \$ paraFoam -touch

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