

# *Process Simulation Using Aspen HYSYS V8*

**Experience the New Aspen HYSYS®.**

V8

The best process simulation software is now easier to use and faster to learn than ever!

Aspen HYSYS is a comprehensive process modeling system used by the world's leading oil & gas producers, refineries, and engineering companies to optimize process design and operations.



Why Aspen HYSYS?

***Eng. Ahmed Deyab Fares  
Process Simulation Consultant***

Aspen HYSYS is a market-leading process modeling tool for conceptual design, optimization, business planning, asset management, and performance monitoring for oil & gas production, gas processing, petroleum refining, and air separation industries. Aspen HYSYS is a core element of AspenTech's aspenONE™ Process Engineering applications.

## Objectives

- Learn to build, navigate and optimize process simulations using Aspen HYSYS
- Learn the efficient use of different HYSYS functions to build steady state process simulations

## Who Should Attend

- New engineering graduates/technologists who will be using Aspen HYSYS in their daily work
- Process engineers doing process design and optimization projects and studies
- Plant engineers checking plant performance under different operating conditions
- R&D engineers and researchers using Aspen HYSYS for process synthesis

## Prerequisites

- A background in chemical engineering or industrial chemistry

## **Content:**

- ✓ *Getting Started*
- ✓ *Propane Refrigeration Loop*
- ✓ *Refrigerated Gas Plant*
- ✓ *Oil Characterization*
- ✓ *Pre-Heat Train*
- ✓ *Atmospheric Towers & Side operations*
- ✓ *Gas Gathering*
- ✓ *Optimization*
- ✓ *NGL Fractionation Train*
- ✓ *Oil Stabilization Optimization*

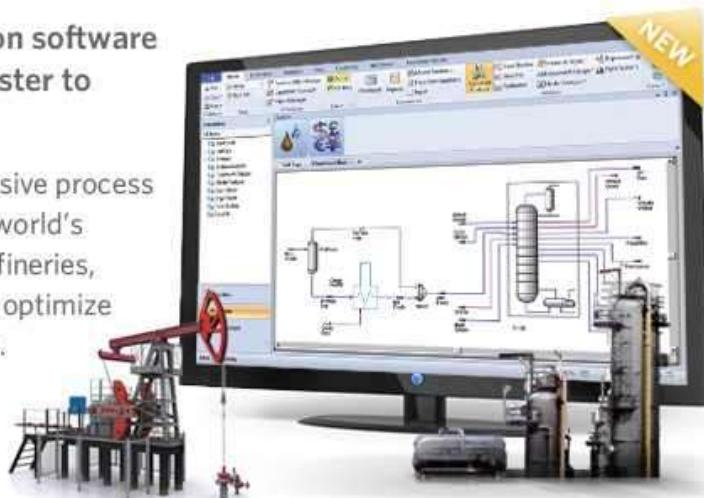
# Getting Started

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1

## ***Workshop***

The Getting Started module introduces you to some of the basic concepts necessary for creating simulations in HYSYS. You will use HYSYS to define three gas streams to be used as feeds to a gas plant. In addition, you will learn how to determine properties of these streams by using the Phase Envelope and the Property Table utilities.

## ***Learning Objectives***

- Define a fluid package (property package, components, hypotheticals).
- Add streams.
- Understand flash calculations.
- Attach stream utilities.
- Customize the Workbook.

Example:

We have a stream containing 15% ethane, 20% propane, 60% i-butane and 5% n-butane at 50°F and atmospheric pressure, and a flow rate of 100lbmole/hr. This stream is to be compressed to 50 psia and then cooled to 32°F. The resulting vapor and liquid are to be separated as the two product streams. Neglect the pressure drop inside the condenser.

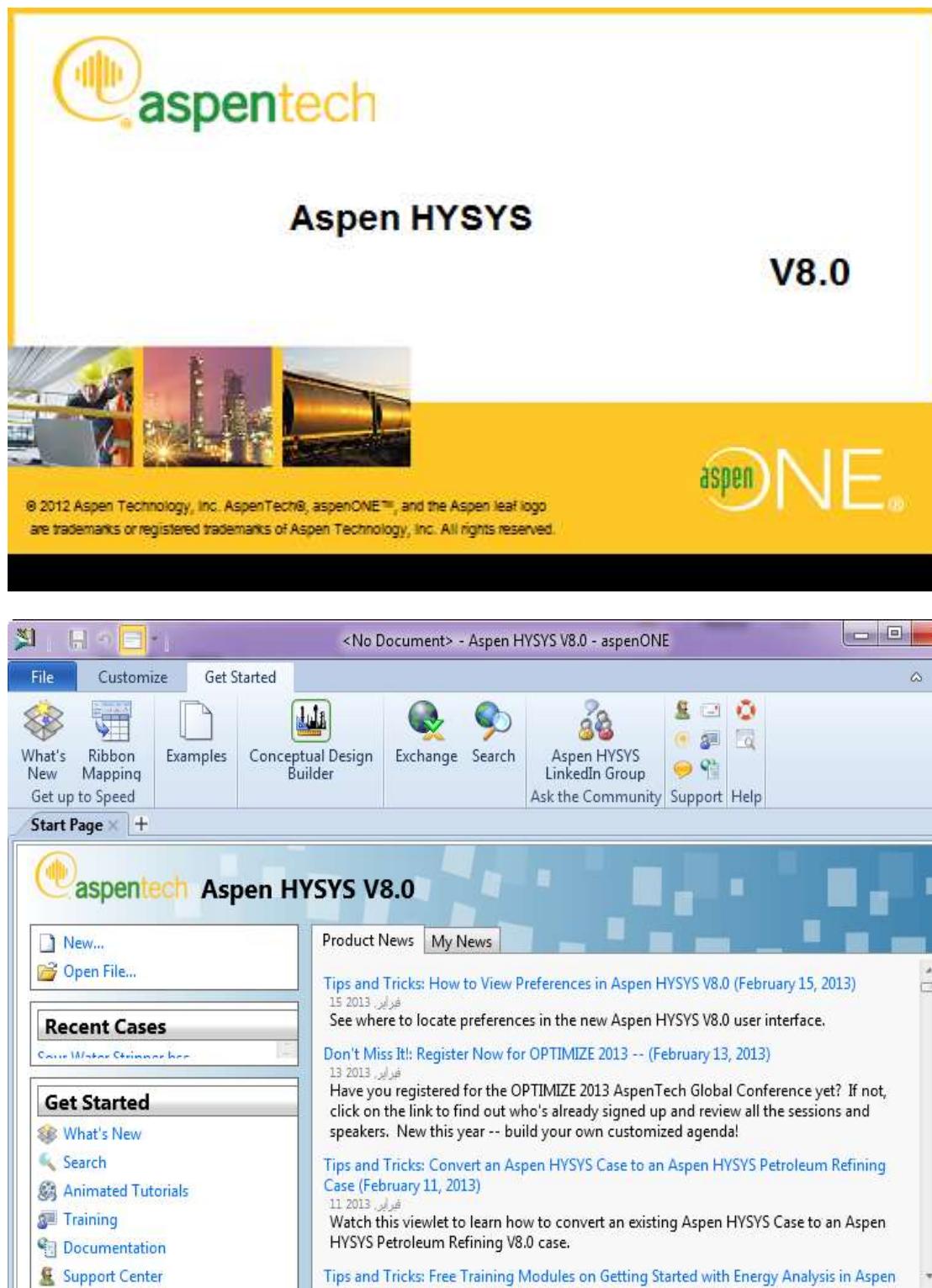
- Fluid pkg: Peng Robinson

\* What are the flow rates and molar compositions of two product streams?

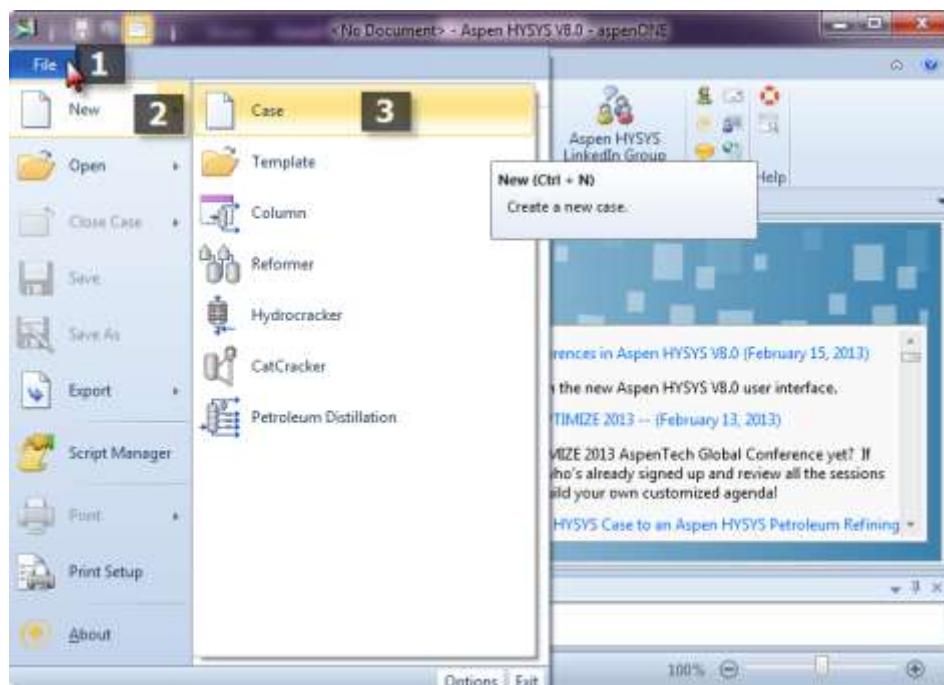
Component	Vapor	Liquid
Ethane		
Propane		
I-Butane		
N-Butane		
Total Flow rate		

\* Create a case study to see the effect of changing temperature of the cooler out stream on the molar flow of the liquid product stream, and write your comment.

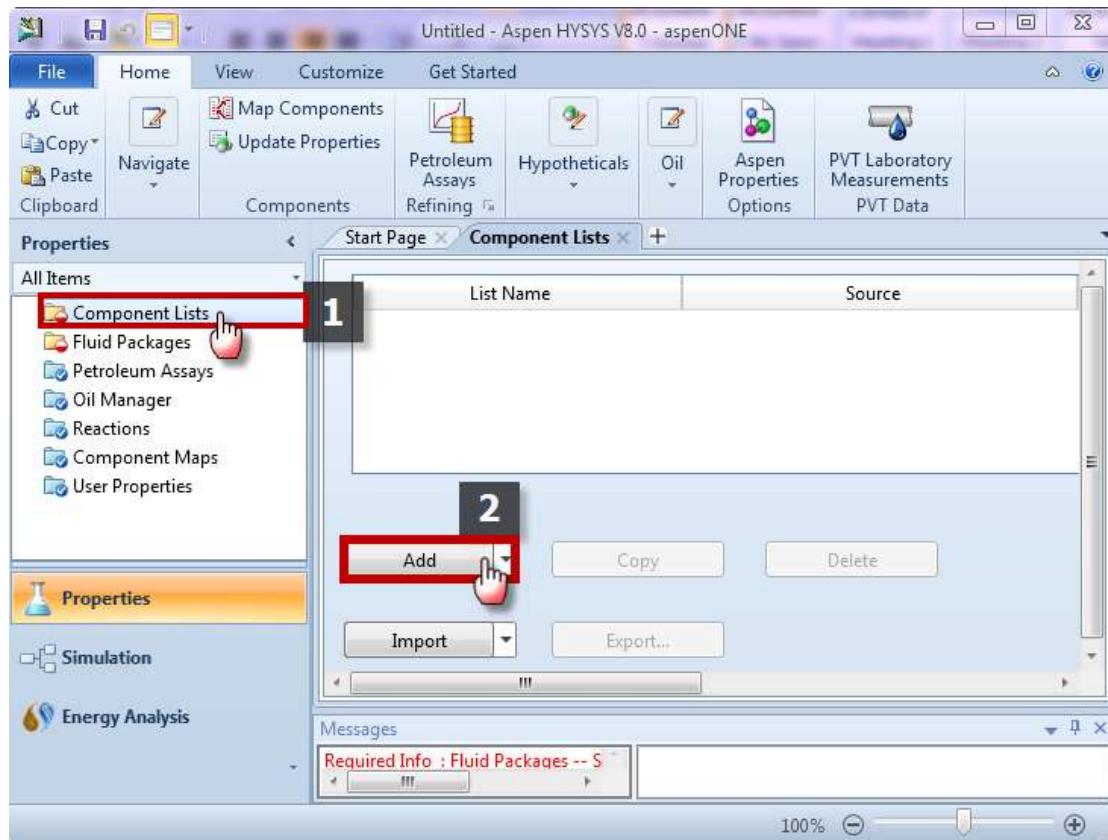
To start the program, From Start Menu, Select All Programs >>  
Aspen Tech >> Process Modeling V8.x >>> Aspen HYSYS >>  
Aspen HYSYS



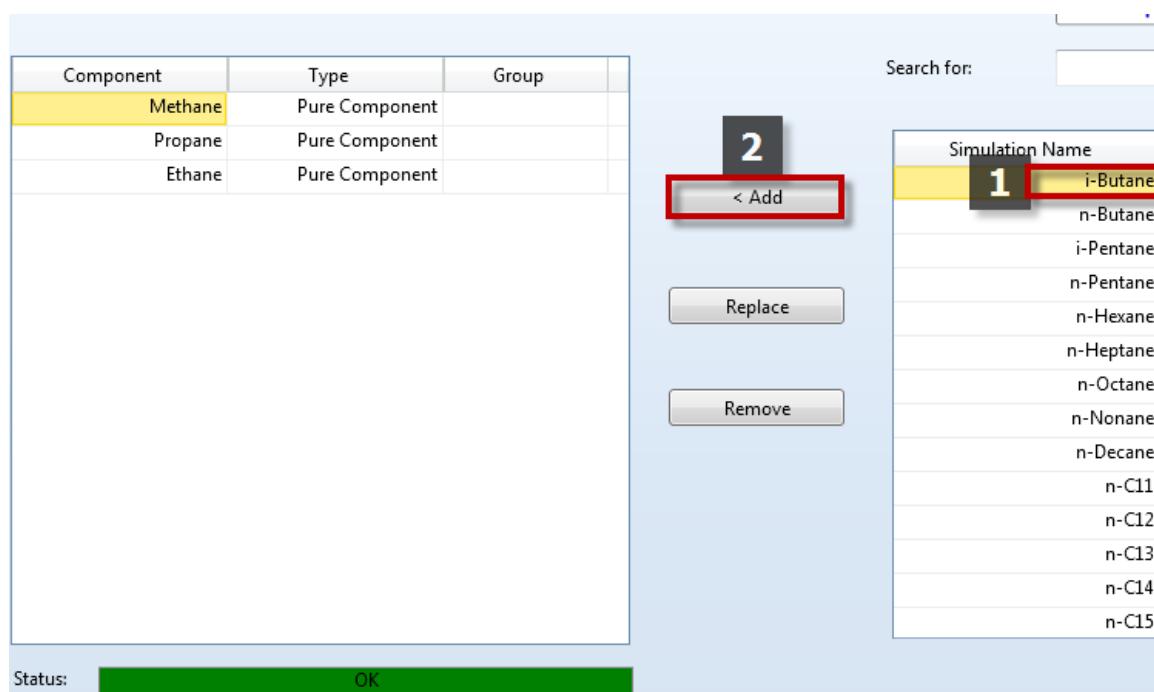
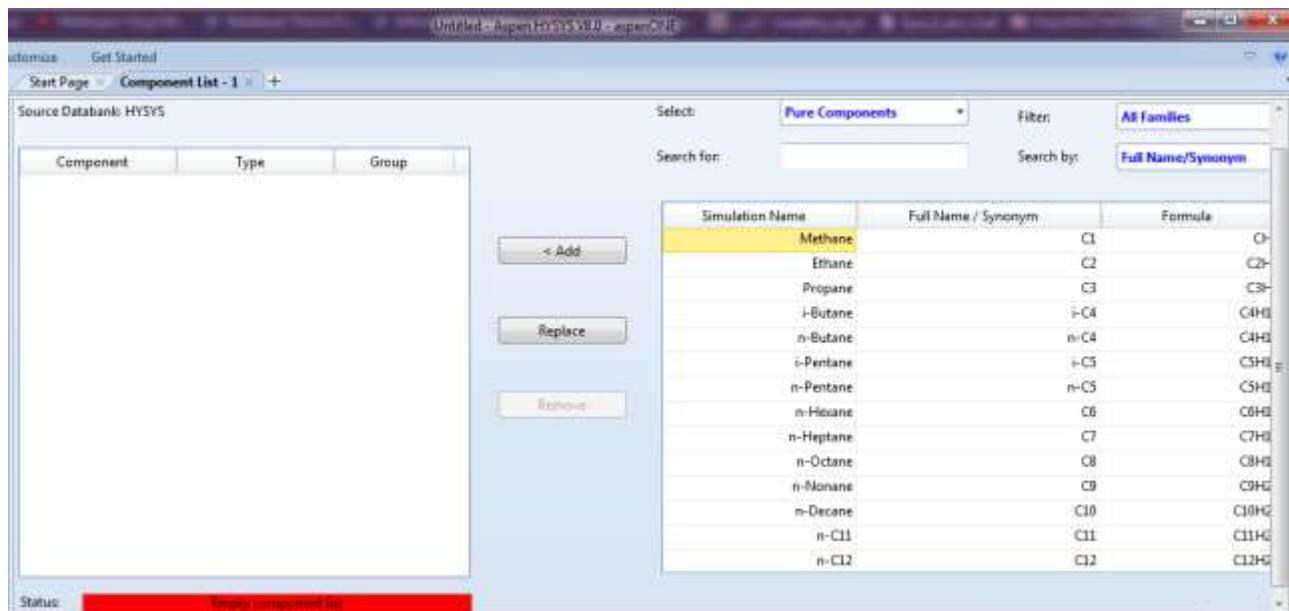
## 1- First, Start a new case



## 2- Add the Components



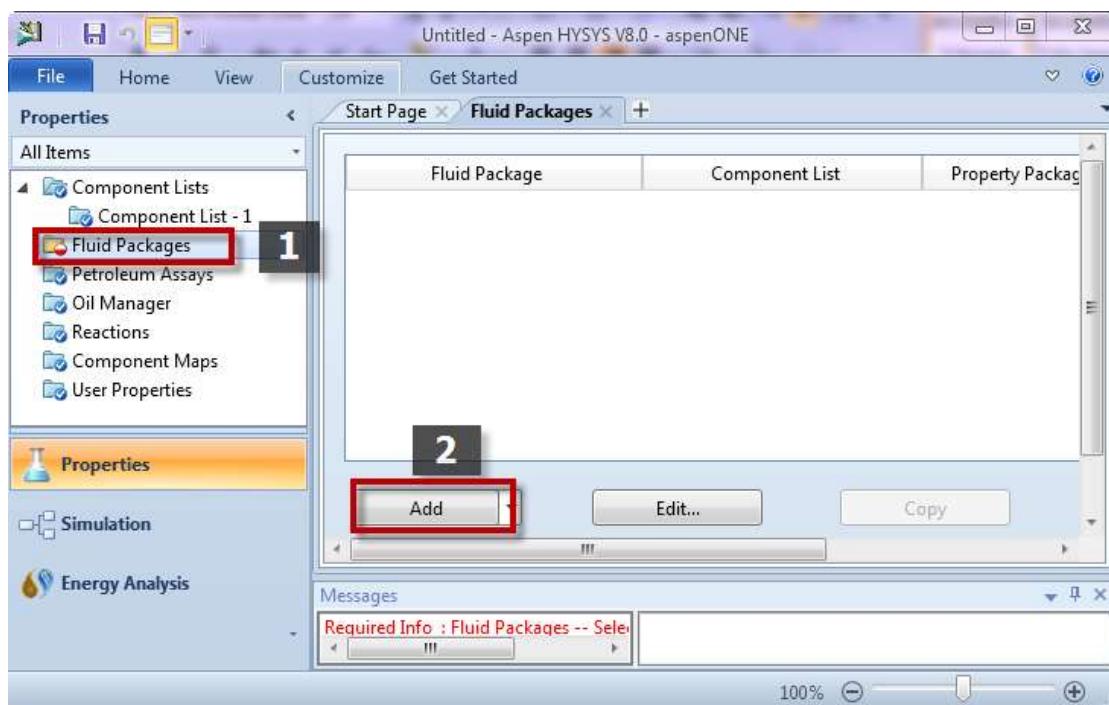
### 3- Choose the system components from the databank:



Now, select the suitable fluid package

When you have established a component list, you combine the component list with a property package. The property package is a collection of methods for calculating the properties of the selected components. The

combination of the component list and the property package, along with other simulation settings, is called the fluid package.



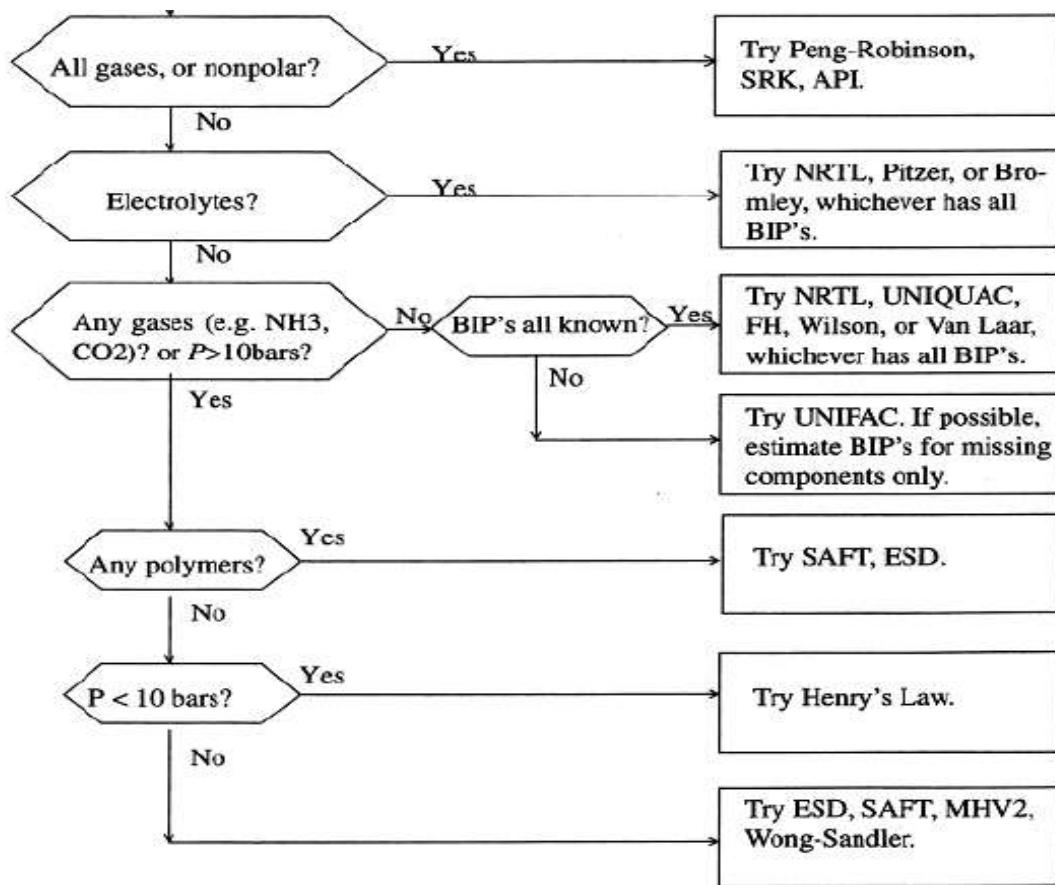
The built-in property packages in HYSYS provide accurate thermodynamic, physical and transport property predictions for hydrocarbon, non-hydrocarbon, petrochemical and chemical fluids.

The database consists of an excess of 1500 components and over 16000 fitted binary coefficients. If a library component cannot be found within the database, a comprehensive selection of estimation methods is available for creating fully defined hypothetical components.

There are about 33 property packages inside HYSYS database; the question now is **HOW TO SELECT THE SUITABLE FLUID PACKAGE?**

We can select the suitable one by specifying:

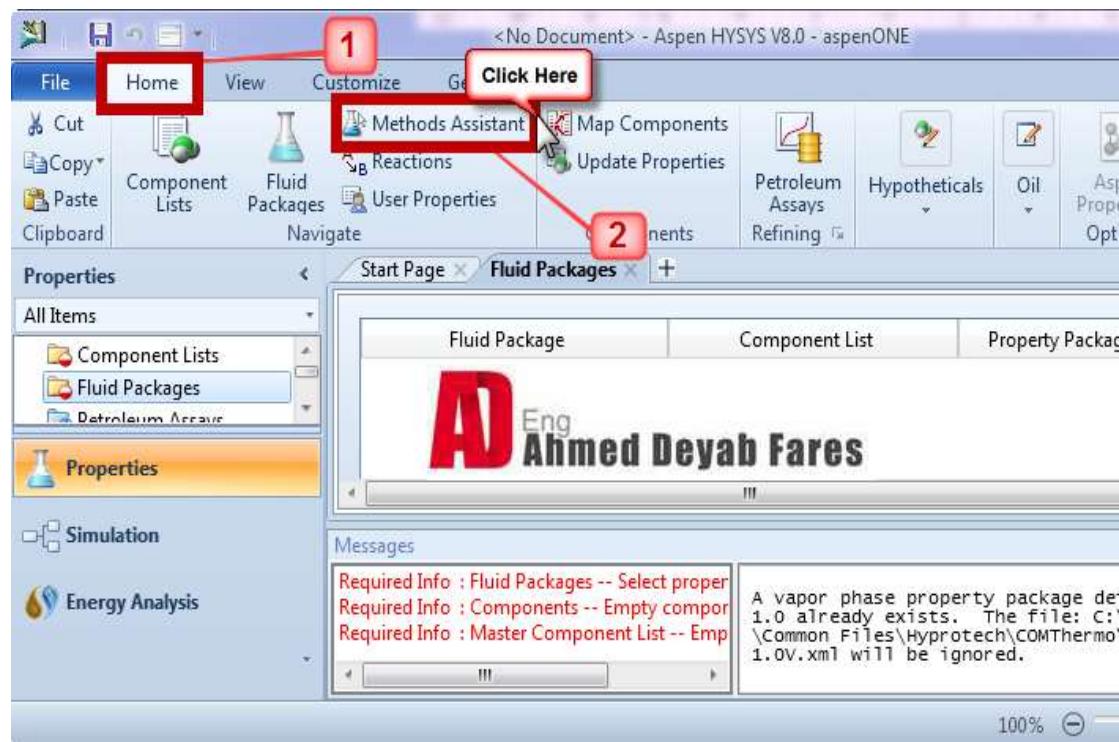
- 1- Process / Application type
- 2- Temperature and Pressure Range



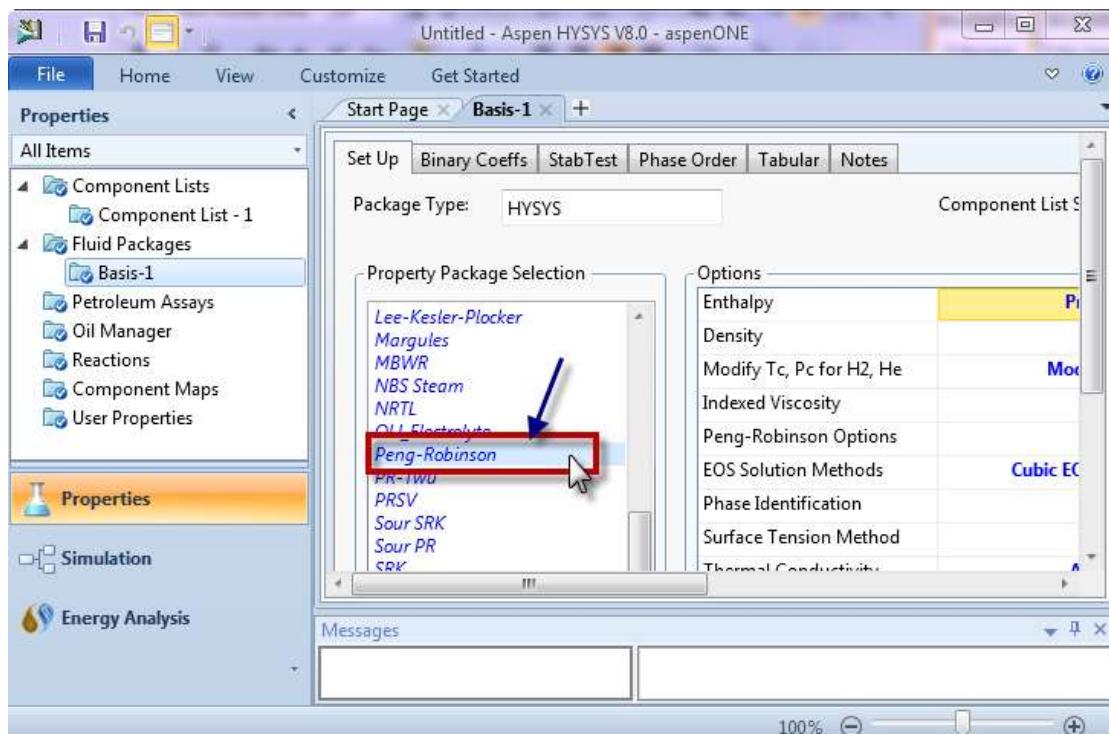
Type of System	Recommended Property Method
TEG Dehydration	PR
Sour Water	PR, Sour PR
Cryogenic Gas Processing	PR, PRSV
Air Separation	PR, PRSV
Atm. Crude Towers	PR, PR Options, GS
Vacuum Towers	PR, PR Options, GS (<10 mmHg), Braun K10, Esso K
Ethylene Towers	Lee Kesler Plocker
High H <sub>2</sub> Systems	PR, ZJ or GS
Reservoir Systems	Steam Package, CS or GS
Hydrate Inhibition	PR
Chemical Systems	Activity Models, PRSV
HF Alkylation	PRSV, NRTL

TEG Dehydration with Aromatics	PR
Hydrocarbon systems where H <sub>2</sub> O solubility in HC is important	Kabadi Danner
Systems with select gases and light HC	MBWR

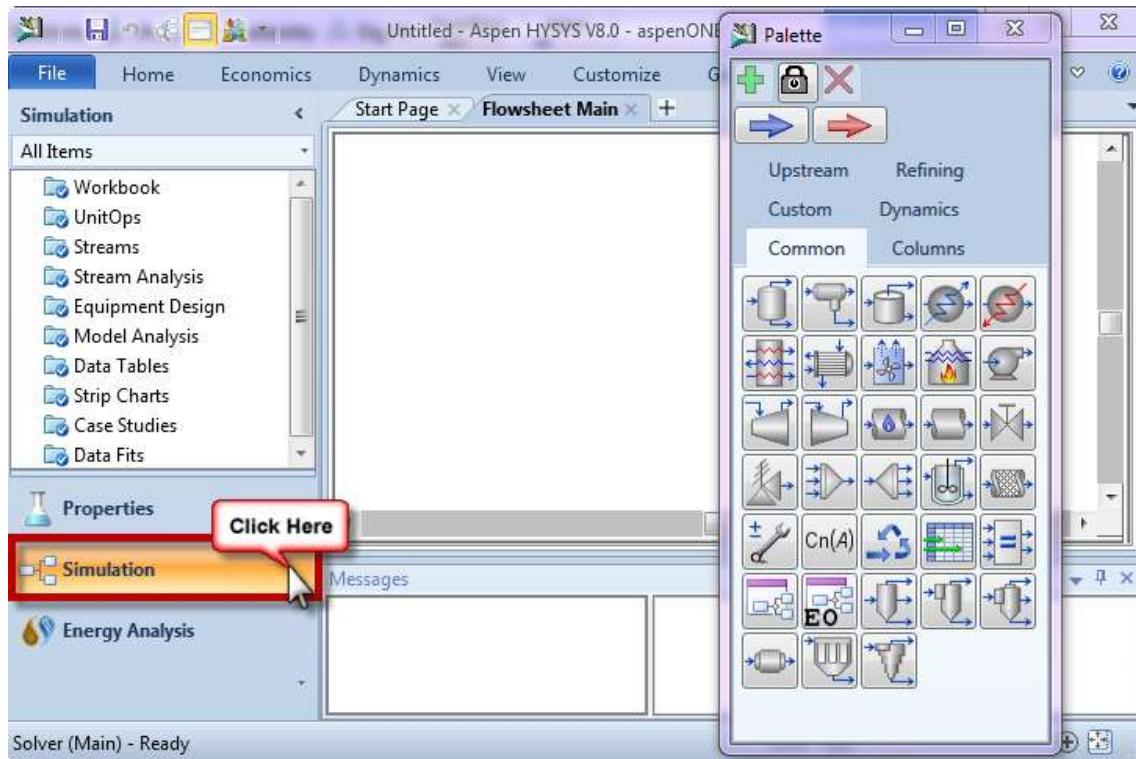
ASPEN HYSYS contains an assistant tool to help you in the selection of the suitable FP, called Methods Assistant:



In this case, select Peng-Robinson

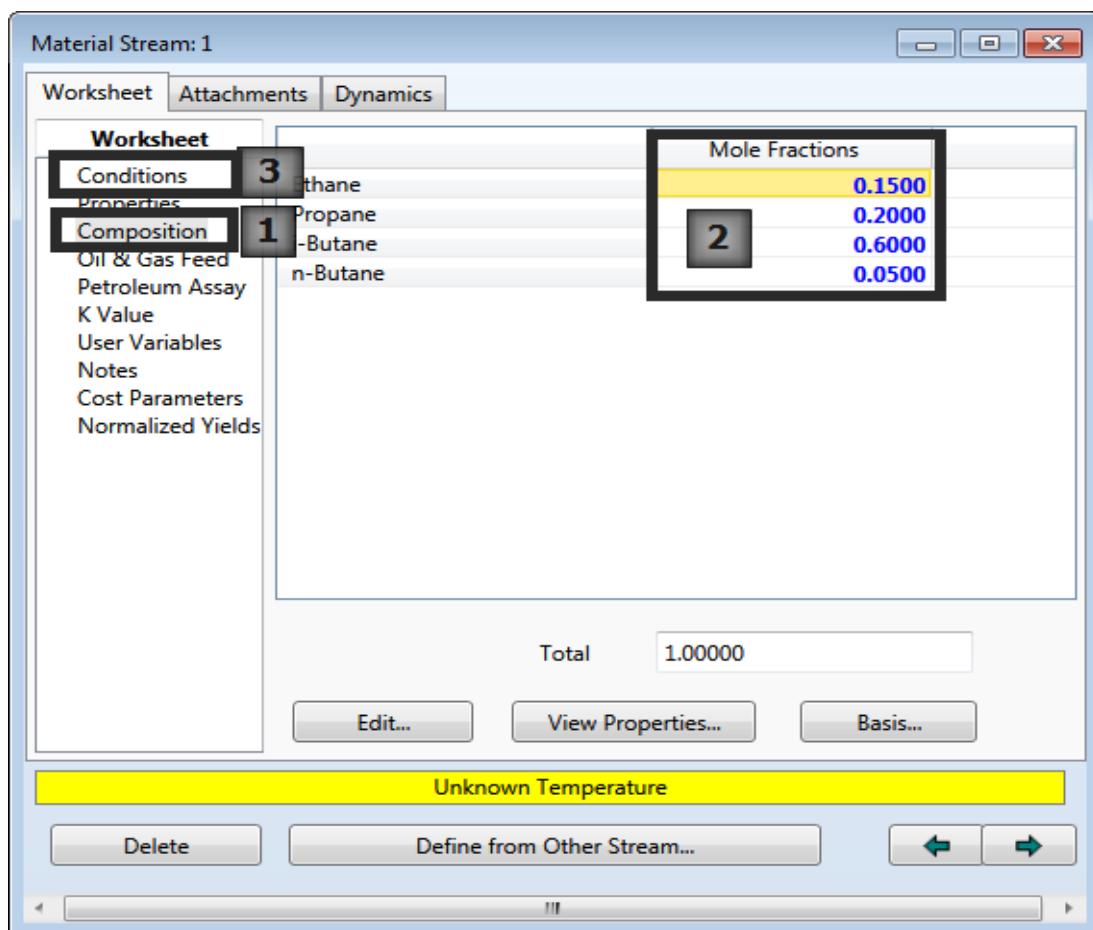
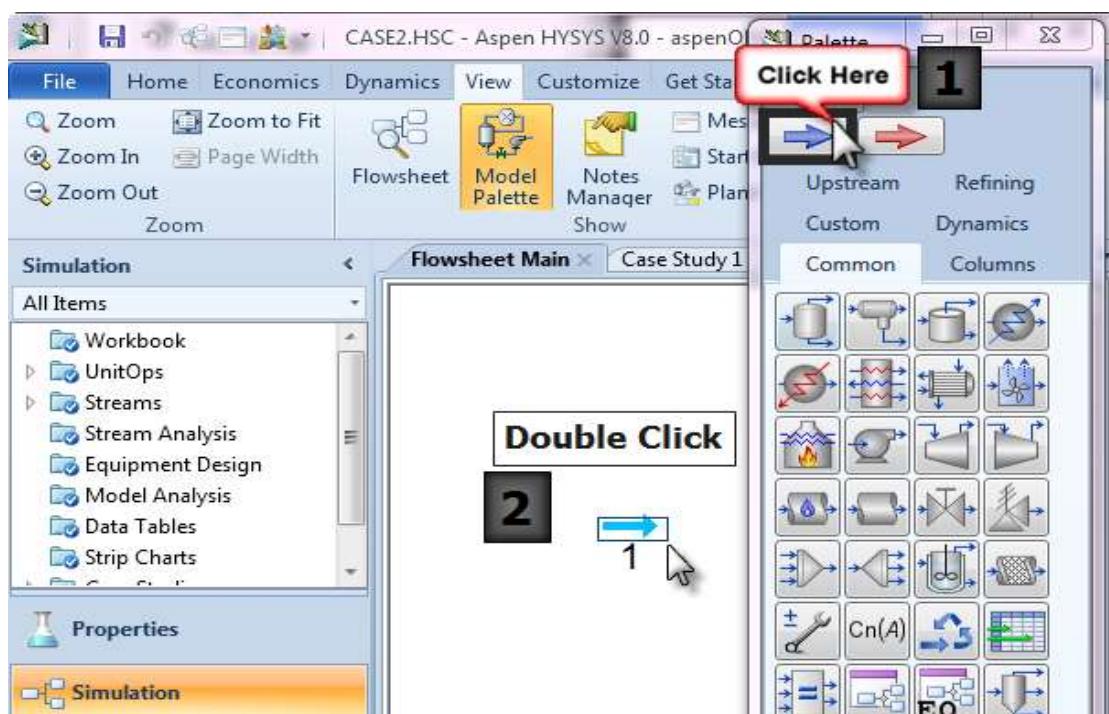


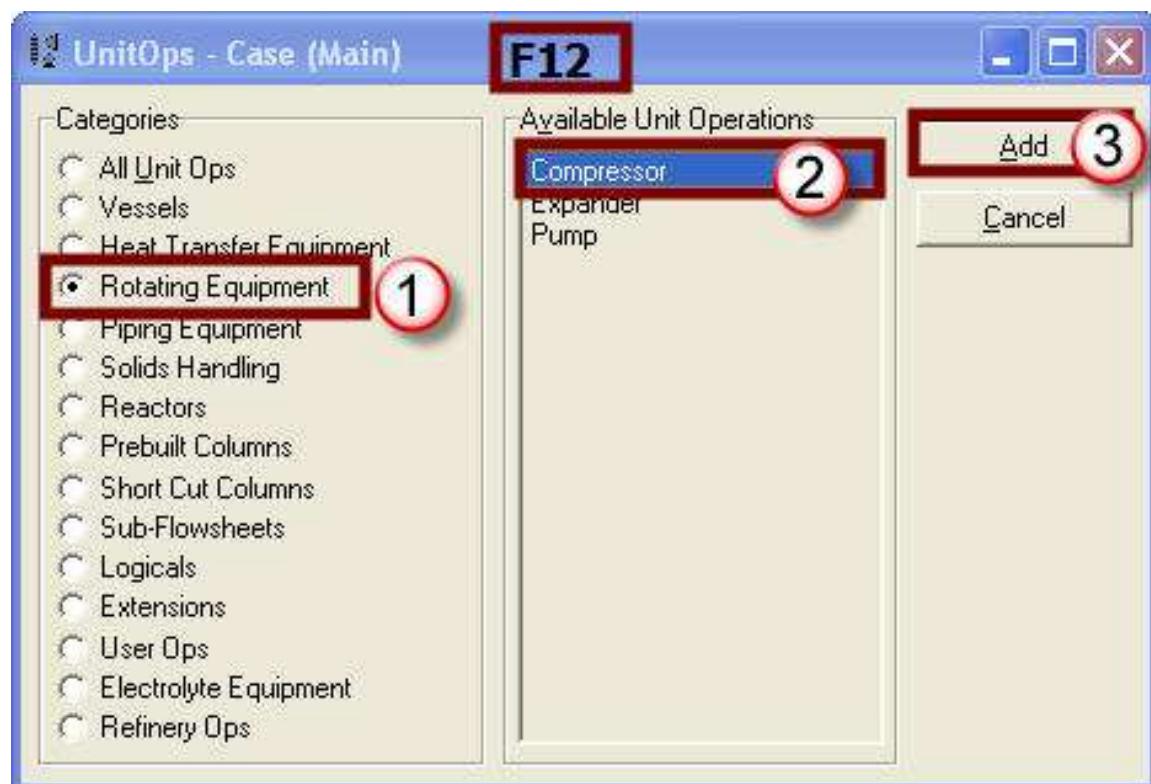
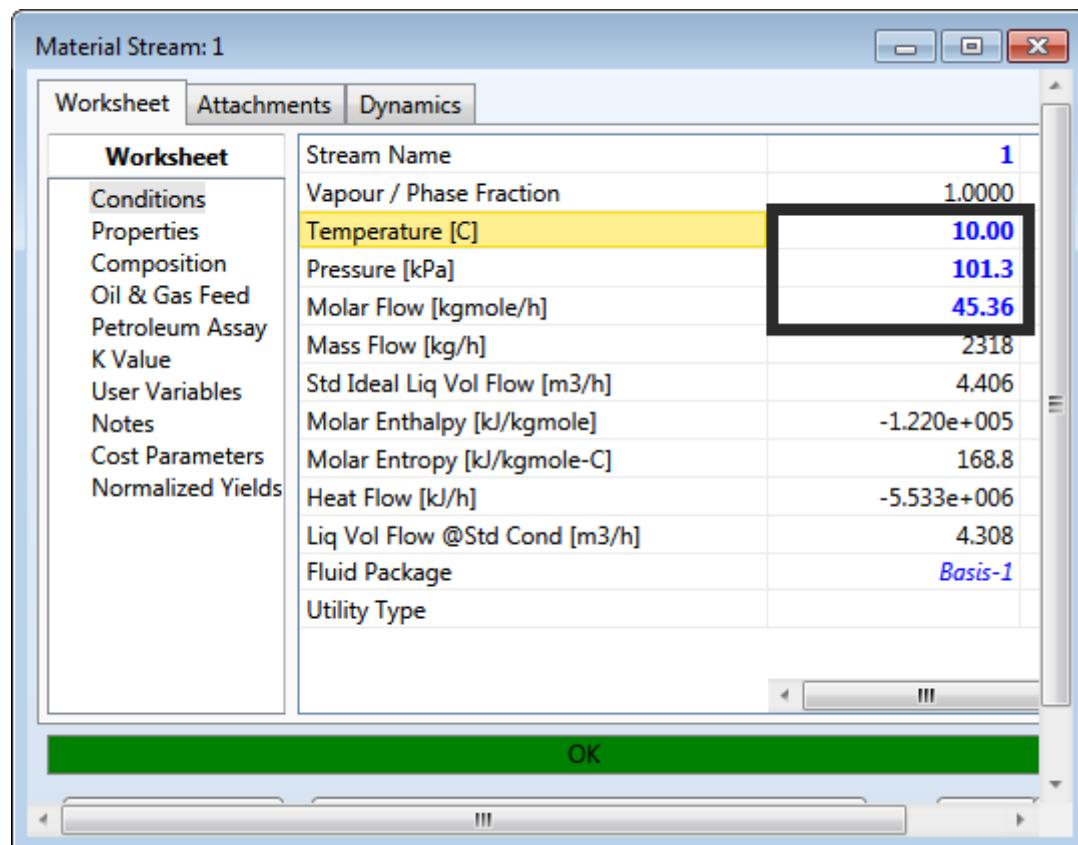
Now you can start drawing the flow sheet for the process by clicking the Simulation button:

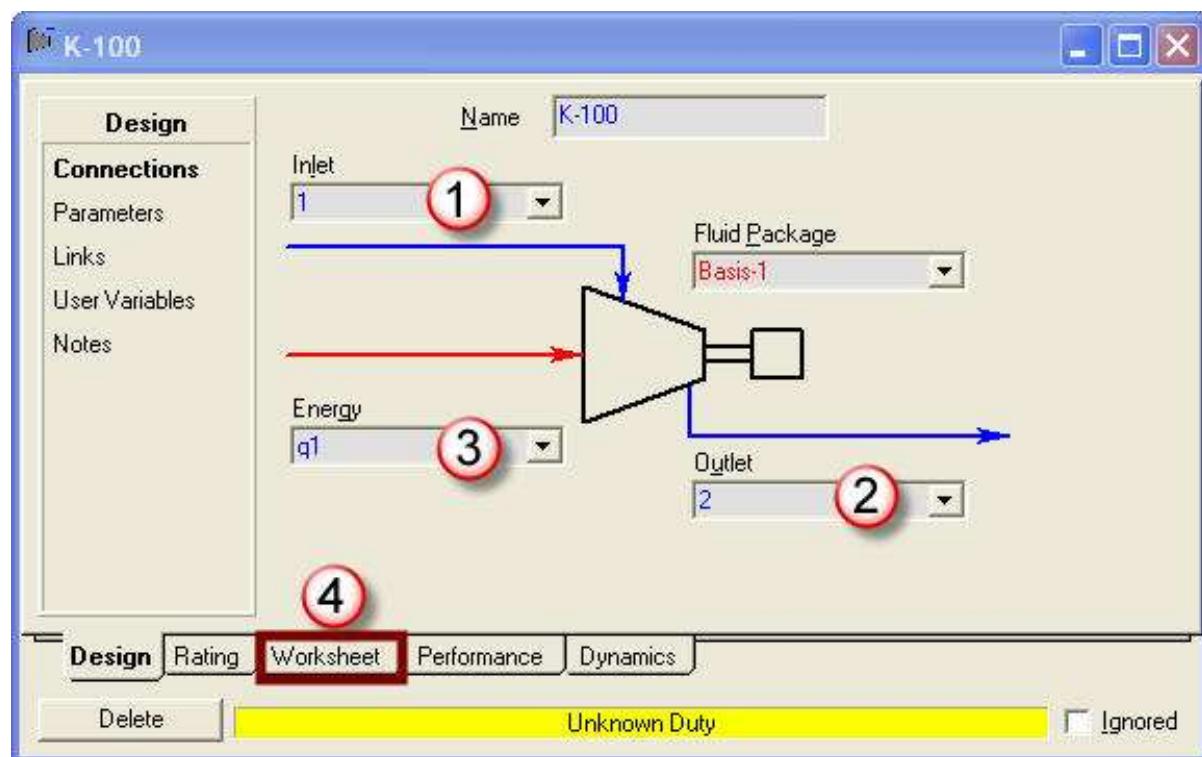


Now add a material stream to define the composition and the conditions of the feed stream

From the palette:







**K-100**

**Worksheet** tab is selected.

Name	1	2	q1
Vapour	1.0000	<empty>	<empty>
Temperature [C]	10.00	<empty>	<empty>
Pressure [kPa]	101.350	50	kPa
Molar Flow [kgmole/h]	45.36	45.36	kg/cm <sup>2</sup>
Mass Flow [kg/h]	2318	2318	psia
LiqVol Flow [m <sup>3</sup> /h]	4.406	4.406	lb/rft <sup>2</sup>
Molar Enthalpy [kJ/kgmole]	-1.220e+005	<empty>	torr
Molar Entropy [kJ/kgmole-C]	168.8	<empty>	mmHg(0C)
Heat Flow [kJ/h]	-5.533e+006	<empty>	inHg(32F)
			inHg(60F)

Buttons: Design, Rating, Worksheet, Performance, Dynamics.

Toolbar: Delete, Unknown Duty, Ignored.

K-100

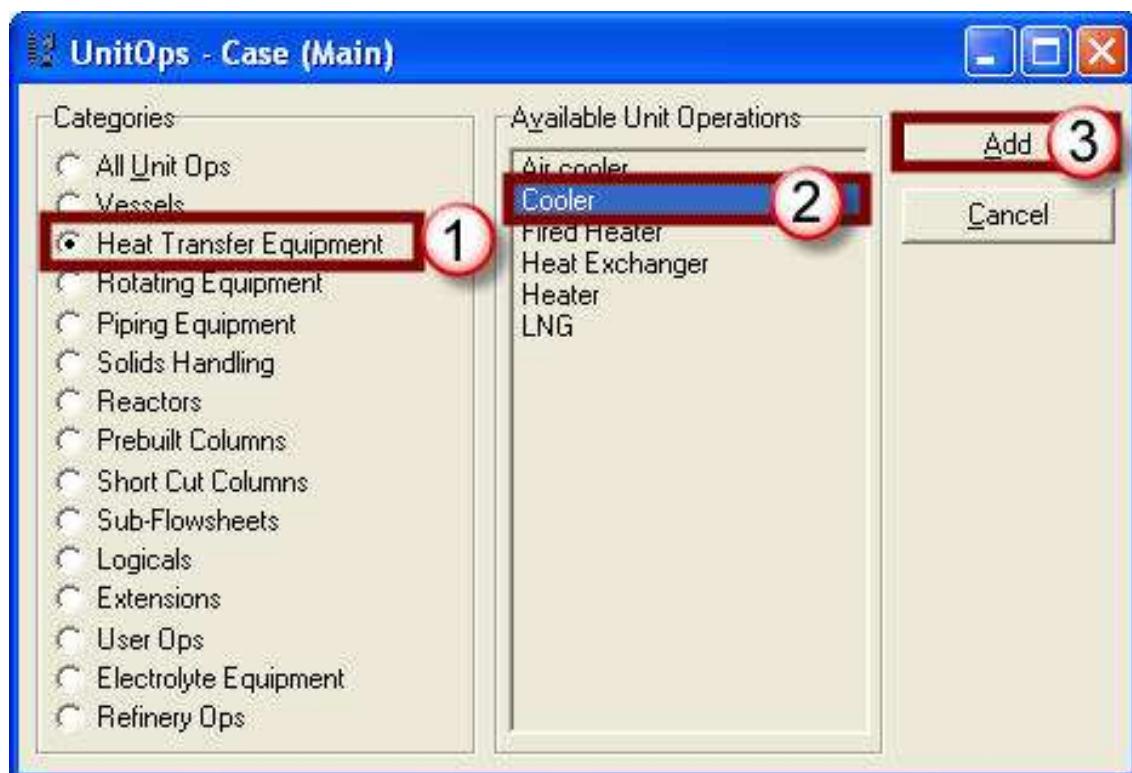
**Worksheet**

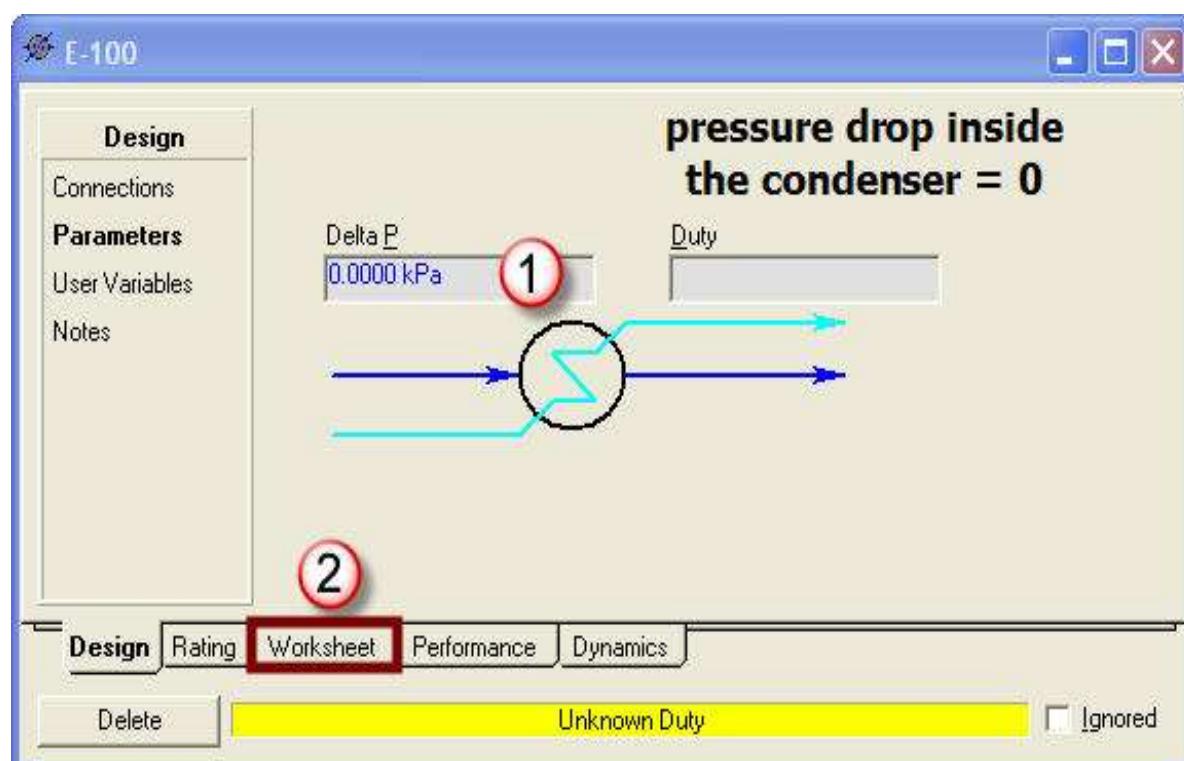
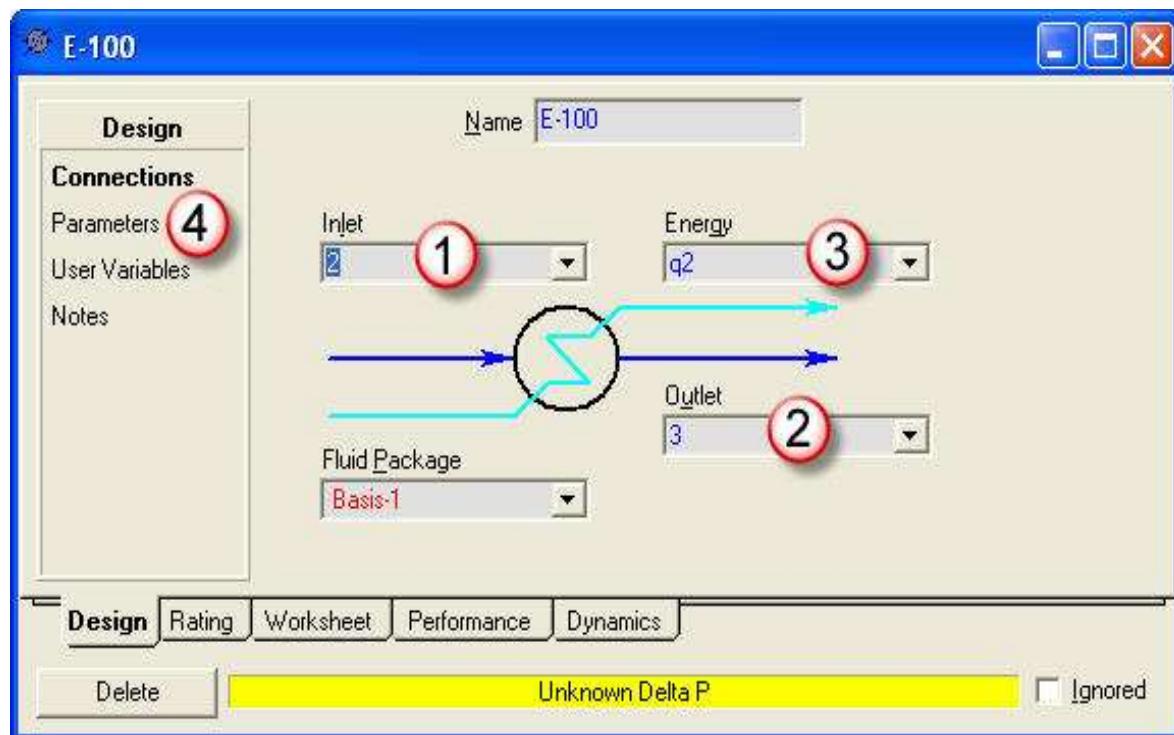
**Conditions**

Name	1	2	q1
Vapour	1.0000	1.0000	<empty>
Temperature [C]	10.00	57.59	<empty>
Pressure [kPa]	101.3	344.7	<empty>
Molar Flow [kgmole/h]	45.36	45.36	<empty>
Mass Flow [kg/h]	2318	2318	<empty>
LiqVol Flow [m <sup>3</sup> /h]	4.406	4.406	<empty>
Molar Enthalpy [kJ/kgmole]	-1.220e+005	-1.180e+005	<empty>
Molar Entropy [kJ/kgmole-C]	168.8	171.8	<empty>
Heat Flow [kJ/h]	-5.533e+006	-5.354e+006	1.783e+005

Design Rating **Worksheet** Performance Dynamics

Delete OK Ignored





E-100

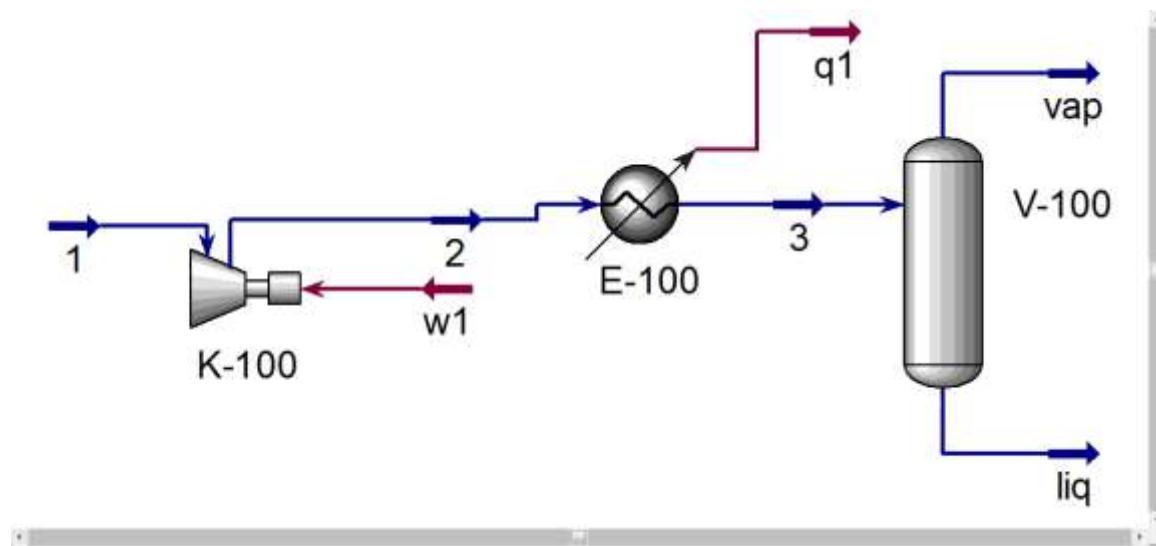
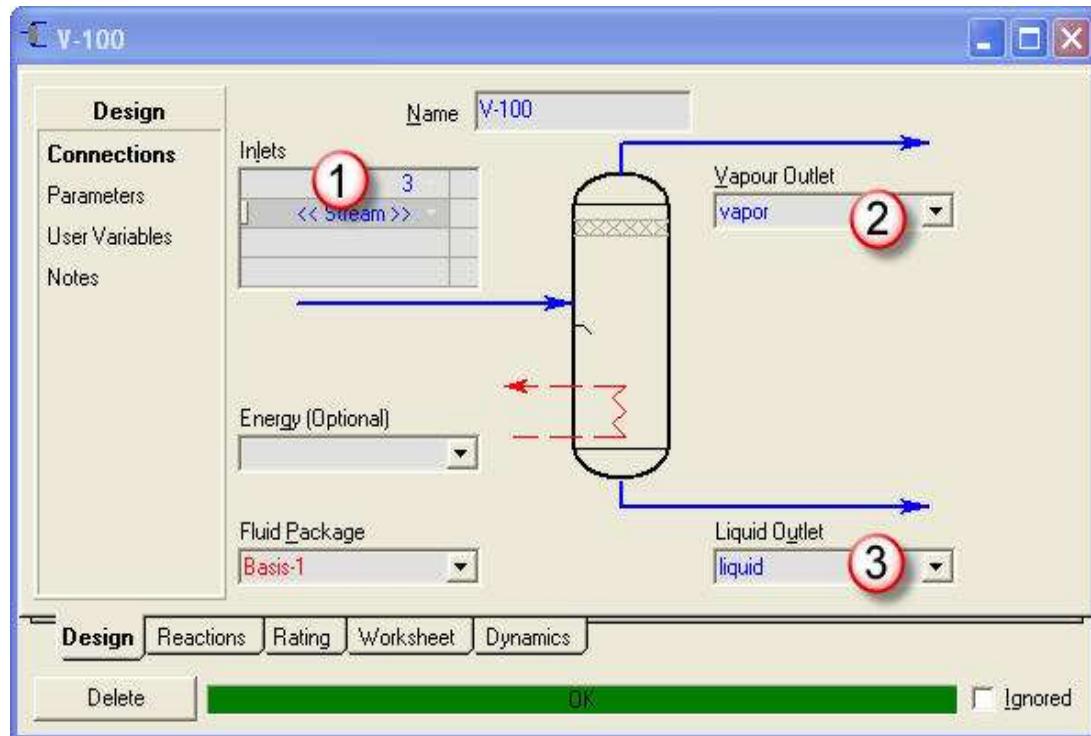
	Name	2	3	q2
Vapour	1.0000	<empty>	<empty>	
Temperature [C]	57.59	32	C	
Pressure [kPa]	344.7	344.7	C	
Molar Flow [kgmole/h]	45.36	45.36	K	
Mass Flow [kg/h]	2318	2318	F	
Std Ideal Liq Vol Flow [m <sup>3</sup> /h]	4.406	4.406	L	
Molar Enthalpy [kJ/kgmole]	-1.180e+005	<empty>	<empty>	
Molar Entropy [kJ/kgmole-C]	171.8	<empty>	<empty>	
Heat Flow [kJ/h]	-5.354e+006	<empty>	<empty>	

Design Rating Worksheet Performance Dynamics

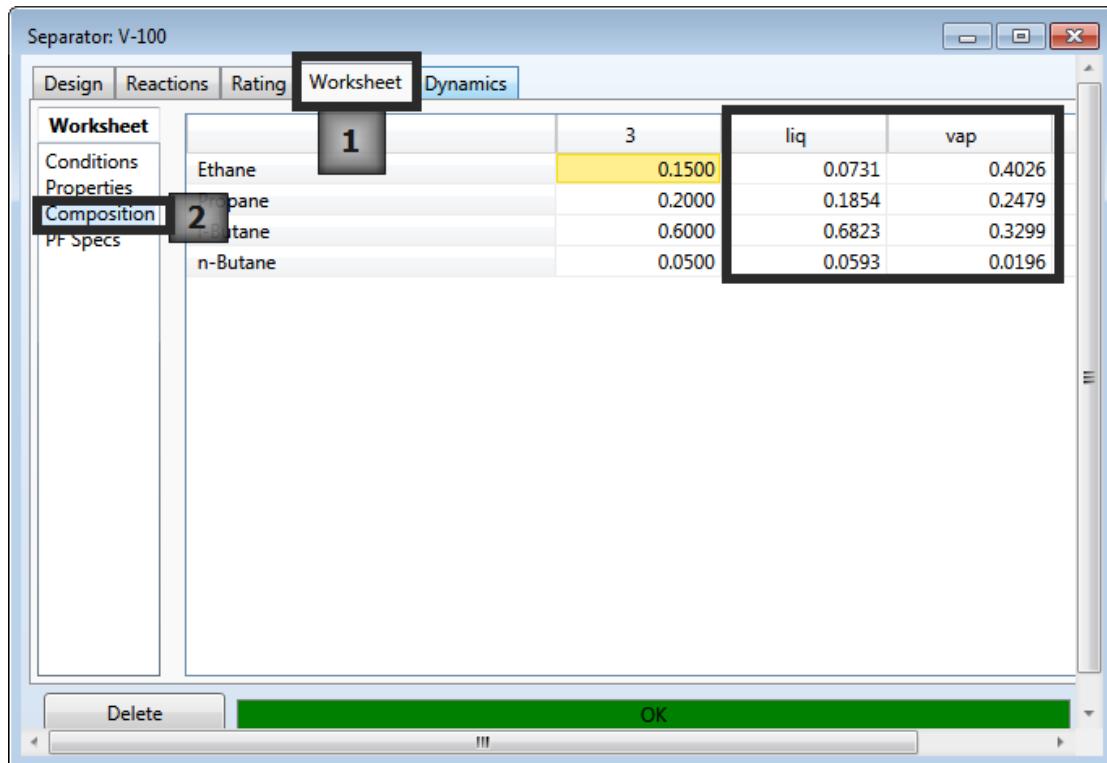
Delete Unknown Duty  Ignored

UnitOps - Case (Main)

Categories	Available Unit Operations	Add
<input checked="" type="radio"/> All Unit Ops <input checked="" type="radio"/> Vessels <input type="radio"/> Heat Transfer Equipment <input type="radio"/> Rotating Equipment <input type="radio"/> Piping Equipment <input type="radio"/> Solids Handling <input type="radio"/> Reactors <input type="radio"/> Prebuilt Columns <input type="radio"/> Short Cut Columns <input type="radio"/> Sub-Flowsheets <input type="radio"/> Logicals <input type="radio"/> Extensions <input type="radio"/> User Ops <input type="radio"/> Electrolyte Equipment <input type="radio"/> Refinery Ops	3 Phase Separator Cont. Stirred Tank Reactor Conversion Reactor Equilibrium Reactor Gibbs Reactor <b>Separator</b> Tank	Add Cancel



Now you can view the results by double clicking on the separator, in the worksheet tab:



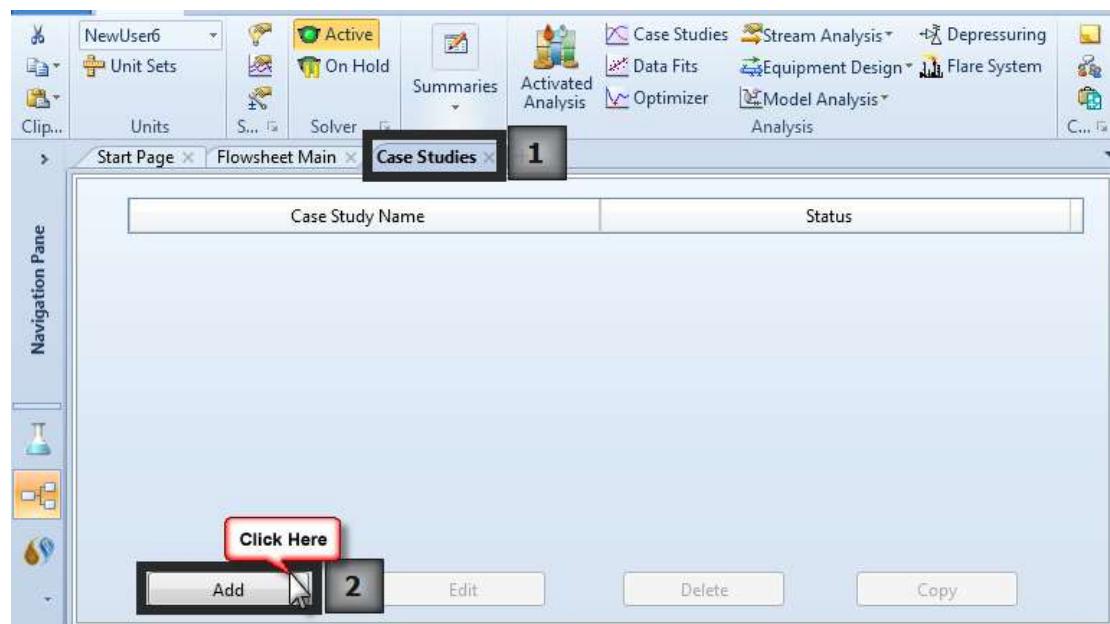
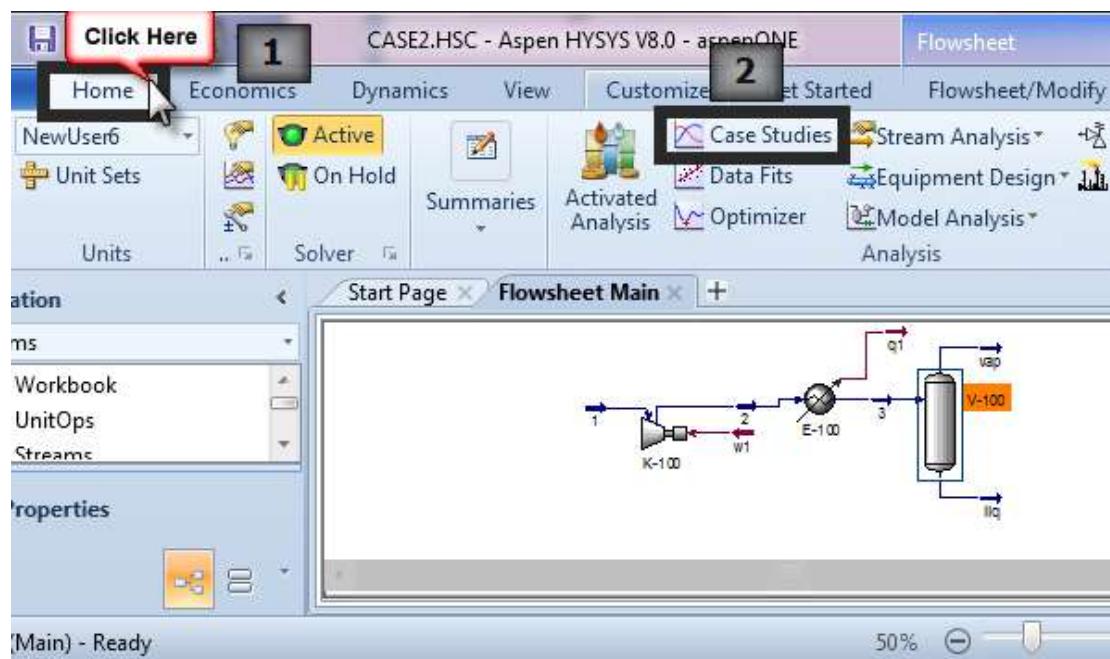
# Save Your Case!

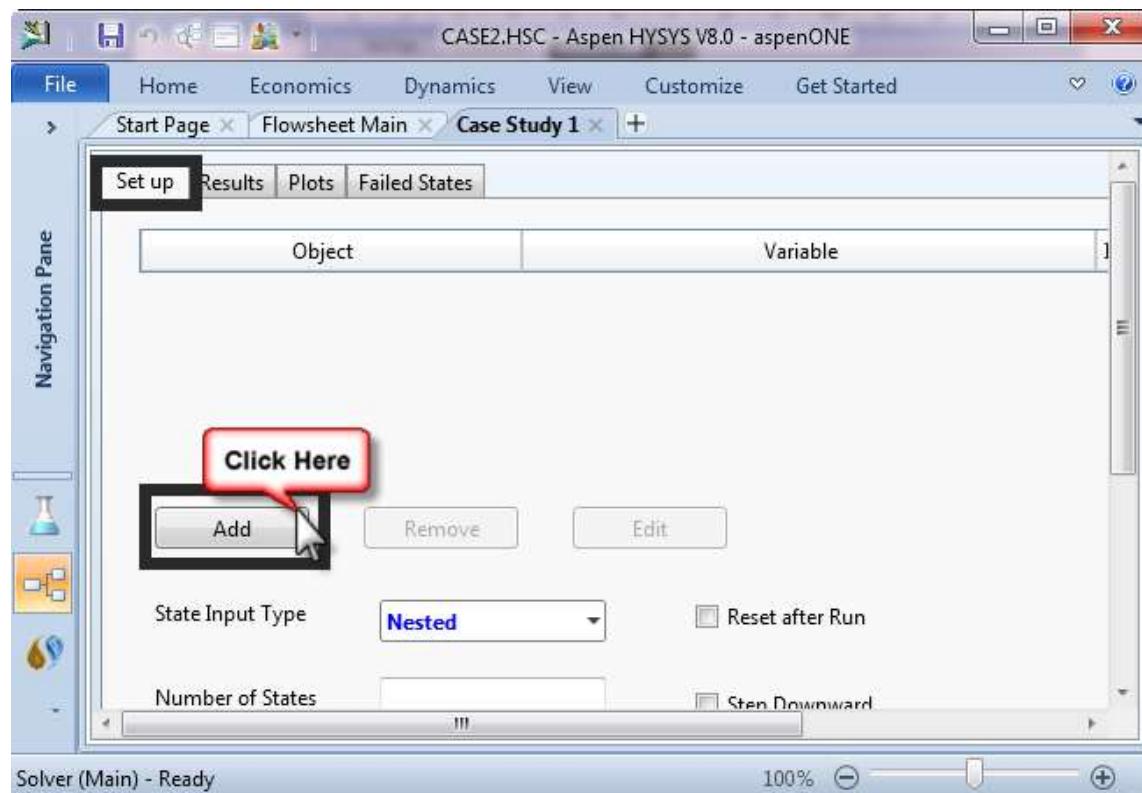
## Case Study

We need to study the effect of changing the Temperature of the cooler out stream (stream no 3) on the flow rate of the liquid product stream.

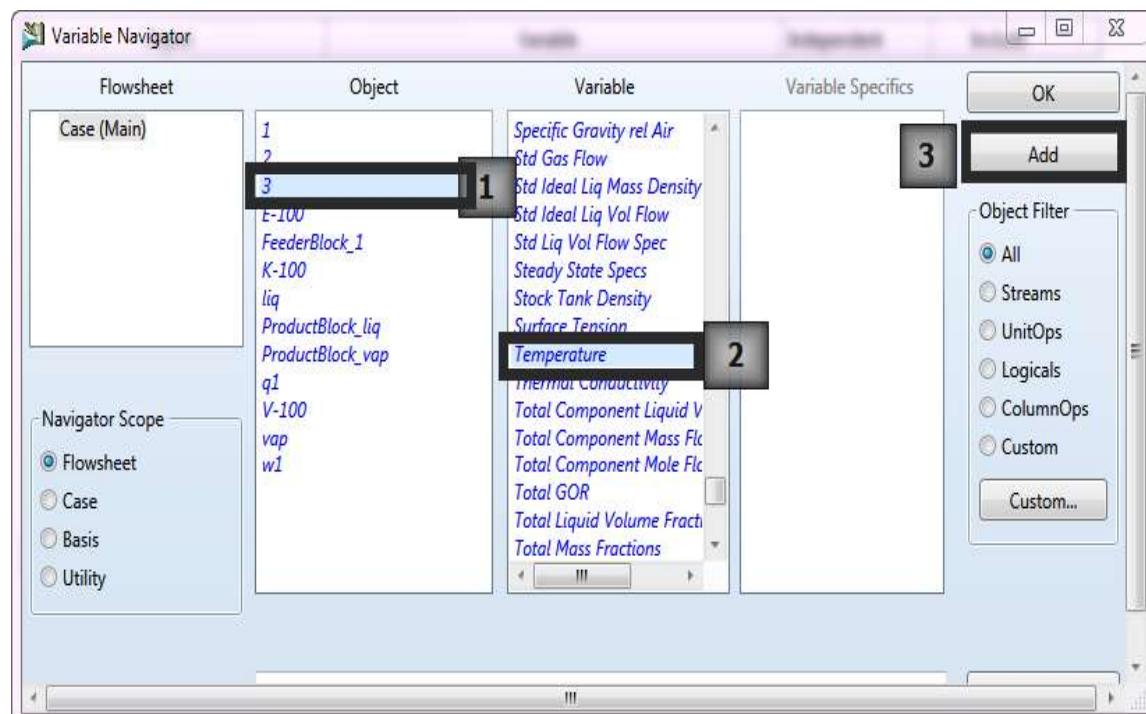
Use range: from -30 to 30 °C with step size =5 °C

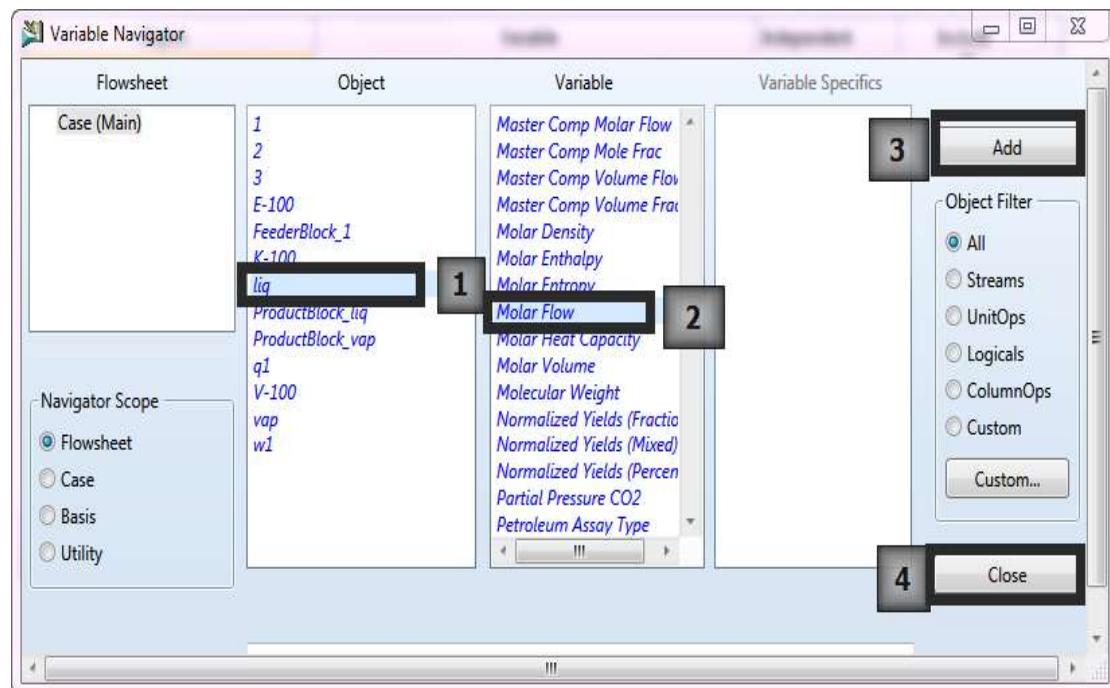
To create a case study in HYSYS you can simply click on Case Studies button on the Home menu:



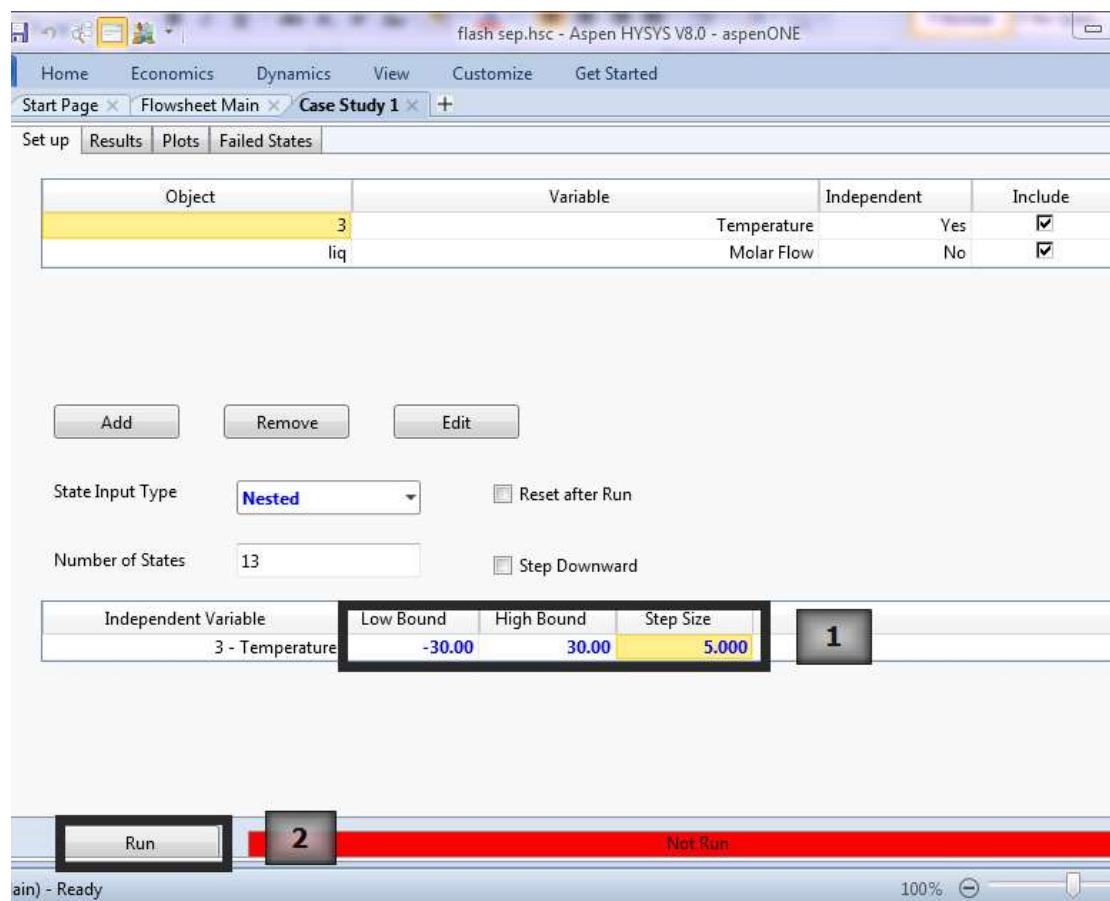


Add the two variables:

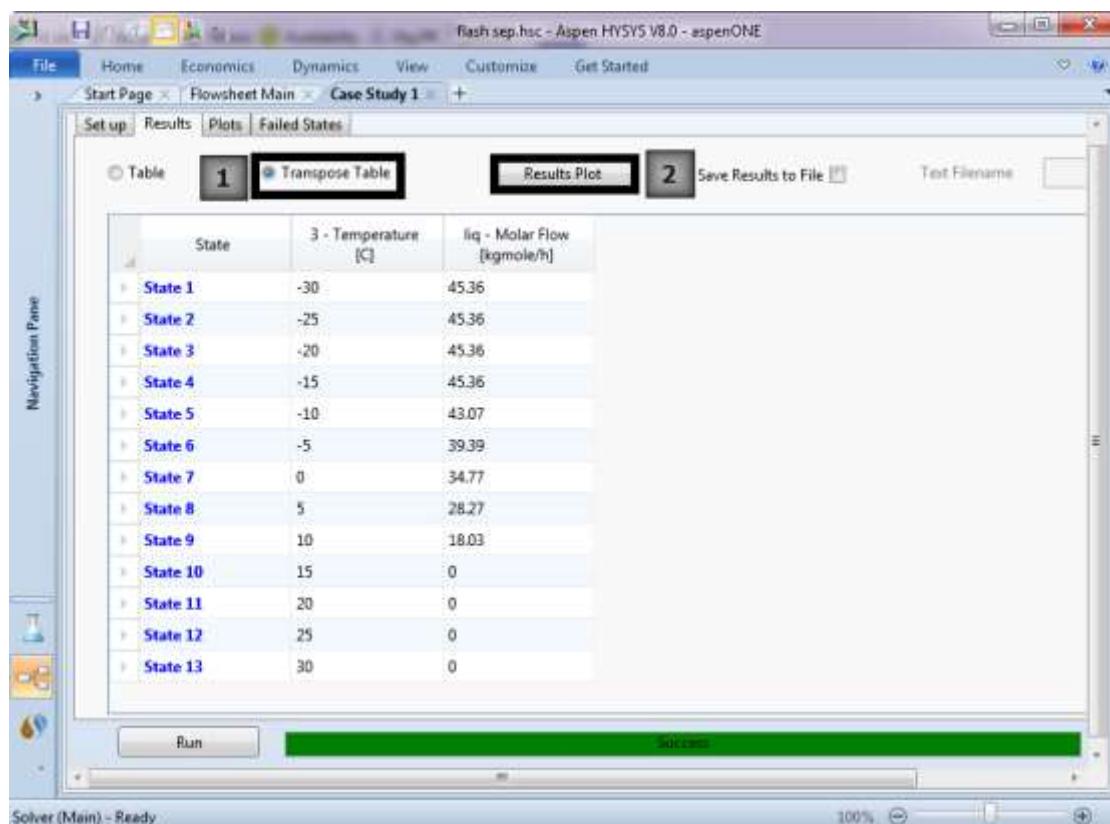
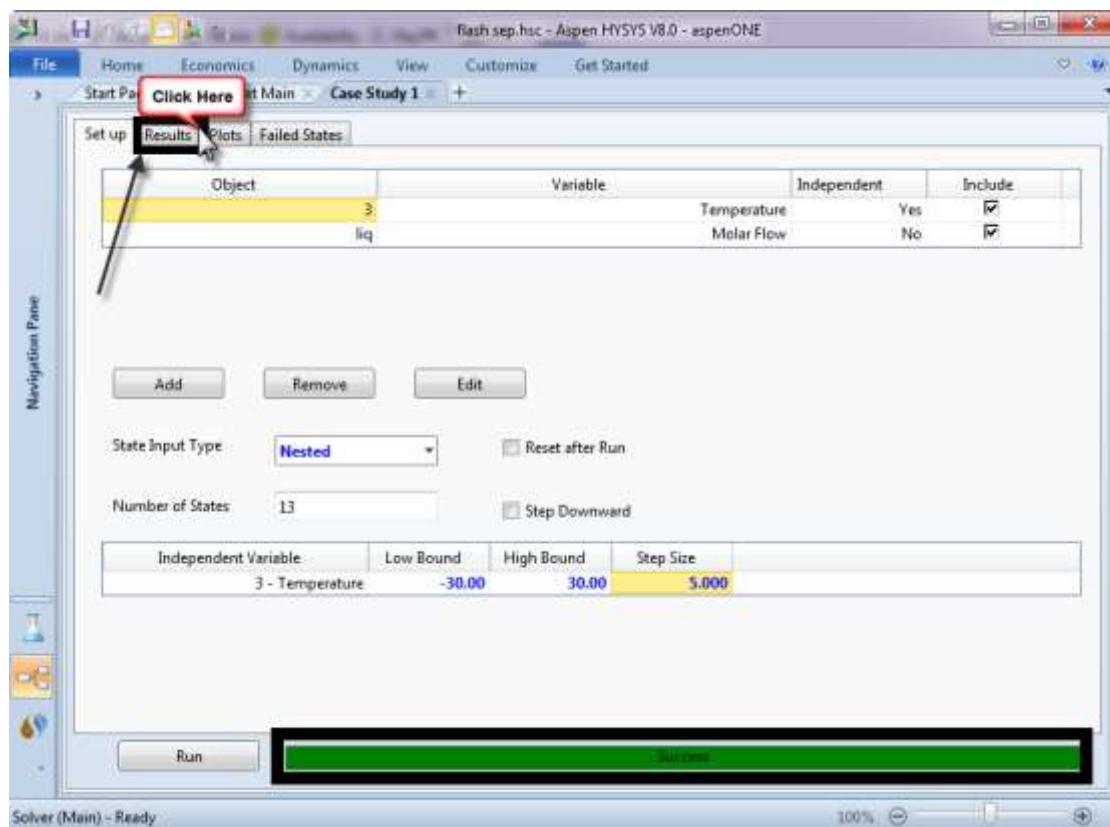




Specify the range of the study:



Click run, and then you can view results from the Results tab



Click Results Plot to view graph

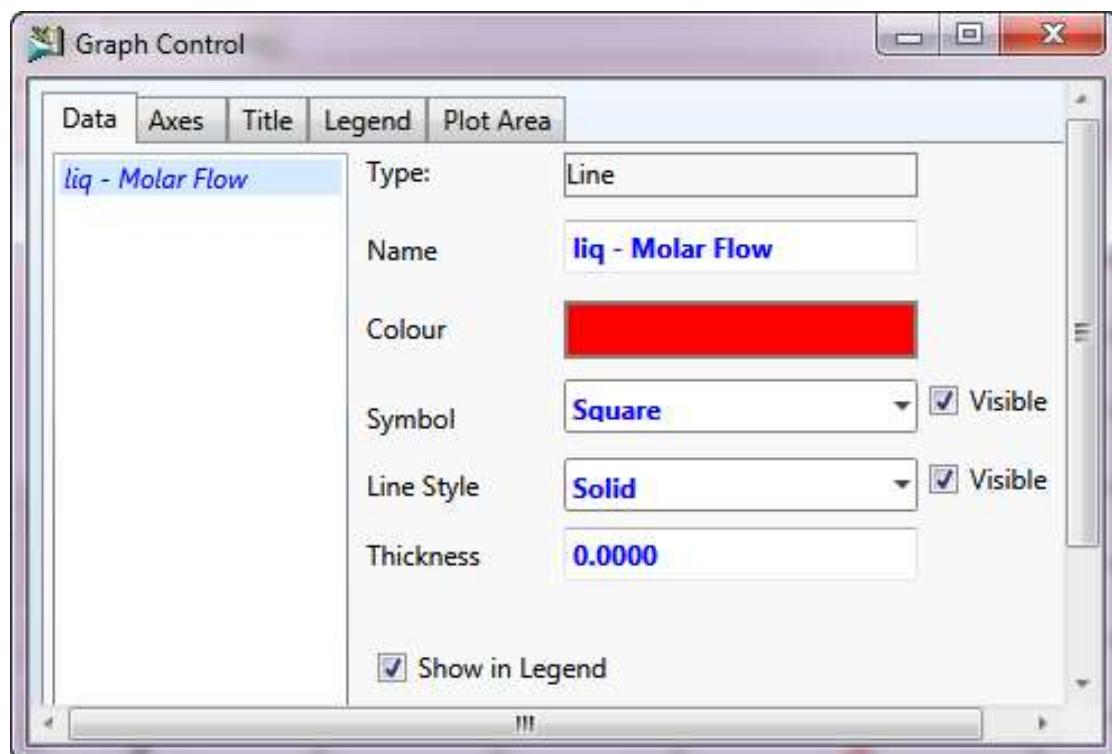
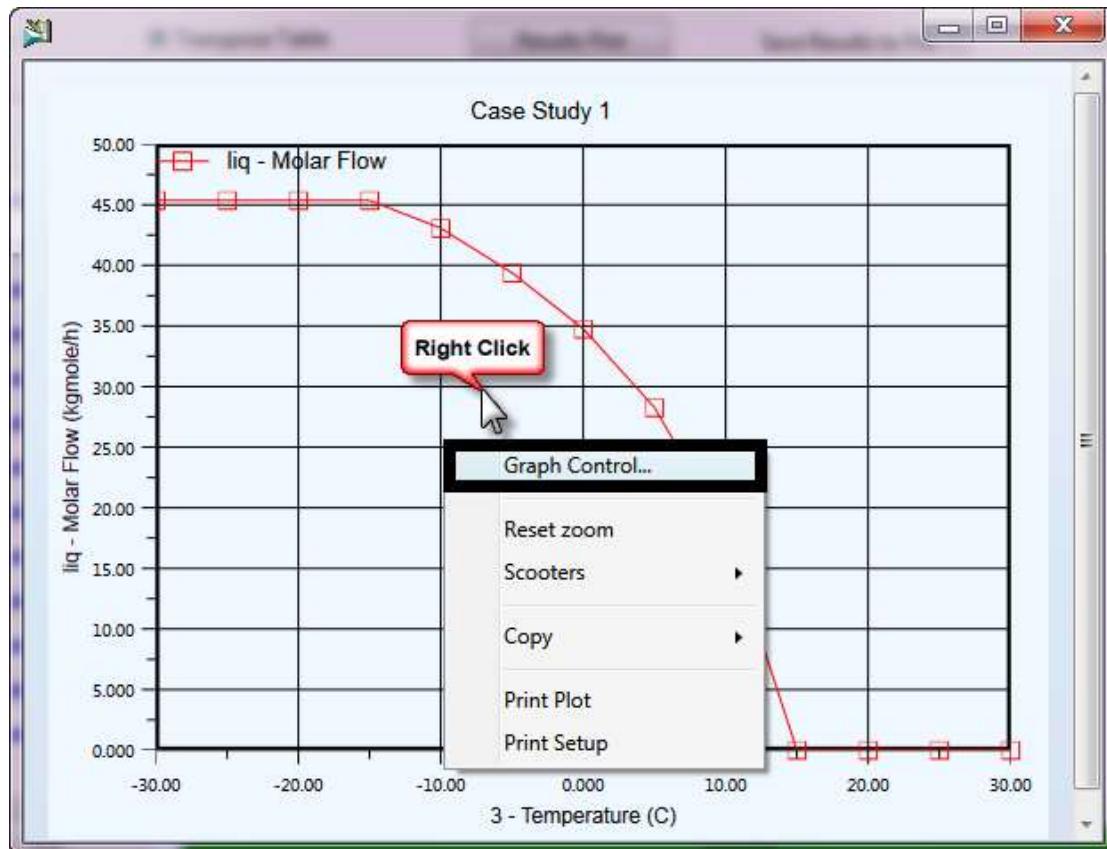


### **Comment:**

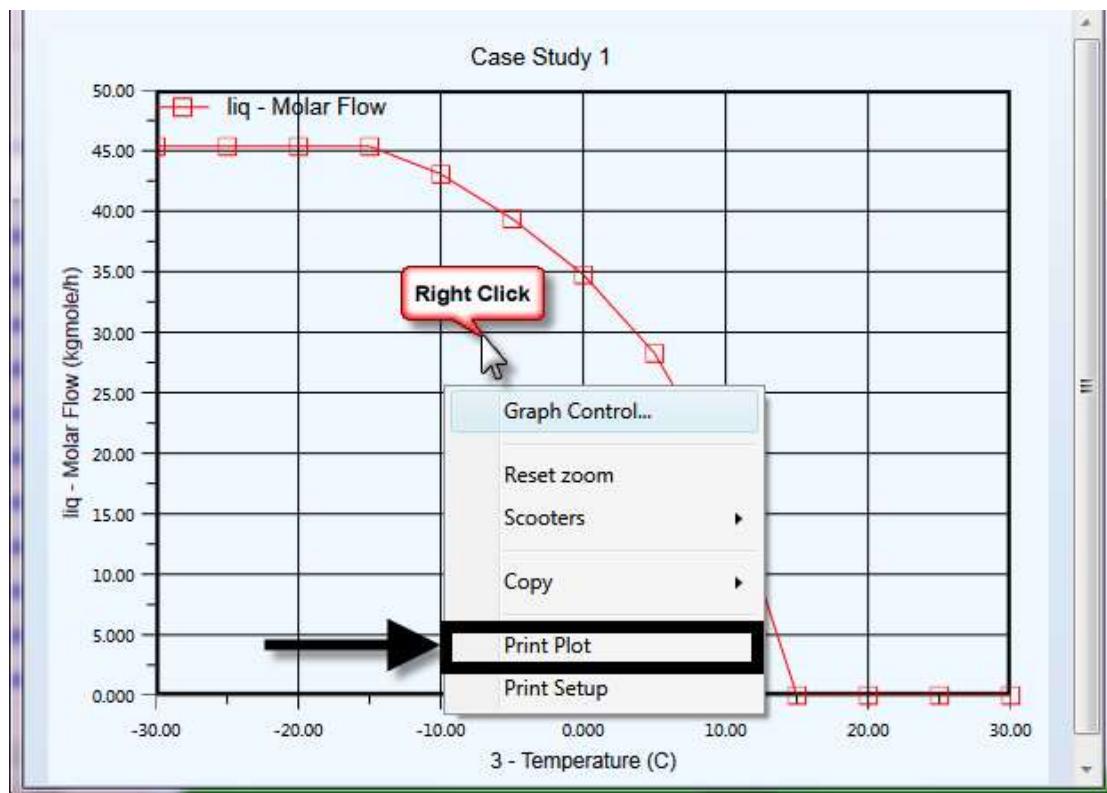
As we see, when the temperature increases the liquid flow rate decrease, the liquid start to decrease @ -15 °C, and @ 15 °C there will be no liquid product and all the product will be vapor.

This is a simple case; you can create your own case study with the same steps.

You can change the scale of axis & the curve color by right click on the plot area and click graph control:



You can also print this plot from the same menu:



# Save Your Case!

# Refrigerated Gas Plant

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2

## ***Workshop***

In this simulation, a simplified version of a refrigerated gas plant is going to be modeled. The purpose is to find the LTS (Low Temperature Separator) temperature at which the hydrocarbon dew point target is met. The Sales Gas hydrocarbon dew point should not exceed -15°C at 6000 kPa. The incoming gas is cooled in two stages—first by exchange with product Sales Gas in a gas-gas exchanger (Gas-Gas) and then in a propane chiller (Chiller), represented here by a Cooler operation. A Balance operation will be used to evaluate the hydrocarbon dew point of the product stream at 6000 kPa.

## ***Learning Objectives***

- Add a hypothetical component
- Install and converge heat exchangers.
- Understand logical operations (Balances and Adjusts).
- Use the Case Study tool to perform case studies on your simulation.

Example:

The feed stream enters an **inlet separator**, which removes the free liquids. Overhead gas from the Separator is fed to the Chiller where it is cooled to  $-20^{\circ}\text{C}$ , which will be modeled simply as a Cooler (Pressure Drop=35 kPa). The cold stream is then separated in a low-temperature separator (LTS). Overhead gas from the LTS is fed to the heater (Pressure drop=5kPa) where it is heated to  $10^{\circ}\text{C}$  to meet Sales Gas Specifications.

