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TRANSFORM Advanced I (60 Minutes)

• Create a model of a force-flow reactor model using TRANSFORM (60 min)





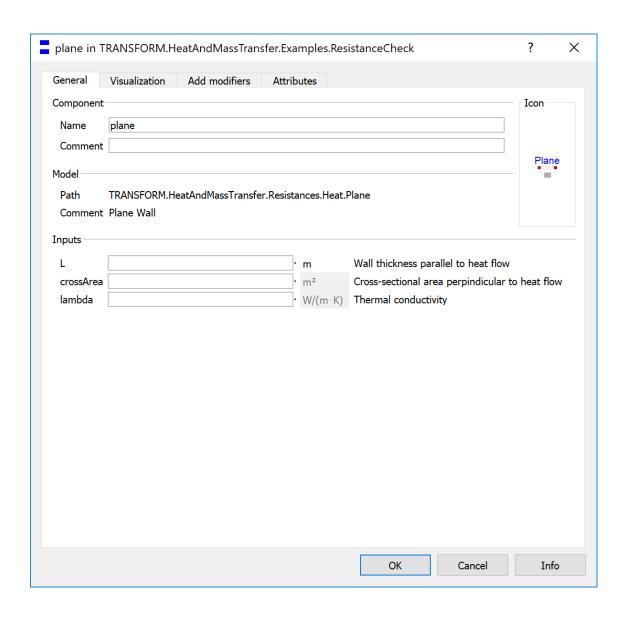
Grab Bag:

• Some random, important information



Parameter GUI: "Inputs" group

- Time dependent inputs use generally use GUI rather than connector
 - To maintain flexibility in application
 - While reducing number of "connections" (lines all over the screen)
 - Allows a wide variety of applications to be supported





Hands-On Examples:

Create a Reactor Model

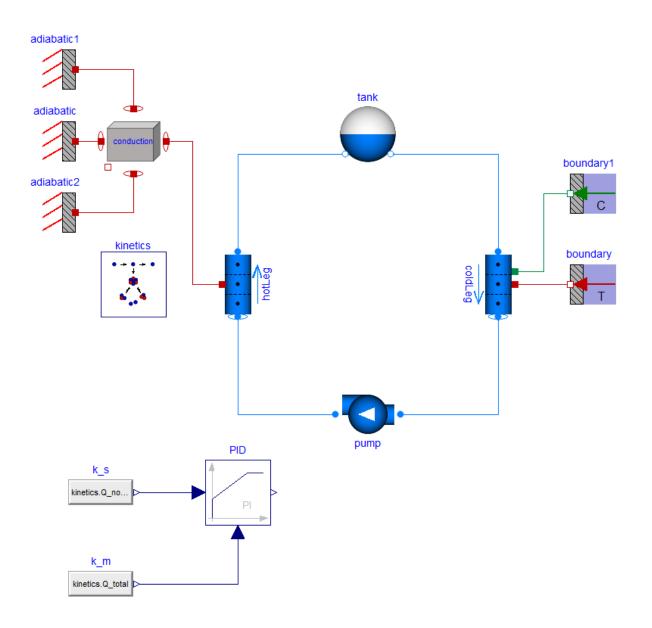
 Steps to create a simple model and gradually make it more complex



Example: Loopty Loop

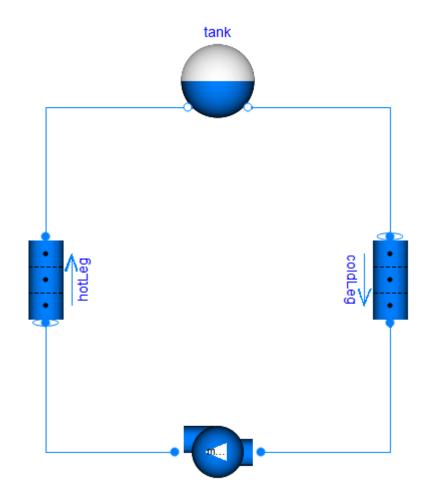
Goal:

- Using TRANSFORM components, create a forced flow reactor model with:
 - Nuclear heating
 - Level indicators
 - Pump curves
 - PID control
 - Heat transfer



Step 1: Add base components

- 1. Add a new model to your packager
- 2. Add 2 pipes to the model
 - TRANSFORM.Fluid.Pipes.GenericPipe_MultiTransferSurface
- 3. Add a tank. Why do we need this?
 - TRANSFORM.Fluid.Volumes.ExpansionTank
- 4. Add a pump
 - TRANSFORM.Fluid.Machines.Pump_SimpleMassFlow
- 5. Connect components like in the picture



Tips:

- While selecting a component, click "H", "V", or "ctrl+R" to flip horizontally, vertically, or rotate.
- In the editor, click "ctrl+A" to select all and then "ctrl+shift+L" to autoformat code. May have to do twice.



