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# Post-processing utilities within ElmerSolver

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# **Postprocessing utilities in ElmerSolver**

- Saving data
  - $\circ$  FEM data
  - $\circ \operatorname{Line} \operatorname{data}$
  - $\circ$  Scalars data
  - $\circ$  Grid data
- Computing data

Derived fields (gradient, curl, divecgence,...)
 Data reduction & filtering
 Creating fields of material properties

- $\circ$  Creating fields of material properties
- The functionality is usually achieved by use of atomic auxialiry solvers

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## **Computing derived fields**

• Many solvers have internal options or dedicated post-processing solvers for computing derived fields

 $\circ\,\text{E.g.}$  stress fields by the elasticity solvers

- $\circ E.g. \ \textbf{MagnetoDynamicsCalcFields}$
- Elmer offers several auxiliary solvers that may be used in a more generic way

• **SaveMaterials**: makes a material parameter into field variable

- $\circ \texttt{StreamlineSolver}: \text{ computes the streamlines of 2D flow}$
- $\circ$  **FluxSolver**: given potential, computes the flux  $q = -c \nabla \phi$

 $\circ$  **VorticitySolver**: computes the vorticity of flow,  $w = \nabla \times \phi$ 

 $\circ$  **PotentialSolver**: given flux, compute the potential -  $c \nabla \phi = q$ 

• FilterTimeSeries: compute filtered data from time series (mean, fourier coefficients,...)

0...

## **Derived nodal data**

• By default Elmer operates on distributed fields but sometimes nodal values are of interest

 Multiphysics coupling may also be performed alternatively using nodal values for computing and setting loads

• Elmer computes the nodal loads from *Ax-b* where *A*, and *b* are saved before boundary conditions are applied

**Calculate Loads = True** 

- This is the most consistant way of obtaining boundary loads
- Note: the nodal data is really pointwise

 $\circ$  expressed in units N, C, W etc.

(rather than N/m<sup>2</sup>, C/m<sup>2</sup>, W/m<sup>2</sup> etc.)

 $\odot$  For comparison with distributed data divided by the ~size of the surface elements

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## Derived lower dimensional data

• Derived boundary data

o SaveLine: Computes fluxes on-the-fly

• Derived lumped (or oD) data

SaveScalars: Computes a large number of different quantities on-the-fly
 FluidicForce: compute the fluidic force acting on a surface
 ElectricForce: compute the electrostatic froce using the Maxwell stress tensor

 Many solvers compute lumped quantities internally for later use (Capacitance, Lumped spring,...)

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#### Exporting FEM data: ResultOutputSolve

- Currently recommend format is VTU
  - $\circ \mathsf{XML}\ \mathsf{based}\ \mathsf{unstructured}\ \mathsf{VTK}$
  - $\circ \mbox{Has}$  the most complete set of features
  - Old ElmerPost format (with suffix .ep) is becoming obsolite
     Simple way to save VTU files: Post File = file.vtu
- ResultOutputSolve offers additionally several formats

   vtk: Visualization tookit legacy format
   vtu: Visualization tookit XML format
   Gid: GiD software from CIMNE: http://gid.cimne.upc.es
   Gmsh: Gmsh software: http://www.geuz.org/gmsh
   Dx: OpenDx software

#### Exporting 2D/3D data: ResultOutputSolve



An example shows how to save data in unstructured XML VTK (.vtu) files to directory "results" in single precision binary format.

```
Solver n
Exec Solver = after timestep
Equation = "result output"
Procedure = "ResultOutputSolve""ResultOutputSolver"
Output File Name = "case"
Output Format = String "vtu"
Binary Output = True
Single Precision = True
End
```

#### Saving 1D data: SaveLine

- Lines of interest may be defined on-the-fly
- Data can either be saved in uniform 1D grid, or where element faces and lines intersect
- Flux computation using integration points on the boundary not the most accurate
- By default saves all existing field variables