Feed Stream:

Temperature	Pressure	Molar Flow Rate
15°C	6200 kPa	1440 kgmole/h

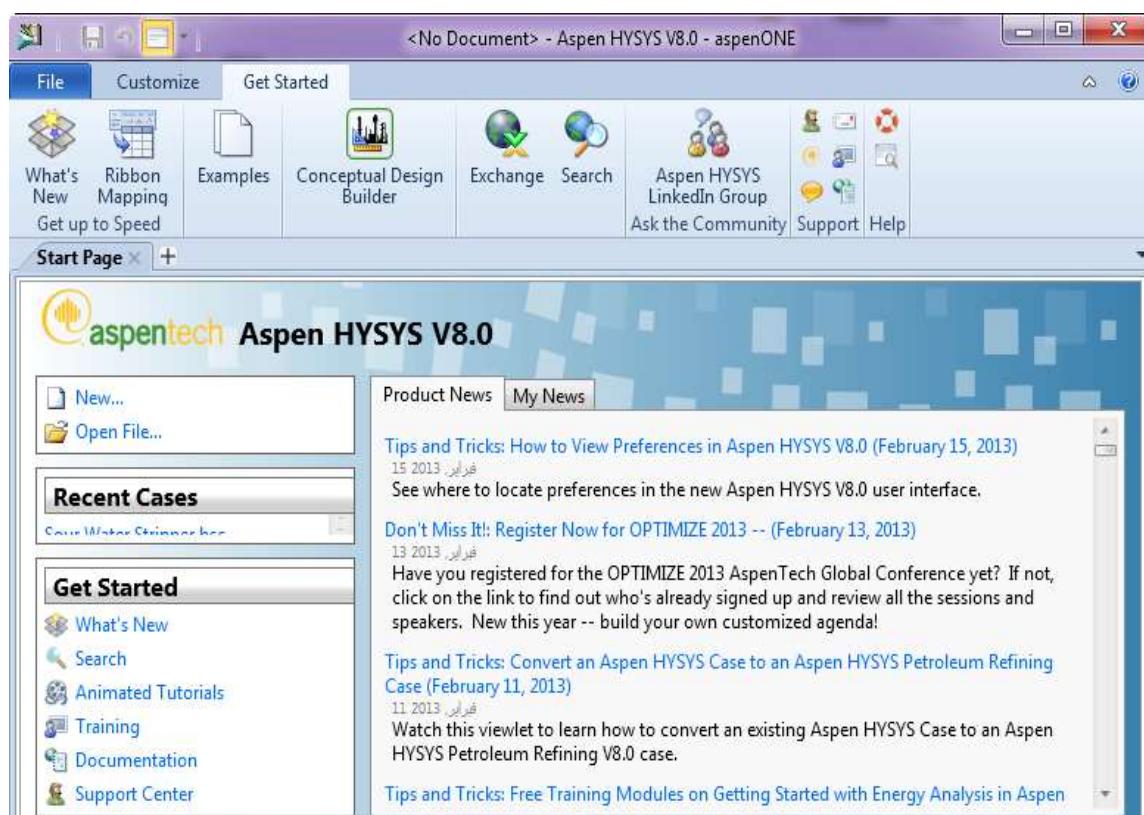
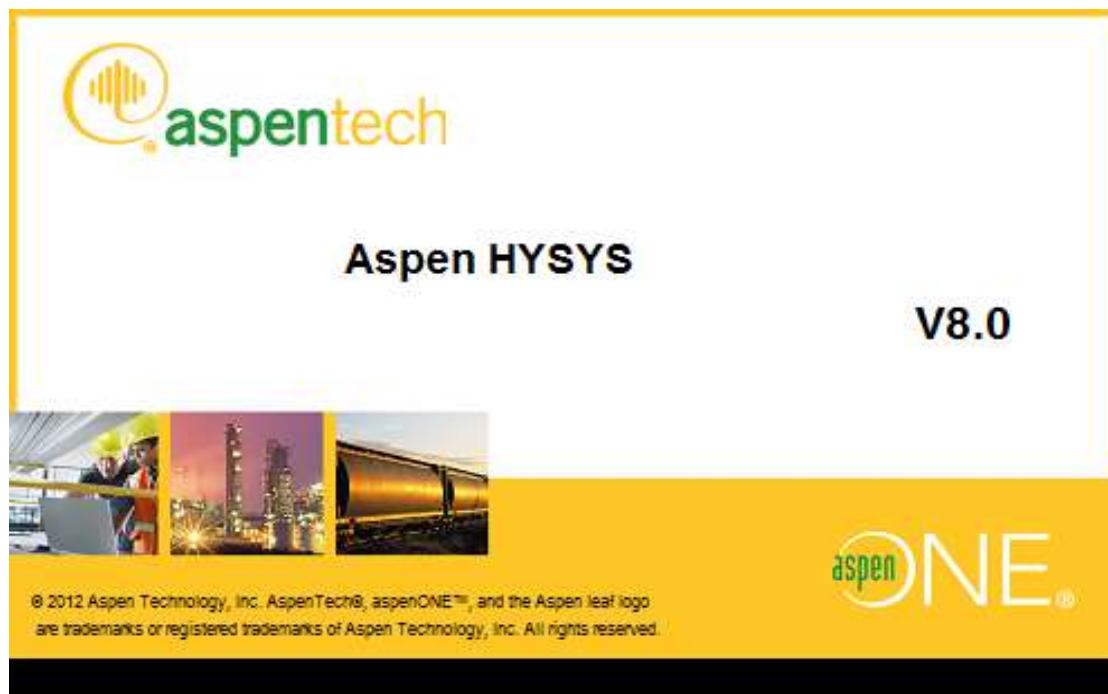
## Composition:

Component	Mole	Component	Mole
N <sub>2</sub>	0.0066	n-Butane	<b>0.0101</b>
H <sub>2</sub> S	0.0003	i-Pentane	<b>0.0028</b>
CO <sub>2</sub>	0.0003	n-Pentane	<b>0.0027</b>
Methane	0.7575	n-Hexane	<b>0.0006</b>
Ethane	0.1709	H <sub>2</sub> O	<b>0.0000</b>
Propane	0.0413	C <sub>7+</sub> (NBP=110°C)	<b>0.0001</b>
i-Butane	0.0068		

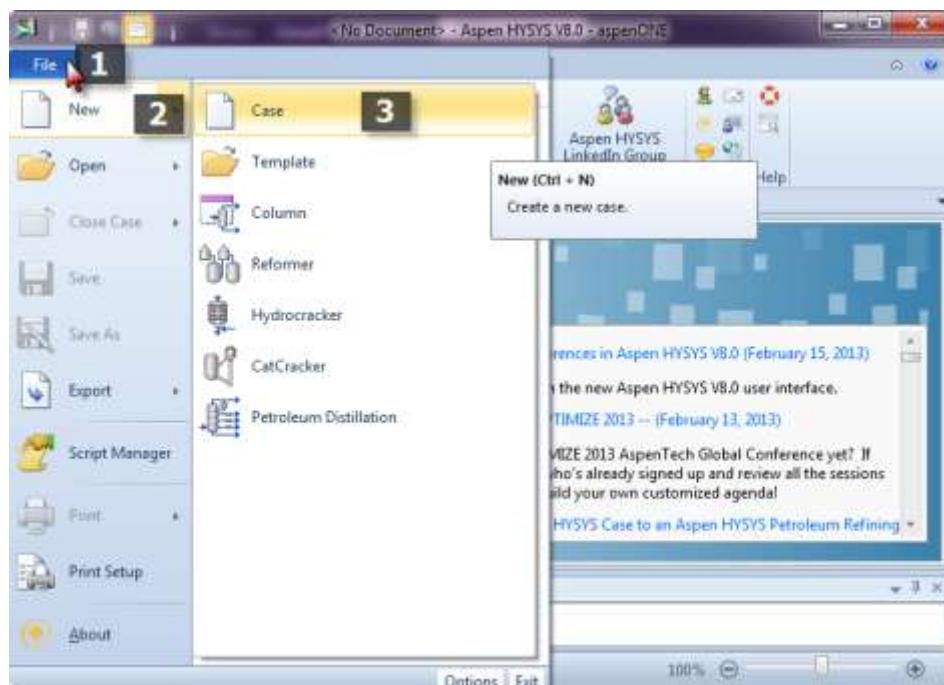
FP: Peng Robinson

- Calculate the duty rejected from the chiller .....
- Calculate the duty Absorbed inside the Heater .....

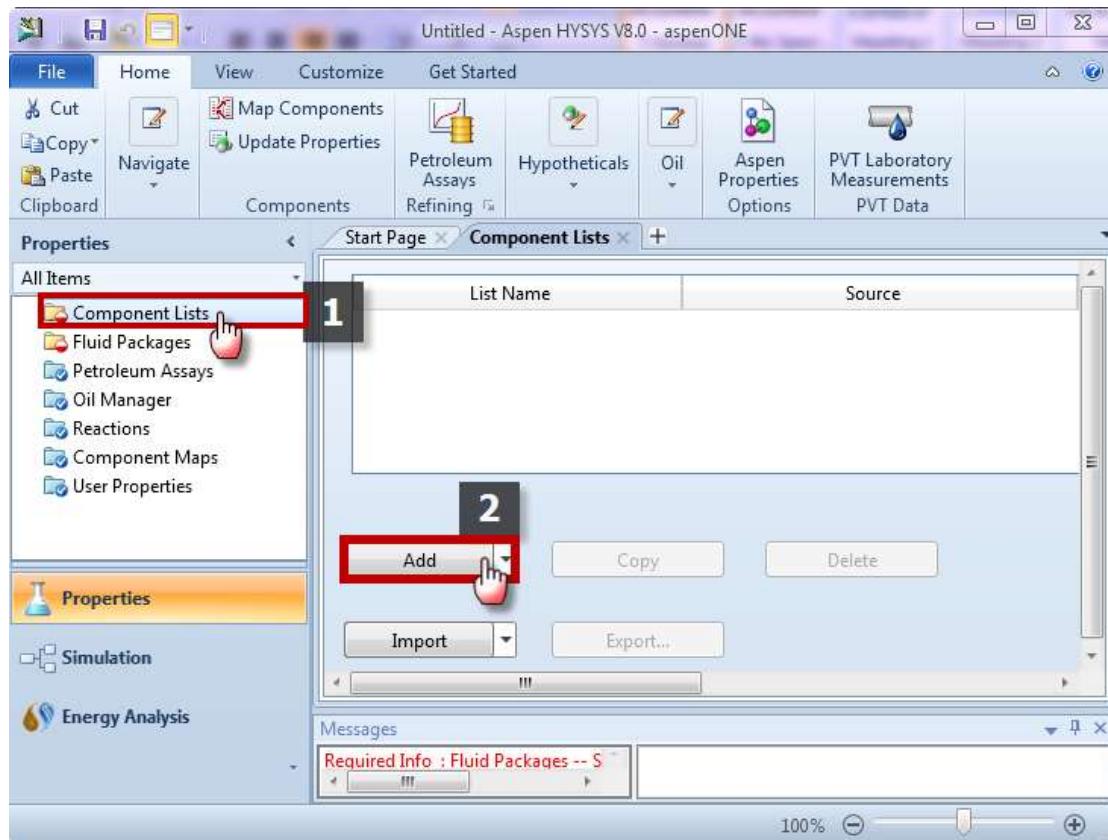
To start the program, From Start Menu, Select All Programs >>  
Aspen Tech >> Process Modeling V8.0 >>> Aspen HYSYS >>  
Aspen HYSYS



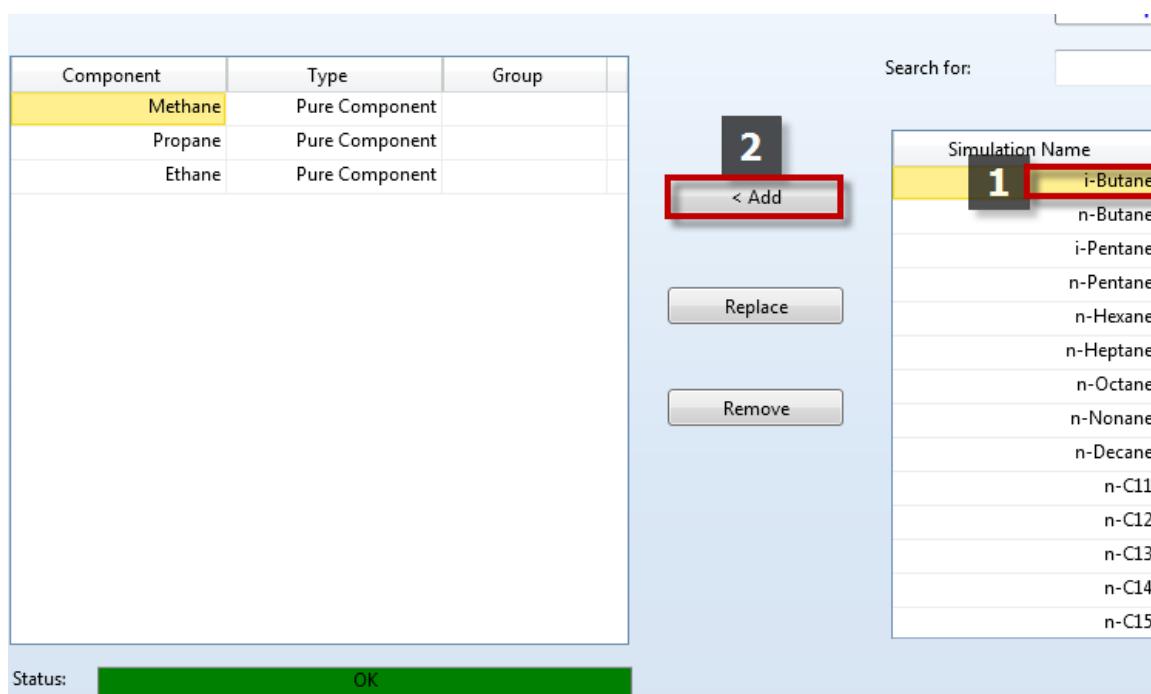
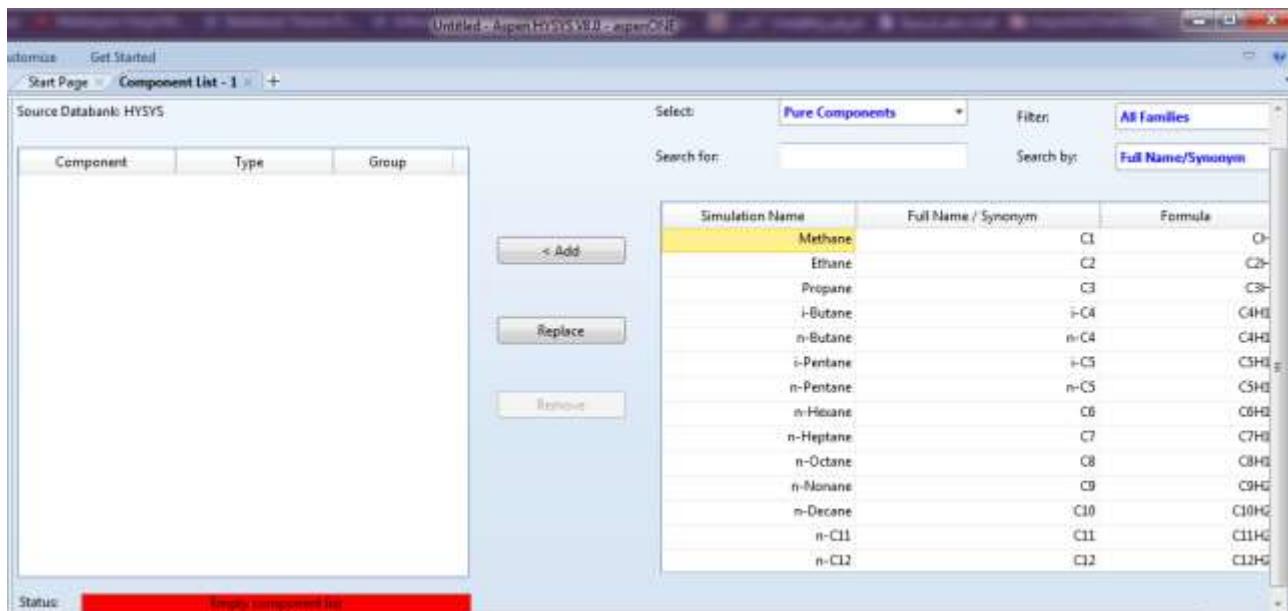
## 4- First, Start a new case



## 5- Add the Components



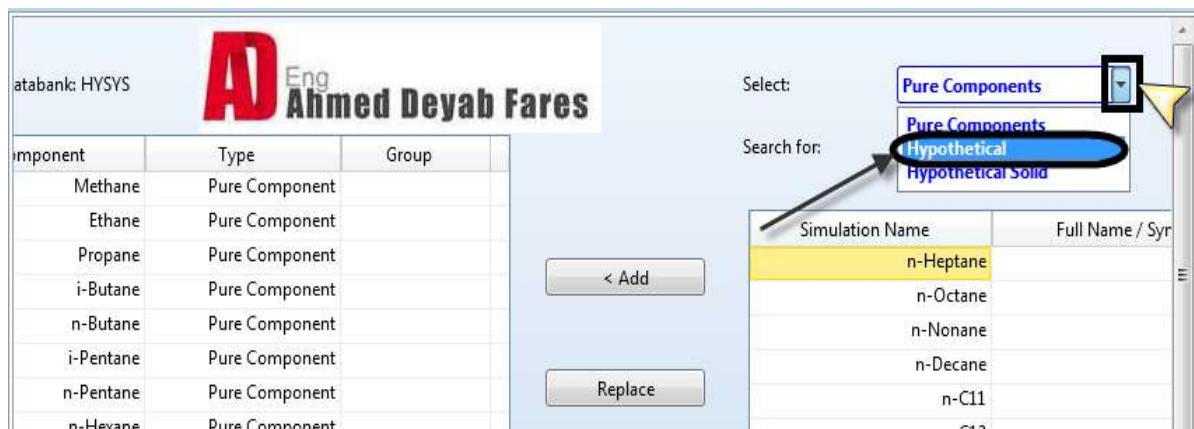
## 6- Choose the system components from the databank:



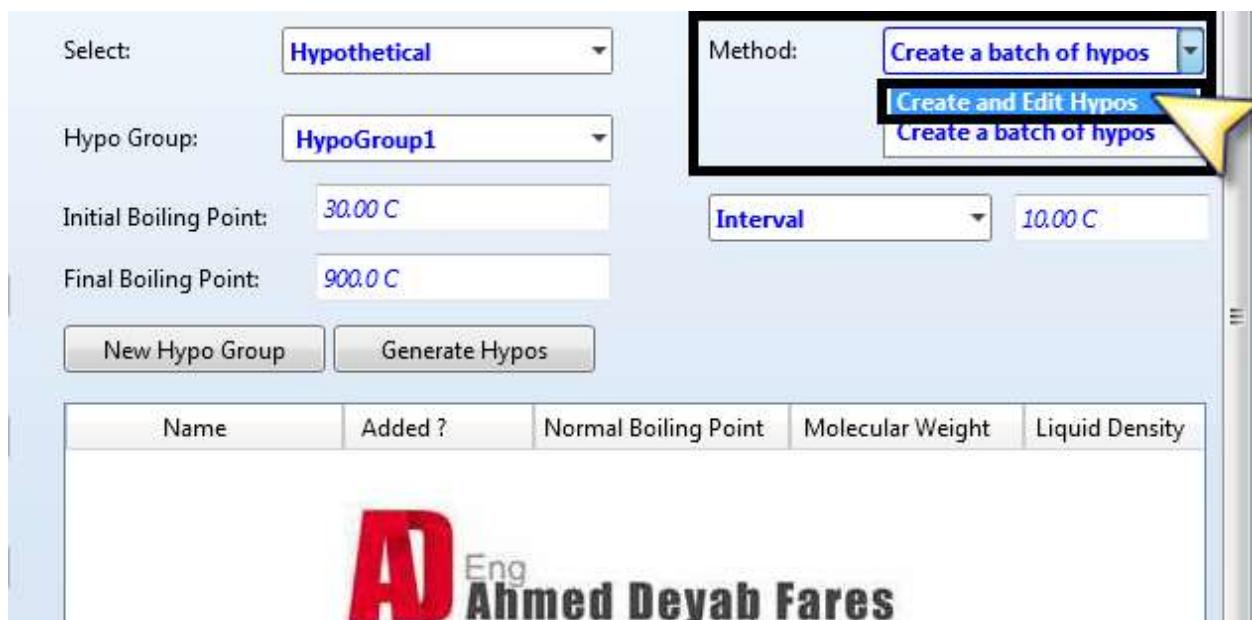
After adding the pure components (N<sub>2</sub>, H<sub>2</sub>S, CO<sub>2</sub>, C1, C2, C3, n-C4, i-C4, n-C5, i-C5, n-C6, H<sub>2</sub>O) we have to add the last component (C7<sup>+</sup>) which is not a pure component as it represents all components above C7 including C7 in the feed.

To define C7+ we have to create it as a hypothetical component as the following:

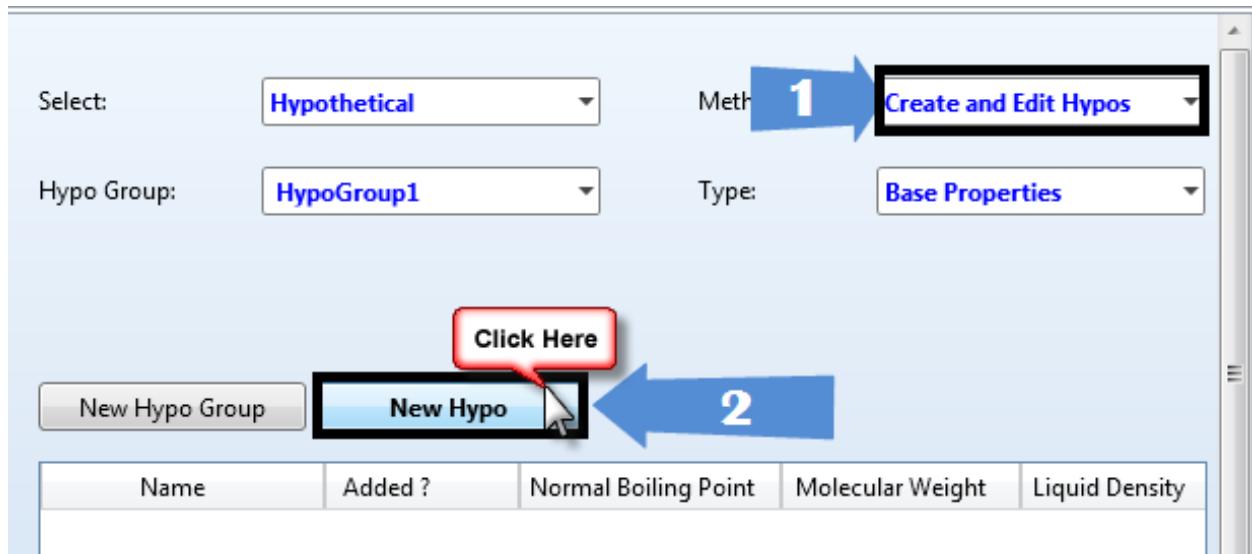
From the drop menu select Hypothetical instead of pure components



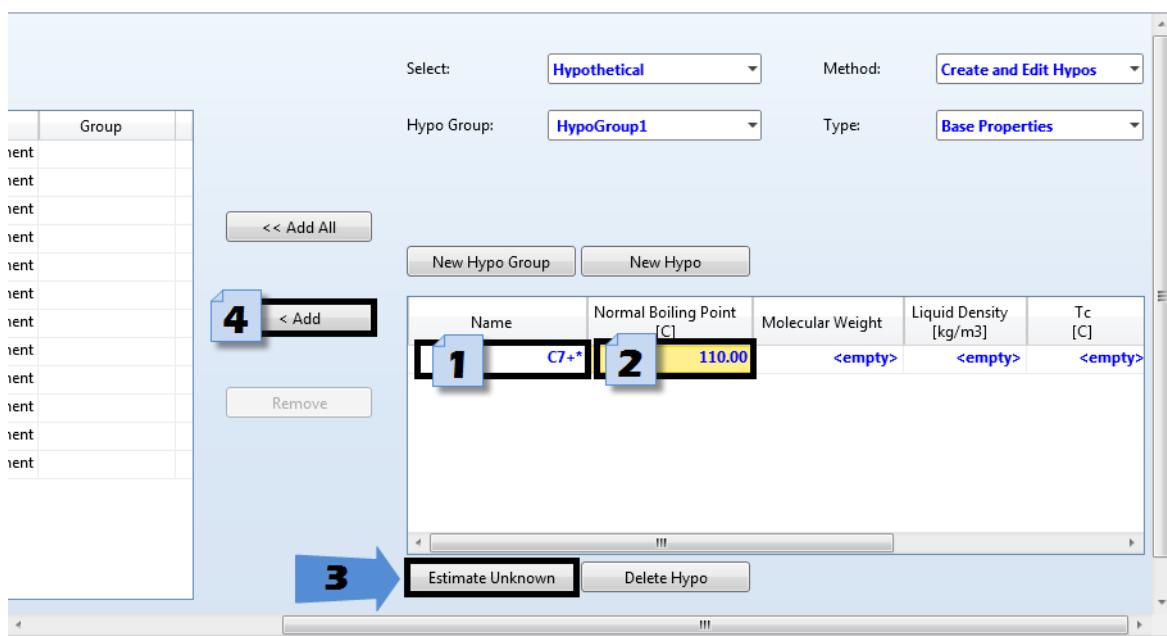
Select create and edit hypos



Click on New Hypo



After adding a hypo component you can edit the name, add the properties you have, and estimate the unknown properties as follows:



Finally add the hypo component to the component list

The screenshot shows the 'Component List - 1' window in HYSYS. On the left, a tree view under 'Properties' shows 'Component Lists' expanded, with 'Component List - 1' selected. Other options like 'Fluid Packages', 'Petroleum Assays', 'Oil Manager', 'Reactions', 'Component Maps', and 'User Properties' are also listed. The main area displays a table titled 'Source Databank: HYSYS' with the following data:

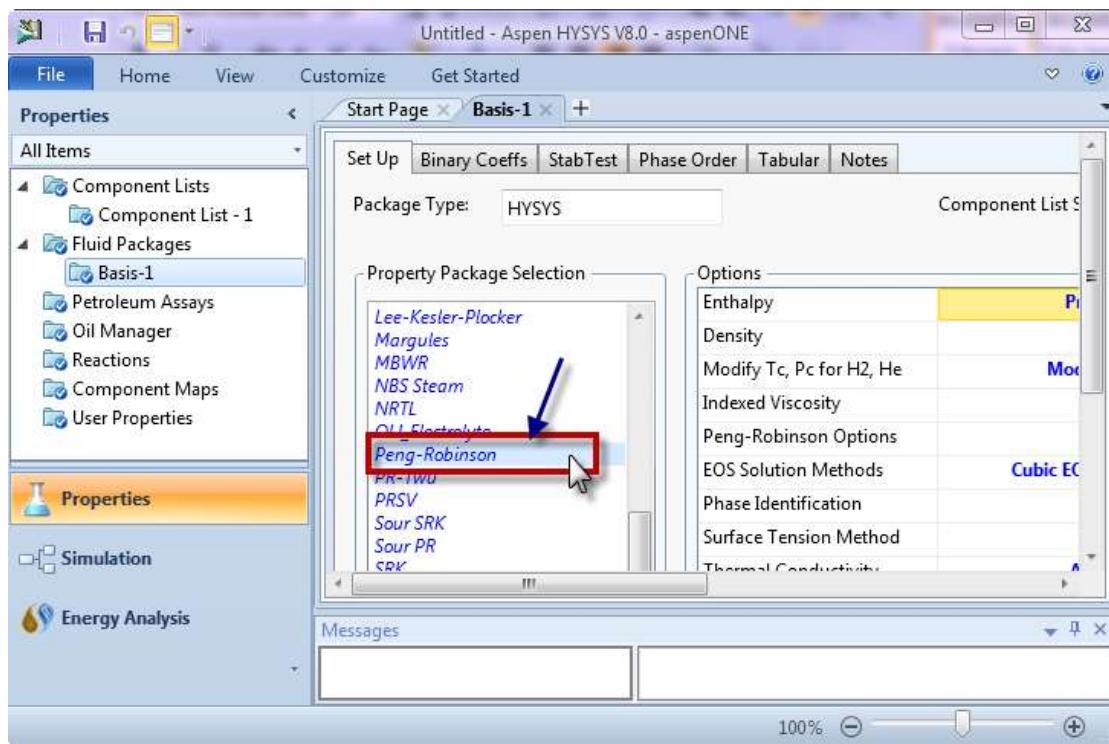
Component	Type	Group
Methane	Pure Component	
Ethane	Pure Component	
Propane	Pure Component	
i-Butane	Pure Component	
n-Butane	Pure Component	
i-Pentane	Pure Component	
n-Pentane	Pure Component	
n-Hexane	Pure Component	
CO <sub>2</sub>	Pure Component	
H <sub>2</sub> S	Pure Component	
Nitrogen	Pure Component	
H <sub>2</sub> O	Pure Component	

At the bottom of the table, there are buttons for 'C7+\* User Defined Hypotheticals' and 'HypoGroup1'. To the right of the table are buttons for '<< Add All', '< Add', and 'Remove'.

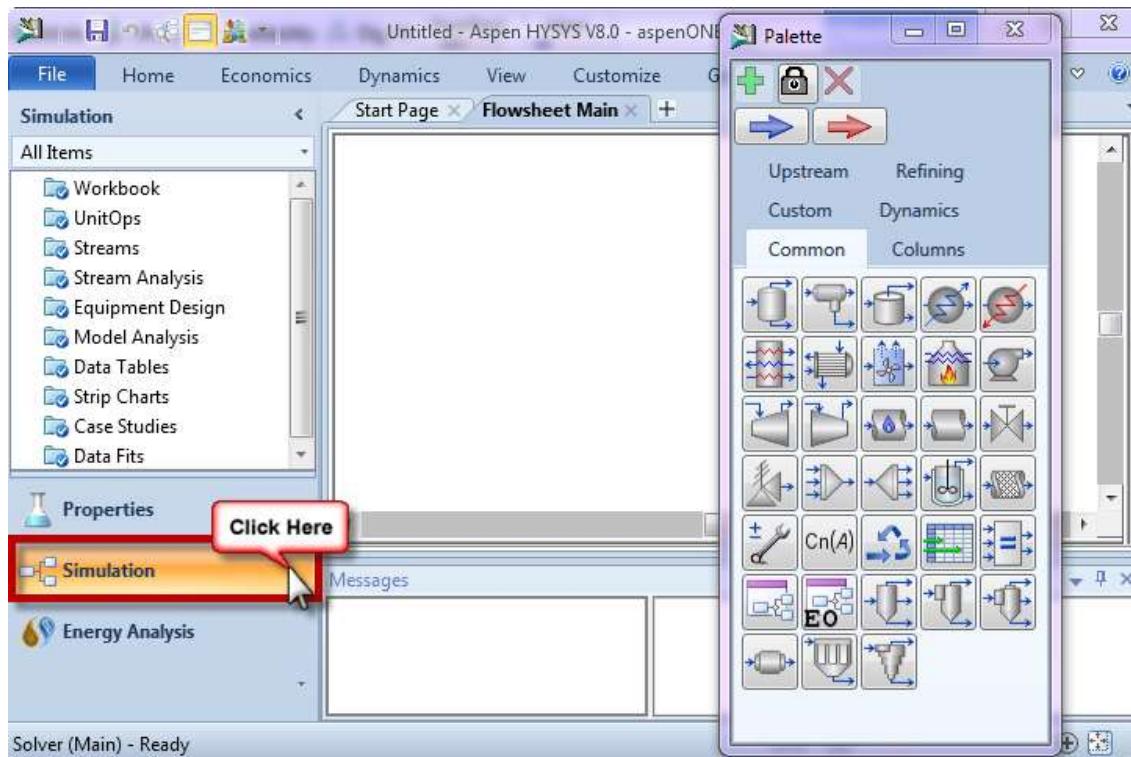
Now, select the suitable fluid package

The screenshot shows the 'Fluid Packages' window in Aspen HYSYS. On the left, a tree view under 'Properties' shows 'Fluid Packages' selected (marked with a red box labeled '1'). Other options like 'Component Lists', 'Petroleum Assays', 'Oil Manager', 'Reactions', 'Component Maps', and 'User Properties' are also listed. The main area has three tabs: 'Fluid Package', 'Component List', and 'Property Package'. Below the tabs, there is a button labeled 'Add' (marked with a red box labeled '2'). A message bar at the bottom says 'Required Info : Fluid Packages -- Select!!!'. The status bar at the bottom right shows '100%'.

In this case, select Peng-Robinson

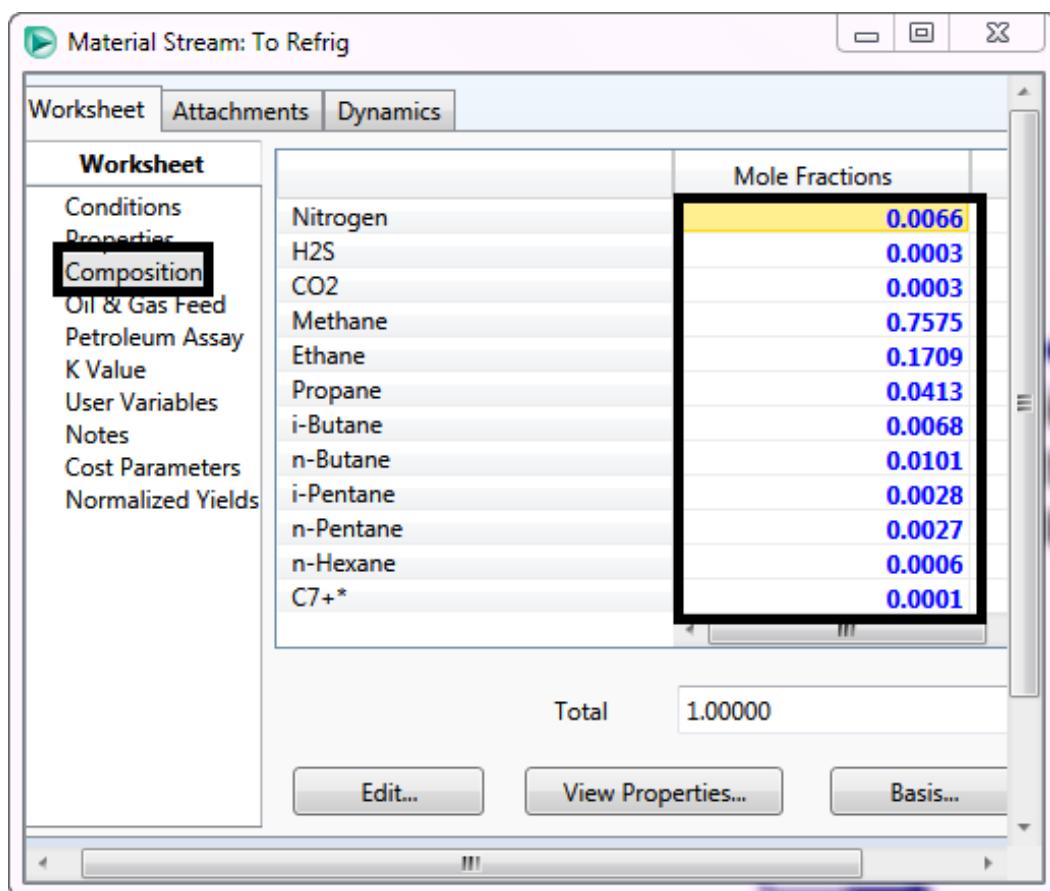
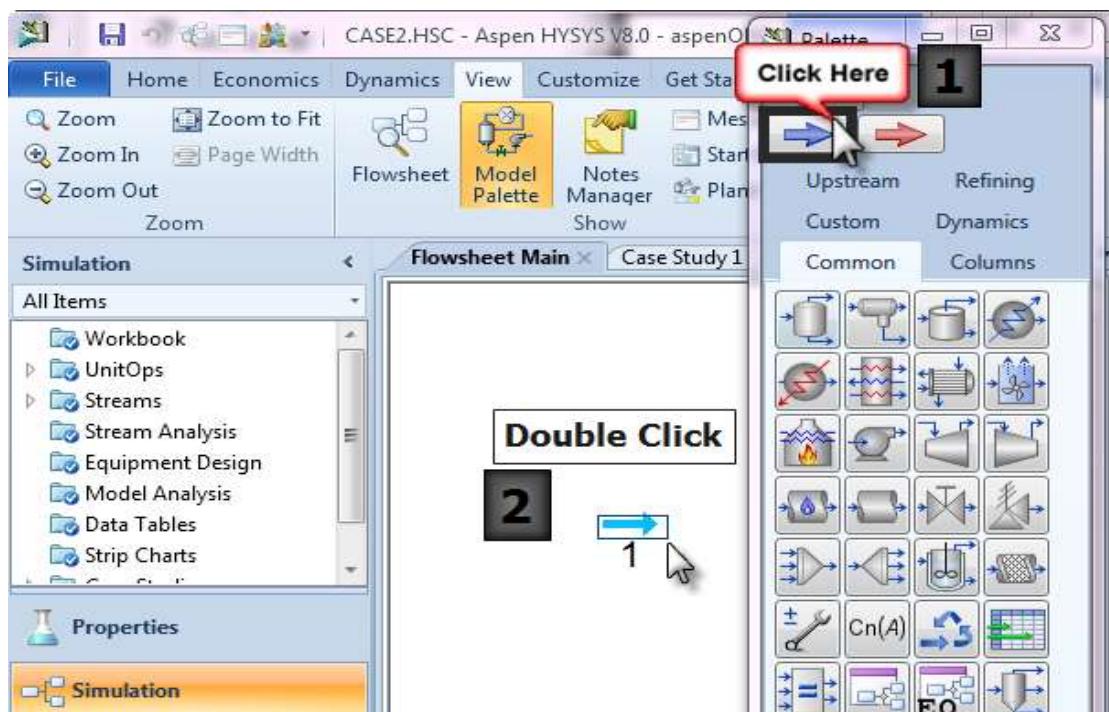


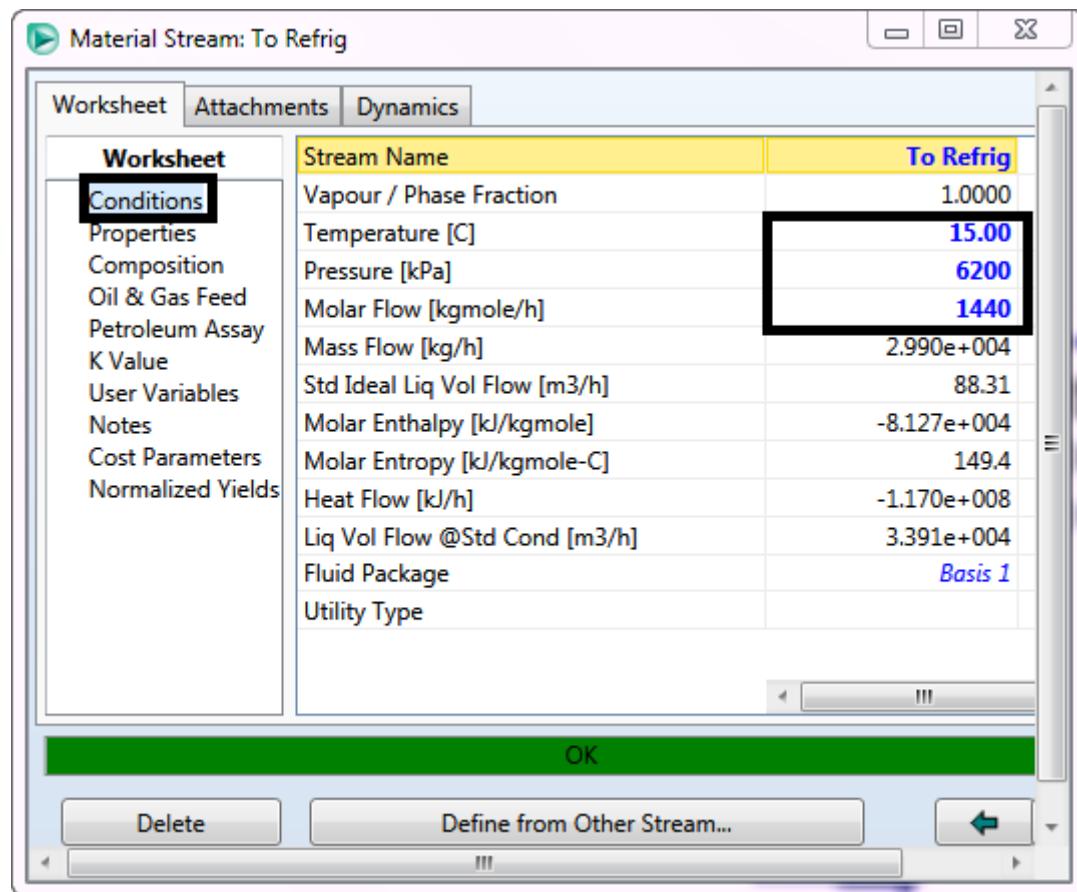
Now you can start drawing the flow sheet for the process by clicking the Simulation button:



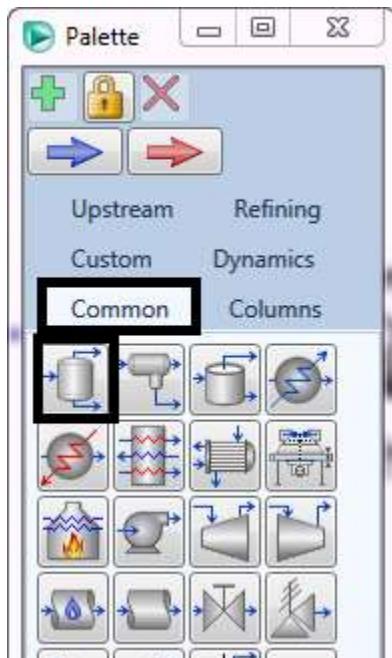
Now add a material stream to define the composition and the conditions of the feed stream

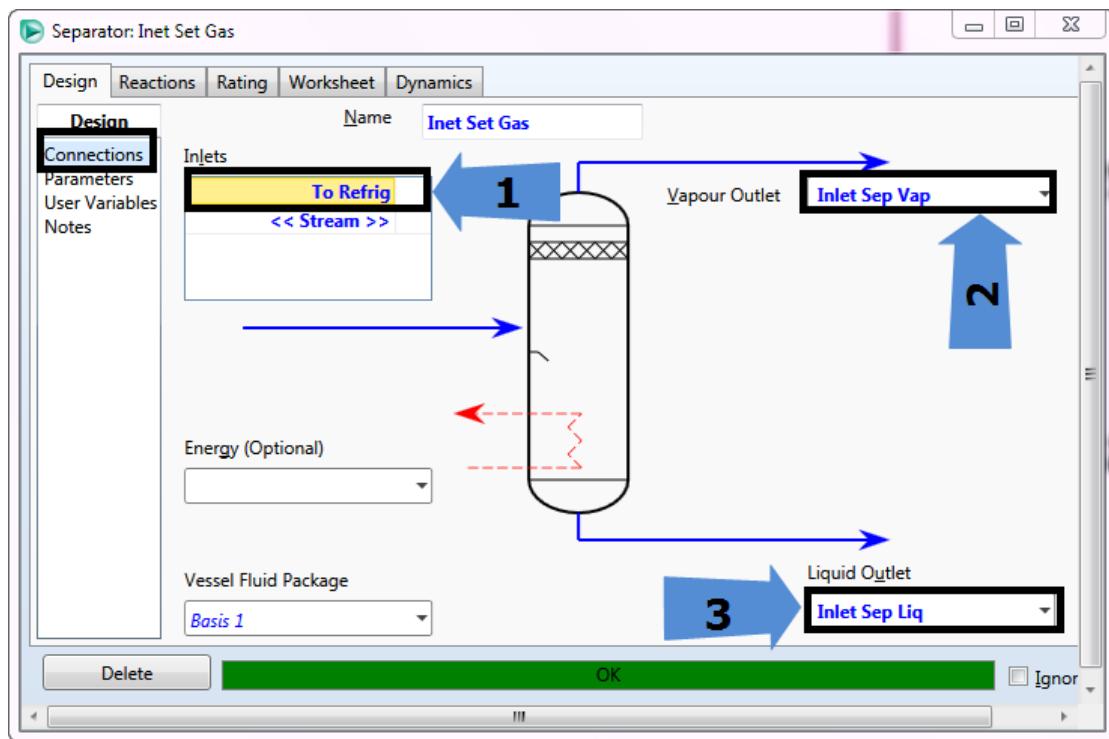
From the palette:



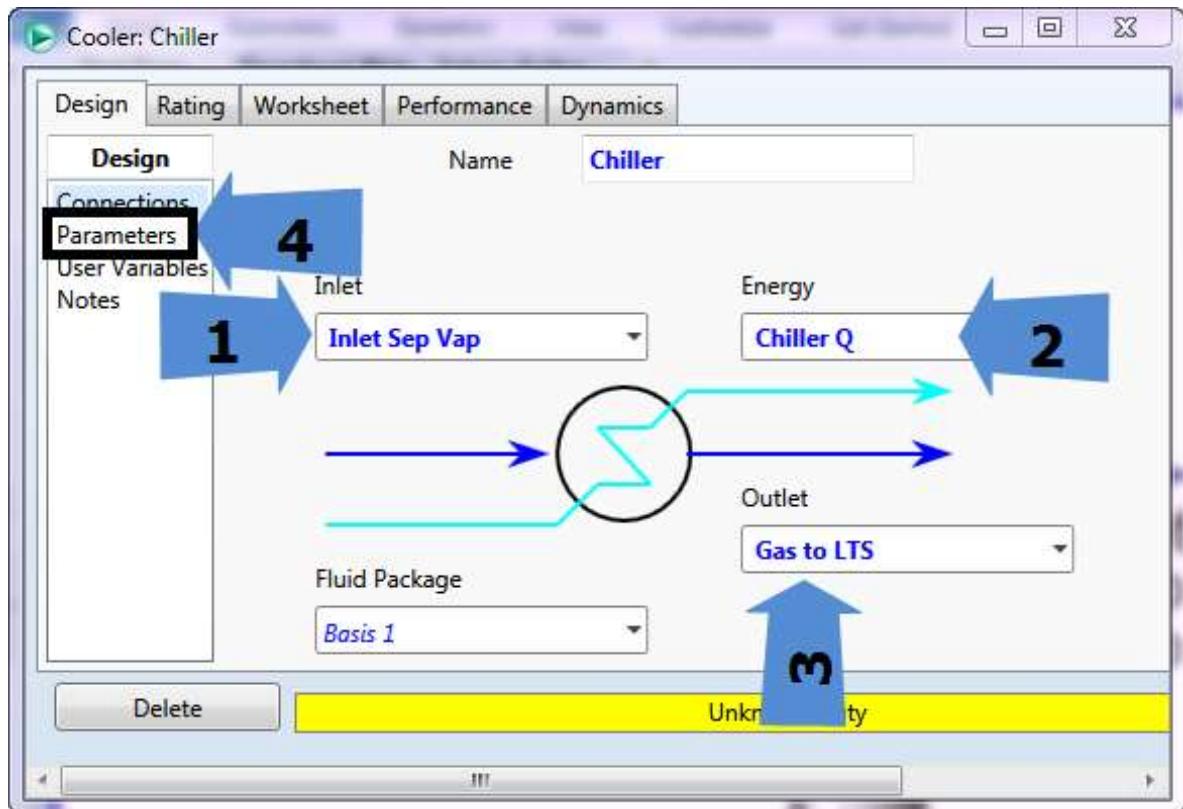


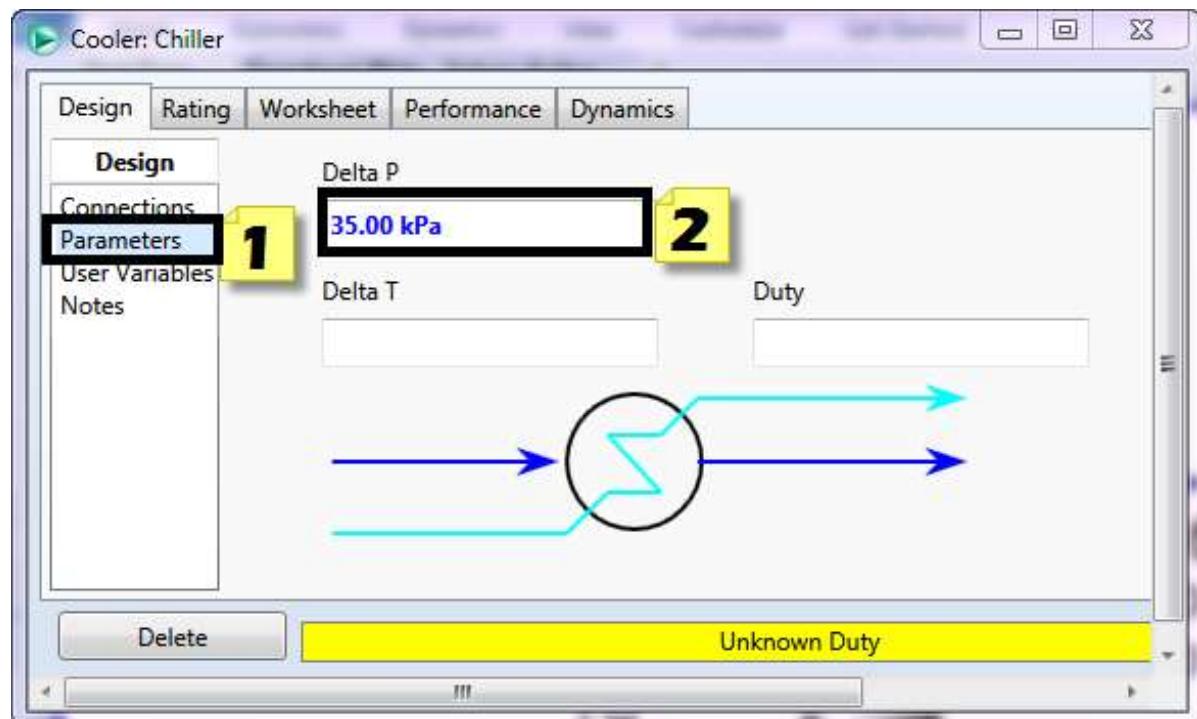
From the palette select the separator:





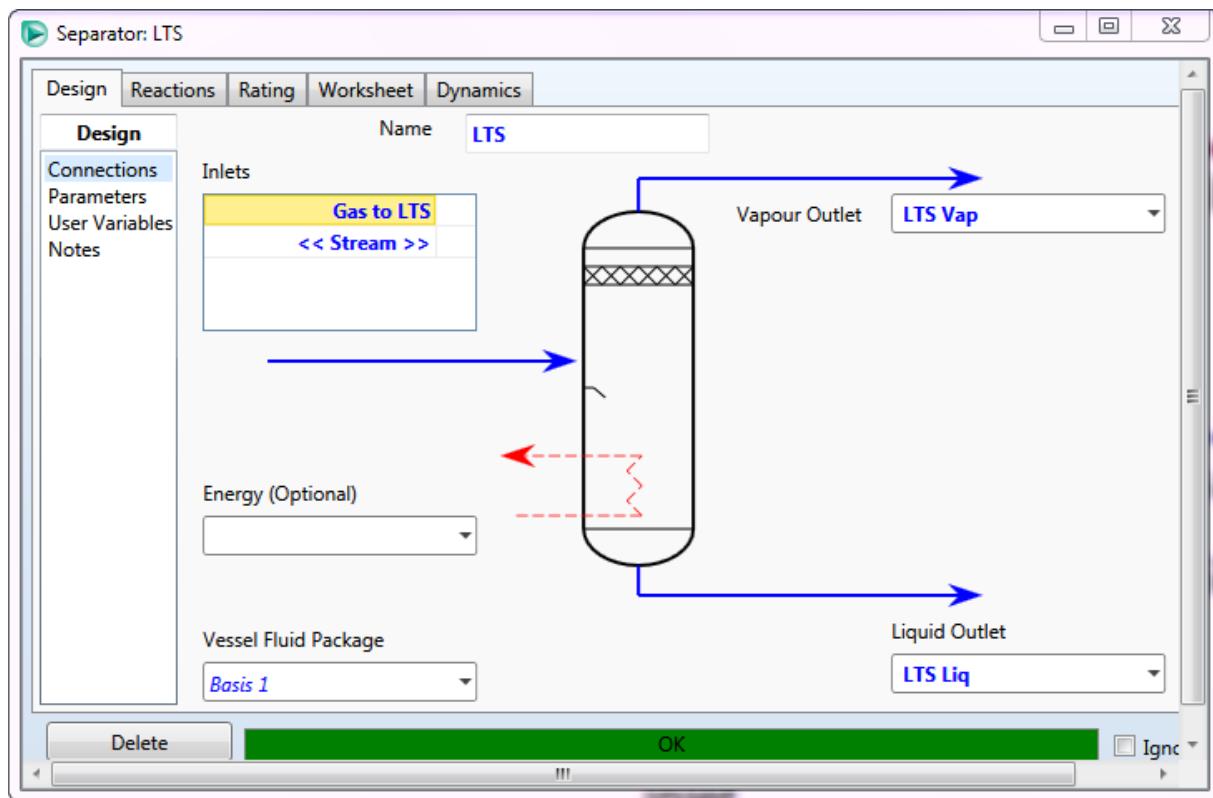
Add a cooler:



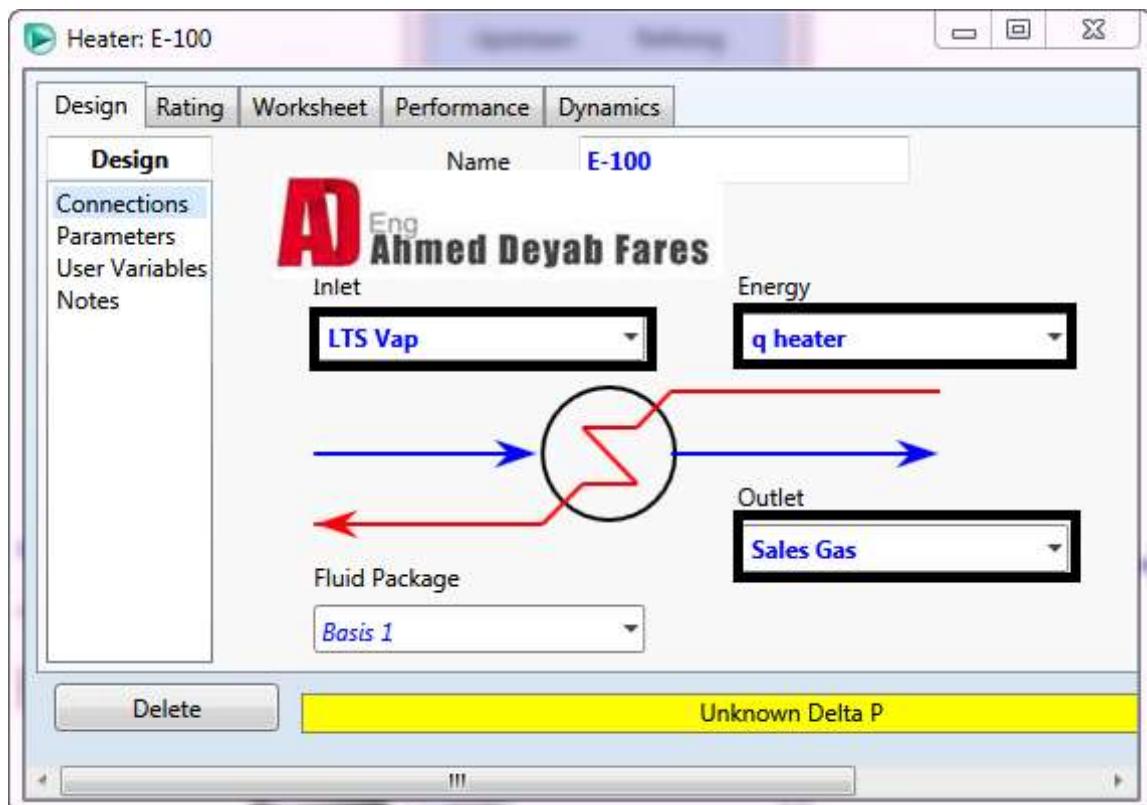


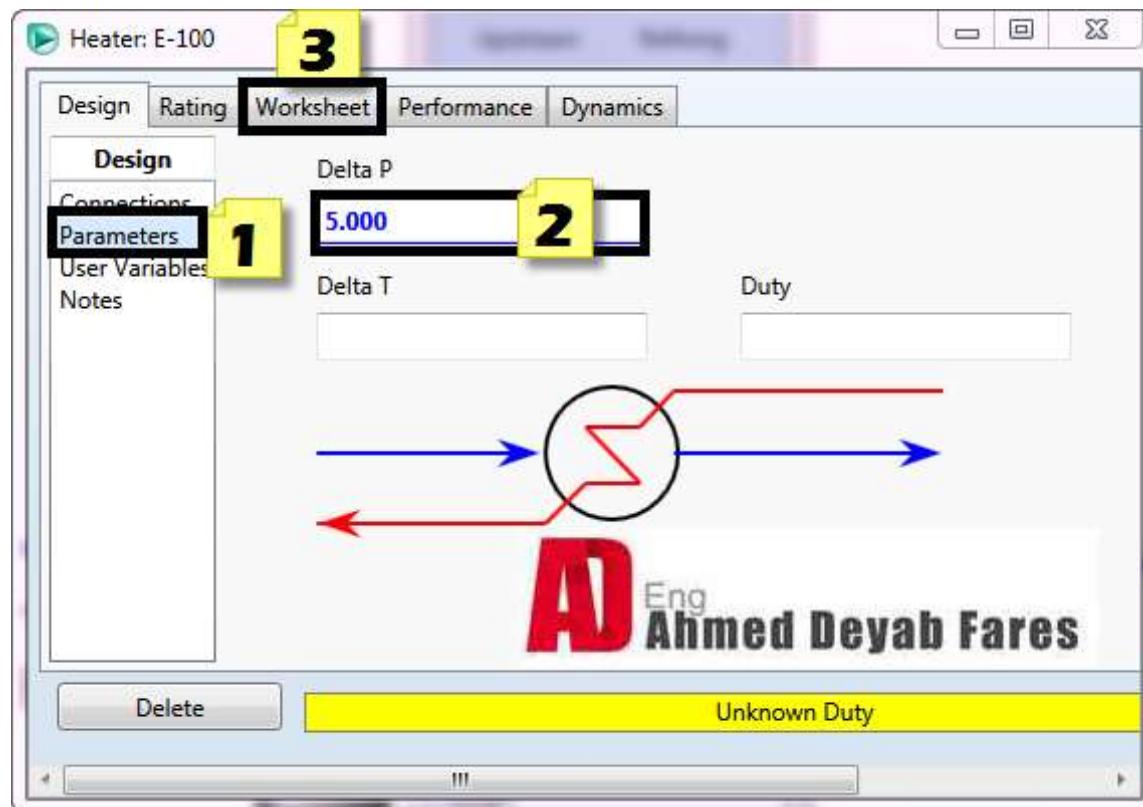
Worksheet				
Conditions	Name	Inlet Sep Vap	Gas to LTS	Chiller Q
Vapour	1.0000	0.8735	<empty>	<empty>
Temperature [C]	15.00	6165	-20.00	<empty>
Pressure [kPa]	6200	1440	6165	<empty>
Molar Flow [kgmole/h]	1440	2.990e+004	1440	<empty>
Mass Flow [kg/h]	2.990e+004	2.990e+004	2.990e+004	<empty>
Std Ideal Liq Vol Flow [m <sup>3</sup> /h]	88.31	88.31	88.31	<empty>
Molar Enthalpy [kJ/kgmole]	-8.127e+004	-8.417e+004	-8.417e+004	<empty>
Molar Entropy [kJ/kgmole-C]	149.4	138.6	138.6	<empty>
Heat Flow [kJ/h]	-1.170e+008	-1.212e+008	-1.212e+008	4.186e+006

Add the LTS Separator:

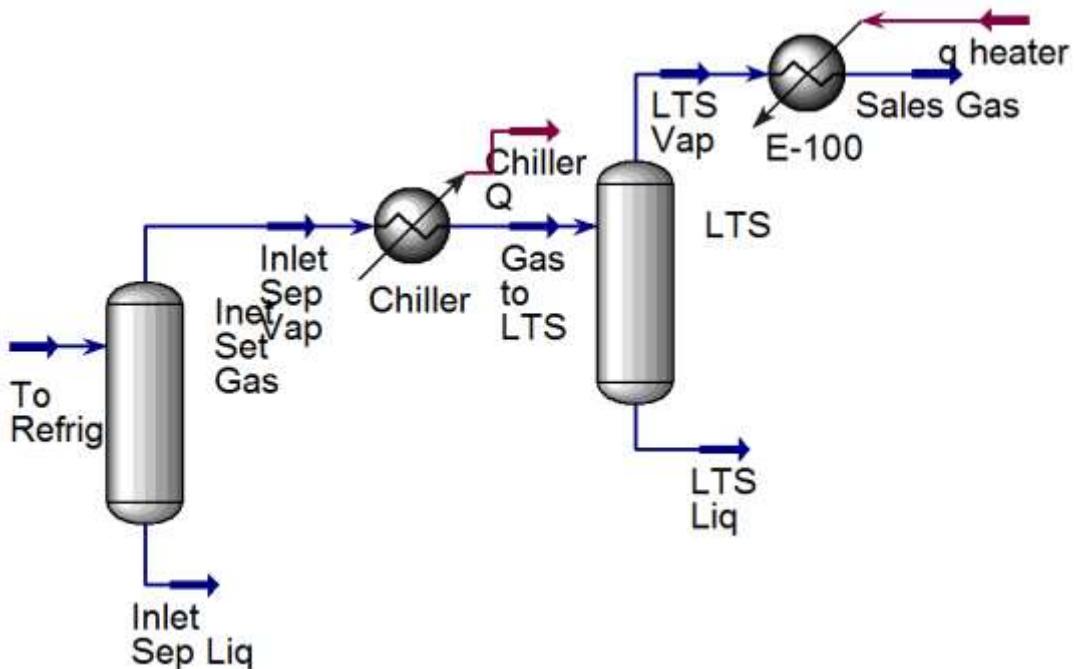


Add a heater:





Name	LTS Vap	Sales Gas
Vapour	1.0000	1.0000
Temperature [C]	-20.00	20.00
Pressure [kPa]	6165	6160
Molar Flow [kgmole/h]	1258	1258
Mass Flow [kg/h]	2.444e+004	2.444e+004
Std Ideal Liq Vol Flow [m <sup>3</sup> /h]	74.84	74.84
Molar Enthalpy [kJ/kgmole]	-8.149e+004	-7.967e+004
Molar Entropy [kJ/kgmole-C]	141.9	148.7
Heat Flow [kJ/h]	-1.025e+008	-1.002e+008



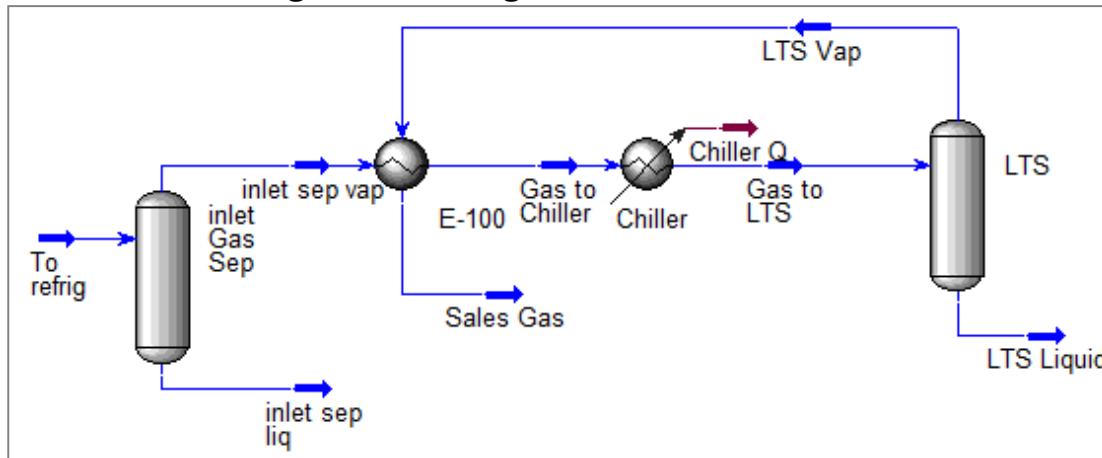
- The duty rejected from the chiller = **4.186 e6** ( $4.186 \times 10^6$ ) kJ/hr
- The duty Absorbed inside the Heater = **2.287 e6** ( $2.287 \times 10^6$ ) kJ/hr

## Heat Exchanger

The design is modified to reduce the operating cost represented in Chiller & Heater duties, by adding a Heat Exchanger before the Chiller where the overhead from the inlet separator is pre-cooled by already refrigerated gas from LTS.

### Heat Exchanger Design Specifications:

- Sales Gas Temperature= 10°C
- Tube side Pressure drop=35kPa
- Shell side Pressure drop= 5kPa
- No heat losses inside the heat exchanger.
- Choose Weighted Exchanger as Model

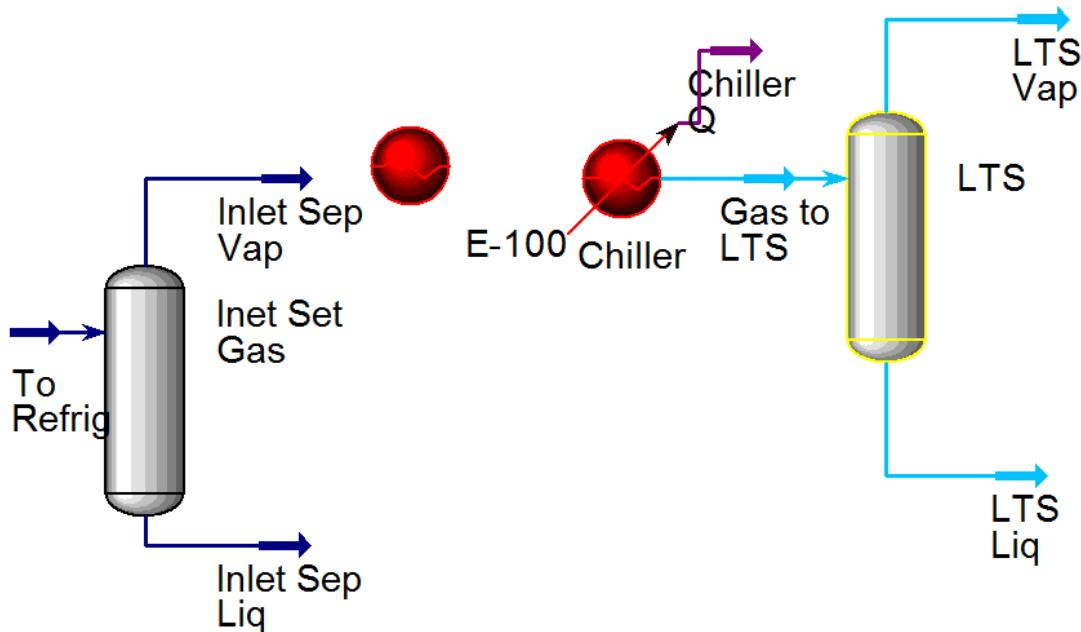


Calculate the duty rejected from the chiller after this modification .....

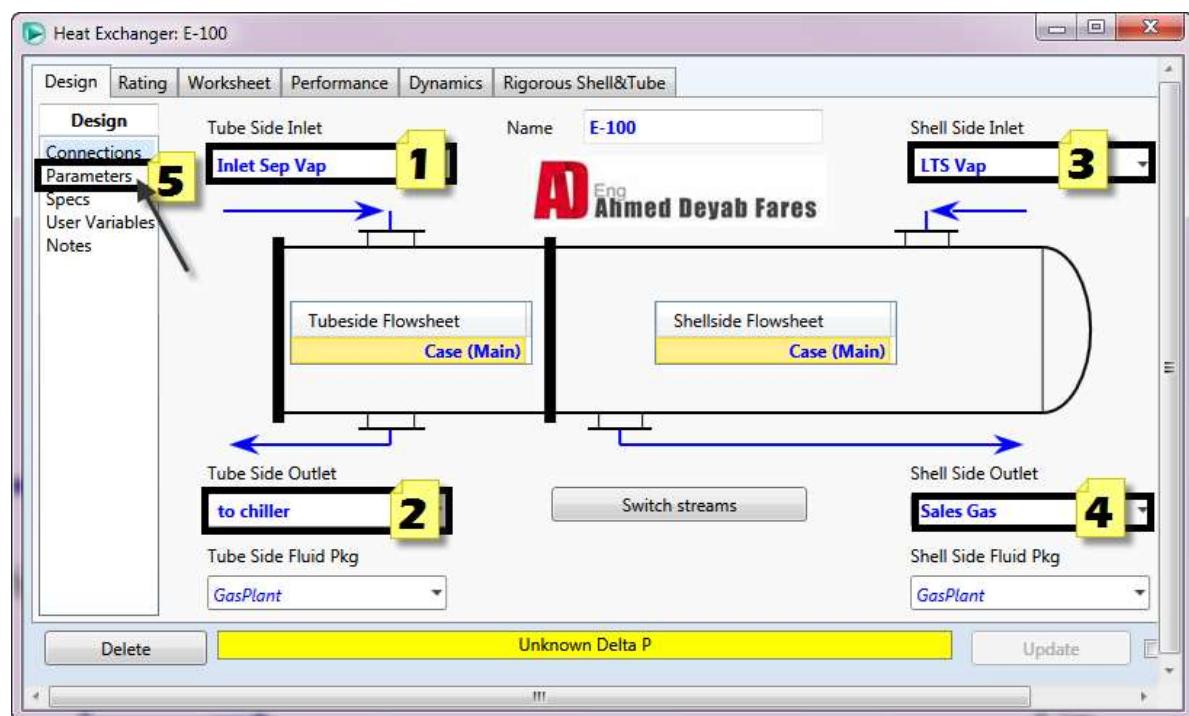
Calculate the Overall Heat Transfer Coefficient (UA) for the HX .....

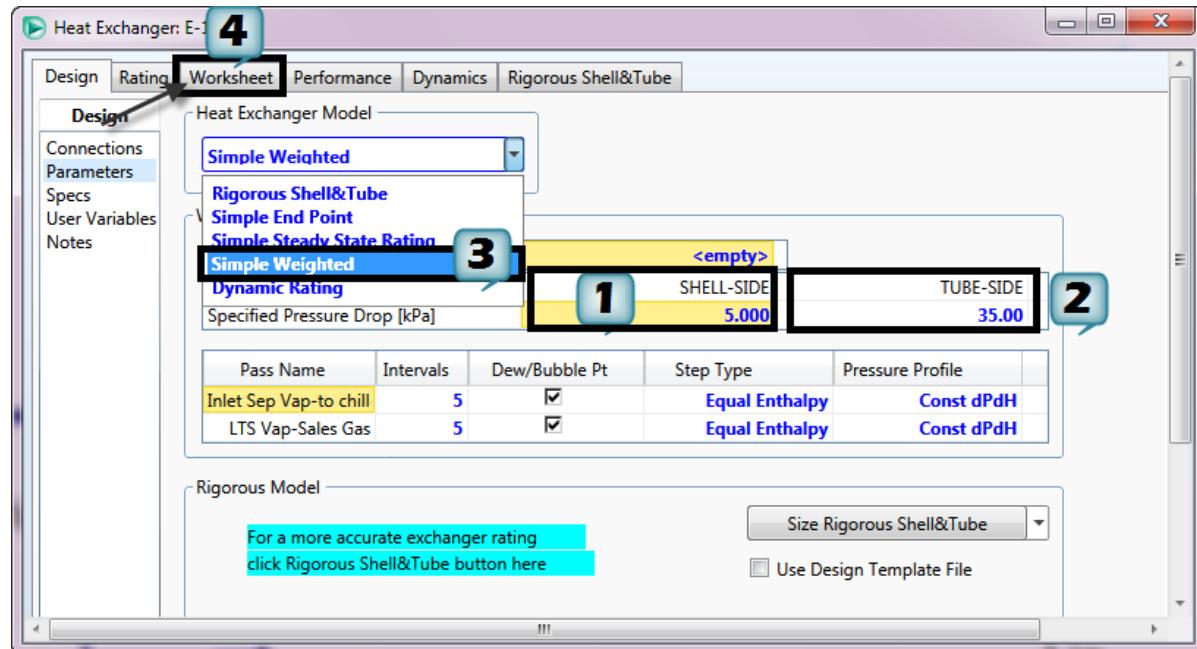
Solution:

- 1- Remove the heater, outlet stream and heater energy stream.
- 2- Disconnect the chiller inlet from the chiller (cooler).
- 3- Add a heat exchanger from the palette:



- 4- Open the heat exchanger and complete the required data:

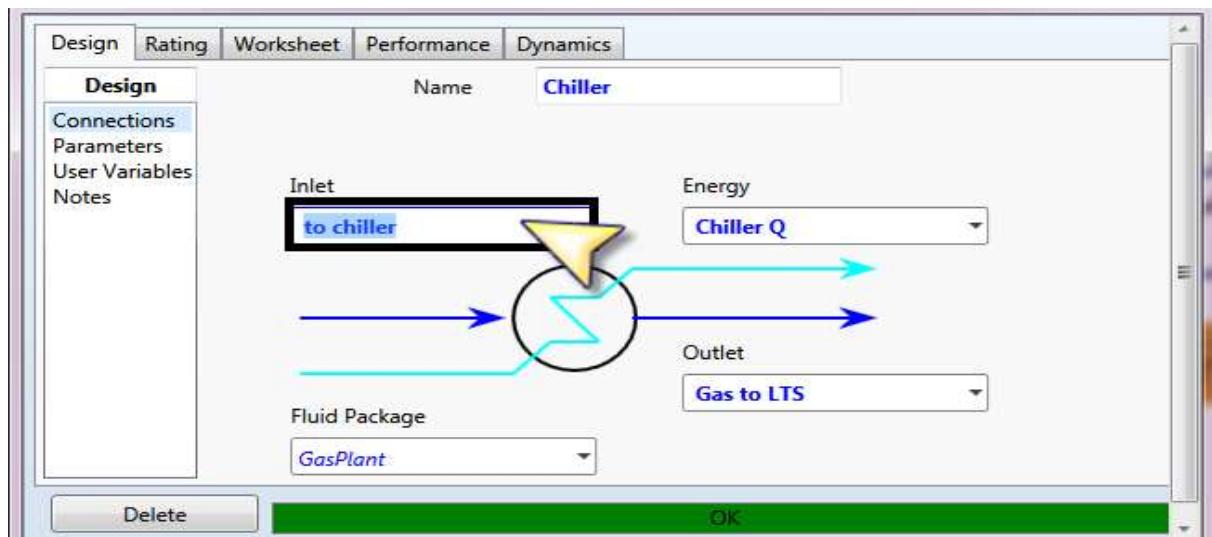




- **Weighted.** The heating curves are broken into intervals, which then exchange energy individually. An LMTD and UA are calculated for each interval in the heat curve and summed to calculate the overall exchanger UA. The Weighted method is available only for Counter-Current exchangers.
- **Endpoint.** A single LMTD and UA are calculated from the inlet and outlet conditions. For simple problems where there is no phase change and  $C_p$  is relatively constant, this option may be sufficient.

Name	Inlet Sep Vap	to chiller	LTS Vap	Sales Gas
Vapour	1.0000	<empty>	1.0000	<empty>
Temperature [C]	15.00	<empty>	-20.00	10.00
Pressure [kPa]	6200	6165	<empty>	<empty>
Molar Flow [kgmole/h]	1440	1440	<empty>	<empty>
Mass Flow [kg/h]	2.990e+004	2.990e+004	<empty>	<empty>
Std Ideal Liq Vol Flow [m3/h]	88.31	88.31	<empty>	<empty>
Molar Enthalpy [kJ/kgmole]	-8.127e+004	<empty>	<empty>	<empty>
Molar Entropy [kJ/kgmole-C]	149.4	<empty>	<empty>	<empty>
Heat Flow [kJ/h]	-1.170e+008	<empty>	<empty>	<empty>

5- Open the chiller and re-connect the tube side outlet to the chiller inlet



The duty rejected from the chiller after this modification = **1.878 e6** kJ/hr

The Overall Heat Transfer Coefficient (UA) for the HX= **2.786 e5** kJ/C-h

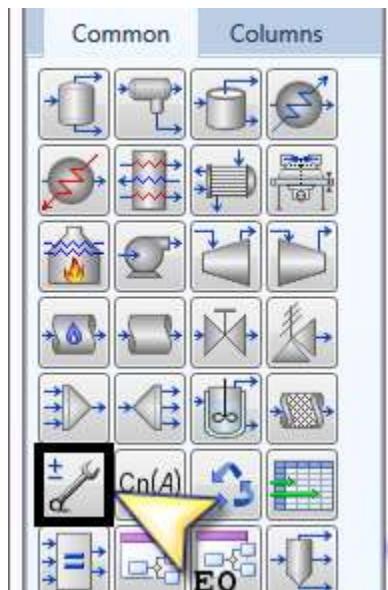
	Before Modification	After Modification
<b>Chiller Duty</b>	<b><math>4.186 \times 10^6</math></b>	<b><math>1.878 \times 10^6</math></b>
<b>Heater Duty</b>	<b><math>2.287 \times 10^6</math></b>	<b>0</b>

# Adjust

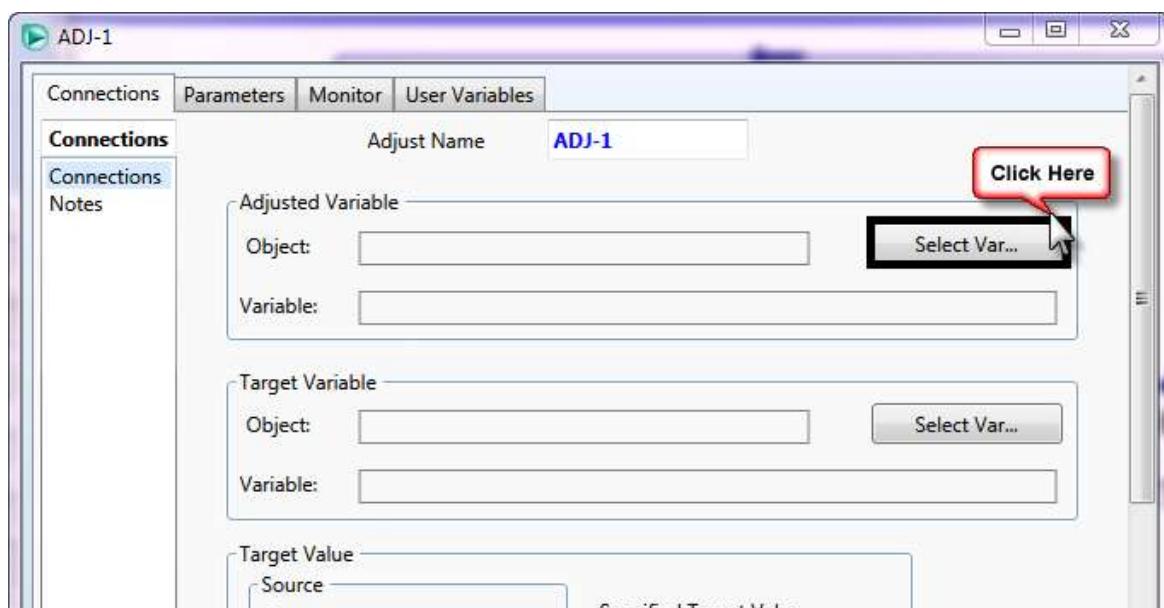
Adjust the LTS feed temperature to ensure the LTS vapor rate of 1200 kgmole/hr using Adjust operation.

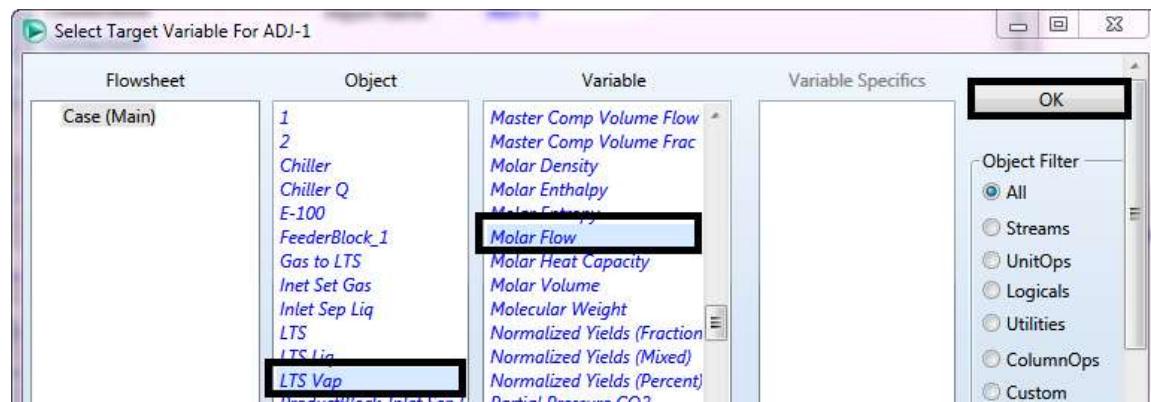
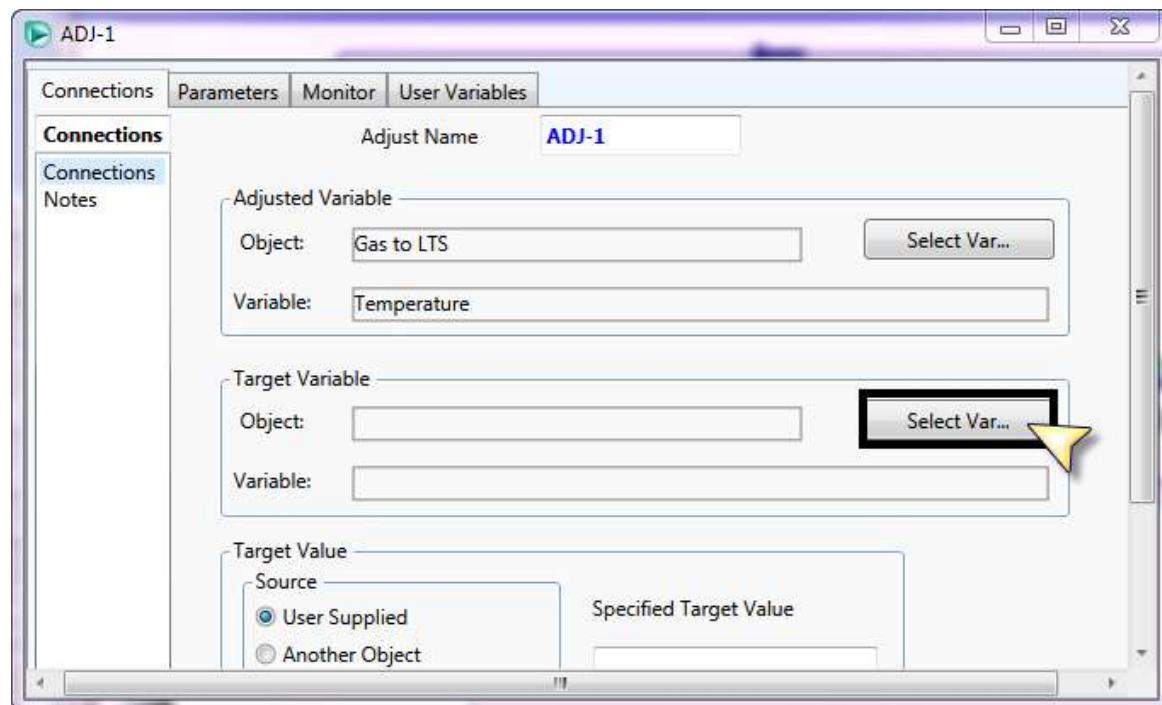
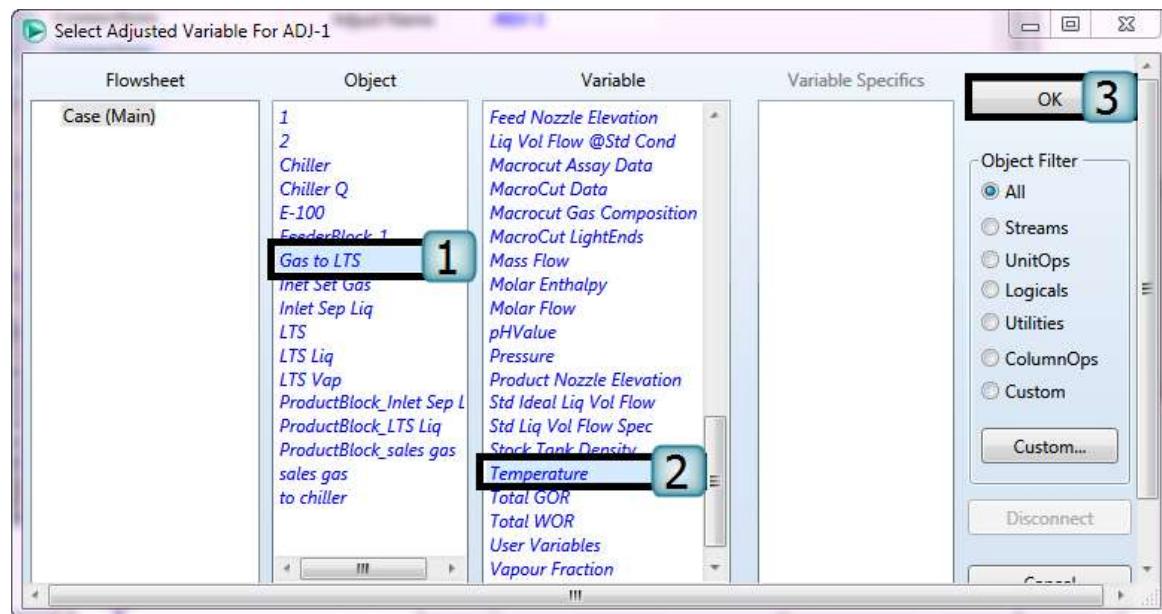
Calculate the temperature of LTS feed .....

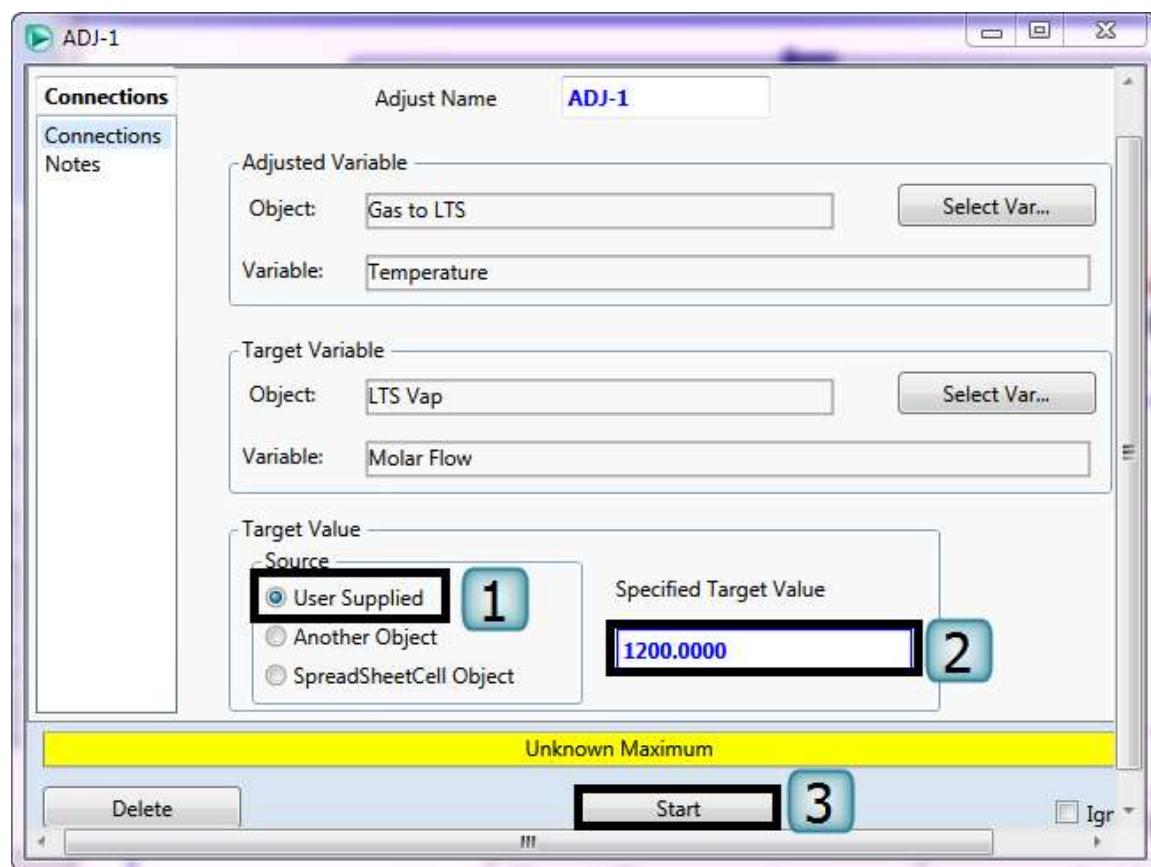
- 1- Select an adjust operation from the palette:



- 2- Open the adjust operation and select the adjusted variable (LTS feed Temperature) and the Target variable (LTS Vapor molar flow).







You can see the total number of iterations from the monitor tab:

Iteration History			
Total Iterations 7			
Iter	Adjusted Value [C]	Target Value [kgmole/h]	Residual [kgmole/h]
1	-20.000	1259.243	59.24
2	-21.000	1247.547	47.55
3	-21.445	1242.197	42.20
4	-31.178	1099.027	-101.0
5	-23.008	1222.687	22.69
6	-24.507	1202.861	2.861
7	-24.723	1199.905	-9.471e-002

# Propane Refrigeration Loop

Experience the New Aspen HYSYS®.

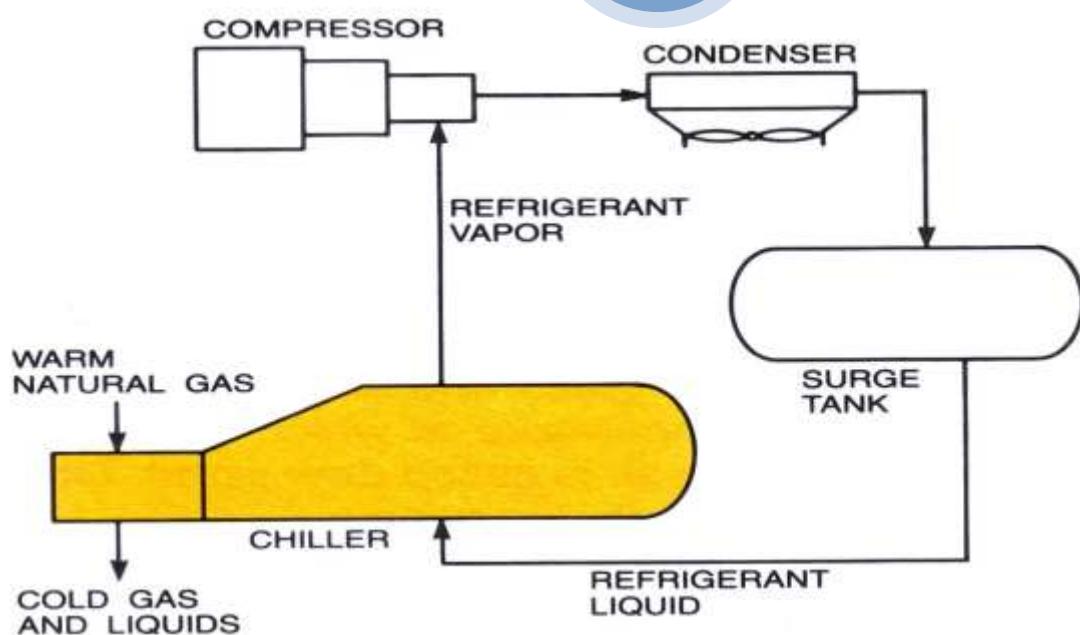
V8

The best process simulation software is now easier to use and faster to learn than ever!

Aspen HYSYS is a comprehensive process modeling system used by the world's leading oil & gas producers, refineries, and engineering companies to optimize process design and operations.



3



## **Workshop**

*Refrigeration systems are commonly found in the natural gas processing industry and in processes related to the petroleum refining, petrochemical, and chemical industries. Refrigeration is used to cool gas to meet a hydrocarbon dewpoint specification and to produce a marketable liquid.*

*In this module you will construct, run, analyze and manipulate a propane refrigeration loop simulation. You will convert the completed simulation to a template, making it available to connect to other simulations.*

## **Learning Objectives**

*Once you have completed this module, you will be able to:*

- *Add and connect operations to build a flowsheet.*
- *Understand how to simulate the vapor compression loop.*
- *Understand forward-backward information propagation in HYSYS.*
- *Using the spread sheet to calculate the COP (Coefficient Of Performance) for the loop.*

Example:

A Refrigeration cycle utilizes propane as the working fluid is used in the liquefaction of the NG. Propane is fed to an evaporator (Heater) the pressure drop=5 kPa, where it absorbed 1.50e+6 kJ/hr from the NG and leaves at the dew point (Vapor Fraction=1.0) at T= -15°C. The output of the evaporator is then compressed adiabatically with efficiency of 75%, and then it's condensed to reject heat. Inside the Condenser there is a pressure drop of 30 kPa, and leaves as saturated liquid at 45°C. Finally, the propane passes through a valve to return the pressure of the Evaporator.

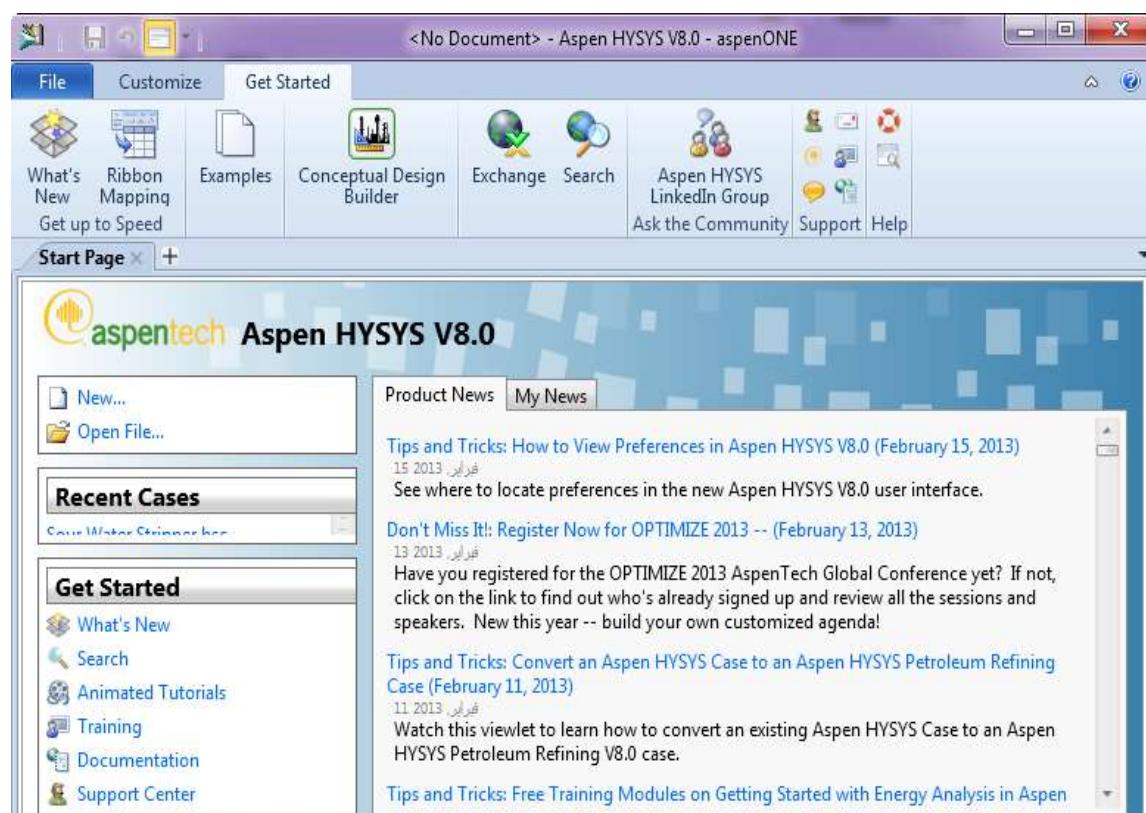
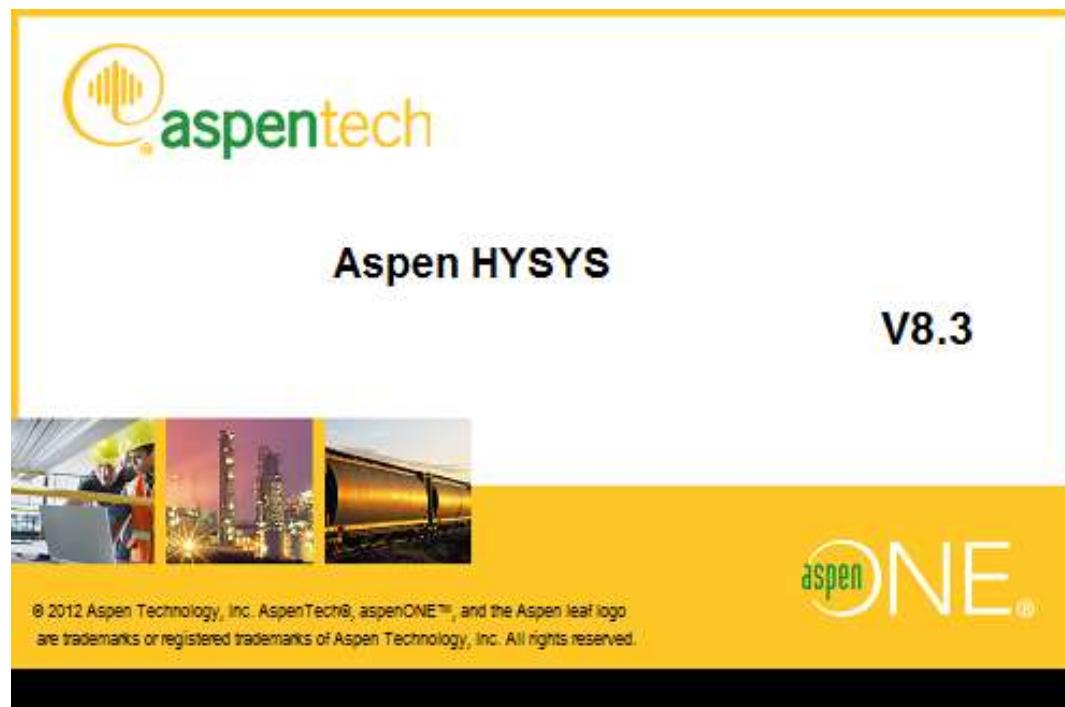
Fluid Pkg: Peng Robinson

Calculate:

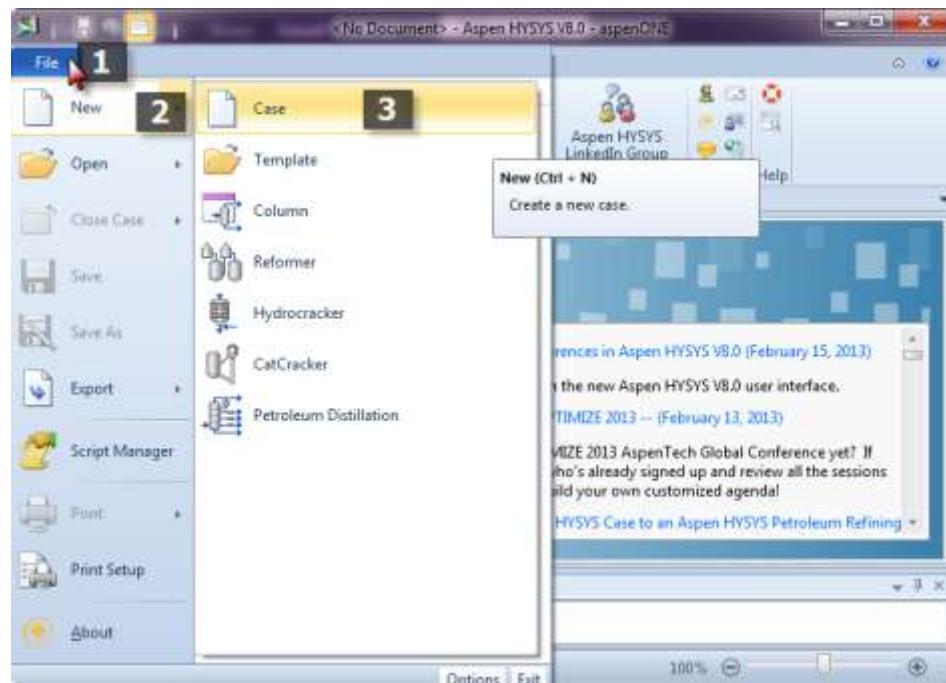
Pressure of the evaporator fed in kPa.	
Flow rate of propane in kmol/hr.	
Valve pressure drop in kPa.	
Temperature of the valve outlet in °C.	
Compressor duty in hp.	
Condenser duty in kJ/hr.	