Step 2: Specify parameters

- 1. Select both pipes.
 - Specify diameter = 2" with nV = 4
 - Under "Initialization" tab set p_start = 1 bar, T_a_start = 50°C, and m_flow_a_start = 1 kg/s
- 2. For "coldLeg" pipe only, under "Advanced" tab specify
 - exposeState a = false
 - exposeState b = true
- Select the tank.
 - Set tank area based on a 1' pipe ($A=\pi r^2$)
 - Set level_start = 1 m, p_start = 1 bar, and h_start based on p = 1 bar and T = 50°C
 - h_start = tank.Medium.specificEnthalpy_pT(tank.p_start, 50 + 273.15)
- 4. Select the pump
 - Set m_flow_nominal = 1 kg/s
- 5. Select all components.
 - Set Medium = Modelica.Media.Water.StandardWater



```
Statistics
     (i) Original Model
        Number of components: 189
        Variables: 1641
        Constants: 62 (62 scalars)
        Parameters: 298 (357 scalars)
        Unknowns: 1281 (1761 scalars)
        Differentiated variables: 44 scalars
        Equations: 823
        Nontrivial: 628

    Translated Model

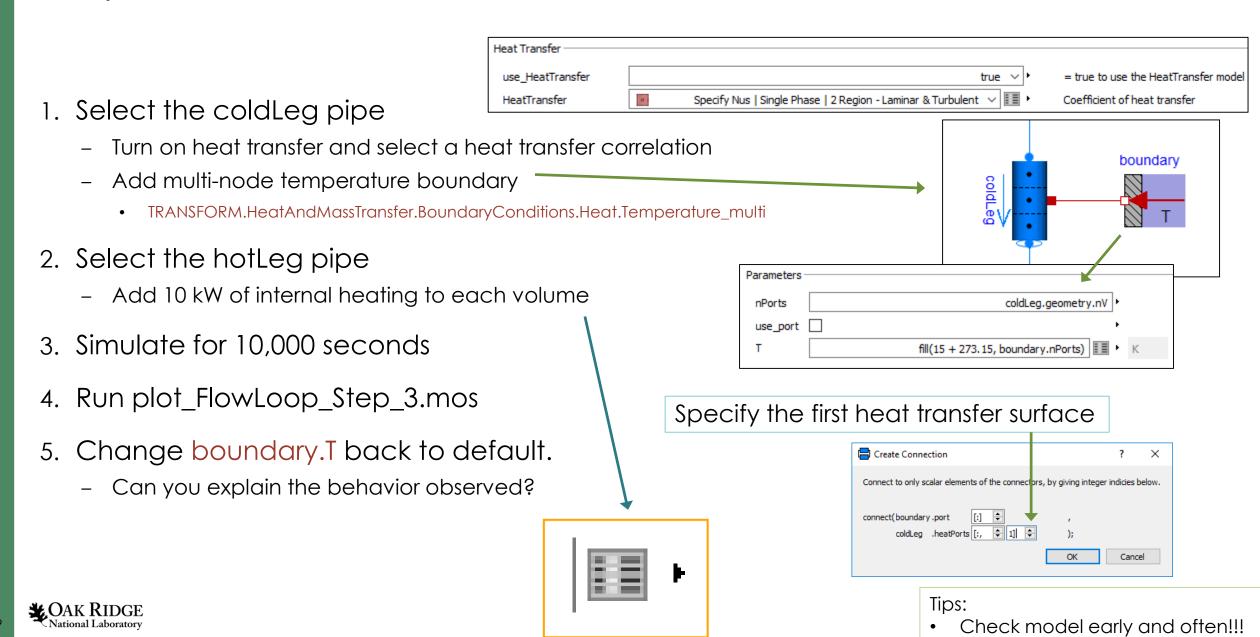
        Constants: 881 scalars
        Free parameters: 58 scalars
        Parameter depending: 102 scalars
        Continuous time states: 26 scalars
        Time-varying variables: 334 scalars
         Alias variables: 805 scalars
         Assumed default initial conditions: 18
        Number of mixed real/discrete systems of equations: 0
        Sizes of linear systems of equations: {2, 2, 2, 2, 2, 2, 2, 2, 2}
        Sizes after manipulation of the linear systems: {0, 0, 0, 0, 0, 0, 0, 0, 0}
        Sizes of nonlinear systems of equations: {}
        Sizes after manipulation of the nonlinear systems: {}
        Number of numerical Jacobians: 0
```

```
Integration started at T = 0 using integration method DASSL
(DAE multi-step solver (dassl/dasslrt of Petzold modified by Das
Integration terminated successfully at T = 100
  CPU-time for integration
                                            : 0.168 seconds
                                            : 0.336 milliseconds
  CPU-time for one grid interval
  CPU-time for initialization
                                            : 0.002 seconds
  Number of result points
  Number of grid points
                                            : 501
  Number of accepted steps
                                            : 76
  Number of f-evaluations (dynamics)
                                            : 158
  Number of crossing function evaluations : 576
  Number of Jacobian-evaluations
                                            : 54
  Number of model time events
  Number of input time events
                                            : 0
  Number of state events
                                            : 0
                                            : 0
  Number of step events
  Minimum integration stepsize
                                            : 1.69e-010
  Maximum integration stepsize
                                            : 37.6
  Maximum integration order
                                            : 1
```