#### Saving 1D data: SaveLine...

```
Solver n
Equation = "SaveLine"
Procedure = File "SaveData" "SaveLine"
Filename = "g.dat"
File Append = Logical True
Polyline Coordinates(2,2) = Real 0.0 1.0 0.0 2.0
End
```

```
Boundary Condition m
Save Line = Logical True
End
```



#### Computing and saving oD data: SaveScalars



Operators on bodies

- Statistical operators
  - $\circ$  Min, max, min abs, max abs, mean, variance, deviation, rms
- Integral operators (quadratures on bodies)

   volume, int mean, int variance, int rms
   Diffusive energy, convective energy, potential energy

Operators on boundaries

- Statistical operators
  - Boundary min, boundary max, boundary min abs, max abs, mean, boundary variance, boundary deviation, boundary sum, boundary rms
  - $\circ$  Min, max, minabs, maxabs, mean
- Integral operators (quadratures on boundary)

   area

 $\circ$  Diffusive flux, convective flux

#### Other operators

 $\circ$  nonlinear change, steady state change, time, timestep size,...

#### Saving oD data: SaveScalars...

```
Solver n
 Exec Solver = after timestep
  Equation = String SaveScalars
  Procedure = File "SaveData" "SaveScalars"
  Filename = File "f.dat"
 Variable 1 = String Temperature
  Operator 1 = String \max
 Variable 2 = String Temperature
  Operator 2 = String min
 Variable 3 = String Temperature
 Operator 3 = String mean
End
```

```
Boundary Condition m
Save Scalars = Logical True
End
```



## Slots for executing postprocessing solvers

- Often the postprocessing solver need to computed only at desired slots, not at every time-step or coupled system iteration
- The execution is controlled by the "Exec Solver" keyword

   Exec Solver = before simulation
   Exec Solver = after simulation
   Exec Solver = before timesteo
   Exec Solver = after timestep
   Exec Solver = before saving
   Exec Solver = after saving
  - The before/after saving slot is controlled by the output intervals

Derived solvers often use the "before saving" slot
Data is often saved with the "after saving" slot



#### Case: TwelveSolvers

# Natural convection with ten auxialiary solvers

#### **Case: Motivation**

- The purpose of the example is to show the flexibility of the modular structure
- The users should not be afraid to add new atomistic solvers to perform specific tasks
- A case of 12 solvers is rather rare, yet not totally unrealitistic



#### **Case:** preliminaries

- Square with hot wall on right and cold wall on left
- Filled with viscous fluid
- Bouyancy modeled with Boussinesq approximation
- Temperature difference initiates a convection roll

Cold wall

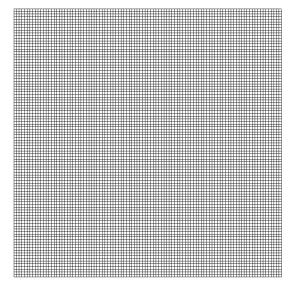






#### Case: 12 solvers

- **1**. HeatSolver
- 2. FlowSolver
- 3. FluxSolver: solve the heat flux
- 4. StreamSolver: solve the stream function
- 5. VorticitySolver: solve the vorticity field (curl of vector field)
- 6. DivergenceSolver: solve the divergence
- 7. ShearrateSolver: calculate the shearrate
- 8. IsosurfaceSolver: generate an isosurface at given value
- 9. ResultOutputSolver: write data
- 10. SaveGridData: save data on uniform grid
- **11. SaveLine**: save data on given lines
- **12. SaveScalars**: save various reductions