\* Calculate the COP (Coefficient of Performance) for the cycle

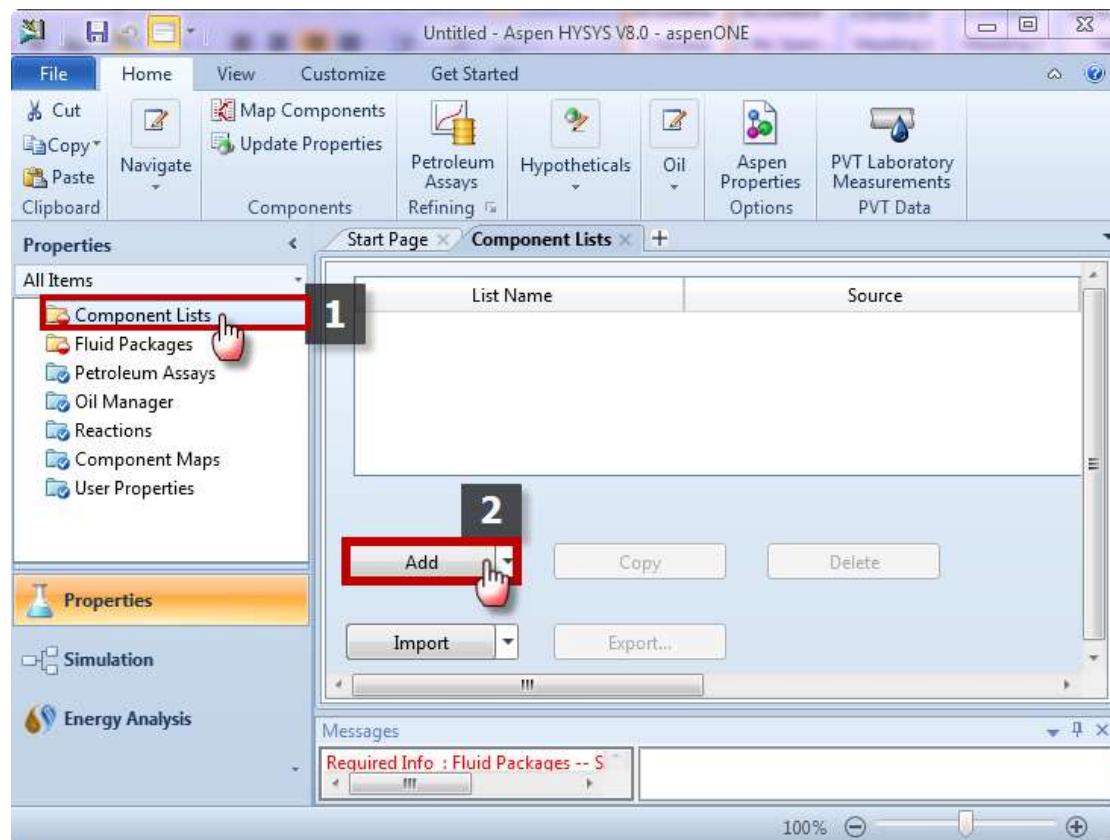
To start the program, From Start Menu, Select All Programs >>  
Aspen Tech >> Process Modeling V8.3 >>> Aspen HYSYS >>  
Aspen HYSYS



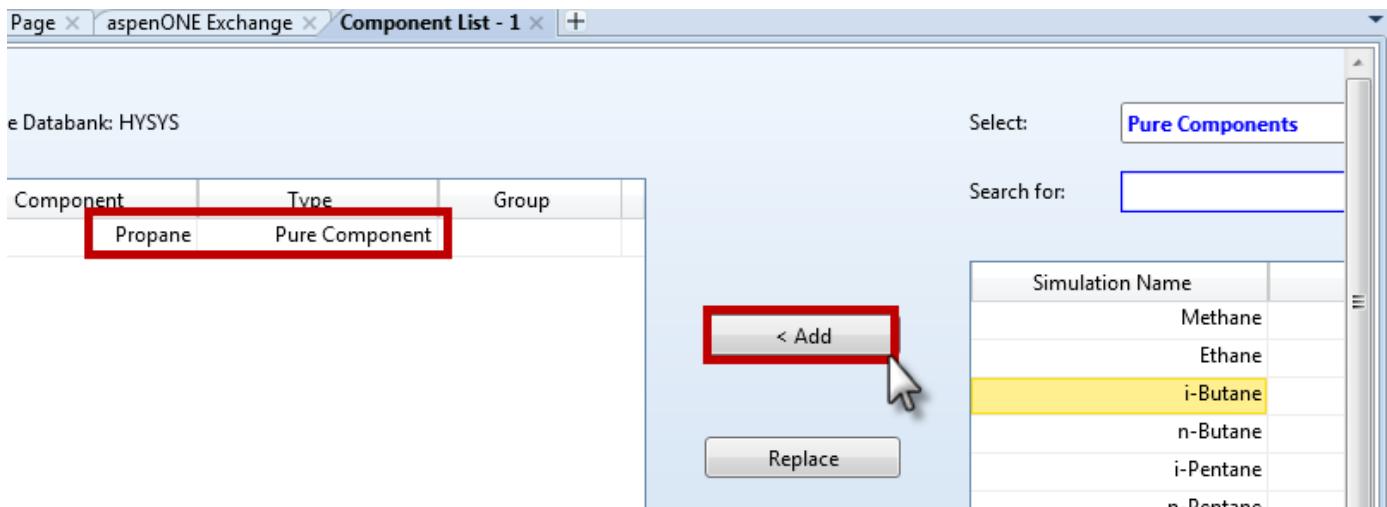
## 1- First, Start a new case



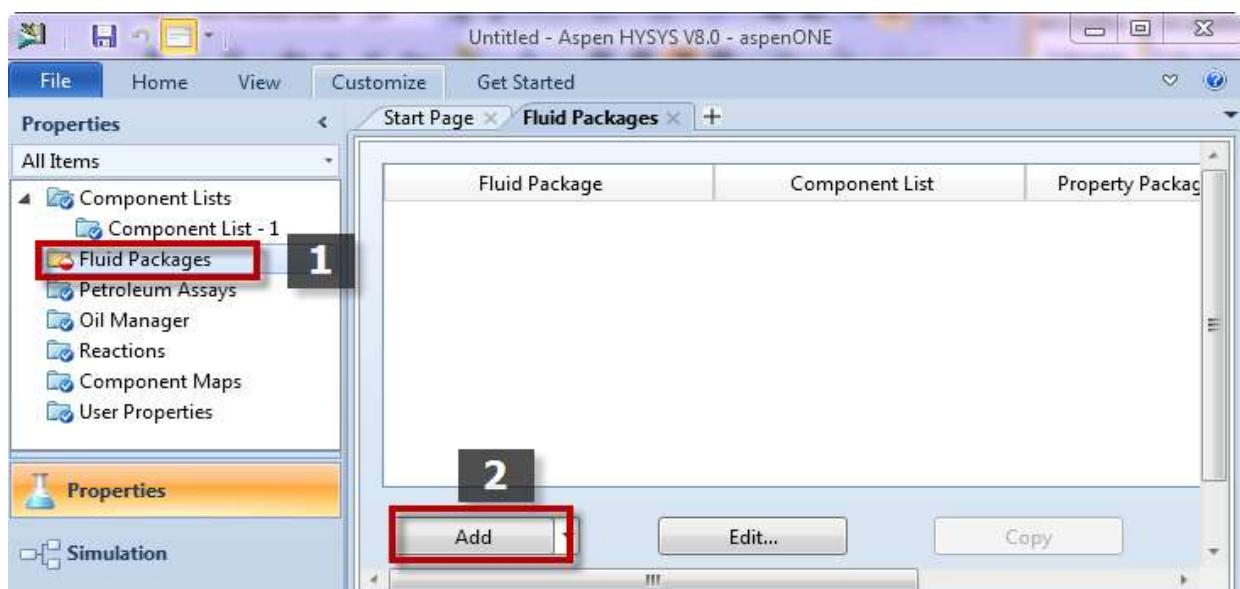
## 2- Add the Components



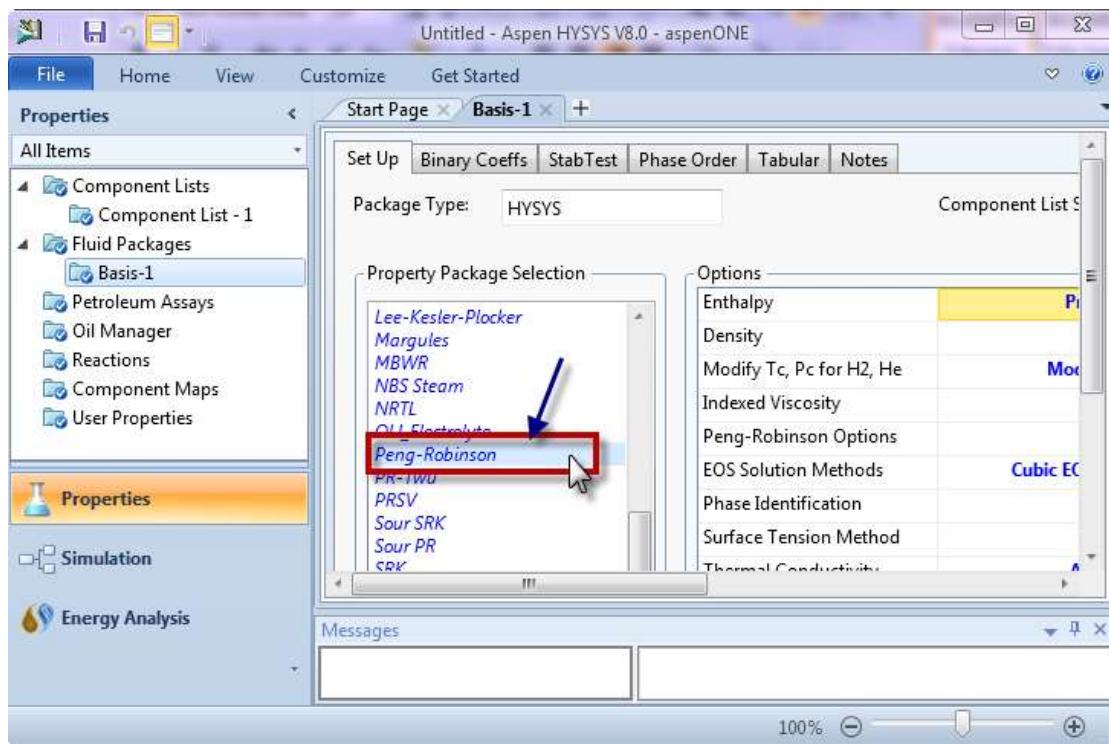
3- Choose the system components from the databank:



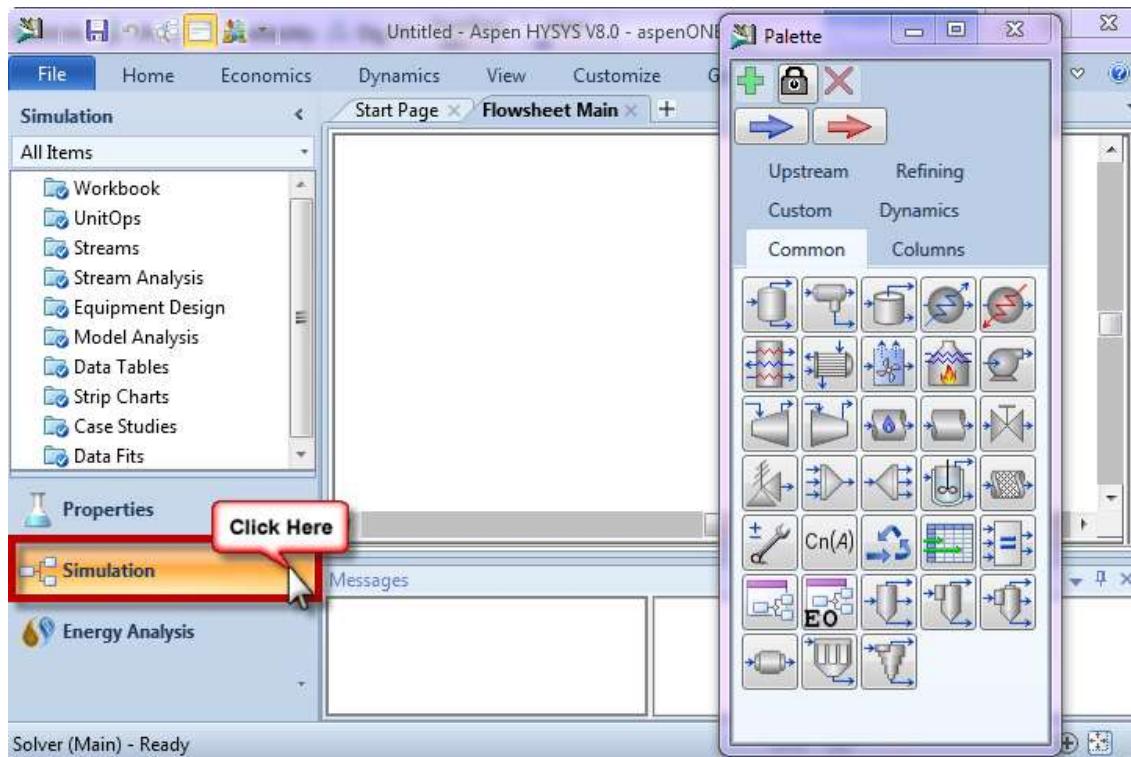
Now, select the suitable fluid package



In this case, select Peng-Robinson

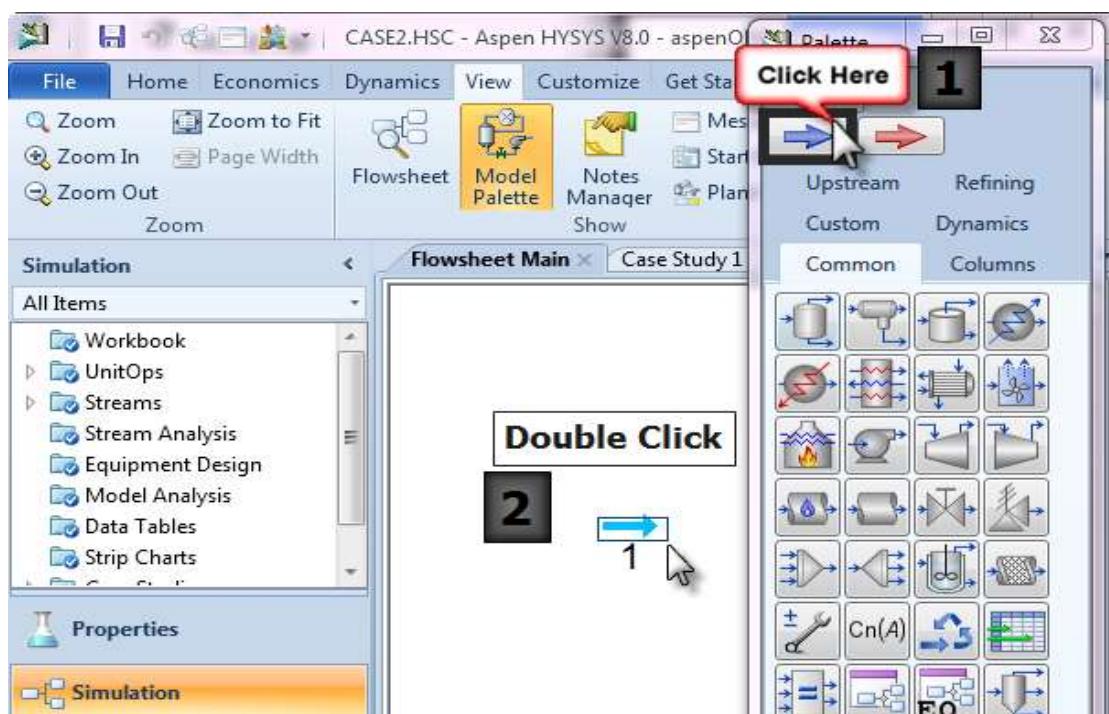


Now you can start drawing the flow sheet for the process by clicking the Simulation button:

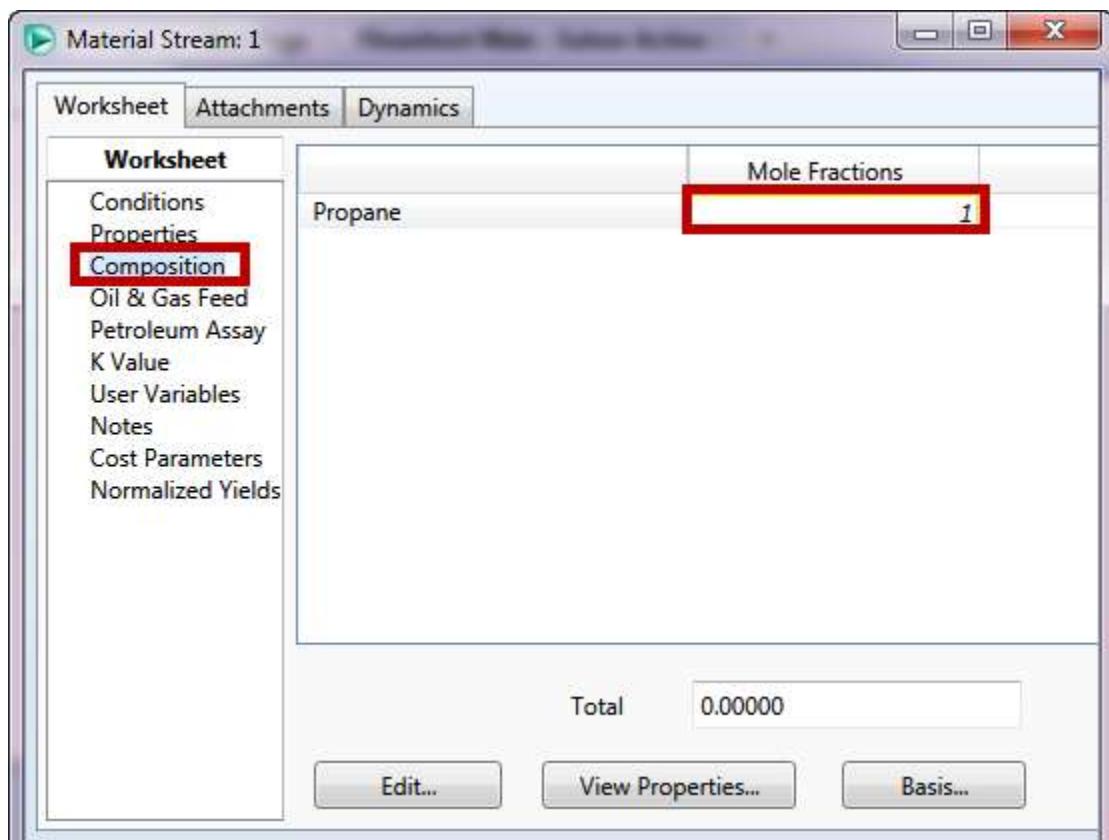


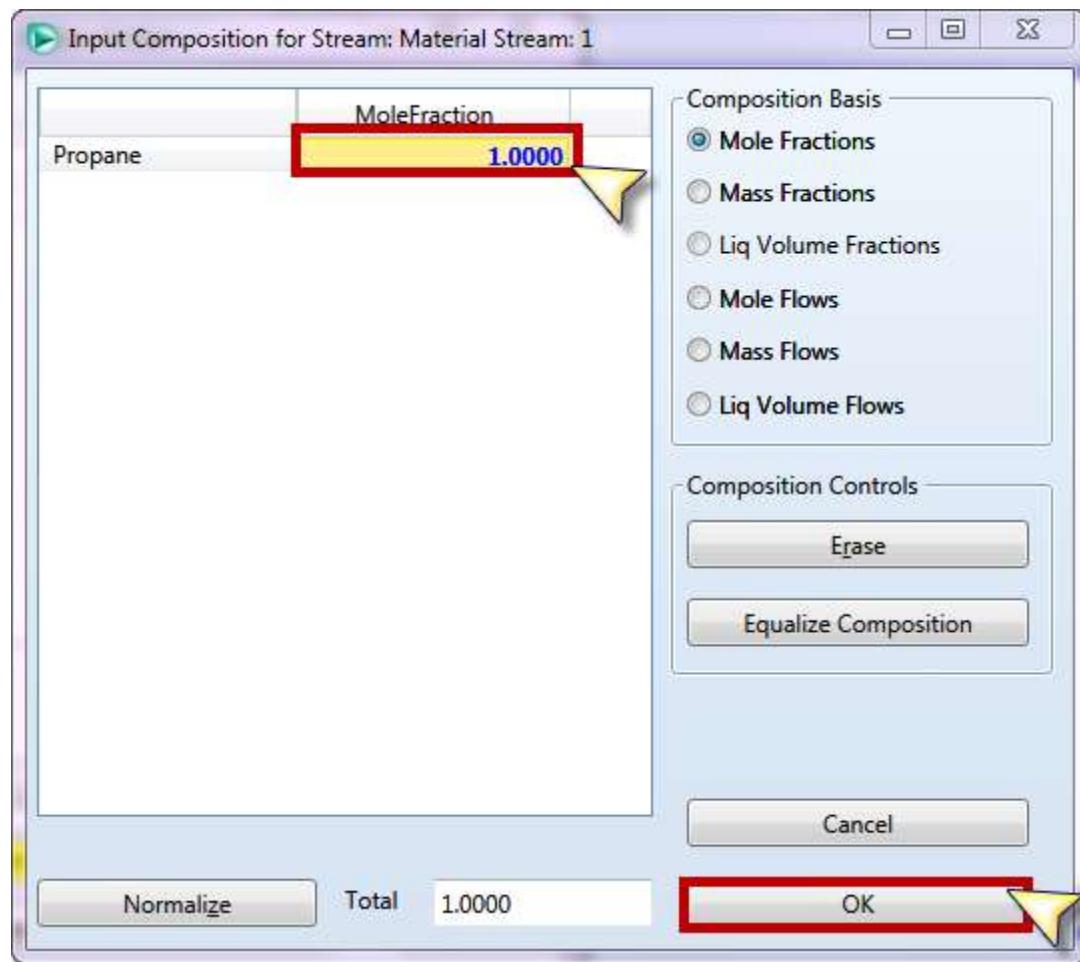
Now add a material stream to define the composition and the conditions of the feed stream

From the palette:



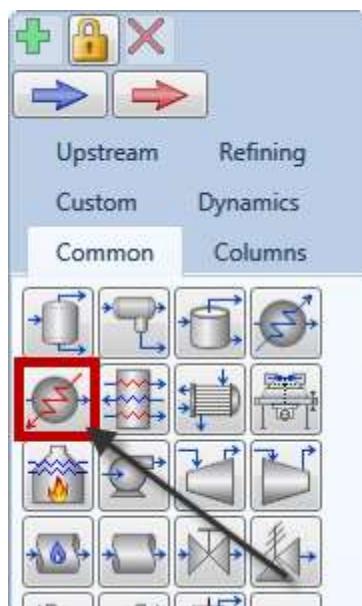
Add the mole fraction for the inlet stream (Propane =1)



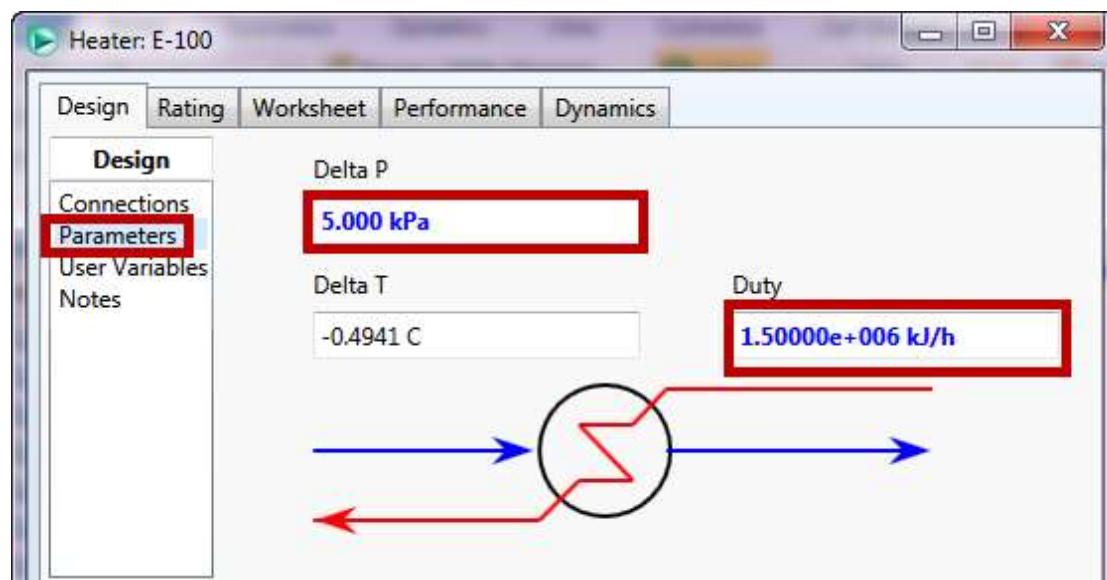
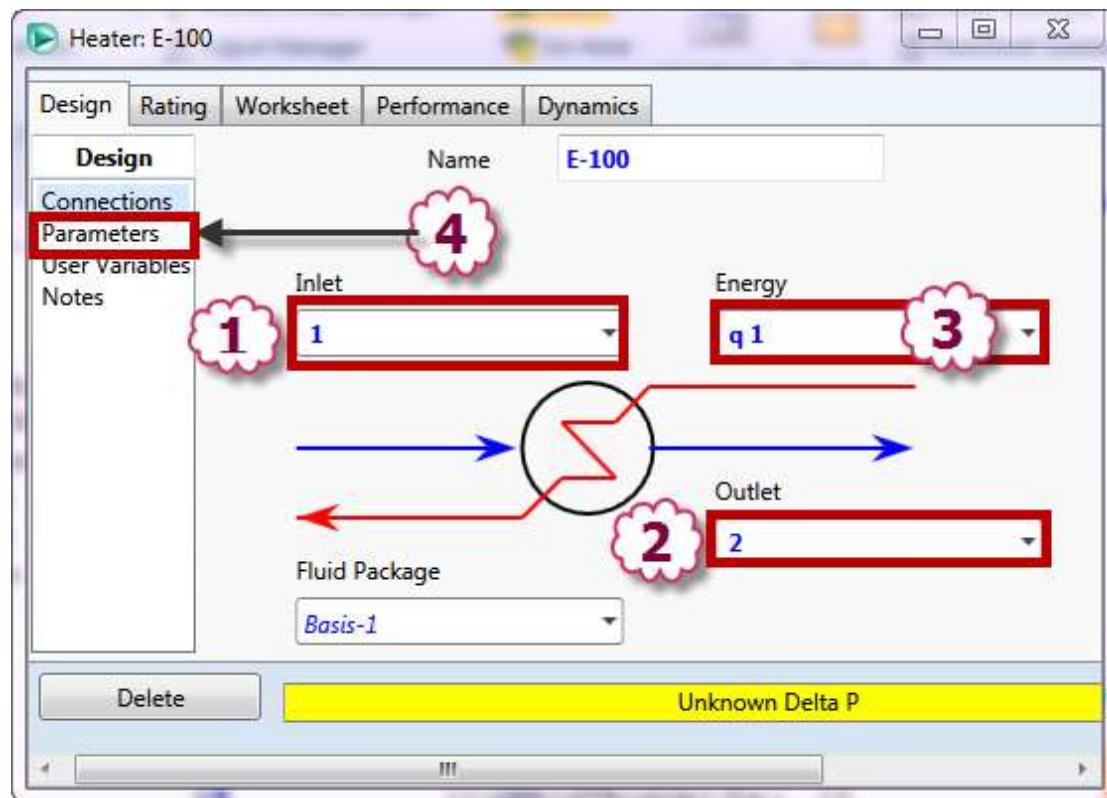


Then leave the stream not solved till the loop is closed

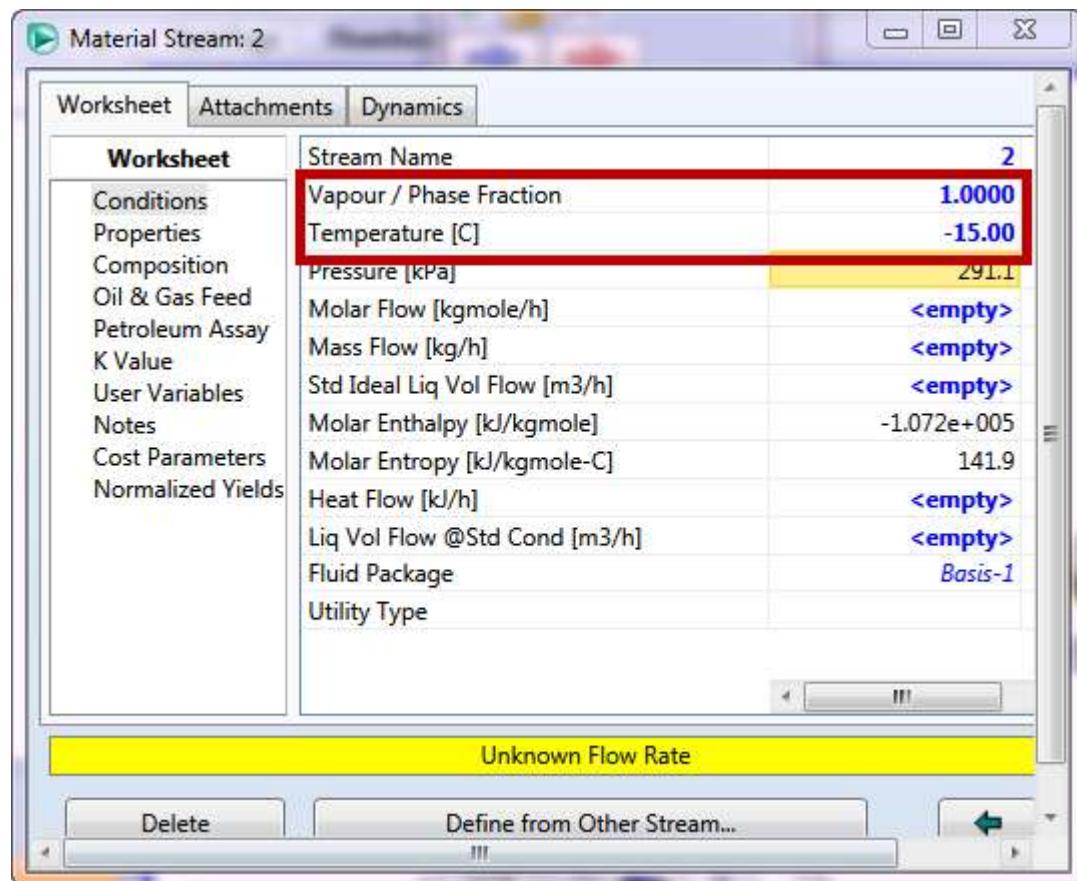
Add the evaporator (heater)



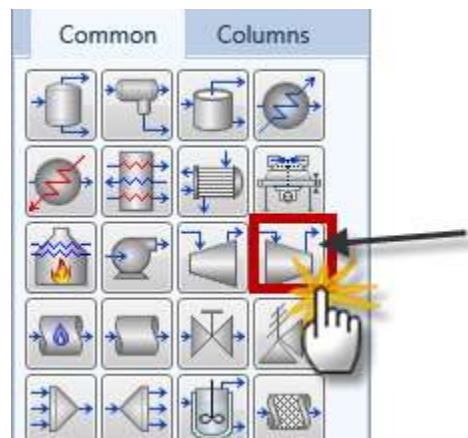
Complete the connections and then go to parameters page to add the pressure drop and the duty rejected

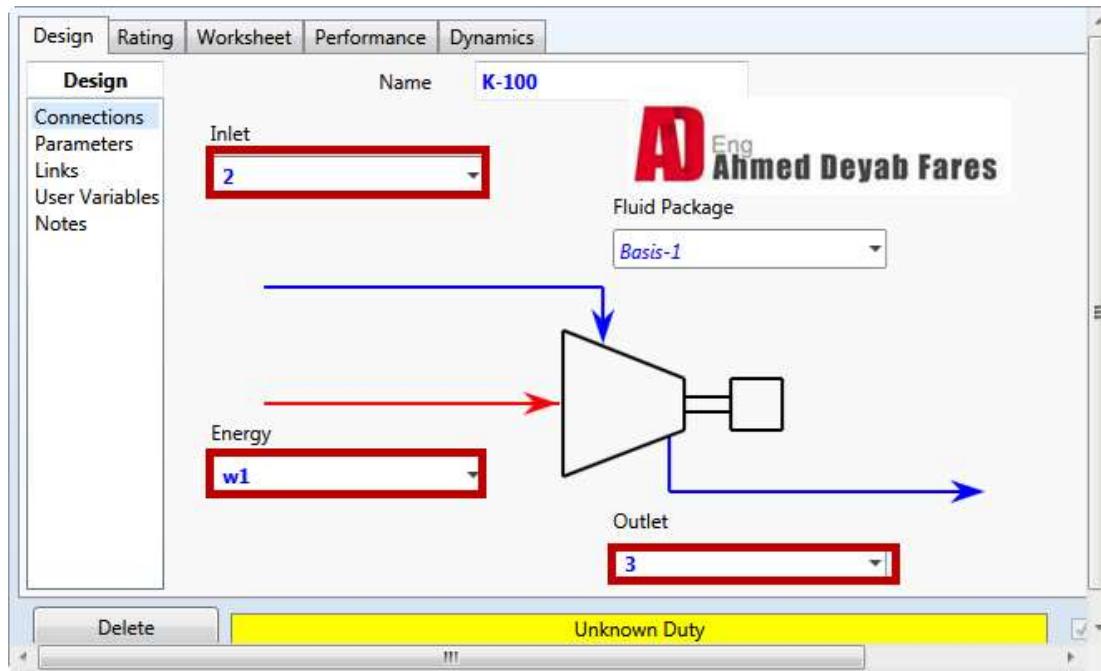


Go to stream 2 and complete the vapor fraction & temperature

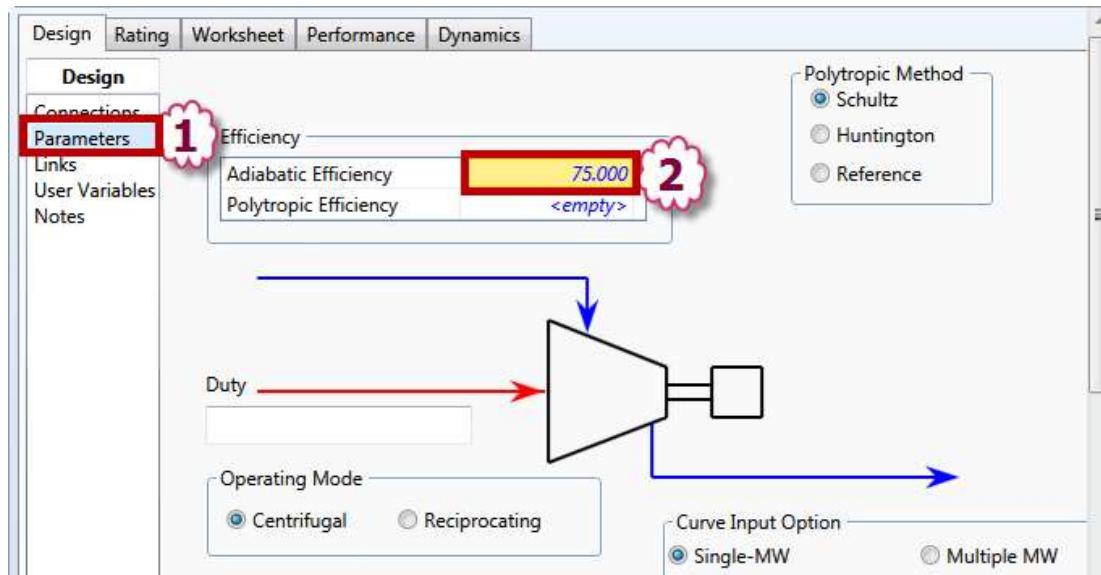


Then add a compressor to raise the pressure of the vapor out from the heater

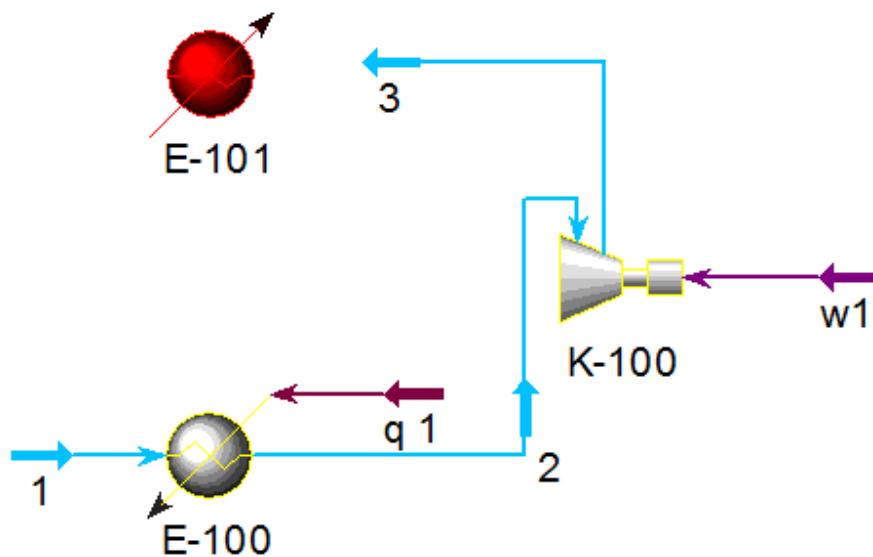
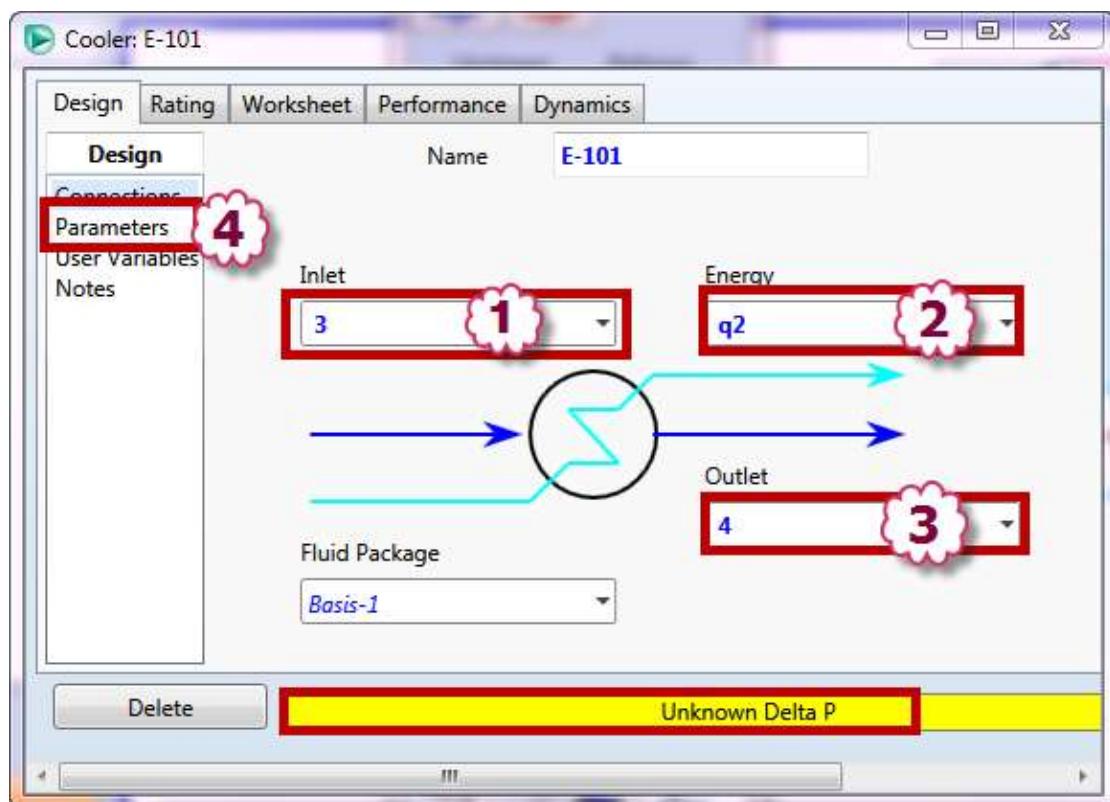




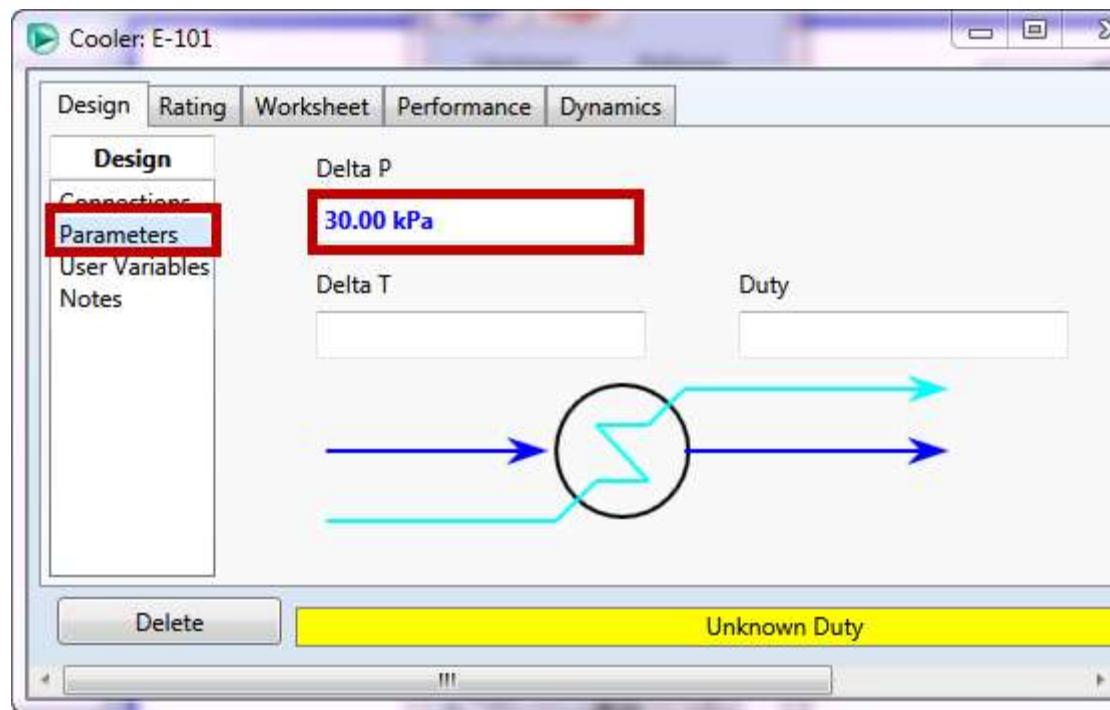
Then go to parameters to make sure that the adiabatic efficiency is 75%



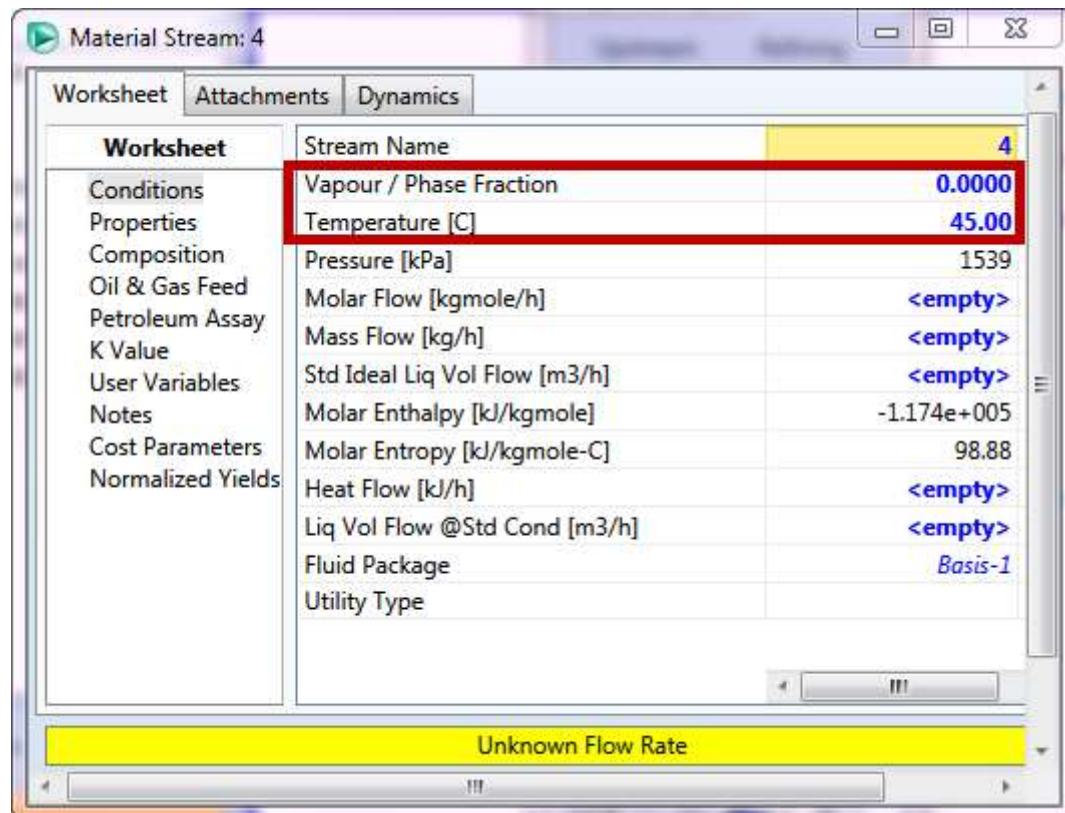
Leave the compressor not solved till the loop is closed then add a cooler


**A** Eng  
**Ahmed Deyab Fares**


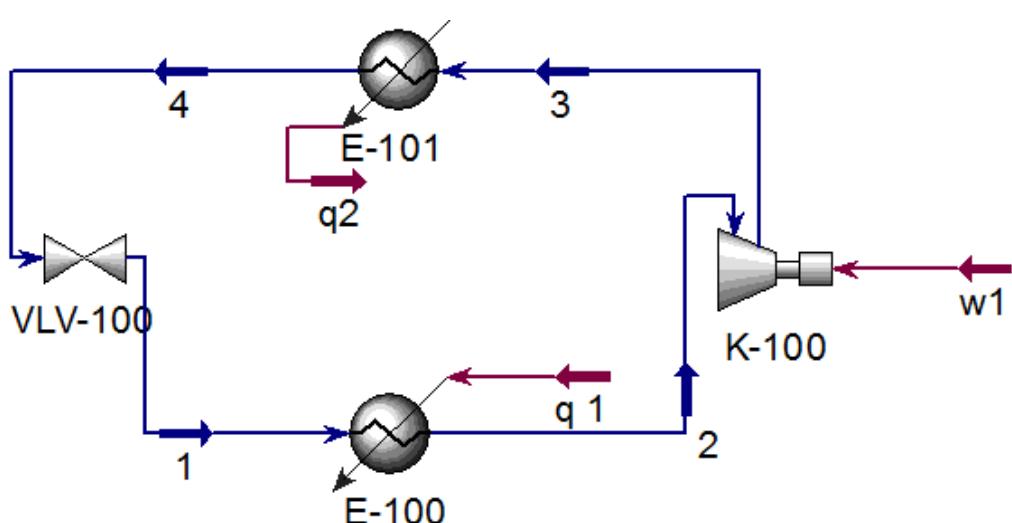
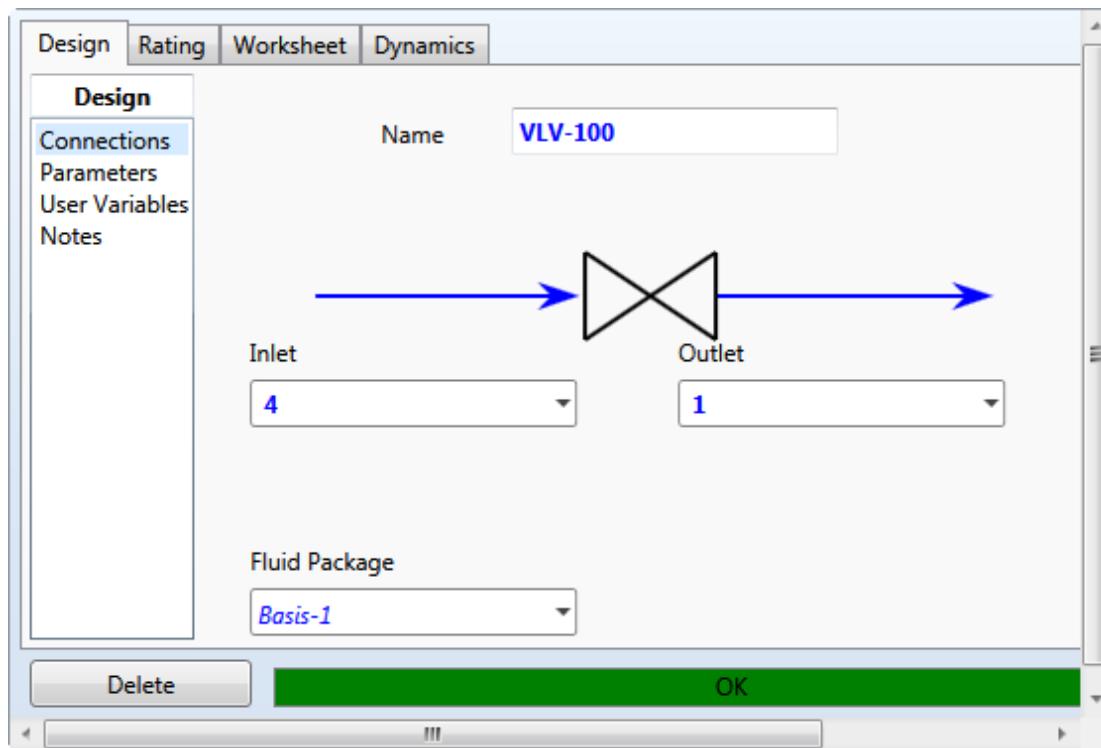
Complete the connections and then go to parameters page to add the pressure drop =30 kPa



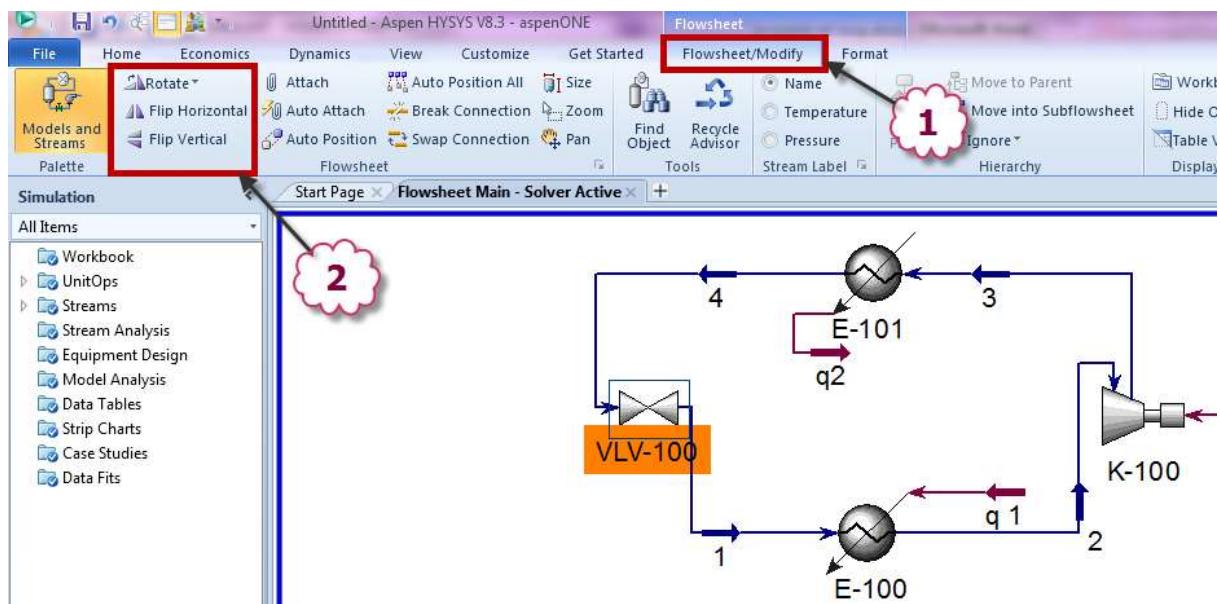
Go to stream 4 and complete the vapor fraction (Saturated liquid=0.0) & temperature (45°C)



Then add a valve to close the loop



You can adjust the Flowsheet and rotate the streams and equipment from the above menu (Flowsheet/Modify)



Results:

Pressure of the evaporator fed in kPa.	296.1
Flow rate of propane in kmol/hr.	146.4
Valve pressure drop in kPa.	1243.34
Temperature of the valve outlet in °C.	-14.51
Compressor duty in hp.	252.8
Condenser duty in kJ/hr.	2.179 e+6

## Coefficient of Performance (COP)

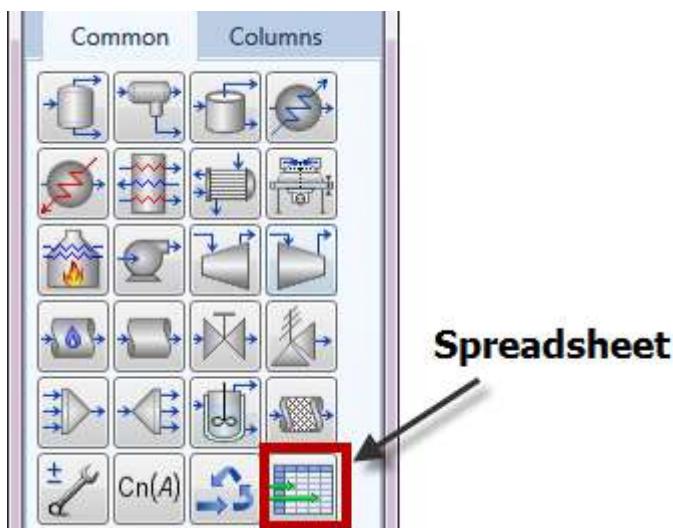
Is a measure of the efficiency of a refrigeration cycle is the *coefficient of performance*, COP

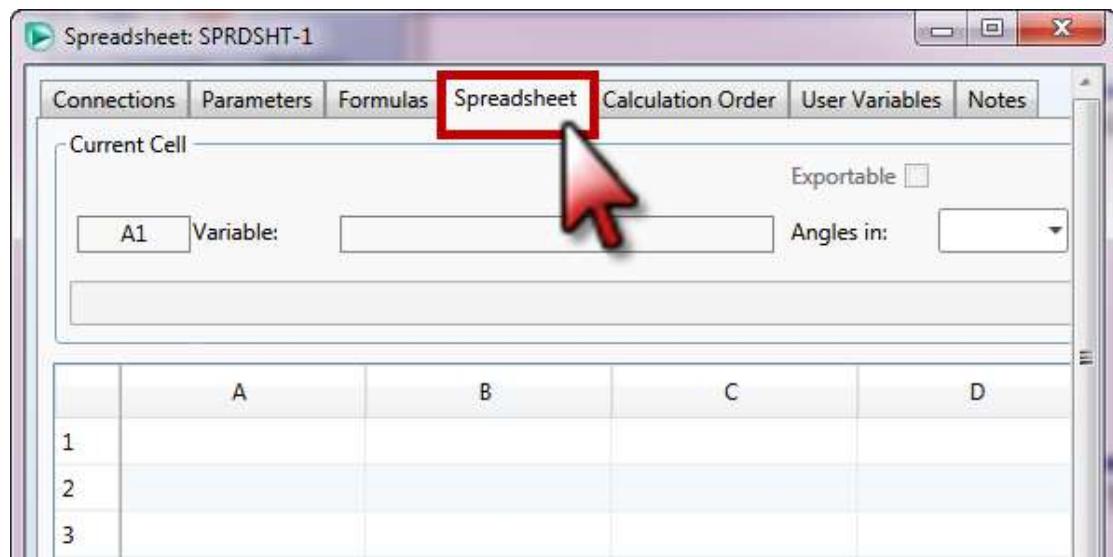
It is the ratio of desired output divided by the required input.

In the vapor-compression system, the net power input is equal to the compressor power, since the expansion valve involves no power input or output.

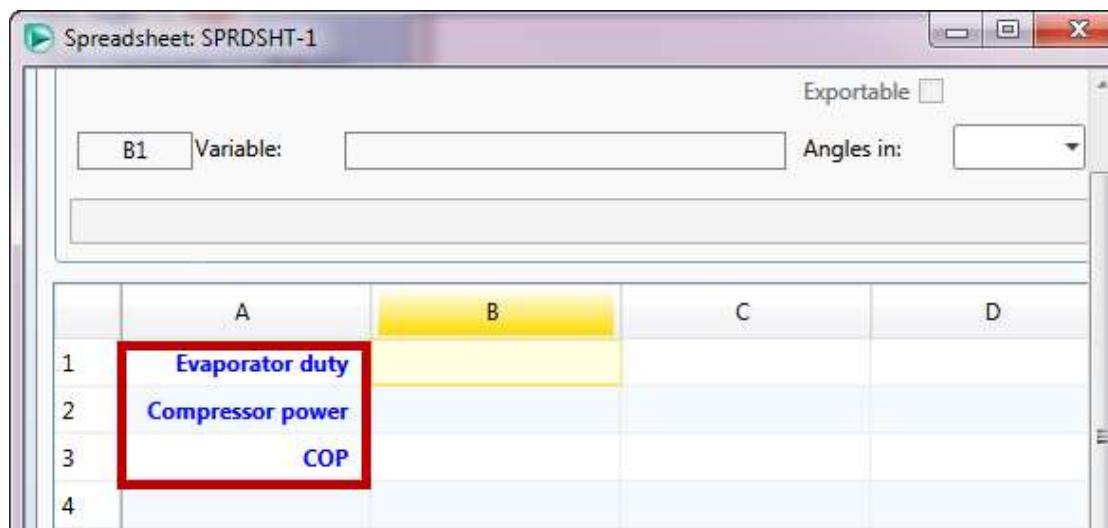
$$\text{COP} = \text{Evaporator Duty} / \text{Compressor Power}$$

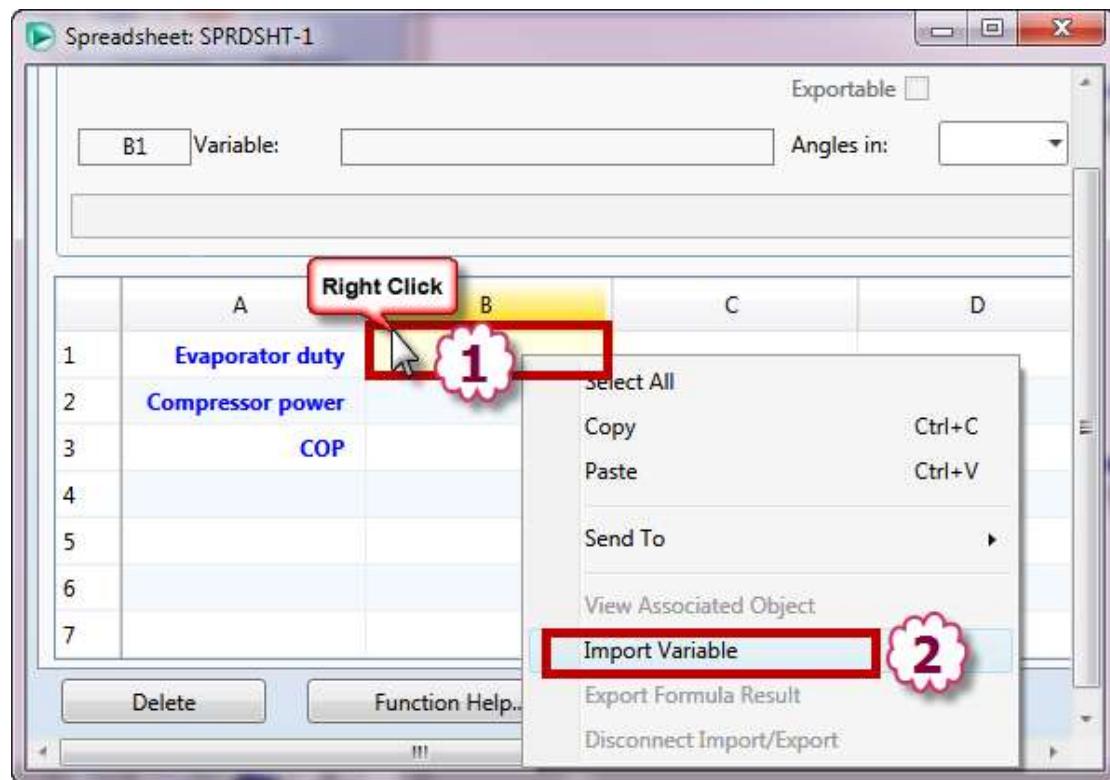
We can use the spreadsheet operation in HYSYS to calculate the COP of the cycle:



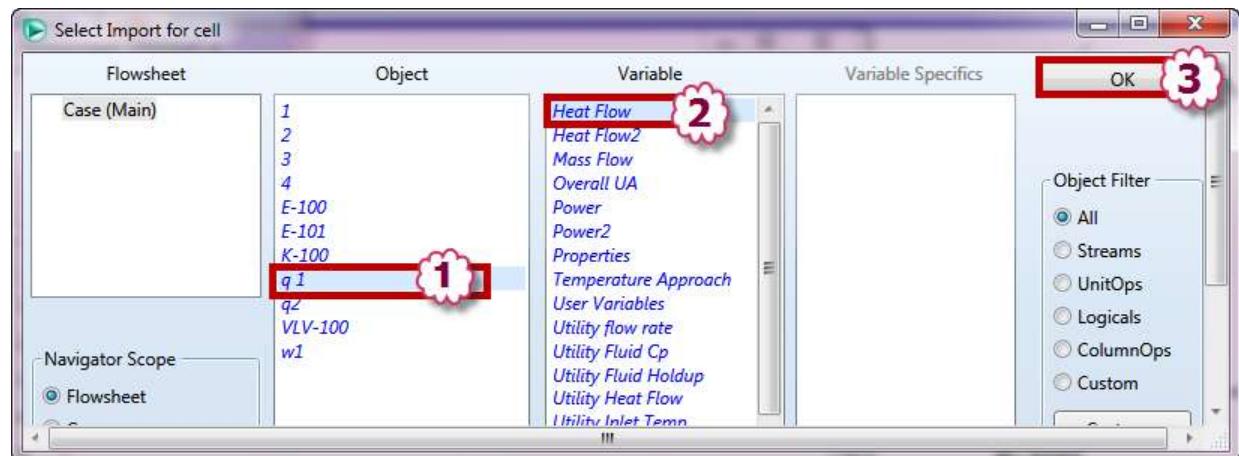


Now we can calculate the COP by importing the two variables;  
Evaporator duty & Compressor power

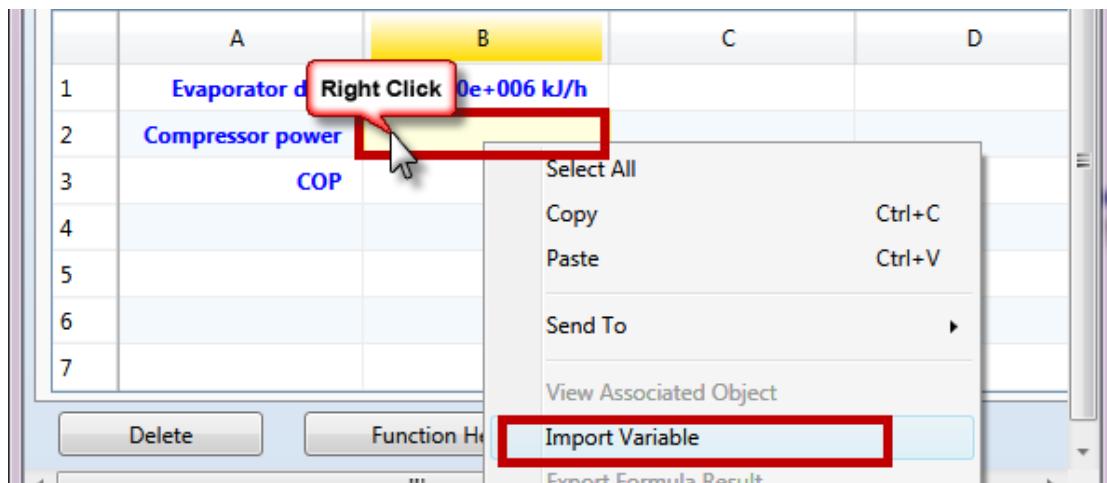




Right click on the cell B1 and select import variable to import the duty of the evaporator



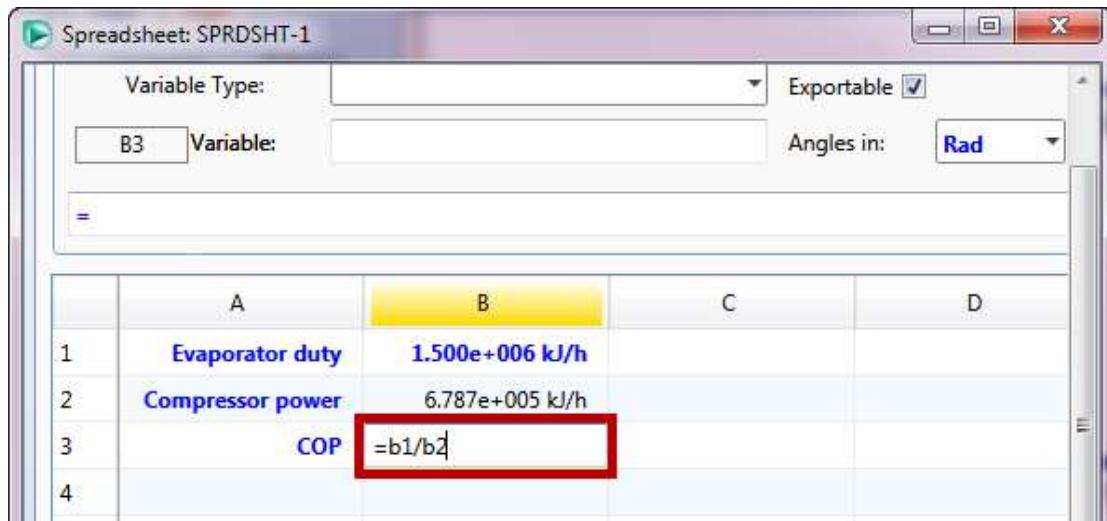
Right click on the cell B2 and select import variable to import the power of the Compressor



Note that the 2 variables must be in the same units (kJ/hr or KW)



Now, divide the two variables in the cells b1 & b2 to calculate the COP in b3



The result will be 2.2

	A	B	C	D
1	Evaporator duty	1.500e+006 kJ/h		
2	Compressor power	6.787e+005 kJ/h		
3	COP	2.210		
4				
5				

**Challenge:**

Now change the duty of the evaporator to 3 e6 kJ/hr (in the evaporator not in the spreadsheet), then open the spreadsheet to calculate the COP & explain the results.

.....  
.....  
.....  
.....  
.....



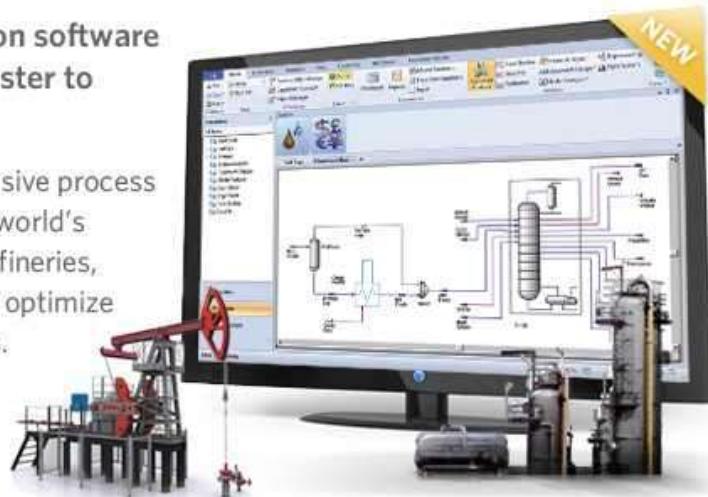
# Distillation Column 4

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is now easier to use and faster to  
learn than ever!

Aspen HYSYS is a comprehensive process  
modeling system used by the world's  
leading oil & gas producers, refineries,  
and engineering companies to optimize  
process design and operations.



## **Workshop**

*Separation of light products is present in any Hydrocarbons operations. In this module, a column will be modeled to separate Light and heavy components from each other using a distillation column with 12 trays.*

## **Learning Objectives**

*Once you have completed this section, you will be able to:*

- *Add columns using the Input Experts.*
- *Add extra specifications to columns.*

**Example:**

We need to separate a mixture of five paraffins into light and heavy fraction by using a distillation column with 12 trays, a full reflux condenser, and a Kettle reboiler.

The feed stream (1000 lbmol/hr) consists of 3% (mole %) ethane, 20% propane, 37% n-butane, 35% n-pentane and 5% n-hexane at 225 °F and 250 psia, which enters the column on the sixth tray, counting from the top. The condenser and reboiler pressures are 248 and 252 psia, respectively. The preliminary design specifications require a **reflux ratio of 6.06** and a **vapor overhead product of 226 lbmol/hr**. Subsequently, the design is modified to ensure propane overhead flow of 191 lb<sub>mol</sub>/hr and n-butane bottom flow of 365 lb<sub>mol</sub>/hr.

**Use SRK Fluid Pkg****Calculate:**

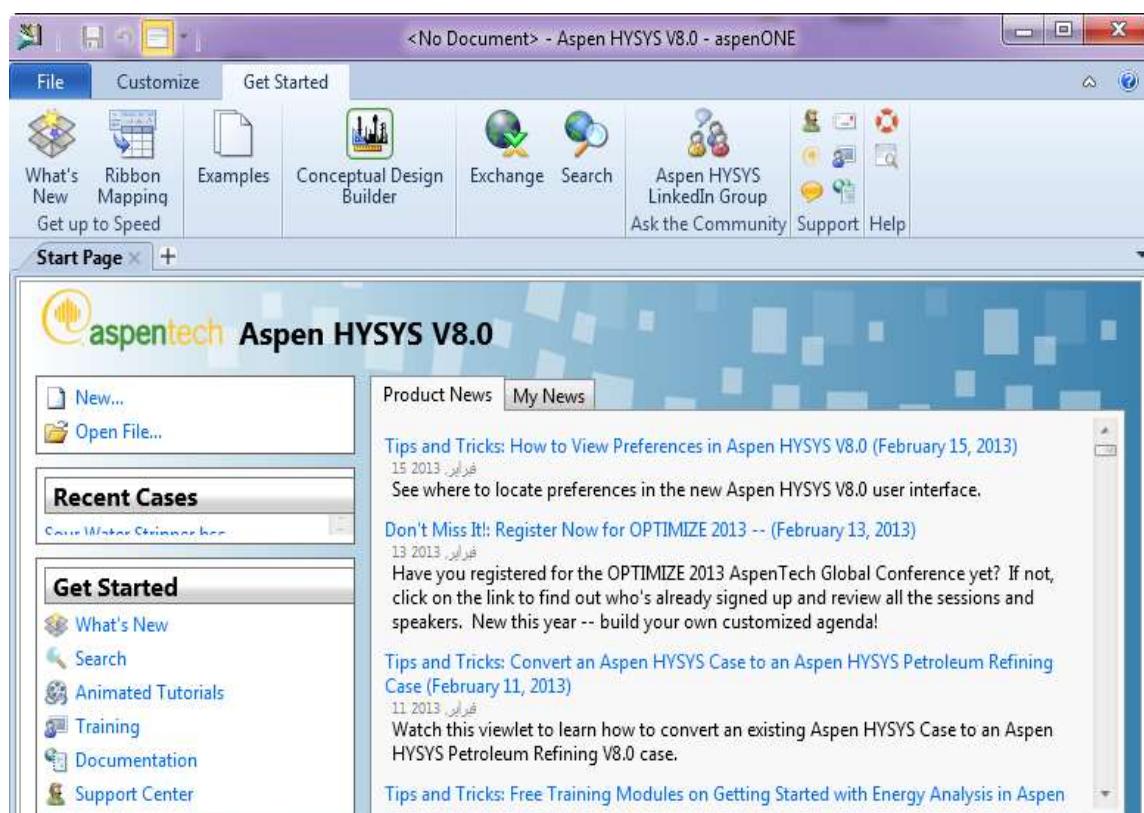
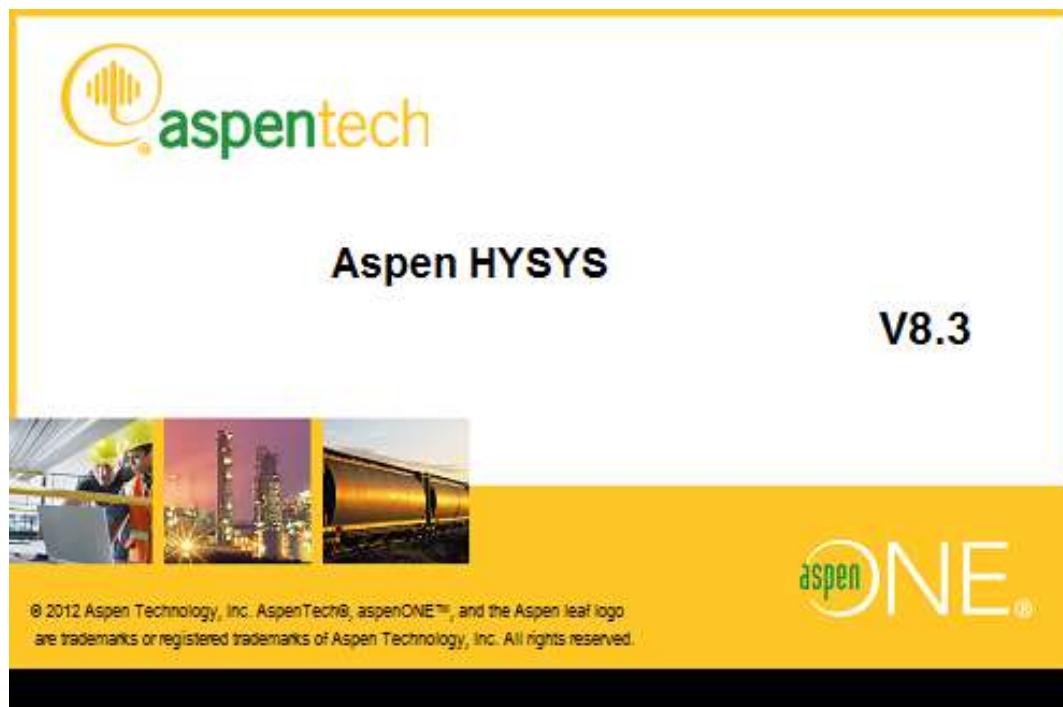
The Condenser, the Reboiler Temperatures & the Reflux Ratio **after modification**

**Condenser Temp** ..... °C

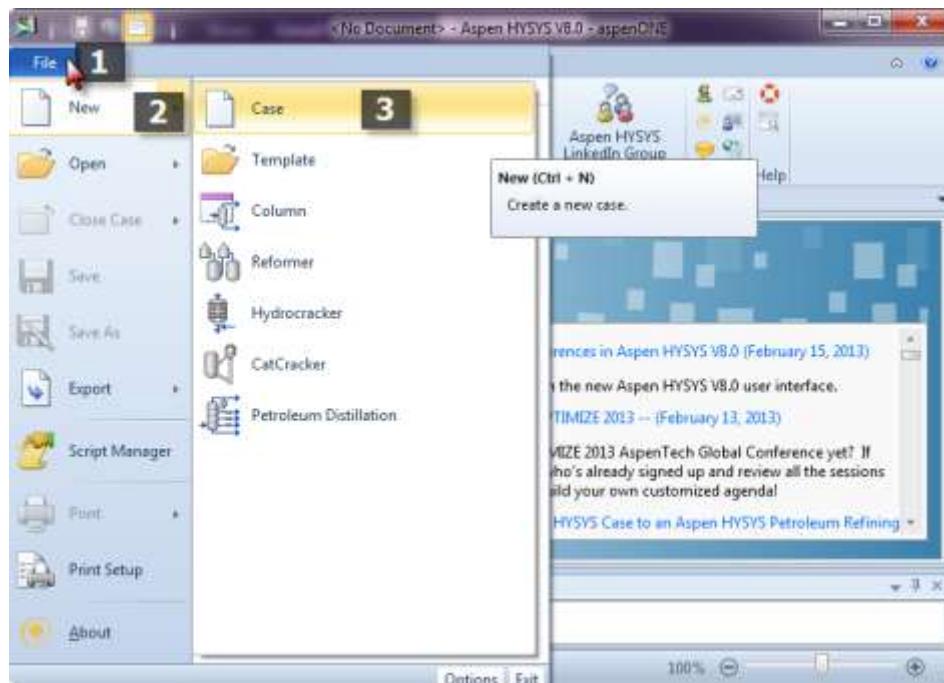
**Reboiler Temp** ..... °C

**Reflux Ratio** .....

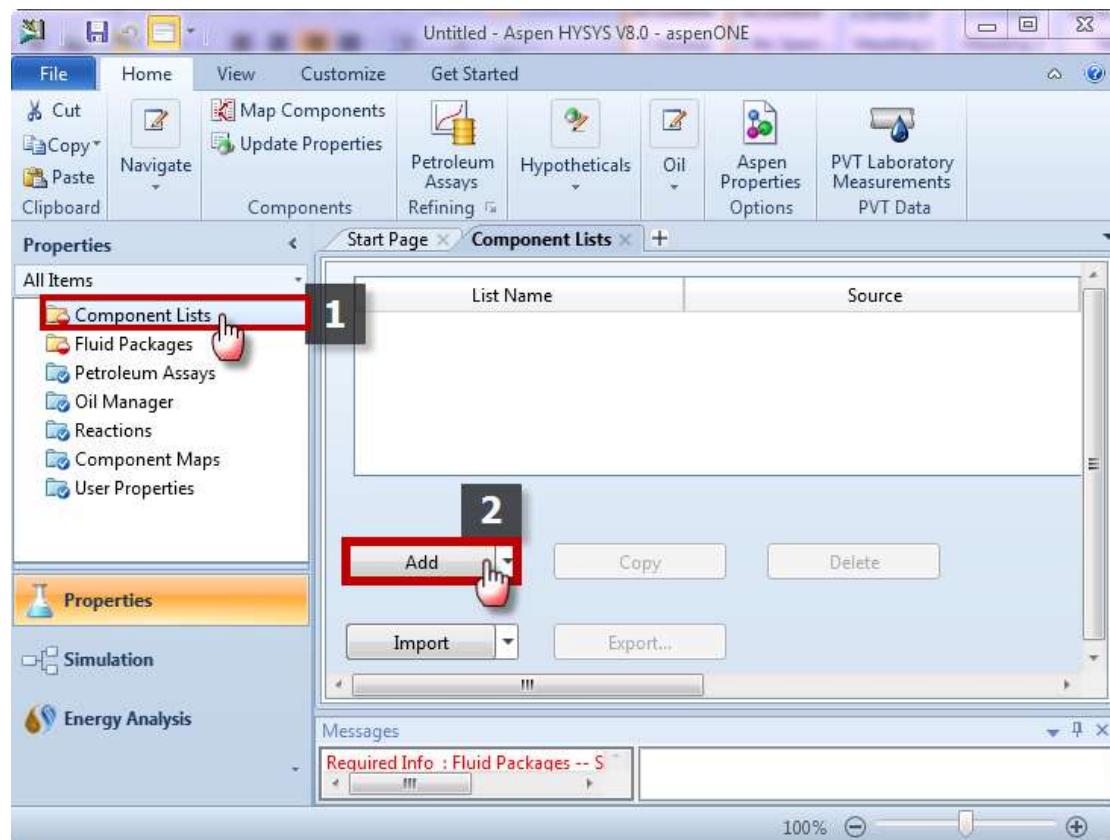
To start the program, From Start Menu, Select All Programs >>  
Aspen Tech >> Process Modeling V8.3 >>> Aspen HYSYS >>  
Aspen HYSYS



## 7- First, Start a new case



## 8- Add the Components



9- Choose the system components from the databank:

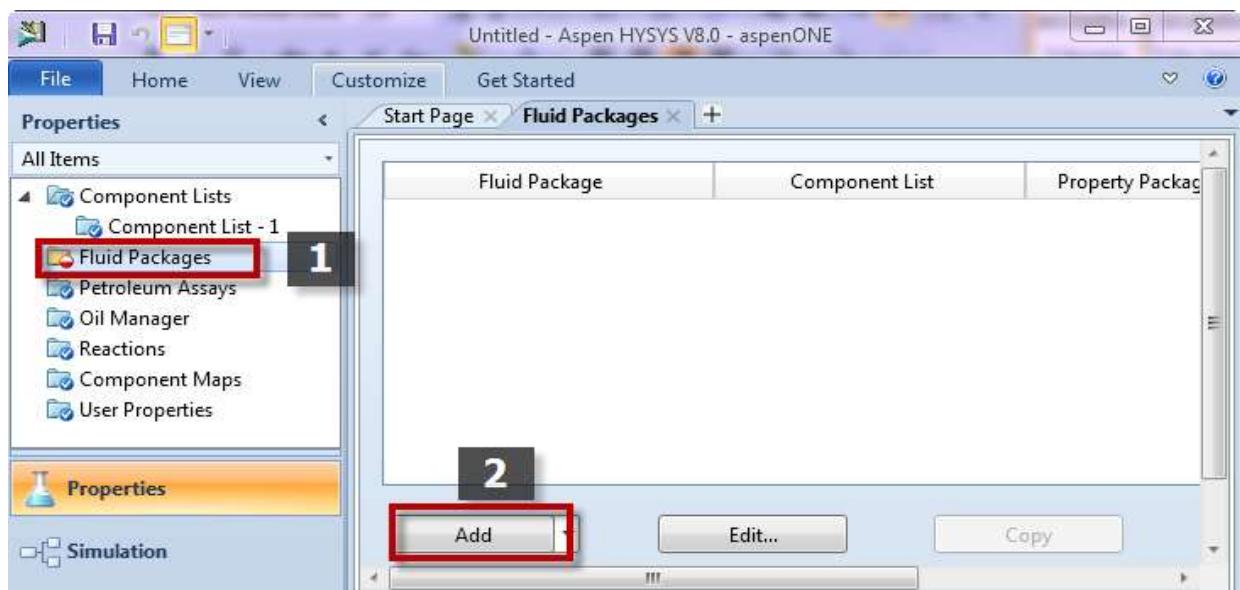
Component	Type	Group
Ethane	Pure Component	
Propane	Pure Component	
n-Butane	Pure Component	
n-Pentane	Pure Component	
n-Hexane	Pure Component	

Select: Pure Components

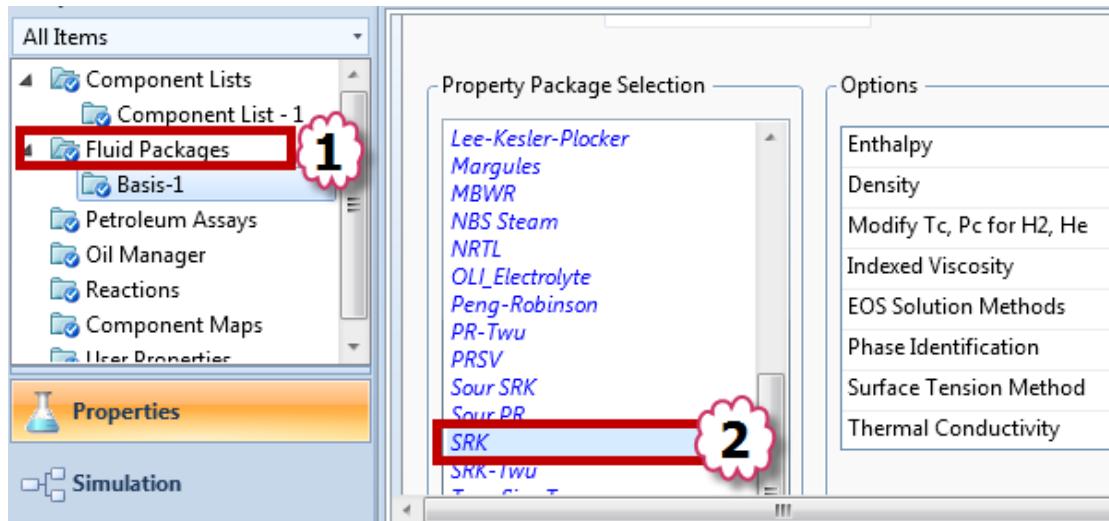
Search for:

Simulation Name
Methane
i-Butane
i-Pentane

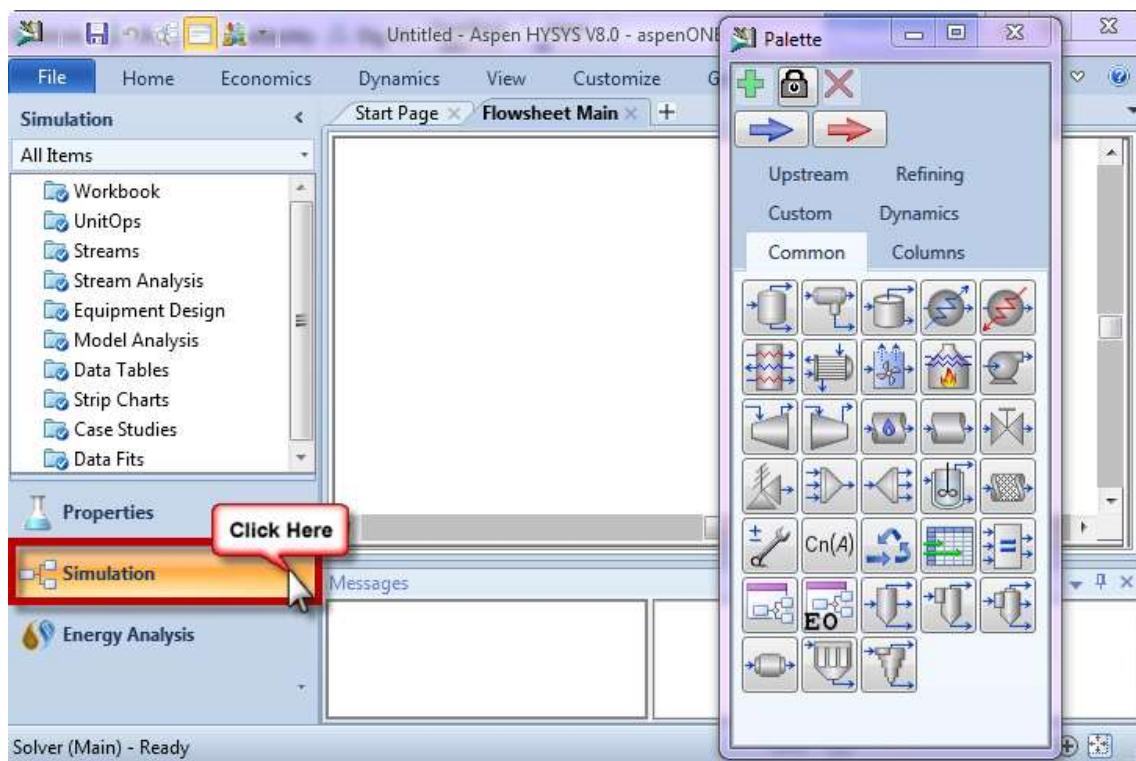
Now, select the suitable fluid package



In this case, select SRK

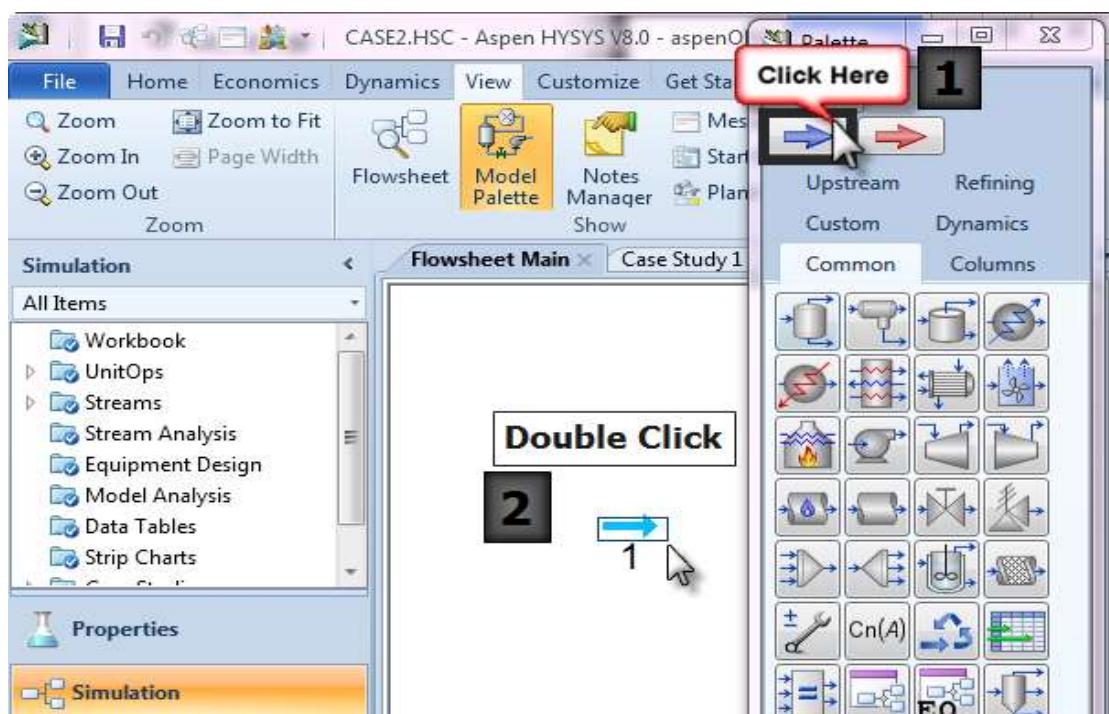


Now you can start drawing the flow sheet for the process by clicking the Simulation button:

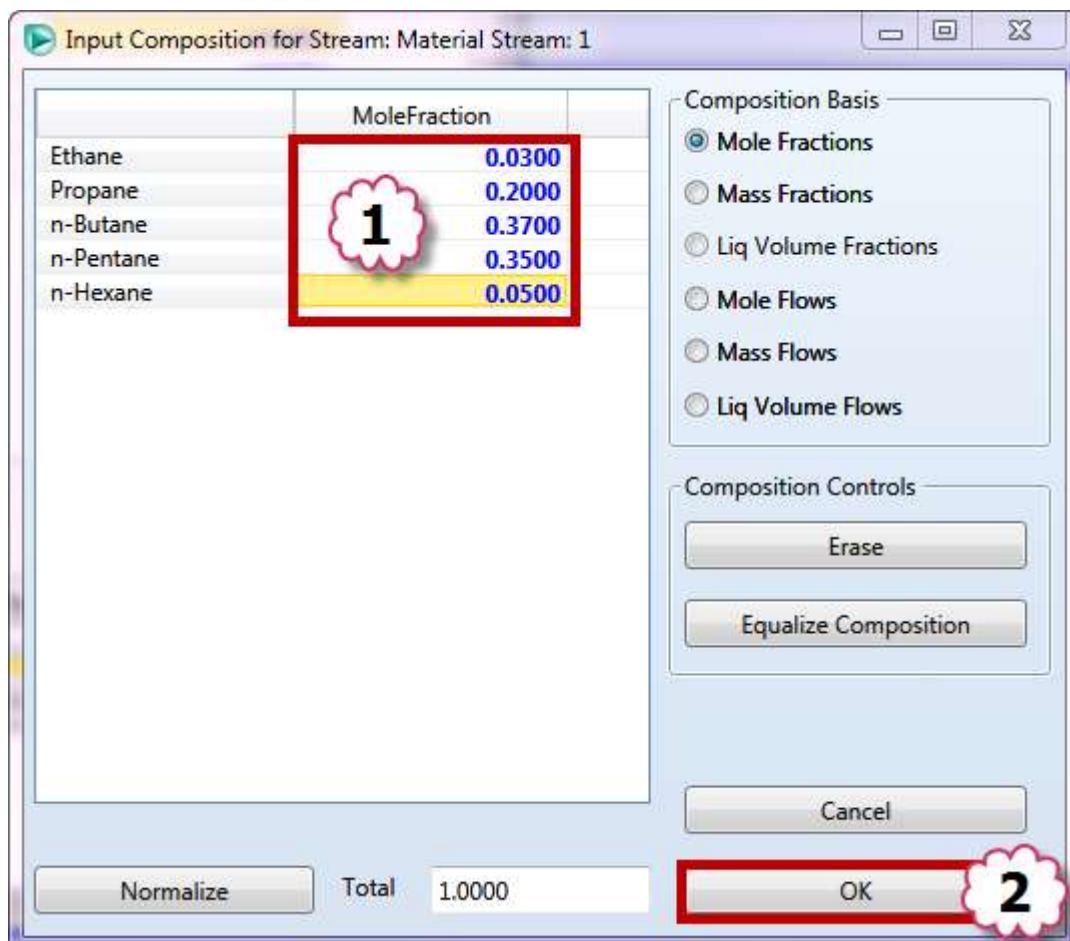


Now add a material stream to define the composition and the conditions of the feed stream

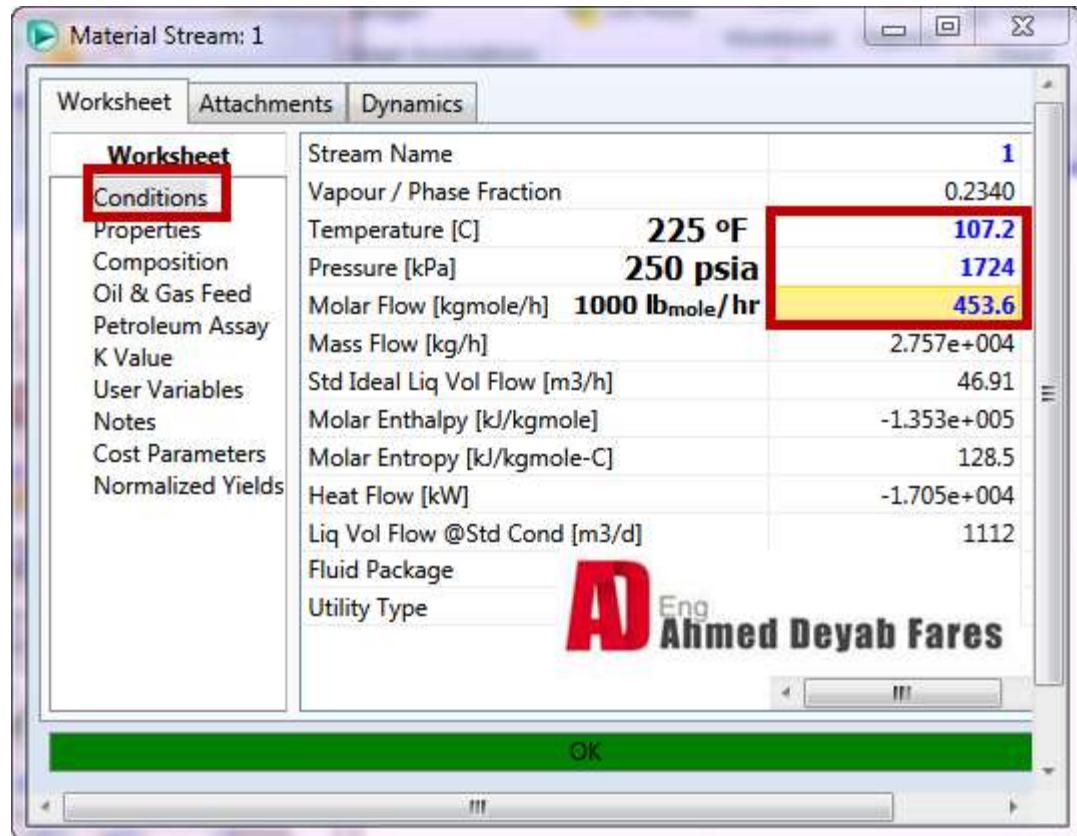
From the palette:



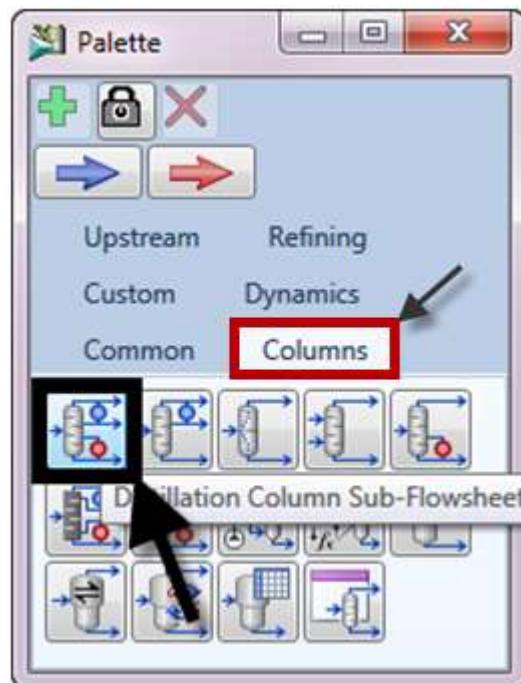
Add the mole fraction for the inlet stream