Tips:

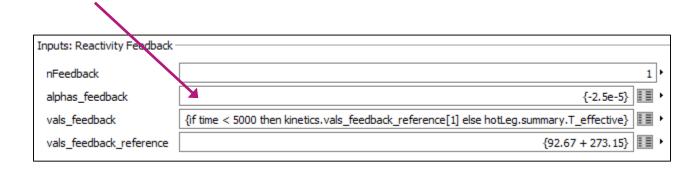
 While in code editor or in a parameter GUI, click "ctrl+SPACE" to get autocomplete

Step 3: Add heat transfer



Step 4: Add nuclear kinetics

- 1. Add a nuclear kinetics component to the model
 - TRANSFORM.Nuclear.ReactorKinetics.PointKinetics_L1_powerBased
 - Set Q_nominal = 40 kW
 - Set a temperature feedback ("kinetics" tab) based on hotLeg effective temperature
- 2. Change the internal heat generation in the hotLeg
 - Set Q_gen = kinetics.Q_total/4
- 3. Simulate. What just happened!?
- 4. Try easing the reactor with some simple logic
 - Modify vals_feedback to have a time delay
- 5. Simulate. Is the issue fixed?



Inputs: Reactivity Feedback

vals_feedback_reference

{-2.5e-5}

{92.67 + 273.15}

{hotLeg.summary.T_effective}

nFeedback

alphas_feedback

vals feedback



Tips:

• Autocomplete in the GUI only works if no other text exists beyond the cursor

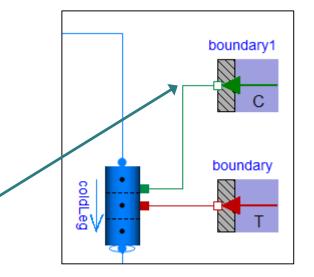
Step 5: Add and remove a trace component

model FlowLoop_Step_5 "Add nuclear heating"

package Medium = Modelica.Media.Water.StandardWater(extraPropertiesNames={"Tritium"});

- 1. Need to modify the media to indicate an extra substance
 - Add to the code side the following line
 - package Medium = Modelica.Media.Water.StandardWater(extraPropertiesNames={"Tritium"})
 - Select all fluid components and change media to "Medium"
- 2. In the hotLeg InternalTraceGen model
 - Set mC_gen = {1e-4*kinetics.Q_total} -/s
 - Trace components are unitless. "-" can mean whatever the user desires (e.g., atoms)
- 3. In the coldLeg, turn on mass transfer
 - Set mass transfer to specify alphaM
 - Set $D_ab0 = 1$ and $alpha M0 = \{1000\}$
- 4. Add a multi-node concentration boundary
 - TRANSFORM.HeatAndMassTransfer.BoundaryConditions.Mass.Concentration_multi
 - Default values are fine
- 5. Simulate. Can you see your substance in the system?
 - Search in variable browser for mC





Tips:

• The Dymola simulation search permits the use of wild cards "*". Use it!

Step 6: Change pump component

- 1. Right click the pump and select "Change Class"
- 2. Select "Pump_Controlled"
- 3. In the pump component
 - Set p_a_start = 1 bar, p_b_start = 1.1 bar
 - Set T_a_start = 50°C and m_flow_start = 1 kg/s



- 5. In the pump change "controlType" from "RPM" to "m_flow"
- 6. Simulate. What changed and why?