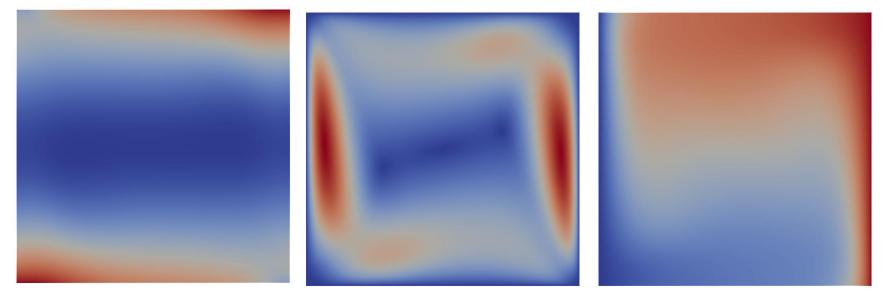


Mesh of 10000 bilinear elements



#### Primary fields for natural convection





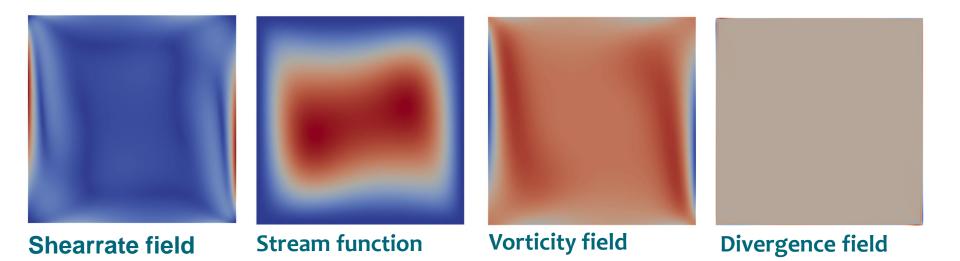
#### Pressure

#### Velocity

Temperature

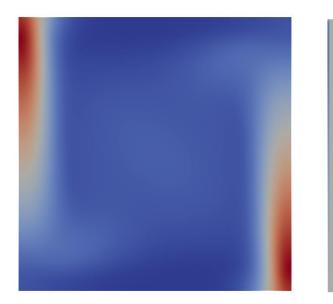


#### **Derived fields for Navier-Stokes solution**





#### **Derived fields for heat equation**



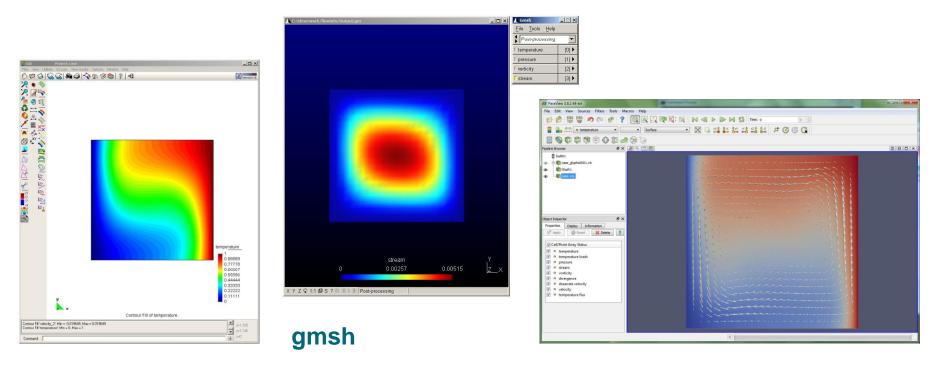


- Nodal loads only occur at boundaries (nonzero heat source)
- Nodal loads are associated to continuous heat flux by element size factor

#### **Heat flux**

**Nodal heat loads** 

#### Visualization in differen postprocessors



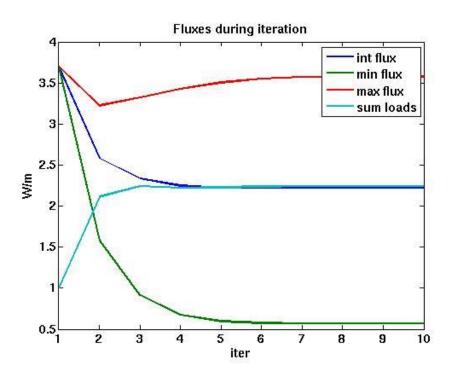




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#### **Example: total flux**