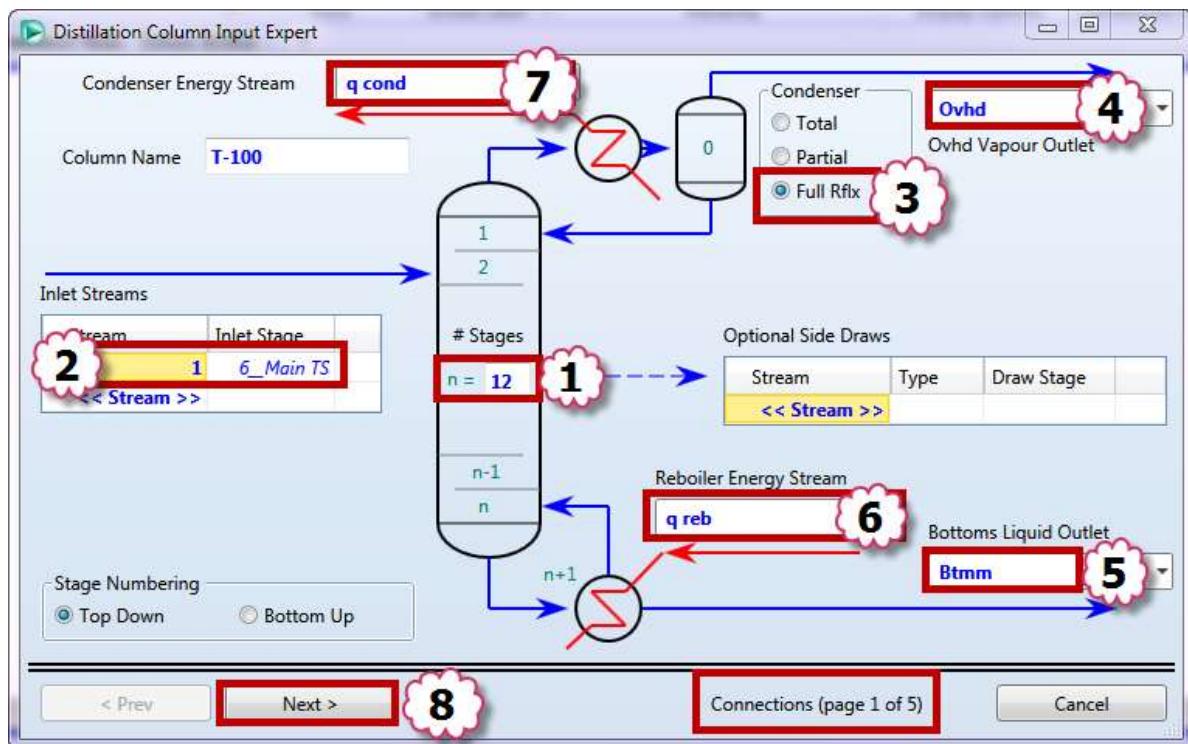
Then go to the conditions page to complete the feed stream conditions:



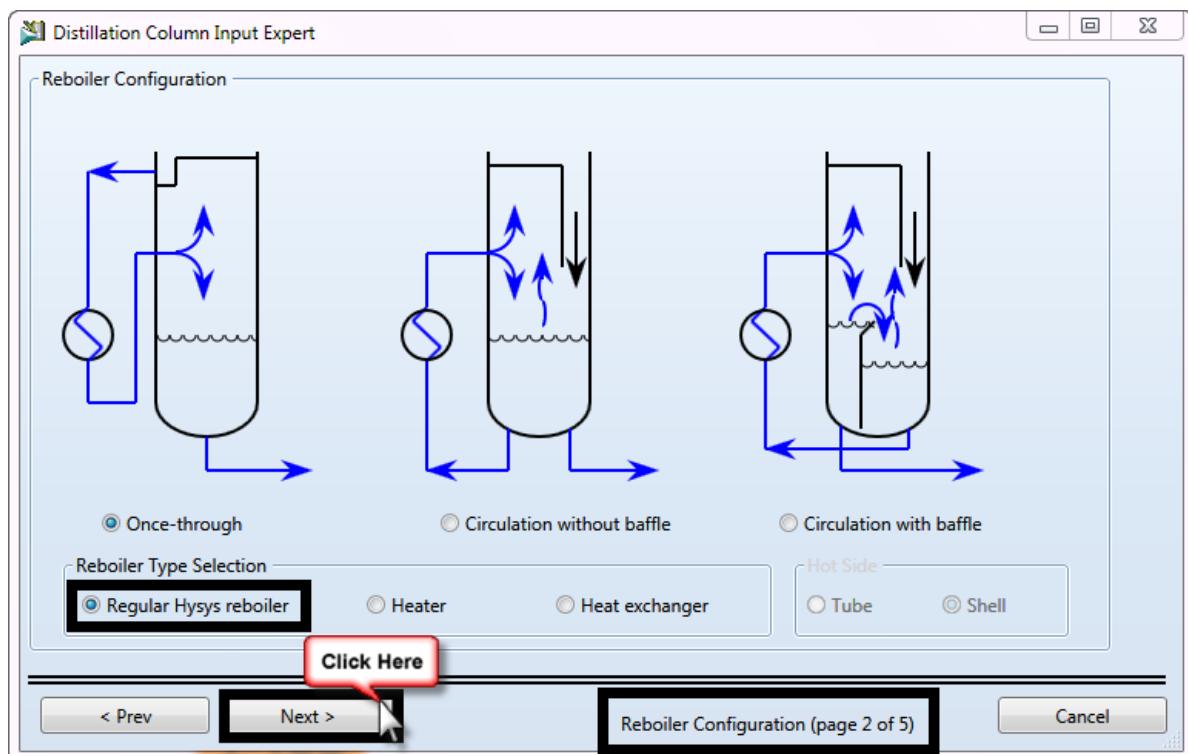
Add a distillation tower (with condenser & Reboiler):

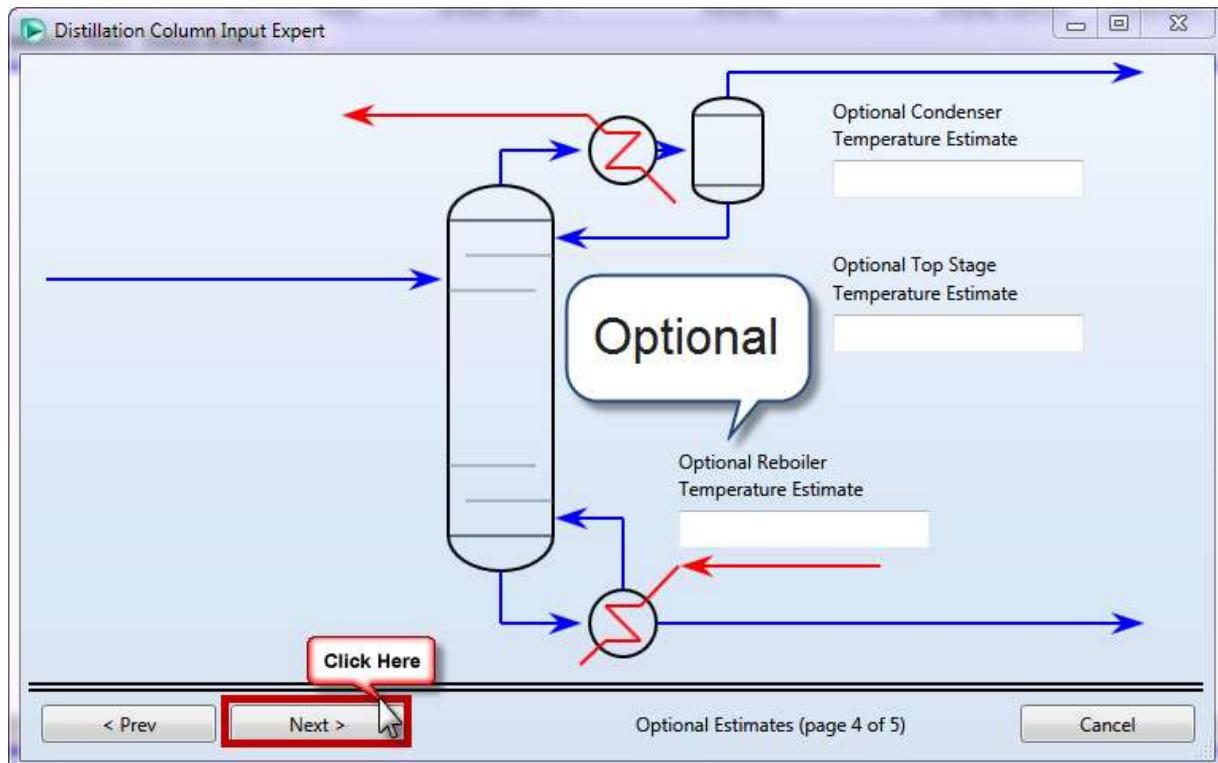
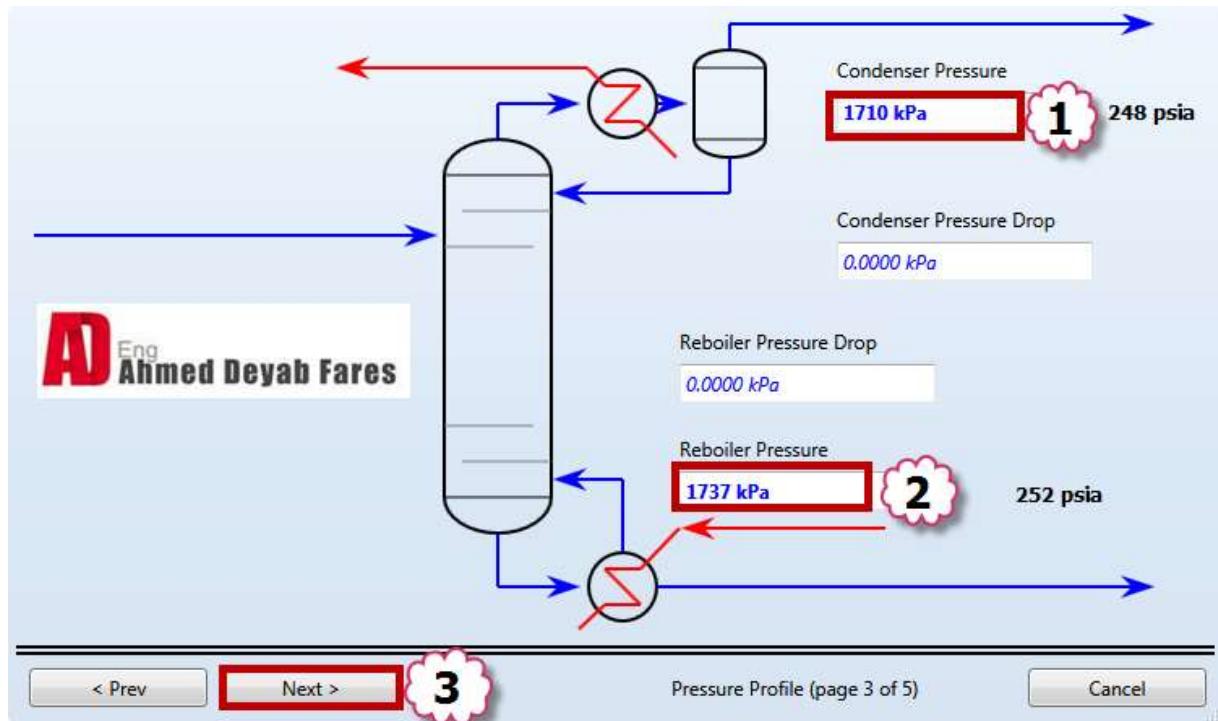


Now, start building the column:



Select Regular Hysys Reboiler in the reboiler configuration page for Kettle Reboiler as follows:

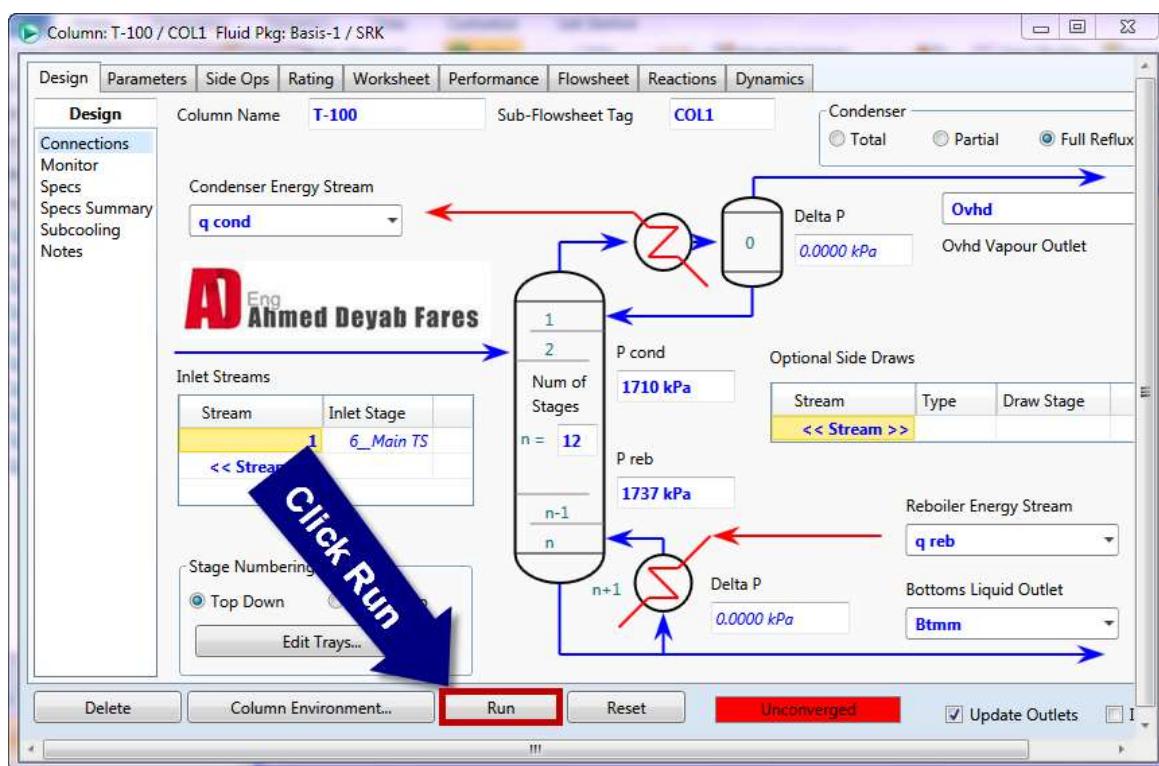
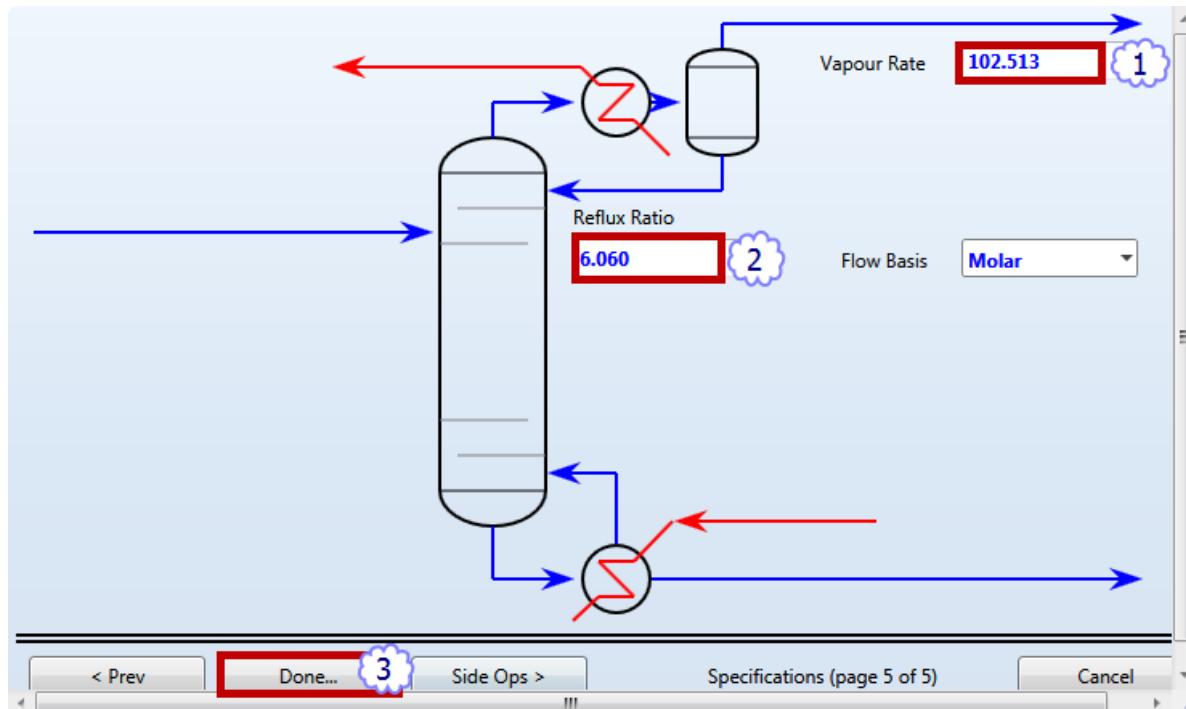




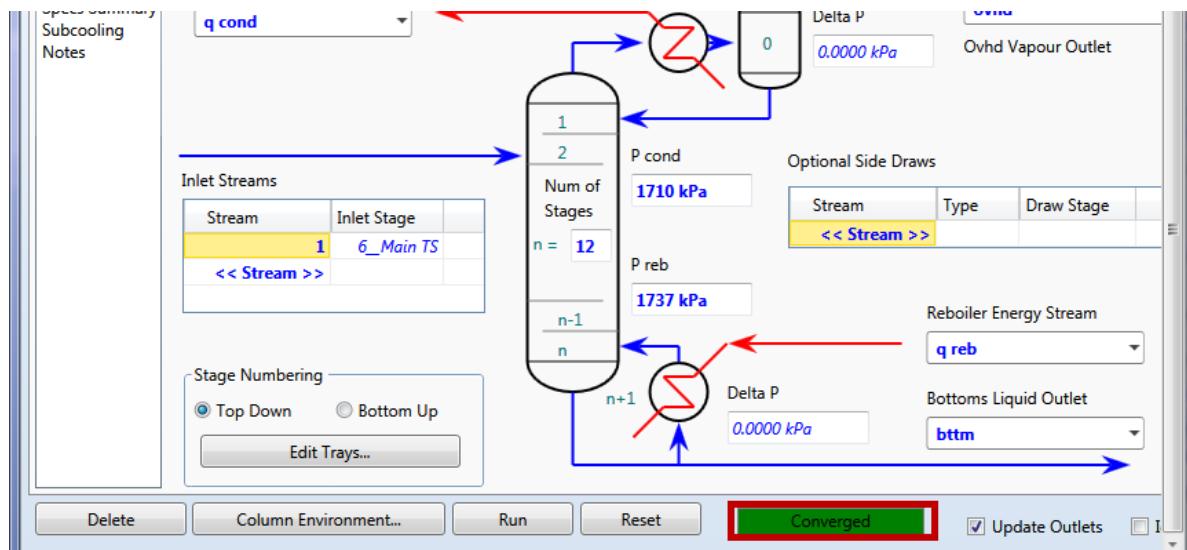
The target is to run the column with 2 specifications:

1- Reflux Ratio = 6.06

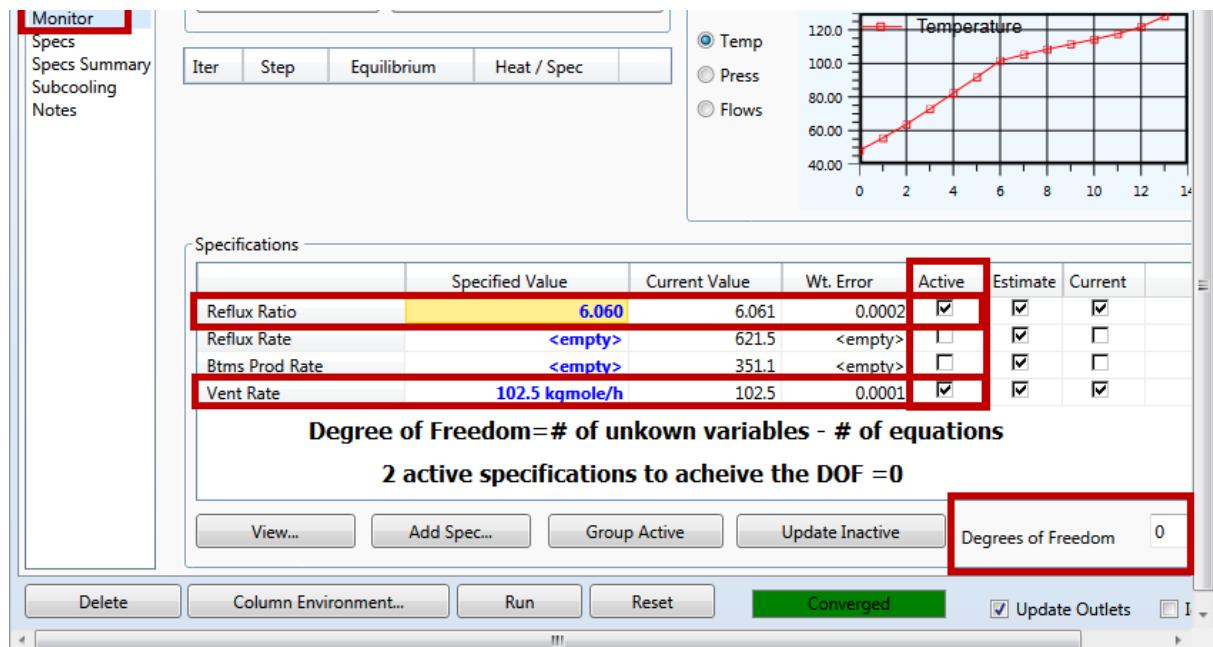
2- Overhead vapor rate (Vent rate) = 226 lb<sub>mole</sub>/hr



The column status bar (Red bar) is now un converged till clicking the RUN button to converge the column.



Let's go to the monitor page and see the current specifications:



We must make 2 specifications active to make the DOF=0.0 and to converge the column.

The two specifications are the target that you want to achieve from the column.

Although the column is converged, it is not always practical to have vapor rate & reflux ratio specifications. These specifications can result in columns which cannot be converged or that produce product streams with undesirable properties if the column feed conditions change.

An alternative approach is to specify either component fractions or component flow rates for the column product streams.

Now we have to give the column another 2 new specifications to run with (more practical)

- 1- Propane overhead flow of 191 lb<sub>mol</sub>/hr
- 2- Butane bottom flow of 365 lb<sub>mol</sub>/hr.

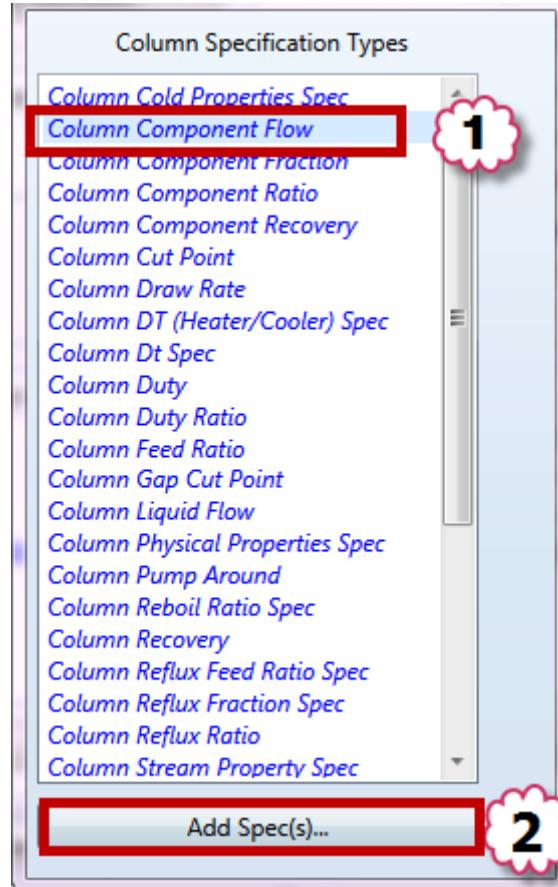
Specifications

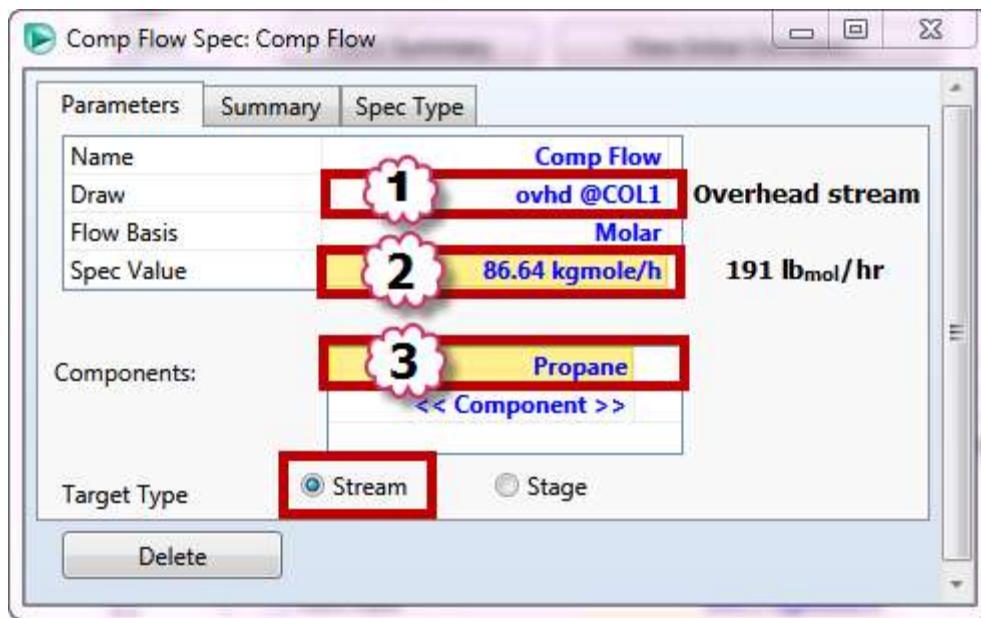
	Specified Value	Current Value	Wt. Error	Active	Estimate	Current	
Reflux Ratio	<b>6.060</b>	6.061	0.0002	<input checked="" type="checkbox"/>	<input checked="" type="checkbox"/>	<input checked="" type="checkbox"/>	
Reflux Rate	<empty>	621.7	<empty>	<input type="checkbox"/>	<input checked="" type="checkbox"/>	<input type="checkbox"/>	
Btms Prod Rate	<empty>	351.0	<empty>	<input type="checkbox"/>	<input checked="" type="checkbox"/>	<input type="checkbox"/>	
Vent Rate	<b>102.5 kgmole/h</b>	102.6	0.0006	<input checked="" type="checkbox"/>	<input checked="" type="checkbox"/>	<input checked="" type="checkbox"/>	

**Click Here**

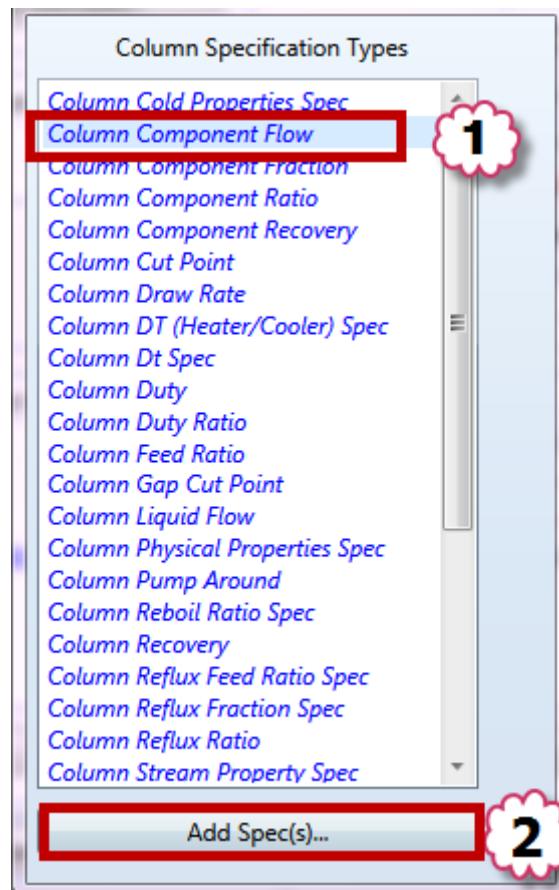
Add Spec...   Degrees of Freedom

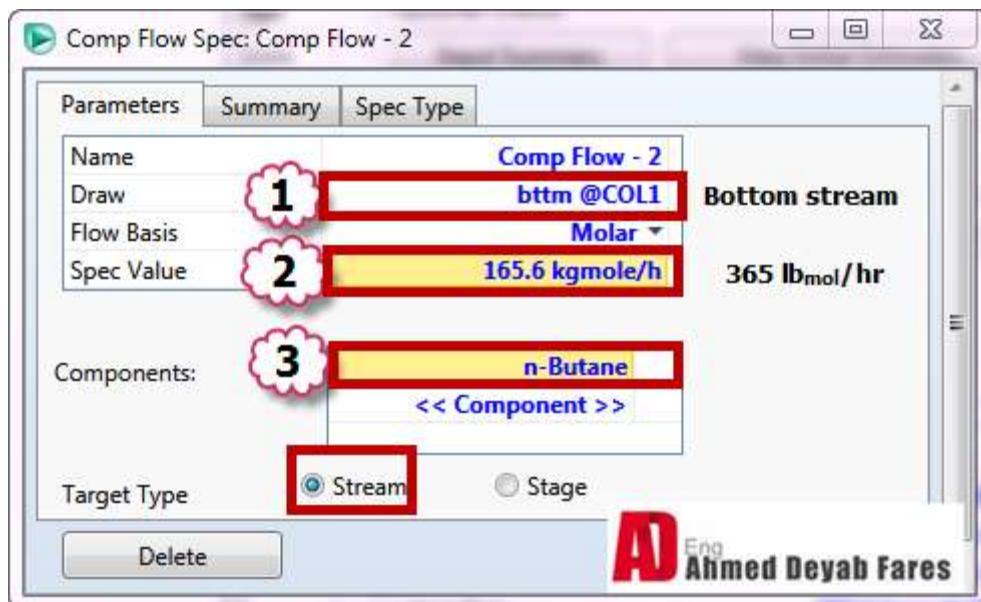
Converged  Update Outlets  I

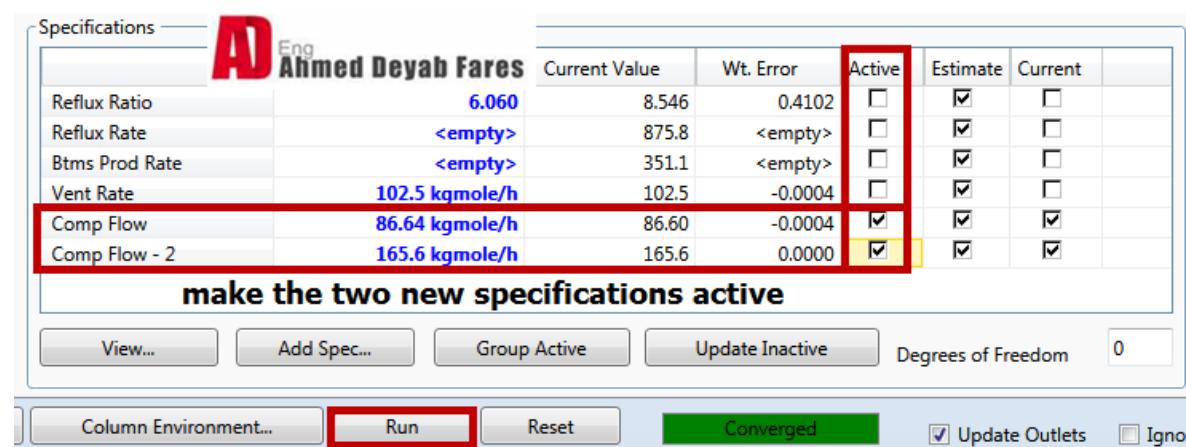


Add the other specification:





You can change the specifications by marking the Active check box on the 2 new specifications



The Results is always inside the performance page:



Go to column profile to see the temperature & Pressure profile across the tower.

	Design	Parameters	Side Ops	Rating	Worksheet	Performance	Flowsheet	Reactions	Dynamics
	Performance				Reflux Ratio	8.546			
	Summary				Boilup Ratio	2.121	Flows	Energy	Basis
	Column Profiles								
	Feeds / Products								
	Plots								
	Cond./Reboiler								
					Temperature [C]	Pressure [kPa]	Net Liquid [kgmole/h]	Net Vapour [kgmole/h]	
					Condenser	46.06	1710	875.775	
					1_Main TS	51.63	1710	854.449	978.252
					2_Main TS	57.72	1712	816.478	956.925
					3_Main TS	65.82	1715	772.861	918.954
					4_Main TS	75.93	1717	733.714	875.337
					5_Main TS	87.10	1720	698.895	836.190
					6_Main TS	98.81	1722	1044.86	801.372
					7_Main TS	103.5	1725	1064.36	693.744
					8_Main TS	107.6	1727	1081.04	713.240
					9_Main TS	111.1	1730	1094.06	729.924
					10_Main TS	114.4	1732	1102.07	742.942
					11_Main TS	117.8	1735	1103.22	750.946
					12_Main TS	122.3	1737	1095.84	752.095
					Reboiler	128.8	1737		744.719

Condenser Temp	46 °C
Reboiler Temp	128.8 °C
Reflux Ratio	8.54

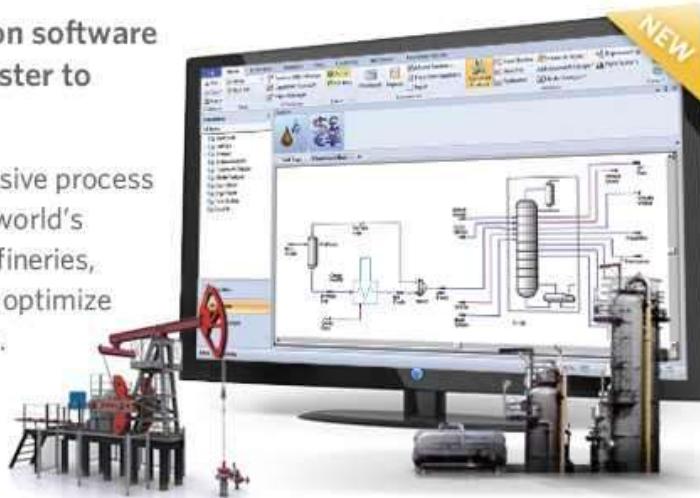
# Oil Characterization

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5

## Workshop

The petroleum characterization method in HYSYS will convert laboratory analyses of condensates, crude oils, petroleum cuts and coaltar liquids into a series of discrete hypothetical components. These petroleum hypo components provide the basis for the property package to predict the remaining thermodynamic and transport properties necessary for fluid modeling.

HYSYS will produce a complete set of physical and critical properties for the petroleum hypo components with a minimal amount of information. However, the more information you can supply about the fluid, the more accurate these properties will be, and the better HYSYS will predict the fluid's actual behavior.

In this example, the Oil Characterization option in HYSYS is used to model a crude oil. The crude is the feed stock to a Pre-heat Train, followed by the Atmospheric Crude Column, which will be modelled in a subsequent module.

## Learning Objectives

Once you have completed, you will be able to use the Oil Characterization option in HYSYS.

### Oil Characterization

The petroleum characterization method in HYSYS will convert laboratory analyses of condensates, crude oils, petroleum cuts and coal-tar liquids into a series of discrete hypothetical components. These petroleum hypocomponents provide the basis for the property package to predict the remaining thermodynamic and transport properties necessary for fluid modeling.

In this example, the Oil Characterization option in HYSYS is used to model a crude oil. The crude is the feed stock to oil refining process. (FPkg=PR)

#### **Bulk Properties of the crude:**

API Gravity of 29 for the crude

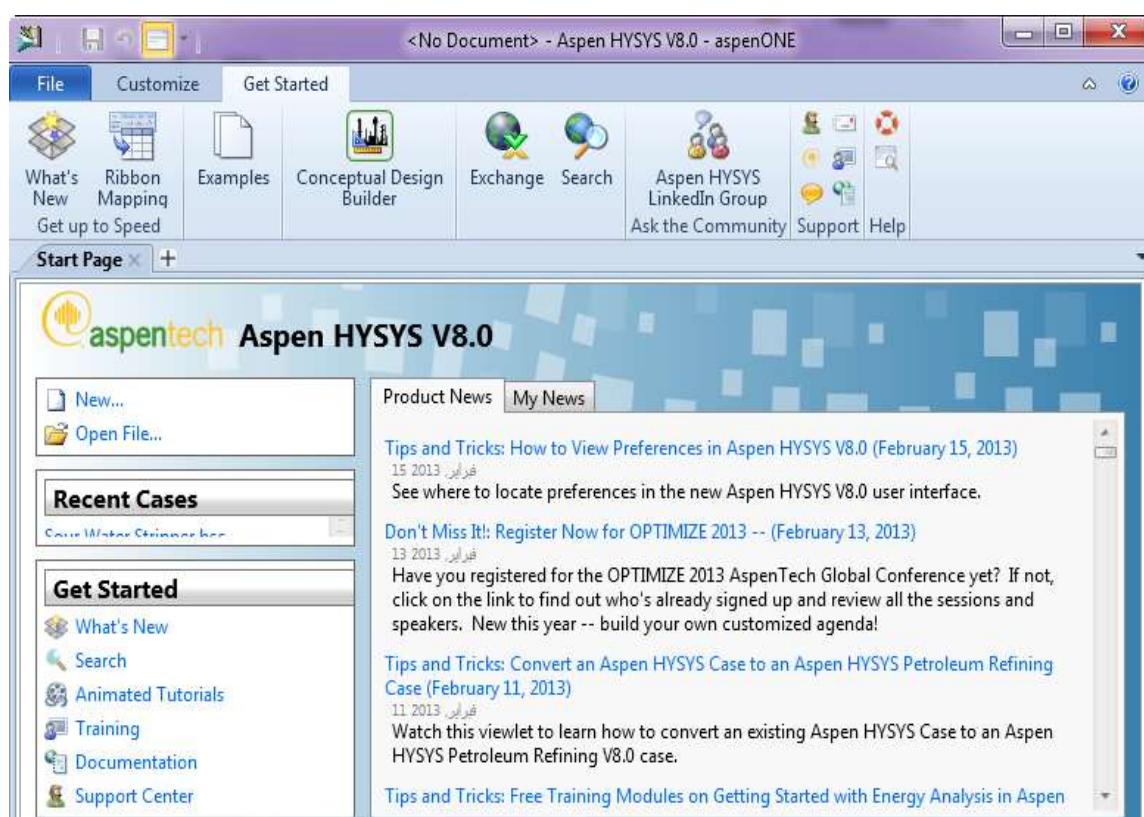
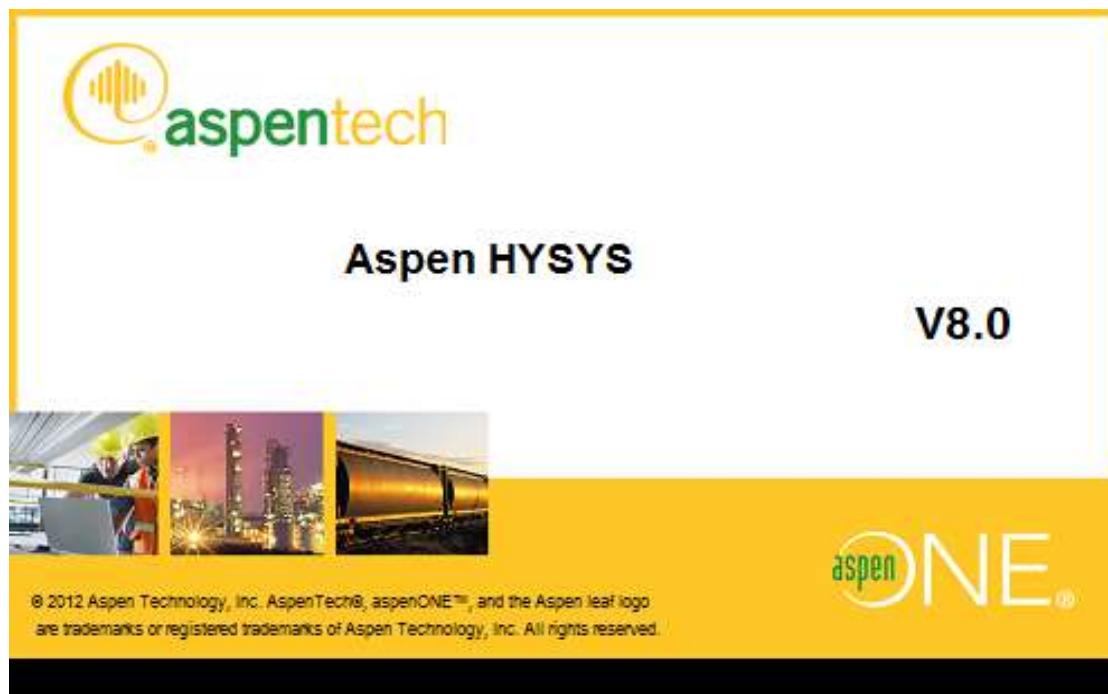
#### **Light Ends (Liquid Volume):**

<i>Light Ends</i>	<i>Liquid Vol %</i>
Methane	0.0065
Ethane	0.0225
Propane	0.3200
i-butane	0.2400
n-butane	1.7500
i-pentane	1.6500
n-pentane	2.2500

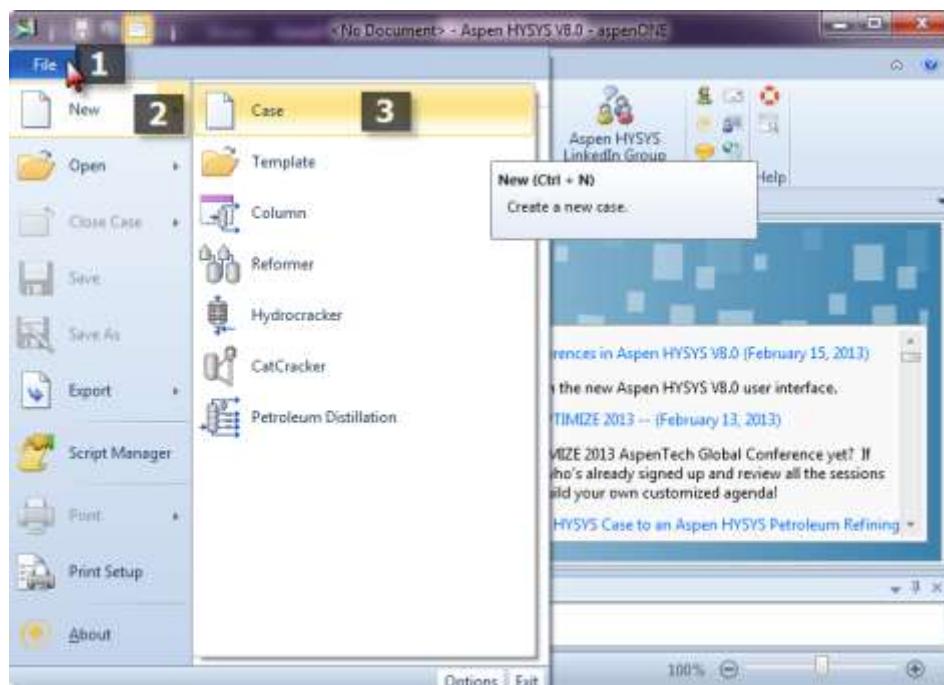
#### **TBP distillation data (Liquid Volume %):**

<b>Assay Percent</b>	<b>Temperature (°C)</b>
0	-12
4	32
9	74
14	116
20	154
30	224
40	273
50	327
60	393
70	450
76	490
80	516

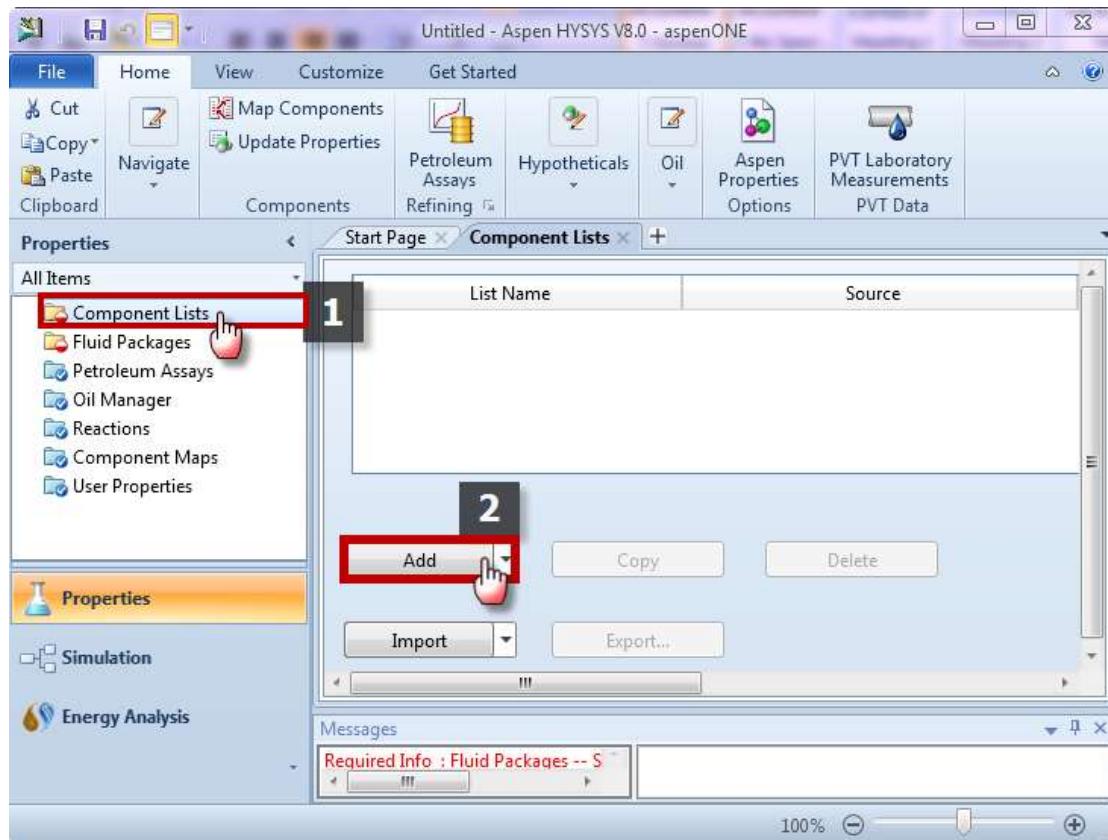
To start the program, From Start Menu, Select All Programs >>  
Aspen Tech >> Process Modeling V8.0 >>> Aspen HYSYS >>  
Aspen HYSYS



## 1- First, Start a new case



## 2- Add the Components

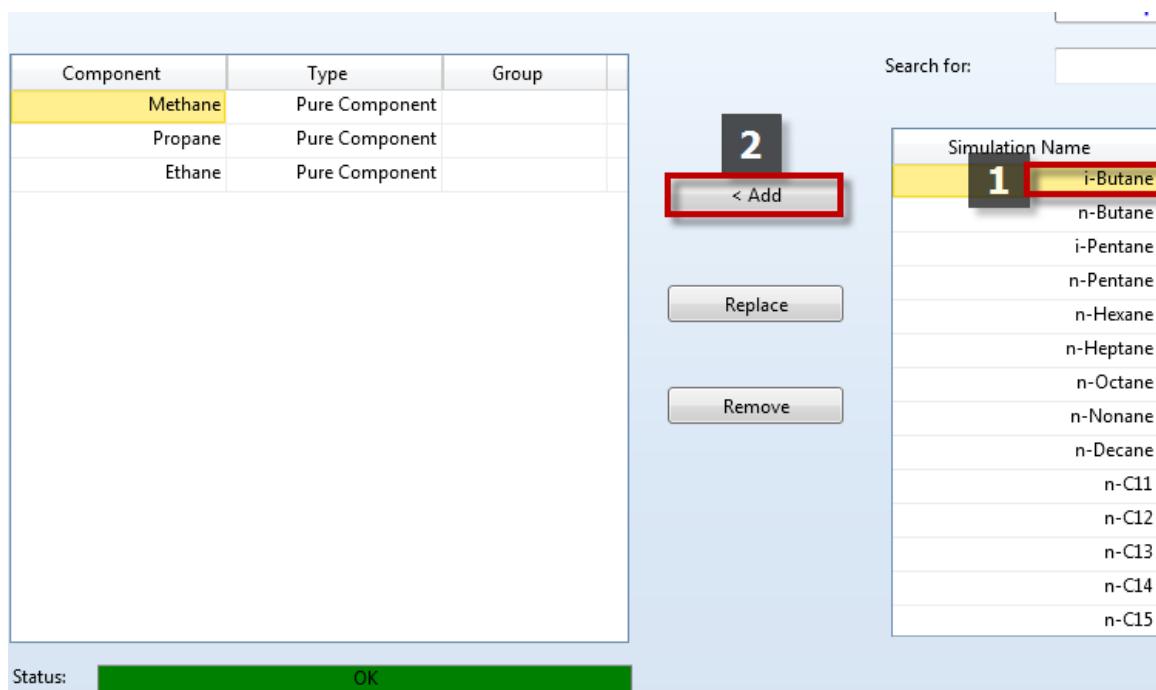
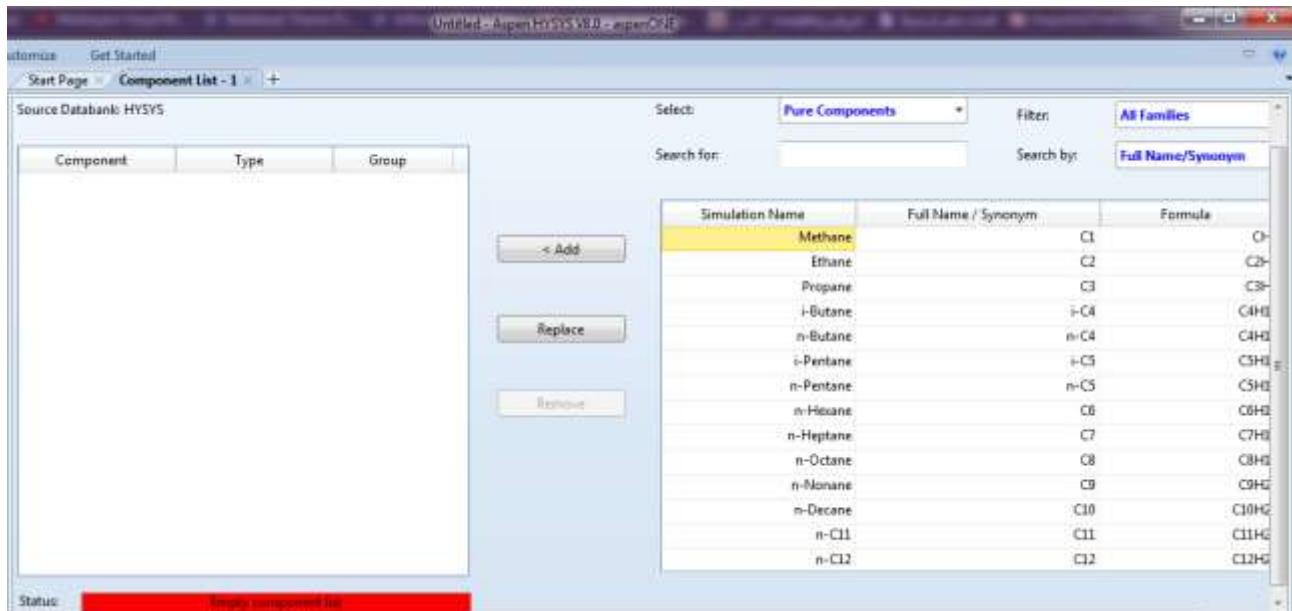


3- Choose the system components from the databank:

4- First we should add the pure components first (Light Ends)

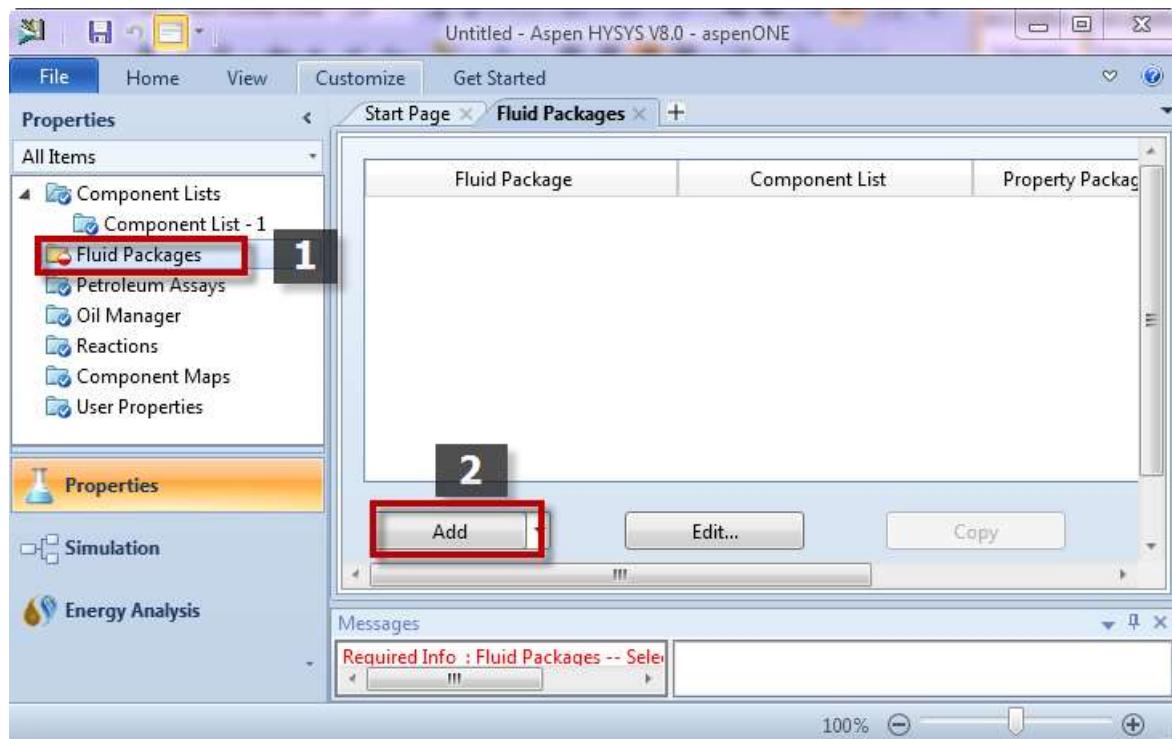
Light Ends are defined as pure components with low boiling points.

Components in the boiling range of C1 to n-C5 are most commonly of interest.

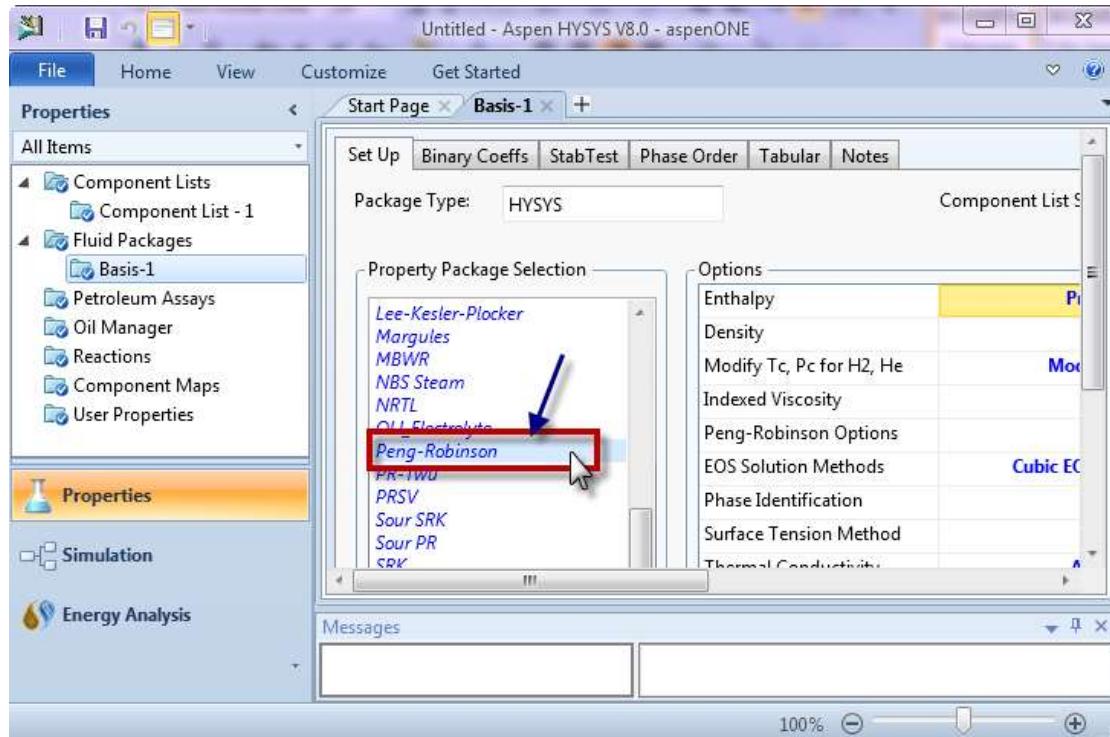


Add the pure components (C1, C2, C3, n-C4, i-C4, n-C5, i-C5, H<sub>2</sub>O)

Now, select the suitable fluid package



In this case, select Peng-Robinson

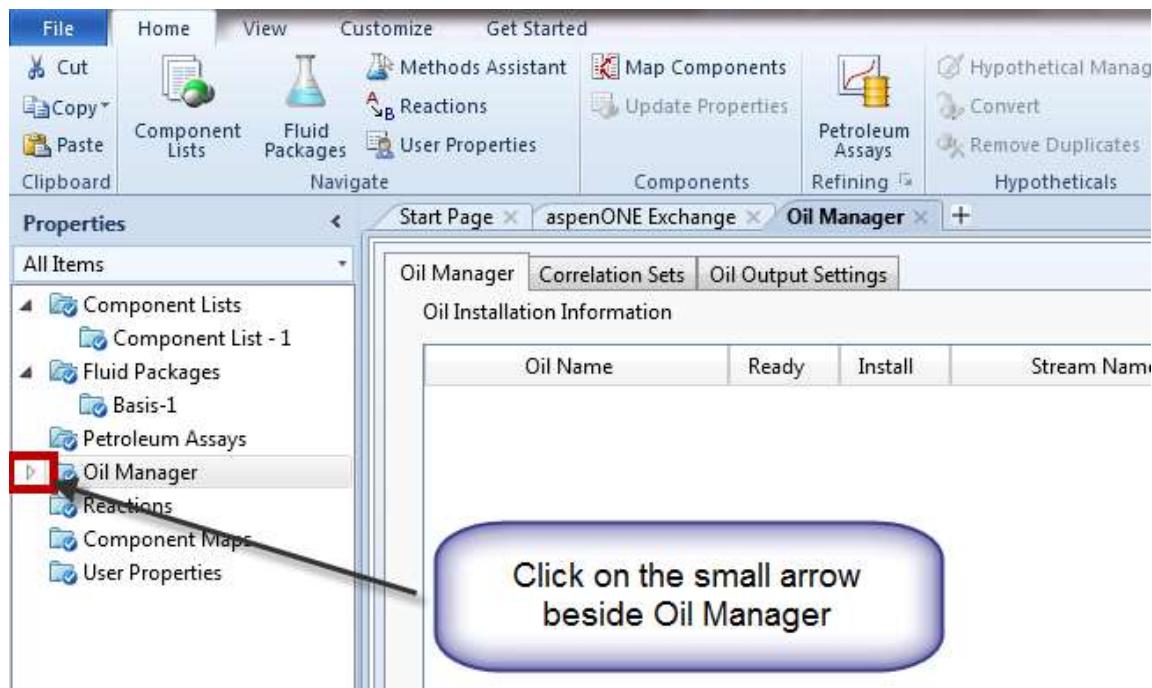


### Characterize the Assay

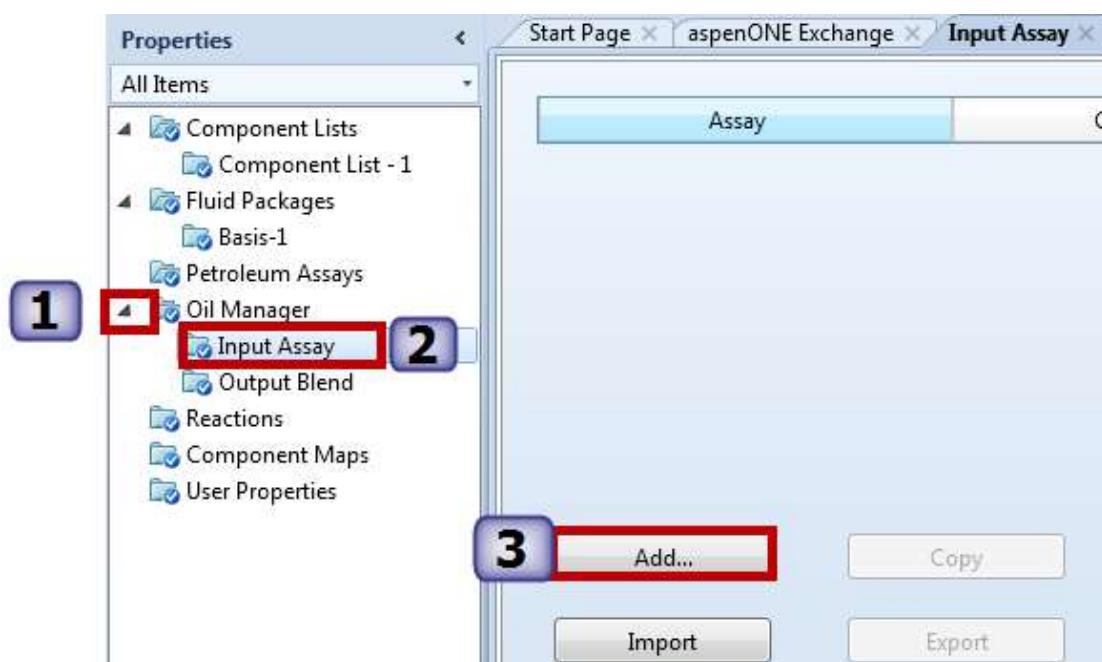
The assay contains all of the petroleum laboratory data, boiling point curves, light ends, property curves and bulk properties. HYSYS uses the supplied Assay data to generate internal TBP, molecular weight, density and viscosity curves, referred to as Working Curves.

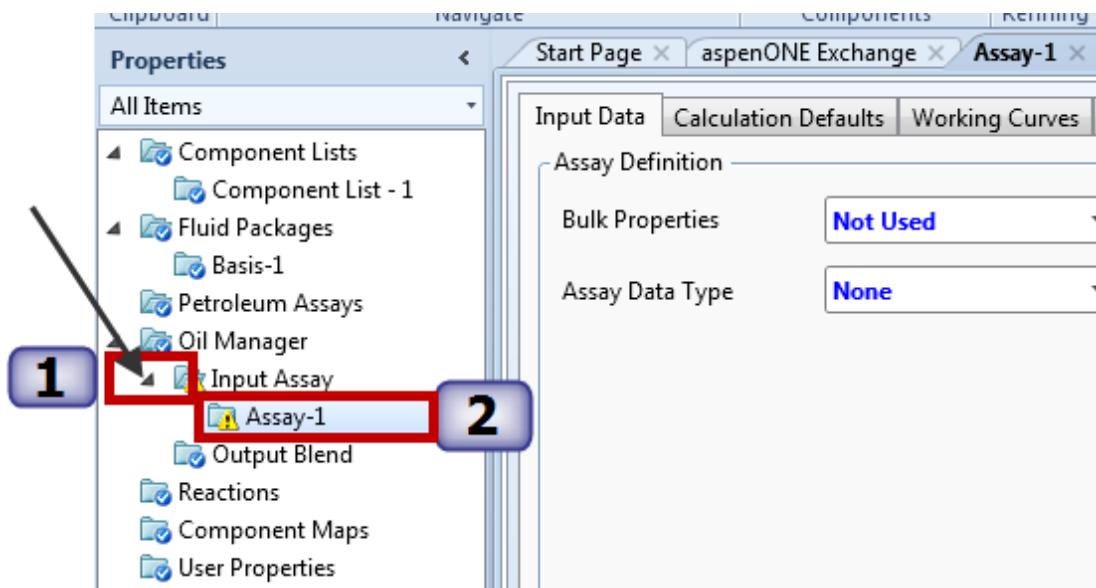
To characterize the assay follow the following steps:

- 1- Go to Oil Manager.



- 2- Click on Input Assay & then Click Add button:



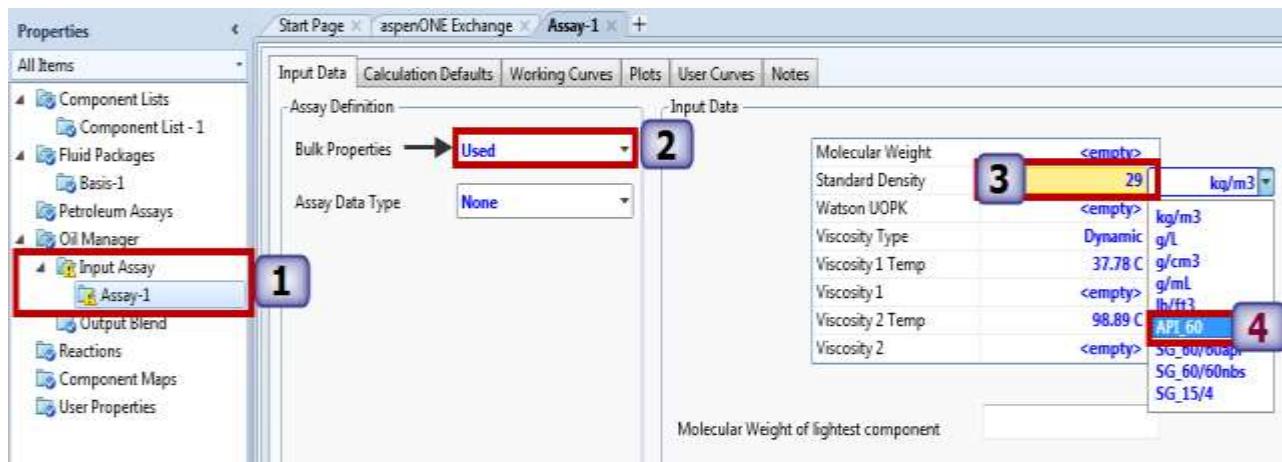


### Bulk Properties

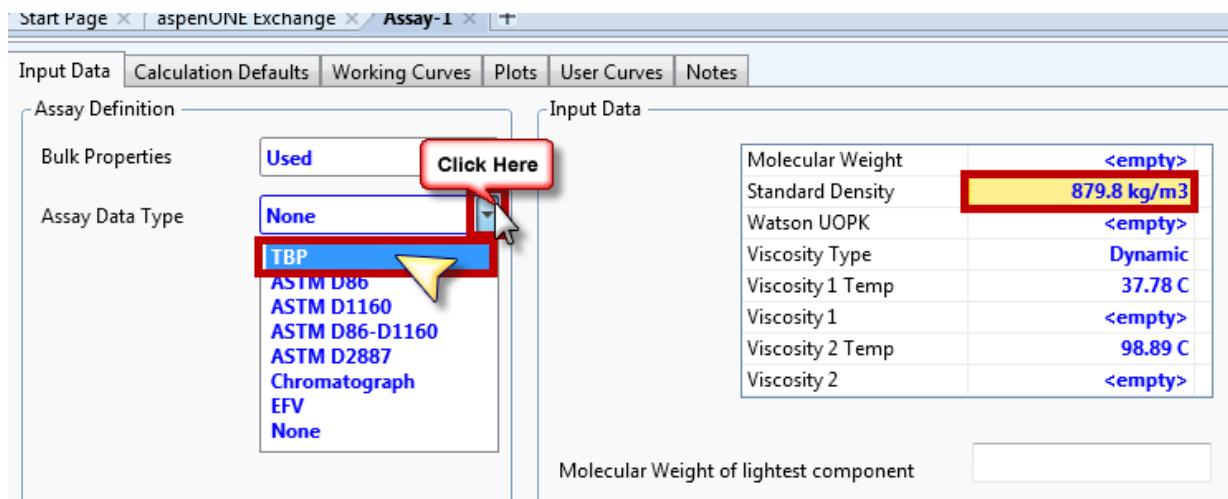
Bulk Properties for the sample may also be supplied. The bulk properties are optional if a distillation curve or chromatograph have been supplied  
Change the bulk properties from **Not used** to **Used** and add the value for standard Density = 29 API\_60

Note:

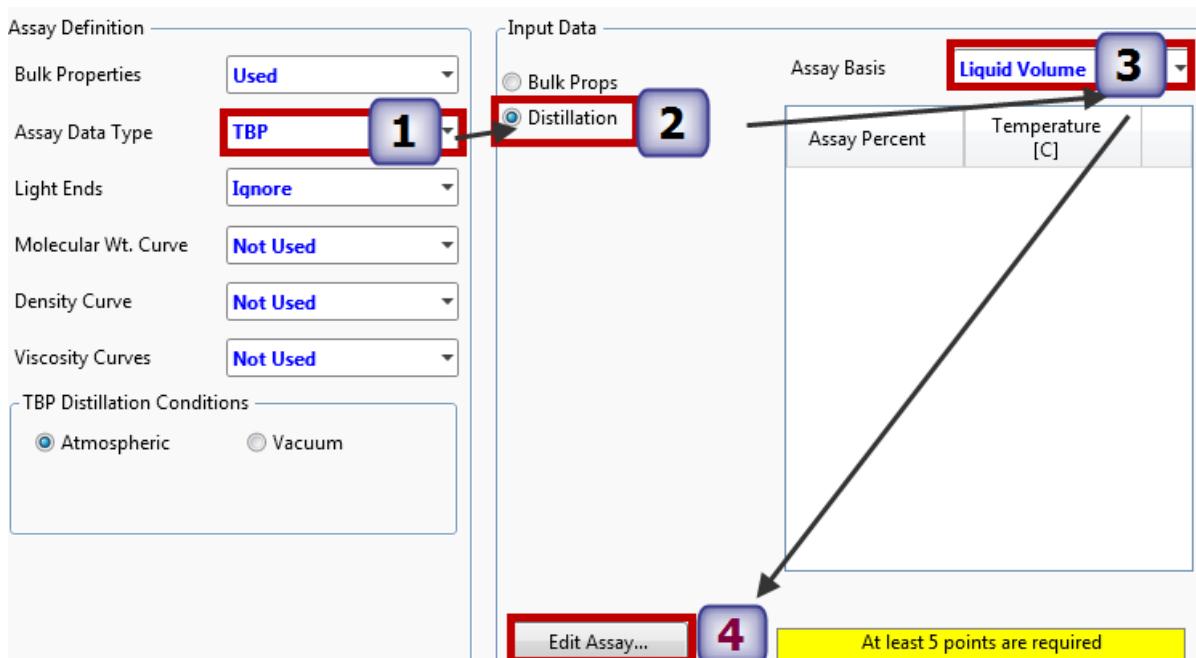
$$\text{API} = \frac{141.5}{SG} - 131.5$$



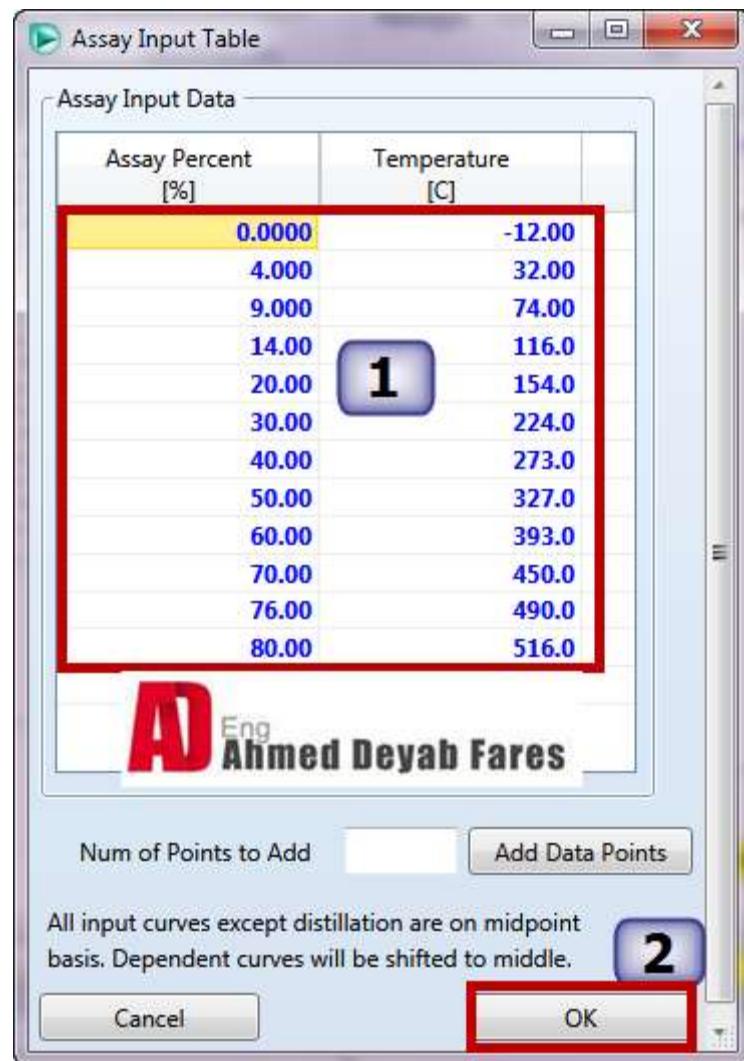
Select the Assay Data Type >> TBP



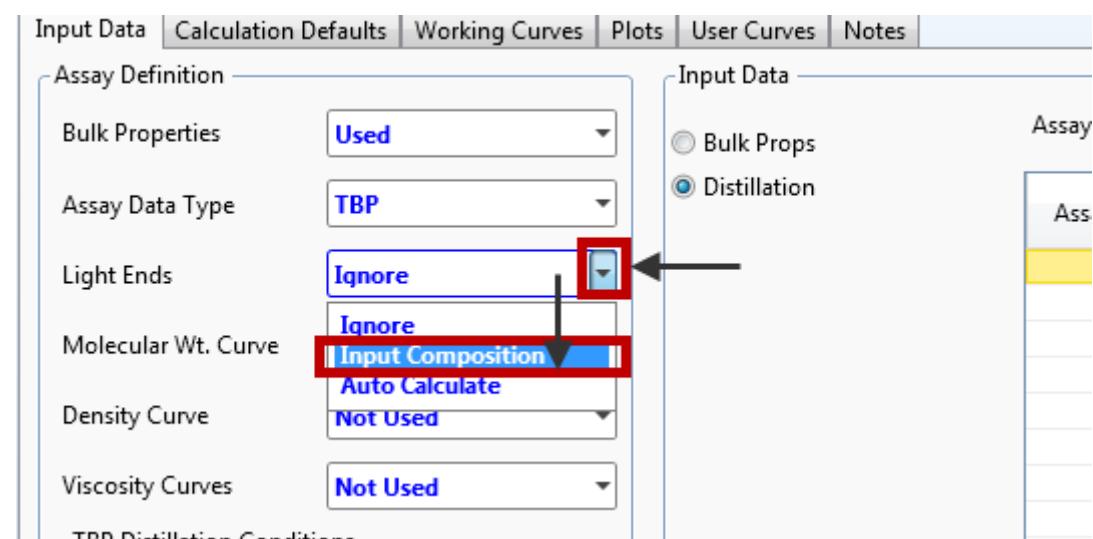
- Select the Distillation radio button in the Input Data group box.
- Select the Assay Basis as Liquid Volume (use the drop-down menu).
- Click the Edit Assay button; this will allow you to enter the assay information below.



Add the assay data:



Use the drop-down lists to select **Input Composition** for Light Ends



Select the **Light Ends** radio button and enter the data given from **Input Data**

put Data | Calculation Defaults | Working Curves | Plots | User Curves | Notes

**Assay Definition**

Bulk Properties	Used
Assay Data Type	TBP
Light Ends	Input Composition
Molecular Wt. Curve	Not Used
Density Curve	Not Used
Viscosity Curves	Not Used
TBP Distillation Conditions	

**Input Data**

<input type="radio"/> Bulk Props	<input checked="" type="radio"/> Light Ends	<input type="radio"/> Distillation
		Light Ends Basis
		Liquid Volume %
Light Ends	Composition	NBP [C]
Methane	6.500e-003	-161.5
Ethane	2.250e-002	-88.60
Propane	0.3200	-42.10
i-Butane	0.2400	-11.73
n-Butane	1.750	-0.5020
i-Pentane	1.650	27.88
n-Pentane	2.250	36.06
H <sub>2</sub> O	0.0000	100.0

Once you have entered all of the data, click the Calculate button. The status message at the bottom of the Assay view will display Assay Was Calculated.

**Assay Definition**

TBP
Input Composition
Not Used
Not Used
Not Used
conditions
: Vacuum

**Input Data**

<input checked="" type="radio"/> Light Ends	<input type="radio"/> Distillation	
Light Ends	Composition	NBP [C]
Methane	6.500e-003	-161.5
Ethane	2.250e-002	-88.60
Propane	0.3200	-42.10
i-Butane	0.2400	-11.73
n-Butane	1.750	-0.5020
i-Pentane	1.650	27.88
n-Pentane	2.250	36.06
H <sub>2</sub> O	0.0000	100.0

**Click Calculate**

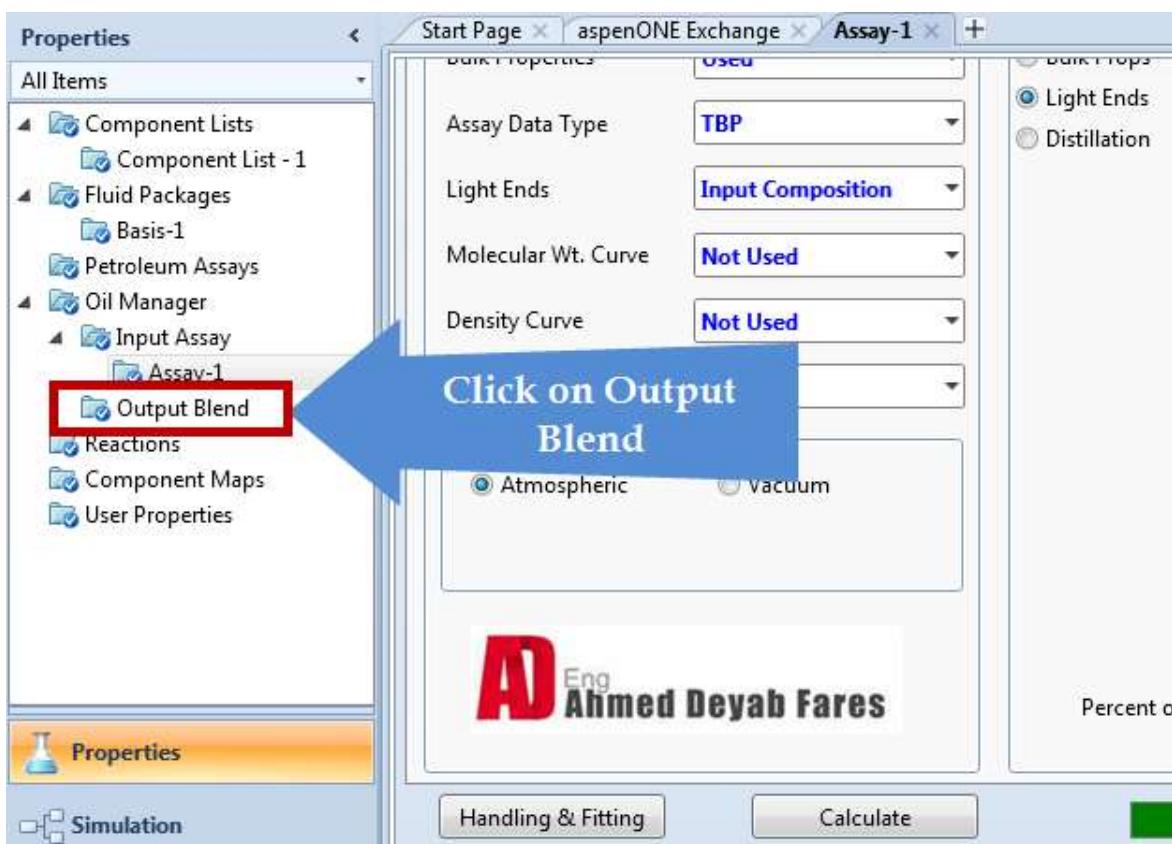
Light Ends in Assay: 6.2390

**Calculate**

Assay Was Calculated

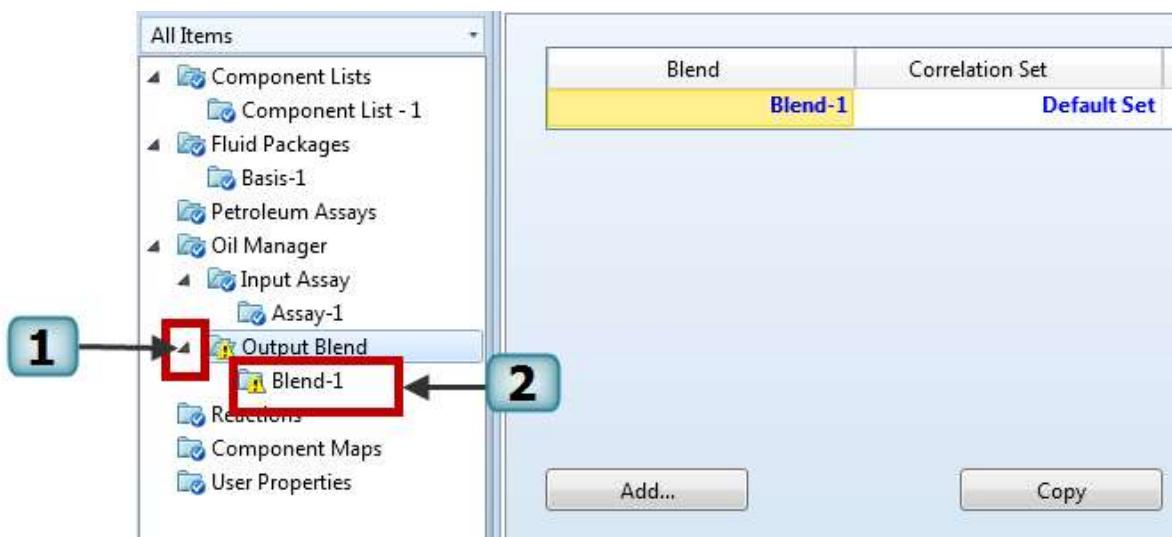
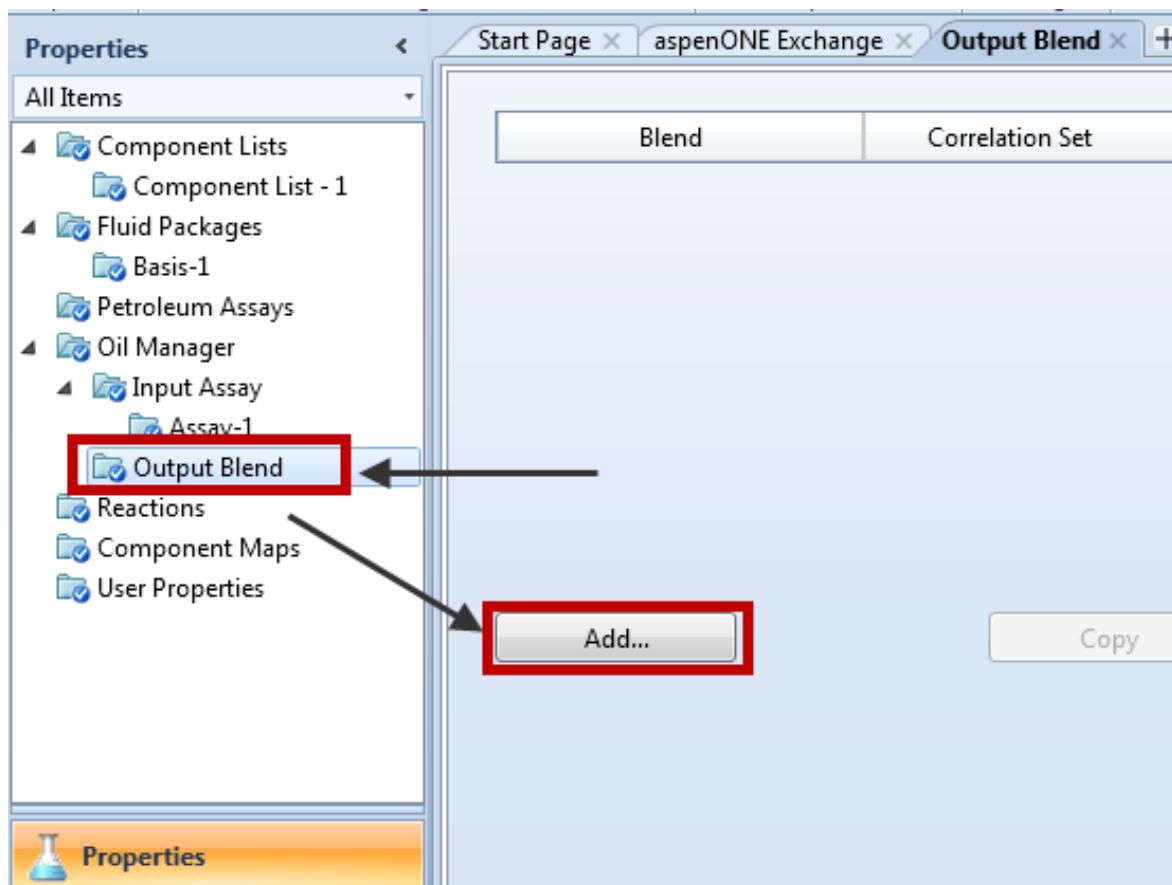
Once the Assay is calculated, the working curves are displayed on the **Plots** and **Working Curves** tabs. The working curves are regressed and extrapolated from the Assay input. From the user-supplied data, HYSYS generates curves for NBP, molecular weight, mass density, and viscosity. These working curves are used in determining the properties of the hypo components generated in the Blend step.

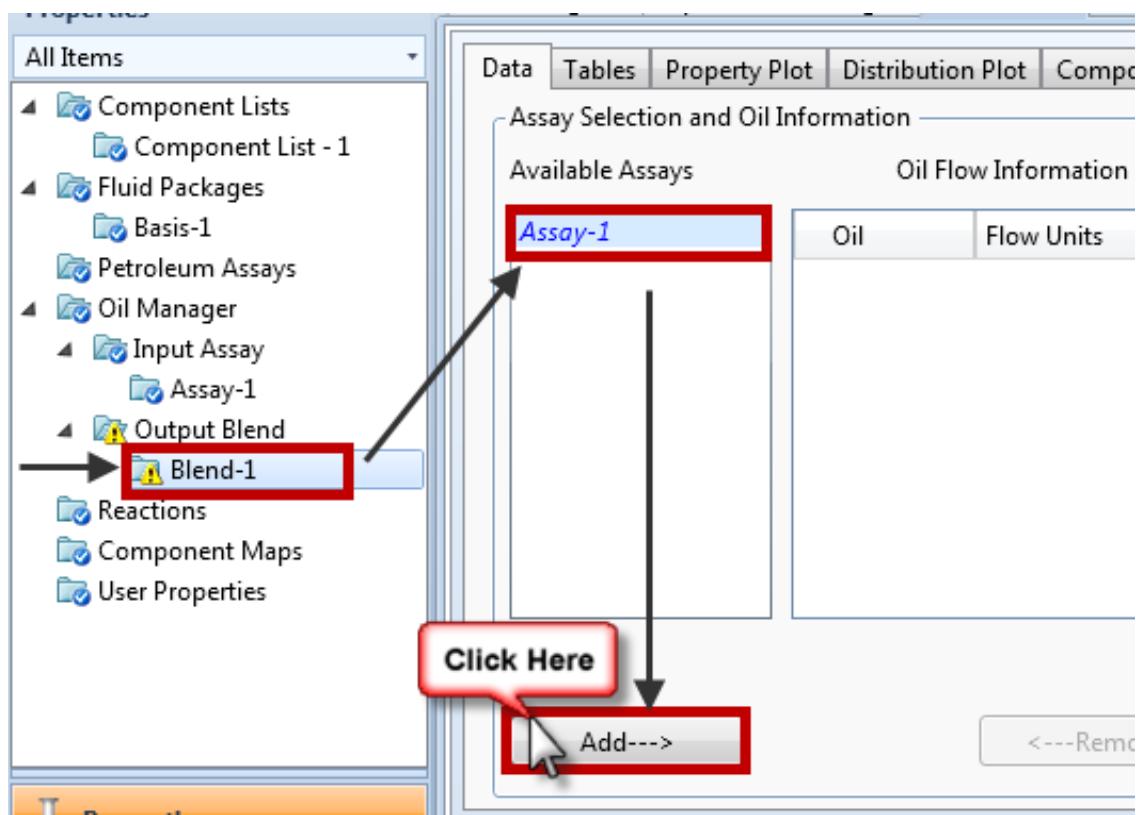
Assay Working Curves							
Point #	Moles	Cum. Moles	NBP [C]	Mole Wt	Mass Density [kg/m³]	Viscosity 1 [cP]	Viscosity 2 [cP]
0	0.00000	0.00000	47.57	71.49	724.7	0.407	0.210
1	0.02524	0.02524	55.42	74.83	730.6	0.404	0.214
2	0.02436	0.04960	63.01	78.13	736.2	0.311	0.188
3	0.02355	0.07315	70.44	81.43	741.6	0.326	0.197
4	0.02274	0.09589	78.20	84.93	747.1	0.345	0.207
5	0.02214	0.11803	84.34	87.77	751.5	0.362	0.216
6	0.02161	0.13964	89.76	90.35	755.2	0.377	0.224
7	0.02103	0.16067	95.80	93.37	759.4	0.396	0.234
8	0.02032	0.18099	103.4	97.27	764.5	0.421	0.246
9	0.01965	0.20064	111.1	101.3	769.8	0.450	0.261
10	0.01907	0.21971	118.3	105.0	774.5	0.479	0.275
11	0.04476	0.26446	134.1	113.3	784.8	0.550	0.309
12	0.04236	0.30682	148.3	121.1	793.8	0.623	0.343
13	0.03994	0.34676	163.7	130.0	803.3	0.714	0.385
14	0.03748	0.38424	180.5	140.3	813.5	0.842	0.441
15	0.03521	0.41945	197.2	151.2	823.4	1.024	0.505
16	0.03313	0.45258	213.7	162.5	832.9	1.252	0.580



The screenshot shows the HYSYS software interface. On the left, the 'Properties' pane is open, displaying a tree structure of items. A blue arrow points from the 'Output Blend' folder under 'Assay-1' to the 'Output Blend' section in the main properties pane. The main pane shows various characterization parameters: Assay Data Type set to 'TBP', Light Ends set to 'Input Composition', Molecular Wt. Curve and Density Curve both set to 'Not Used'. At the bottom, there are buttons for 'Handling & Fitting' and 'Calculate'.

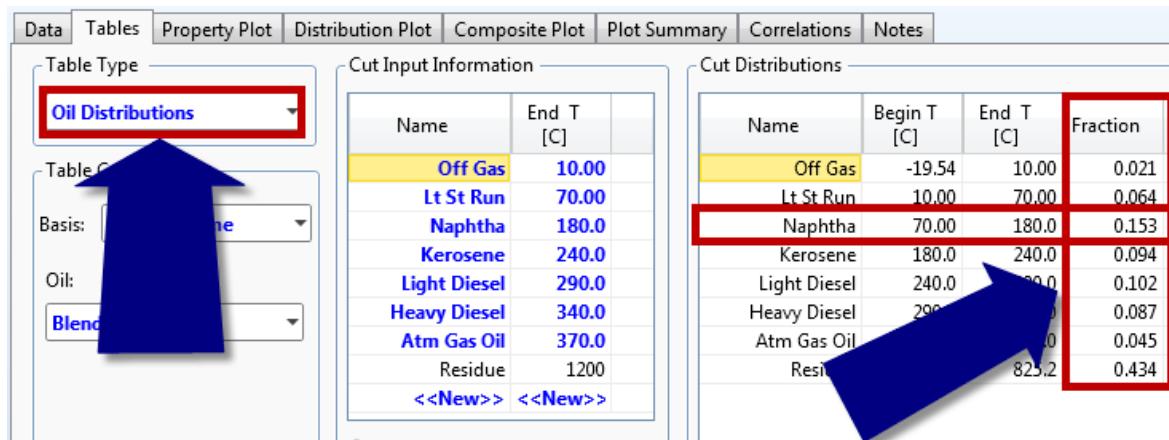
The Output Blend characterization in HYSYS splits the internal working curves for one or more assays into hypo components. The Blend tab of the Oil Characterization view provides two functions, cutting the Oil into Hypo components and Blending two or more Assays into one set of hypo components.



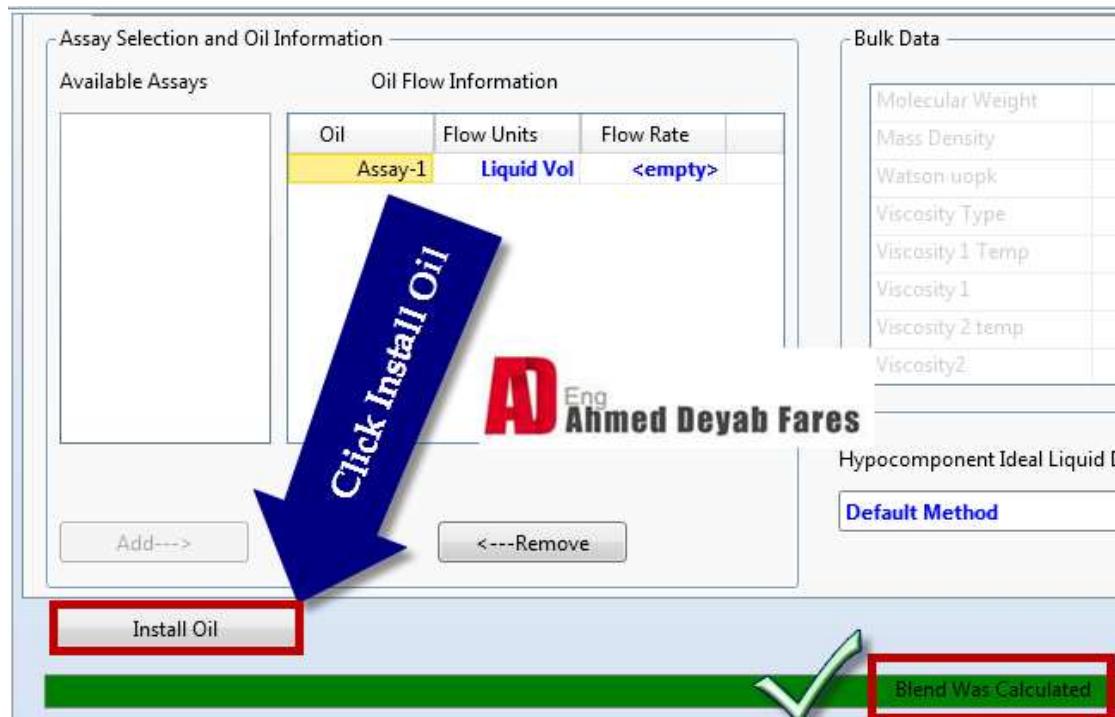


The results of the calculation can be viewed on the Tables tab

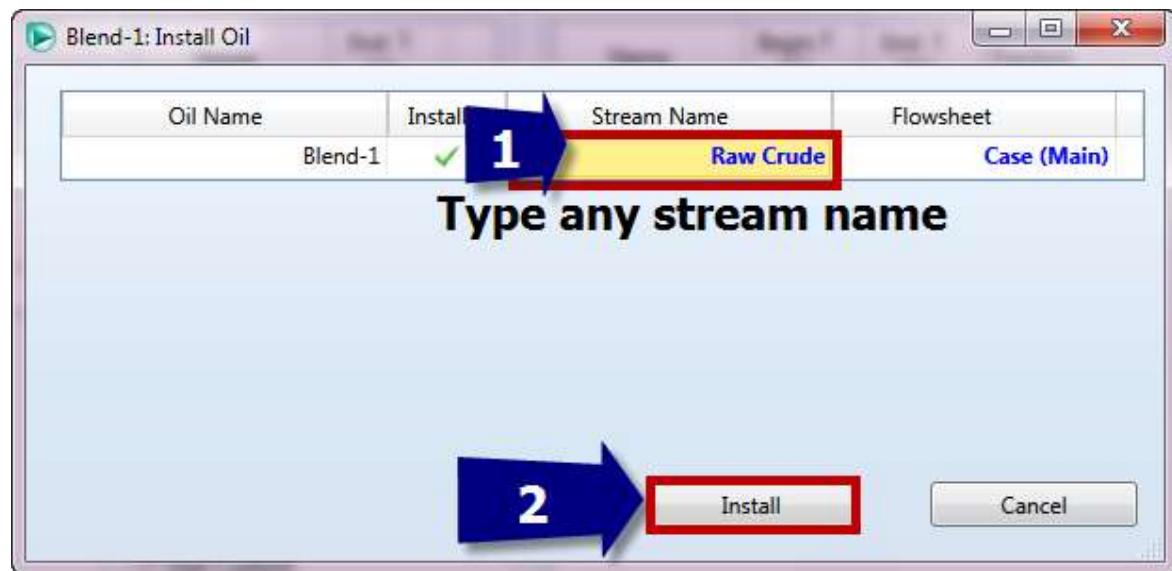
Component Physical Properties					
Comp Name	NBP [C]	Mole Wt.	Density [kg/m <sup>3</sup> ]	Viscos [cP]	
NBP_57	56.57	73.71	728.7	(	
NBP_68	68.42	78.92	737.5	(	
NBP_83	83.35	85.95	748.7	(	
NBP_96	95.96	91.89	757.4	(	
NBP_111	111.1	99.31	767.2	(	
NBP_125	124.6	105.8	775.5	(	
NBP_140	139.9	112.3	783.5	(	
NBP_153	152.7	119.6	792.1	(	
NBP_167	166.9	127.2	800.5	(	
NBP_181	181.1	135.5	808.9	(	
NBP_195	195.1	144.3	817.3	(	
NBP_209	209.3	153.8	825.7	(	
NBP_224	223.5	164.3	834.4	(	
NBP_237	237.5	175.5	843.0	(	
NBP_251	251.4	186.7	852.0	(	



The final step of the characterization is to transfer the hypo component information into the Flowsheet.

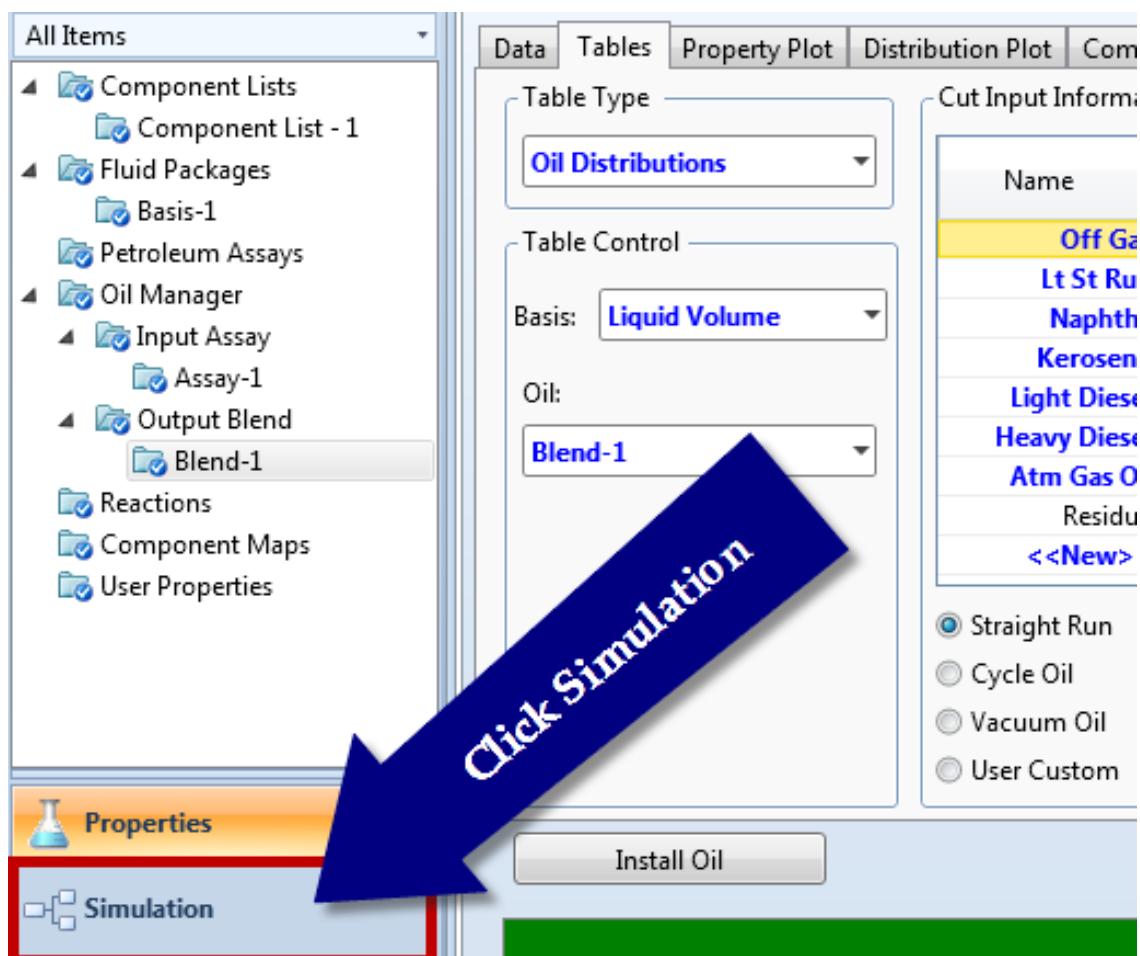


In the Stream Name column, enter the name **Raw Crude** to which the oil composition will be transferred.