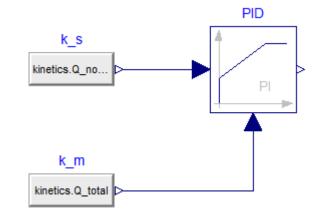


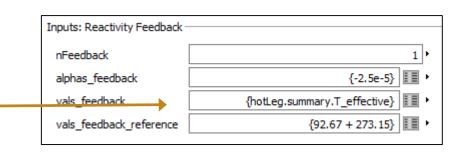
- Using GUI to change component names will update all references to it.
- Update variable using the Edit>Variables for all uses to be updated as well.



Step 7: Use controller to alter power behavior

- Drag a PID component in the model
 - TRANSFORM.Controls.LimPID
 - Set controllerType = PI
 - Set k_s and k_m = 1/kinetics.Q_nominal
 - Set k = 1e-5
- 2. Add 2 real expression inputs into the model
 - Modelica.Blocks.Sources.RealExpression
 - Connect them to the PID
 - Set the one connected to k_s equal to kinetics.Q_nominal
 - Set the one connected to k m equal to kinetics. Q total
- 3. Set kinetics.rho = PID.y
- 4. Simulate. What happened?
- 5. Remove the "if then" logic from vals_feedback
- 6. Simulate. Much better now, yes?





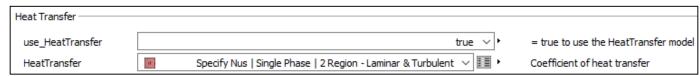
Tips:

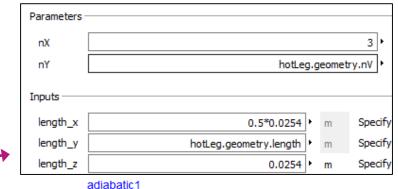
Hover over connections to get connector information

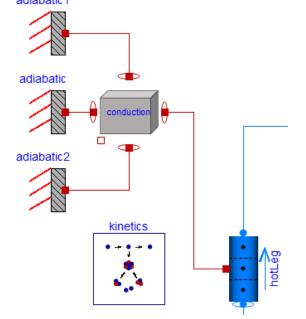


Step 8: Include a solid heating element (e.g., fuel)

- 1. Turn on heat transfer in the hotLeg
- 2. Turn off internal heat generation
- Drag a 2D Conduction component into the model
 - TRANSFORM.HeatAndMassTransfer.DiscritizedModels.Conduction 2D
 - Under "Advanced" tab set all exposeStates* = true
 - Under "Initialization" tab and set temperatures 50°C
 - Set geometry as shown
 - Under "InternalHeatModel"
 - Set Q_gen = kinetics.Q_total/(conduction.geometry.nX*conduction.geometry.nY)
 - Set Material to SS316
- 4. Drag 3 adiabatic multi-node boundary conditions into the model
 - TRANSFORM.HeatAndMassTransfer.BoundaryConditions.Heat.Adiabatic_multi
 - Set "adiabatic" nPorts = conduction.geometry.nY and the other two to nX
- 5. Connect models as shown









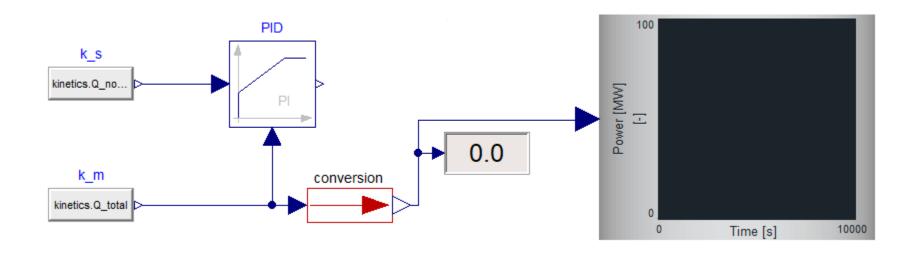
Tips:

Look into sigmoid or easing functions to smooth transitions

Step 9: Add Visual Components

- 1. Add the following to the model and connect
 - TRANSFORM.Utilities.Visualizers.DynamicGraph
 - TRANSFORM.Utilities.Visualizers.displayReal
 - TRANSFORM.Units.Conversions.Models.Conversion
- 2. For conversion, set to_kilo

- 3. In plot, set $y_max = 100$ and $t_end = 10,000$
- 4. Under the coldLeg and hotLeg components
 - Under "Visualization" tab turn on show_colors
 - Set val_min = 40°C and val_max = 100°C
 - Careful with the units!