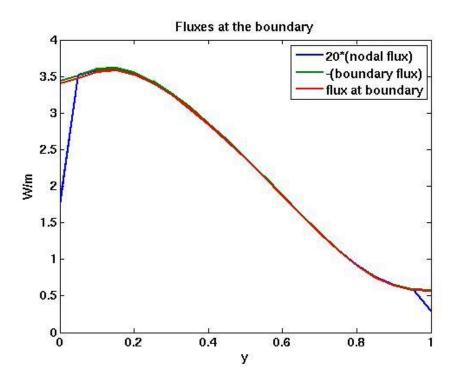
- Saved by SaveScalars
- Two ways of computing the total flux give different approximations
- When convergence is reached the agreement is good



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#### **Example: boundary flux**

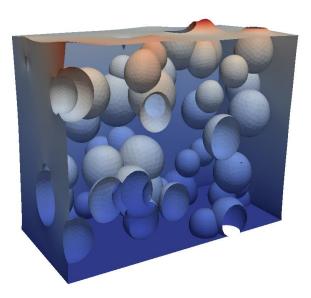
- Saved by SaveLine
- Three ways of computing the boundary flux give different approximations
- At the corner the nodal flux should be normalized using only h/2



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#### Example, saving boundaries in .sif file

Solver 2
Exec Solver = Always
Equation = "result output"
Procedure = "ResultOutputSolve"
"ResultOutputSolver"
Output File Name = case
Vtu Format = Logical True
Save Boundaries Only = Logical True
End

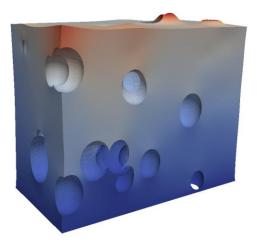




## Example, File size in Swiss Cheese

- Memory consumption of vtu-files (for Paraview) was studied in the "swiss cheese" case
- The ResultOutputSolver with different flags was used to write output in parallel
- Saving just boundaries in single precision binary format may save over 90% in files size compared to full data in ascii
- With larger problem sizes the benefits are amplified

Binary output	Single Prec.	Only bound.	Bytes/node
-	Х	-	376.0
Х	-	-	236.5
Х	Х	-	184.5
Х	-	Х	67.2
Х	Х	Х	38.5





## Manually editing the command files

- Only the most important solvers and features are supported by the GUI
- Minor modifications are most easily done by manual manipulation of the files
- The tutorials, test cases and documentation all include usable sif file pieces
- Use your favorite text editor (emacs, notepad++,...) and copy-paste new definitions to your .sif file
- If your additiones were sensible you can rerun your case
- Note: you cannot read in the changes made in the .sif file

## Exercise

- Study the command file with 12 solvers
- Copy-paste an appropriate solver from there to some existing case of your own

○ResultOutputSolverforVTU output

o StreamSolver, VorticitySolver, FluxSolver,...

- Note: Make sure that the numbering of Solvers is consistant • Solvers that involve finite element solution you need to activate by Active Solvers
- Run the modified case
- Visualize results in Paraview in different ways

#### Using tests as a starting point

- There are over 500 consistancy tests that come with the Elmer distribution
  - $\circ$   $\;$  The hope is to minimize the propability of new bugs  $\;$
- The tests are small for speedy computation
- Step-by-step instructions
  - 1. Go to tests at \$ELMER\_HOME/tests
  - 2. Choose a test case relevant to you (by name, or by grep)
    - Look in Models manual for good search strings
  - 3. Copy the tests to your working directory
  - 4. Edit the sif file
    - Activate the output writing: Post File
    - Make the solver more verbose: Max Output Level
  - 5. Run the case (see runtest.cmake for the meshing procedure)
    - Often just: ElmerSolver
  - 6. Open the result file to see what you got
  - 7. Modify the case and rerun etc.



## Conclusions



- It is good to think in advance what kind of data you need
   3D volume and 2D surface data
   Derived fields
   1D line data
   0D lumped data
- Internal strategies may allow better accuracy than doing the analysis with external postprocessing software

   Consistent use of basis functions to evaluate the data
- Often the same reduction operations may be done also at later stages but with significantly greater effort