HYSYS will assign the composition of your calculated Oil and Light Ends into this stream, completing the characterization process.

Now you can return to the Simulation environment to see the stream (Raw Crude) with a full composition:



All Items

- Component Lists
- Fluid Packages
- Oil Manager
- Reactions
- Component Maps
- User Properties

Component List - 1

Basis-1

Assay-1

Blend-1

Properties

Simulation

Data Tables Property Plot Distribution Plot Com

Table Type: Oil Distributions

Table Control: Basis: Liquid Volume

Oil: Blend-1

Cut Input Information

Name

- Off Gas
- Lt St Ru
- Naphth
- Kerosen
- Light Diesel
- Heavy Diesel
- Atm Gas O
- Residu
- <<New>>

Straight Run

Cycle Oil

Vacuum Oil

User Custom

Install Oil

Click Simulation

The screenshot shows the HYSYS interface. On the left, the Stream Map palette is open, displaying various process icons categorized into Upstream, Refining, Custom, Dynamics, Common, and Columns sections. A stream labeled "Raw Crude" is highlighted with a red border and a blue arrow pointing to it.

The main window displays the "Material Stream: Raw Crude" worksheet. The left sidebar contains tabs for Worksheet, Attachments, and Dynamics, with "Worksheet" selected. Under the "Composition" tab, a table lists the mole fractions of various components:

	Mole Fractions
Methane	0.0003
Ethane	0.0006
Propane	0.0086
i-Butane	0.0054
n-Butane	0.0410
i-Pentane	0.0333
n-Pentane	0.0458
H <sub>2</sub> O	0.0000
NBP[0]57*	0.0282
NBP[0]68*	0.0380
NBP[0]83*	0.0445
NBP[0]96*	0.0371
NBP[0]111*	0.0319

The total value is listed as 1.00000. Below the table are buttons for Edit..., View Properties..., and Basis... . A yellow bar at the bottom indicates "Unknown Temperature".

*Save Your Case!*

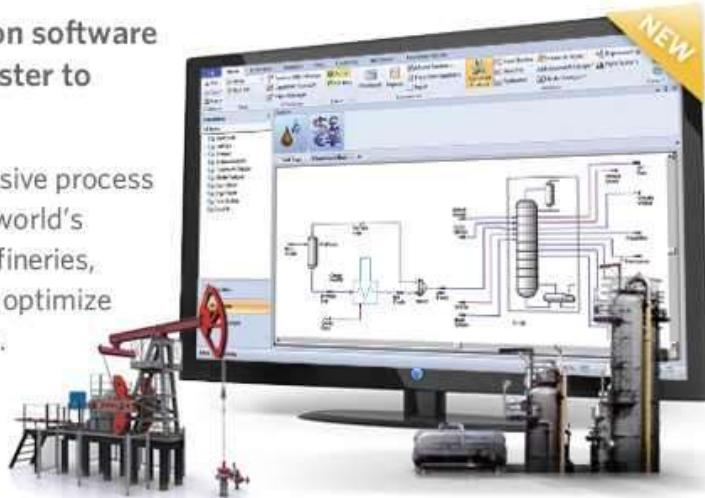
# Atmospheric Distillation

Experience the New Aspen HYSYS®.

V8

The best process simulation software is now easier to use and faster to learn than ever!

Aspen HYSYS is a comprehensive process modeling system used by the world's leading oil & gas producers, refineries, and engineering companies to optimize process design and operations.



6

## Workshop

*Atmospheric Crude Columns are one of the most important pieces of equipment in the petroleum refining industry. Typically located after the Desalter and the Crude Furnace, the Atmospheric Tower serves to distil the crude oil into several different cuts. These include naphtha, kerosene, light diesel, heavy diesel and AGO.*

*In this module, you will construct, run, analyze and manipulate an Atmospheric Crude Column simulation. You will begin by building a simple column and continue by adding side operations to the column.*

## Learning Objectives

- Build and converge an Atmospheric Crude Column.
- Use HYSYS to analyze and predict the behavior of a simulated column.
- Add side operations to a column to improve operation and efficiency.
- Add cut point specifications to increase side product quality and quantity.

### Pre-Heat Train

A crude stream at 15°C, 1000 kPa and flowrate of  $6 \times 10^5$  kg/hr is mixed with a stream of water at 15°C, 1000 kPa and flowrate of 21600 kg/hr using a Mixer, the outlet from the mixer is then heated to 65°C through a Heater ( $\Delta p=50$  kPa), the heater outlet is fed to the tube side of a Shell & Tube Heat Exchanger, where it's heated using a Shell inlet stream having the same composition as the crude feed stream and enters the shell of the heat exchanger at 180°C, 200 kPa and flowrate of 175m<sup>3</sup>/hr. The pressure drops for the Tube and Shell sides, will be 35 kPa and 5 kPa, respectively. The tube outlet from the HX is then sent to a desalter which is simply modeled as Three Phase Separator where desalted water, oil and gas is separated. The oil stream (light liquid) from the desalter is then heated to 175°C through a Heater (Pressure drop=375 kPa) and then sent to a Preflash (Separator) to reduce the light components in the feed. The liquid product from the separator is then heated to 400°C inside a Heater (Pressure drop=250 kPa) before entering the Atmospheric Column.

#### Heat Exchanger Specification:

- Use Simple weighted model
- Min Approach = 30°C (54°F). This is the minimum temperature difference between the hot and the cold stream.

#### Calculate:

- The vapor fraction of the product stream before entering the Atmospheric column.
- The Shell side outlet Temperature.
- The vapor molar flow rate from the Preflash.

## Atmospheric Distillation

A feed stream from the pre-heat train is fed to the 28<sup>th</sup> tray of a **Refluxed Absorber** with 29 trays and a partial condenser to separate Off Gases, Naphtha and Bottom Residue. A steam stream (vapor fraction =1.0, pressure =1380 kPa and flowrate=3400 kg/hr) is fed to the bottom of the tower to provide the necessary heat. A water draw stream is required to remove the condensed steam from the overhead condenser. The tower is operated with the following conditions:

Condenser Pressure	140 kPa
Condenser Pressure Drop	60 kPa
Bottom Stage Pressure	230 kPa

Condenser Temperature	40 °C
Top Stage Temperature	120 °C
Bottom Stage Temperature	340 °C

Specifications:

Overhead vapor rate	0.0 kgmole/hr
Distillate rate (Naphtha rate)	.....

What is the flow rate of?

Naphtha \_\_\_\_\_ Residue \_\_\_\_\_

Wastewater \_\_\_\_\_

## Side Strippers & Pump Arounds

**Side Strippers** are added to the column in order to improve the quality of the three main products (Kerosene, Diesel, and AGO). There are two types of side strippers available in HYSYS: Reboiled and Steam Stripped. We will install one reboiled side stripper and two steam stripped.

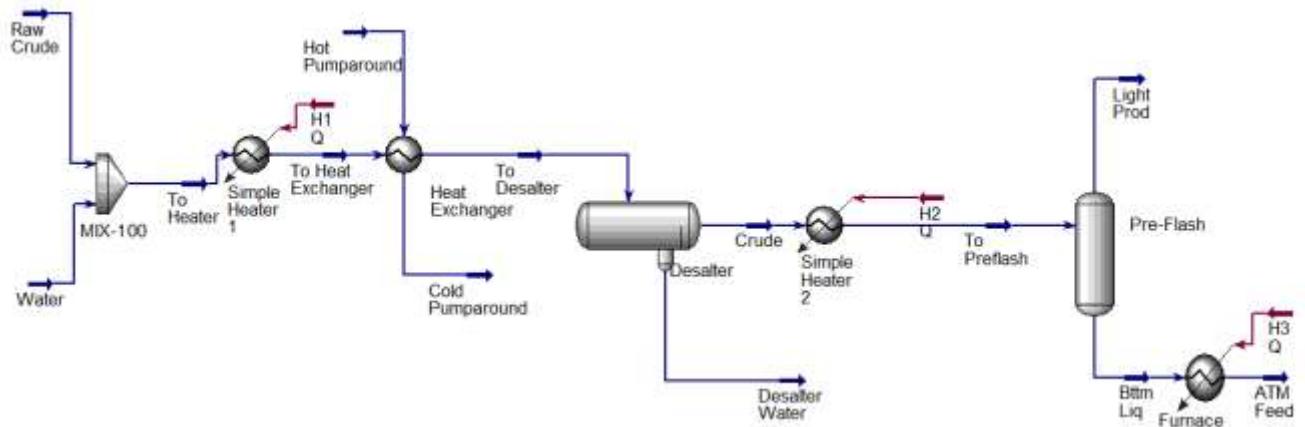
	AGO SS	Diesel SS	Kerosene SS
Draw Stage	22	17	9
Return Stage	21	16	8
Configuration	Steam Stripped	Steam Stripped	Reboiled
Product flow	.....	.....	.....

	Temperature	Pressure	Flowrate
AGO Steam	150 °C	350 kPa	1150 kg/hr
Diesel Steam	150 °C	350 kPa	1350 kg/hr

**Pump Arounds** help to improve the column's efficiency. They operate by drawing a liquid stream from one stage cooling it, and pumping it into a higher stage. In effect, this process adds to the reflux between these two stages.

	AGO PA	Diesel PA	Kerosene PA
Draw Stage	22	17	9
Return Stage	21	16	8
Flowrate	200 m <sup>3</sup> /h	200 m <sup>3</sup> /h	330 m <sup>3</sup> /h
Duty	-3.7e7 kJ/h	-3.7e7 kJ/h	-4.5e7 kJ/h

**Load your Pre-Heat Train case from the Pre-Heat Train module.**



Add a steam stream to provide the heating effect to the bottom of the tower instead of using a reboiler:

Vapor fraction =1.0, pressure =1380 kPa and flowrate=3400 kg/hr

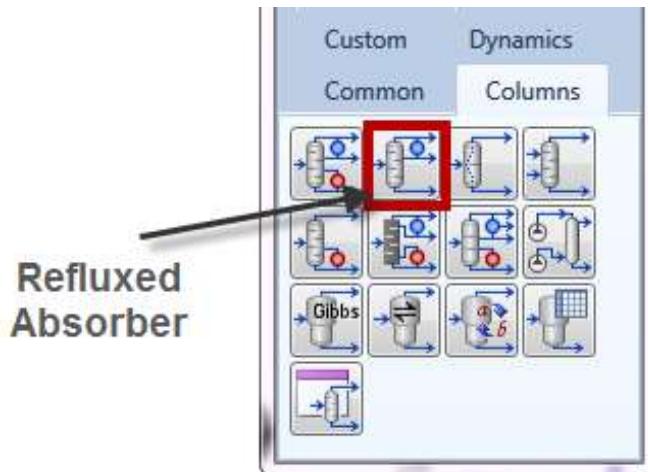
Composition (100% water)

Worksheet	Attachments	Dynamics
<b>Worksheet</b>	Stream Name	<b>Bottom Steam</b>
Conditions	Vapour / Phase Fraction	<b>1.0000</b>
Properties	Temperature [C]	194.6
Composition	Pressure [kPa]	<b>1380</b>
Oil & Gas Feed	Molar Flow [kgmole/h]	188.7
Petroleum Assay	Mass Flow [kg/h]	<b>3400</b>
K Value	Std Ideal Liq Vol Flow [m3/h]	3.407
User Variables	Molar Enthalpy [kJ/kgmole]	-2.359e+005
Notes	Molar Entropy [kJ/kgmole-C]	166.3
Cost Parameters	Heat Flow [kJ/h]	-4.453e+007
Normalized Yields	Liq Vol Flow @Std Cond [m3/h]	3.350
	Fluid Package	<b>Refinery</b>
	Utility Type	

## ADDING THE ATMOSPHERIC COLUMN

The atmospheric column is modeled as a refluxed absorber,

- Double-click on the **Refluxed Absorber** icon on the Object Palette.



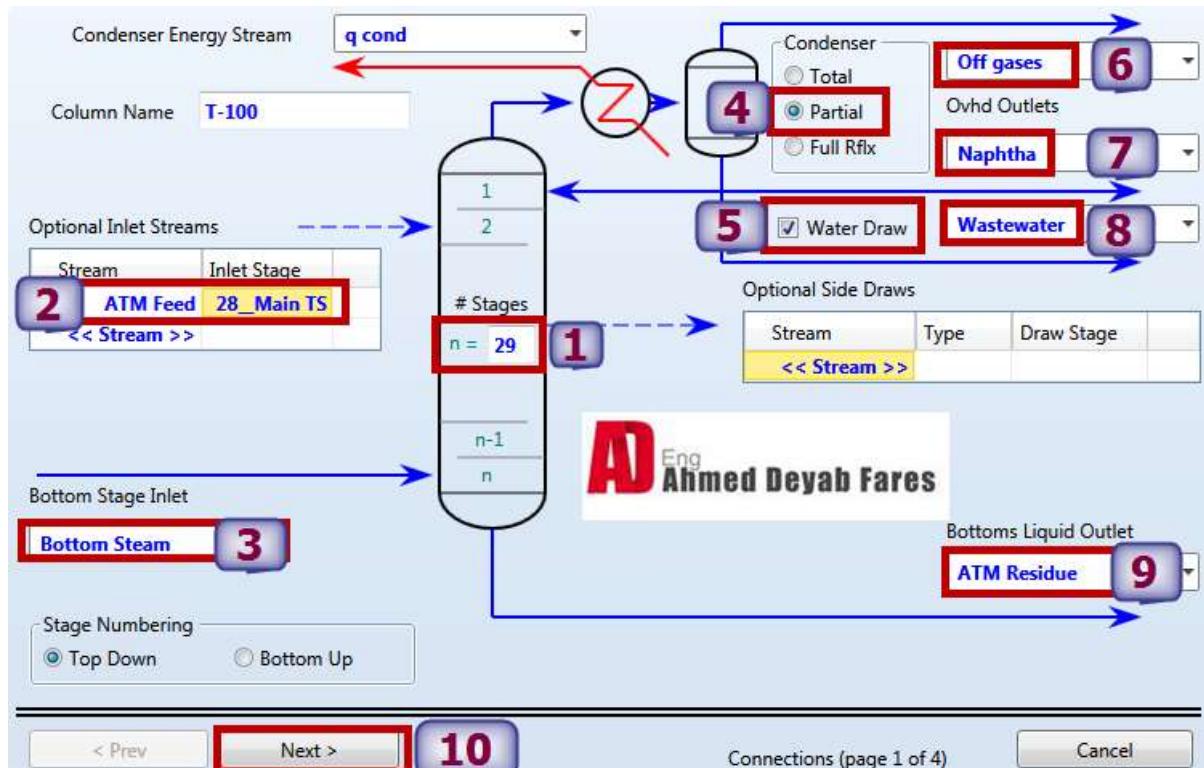
The tower (29 trays) is operated with the following conditions:

Condenser Pressure	140 kPa	Condenser Temperature	40 °C
Condenser Pressure Drop	60 kPa	Top Stage Temperature	120 °C
Bottom Stage Pressure	230 kPa	Bottom Stage Temperature	340 °C

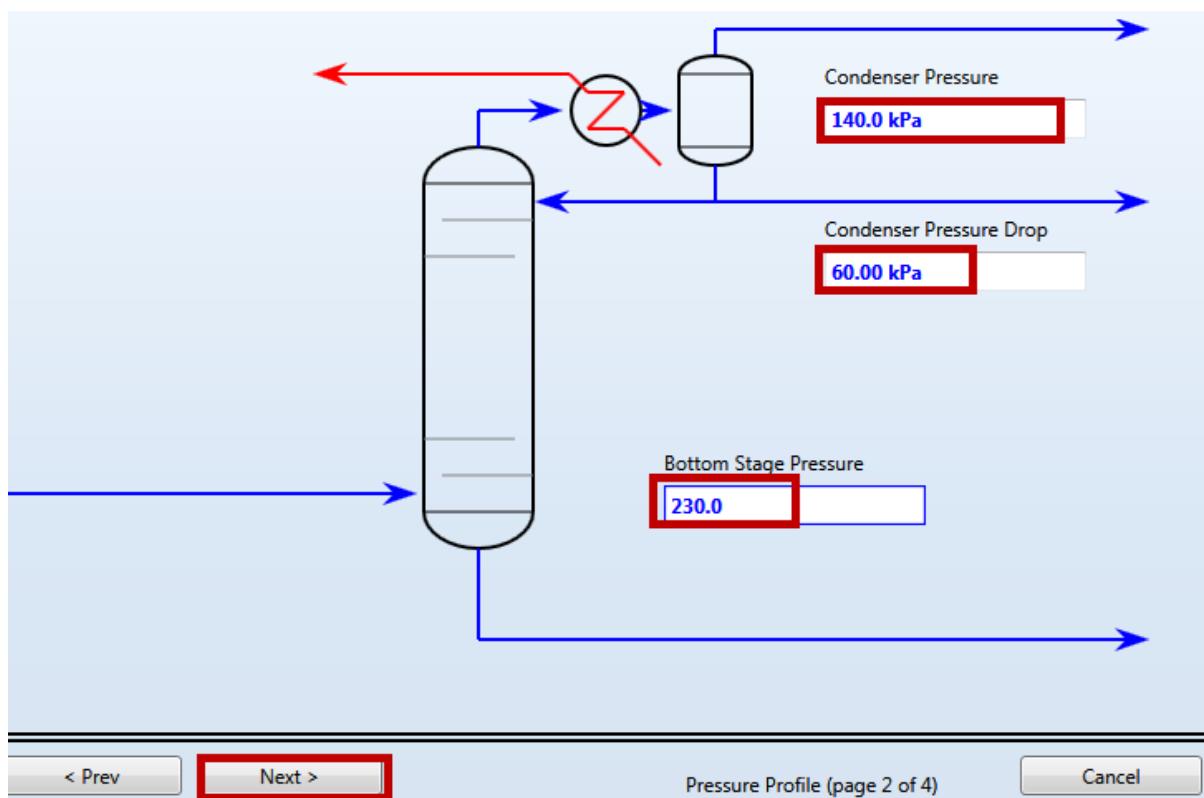
Specifications:

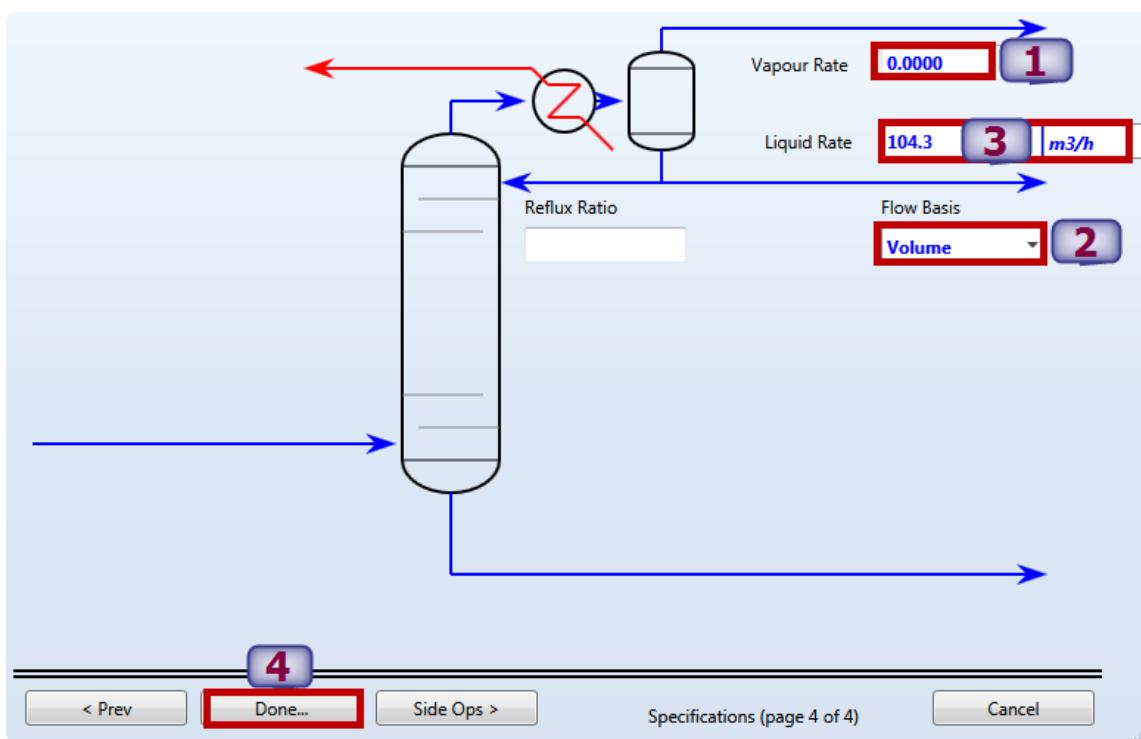
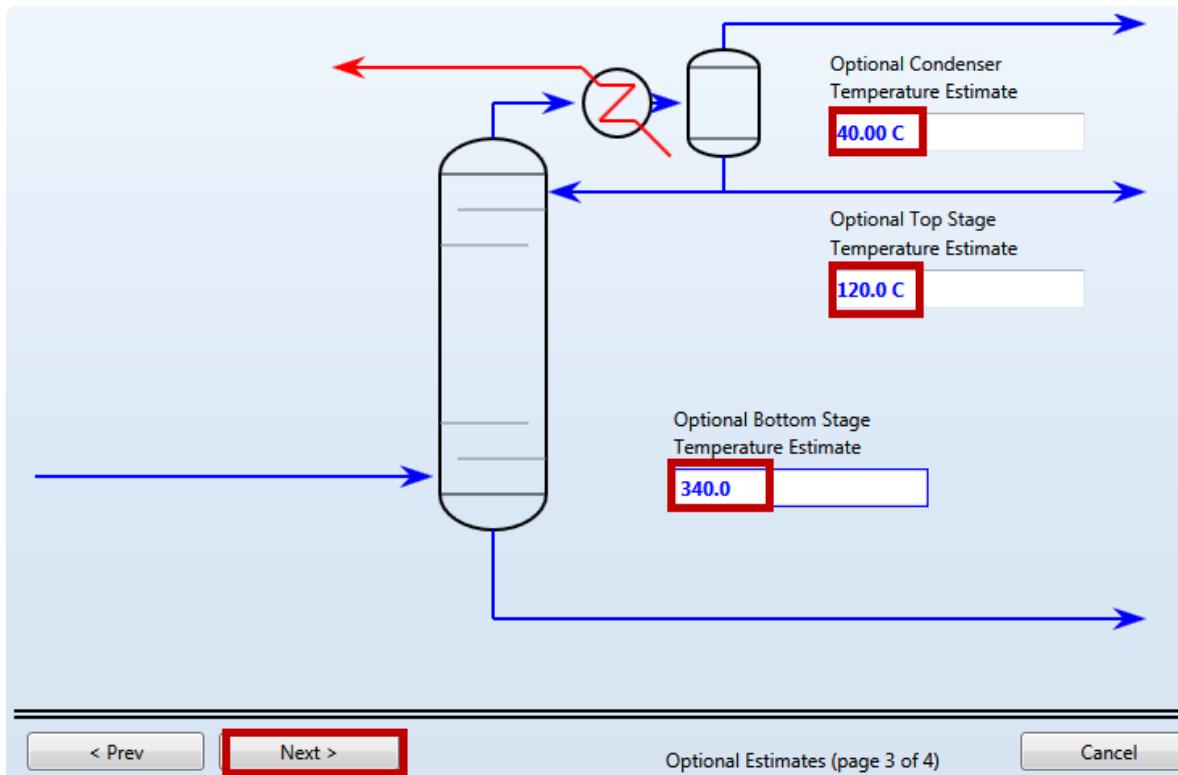
Overhead vapor rate	0.0 kgmole/hr
Distillate rate (Naphtha rate)	$0.153 \times 681.9 \text{ m}^3/\text{hr} = 104.3 \text{ m}^3/\text{hr}$

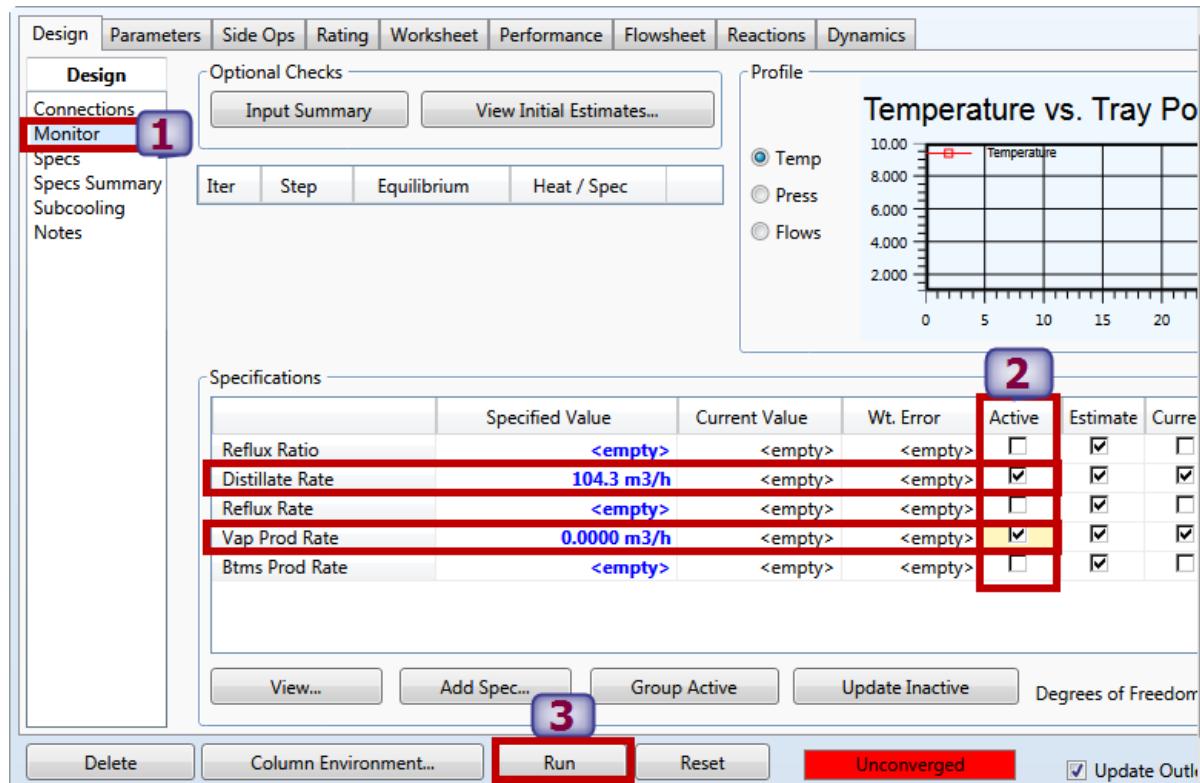
Cut Input Information		Cut Distributions			
Name	End T [C]	Name	Begin T [C]	End T [C]	Fraction
Off Gas	10.00	Off Gas	-19.54	10.00	0.021
Lt St Run	70.00	Lt St Run	10.00	70.00	0.064
Naphtha	180.0	Naphtha	70.00	180.0	0.153
Kerosene	240.0	Kerosene	180.0	240.0	0.094
Light Diesel	290.0	Light Diesel	240.0	290.0	0.102
Heavy Diesel	340.0	Heavy Diesel	290.0	340.0	0.087
Atm Gas Oil	370.0	Atm Gas Oil	340.0	370.0	0.045
Residue	1200	Residue	370.0	825.2	0.434
..	..				



The Water Draw checkbox must be checked to prevent two liquid phases being formed in the column.







Run the column

## Adding Side Strippers

**Side Strippers** are added to the column in order to improve the quality of the three main products (Kerosene, Diesel, and AGO). There are two types of side strippers available in HYSYS: Reboiled and Steam Stripped. We will install one reboiled side stripper and two steam stripped.

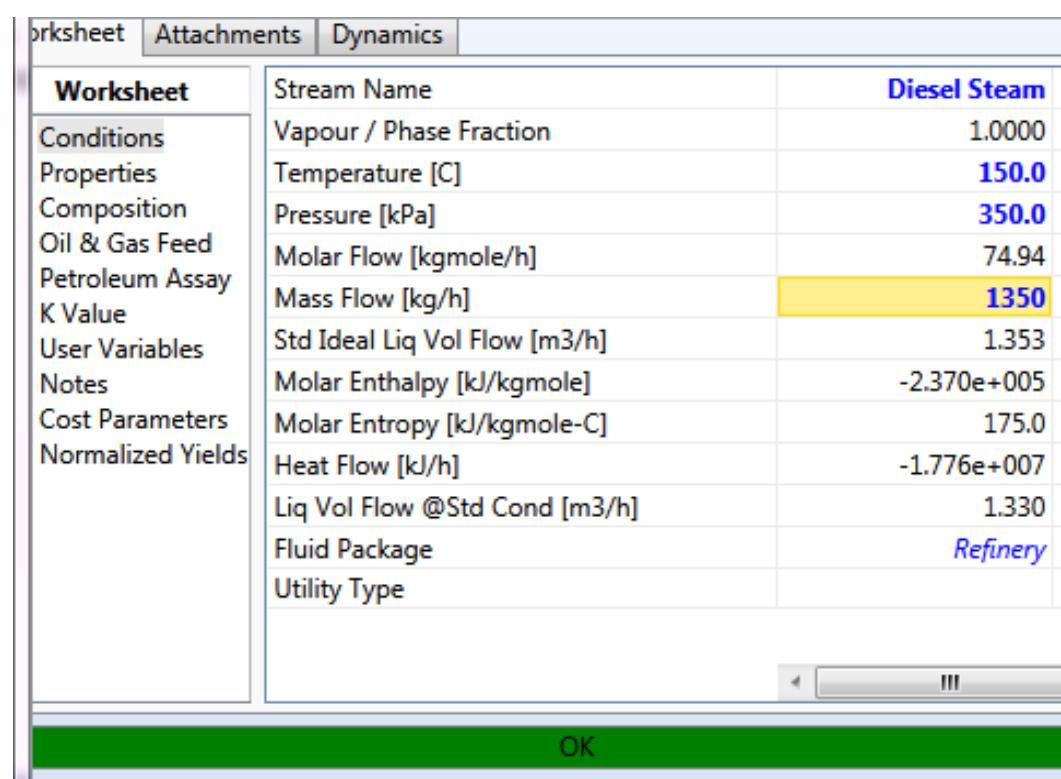
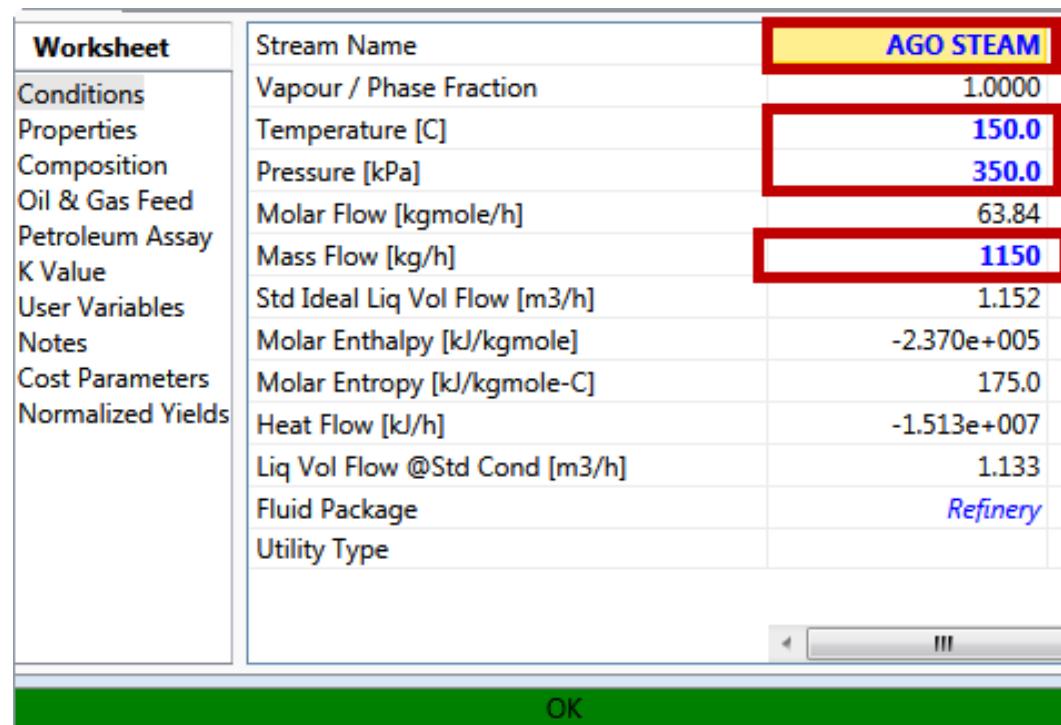
	AGO SS	Diesel SS	Kerosene SS
Draw Stage	22	17	9
Return Stage	21	16	8
Configuration	Steam Stripped	Steam Stripped	Reboiled
Product flow	$0.045 \times 681.9 \text{ m}^3/\text{hr}$ $= 30.69 \text{ m}^3/\text{hr}$	$(0.102+0.087) \times 681.9 \text{ m}^3/\text{hr}$ $= 128.88 \text{ m}^3/\text{hr}$	$0.094 \times 681.9 \text{ m}^3/\text{hr}$ $= 64.1 \text{ m}^3/\text{hr}$

Cut Input Information		Cut Distributions			
Name	End T [C]	Name	Begin T [C]	End T [C]	Fraction
Off Gas	10.00	Off Gas	-19.54	10.00	0.021
Lt St Run	70.00	Lt St Run	10.00	70.00	0.064
Naphtha	180.0	Naphtha	70.00	180.0	0.153
Kerosene	240.0	Kerosene	180.0	240.0	0.094
Light Diesel	290.0	Light Diesel	240.0	290.0	0.102
Heavy Diesel	340.0	Heavy Diesel	290.0	340.0	0.087
Atm Gas Oil	370.0	Atm Gas Oil	340.0	370.0	0.045
Residue	1200	Residue	370.0	825.2	0.434
..	..				

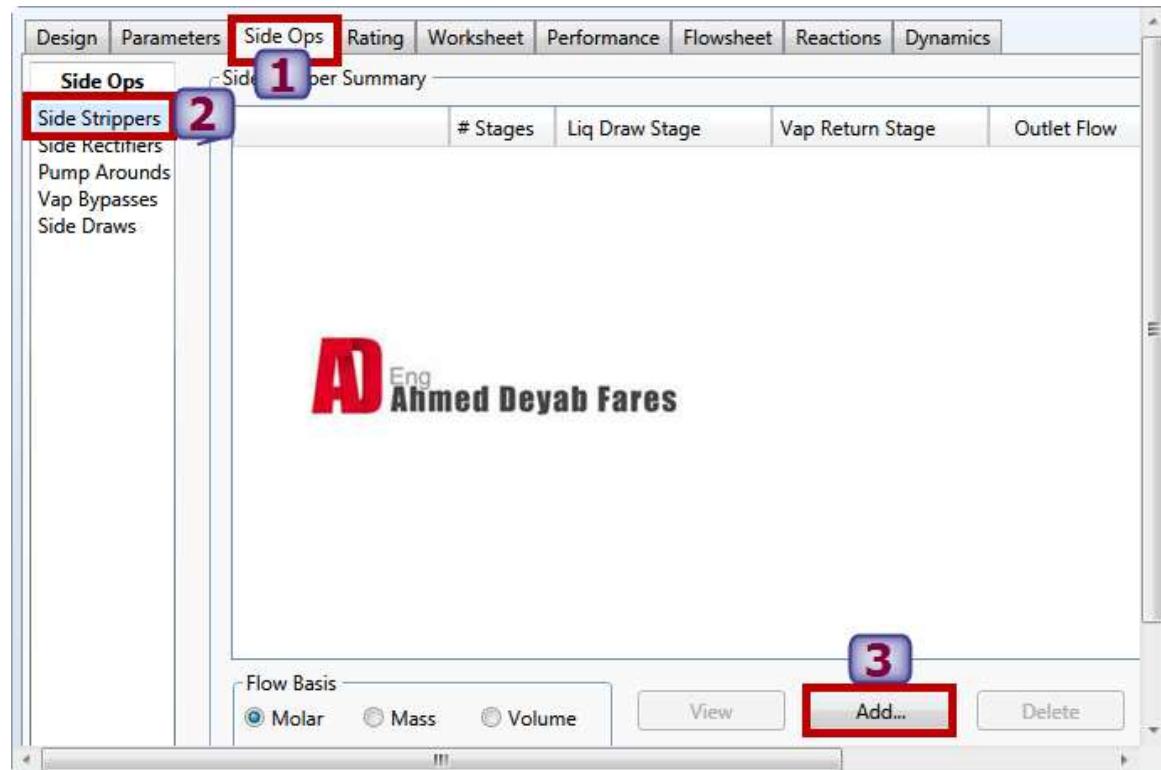
	Temperature	Pressure	Flowrate
AGO Steam	150 °C	350 kPa	1150 kg/hr
Diesel Steam	150 °C	350 kPa	1350 kg/hr

First we need to add two steam streams to provide the heating effect for both AGO & Diesel Side Strippers

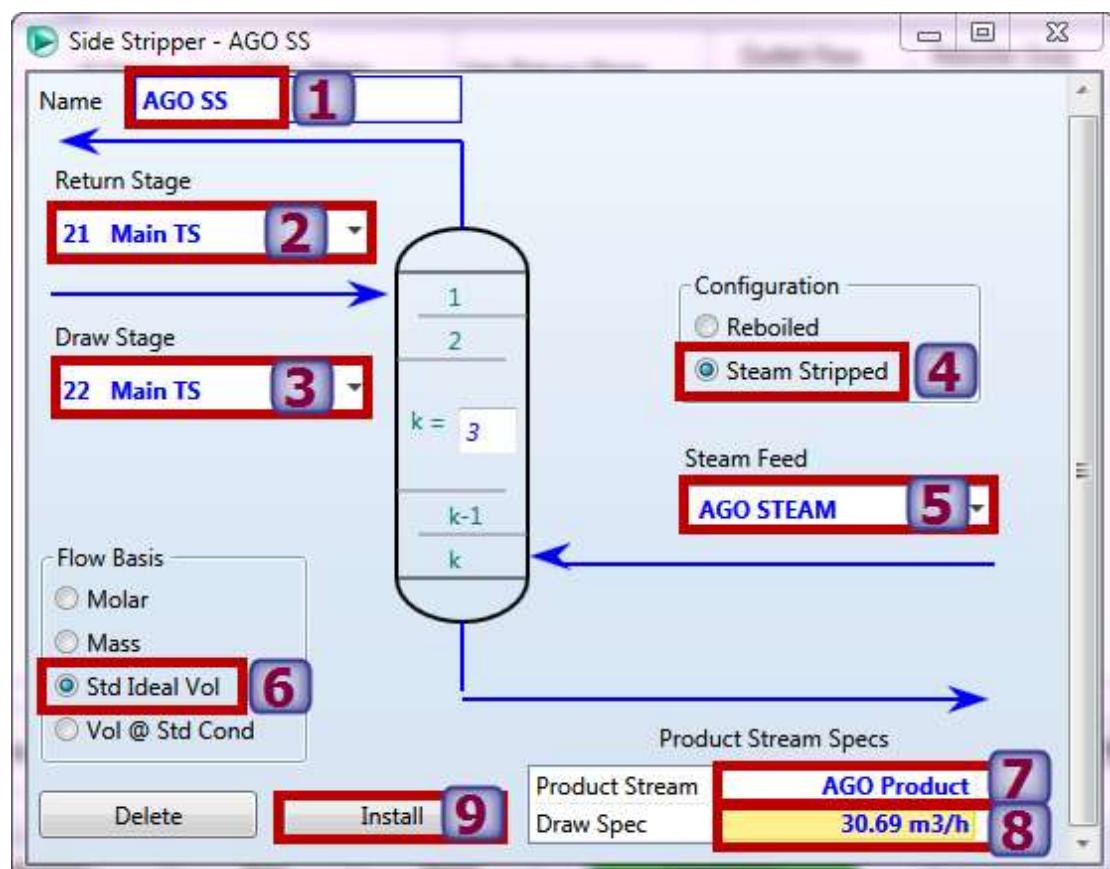
	Temperature	Pressure	Flowrate
AGO Steam	150 °C	350 kPa	1150 kg/hr
Diesel Steam	150 °C	350 kPa	1350 kg/hr

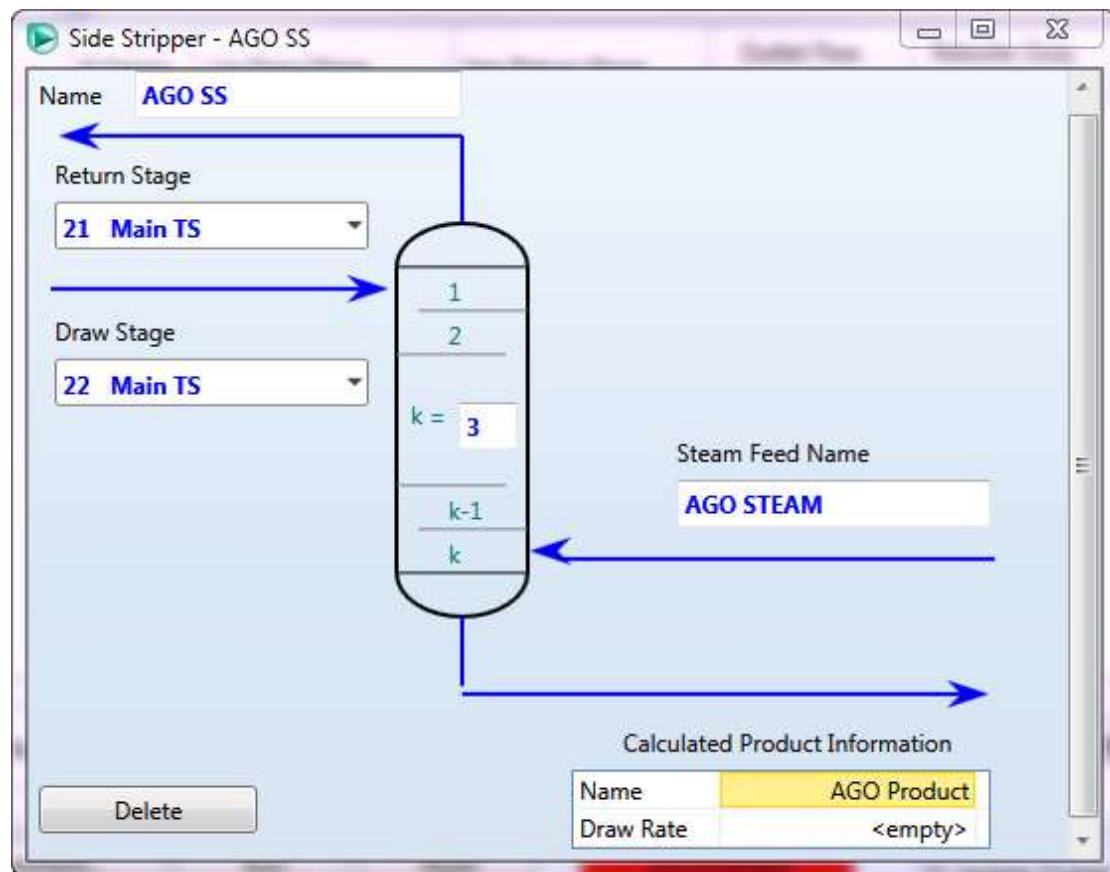


Go to the *Side Ops* tab inside the atmospheric tower to *add* the three side strippers:



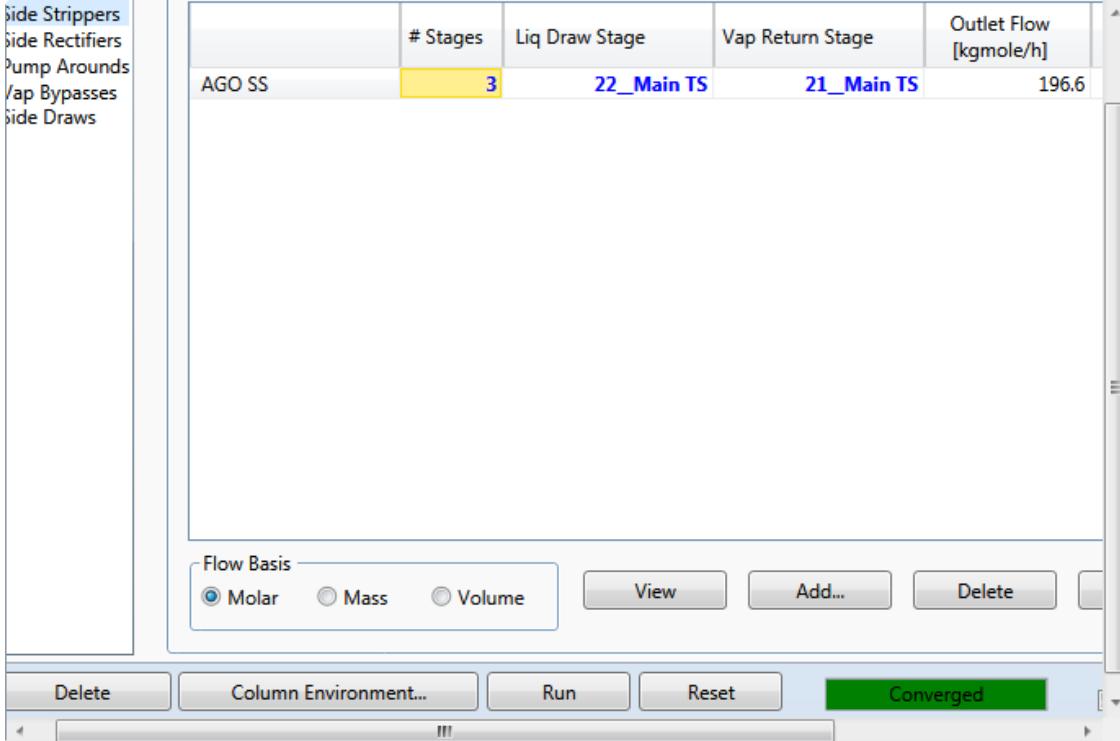
First add the AGO Side Stripper





**Close the window and run the column**

**Make sure that the column is converged**



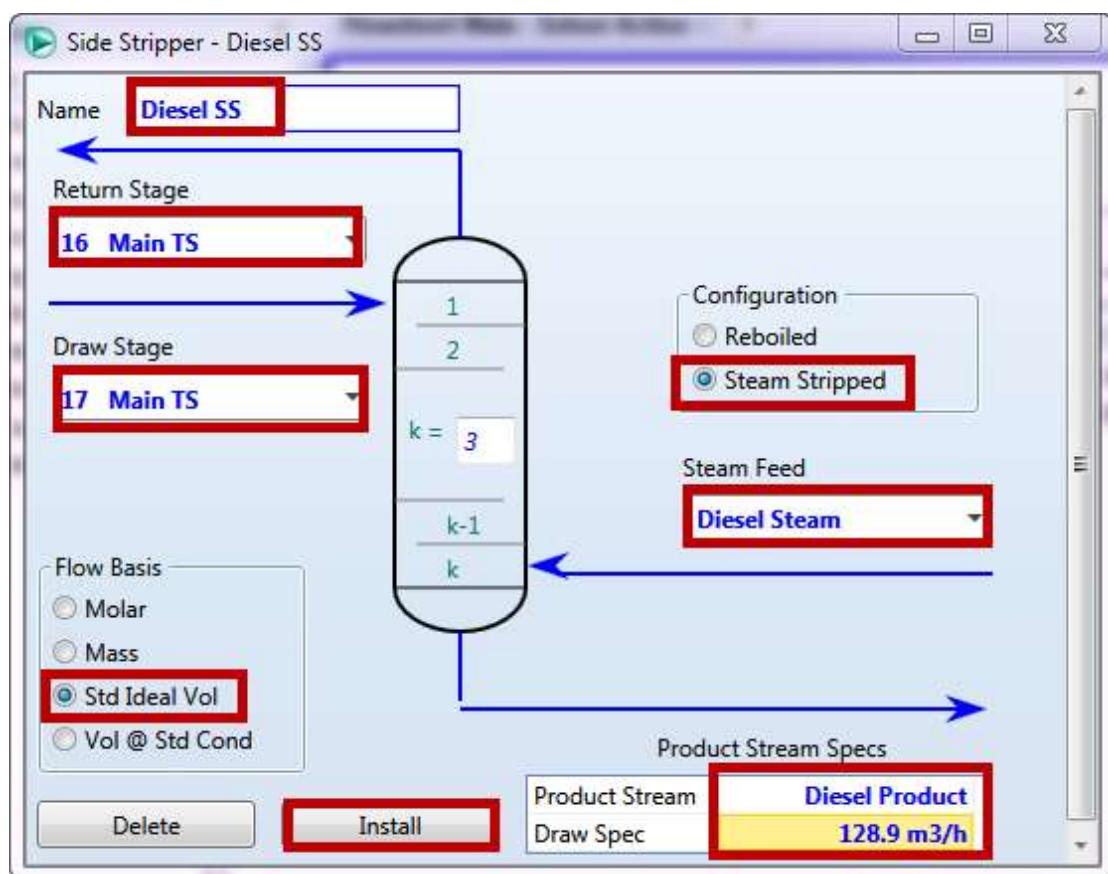
Side Strippers	# Stages	Liq Draw Stage	Vap Return Stage	Outlet Flow [kgmole/h]
AGO SS	3	22_Main TS	21_Main TS	196.6

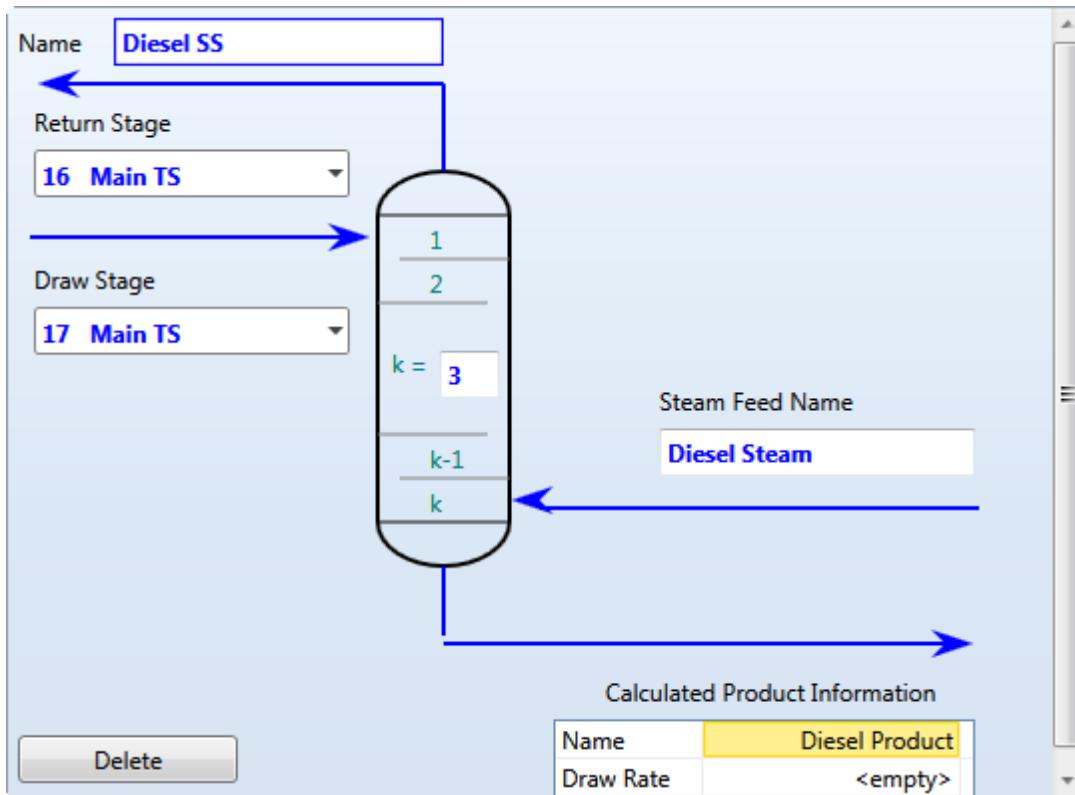
Flow Basis:

Molar    Mass    Volume

View Add... Delete

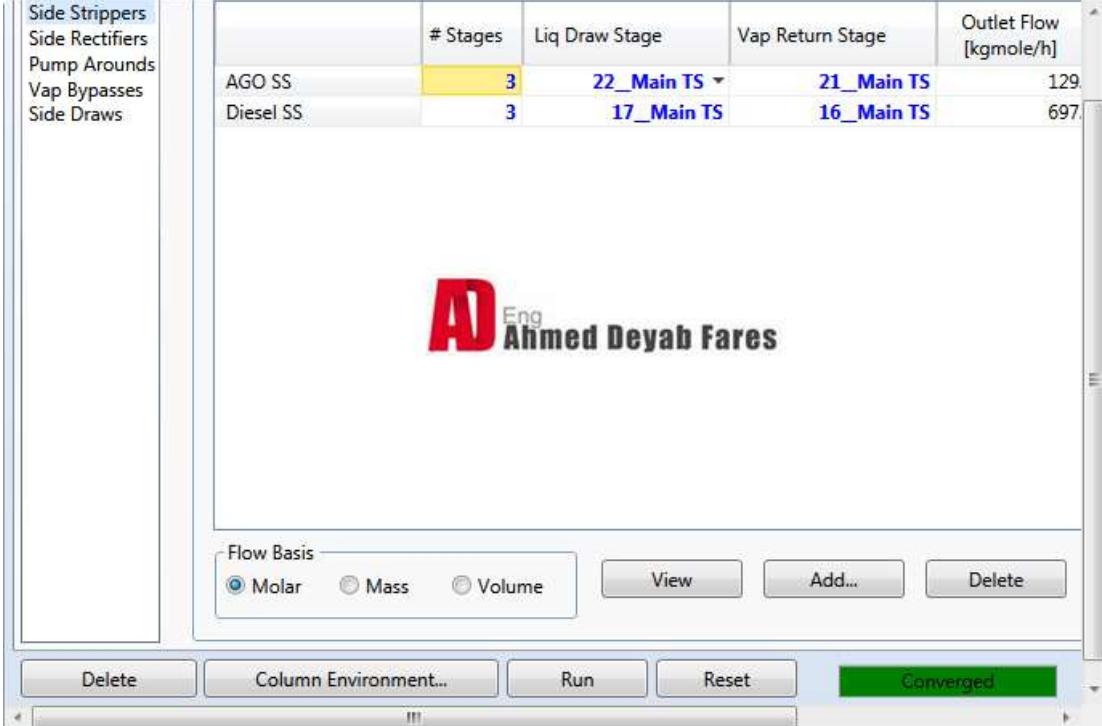
Converged

Add the Diesel Side stripper:



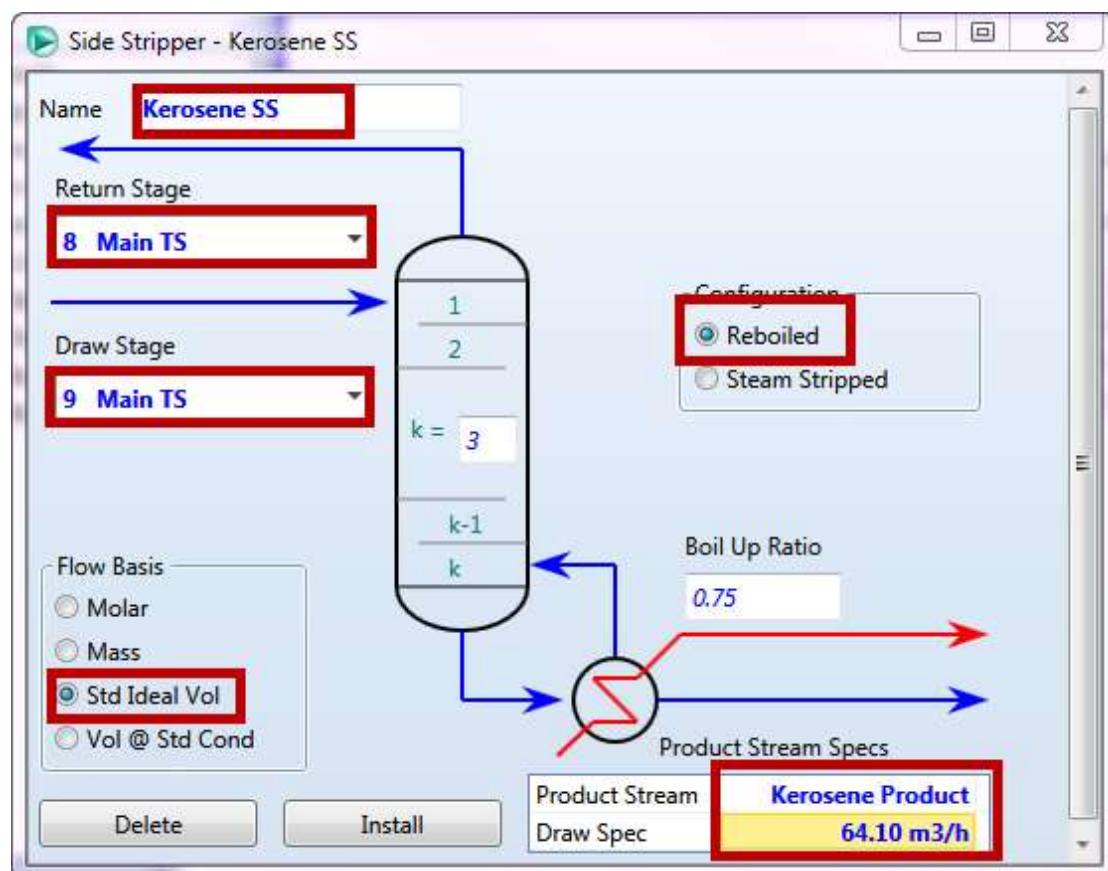
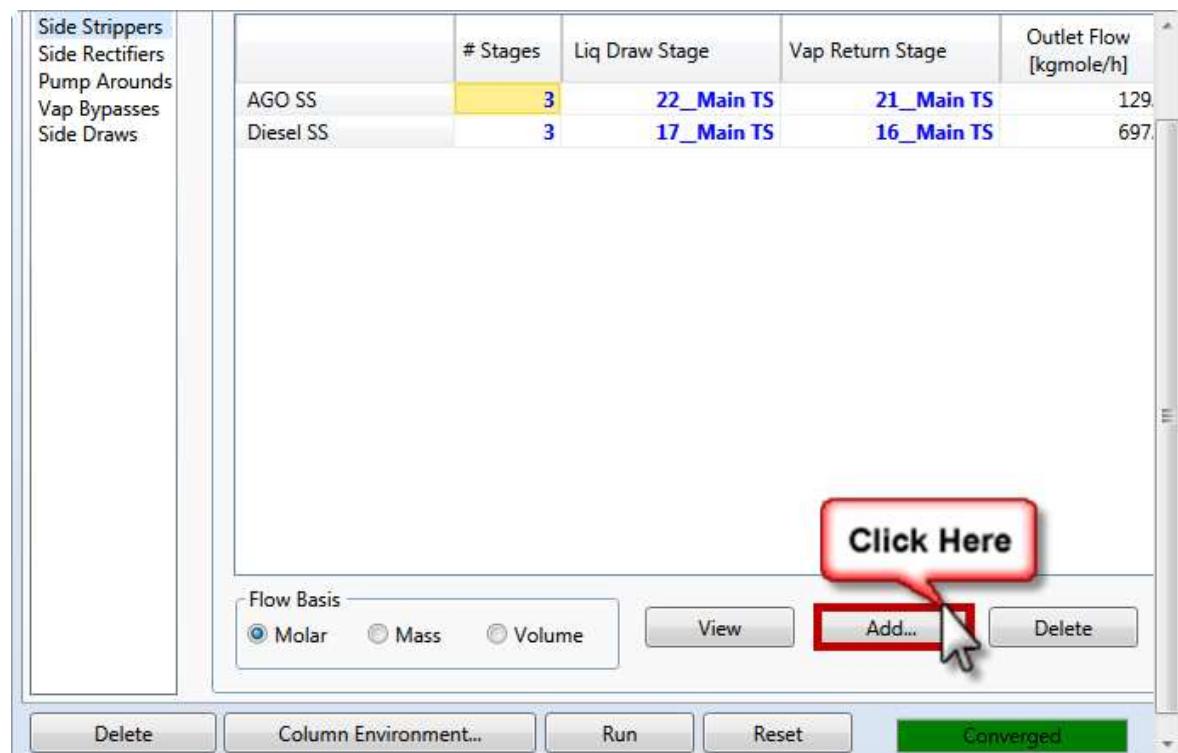
**Close the window and run the column**

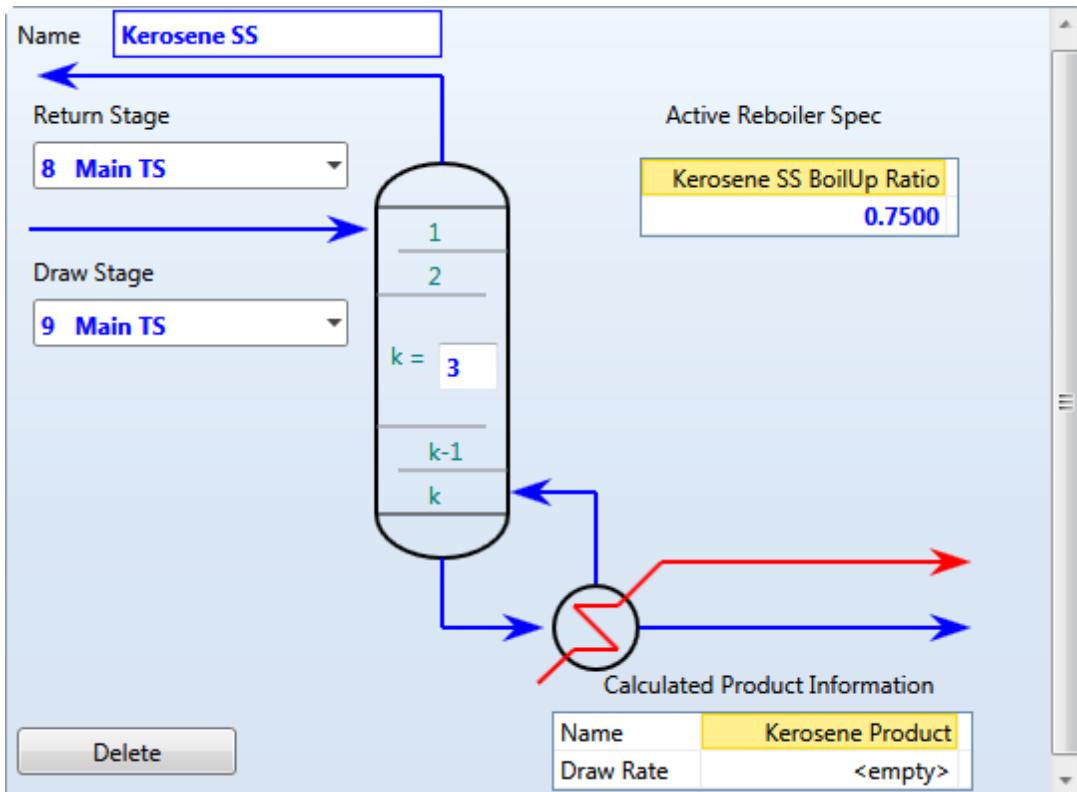
**Make sure that the column is converged**



The table displays the properties of two columns: AGO SS and Diesel SS. Both columns have 3 stages, with AGO SS having a liquid draw stage of 22\_Main TS and a vapor return stage of 21\_Main TS, resulting in an outlet flow of 129 kgmole/h. Diesel SS has a liquid draw stage of 17\_Main TS and a vapor return stage of 16\_Main TS, resulting in an outlet flow of 697 kgmole/h. The 'Converged' button at the bottom right is highlighted in green, indicating successful convergence.

### Add the Kerosene Side stripper:





**Close the window and run the column**

**Make sure that the column is converged**

The table displays the properties of three columns: AGO SS, Diesel SS, and Kerosene SS. All three columns have 3 stages, a Liq Draw Stage of 22\_Main TS, and a Vap Return Stage of 21\_Main TS, 17\_Main TS, and 9\_Main TS respectively. The Kerosene SS column has an outlet flow of 386.2 kgmole/h and a reboiler duty of 1.2177e+007 kJ/h.

	# Stages	Liq Draw Stage	Vap Return Stage	Outlet Flow [kgmole/h]	Reboiler Duty [kJ/h]
AGO SS	3	22_Main TS	21_Main TS	107.9	
Diesel SS	3	17_Main TS	16_Main TS	569.0	
Kerosene SS	3	9_Main TS	8_Main TS	386.2	1.2177e+007

Flow Basis: Molar (radio button selected)

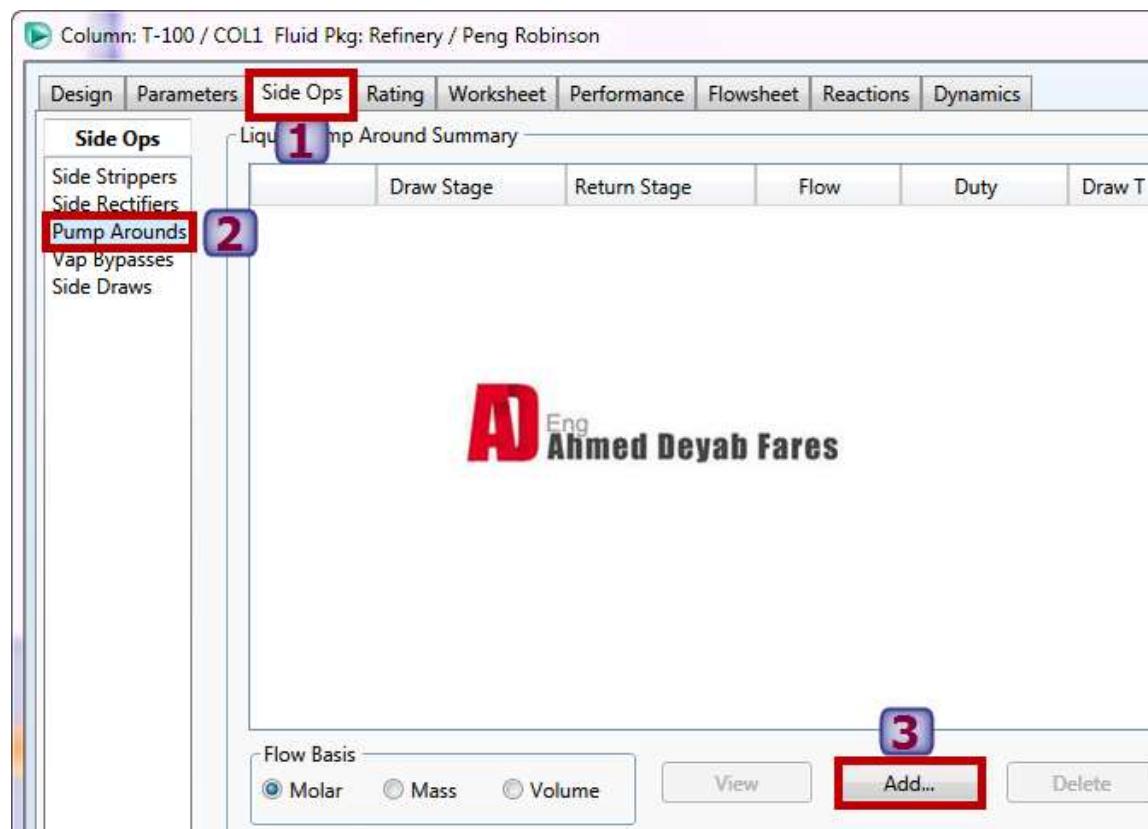
Buttons: View, Add..., Delete, Side Ops Input Export, Delete, Column Environment..., Run, Reset, Converged (green), Update Outlets (checkbox checked).

## Adding Pump Arounds

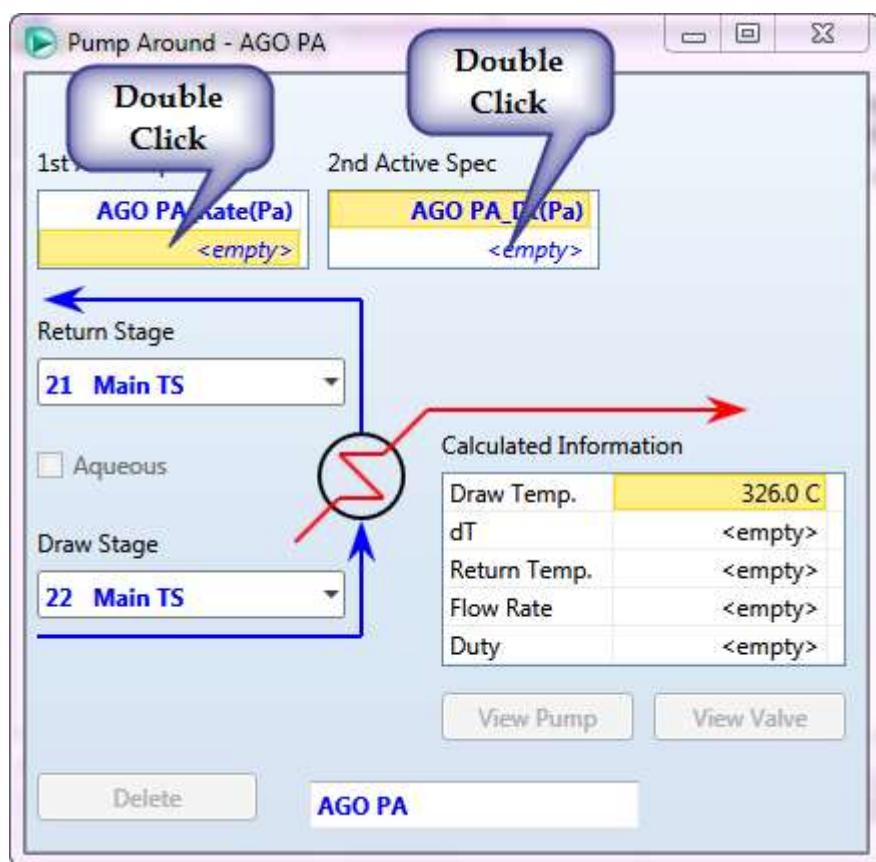
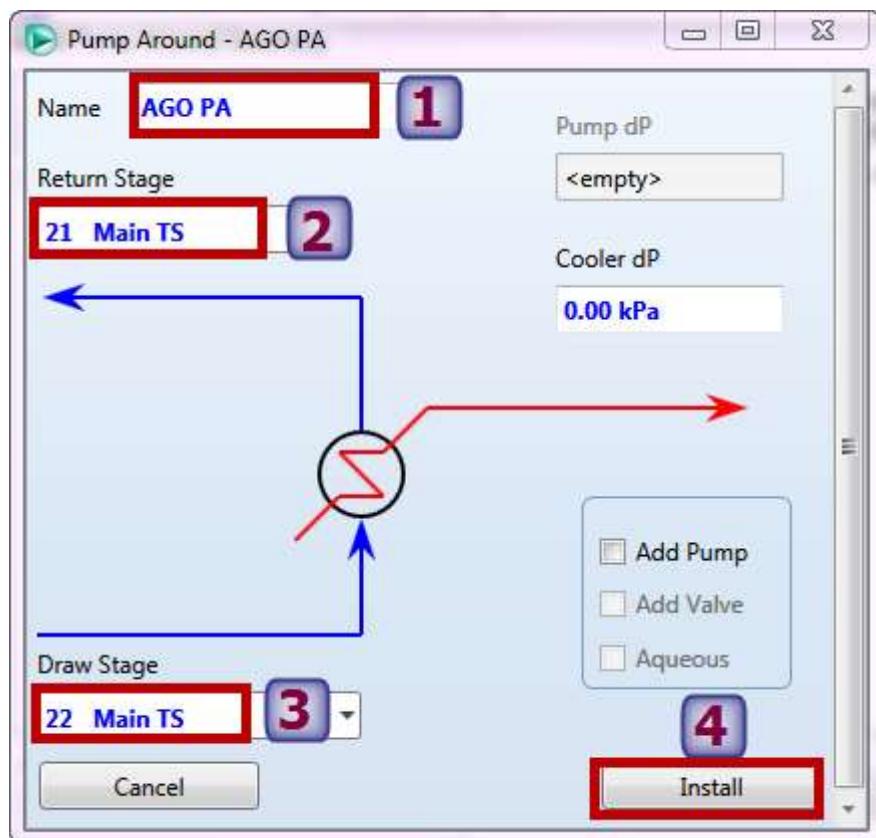
**Pump Arounds** help to improve the column's efficiency. They operate by drawing a liquid stream from one stage cooling it, and pumping it into a higher stage. In effect, this process adds to the reflux between these two stages.

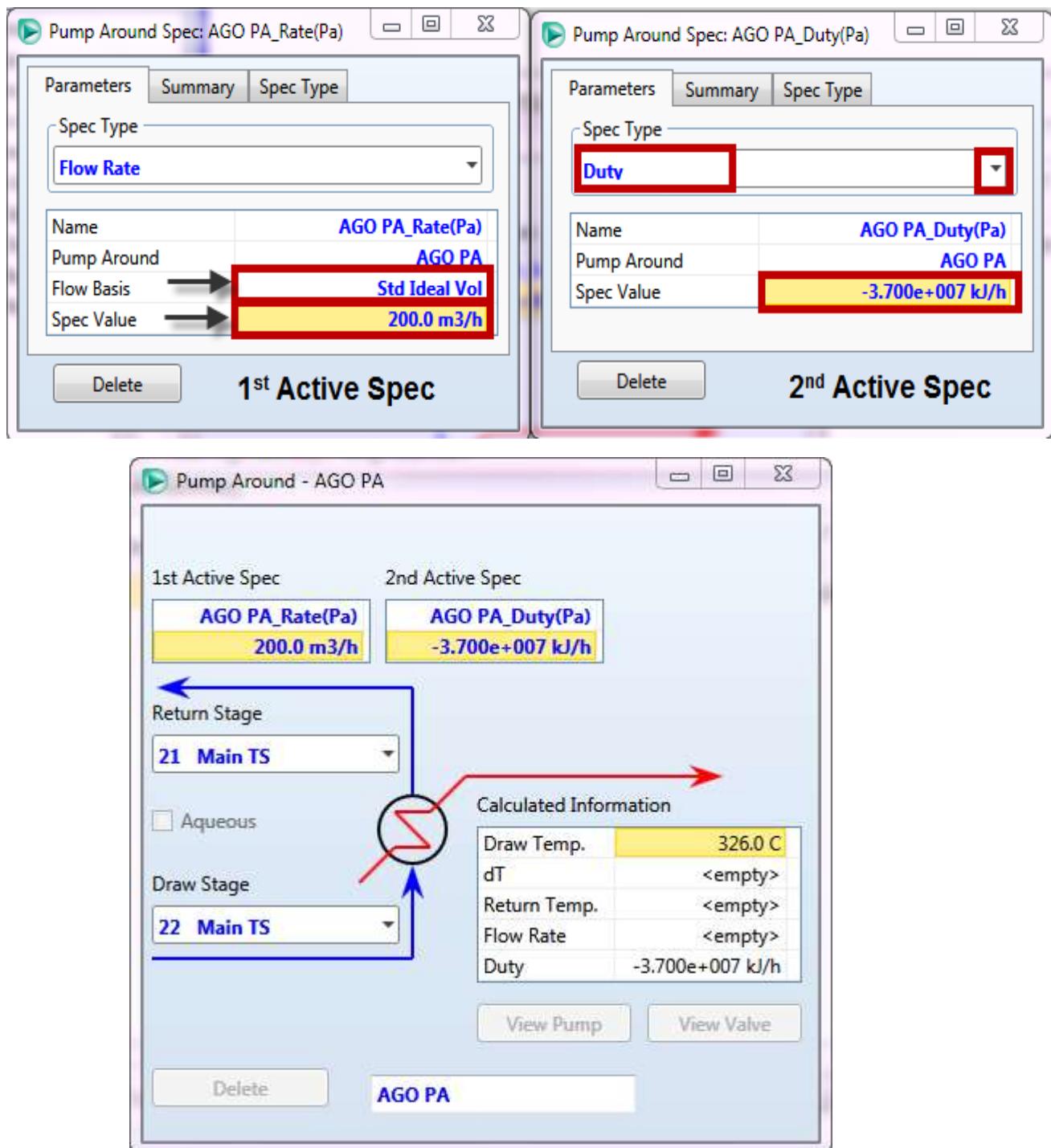
	AGO PA	Diesel PA	Kerosene PA
Draw Stage	22	17	9
Return Stage	21	16	8
Flowrate	200 m <sup>3</sup> /h	200 m <sup>3</sup> /h	330 m <sup>3</sup> /h
Duty	-3.7e7 kJ/h	-3.7e7 kJ/h	-4.5e7 kJ/h

Go to the **Side Ops** tab inside the atmospheric tower to **add** the three Pump Arounds:



## Add the AGO Pump Around (PA)

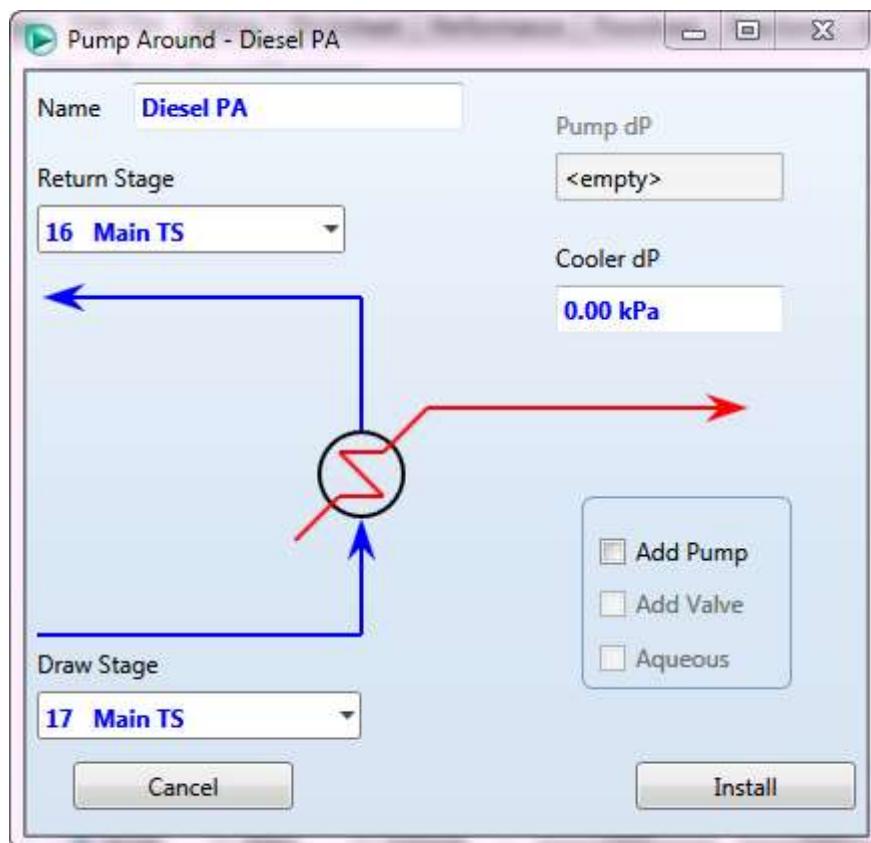
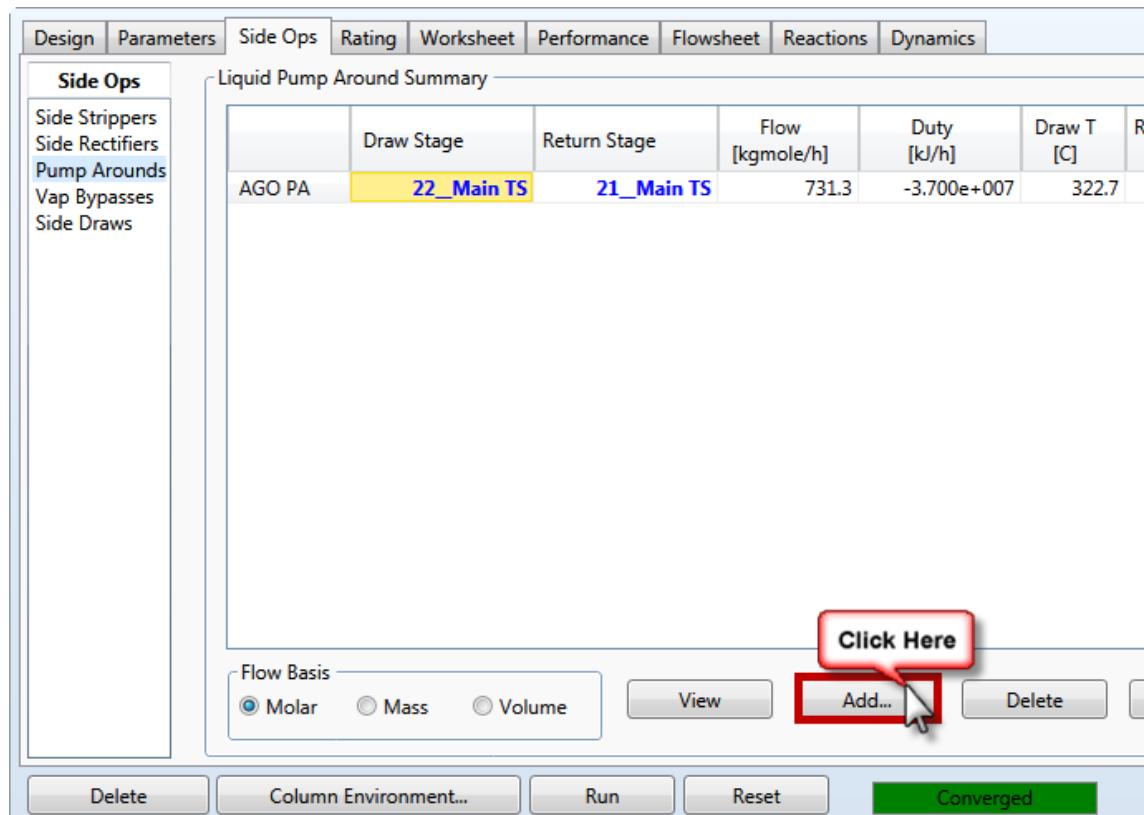


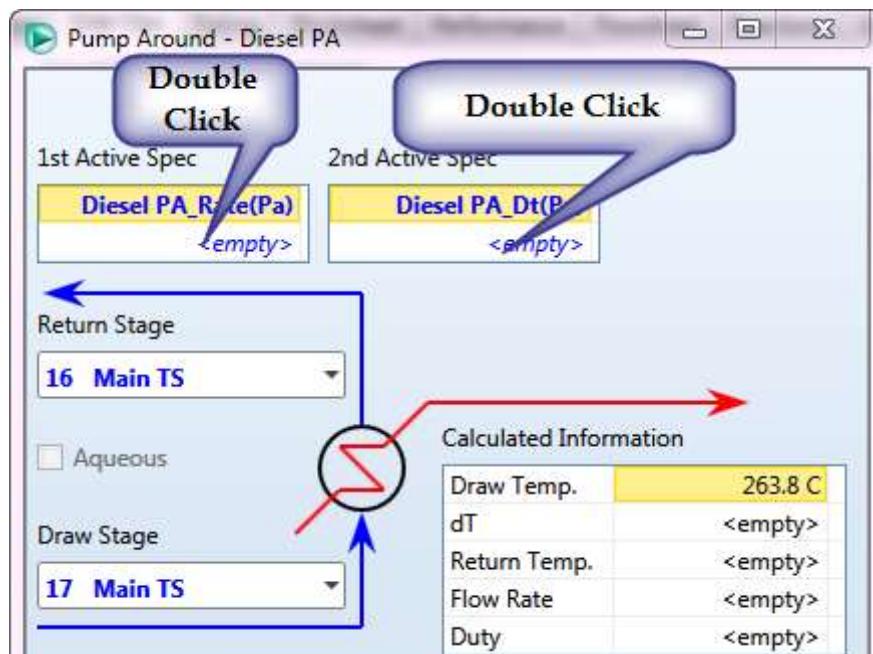
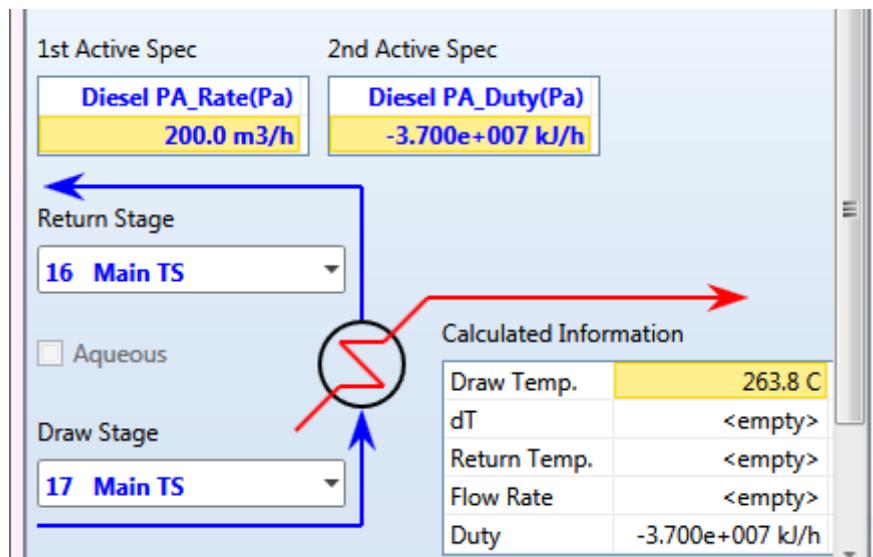


**Close the window and run the column**

**Make sure that the column is converged**

## Add the Diesel Pump Around



**Close the window and run the column  
Make sure that the column is converged**

## Add the Kerosene Pump Around

Liquid Pump Around Summary

	Draw Stage	Return Stage	Flow [kgmole/h]	Duty [kJ/h]	Draw T [C]	Return T [C]	Export
AGO PA	<b>22_Main TS</b>	<b>21_Main TS</b>	<empty>	-3.700e+007	322.7	<empty>	<input type="checkbox"/>
Diesel PA	<b>17_Main TS</b>	<b>16_Main TS</b>	<empty>	-3.700e+007	263.8	<empty>	<input type="checkbox"/>

Click Here

Flow Basis:  Molar  Mass  Volume      View      Add...      Delete      Side Ops Input Expert...

**Kerosene PA**

Return Stage: **8\_Main TS**

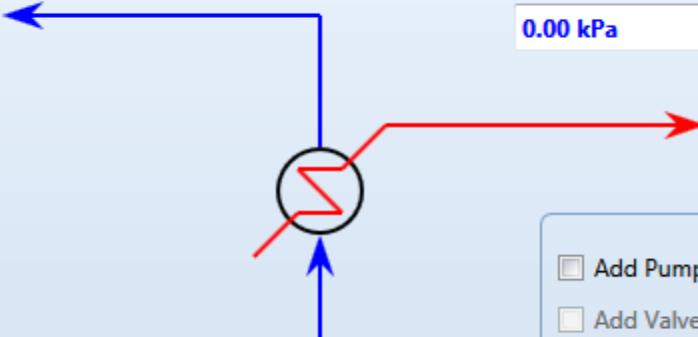
Pump dP: <empty>

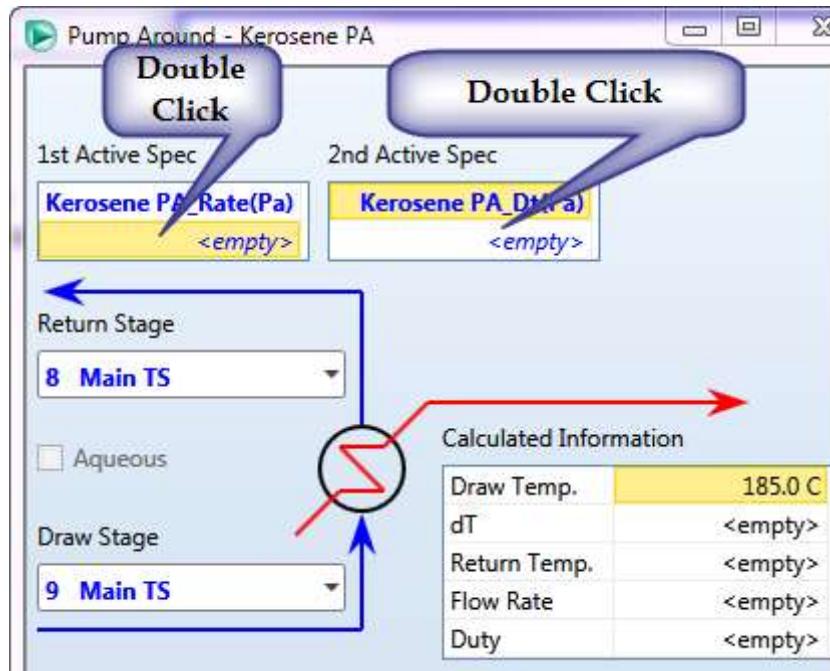
Cooler dP: 0.00 kPa

Draw Stage: **9\_Main TS**

Add Pump   
Add Valve   
Aqueous

Cancel      Install





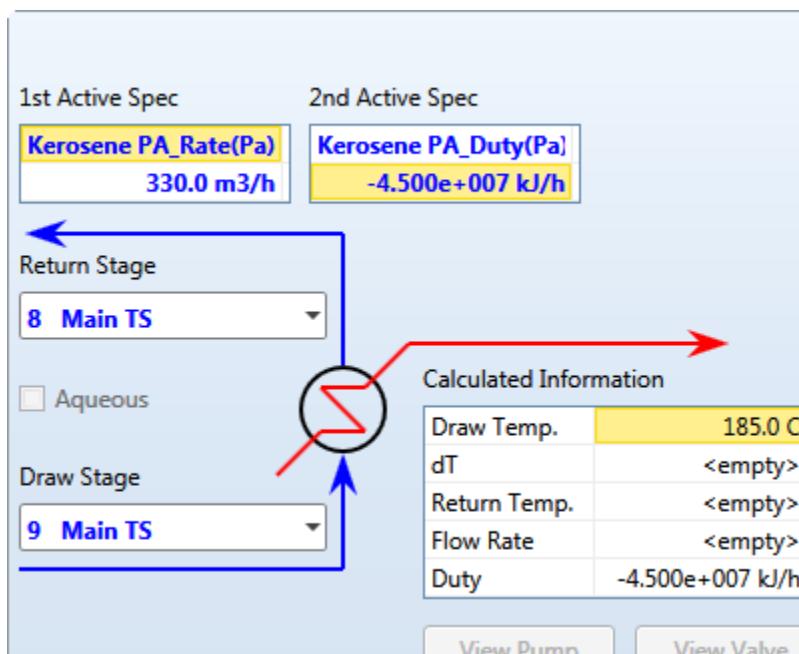
Two separate dialog boxes are shown, each detailing one of the active specifications from the main dialog.

**1st Active Spec (Kerosene PA\_Rate(Pa)):**

- Spec Type: Flow Rate
- Name: Kerosene PA\_Rate(Pa)
- Pump Around
- Flow Basis: Std Ideal Vol
- Spec Value: 330.0 m3/h

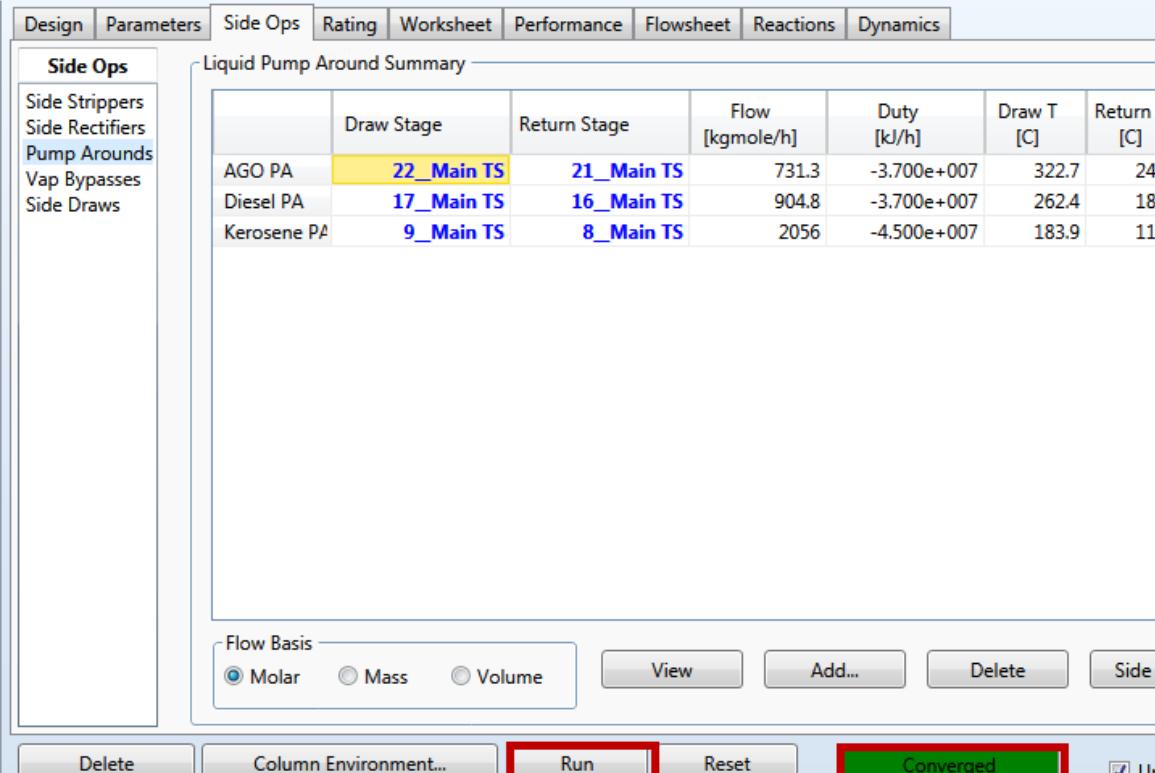
**2nd Active Spec (Kerosene PA\_Duty(Pa)):**

- Spec Type: Duty
- Name: Kerosene PA\_Duty(Pa)
- Pump Around
- Spec Value: -4.500e+007 kJ/h



**Close the window and run the column**

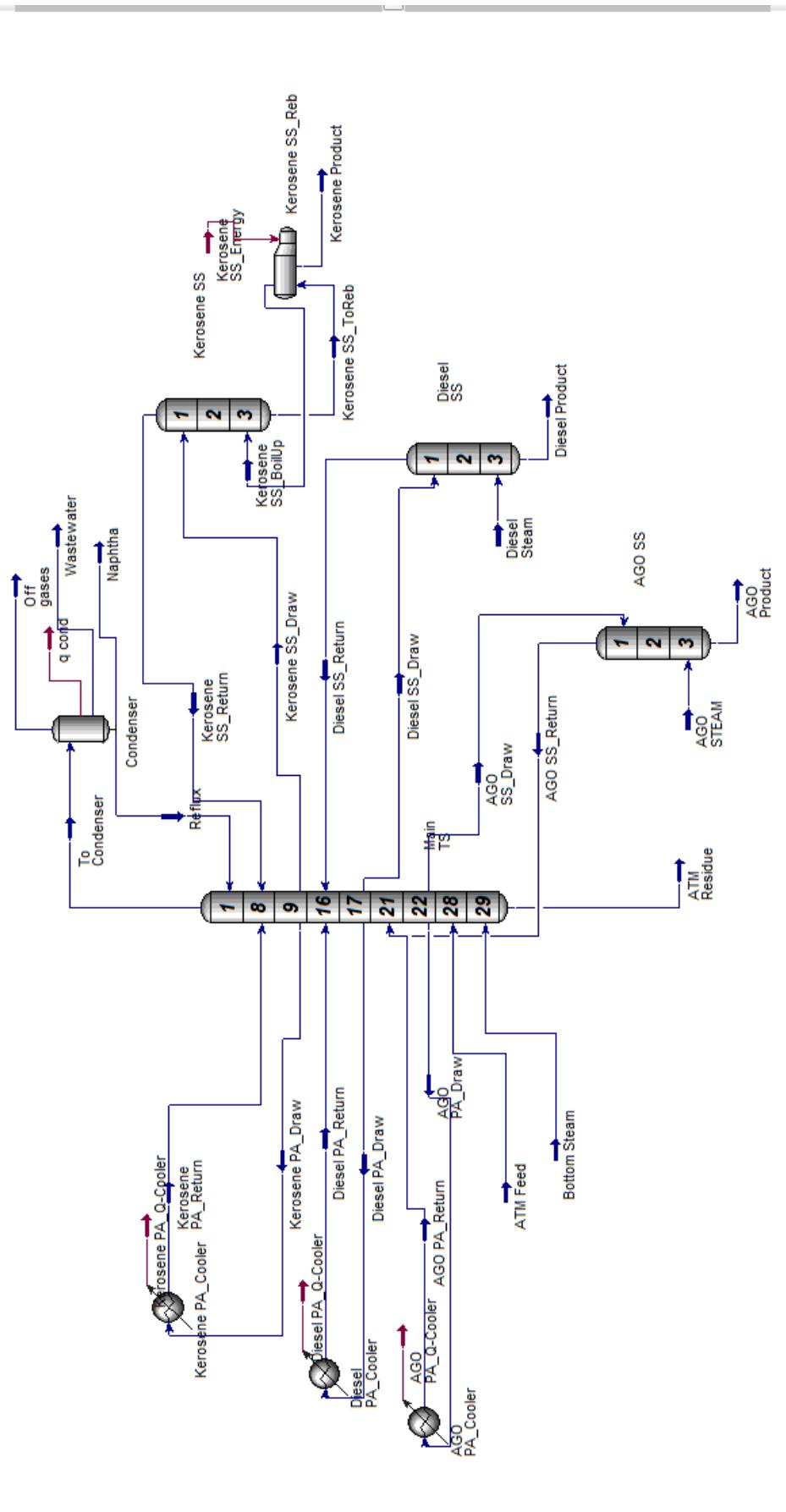
**Make sure that the column is converged**



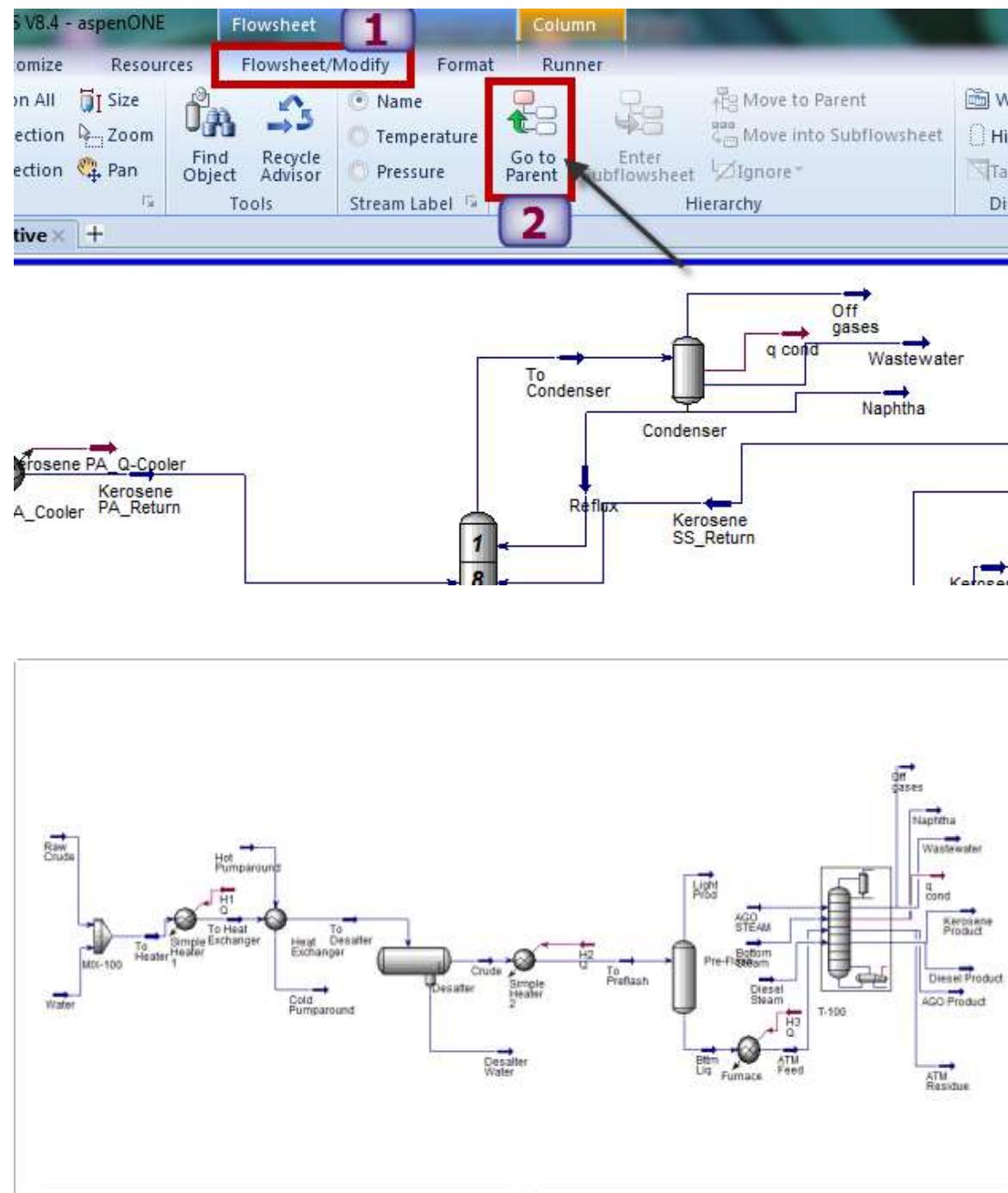
	Draw Stage	Return Stage	Flow [kgmole/h]	Duty [kJ/h]	Draw T [C]	Return T [C]
AGO PA	22_Main TS	21_Main TS	731.3	-3.700e+007	322.7	24
Diesel PA	17_Main TS	16_Main TS	904.8	-3.700e+007	262.4	18
Kerosene PA	9_Main TS	8_Main TS	2056	-4.500e+007	183.9	11

You can find the detailed flow sheet by entering the *Column Environment*



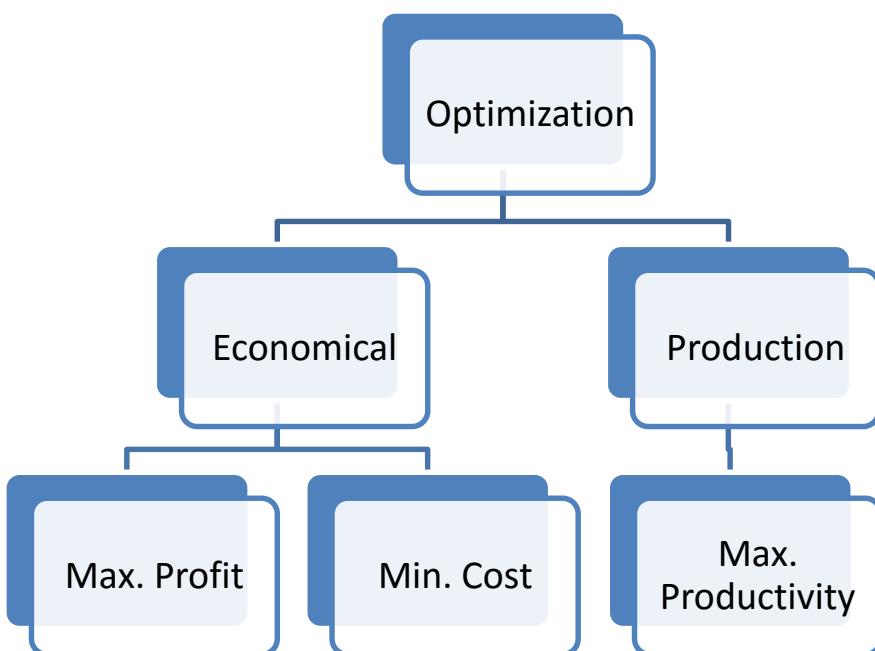


To return to the main environment:



*Save Your Case!*

# Optimization



7

## ***Workshop***

In this example, a simple distillation column to separate Tetrahydrofuran (THF) from Toluene is simulated. The object of the exercise is to select the product specifications such that profit is maximized. A special tool in HYSYS, the Optimizer, will be used to find the optimum operating conditions.