Step 10: Add Realtime User Input

- Add a user input to the model
 - UserInteraction.Inputs.NumericInputIO
- Add the input value to the kinetic model
- 3. Switch to the simulation view
- 4. Open the model viewer
- 5. Under simulation settings
 - "Realtime" tab set settings
 - "Compiler" tab select DDE server
- 6. Simulate and plot kinetics.Q_total and kinetics.rho
- 7. Change the input value and press enter. What happened?

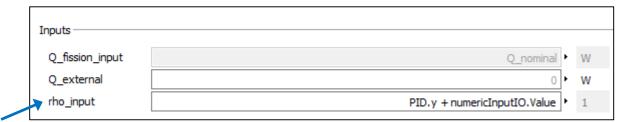
Realtime simulation

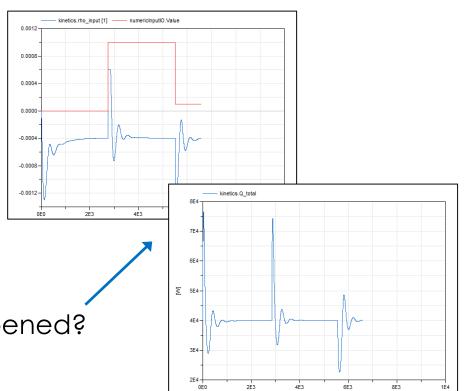
Slowdown factor

✓ Synchronize with realtime

Load result interval [s] 0.5

0.01

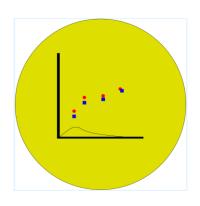






Step 11: Add Unit Test

- Add unit test checker to model
 - TRANSFORM.Utilities.ErrorAnalysis.UnitTests
- 2. In unitTest parameter GUI assign one or more variables to x
- 3. Simulate to ensure everything is working
- 4. Open python/Spyder and
 - Run createUnitScripts.py
 - Run regTestsWin_customBuildPy.py
 - May require custom buildingspy regression script from ModelicaPy (see Session 4)
 - Select "y" if prompted
- 5. Check the Resources folder. What happened?



Step 12: Extend/Modify TRANSFORM – Heat Transfer

- 1. Create a new package in TRANSFORM_Training_2019 called HeatTransfer
- Duplicate the used correlation into the new HeatTransfer package with a modified name and description
 - TRANSFORM.Fluid.ClosureRelations.HeatTransfer.Models.DistributedPipe_1D_MultiTransferSurface.Nus_SinglePhase_2Region
- 3. Replace Nu_DittusBoelter with Nu_SiederTate
 - Check the heat transfer model. What is missing?
- 4. Change the hotLeg correlation to the new correlation
- 5. Simulate. What changed?
 - Use the compare feature to view the difference in effective solid temperature
 - conduction.summary.T_effective



Step 13: Read results in Python/Matlab

Python

- Example in TRANSFORM Resources folder
- Install buildingspy: pip install buildingspy
- Change directory to location of '.mat'
- Create notebook or python file and add:

from buildingspy.io.outputfile import Reader

r = Reader("MATFILE.mat','dymola')

time, variable = r.values('VARNAME')

Matlab

- Dymola provides useful functions
 - C:\Program Files\Dymola 2020\Mfiles\dymtools
- Example found in Resources folder
- Grab dymget.m and dymload.m from
 - C:\Program Files\Dymola 2020\Mfiles\dymtools

% Load the .mat file. If not found see 'err' for error [dymstr,err] = dymload(fullPath_mat);

% Get time variable. Only need to load once as it is the same for all variables t = dymget(dymstr,'Time');



Step 14: Create an FMU

- Generate the FMU from the Dymola GUI
- 2. Open the FMU in FMU Simulator
 - FMPy Github (https://github.com/CATIA-Systems/FMPy)
 - Windows Executable: https://www.3ds.com/productsservices/catia/products/dymola/free-downloads/fmu-simulator/
- 3. Simulate to ensure everything is working



Free Time

- 1. Explore capabilities. Try things such as:
 - Adding time dependent signals
 - Make variables dependent on other variables
 - Create summary of variables in the code side...
 - ... or create a "summary" model that you can drag in to populate
 - Add more trace components
 - Add a source/sink for another fluid stream
 - Try to remove the trace component using an ideal sink
 - Try a difference geometry for the conduction model



Thank you.

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