HYSYS includes additional modelling and decision support tools that can be used to enhance the usability of your models. In this module, you will use the HYSYS optimization tool available in HYSYS to investigate the debottlenecking and optimization of a crude column.

## ***Learning Objectives***

Once you have completed this section, you will be able to:

- Use the Optimizer tool in HYSYS to optimize flowsheets
- Use the Spreadsheet to perform calculations

Example:

3700 kg/hr mixture of tetrahydrofuran & toluene (44 mass% THF) at 10°C and 140 kPa is to be separated by distillation to get each of them with purity of 99.5 mass% of THF & 94 mass% of Toluene (THF is the more volatile component).

- Use Wilson fluid package

**The column specifications are:**

- The condenser & reboiler pressure are 103 kPa & 107 kPa.
- The condenser works on **total condensation** conditions.
- **Number of stages = 10.**
- Feed enters from the 5<sup>th</sup> tray.

Calculate:

The reflux ratio and the distillate rate under the specified conditions.

<b>Reflux Ratio</b>	.....
<b>Distillate Rate</b>	..... kgmol/hr

Data:

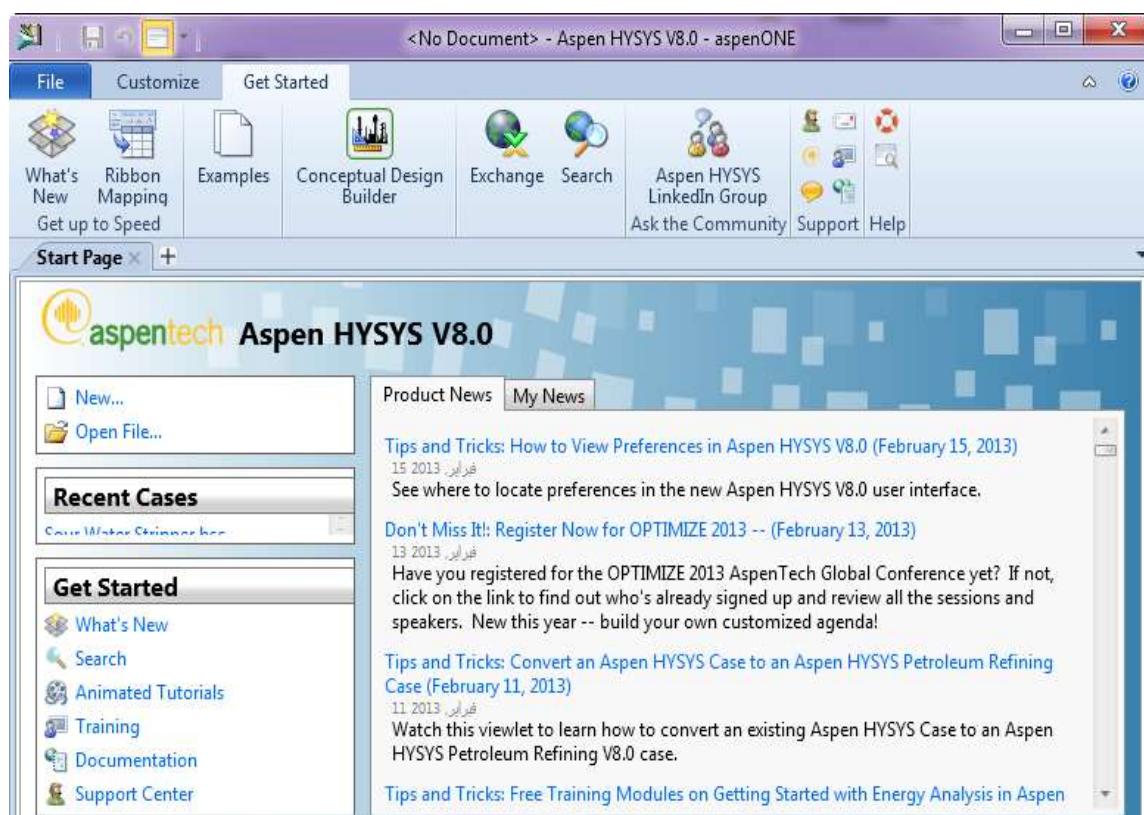
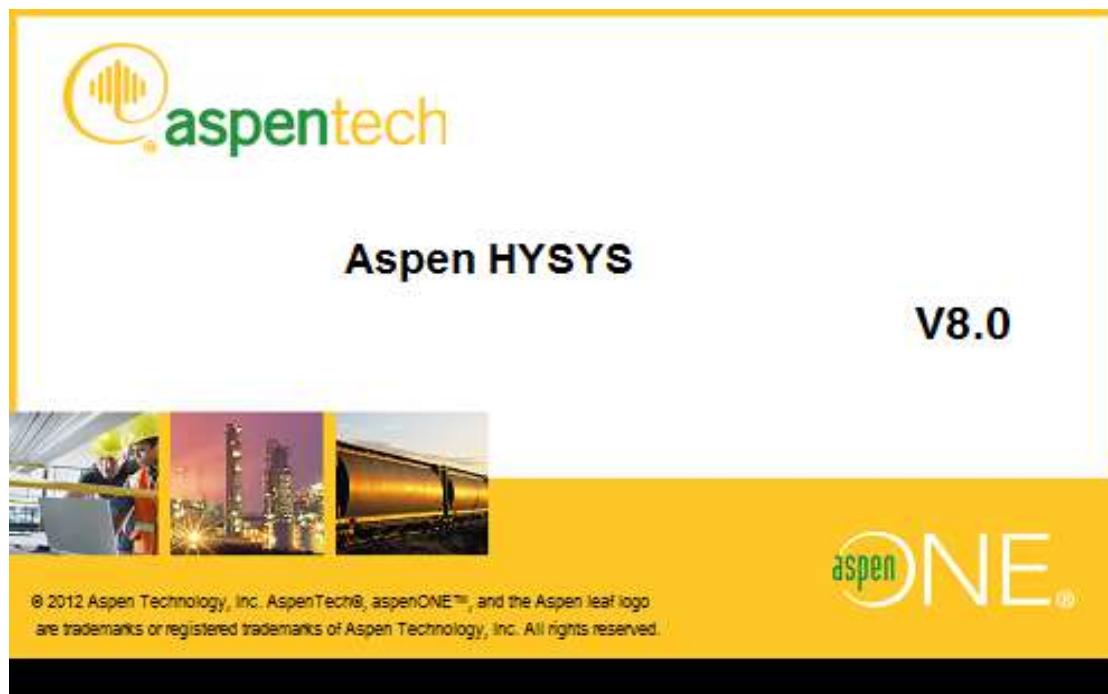
- Feed price= 0.05 \$/kg.
- Pure toluene selling price= 0.136 \$/kg
- Pure THF selling price= 0.333 \$/ kg
- Cooling Cost= 0.471 \$/ kw.hr
- Heating Cost= 0.737 \$/kw.hr

Note:

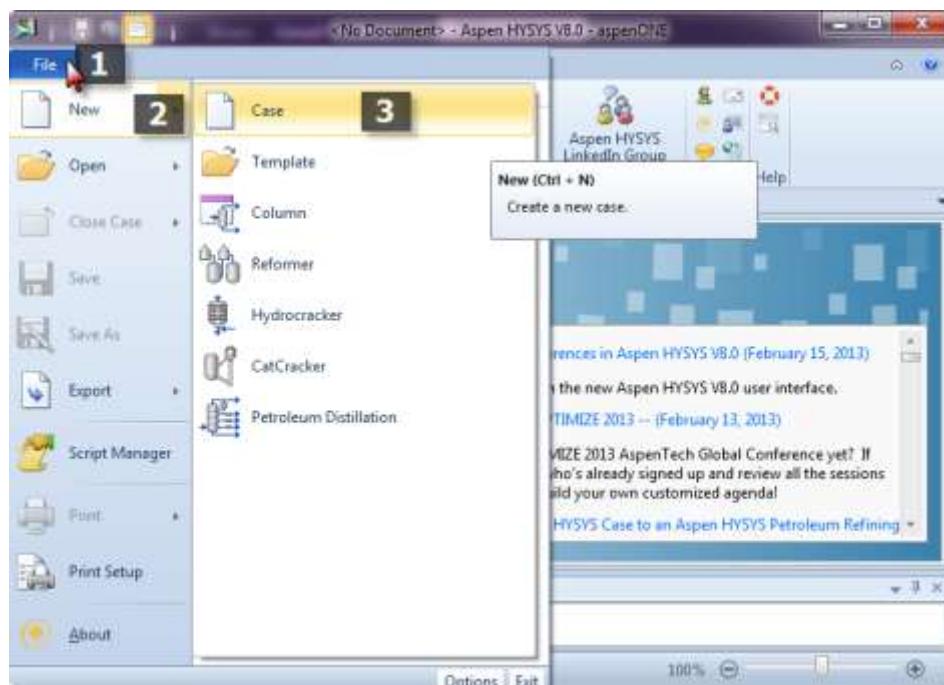
Profit = (Total Toluene selling price + Total THF selling price) - (Feed cost + Heating cost + Cooling Cost)

- Use a range of 0.99 to 0.999 for THF limit & 0.9 to 0.99 for the toluene.

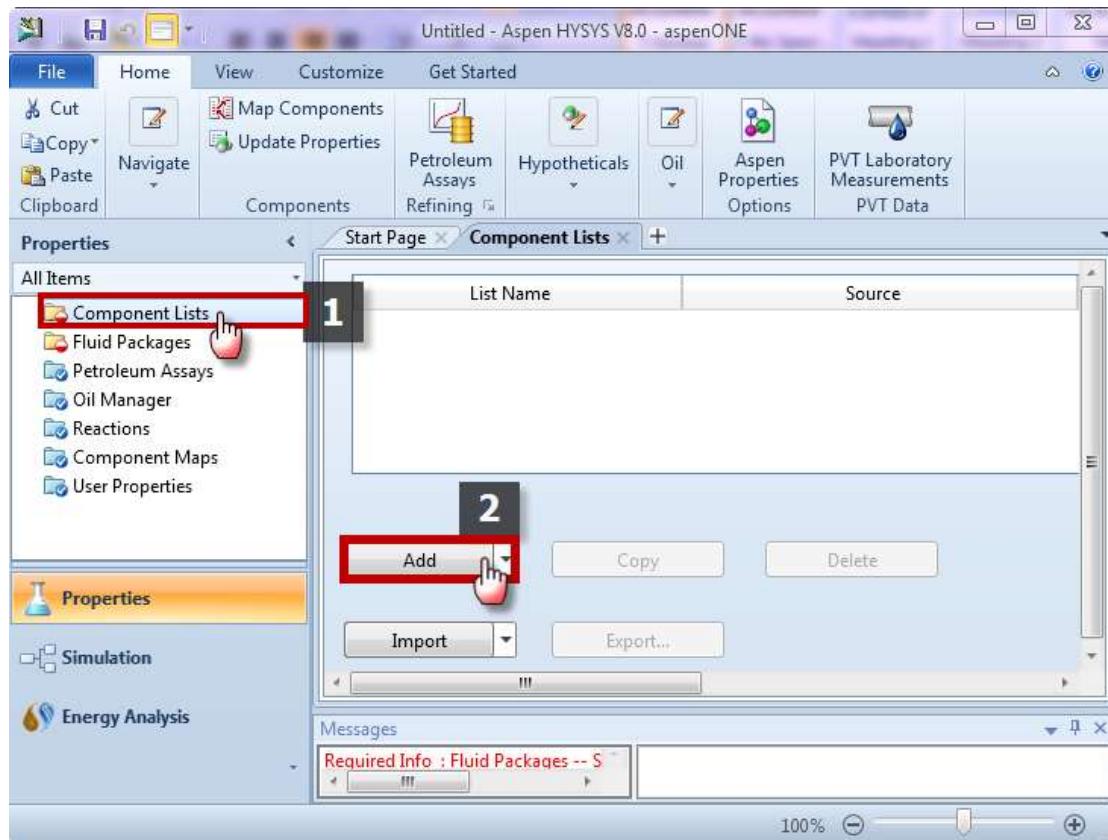
To start the program, From Start Menu, Select All Programs >>  
Aspen Tech >> Process Modeling V8.0 >>> Aspen HYSYS >>  
Aspen HYSYS



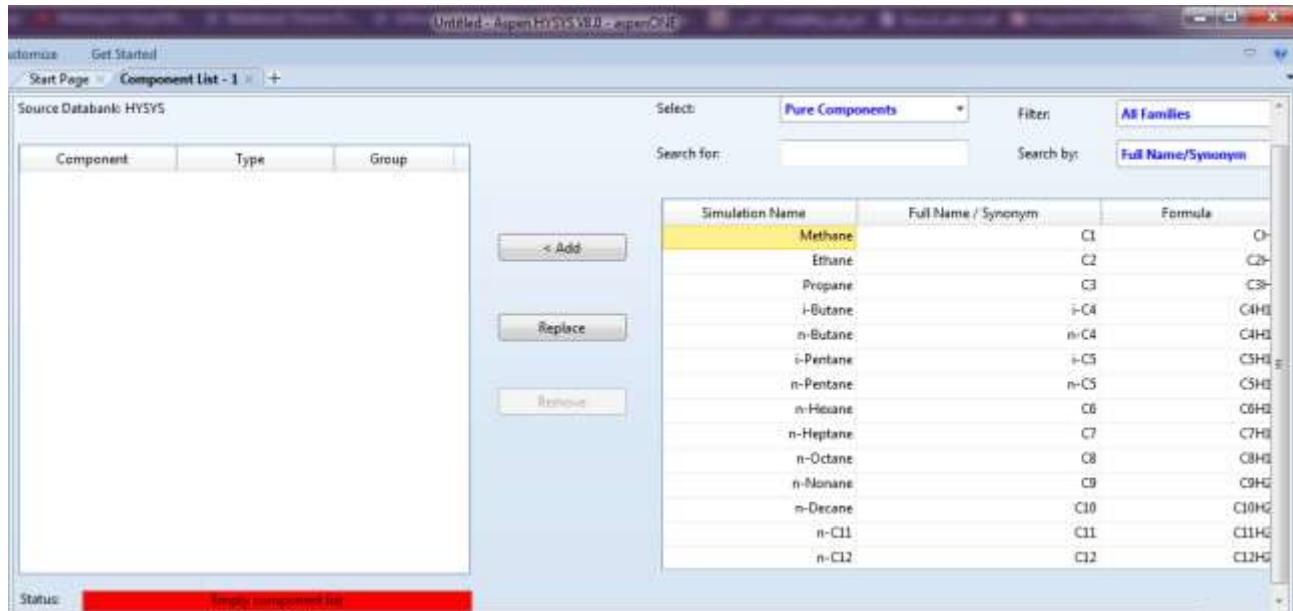
## 5- First, Start a new case



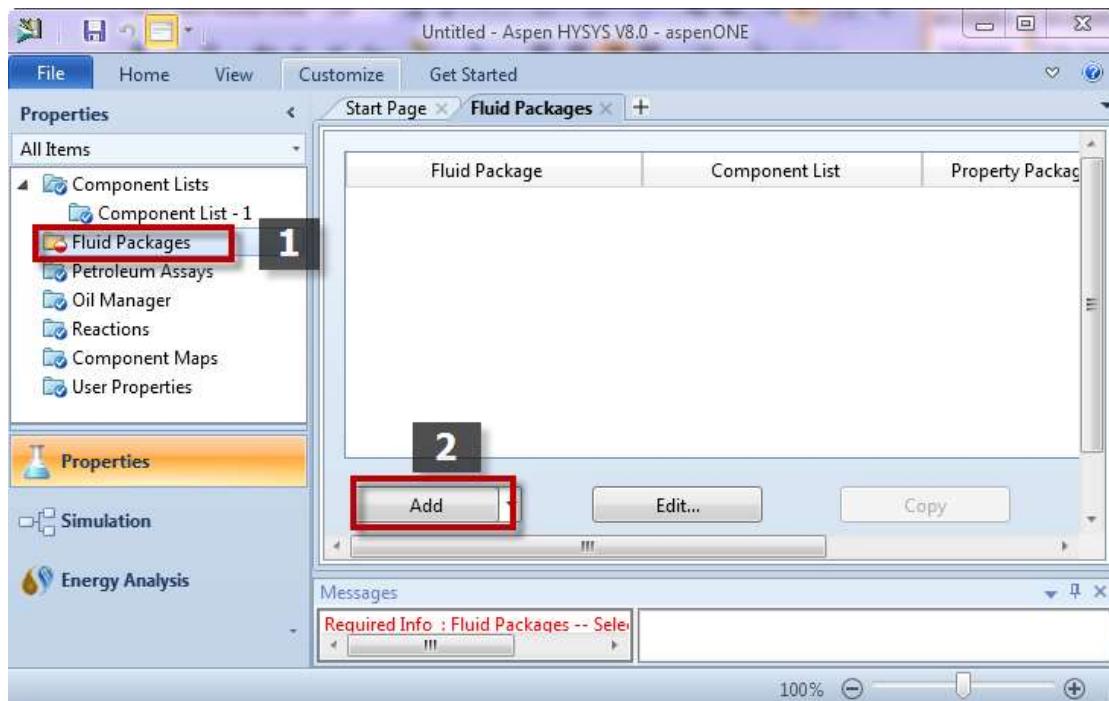
## 6- Add the Components



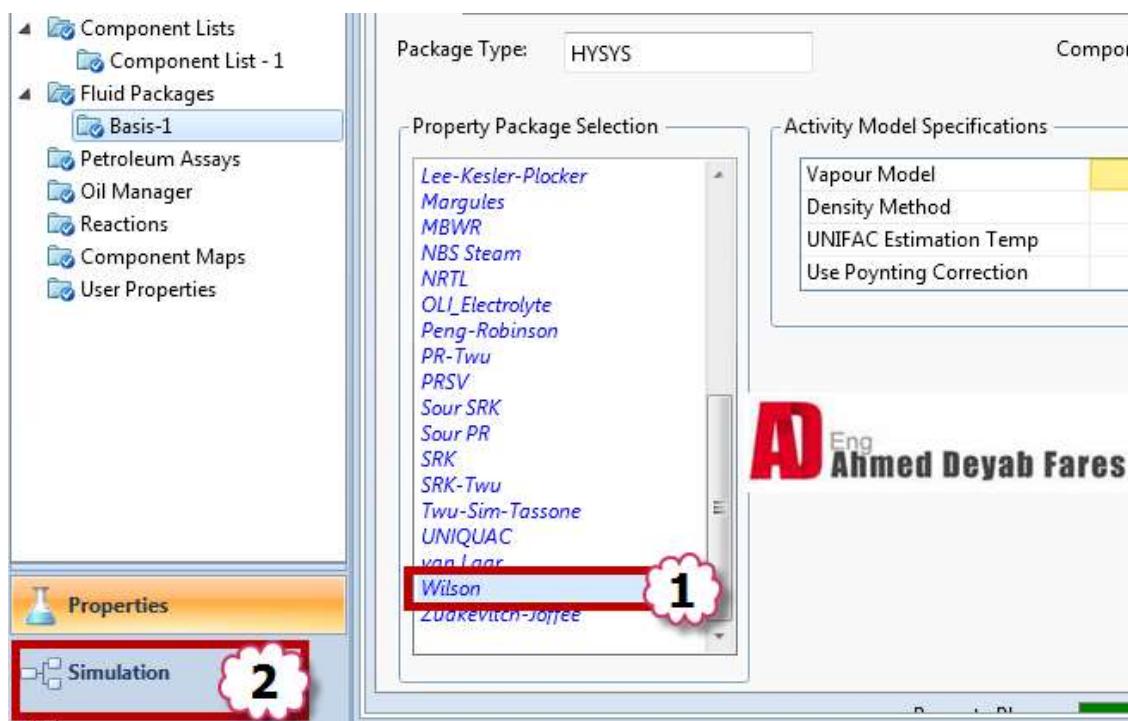
## 7- Choose the system components from the databank:



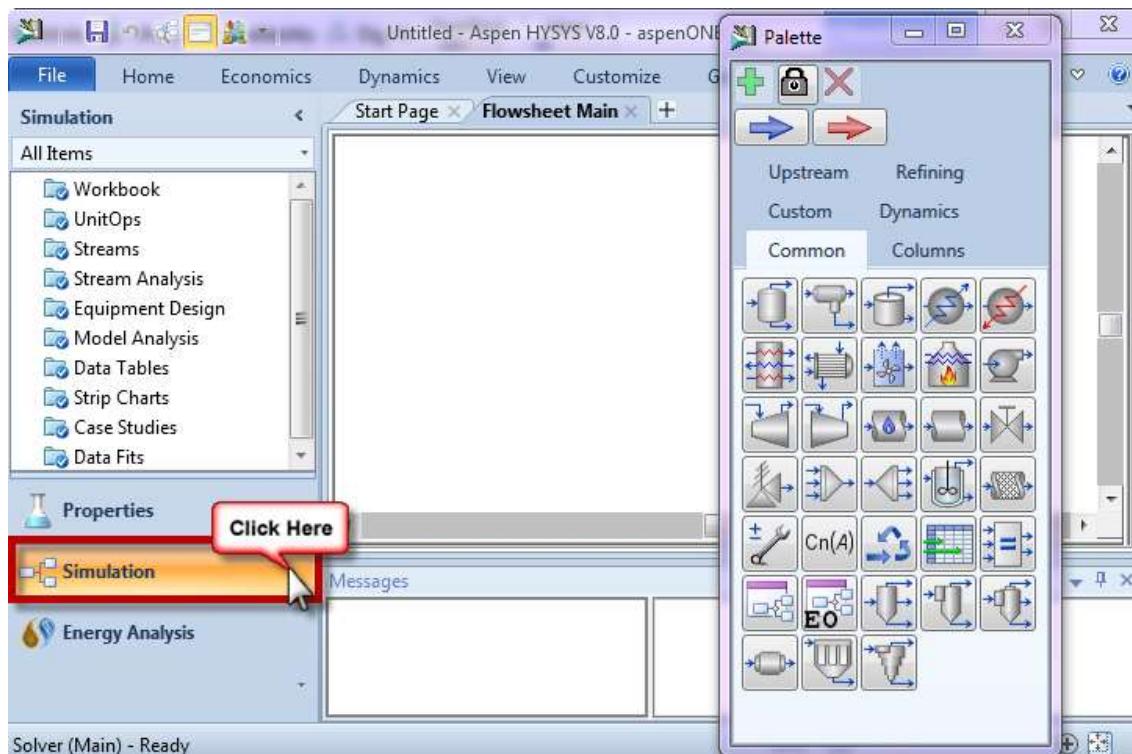
Now, select the suitable fluid package



In this case, select Wilson

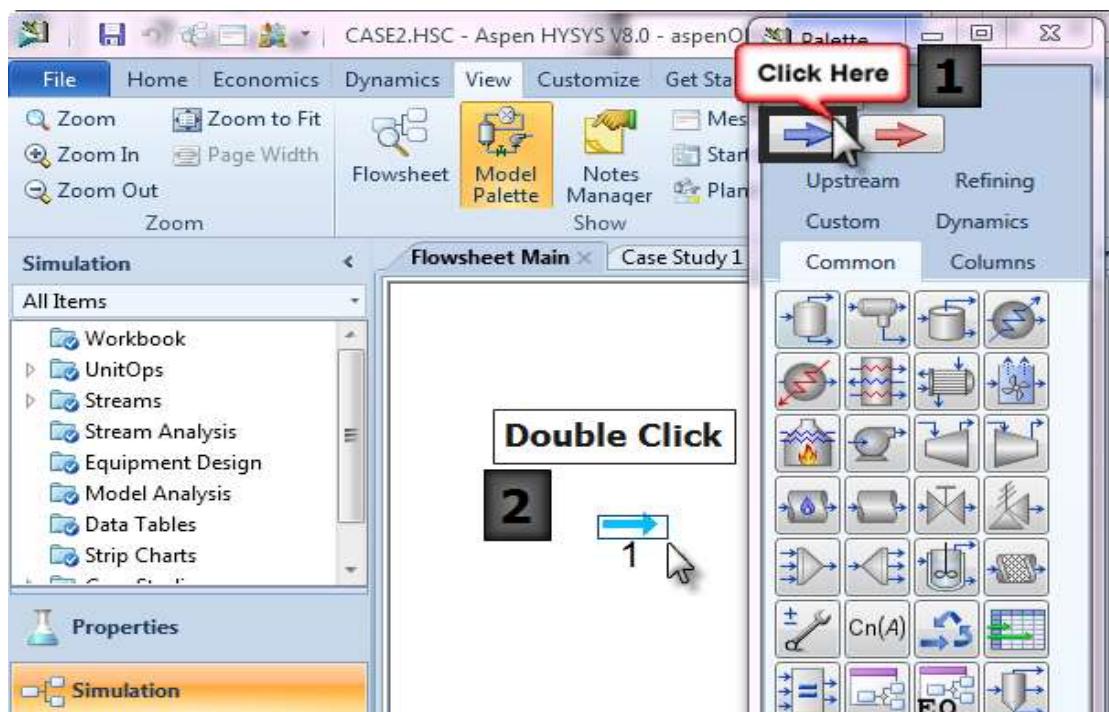


Now you can start drawing the flow sheet for the process by clicking the Simulation button:

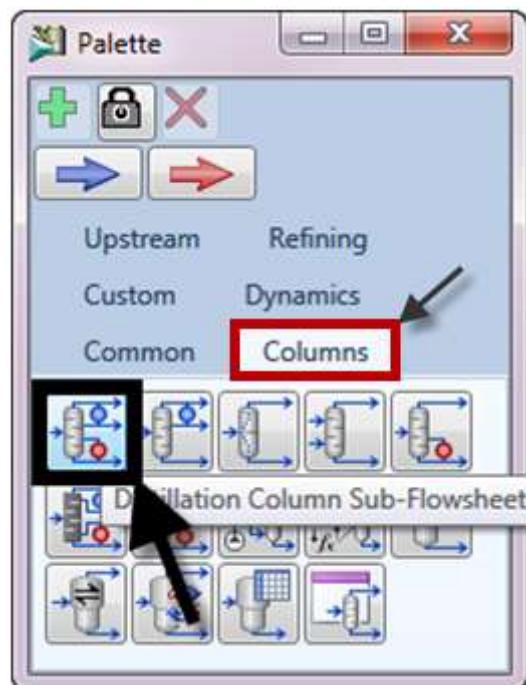


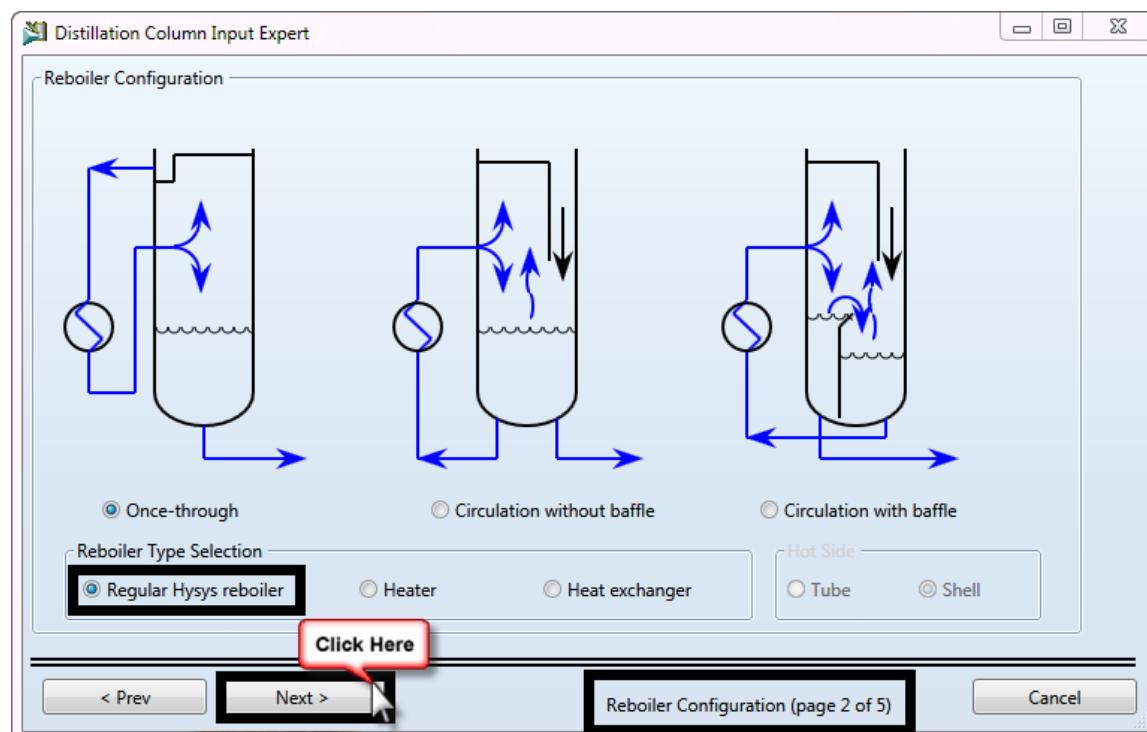
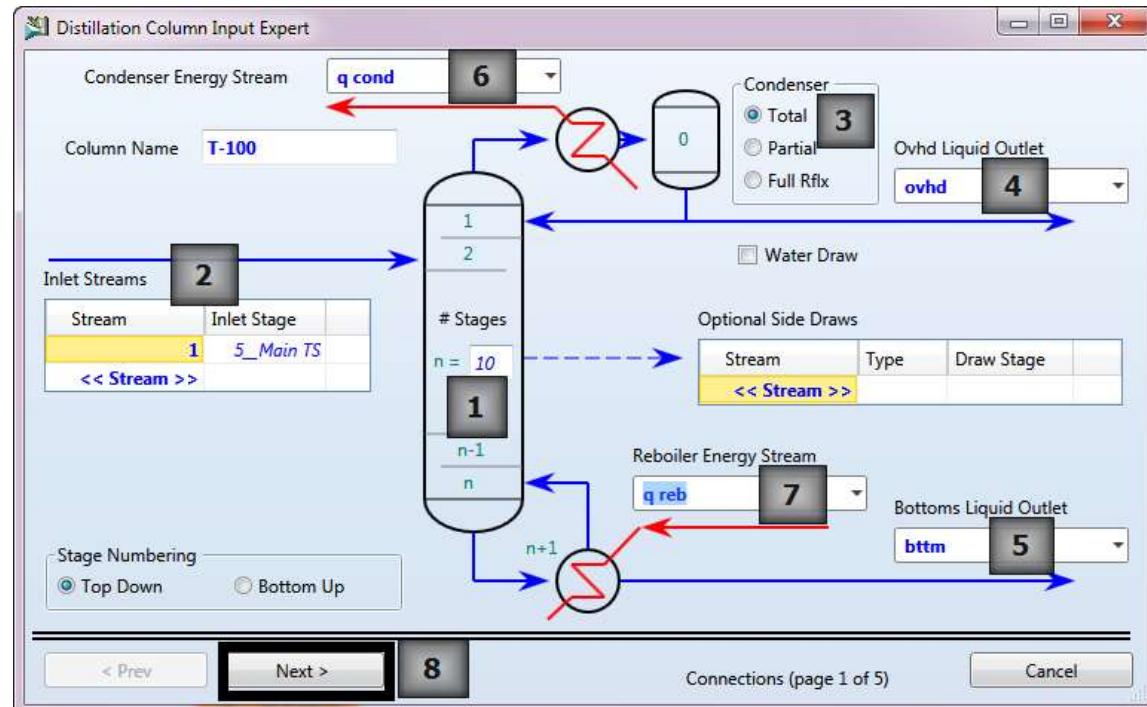
Now add a material stream to define the composition and the conditions of the feed stream

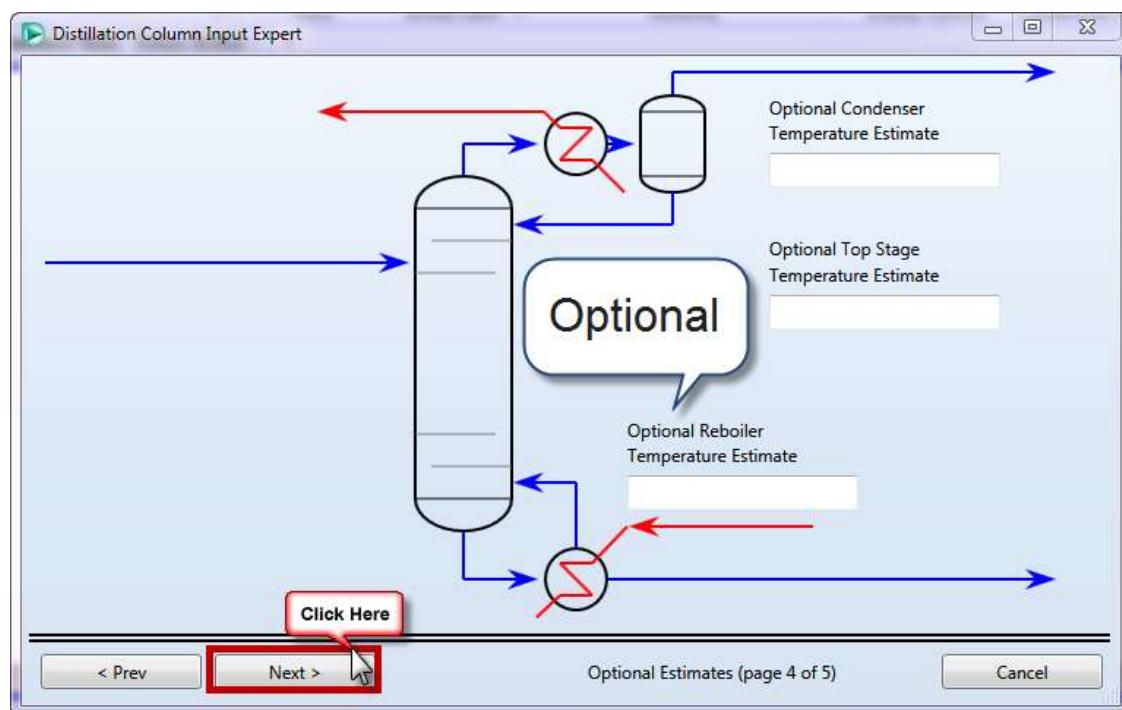
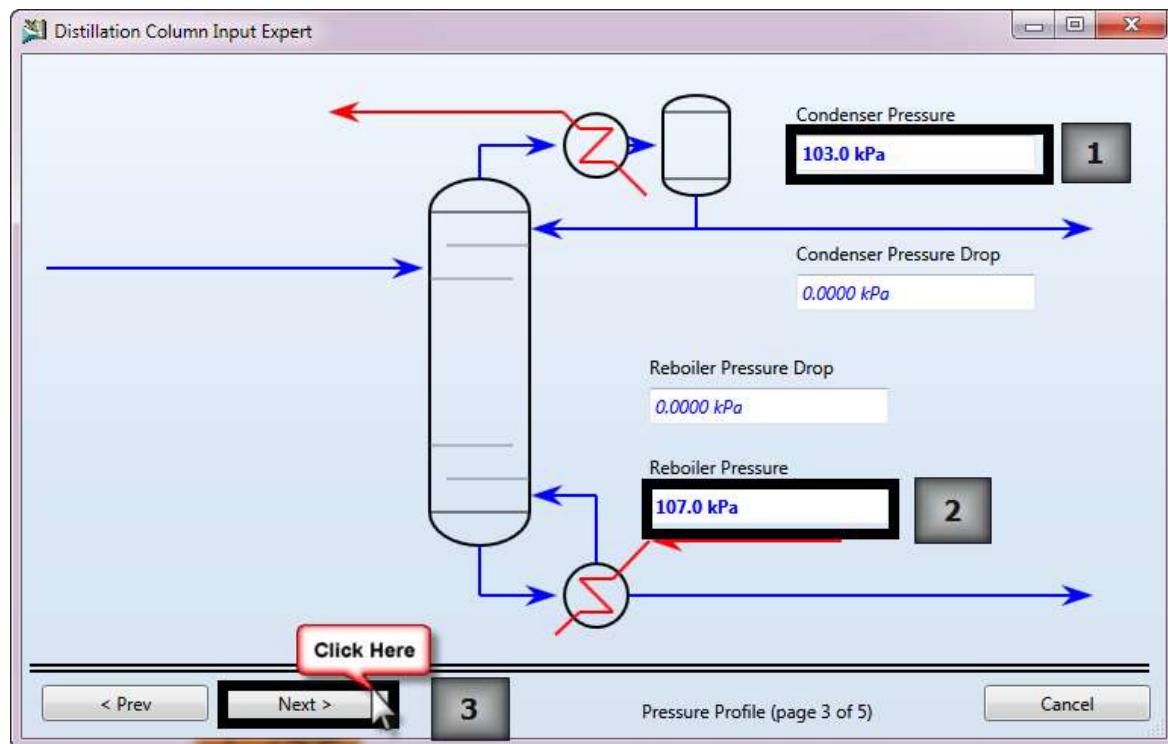
From the palette:

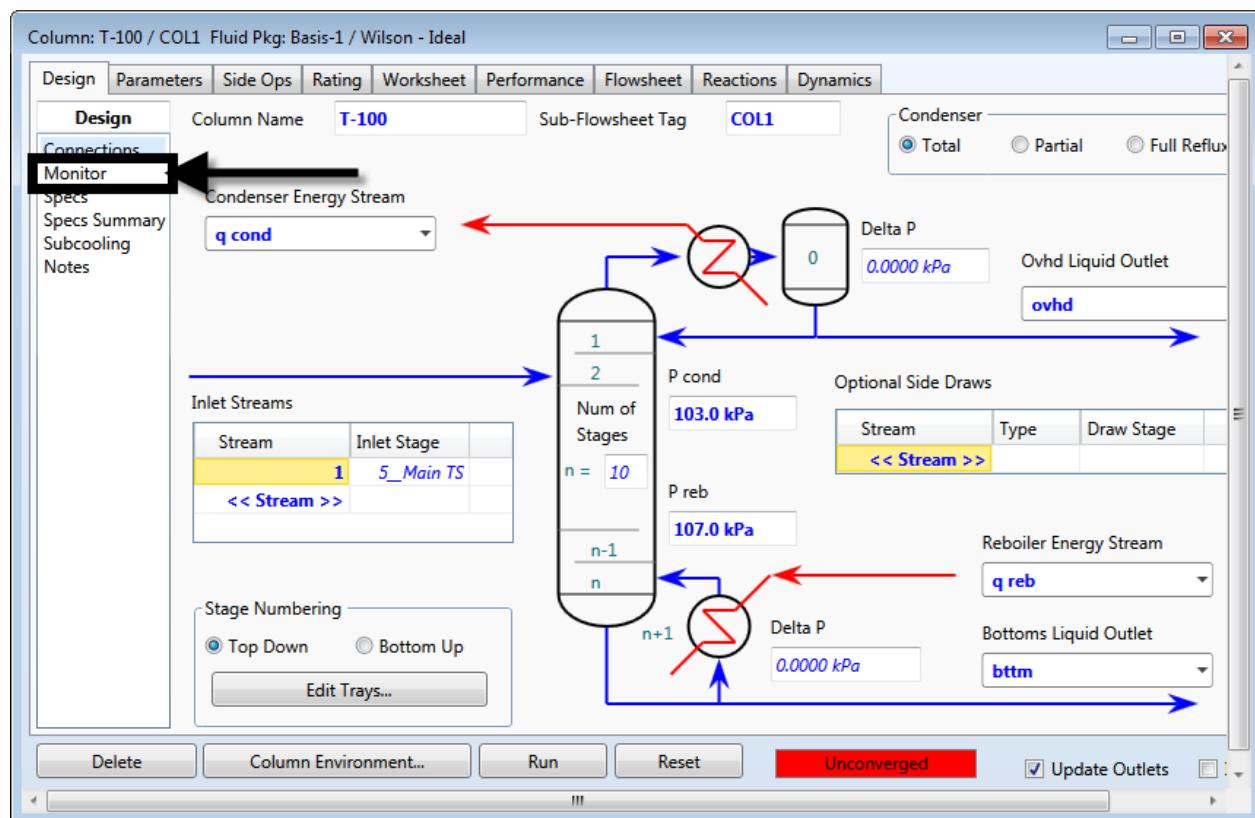
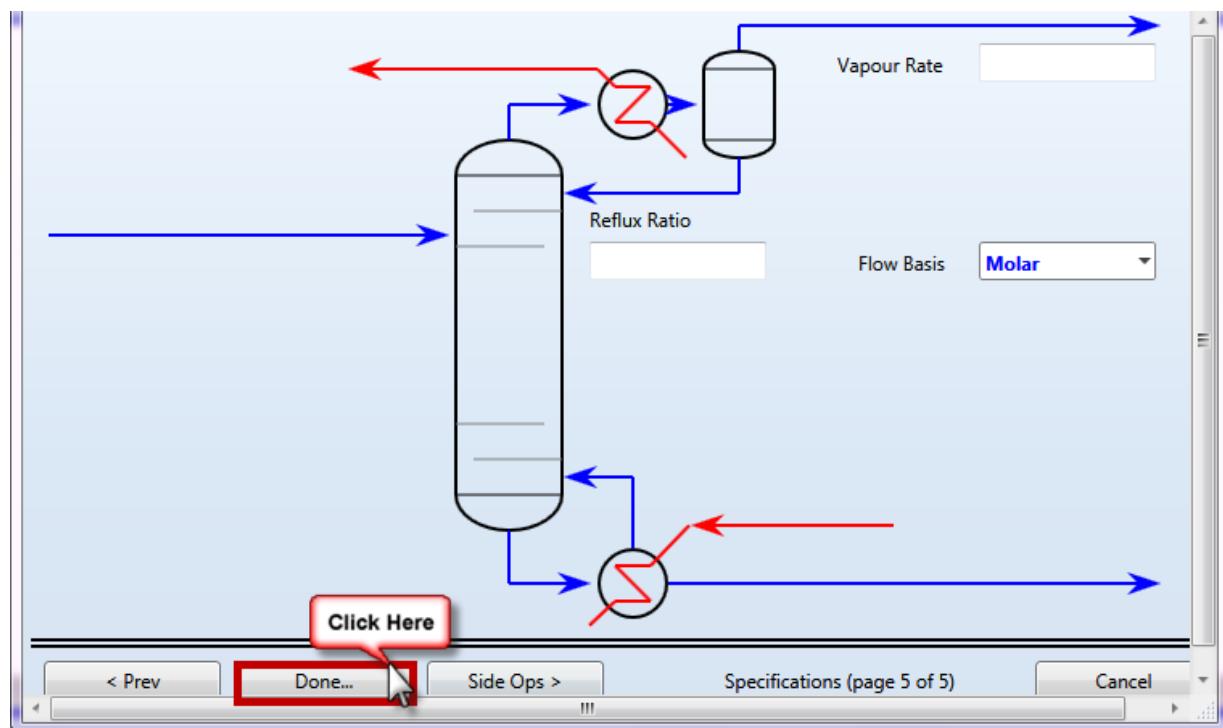


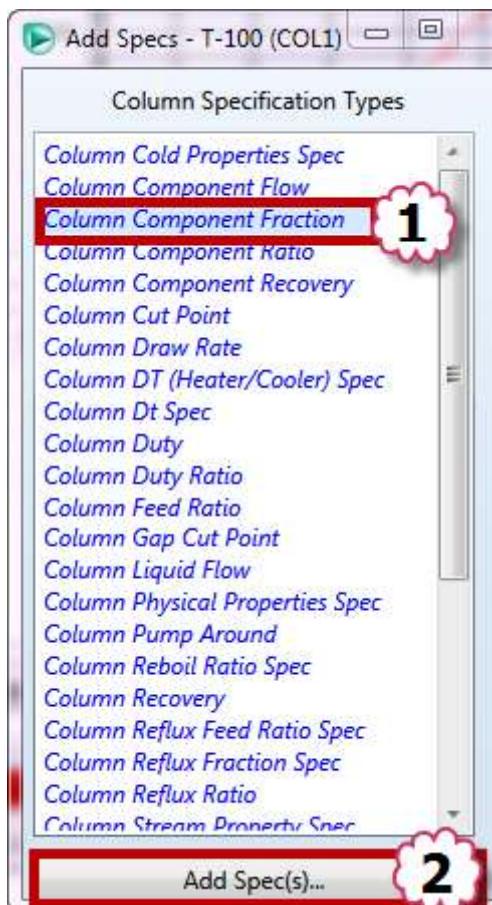
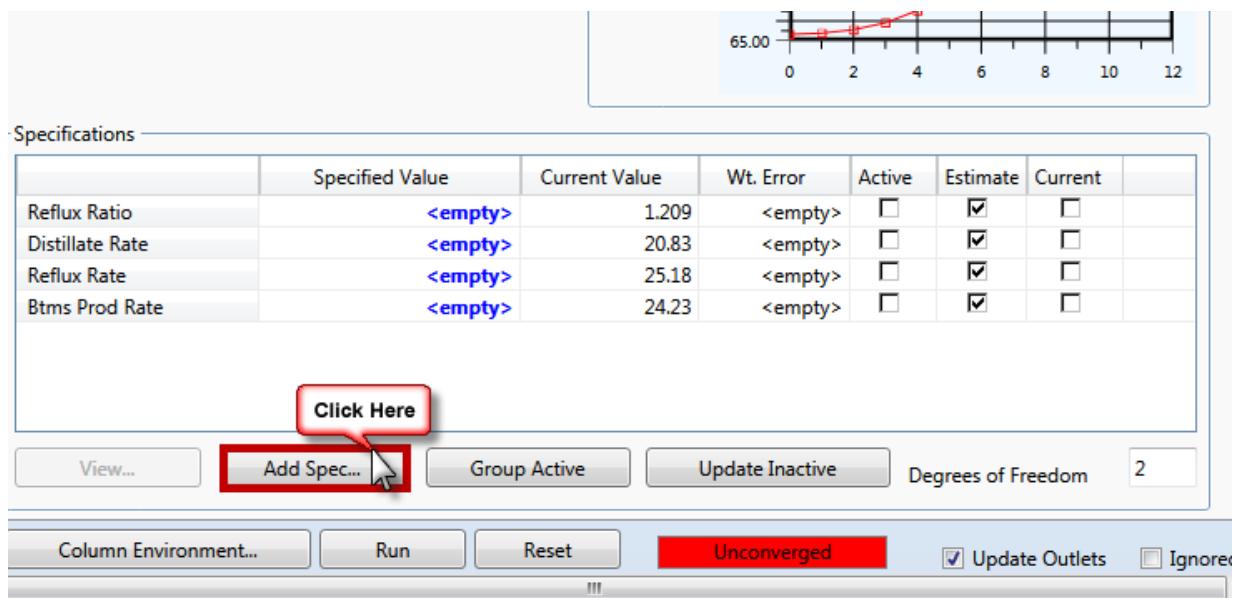
Add a distillation tower:











The image shows two separate windows from the HYSYS software interface. Both windows are titled "Comp Frac Spec".

**Left Window (THF Purity):**

- Parameters:** Name, Stage, Flow Basis, Phase, Spec Value.
- Spec Type:** THF Purity, Condenser, Mass Fraction, Liquid, Spec Value: 0.9950.
- Components:** TetraHyFuran, << Component >>.
- Target Type:** Stream (radio button).
- Buttons:** Delete.

**Right Window (Toluene purity):**

- Parameters:** Name, Stage, Flow Basis, Phase, Spec Value.
- Spec Type:** Toluene purity, Reboiler, Mass Fraction, Liquid, Spec Value: 0.9400.
- Components:** Toluene, << Component >>.
- Target Type:** Stage (radio button).

**A Eng Ahmed Deyab Fares** watermark is visible at the bottom center of the interface.

**Specifications** window:

	Specified Value	Current Value	Wt. Error	Active	Estimate	Current
Reflux Ratio	<empty>	1.209	<empty>	<input type="checkbox"/>	<input checked="" type="checkbox"/>	<input type="checkbox"/>
Distillate Rate	<empty>	20.83	<empty>	<input type="checkbox"/>	<input checked="" type="checkbox"/>	<input type="checkbox"/>
Reflux Rate	<empty>	25.18	<empty>	<input type="checkbox"/>	<input checked="" type="checkbox"/>	<input type="checkbox"/>
Btms Prod Rate	<empty>	24.23	<empty>	<input type="checkbox"/>	<input checked="" type="checkbox"/>	<input type="checkbox"/>
THF Purity	0.9950	0.9950	-0.0001	<input checked="" type="checkbox"/>	<input checked="" type="checkbox"/>	<input type="checkbox"/>
Toluene purity	0.9400	0.9400	-0.0005	<input checked="" type="checkbox"/>	<input checked="" type="checkbox"/>	<input type="checkbox"/>

**Buttons:** View..., Add Spec..., Group Active, Update Inactive, Degrees of Freedom: 0.

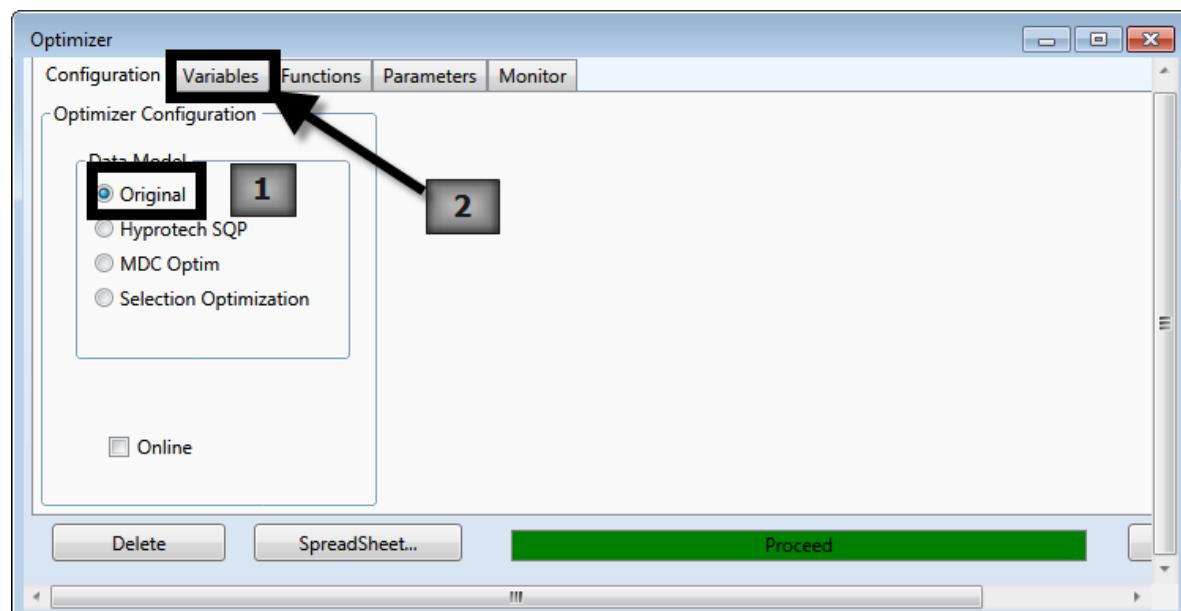
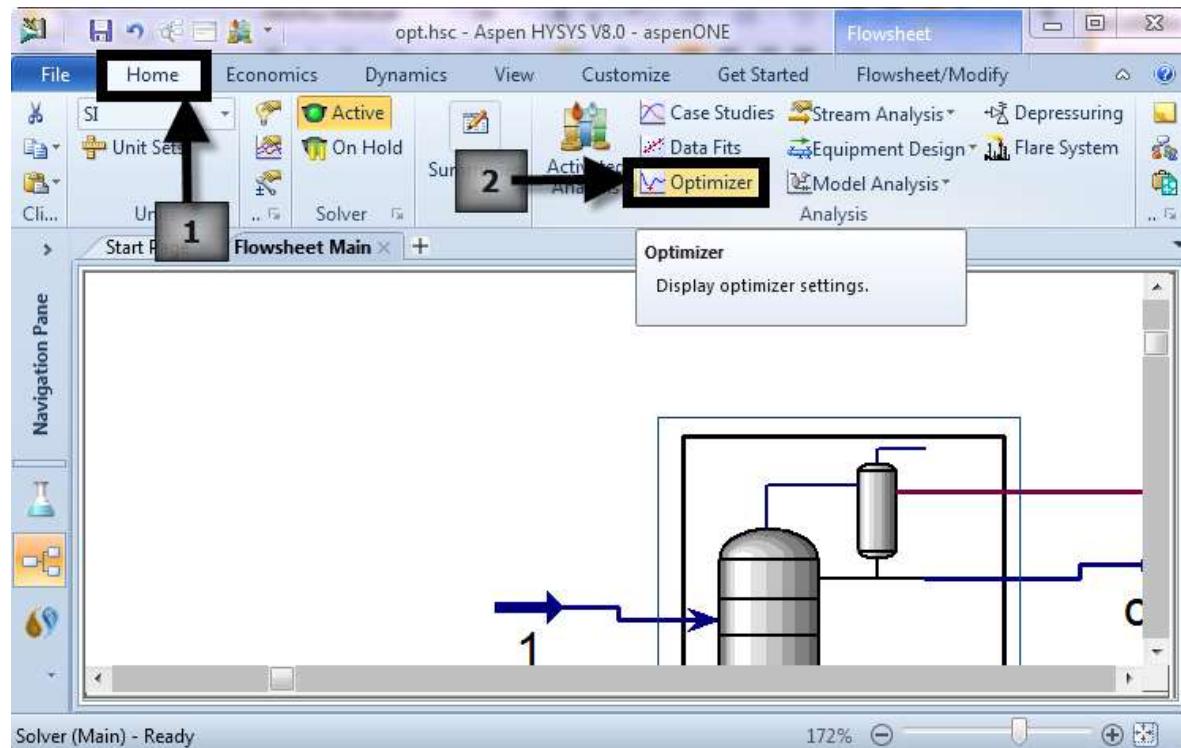
**Control Buttons:** Column Environment (button 2), Run (button highlighted in red), Reset, Unconverged, Update Outlets, Ign.

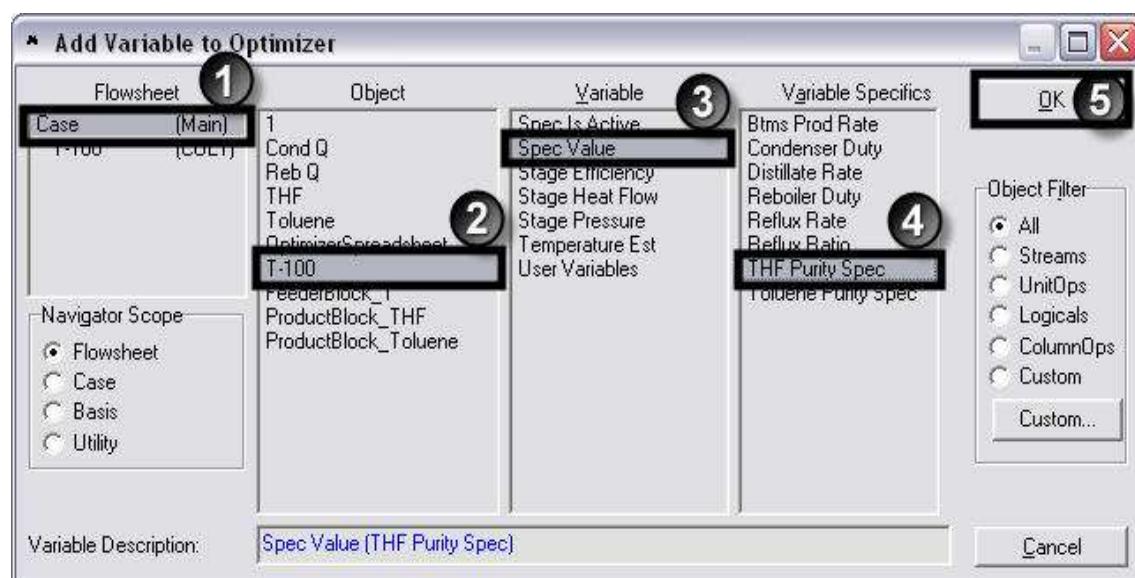
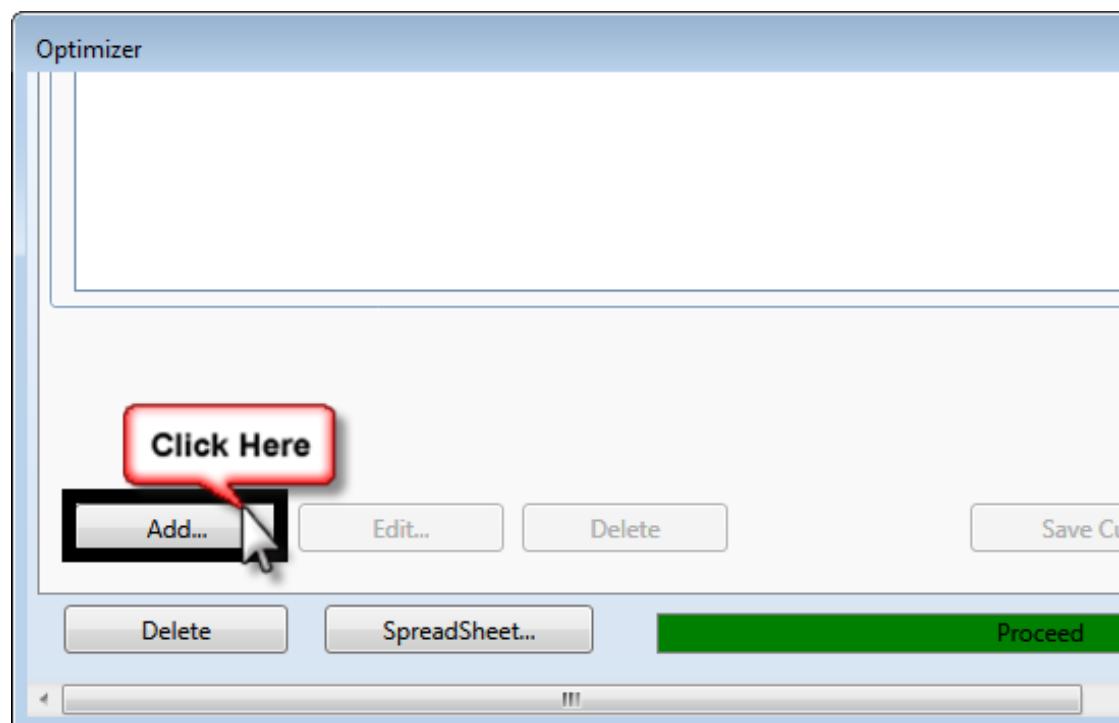
# Optimization

We need to check if the operating conditions are optimum or not.

The Variables to check are: THF Purity & Toluene Purity

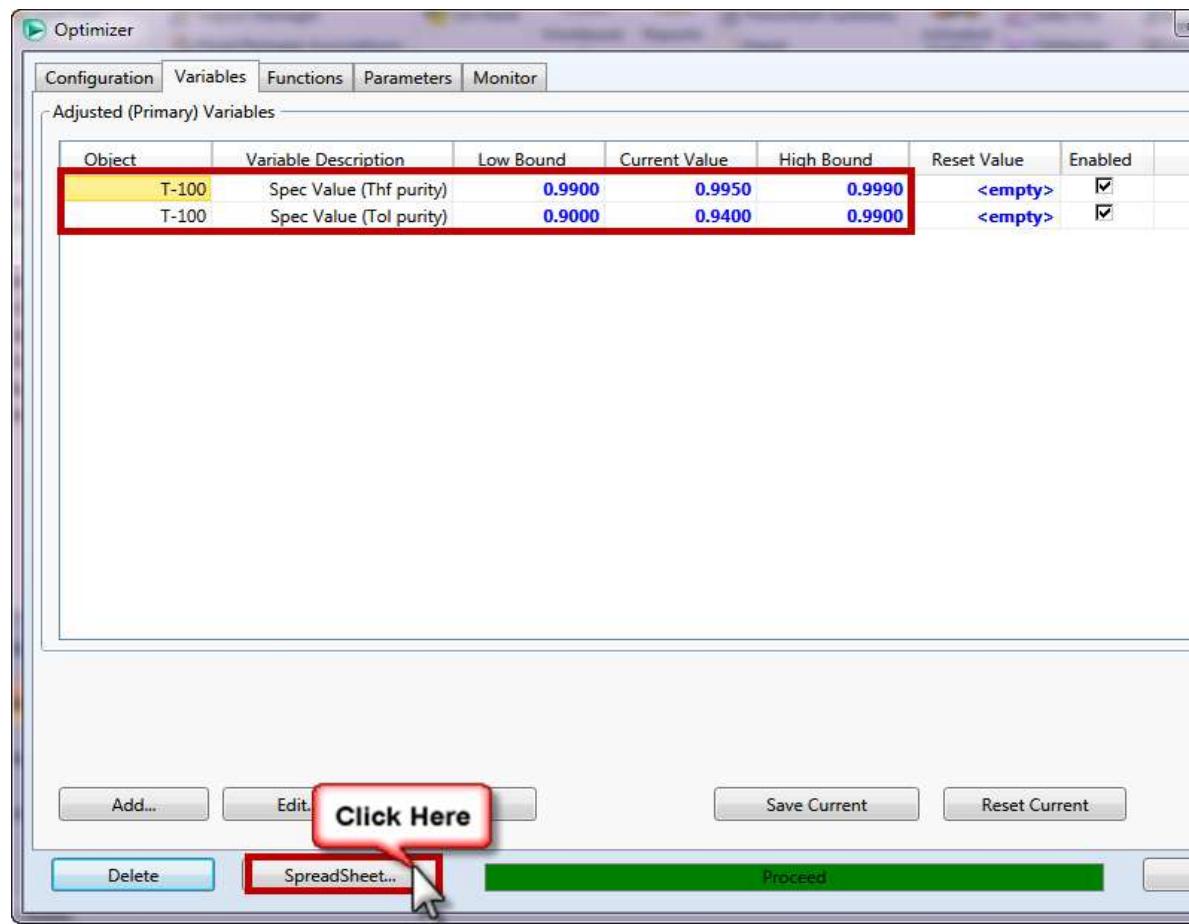
Using the Optimizer tool:



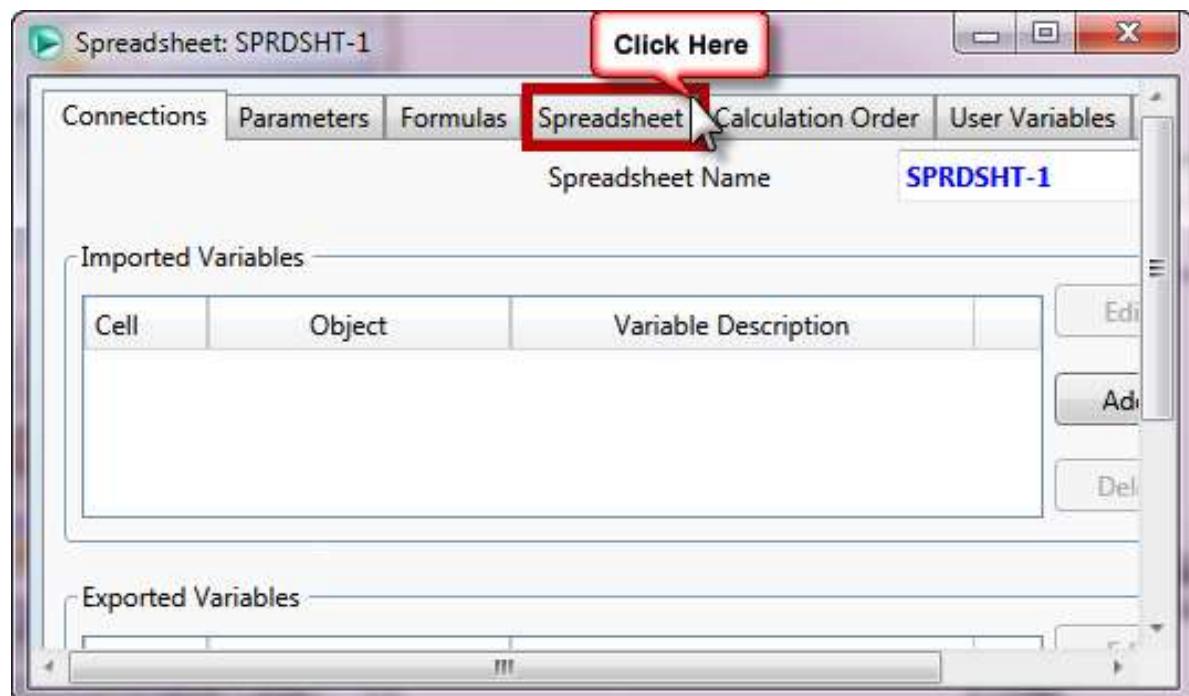


And do the same steps to add the other variable

Use a range of 0.99 to 0.999 for THF limit & 0.9 to 0.99 for the toluene.

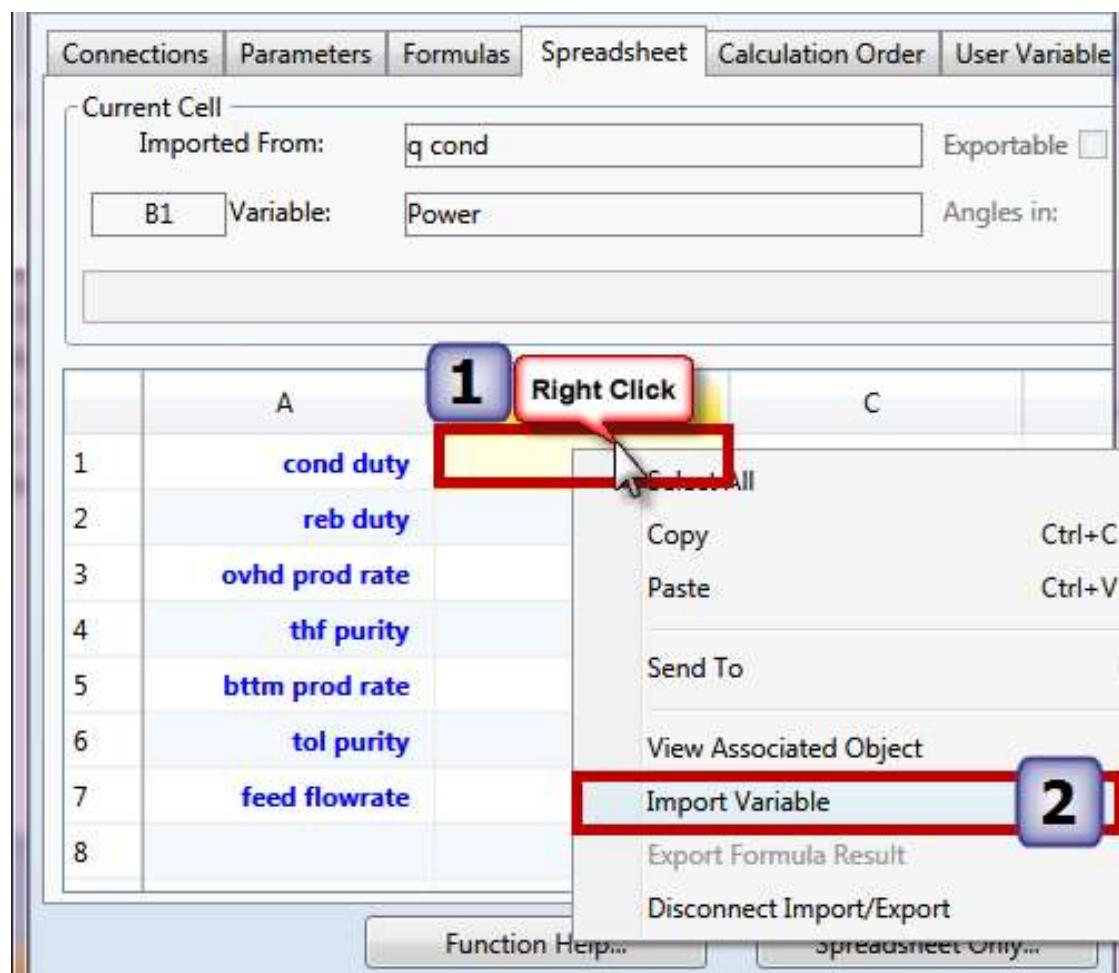
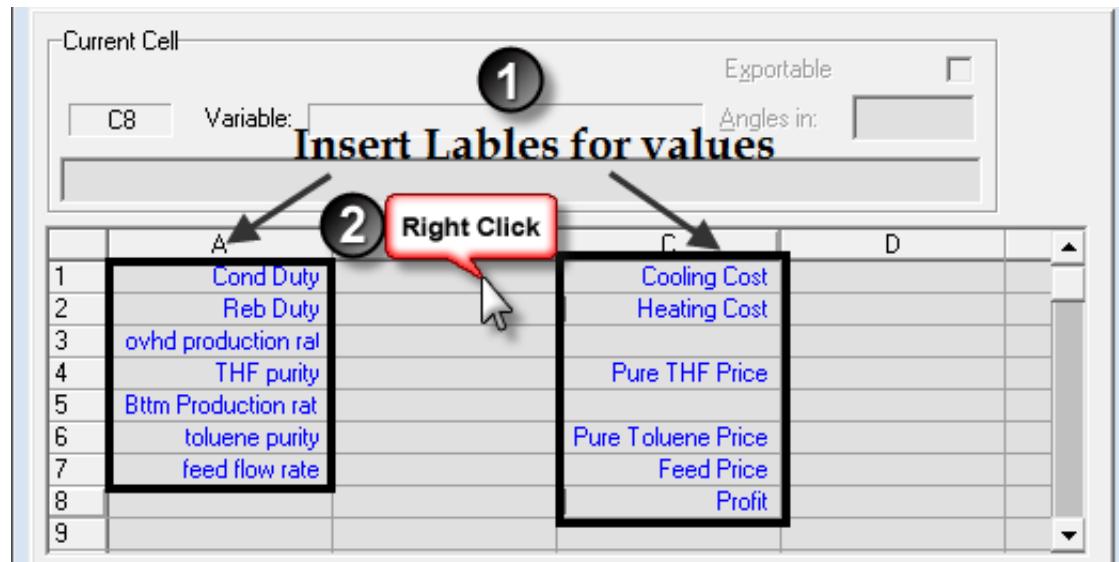


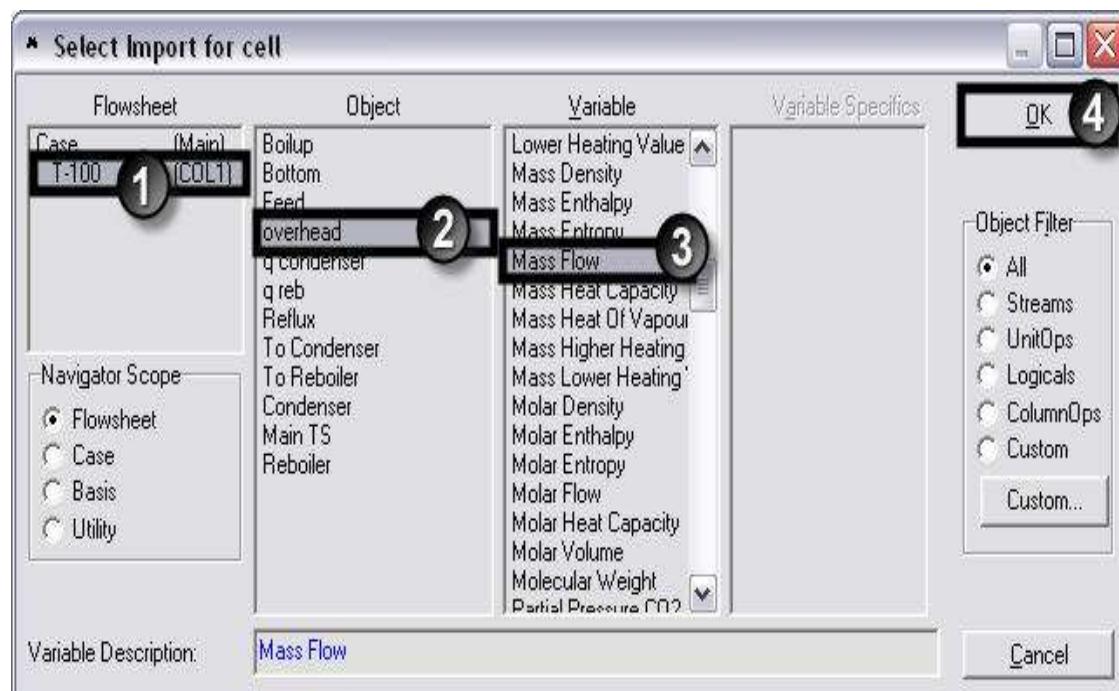
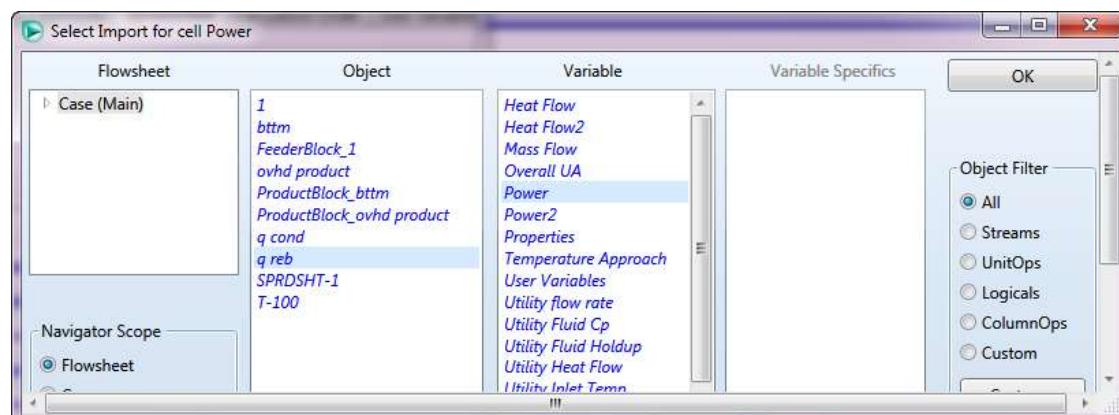
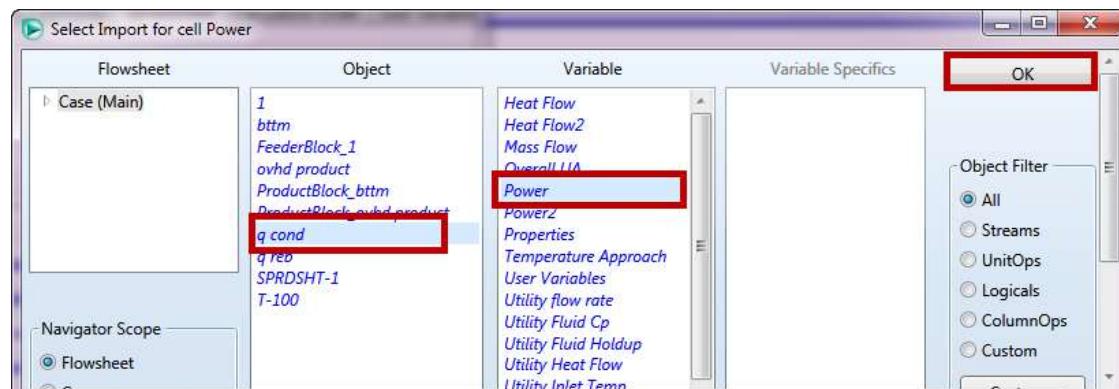
Now we have to start building the profit module using the spreadsheet operation:



Profit= Income - Cost

Profit = (Total Toluene selling price + Total THF selling price) - (Feed cost + Heating cost + Cooling Cost)

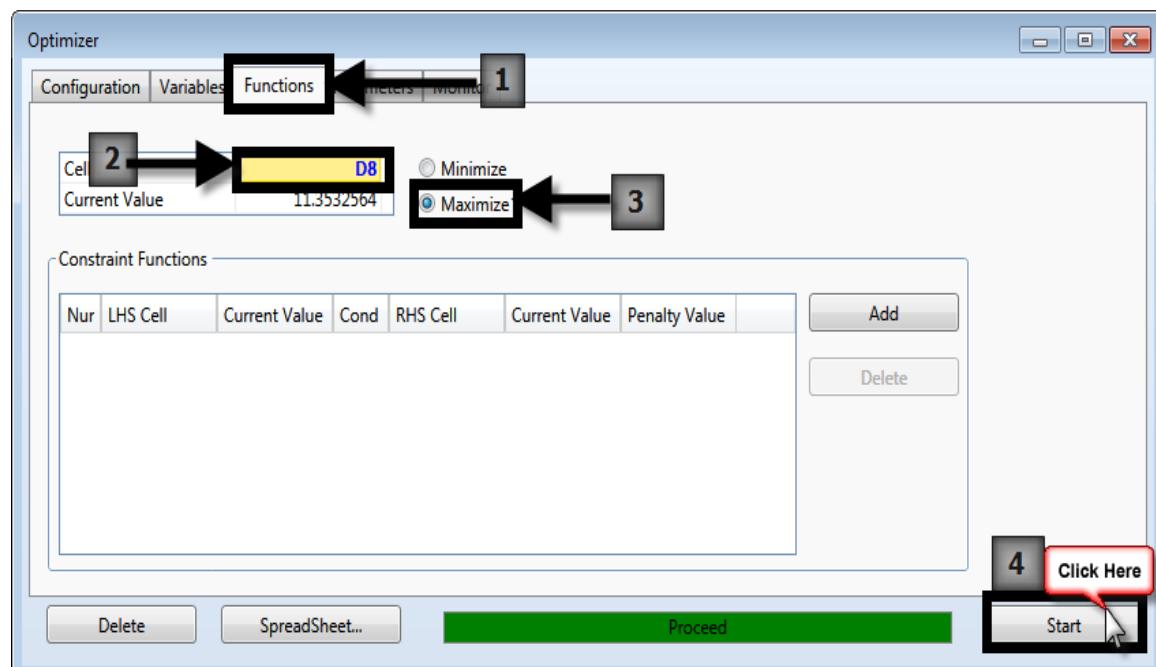


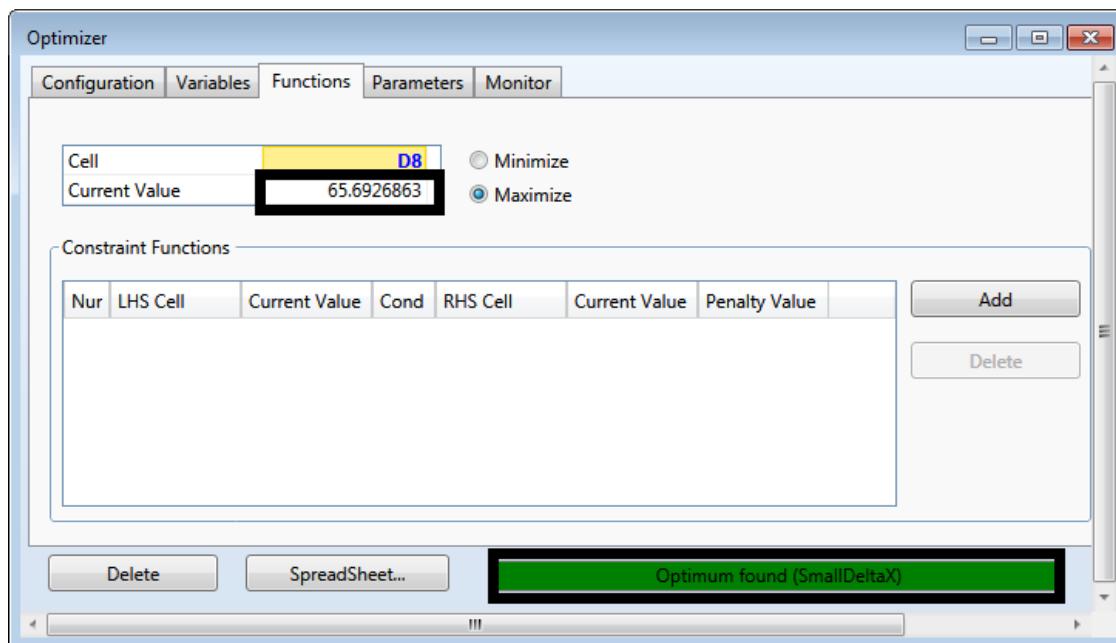


Connections Parameters Formulas Spreadsheet Calculation Order User Variables Notes				
Current Cell Variable Type: <input type="text"/> D8 Variable: <input type="text"/> Angles in: Rad				
$=-(d4*b4*b3+d6*b6*b5)-(d7*b7+d1*b1+d2*b2)$				
1	A cond duty	378.9 kW	C cooling cost	D 0.4710
2	A reb duty	548.4 kW	C heating cost	D 0.7370
3	A ovhd prod rate	1504 kg/h		
4	A thf purity	0.9950	C thf sp	D 0.3330
5	A bttm prod rate	2196 kg/h		
6	A tol purity	0.9400	C tol sp	D 0.1360
7	A feed flowrate	3700 kg/h	C feed price	D 5.000e-002
8			C profit	D 11.35
<a href="#">Function Help...</a> <a href="#">Spreadsheet Only...</a>				

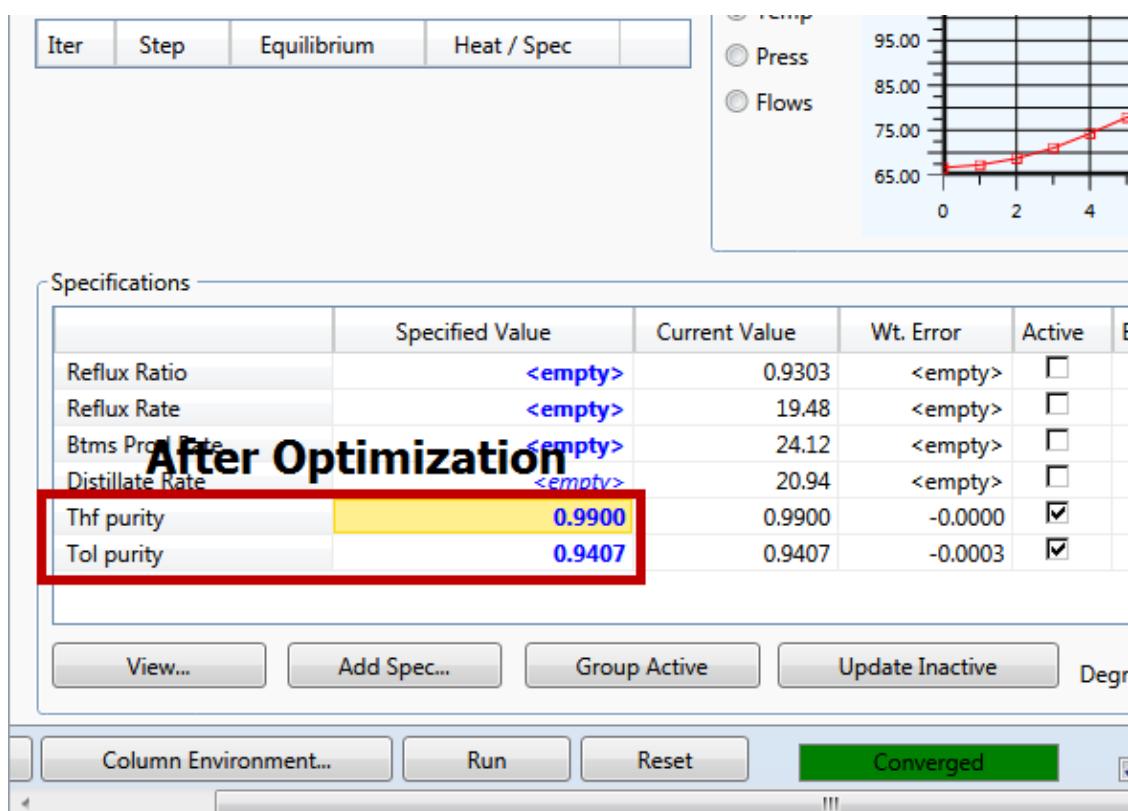
The profit formula will be in D8:

$$= (d3*b3*b4+d5*b5*b6)-(d1*b1+d2*b2+d7*b7)$$





Now you can go to the monitor tab inside the column to see the optimum values for the THF & Toluene purities.



## Save Your Case!

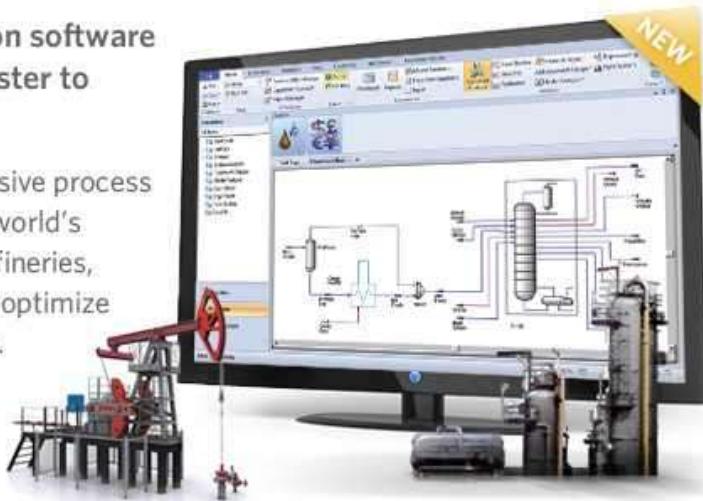
# Gas Gathering

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V8

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Aspen HYSYS is a comprehensive process modeling system used by the world's leading oil & gas producers, refineries, and engineering companies to optimize process design and operations.



8

## ***Workshop***

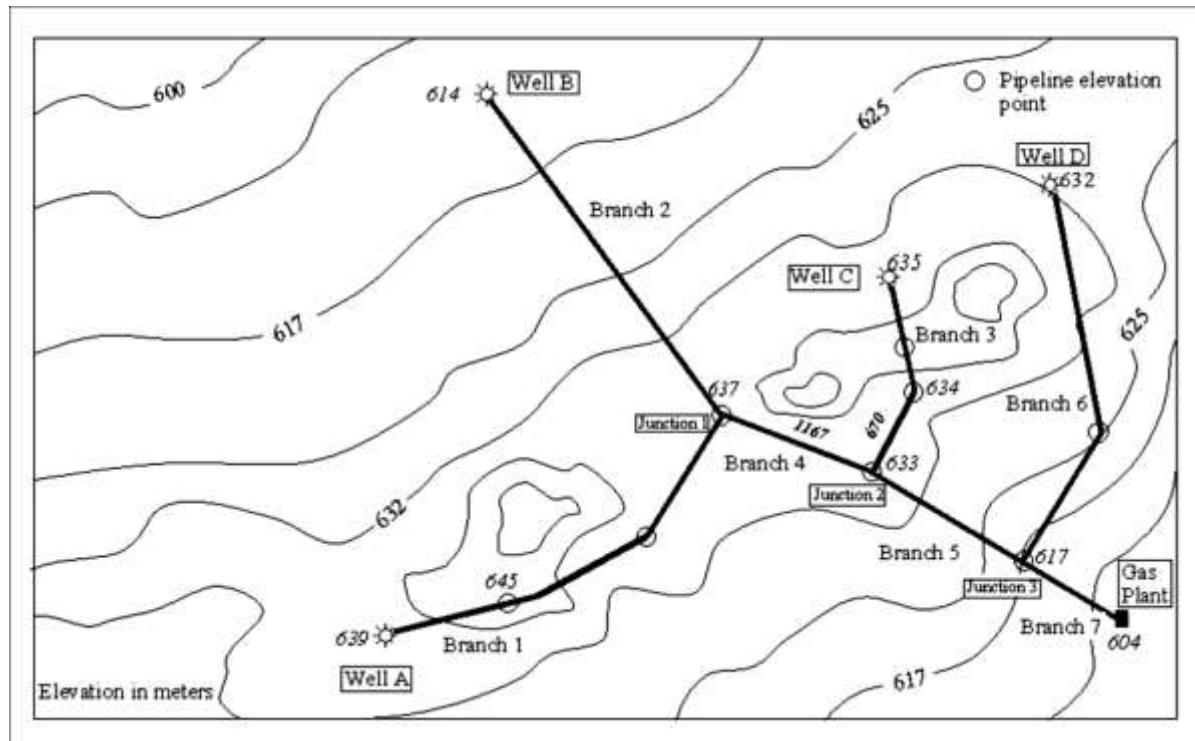
*In this example, a gas gathering system located on varied terrain is simulated using the steady state capabilities of HYSYS. The following figure shows the physical configuration of this system superimposed on a topographic map. The system consists of four wells distributed over an area of approximately 2.0 square km, connected to a gas plant via a network of pipelines.*

## ***Learning Objectives***

*Once you have completed this module, you will be able to use the Pipe Segment in HYSYS to model pipelines.*

### **Example:**

There are 4 gas wells, we need to gather the gas from the wells and transfer it to the plant through pipe lines shown below:



The composition of the four wells is the same:

Component	Mole Fraction	Component	Mole Fraction
Methane	0.6230	n-pentane	0.00405
Ethane	0.2800	n-hexane	0.00659
Propane	0.0163	C7+	0.00992
i-butane	0.00433	N <sub>2</sub>	0.00554
n-butane	0.00821	CO <sub>2</sub>	0.0225
i-pentane	0.00416	H <sub>2</sub> S	0.0154

C7+: MW=122, ρ=47.45 lb/ft<sup>3</sup>

	GasWell 1	GasWell 2	GasWell 3	GasWell 4
Temperature °C (°F)	40 (105)	45 (115)	45 (115)	35 (95)
Pressure kPa (psia)	4135 (600)	3450 (500)	<empty>	<empty>
Flow kgmole/h (lbmole/hr)	425 (935)	375 (825)	575 (1270)	545 (1200)

The pipe segments data are given below:

Branch	Segment	Length meters	Elevation meters	Elevation Change meters
<b>Branch 1</b>	GasWell 1		639	
	1	150	645	6
	2	125	636.5	-8.5
	3	100	637	0.5
<b>Branch 2</b>	GasWell 2			614
	1	200	637	23
<b>Branch 3</b>	GasWell 3		635.5	
	1	160	648	12.5
	2	100	634	-14
	3	205	633	-1
<b>Branch 4</b>	Branch 1 & 2		637	
	1	355	633	-4
<b>Branch 5</b>	Branch 3 & 4		633	633
	1	300	617	-16
<b>Branch 6</b>	GasWell 4		632.5	
	1	180	625	-7.5
	2	165	617	-8
<b>Branch 7</b>	Branch 5 & 6		617	
	1	340	604	-13

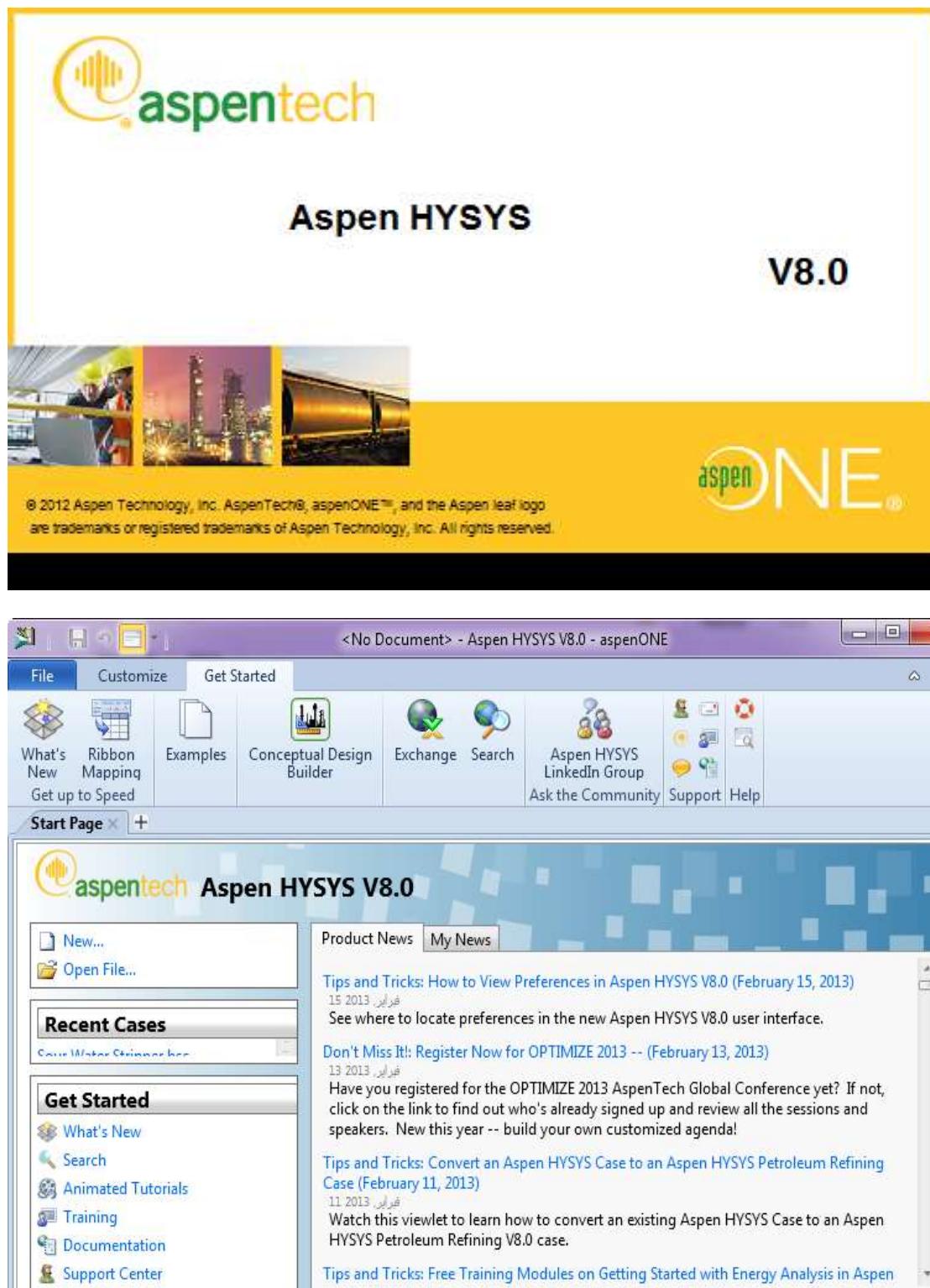
Pipe Branch	Diameter
Branch 1	76.2 mm (3")
Branch 2	101.6 mm (4")
Branch 3	76.2 mm (3")
Branch 4	101.6 mm (4")
Branch 5	152 mm (6")
Branch 6	76.2 mm (3")
Branch 7	152 mm (6")

Schedule 40 steel pipes is used throughout and all branches are buried at a depth of 1 m (3 ft). All pipes are uninsulated

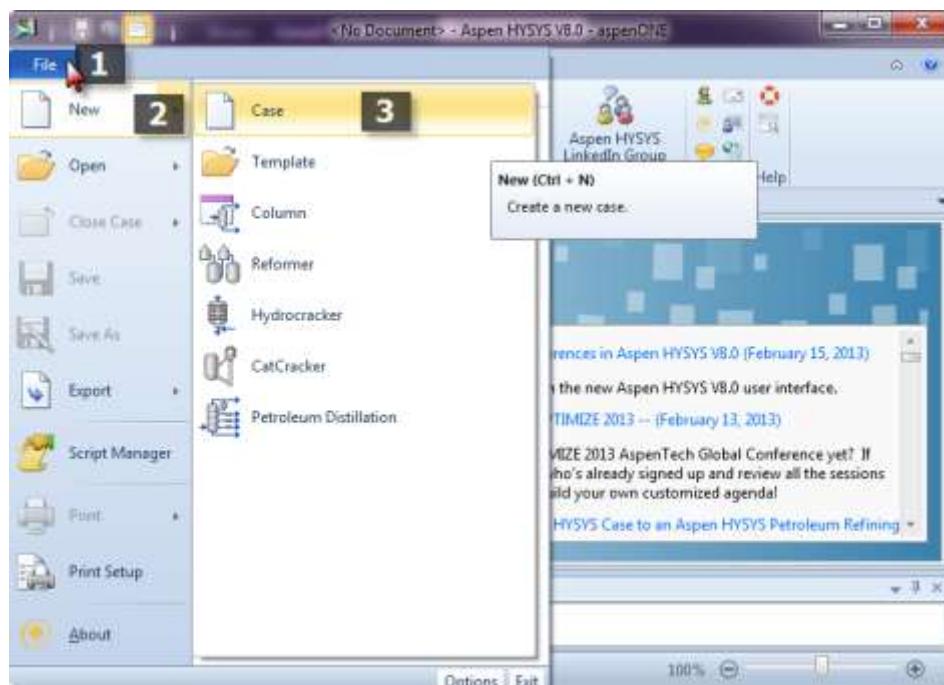
Consider inner and outer HTC and the pipe wall in heat transfer estimation. (Ambient Temperature=5°C)

- Calculate the pressure drop and the heat loses inside each branch.

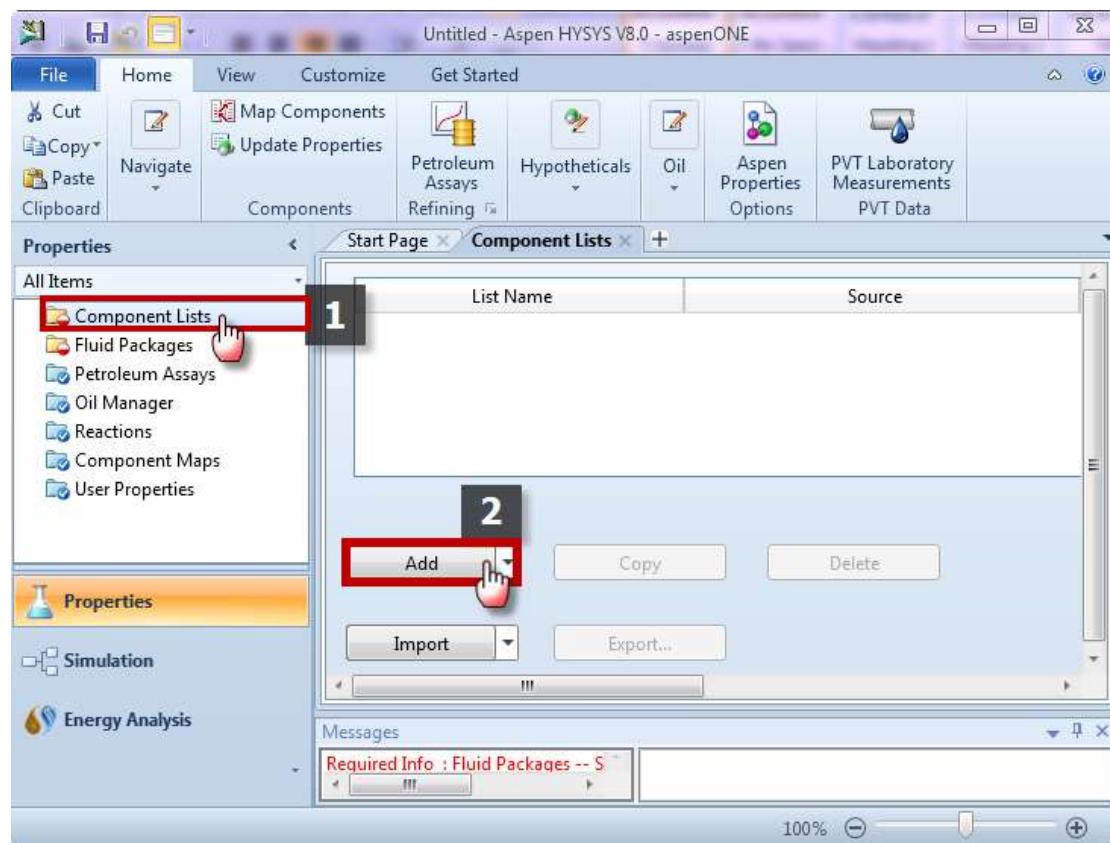
To start the program, From Start Menu, Select All Programs >>  
Aspen Tech >> Process Modeling V8.x >>> Aspen HYSYS >>  
Aspen HYSYS



## 8- First, Start a new case



## 9- Add the Components



10- Choose the system components from the databank:

The screenshot shows the 'Component List' dialog box in Aspen HYSYS V8. On the left, a table lists selected components: Methane, Propane, and Ethane, all categorized as 'Pure Component'. On the right, a larger table lists various hydrocarbons with their simulation names, full names/synonyms, and formulas. The first row, 'Methane', is highlighted in yellow. A red box highlights the '< Add' button, which is labeled with the number '2'. Another red box highlights the 'i-Butane' entry in the list, which is labeled with the number '1'.

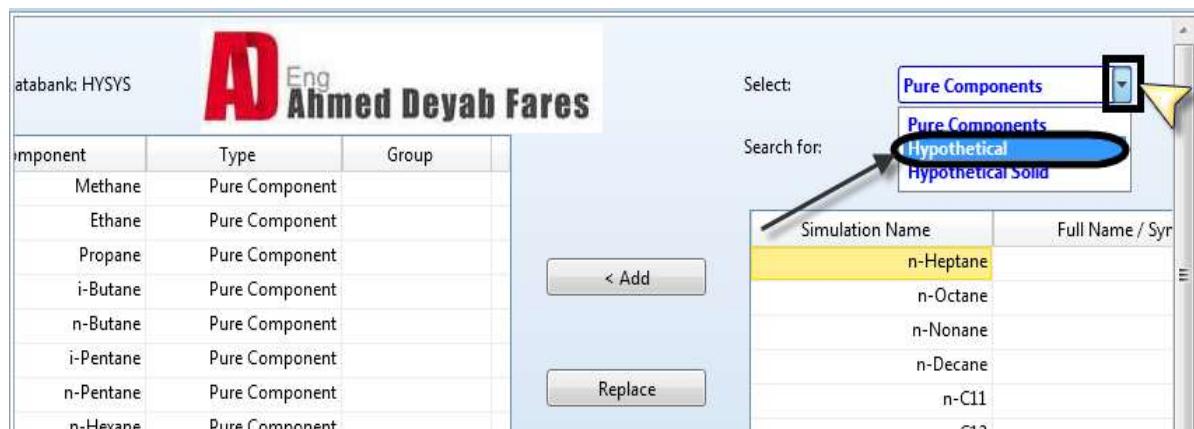
Component	Type	Group
Methane	Pure Component	
Propane	Pure Component	
Ethane	Pure Component	

Simulation Name	Full Name / Synonym	Formula
Methane	C1	CH <sub>4</sub>
Ethane	C2	C <sub>2</sub> H <sub>6</sub>
Propane	C3	C <sub>3</sub> H <sub>8</sub>
i-Butane	i-C4	C <sub>4</sub> H <sub>10</sub>
n-Butane	n-C4	C <sub>4</sub> H <sub>10</sub>
i-Pentane	i-C5	C <sub>5</sub> H <sub>12</sub>
n-Pentane	n-C5	C <sub>5</sub> H <sub>12</sub>
n-Hexane	C6	C <sub>6</sub> H <sub>14</sub>
n-Heptane	C7	C <sub>7</sub> H <sub>16</sub>
n-Octane	C8	C <sub>8</sub> H <sub>18</sub>
n-Nonane	C9	C <sub>9</sub> H <sub>20</sub>
n-Decane	C10	C <sub>10</sub> H <sub>22</sub>
n-C11	C11	C <sub>11</sub> H <sub>24</sub>
n-C12	C12	C <sub>12</sub> H <sub>26</sub>

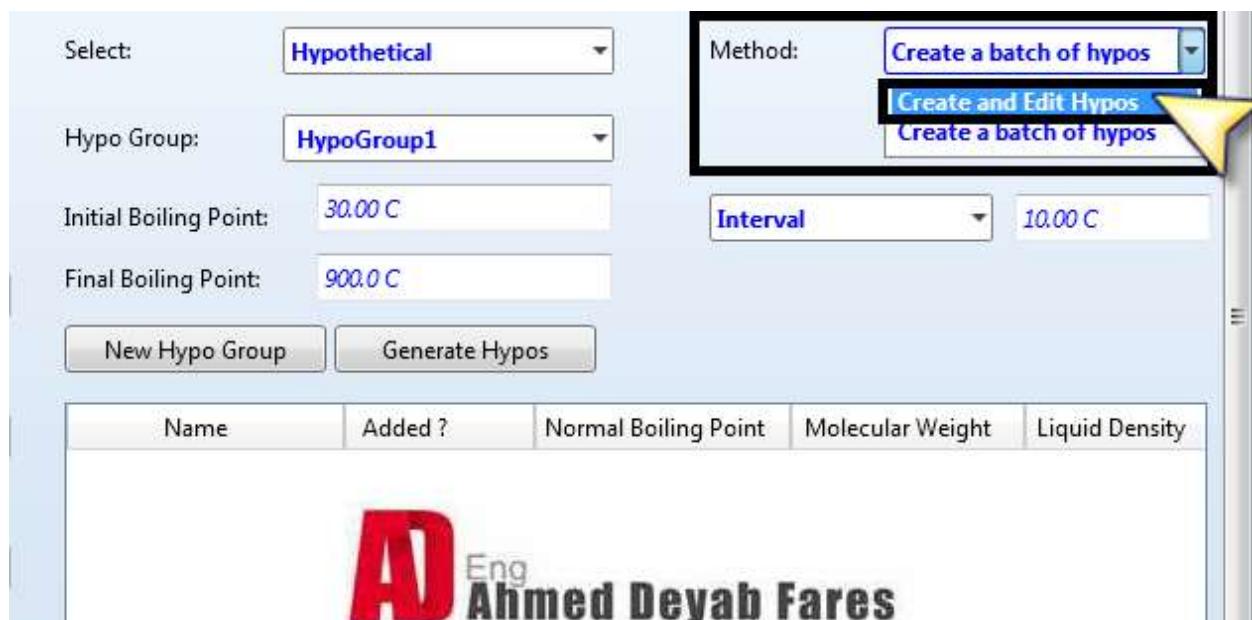
After adding the pure components (N<sub>2</sub>, H<sub>2</sub>S, CO<sub>2</sub>, C1, C2, C3, n-C4, i-C4, n-C5, i-C5, n-C6, H<sub>2</sub>O) we have to add the last component (C7<sup>+</sup>) which is not a pure component as it represents all components above C7 including C7 in the feed.

To define C7<sup>+</sup> we have to create it as a hypothetical component as the following:

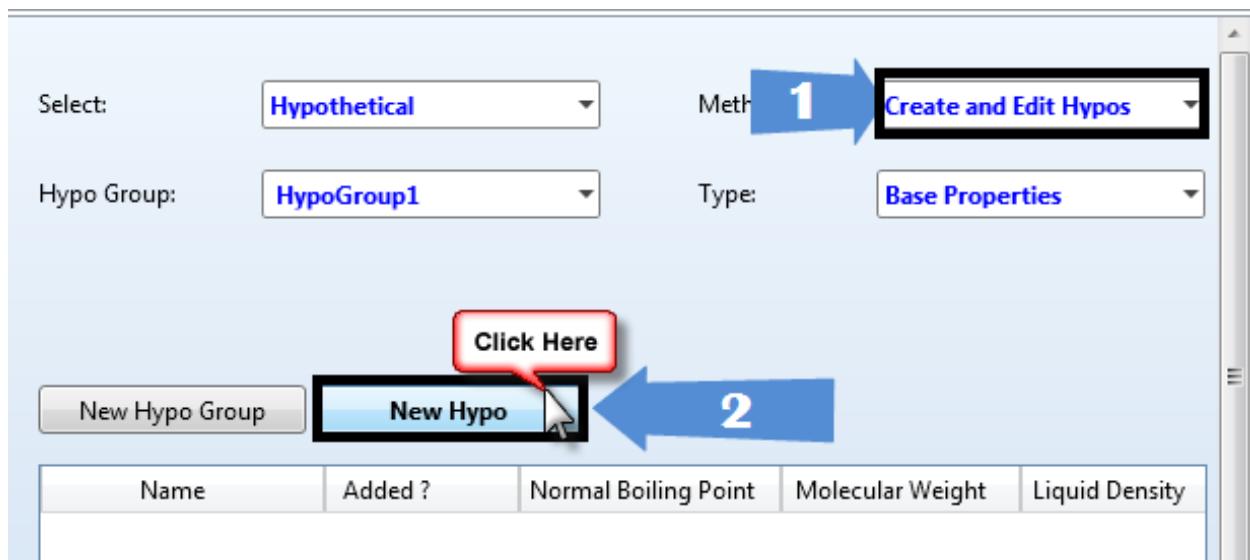
From the drop menu select Hypothetical instead of pure components



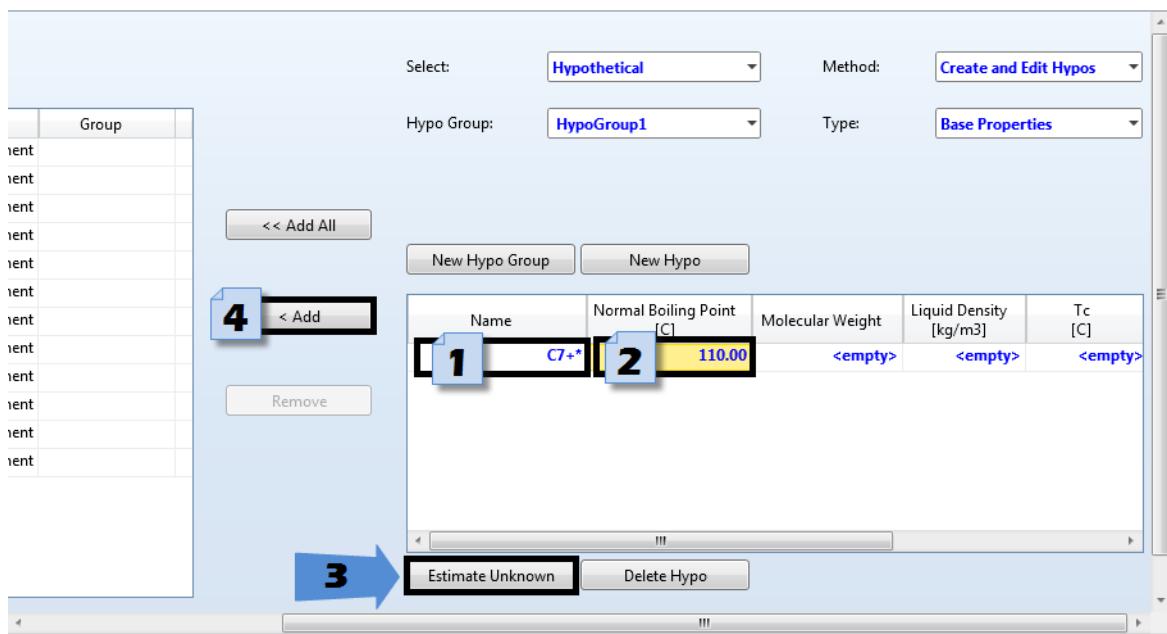
Select create and edit hypos



Click on New Hypo



After adding a hypo component you can edit the name, add the properties you have, and estimate the unknown properties as follows:

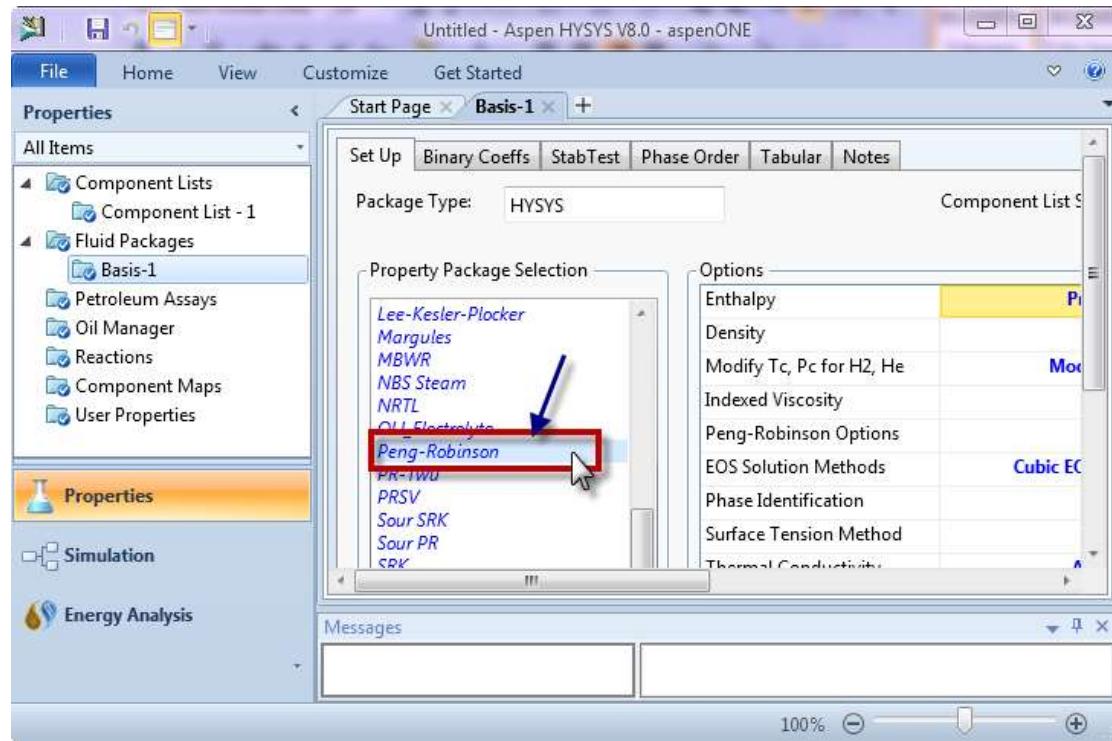


Finally add the hypo component to the component list

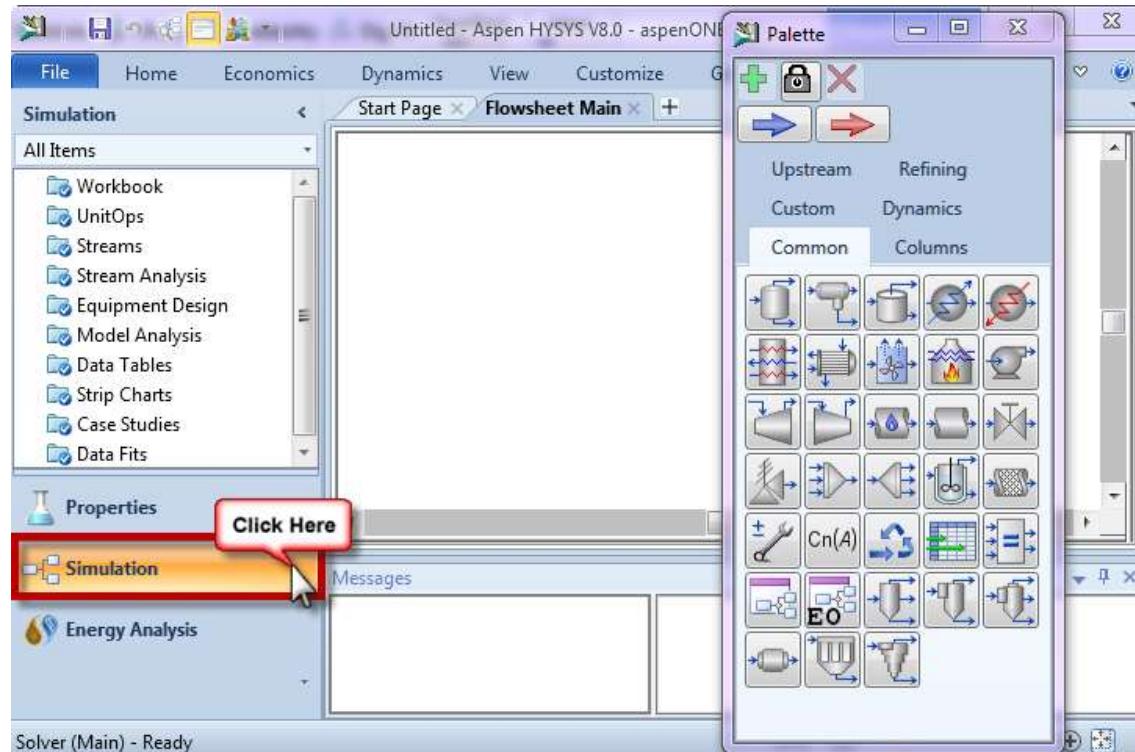
Component	Type	Group
Methane	Pure Component	
Ethane	Pure Component	
Propane	Pure Component	
i-Butane	Pure Component	
n-Butane	Pure Component	
i-Pentane	Pure Component	
n-Pentane	Pure Component	
n-Hexane	Pure Component	
CO <sub>2</sub>	Pure Component	
H <sub>2</sub> S	Pure Component	
Nitrogen	Pure Component	
H <sub>2</sub> O	Pure Component	
C7+* User Defined Hypothetical		HypoGroup1

Now, select the suitable fluid package

In this case, select Peng-Robinson

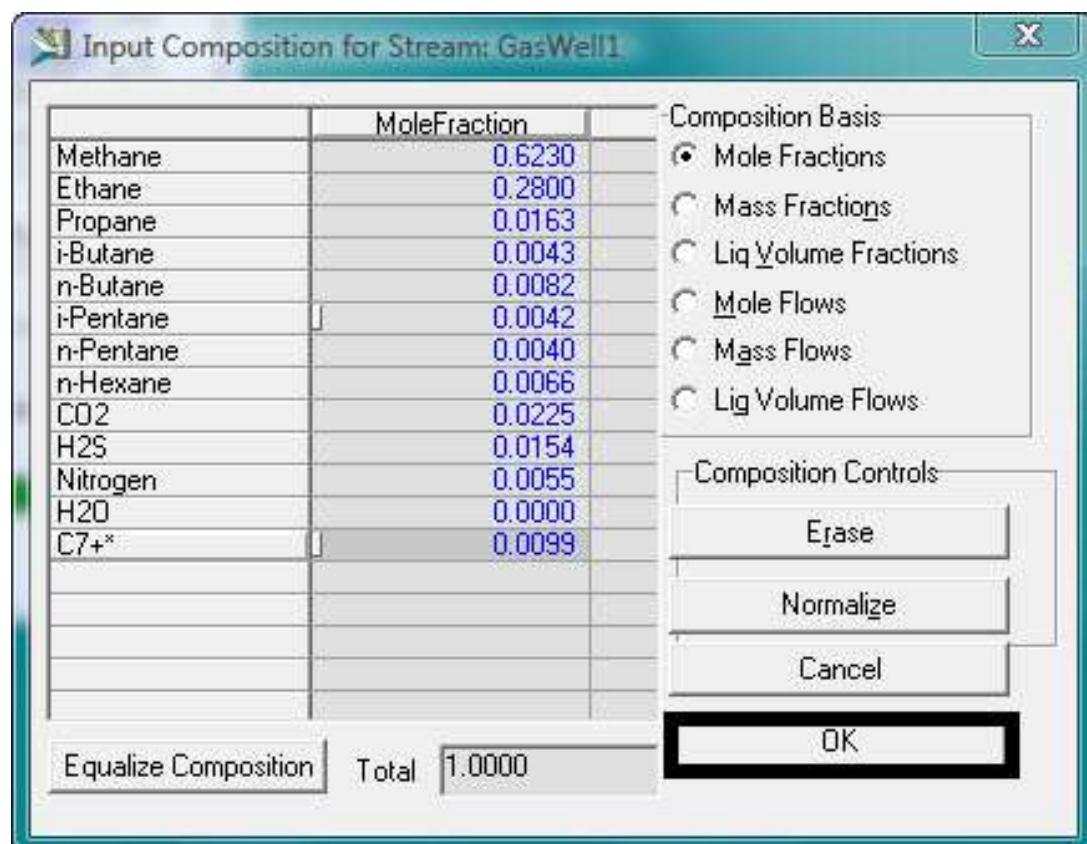
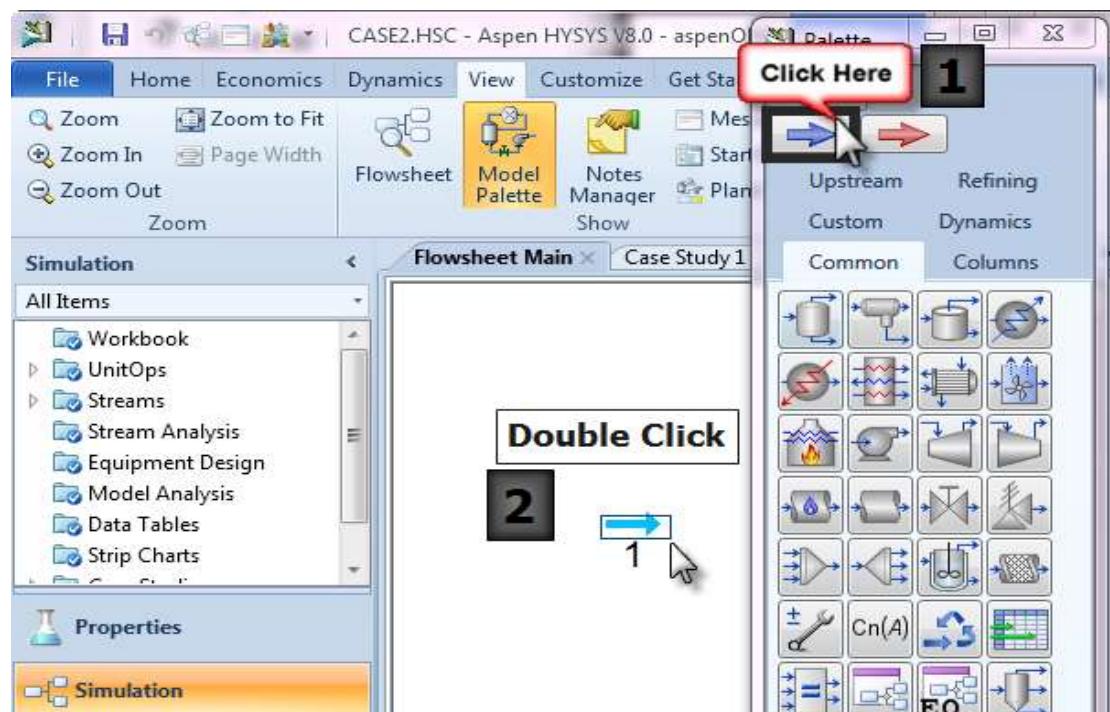


Now you can start drawing the flow sheet for the process by clicking the Simulation button:

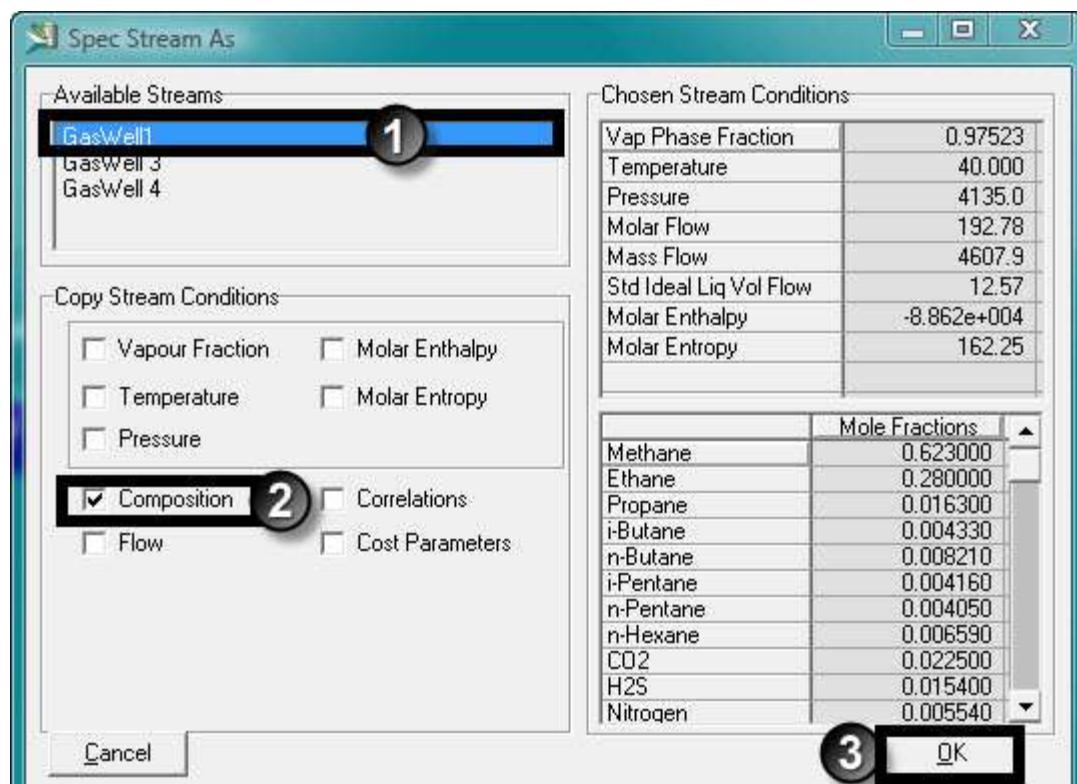
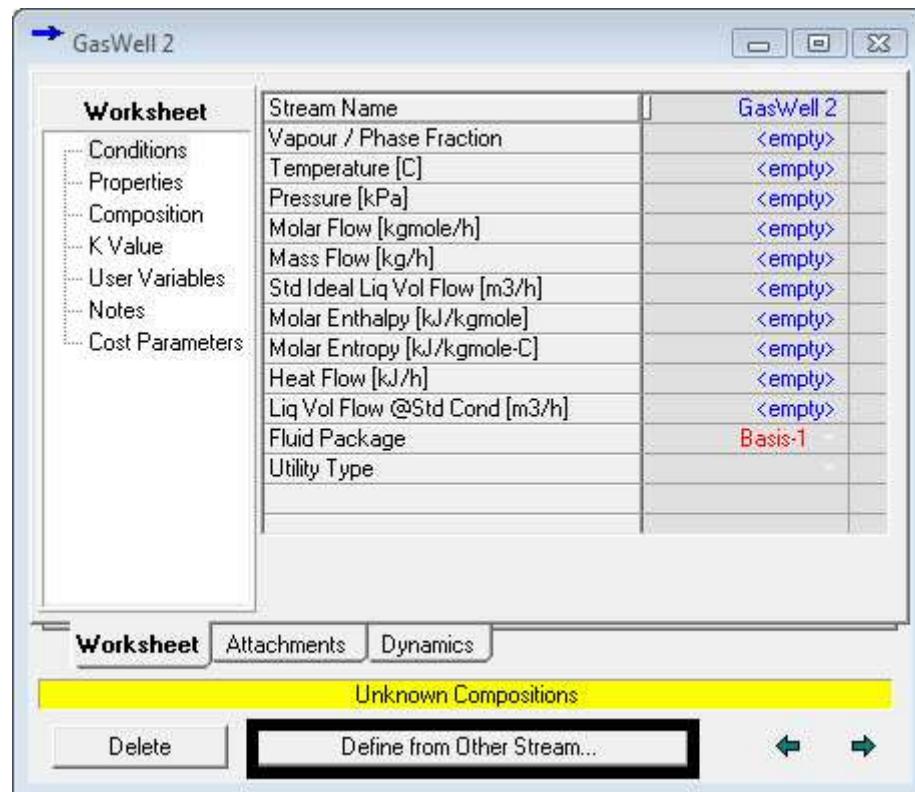


Now add a material stream to define the composition and the conditions of the feed stream

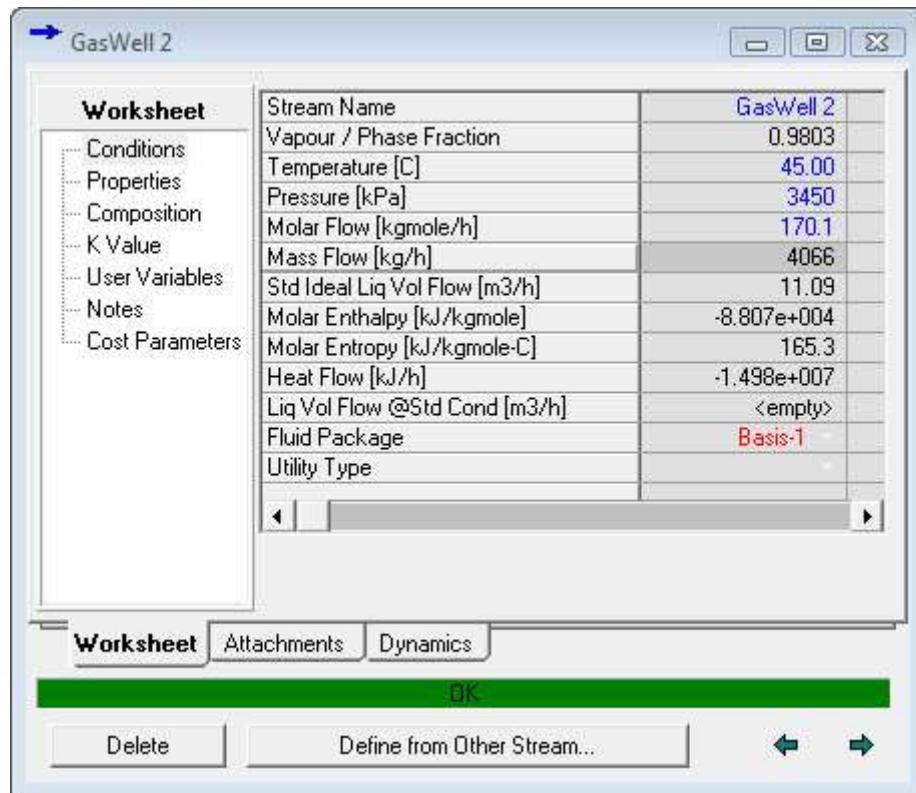
From the palette:



After adding the compositions and the conditions for the first well, add another stream for well 2 and define the composition from the first stream as follows:

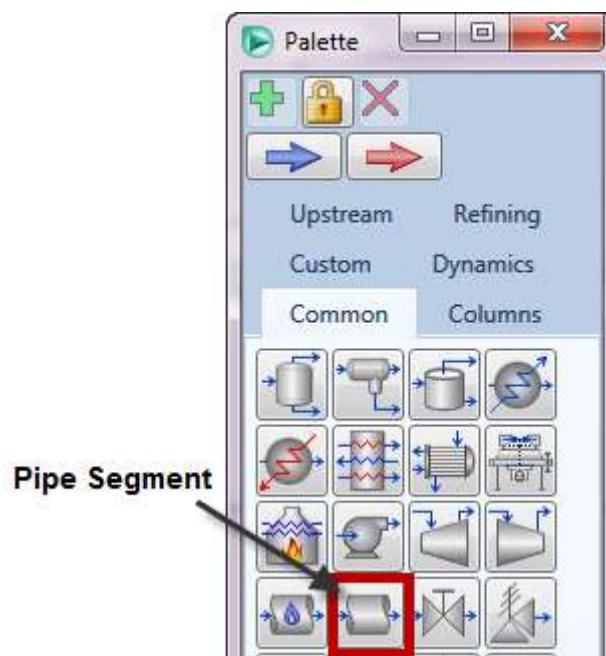


Add the conditions for the second stream

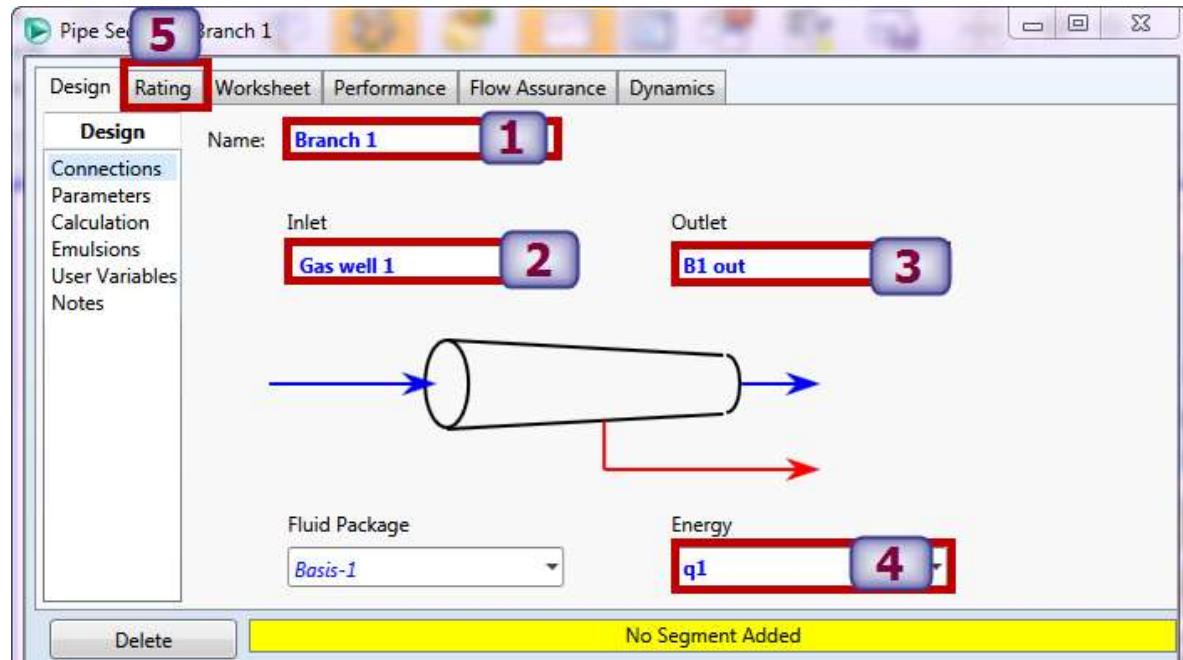


Add the conditions & Compositions for the other two streams as above

Now, Add a **Pipe Segment** from the palette

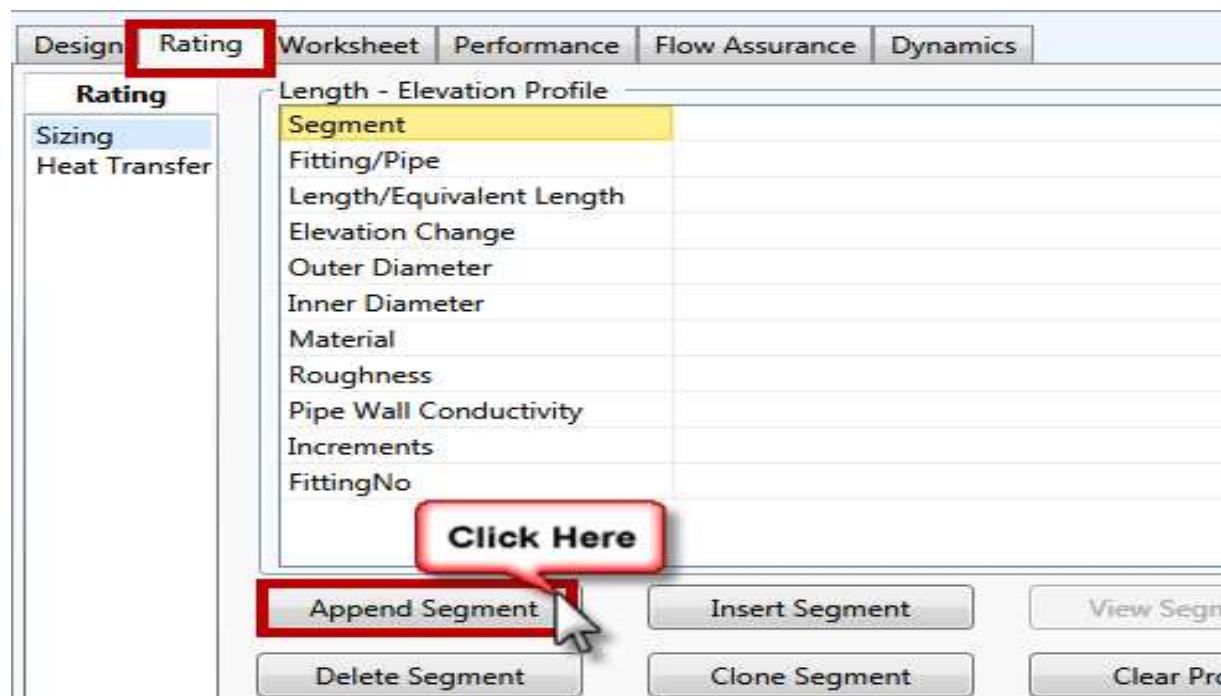


The pipe segment is used to simulate a wide variety of piping situations ranging from single/multiphase plant piping with rigorous heat transfer estimation, to large capacity looped pipeline problems

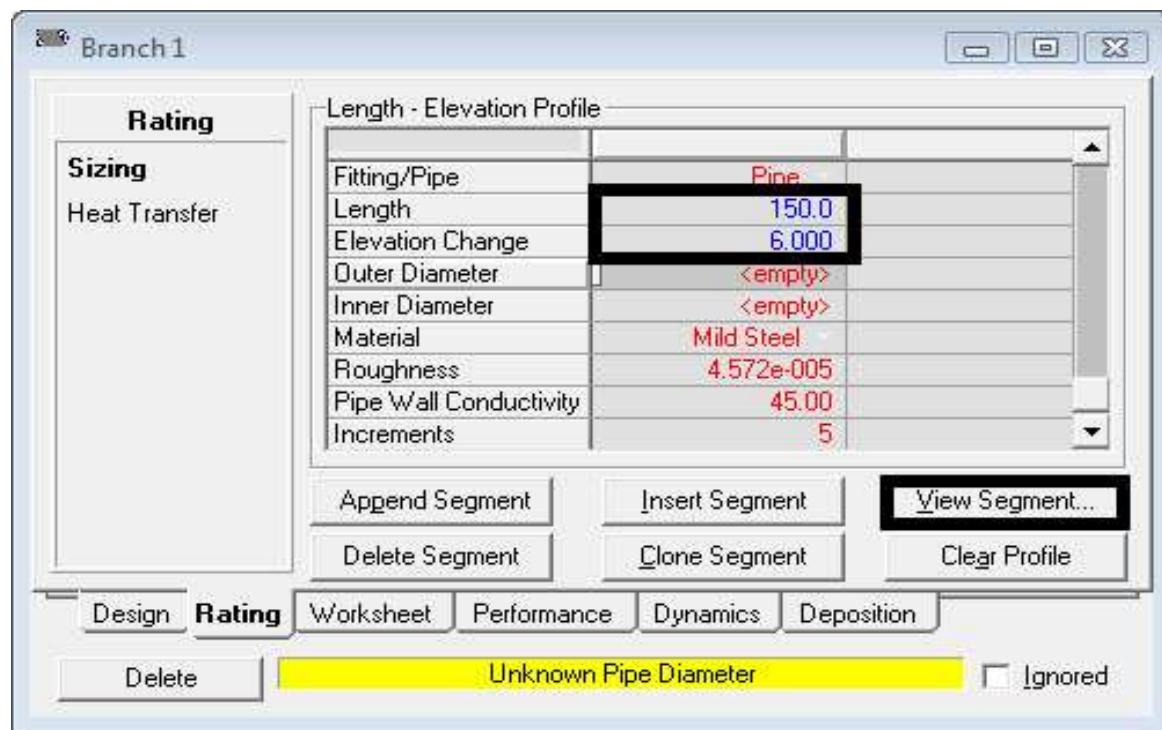


On the **Sizing** page, you construct the length-elevation profile for the Pipe Segment. Each pipe section and fitting is labeled as a segment. To fully define the pipe sections segments, you must also specify pipe schedule, diameters, pipe material and a number of increments.

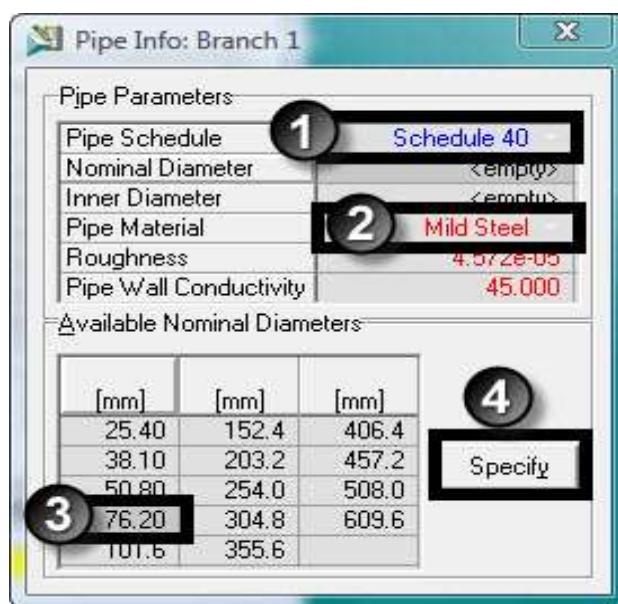
- The first pipe, Branch 1 is broken into three segments.



- Add the first segment to the pipe unit operation by clicking the **Append Segment** button. Specify the following information for the segment.
- To specify the diameter, click the **View Segment** button.

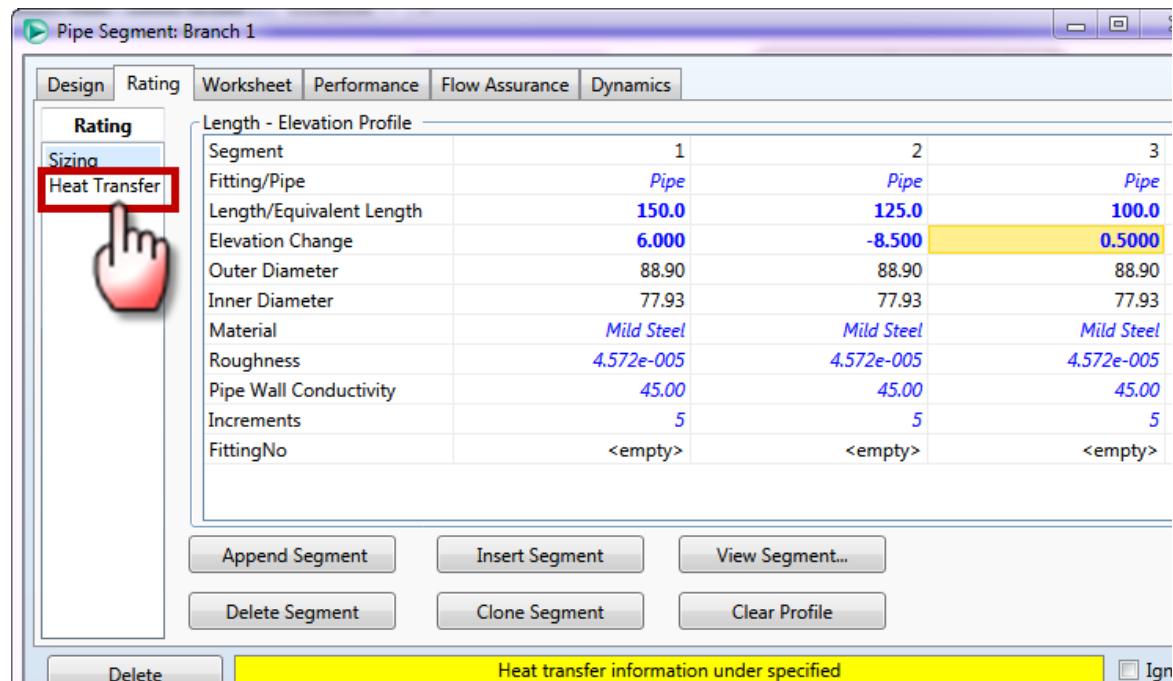


- Select **Schedule 40** as the Pipe Schedule.
- From the Available Nominal Diameters group, select **76.20 mm (3 inch)** diameter pipe and click the **Specify** button. The Outer and Inner Diameter will be calculated by HYSYS.
- Use the default Pipe Material, Mild Steel



- Two more segments are needed to complete the branch.

In this cell...	Enter...	Enter...
Segment	2	3
Fitting/Pipe	Pipe	Pipe
Length	125 m (410 ft)	100 m (325 ft)
Elevation	-6.5 m (-21 ft)	0.5 m (1 ft)
Schedule	40	40
Nominal Diameter	76.2 mm (3 inch)	76.2 mm (3 inch)

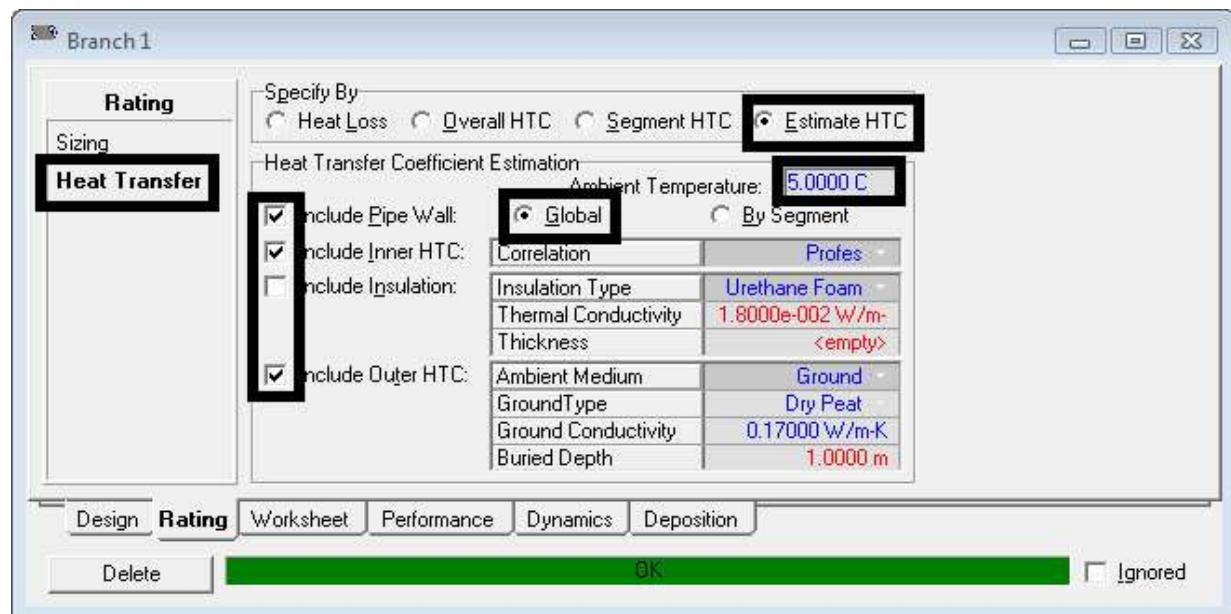


The Pipe Segment is not yet able to solve because we have not specified any information about the heat transfer properties of the pipe.

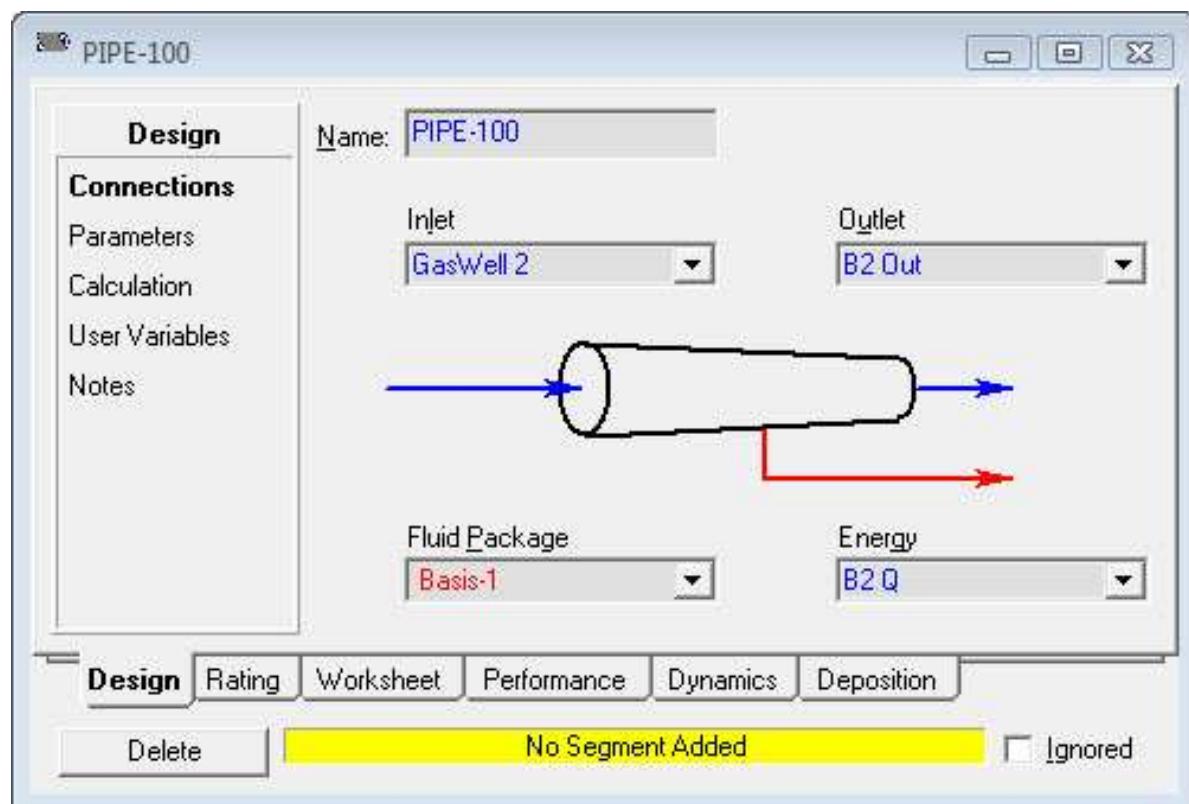
## Heat Transfer page

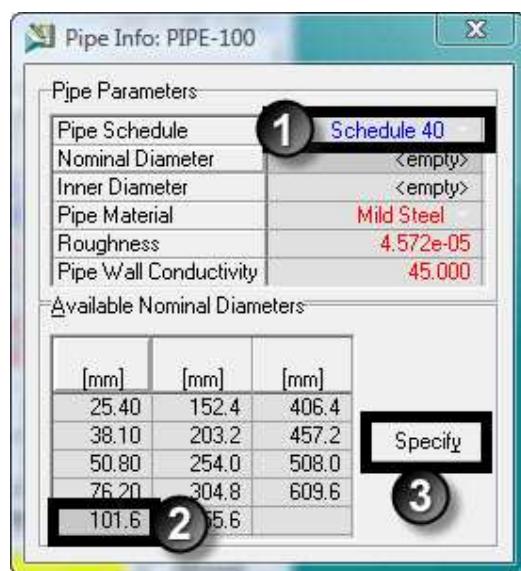
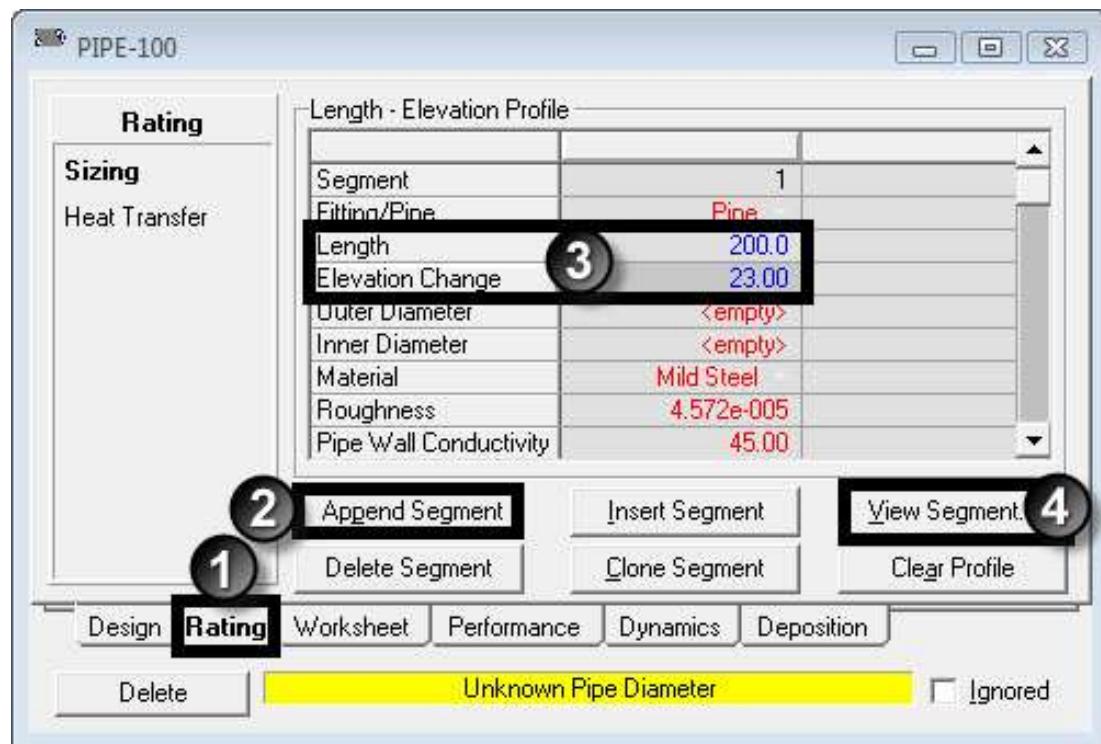
On this page, you select the method that HYSYS will use for the heat transfer calculations.

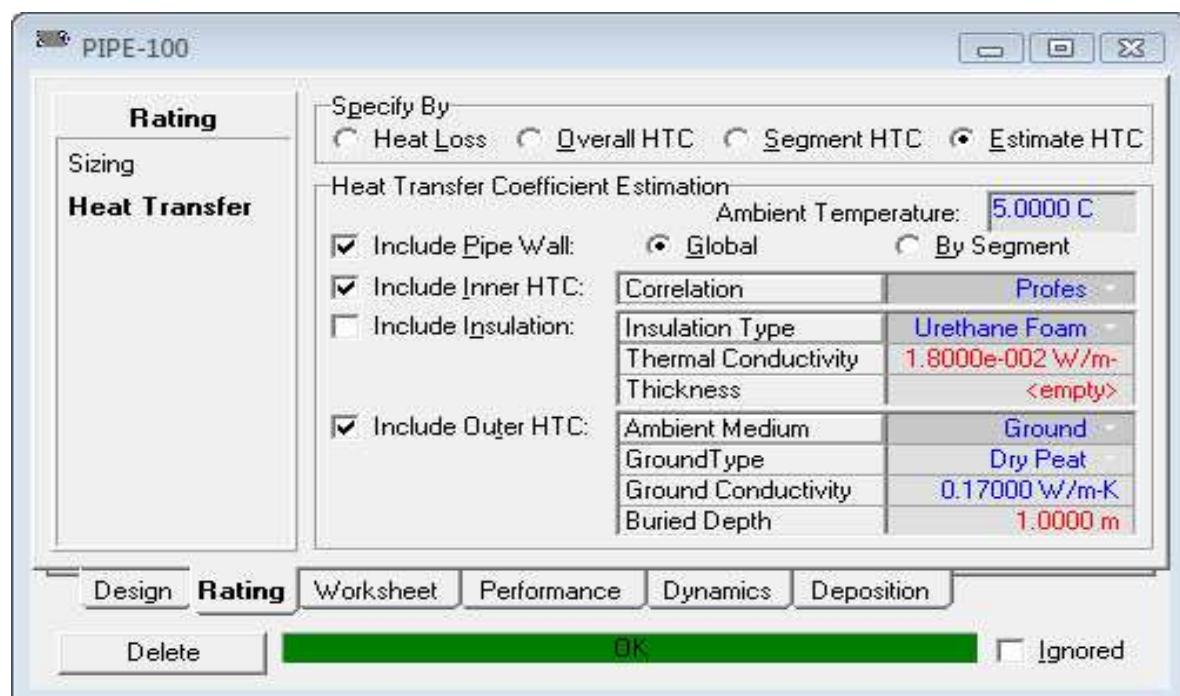
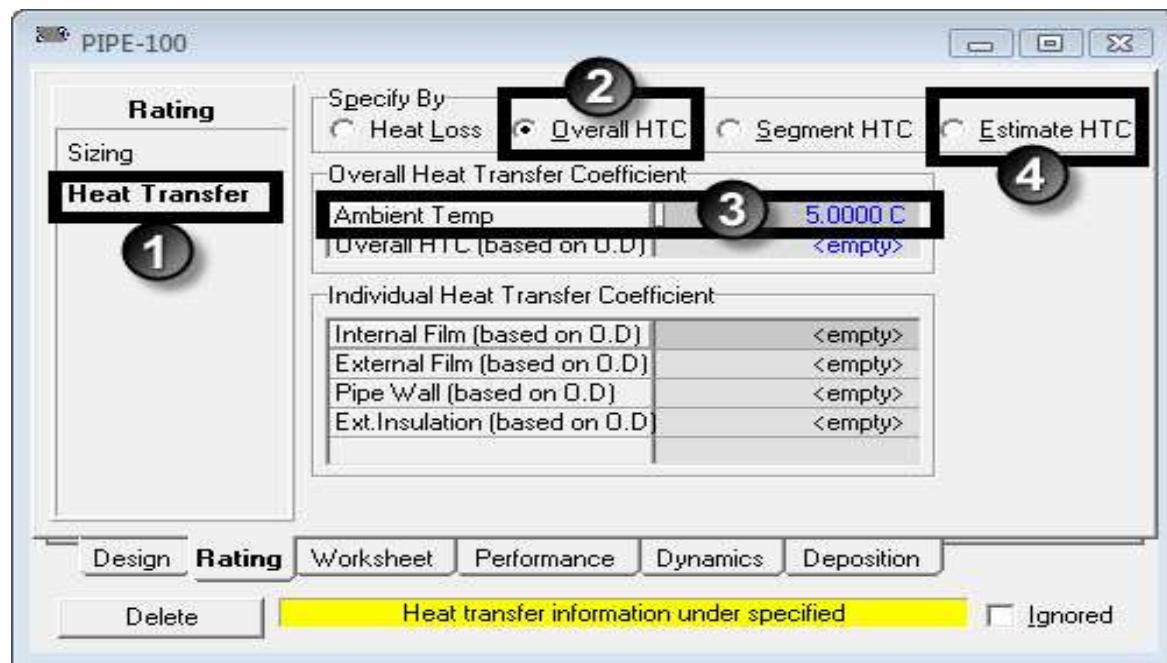
For all pipes in this simulation, use the Estimate HTC method.

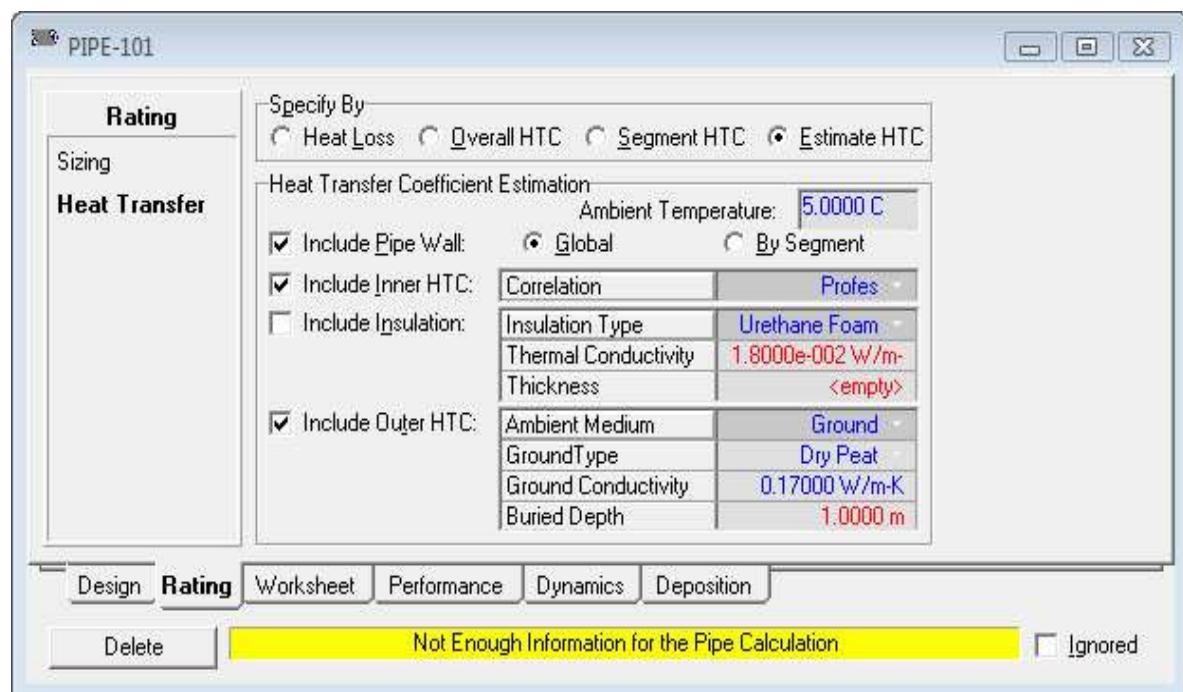
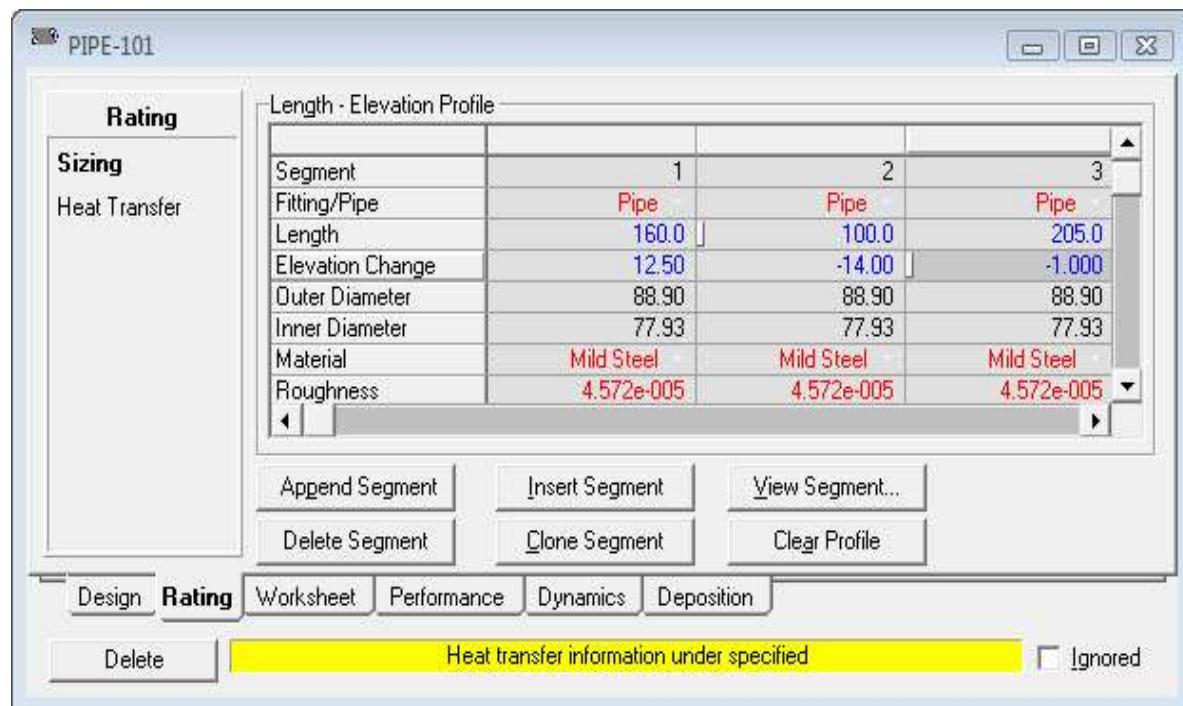


Now add the remaining unit operations to your case.

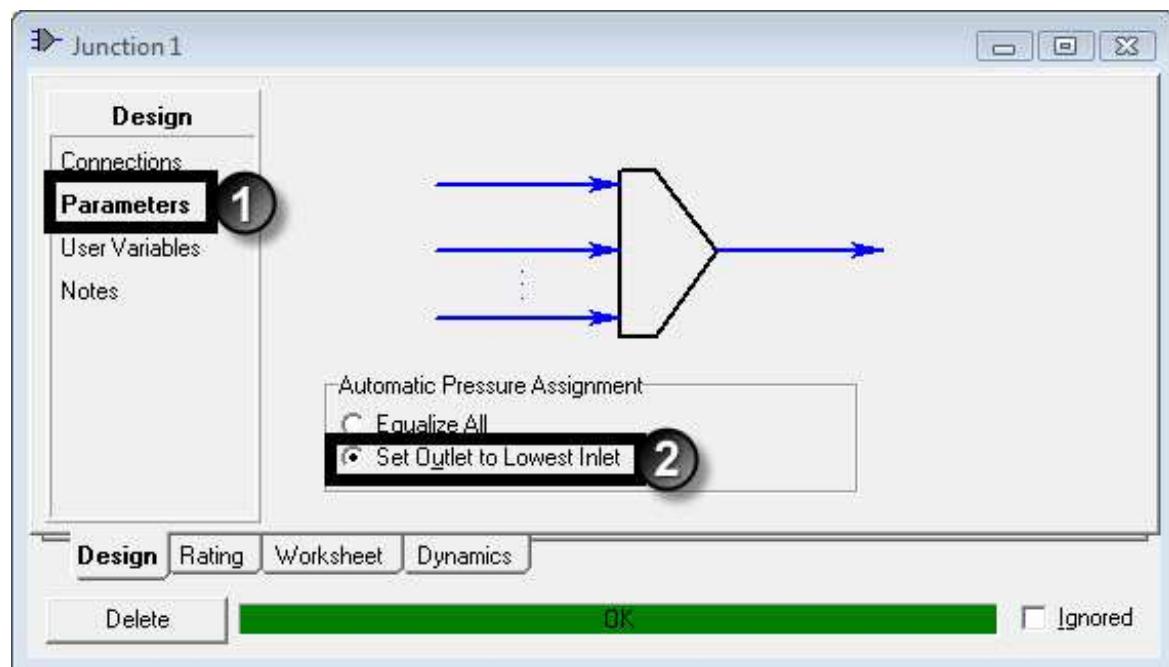
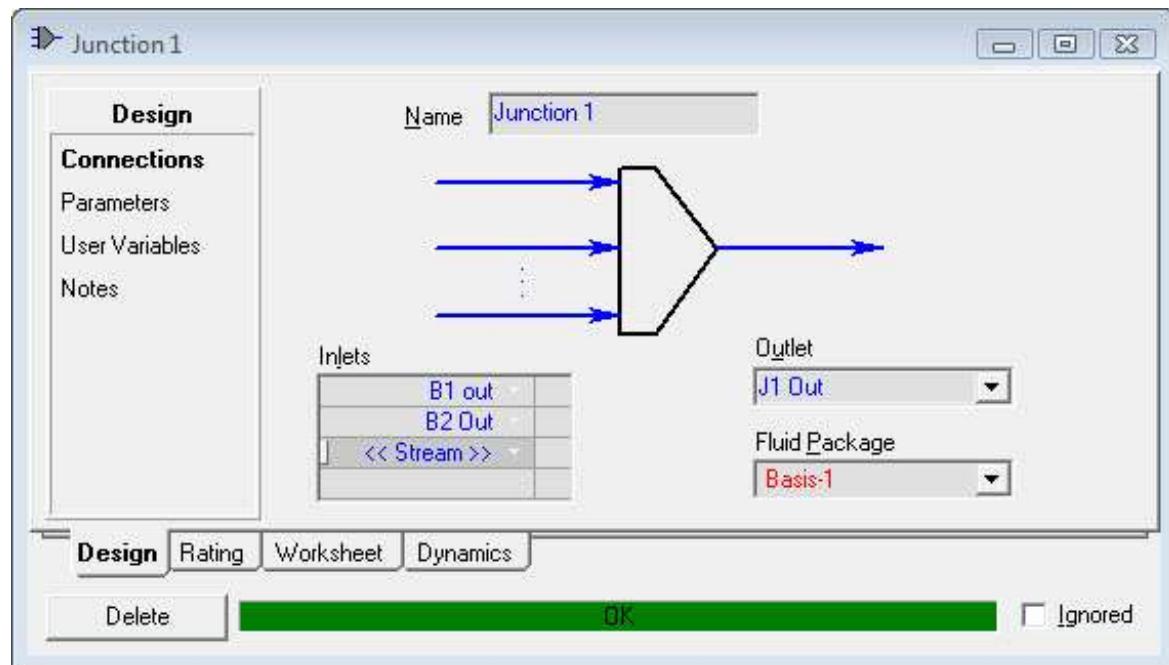


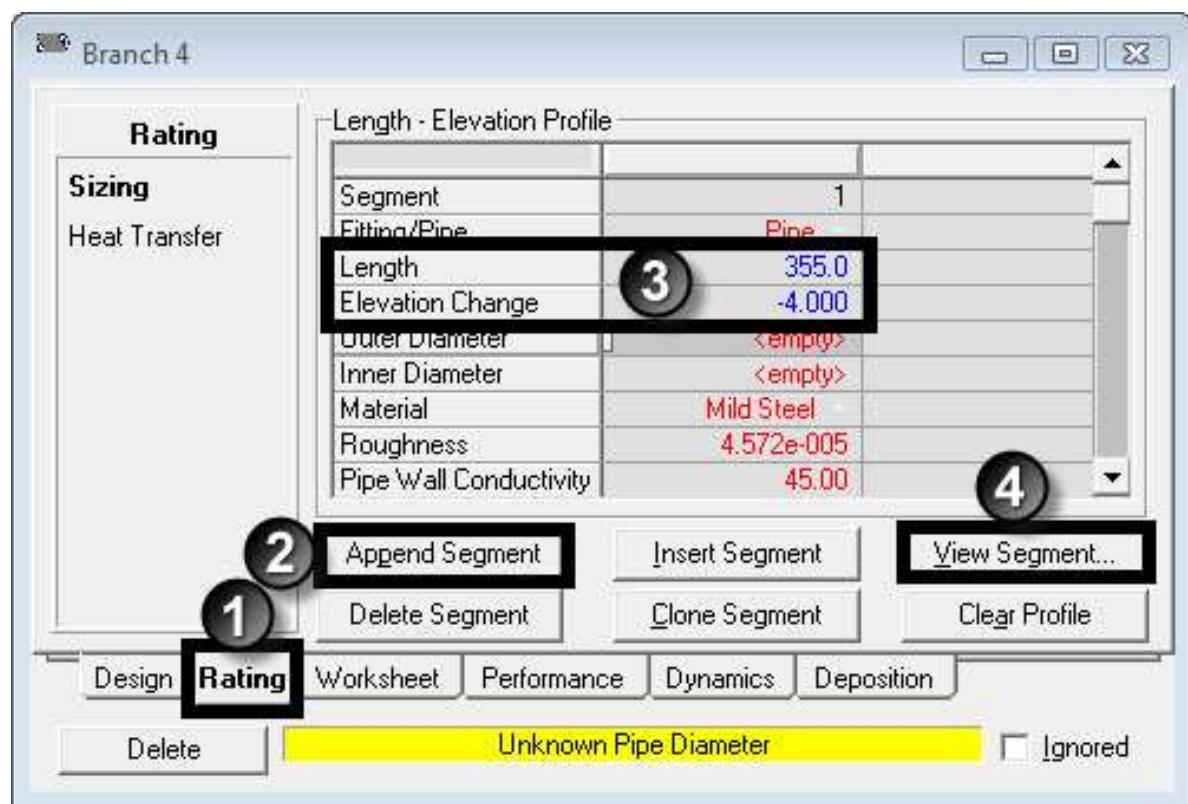
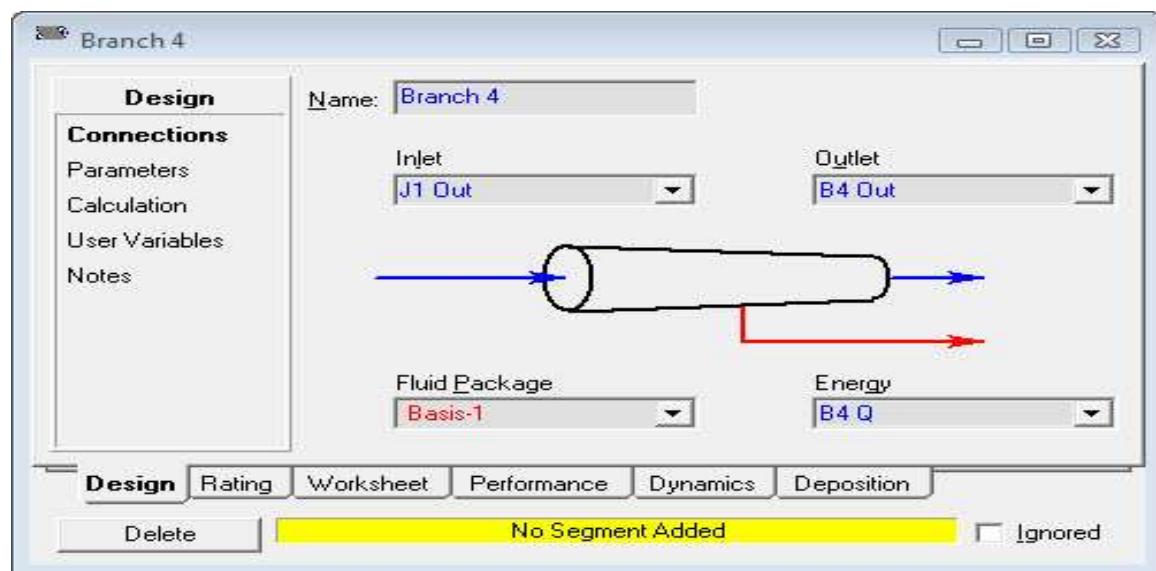


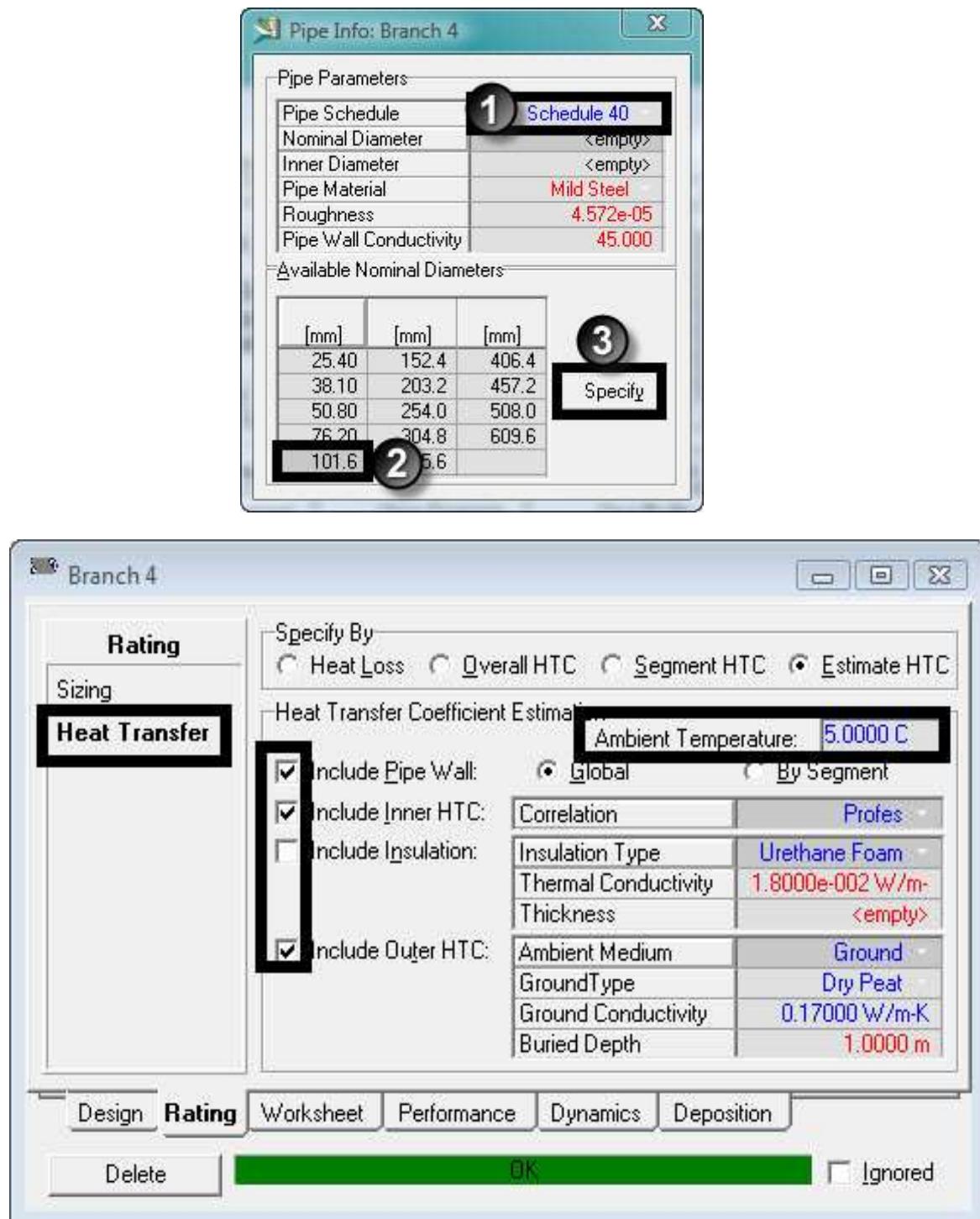


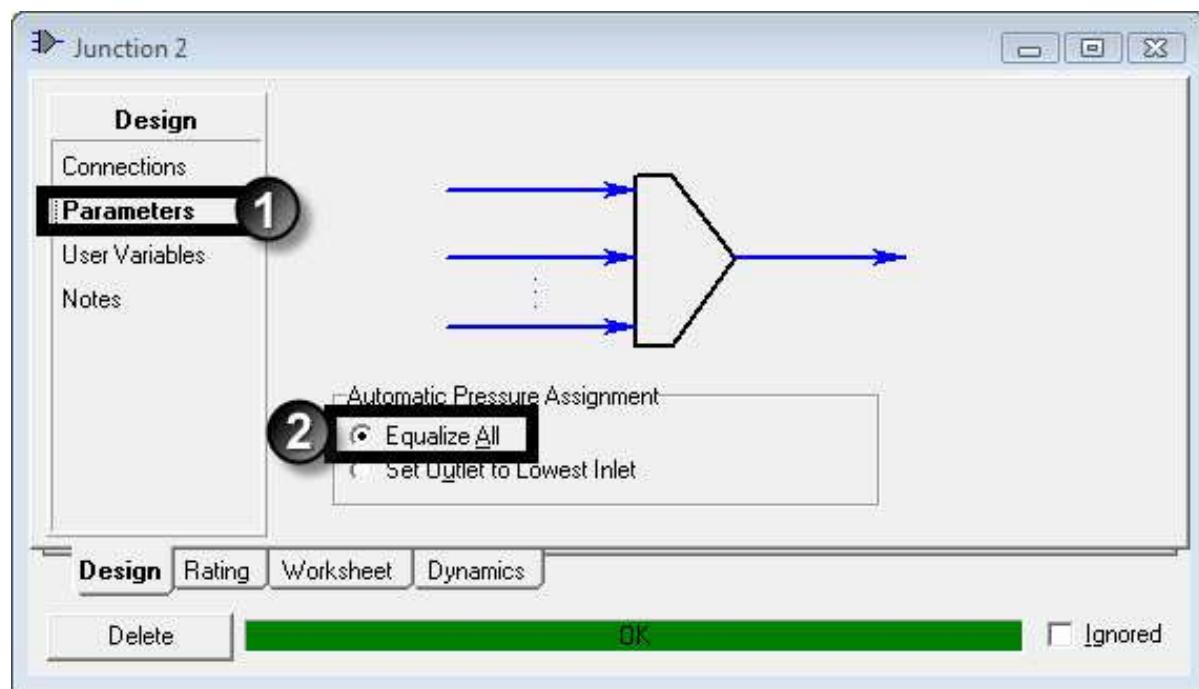
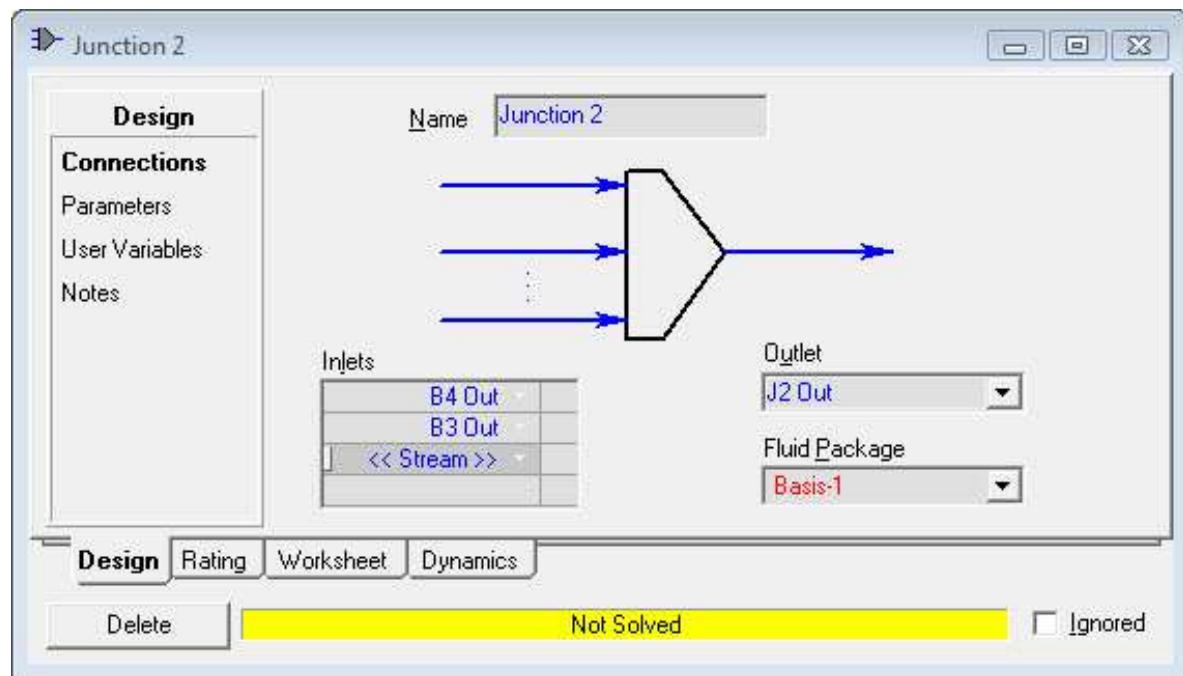


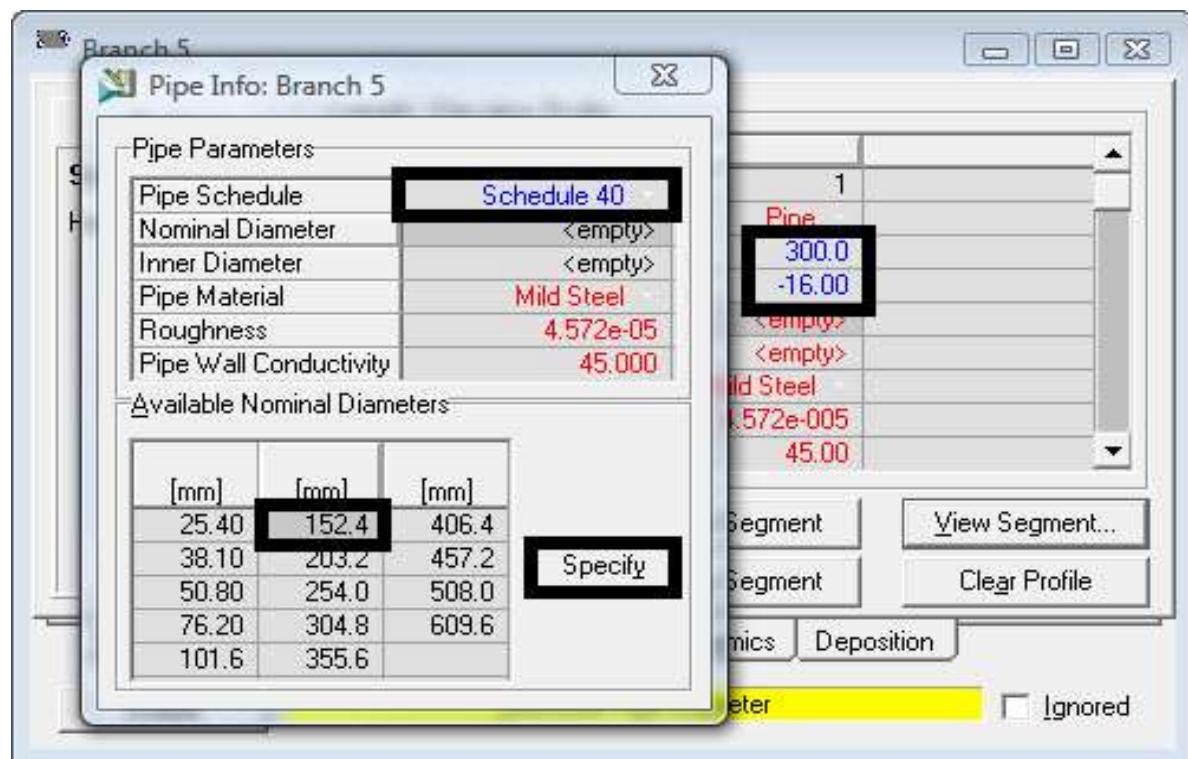
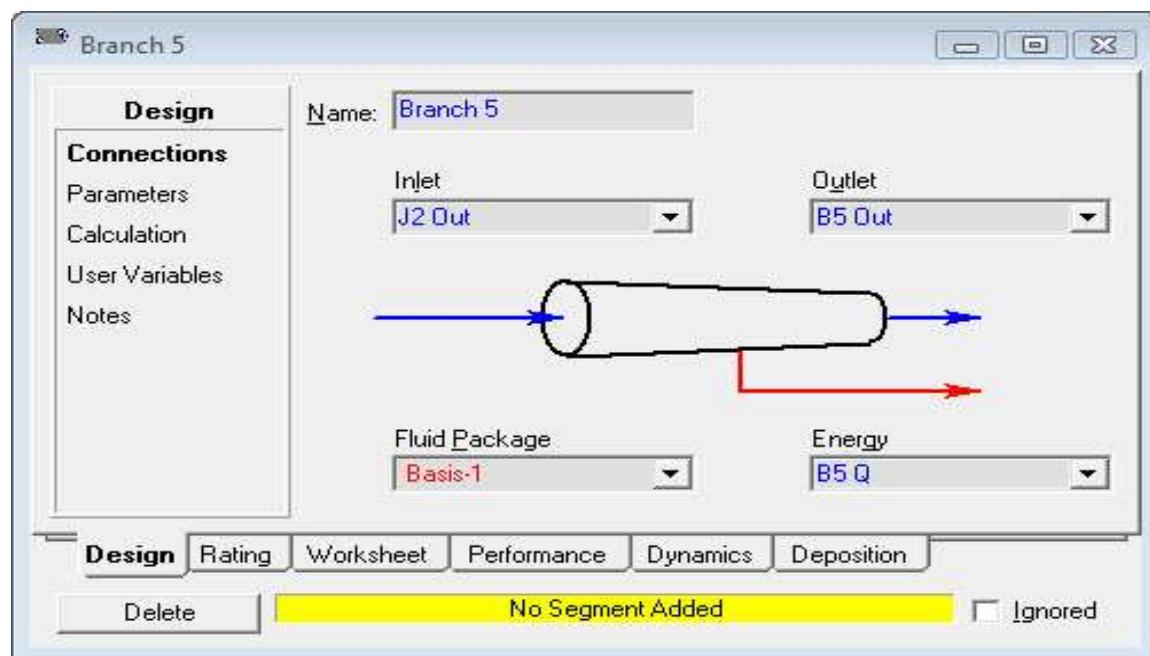
Now, add a mixer operation to mix the outlet streams from both branch 1 & branch 2

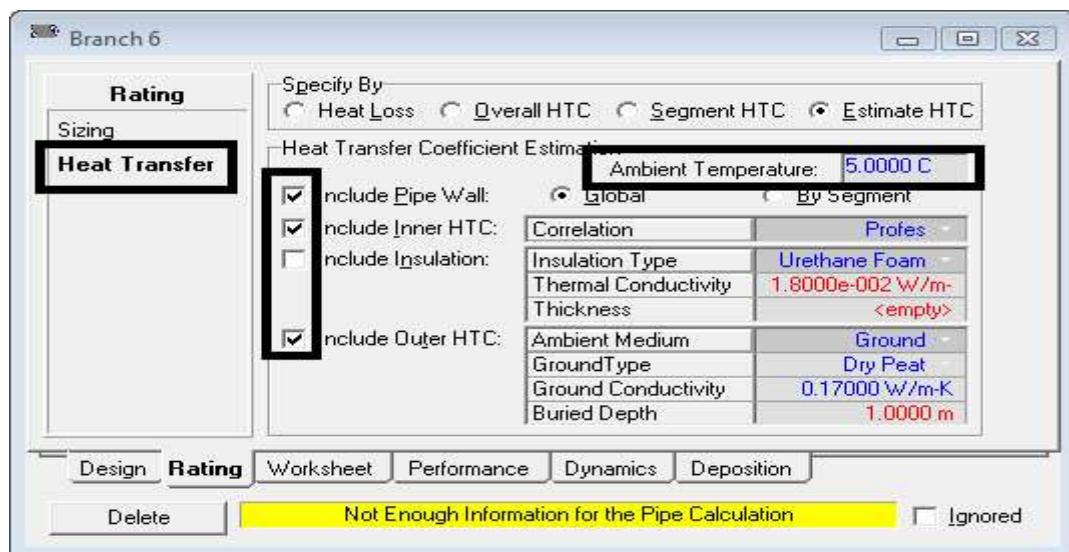
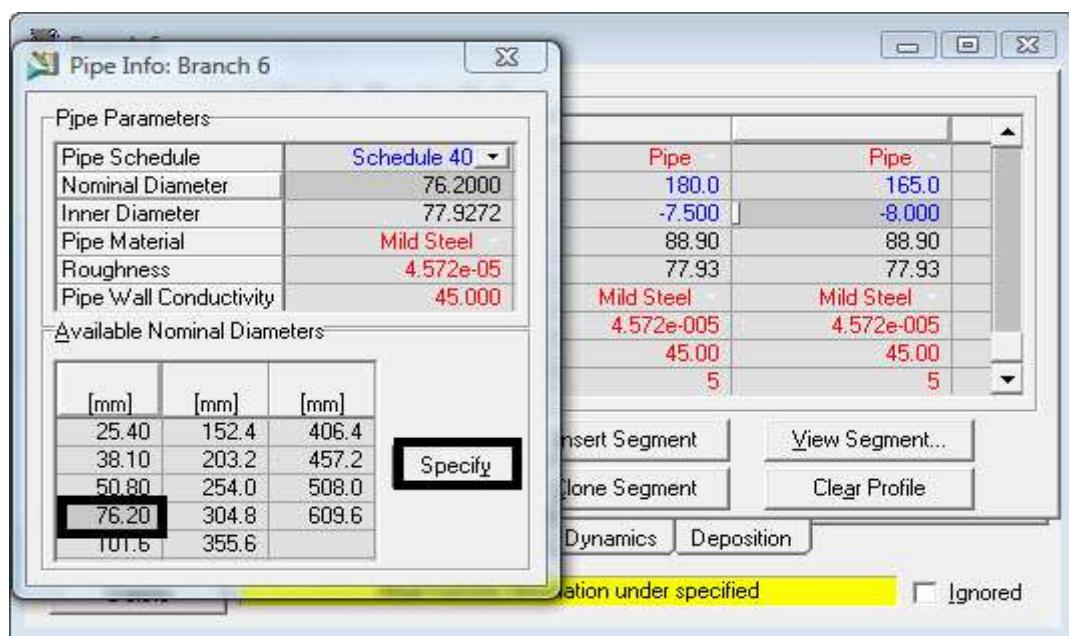
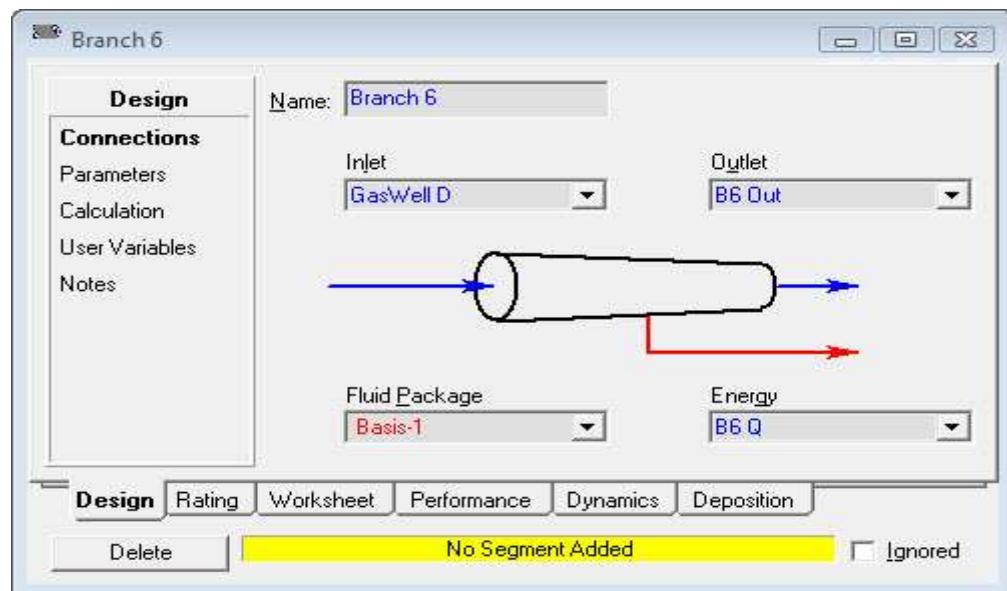




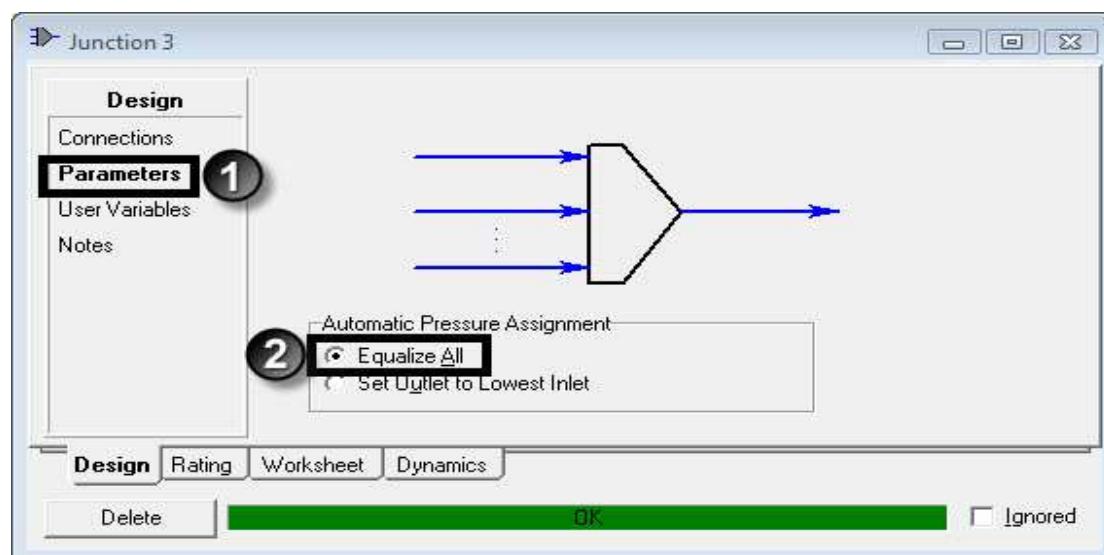
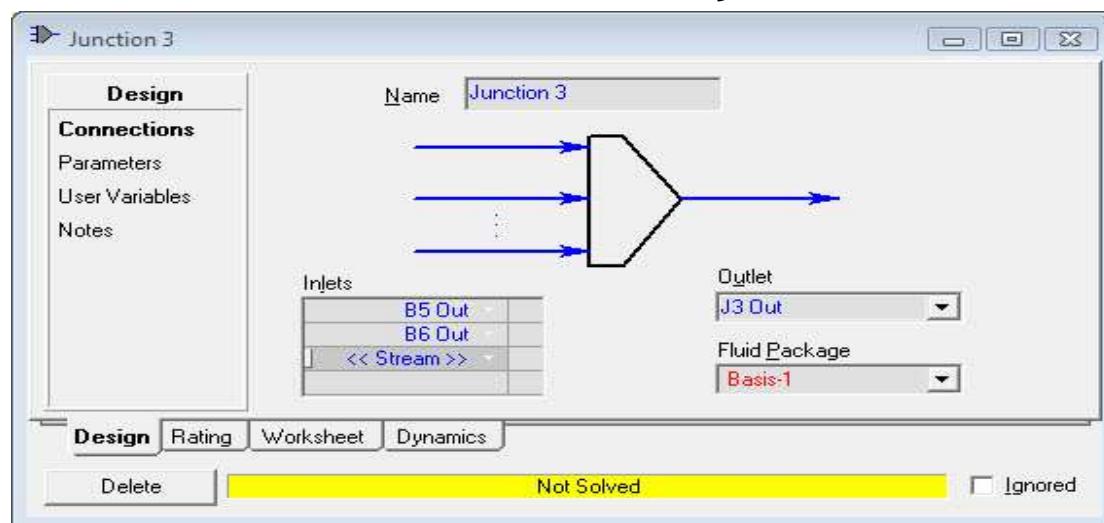




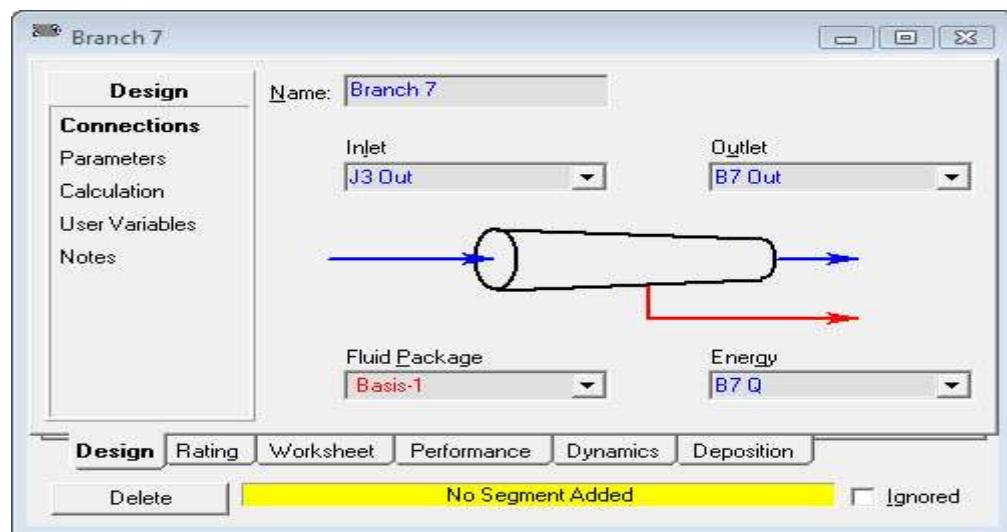


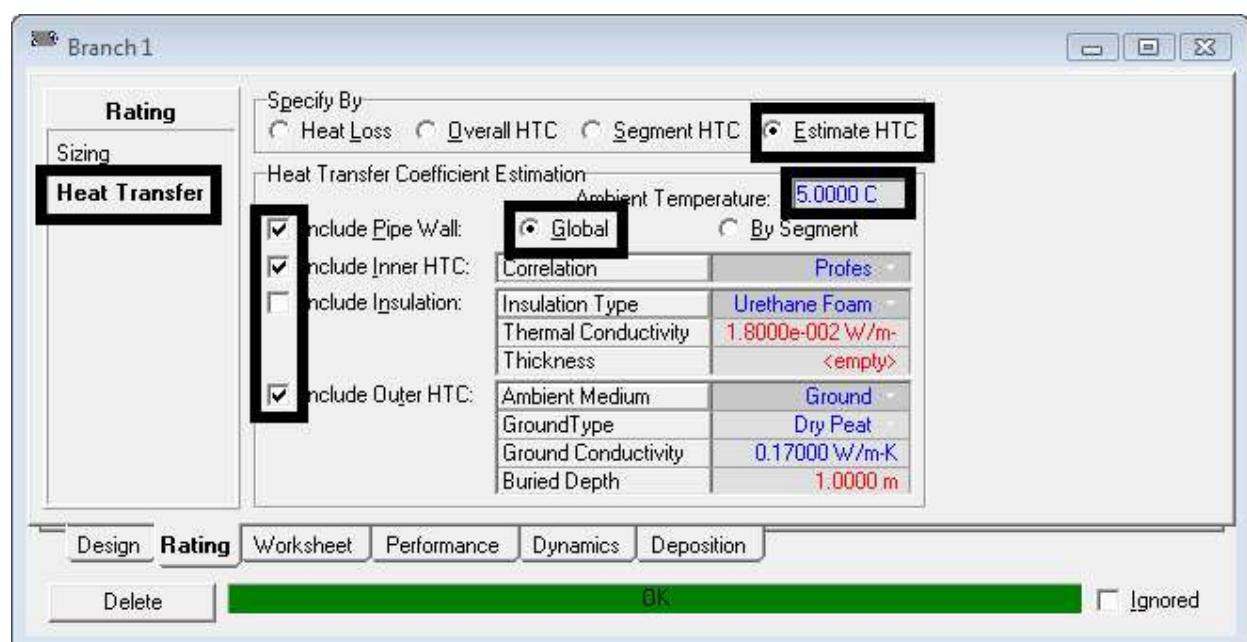
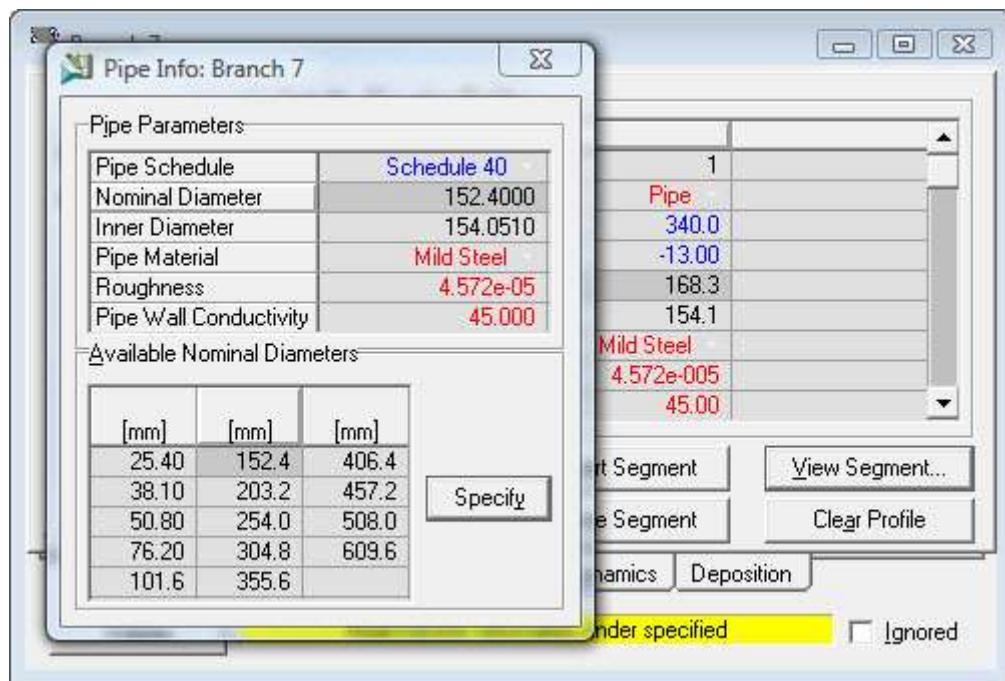


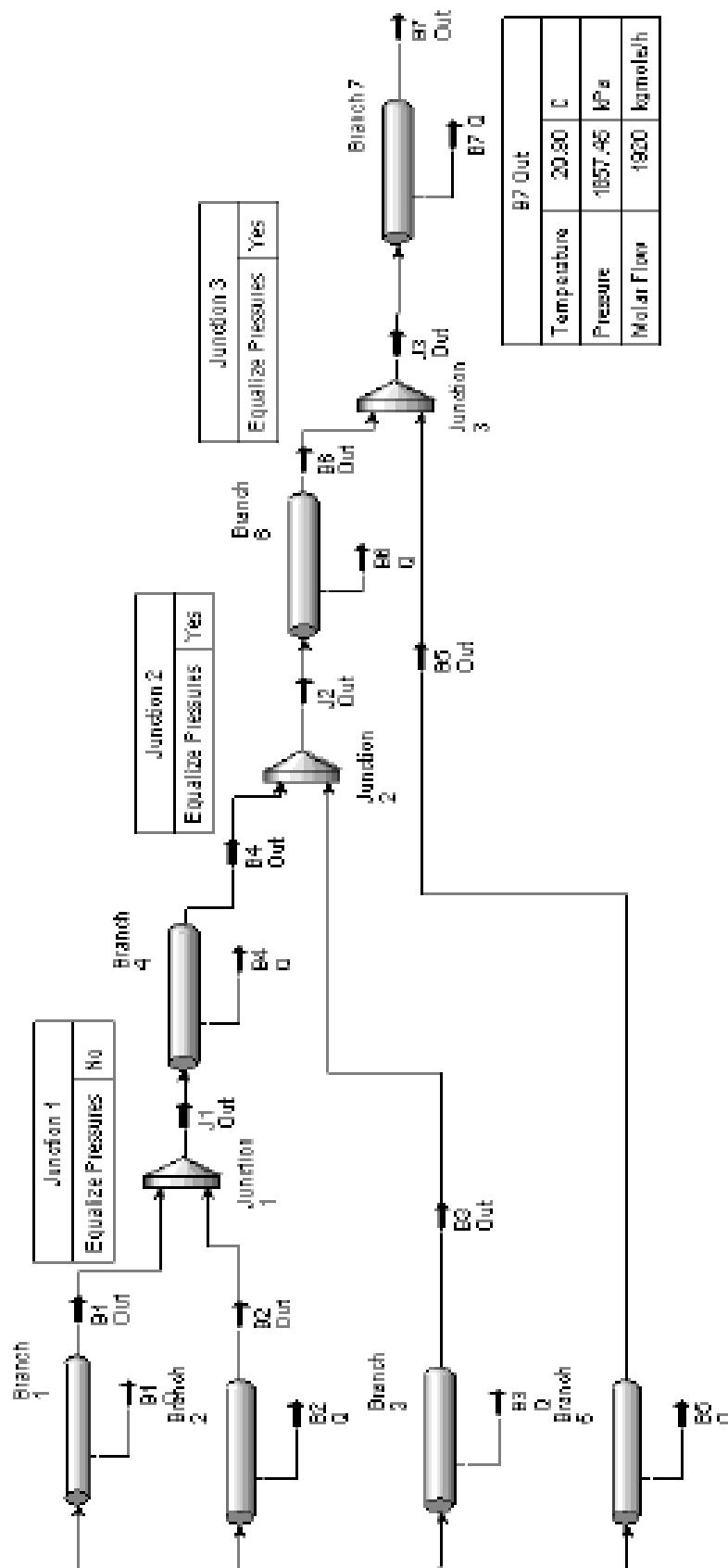
Mix the outlet from both branch 5 & branch 6



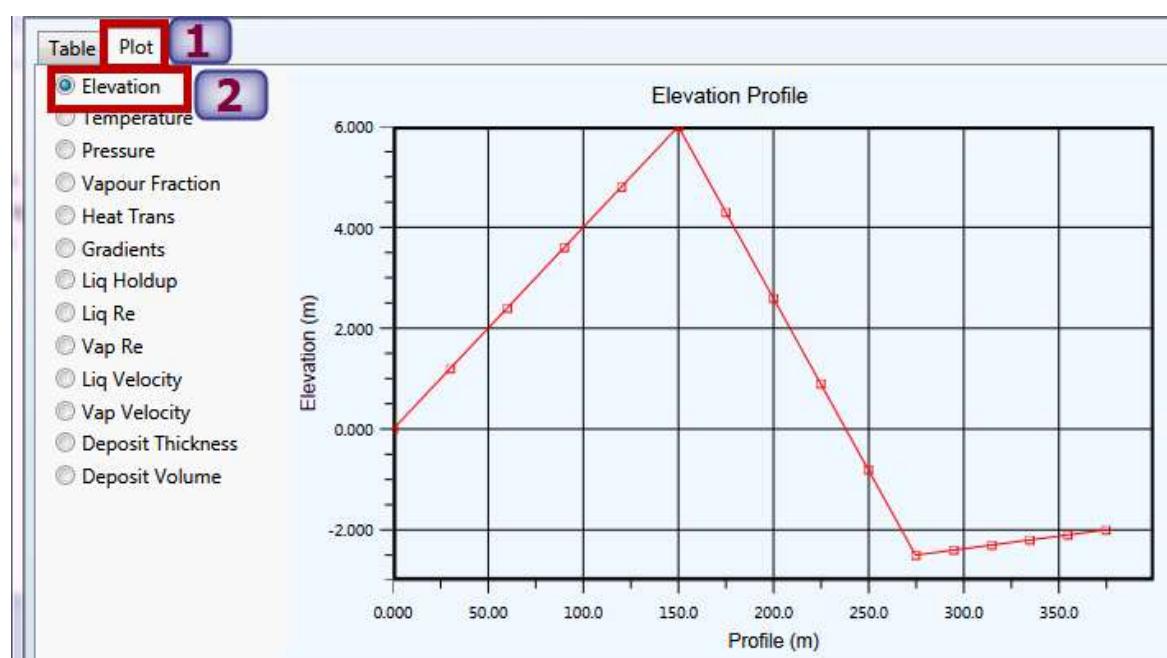
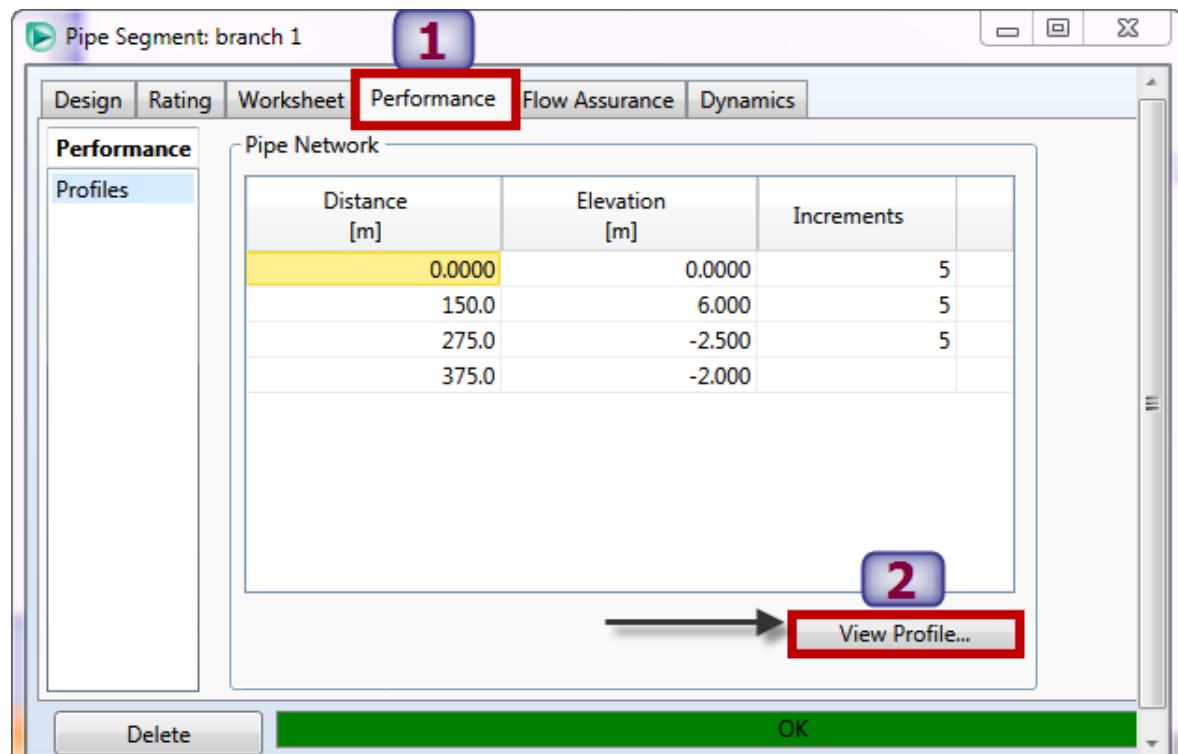
The last branch

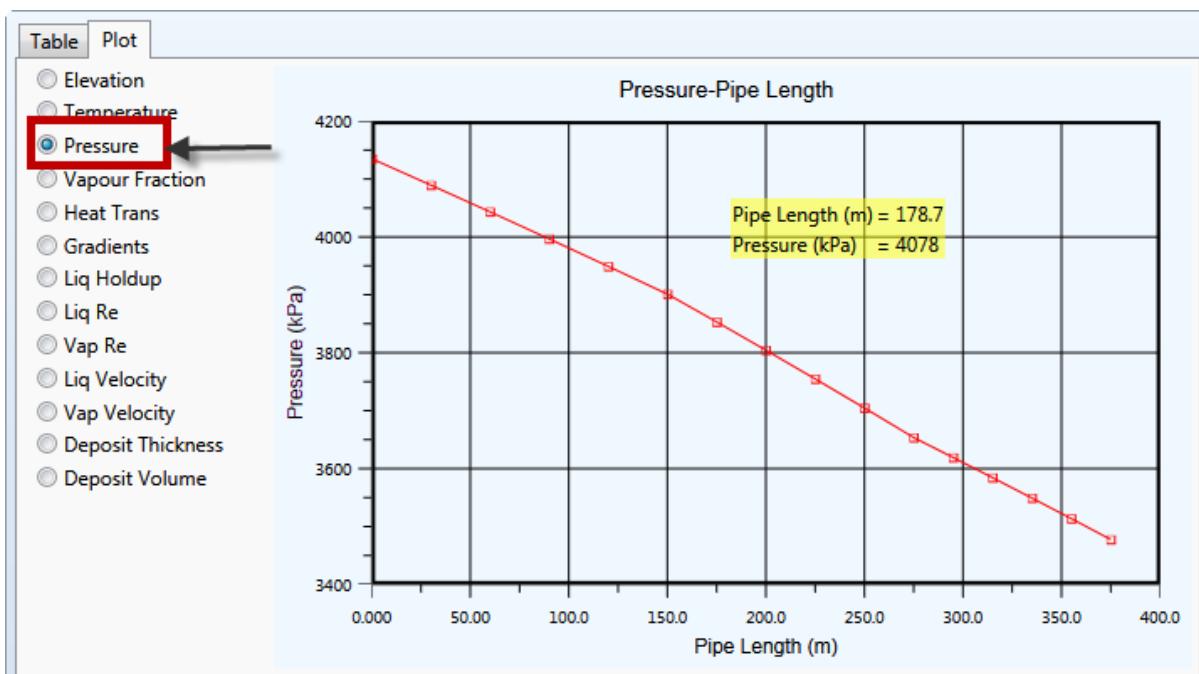
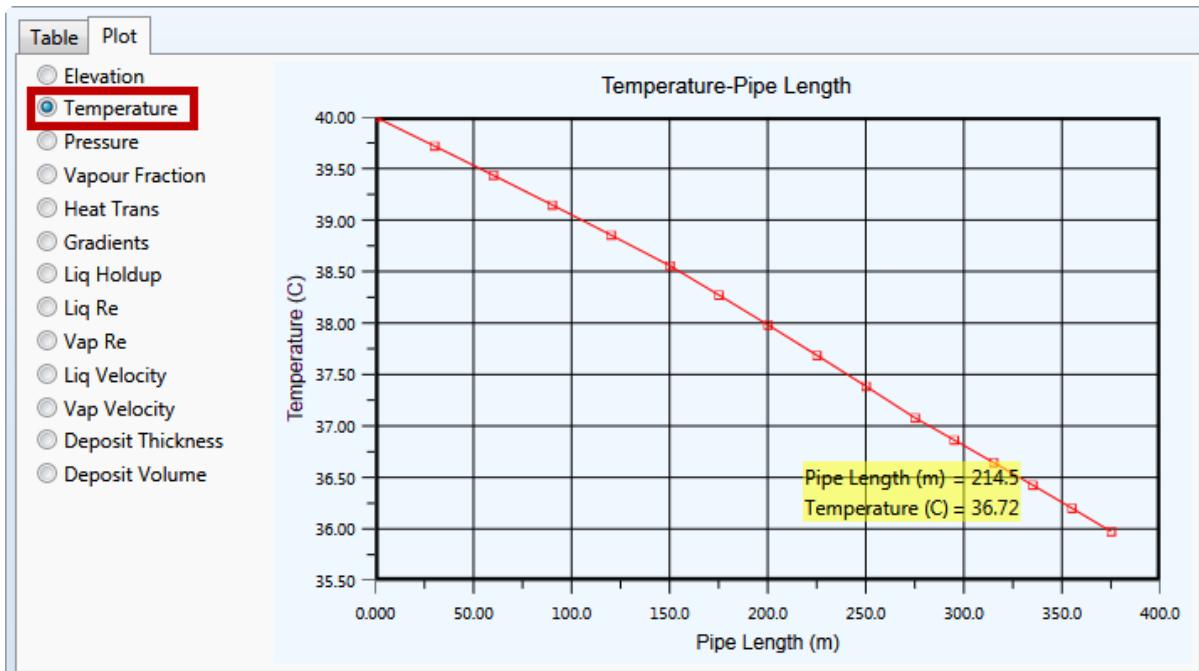




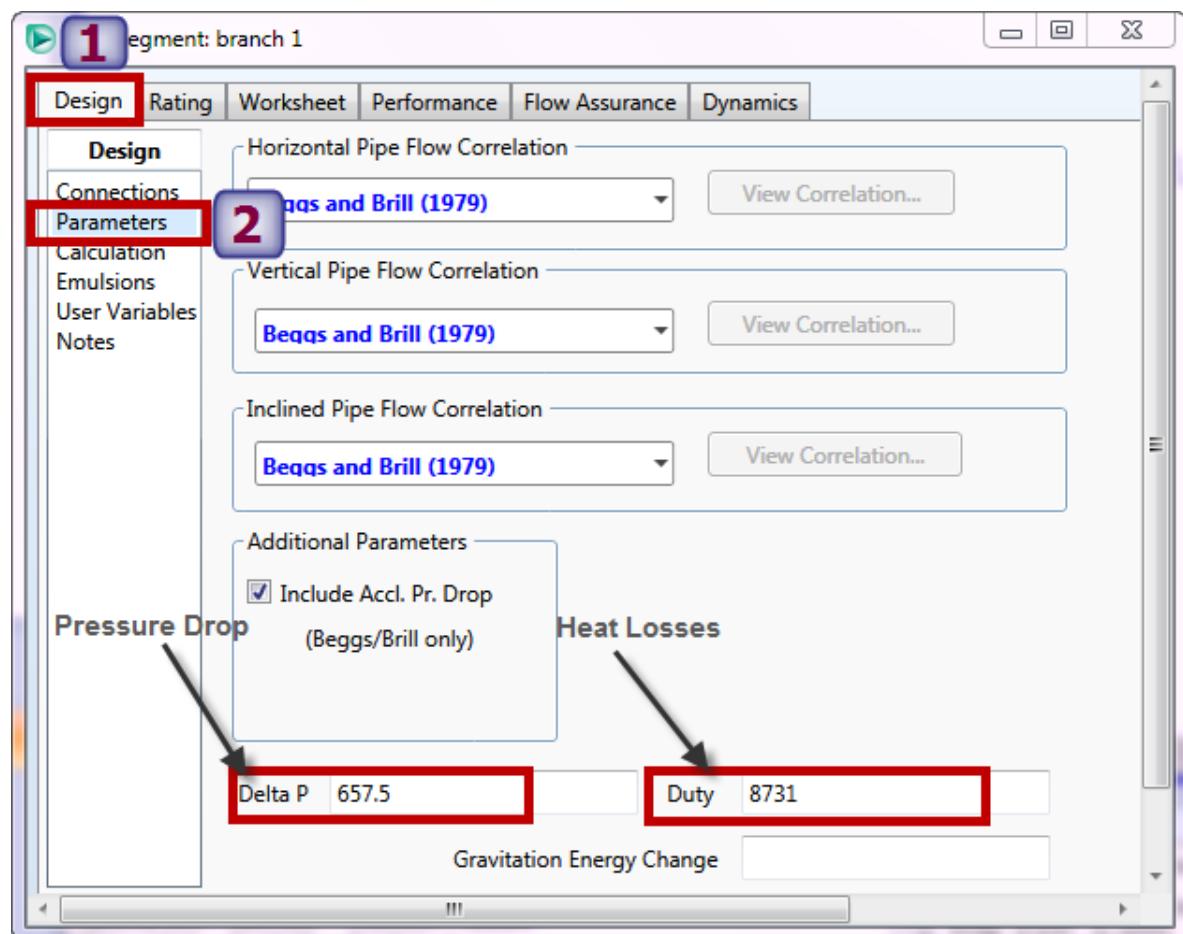


You can go to the performance tab inside each segment to view the profile





You can also go to the parameters page to see the pressure drop & heat losses inside each branch:



# Save Your Case!

# NGL Fractionation

Experience the New Aspen HYSYS®.

V8

The best process simulation software is now easier to use and faster to learn than ever!

Aspen HYSYS is a comprehensive process modeling system used by the world's leading oil & gas producers, refineries, and engineering companies to optimize process design and operations.



9

## Workshop

*Recovery of natural-gas liquids (NGL) from natural gas is quite common in natural gas processing. Recovery is usually done to:*

- *Produce transportable gas (free from heavier hydrocarbons which may condense in the pipeline).*
- *Meet a sales gas specification.*
- *Maximize liquid recovery (when liquid products are more valuable than gas).*

*Aspen HYSYS can model a wide range of different column configurations.*

*In this simulation, an NGL Plant will be constructed, consisting of three columns:*

- *De-Methanizer (operated and modeled as a Reboiled Absorber column)*
- *De-Ethanizer (Distillation column)*
- *De-Propanizer (Distillation column)*

## Learning Objectives

*Once you have completed this section, you will be able to:*

- *Add columns using the Input Experts.*
- *Add extra specifications to columns.*

## NGL Fractionation Train

It's required to process a crude natural gas to remove the heavier hydrocarbons from it thus the composition of it would be suitable for transportation by passing the crude gas which comes from two different wells to three towers: De-Methanizer, De-Ethanizer and De-Propanizer.

The first well at (-140 °F, 330 psia and flow rate of 3575 lbmol/hr) have the following conditions and compositions: (**Fluid Package: Peng Robinson**)

Component	Mol frac	Component	Mol frac
N <sub>2</sub>	0.0025	n-C4	0.0085
CO <sub>2</sub>	0.0048	i-C5	0.0036
C1	0.7041	n-C5	0.002
C2	0.1921	n-C6	0.0003
C3	0.0706	n-C7	0.0002
<b>i-C4</b>	<b>0.0112</b>	n-C8	0.0001

The second one at (-120 °F, 332 psia and flow rate of 475 lbmol/hr) have the following conditions and composition

Component	Mol frac	Component	Mol frac
N <sub>2</sub>	0.0057	n-C4	0.0197
CO <sub>2</sub>	0.0029	i-C5	0.0147
C1	0.7227	n-C5	0.0102
C2	0.1176	n-C6	0.0037
C3	0.075	n-C7	0.0047
<b>i-C4</b>	<b>0.0204</b>	n-C8	0.0027

There's also an energy supplied to the De-Methanizer of (2e6 BTU/hr) which is used to improve the efficiency of the separation.

The De-Methanizer (reboiled absorber) has the following specifications:

- Feed 1 Material Stream enters the column from the top stage inlet
- Feed 2 Material stream enters from the 2nd stage
- Ex-duty Energy Stream enters from the 4th stage
- Number of stages = 10
- Top Stage pressure = 330 psia & Reboiler pressure = 335 psia
- Top Stage temperature = -125°F & Reboiler temperature = 80 °F
- Ovhd Prod Rate = 2950 lbmole/hr (1338 Kg/hr)

**After running the column with the above specifications, the design is then modified to ensure that the overhead molar fraction of the methane is 0.96**

Most of methane is removed from the top of the tower and the bottom stream is pumped to 2790 kpa then it enters to the De-Ethanizer (distillation Column) where most of the ethane in the crude is taken as an overhead product then it is processed to use in different applications.

The de-ethanizer has the following specifications:

<b>Connections</b>	
Name	DC2
No. of Stages	14
Inlet Stream/Stage	DC2 Feed/6
Condenser Type	Partial
Overhead Vapour Product	DC2 Ovhd
Overhead Liquid Product	DC2 Dist
Bottoms Liquid Outlet	DC2 Btm
Reboiler Duty Energy Stream	DC2 Reb Q
Condenser Duty Energy Stream	DC2 Cond Q
<b>Pressures</b>	
Condenser	2725 kPa (395 psia)
Condenser Delta P	35 kPa (5 psi)
Reboiler	2792 kPa (405 psia)
<b>Temperature Estimates</b>	
Condenser	-4°C (25°F)
Reboiler	95°C (200°F)
<b>Specifications</b>	
Overhead Vapour Rate	320 kgmole/h (700 lbmole/hr)
Distillate Rate	0 kgmole/h
Reflux Ratio	2.5 (Molar)

After running the column with the above specifications, the design is modified to ensure that the bottom stream has ratio between  $C_2/C_3 = 0.01$

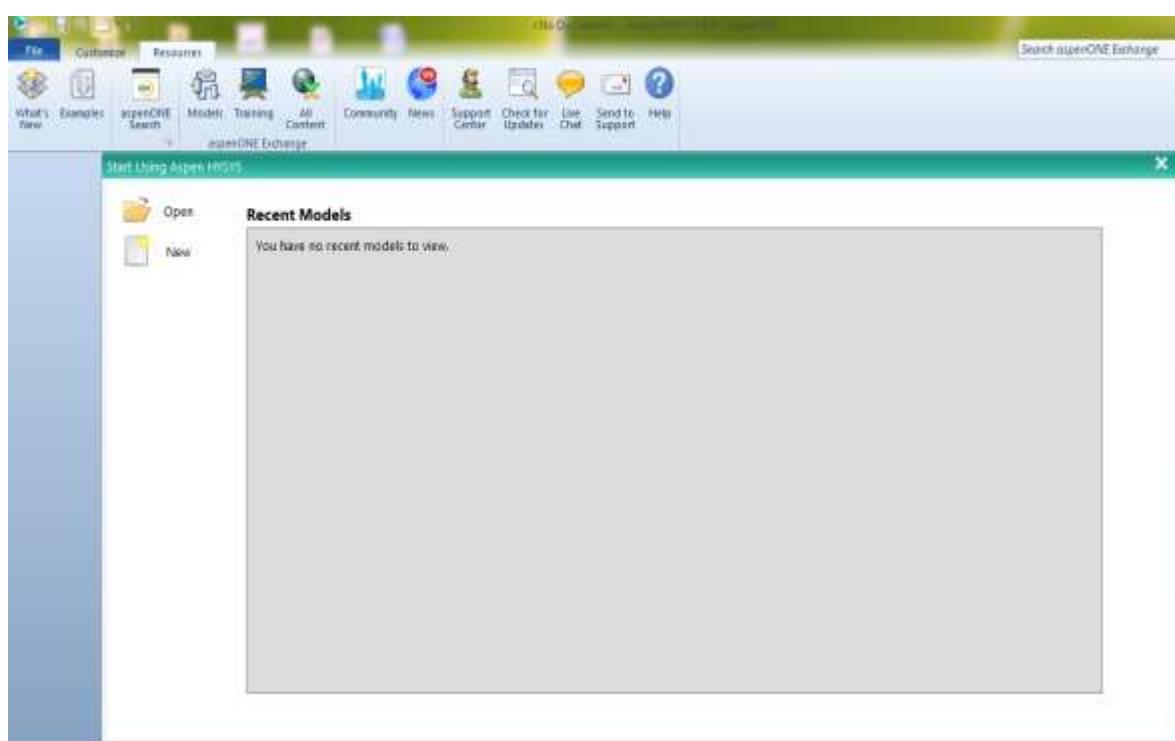
Then the bottom product is sent to a valve where the pressure of it decrease to 1690 kpa, the outlet flow from the valve is sent to a de-propanizer (distillation Column) where most of propane is removed from the top and the heavier hydrocarbons is removed from the bottom

The specifications of the de-propanizer are:

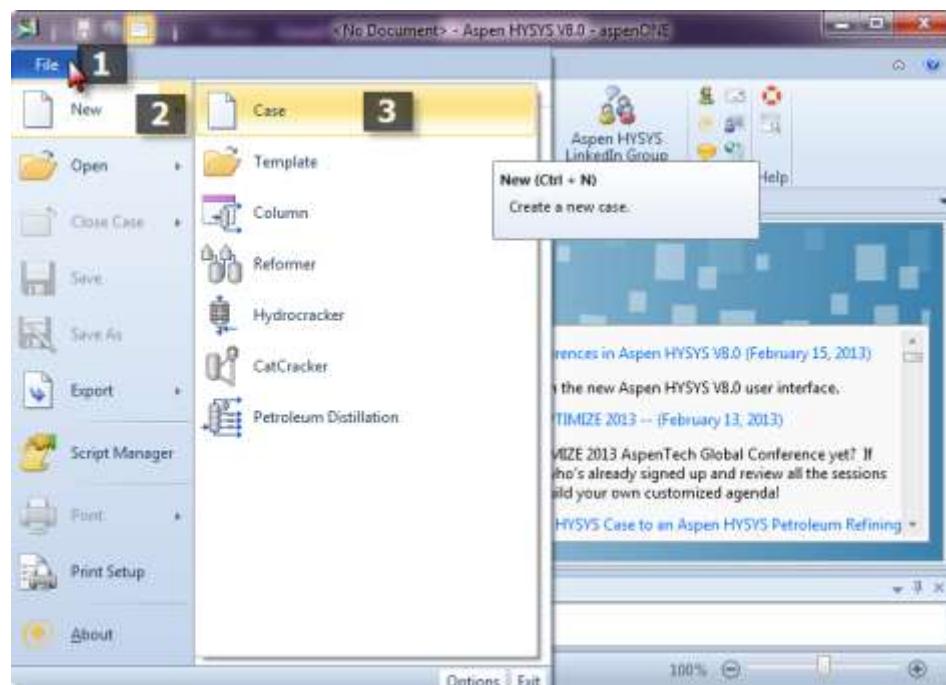
In this cell...	Enter...
<b>Connections</b>	
Name	DC3
No. of Stages	24
Inlet Streams/Stage	DC3 Feed/11
Condenser Type	Total
Ovhd Liquid Outlet	DC3 Dist
Bottom Liquid Outlet	DC3 Btm
Reboiler Duty Energy Stream	DC3 Reb Q
Condenser Duty Energy Stream	DC3 Cond Q
<b>Pressures</b>	
Condenser	1585 kPa (230 psia)
Condenser Delta P	35 kPa (5 psi)
Reboiler	1655 kPa (240 psia)
<b>Temperature Estimates</b>	
Condenser	38°C (100°F)
Reboiler	120°C (250°F)
<b>Specifications</b>	
Liquid Rate	110 kgmole/h (240 lbmole/hr)
Reflux Ratio	1.0 Molar

After running the column with the above specifications, the design is modified to ensure that the overhead molar fraction of the i-C4 & n-C4 =0.15 AND propane bottom product molar fraction = 0.02

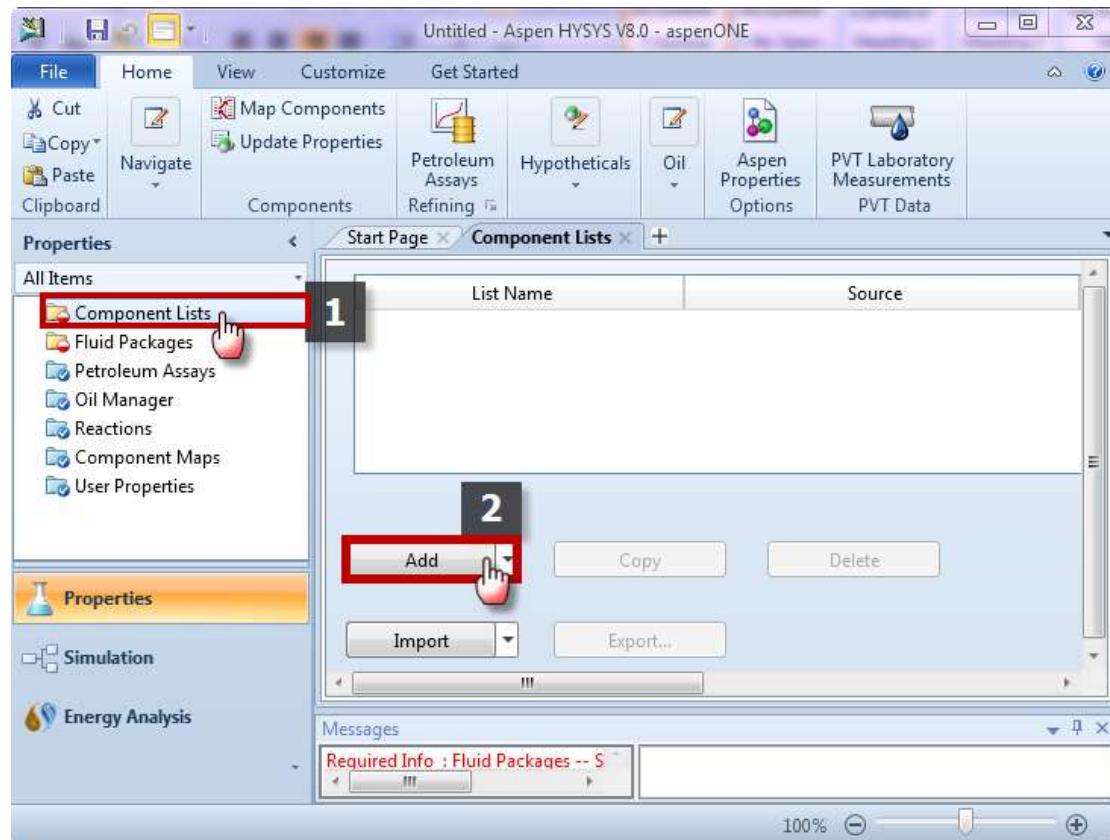
To start the program, From Start Menu, Select All Programs >> Aspen Tech >> Process Modeling V8.x >>> Aspen HYSYS >> Aspen HYSYS



## 11- First, Start a new case



## 12- Add the Components



13- Choose the system components from the databank:

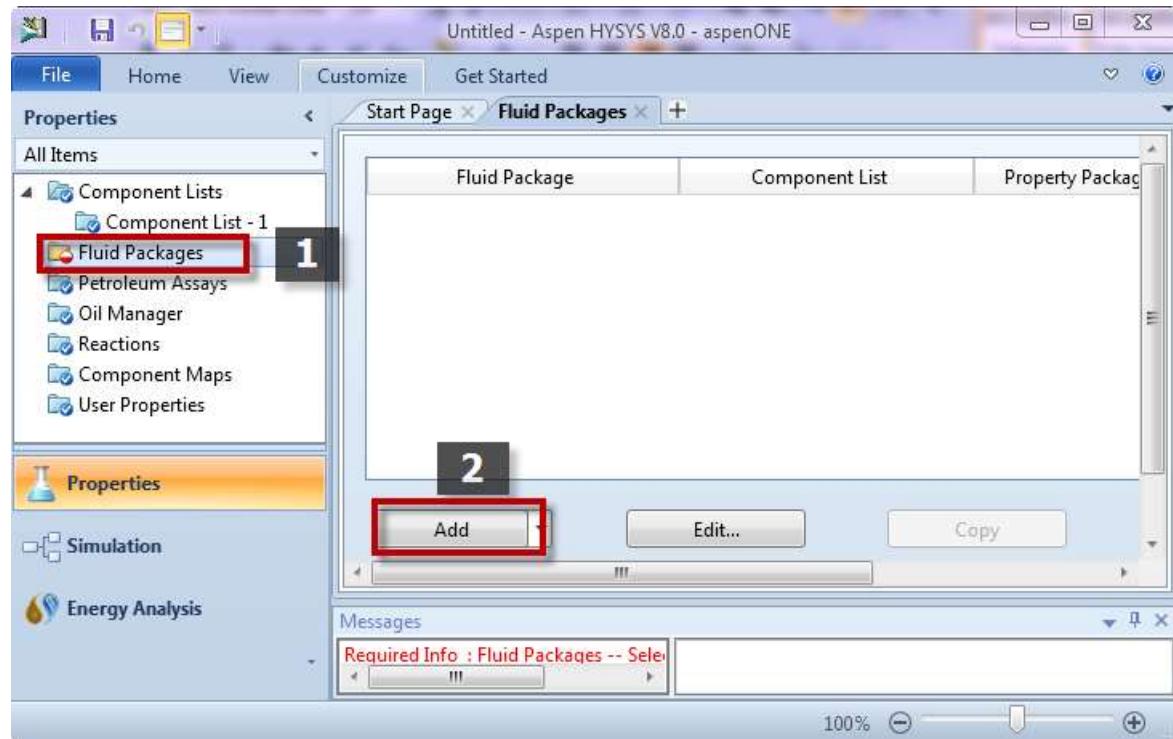
The screenshot shows the 'Component List' dialog in Aspen HYSYS V8. On the left, a table lists selected components: Methane, Propane, and Ethane, all categorized as 'Pure Component'. On the right, a larger table lists various hydrocarbons with their simulation names, full names/synonyms, and formulas. The first row, 'Methane', is highlighted with a yellow background. A red box highlights the '< Add' button, which is labeled with the number '2'. Another red box highlights the 'i-Butane' entry in the list, which is labeled with the number '1'.

Component	Type	Group
Methane	Pure Component	
Propane	Pure Component	
Ethane	Pure Component	

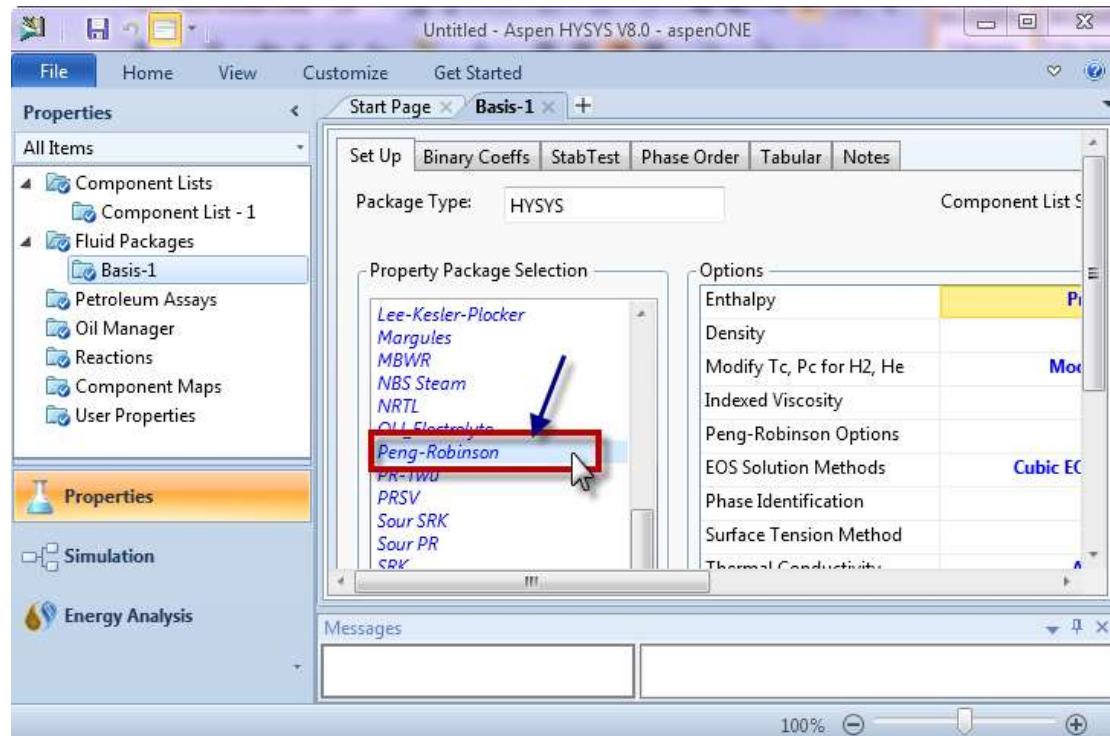
Simulation Name	Full Name / Synonym	Formula
Methane	C1	CH <sub>4</sub>
Ethane	C2	C <sub>2</sub> H <sub>6</sub>
Propane	C3	C <sub>3</sub> H <sub>8</sub>
i-Butane	i-C4	C <sub>4</sub> H <sub>10</sub>
n-Butane	n-C4	C <sub>4</sub> H <sub>10</sub>
i-Pentane	i-C5	C <sub>5</sub> H <sub>12</sub>
n-Pentane	n-C5	C <sub>5</sub> H <sub>12</sub>
n-Hexane	C6	C <sub>6</sub> H <sub>14</sub>
n-Heptane	C7	C <sub>7</sub> H <sub>16</sub>
n-Octane	C8	C <sub>8</sub> H <sub>18</sub>
n-Nonane	C9	C <sub>9</sub> H <sub>20</sub>
n-Decane	C10	C <sub>10</sub> H <sub>22</sub>
n-C11	C11	C <sub>11</sub> H <sub>24</sub>
n-C12	C12	C <sub>12</sub> H <sub>26</sub>

After adding the pure components (N<sub>2</sub>, CO<sub>2</sub>, C1, C2, C3, n-C4, i-C4, n-C5, i-C5, n-C6, C7, C8)

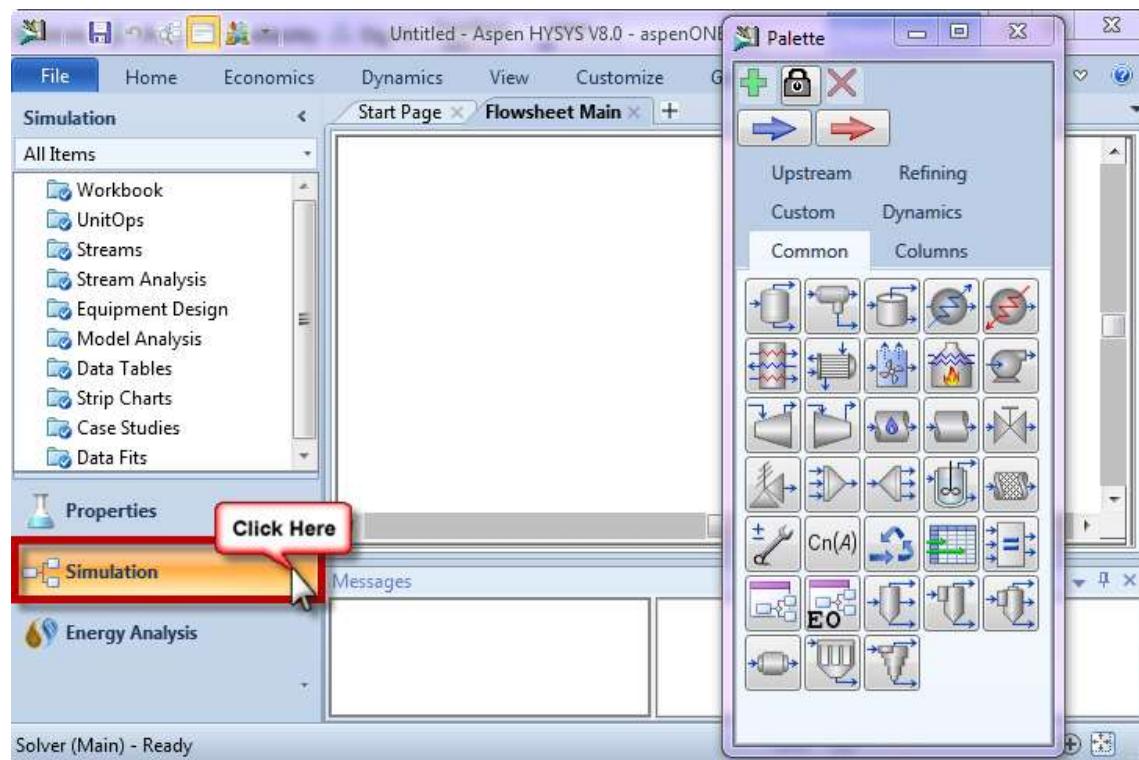
Now, select the suitable fluid package



In this case, select Peng-Robinson

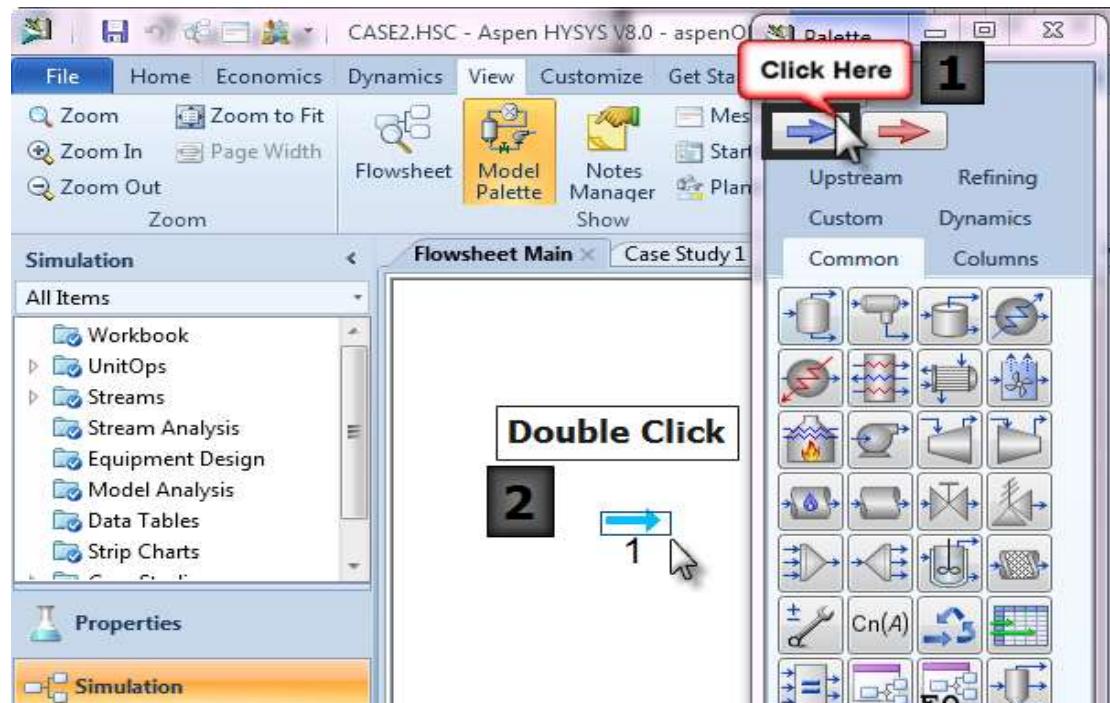


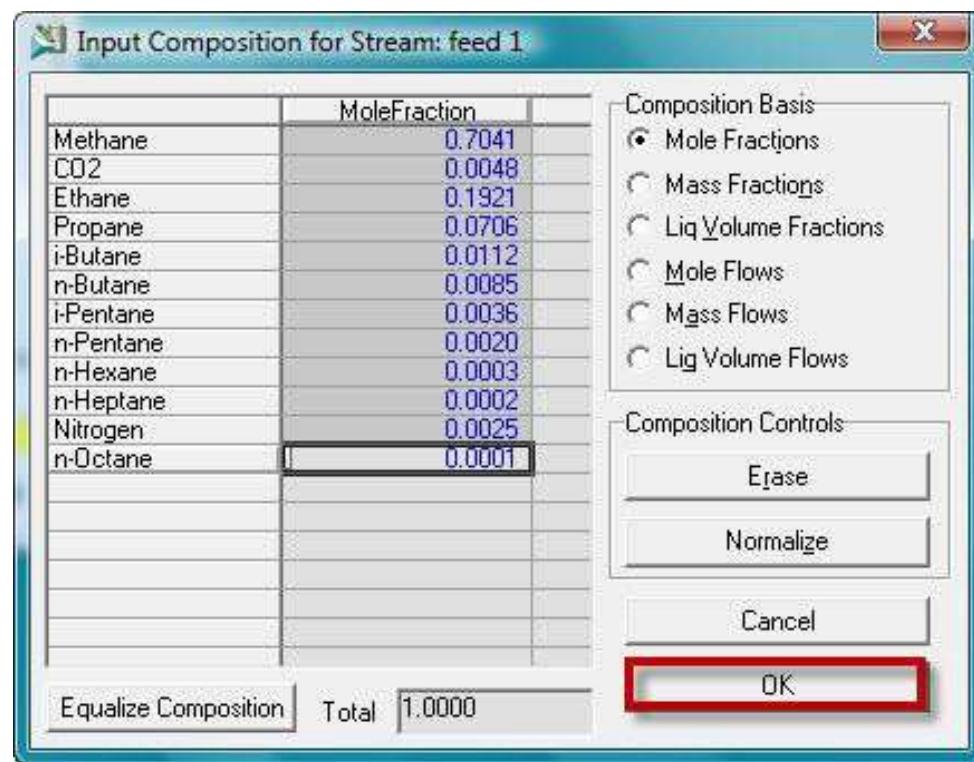
Now you can start drawing the flow sheet for the process by clicking the Simulation button:



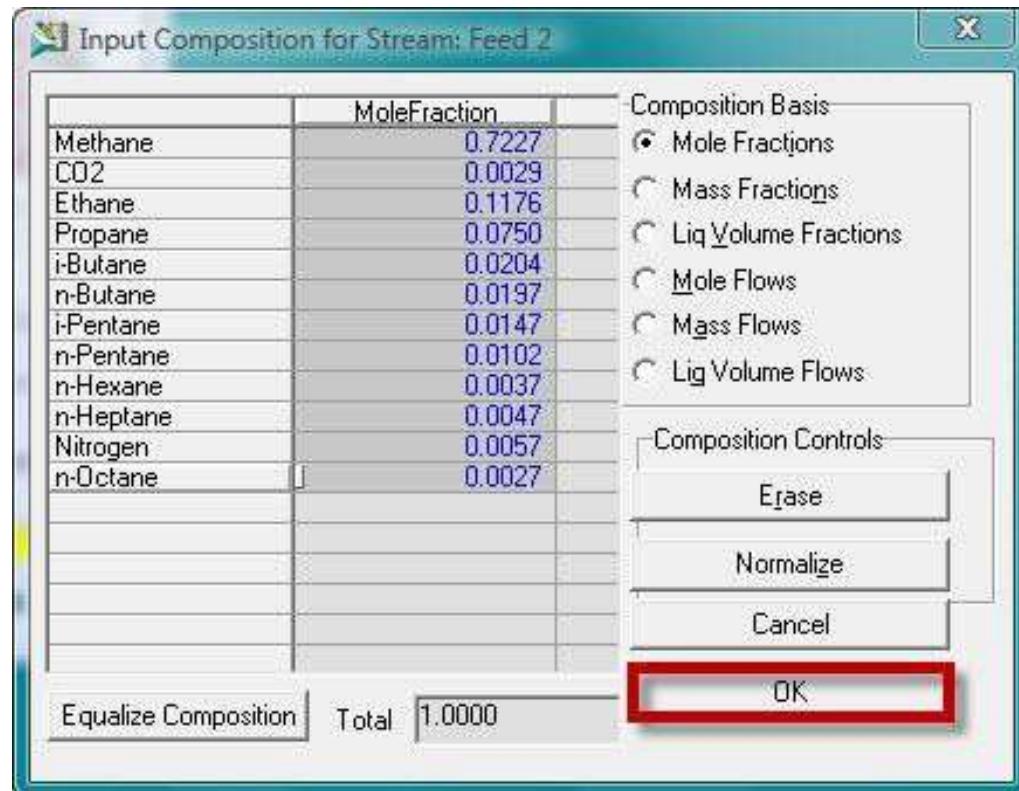
Now add a material stream to define the composition and the conditions of the feed stream

From the palette:





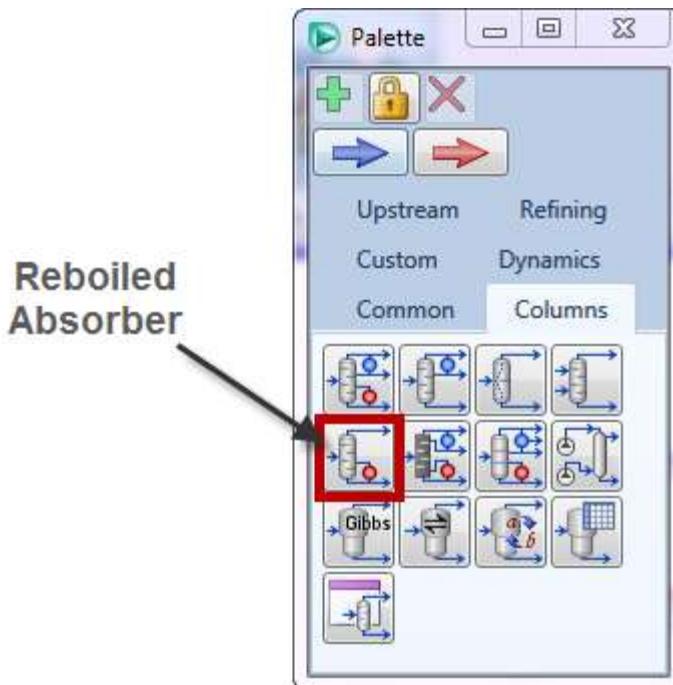
After adding the compositions and the conditions for the first stream, add another stream for feed 2:



## PART 1: DE-METHANIZER COLUMN

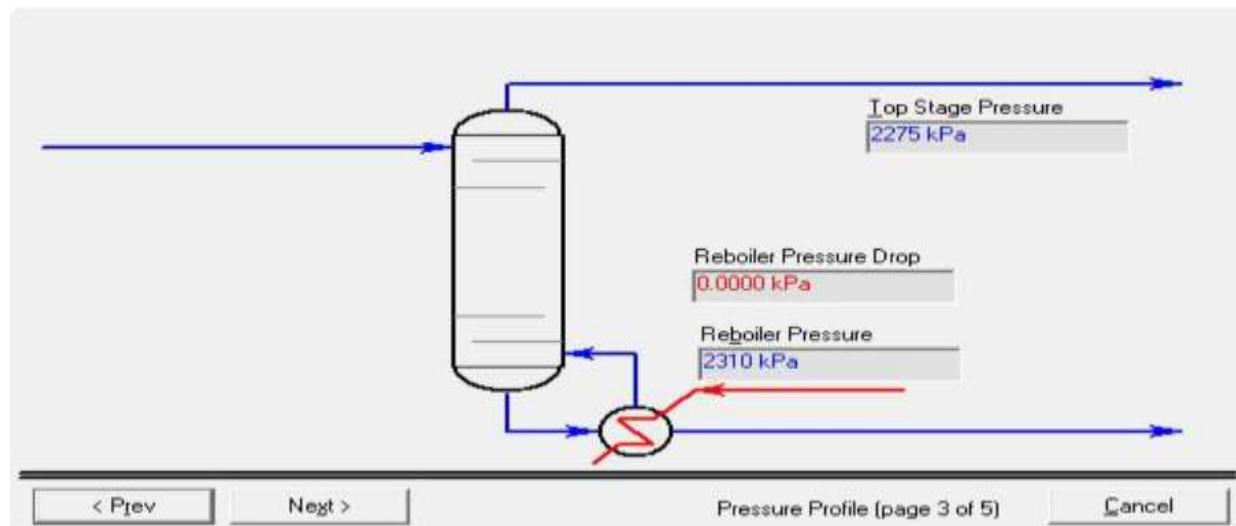
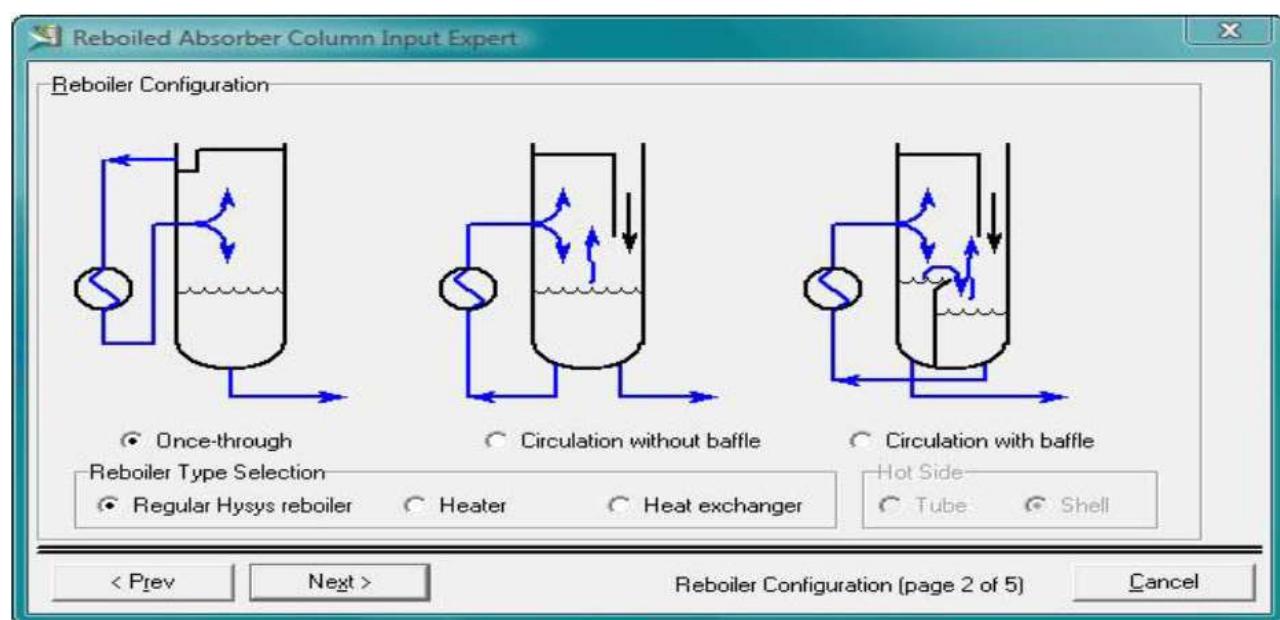
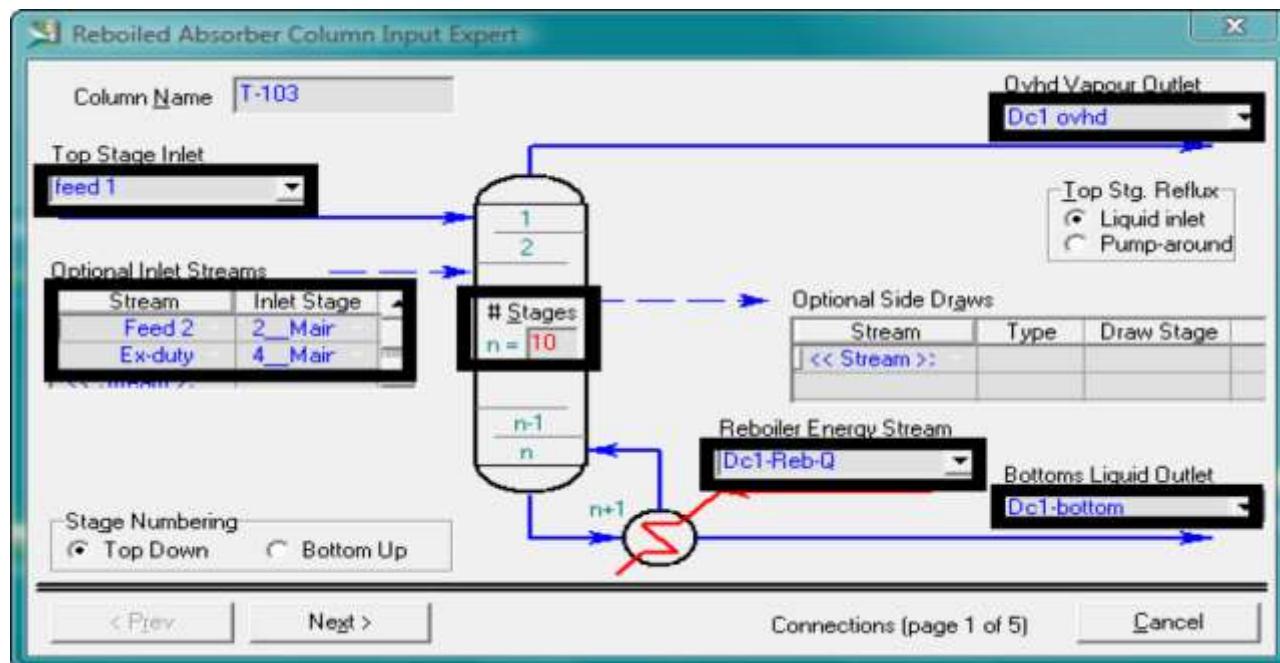
The De-Methanizer is modeled as a reboiled absorber operation, with two feed streams and an energy stream feed, which represents a side heater on the column.

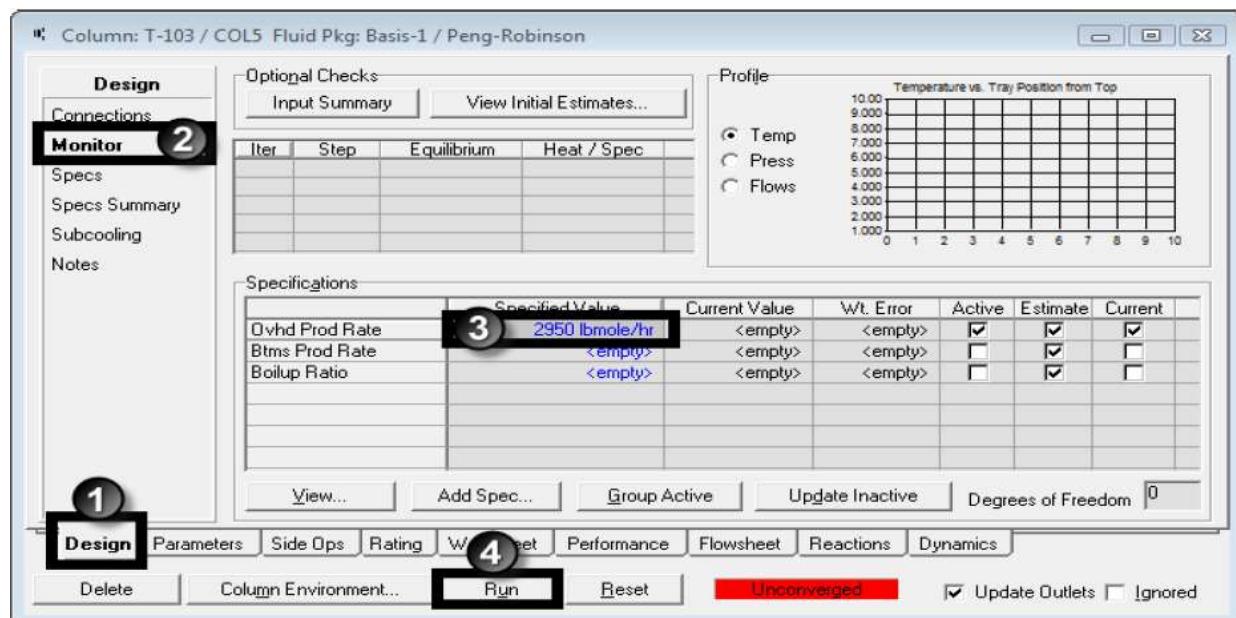
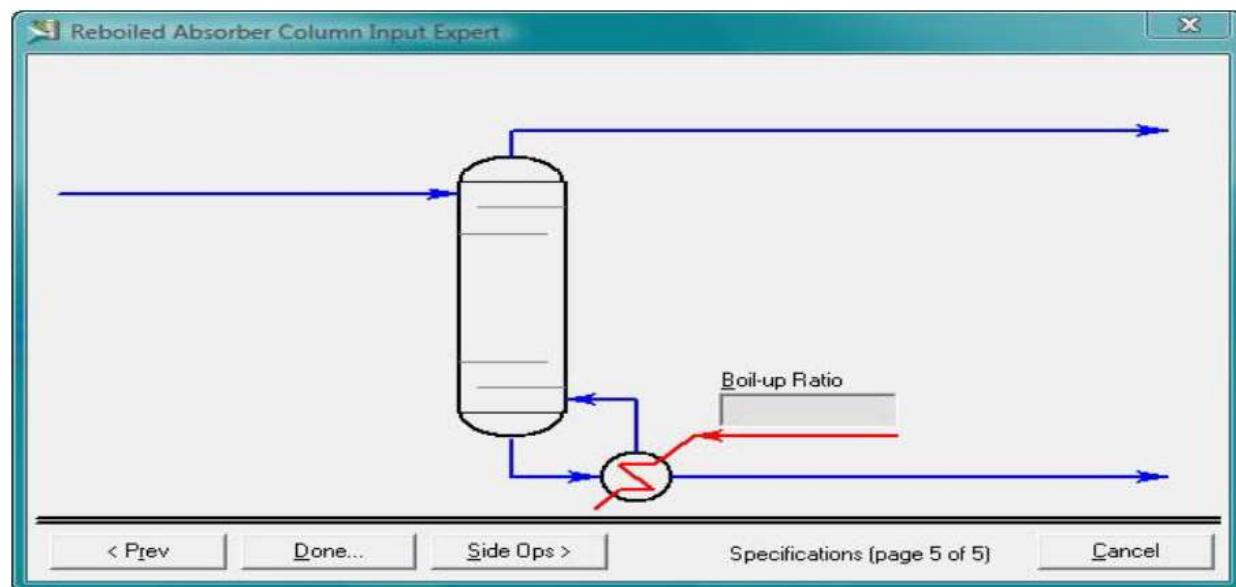
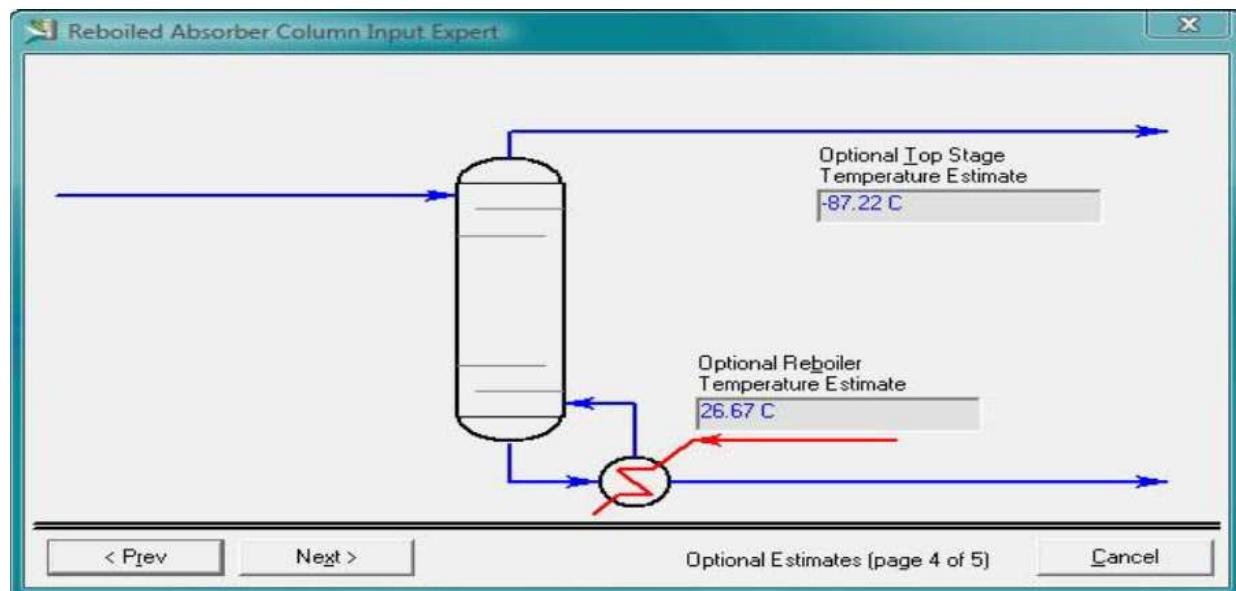
- Add an **Energy** stream with the duty =  $2.1 \text{ e}6 \text{ kJ/hr}$
- Double-click on the **Reboiled Absorber** icon on the Object Palette.



### De-Methanizer Specs:

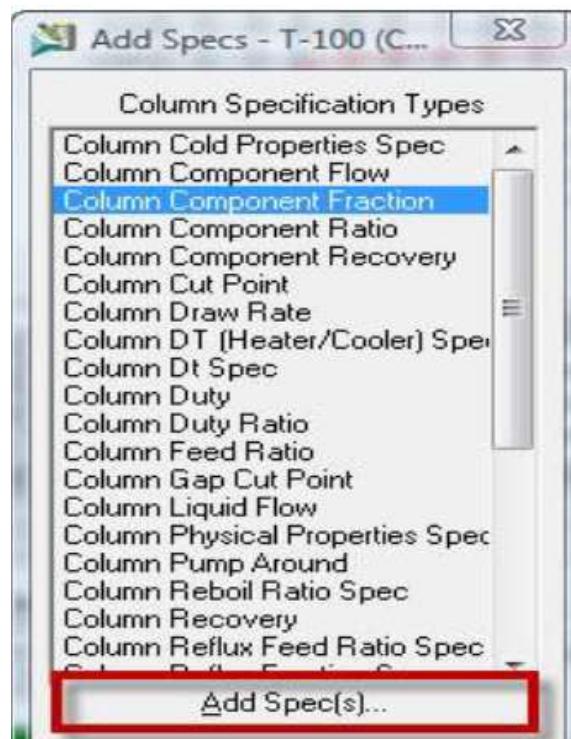
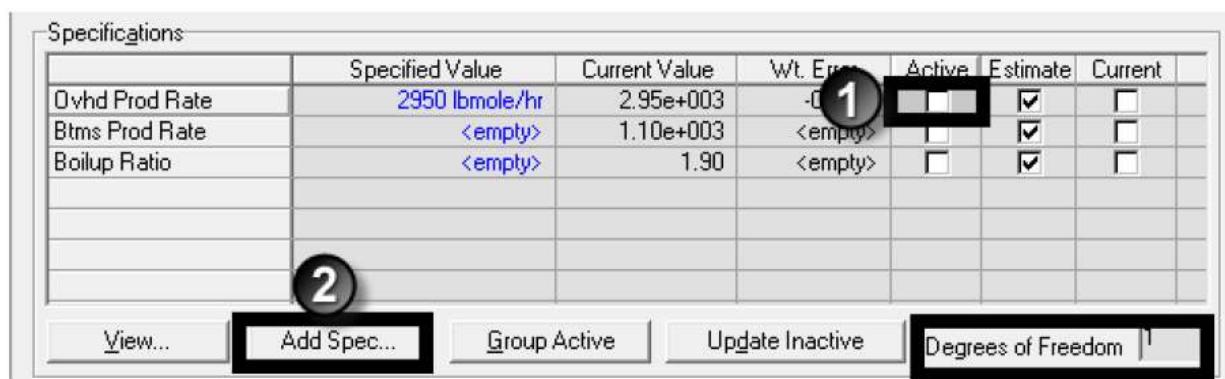
- Feed 1 Material Stream enters the column from the top stage inlet
- Feed 2 Material stream enters from the 2nd stage
- Ex-duty Energy Stream enters from the 4th stage
- Number of stages = 10
- Top Stage pressure = 330 psia
- Reboiler pressure = 335 psia
- Top Stage temperature =  $-125^{\circ}\text{F}$
- Reboiler temperature =  $80^{\circ}\text{F}$
- Ovhd Prod Rate = 2950 lbmole/hr (1338 Kg/hr)
- C1 fraction in the Ovhd stream = 0.96

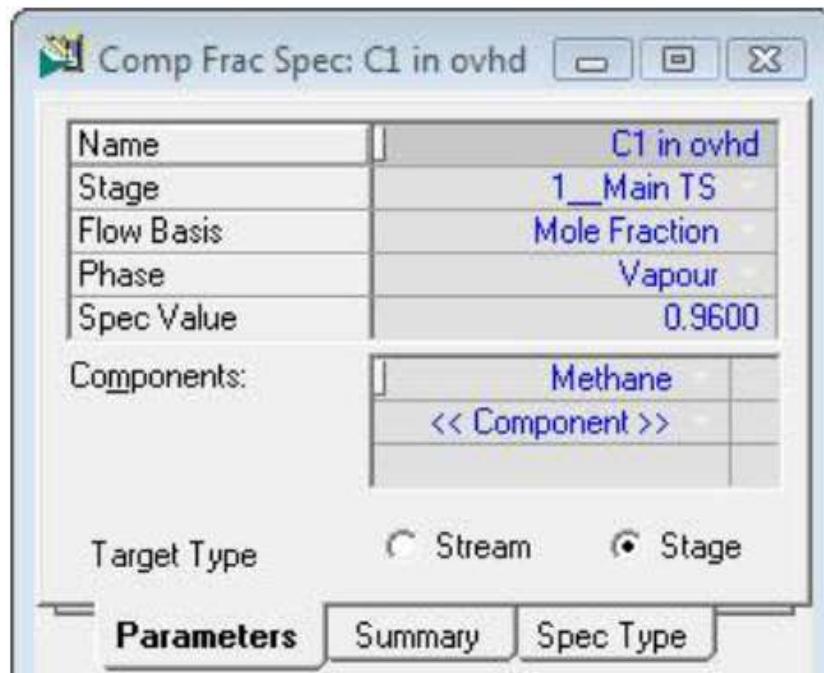




Although the column is converged, it is not always practical to have flow rate specifications. These specifications can result in columns which cannot be converged or that produce product streams with undesirable properties if the column feed conditions change.

An alternative approach is to specify either component fractions or component recoveries for the column product streams.

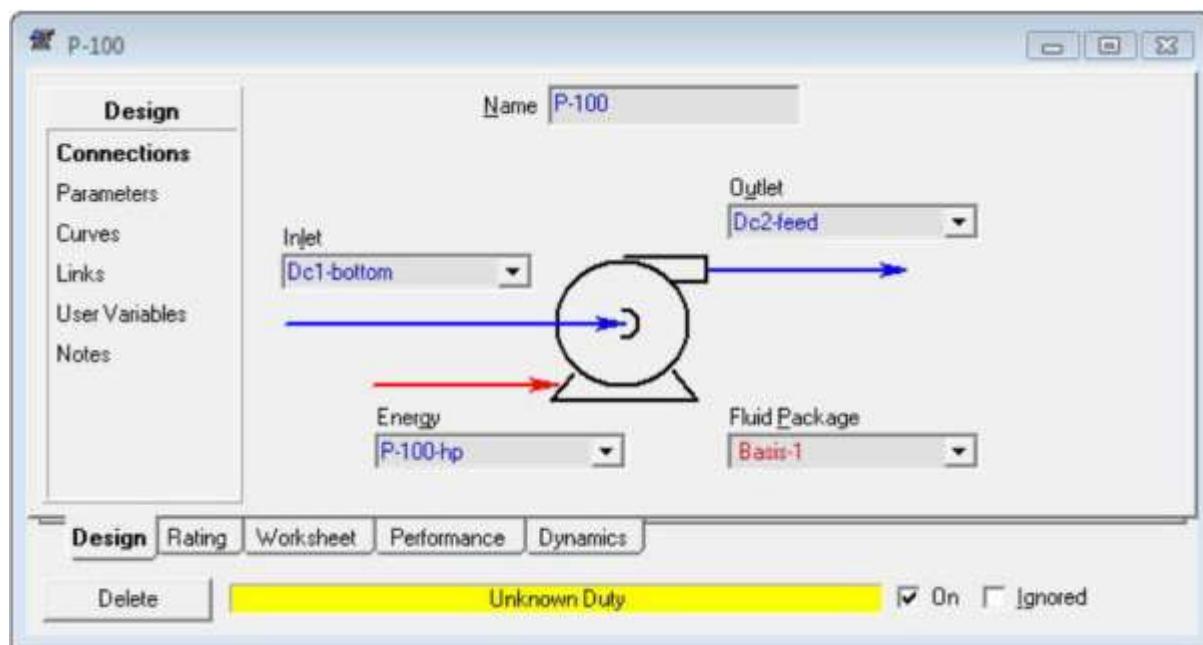




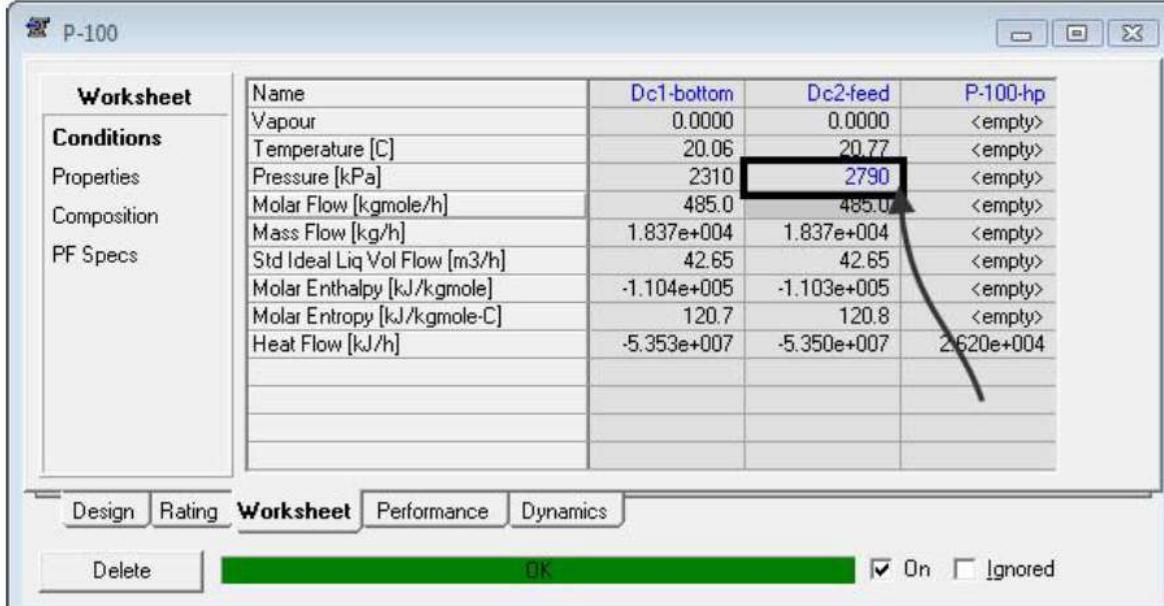
Run the column:



After running the column, add a pump to transfer the bottom liquid to the De-ethanizer:



The pump outlet pressure is 2790 kPa (from Worksheet)

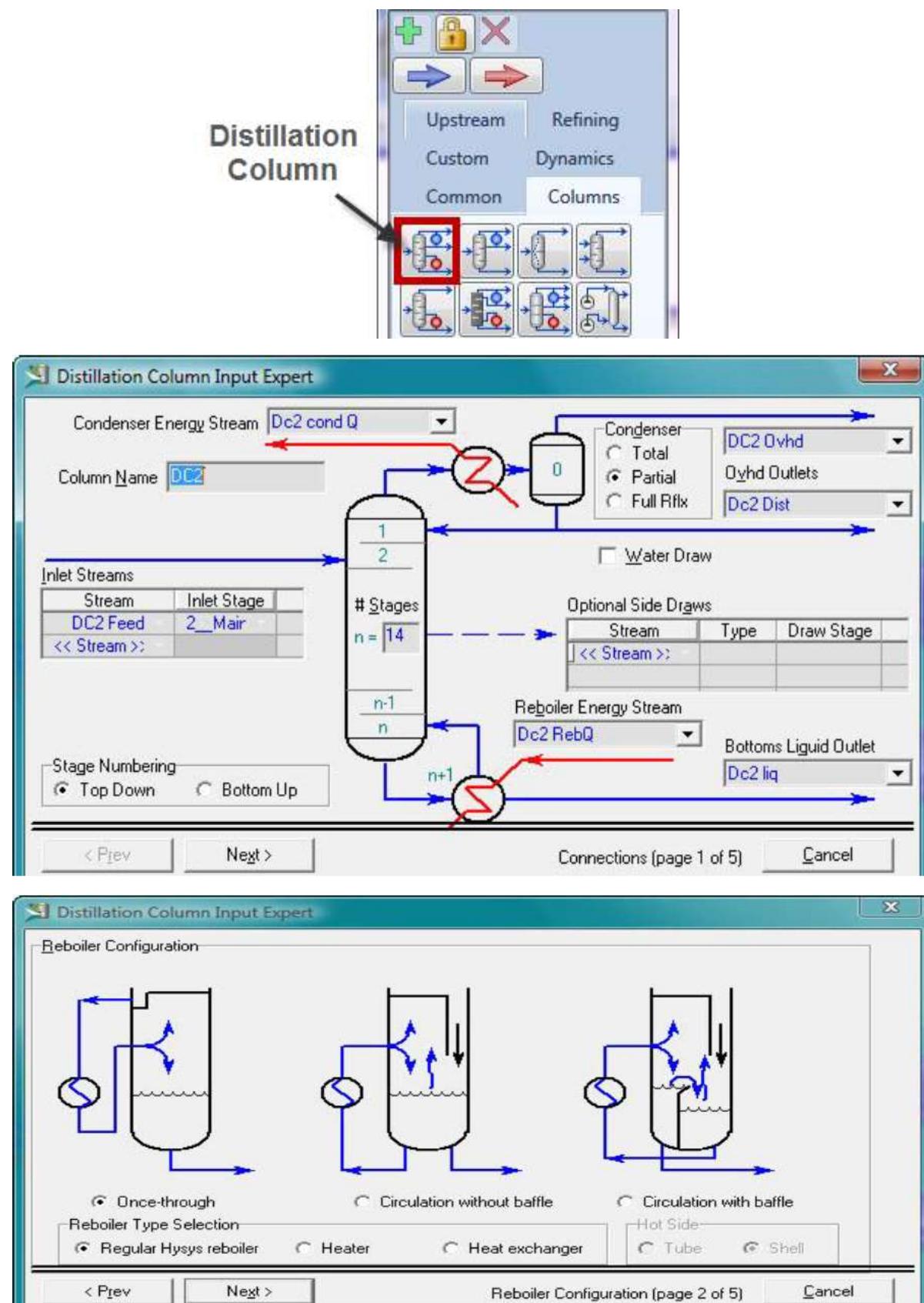


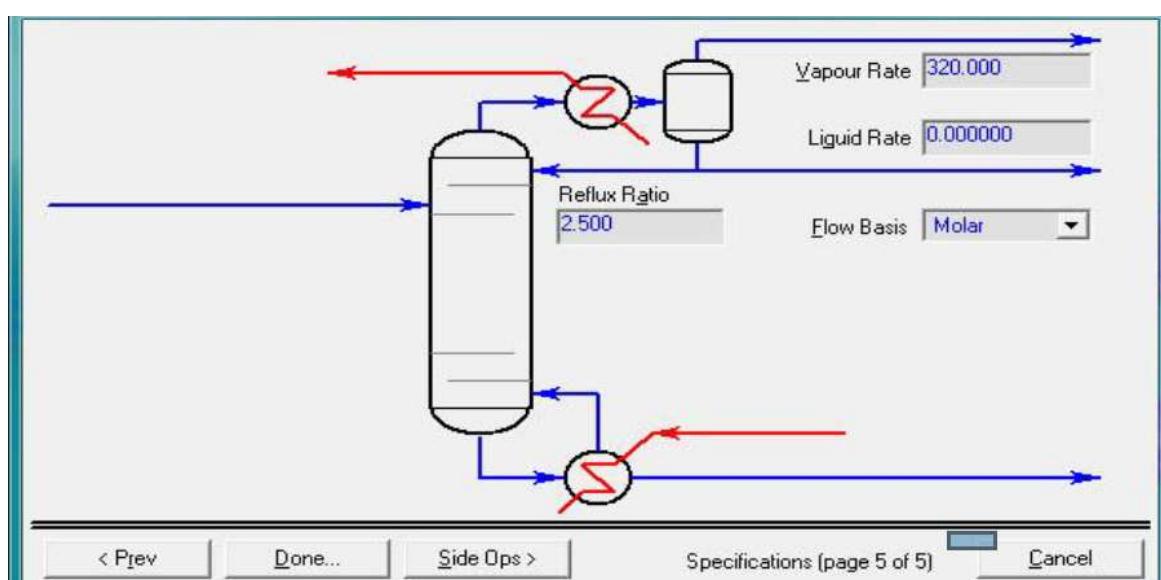
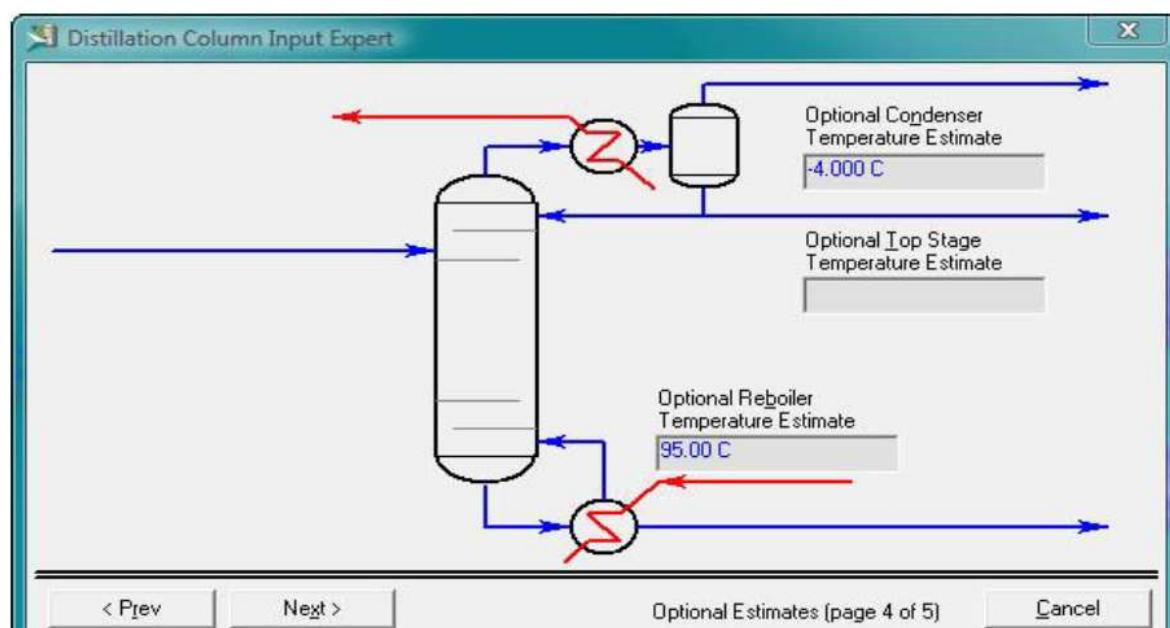
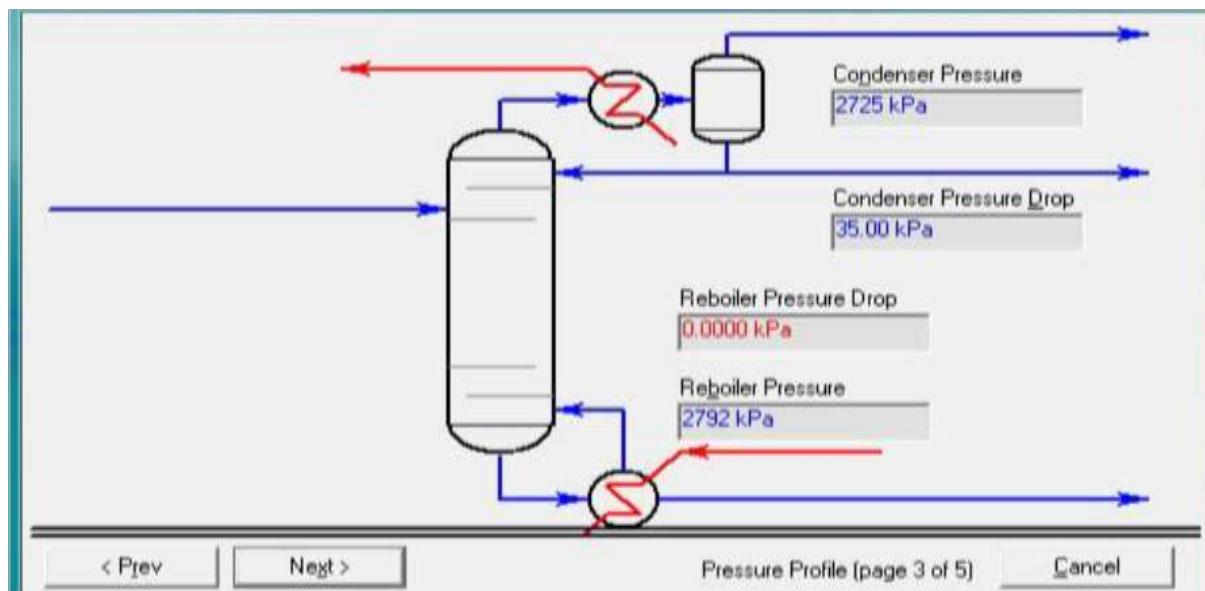
The screenshot shows the HYSYS Worksheet window for the pump. The 'Worksheet' tab is selected. The table displays various conditions and properties. The 'Dc2-feed' column contains values for different properties, and the value for Pressure [kPa] is highlighted in red as 2790. A callout arrow points from the text above to this highlighted value.

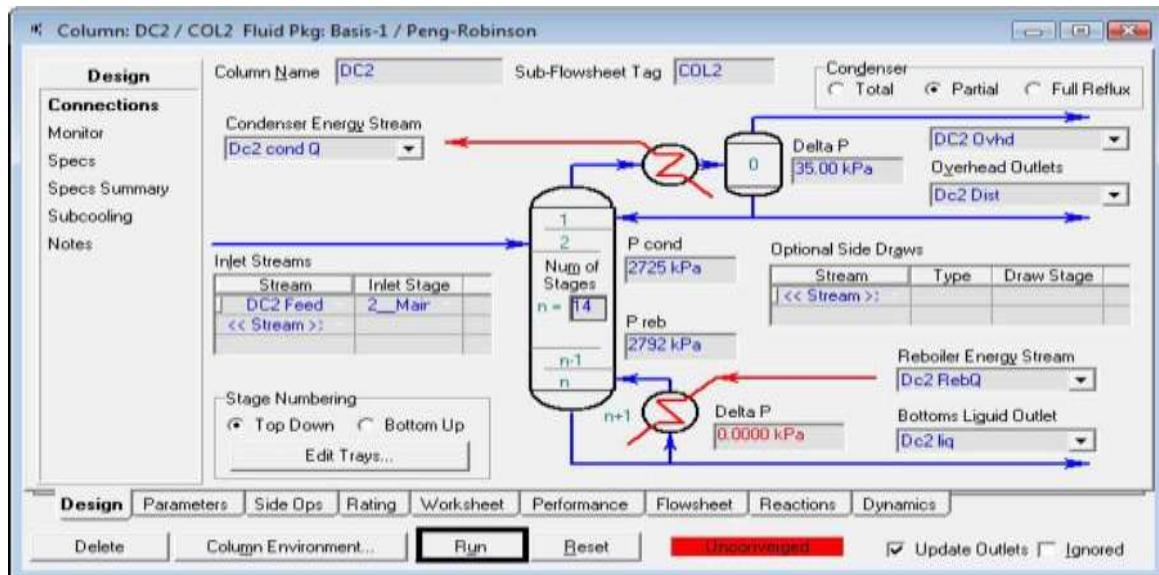
	Name	Dc1-bottom	Dc2-feed	P-100-hp
Vapour	0.0000	0.0000	<empty>	
Temperature [C]	20.06	20.77	<empty>	
Pressure [kPa]	2310	2790	<empty>	
Molar Flow [kgmole/h]	485.0	485.0	<empty>	
Mass Flow [kg/h]	1.837e+004	1.837e+004	<empty>	
Std Ideal Liq Vol Flow [m3/h]	42.65	42.65	<empty>	
Molar Enthalpy [kJ/kgmole]	-1.104e+005	-1.103e+005	<empty>	
Molar Entropy [kJ/kgmole-C]	120.7	120.8	<empty>	
Heat Flow [kJ/h]	-5.353e+007	-5.350e+007	2.620e+004	

## PART 2: DE- ETHANIZER COLUMN

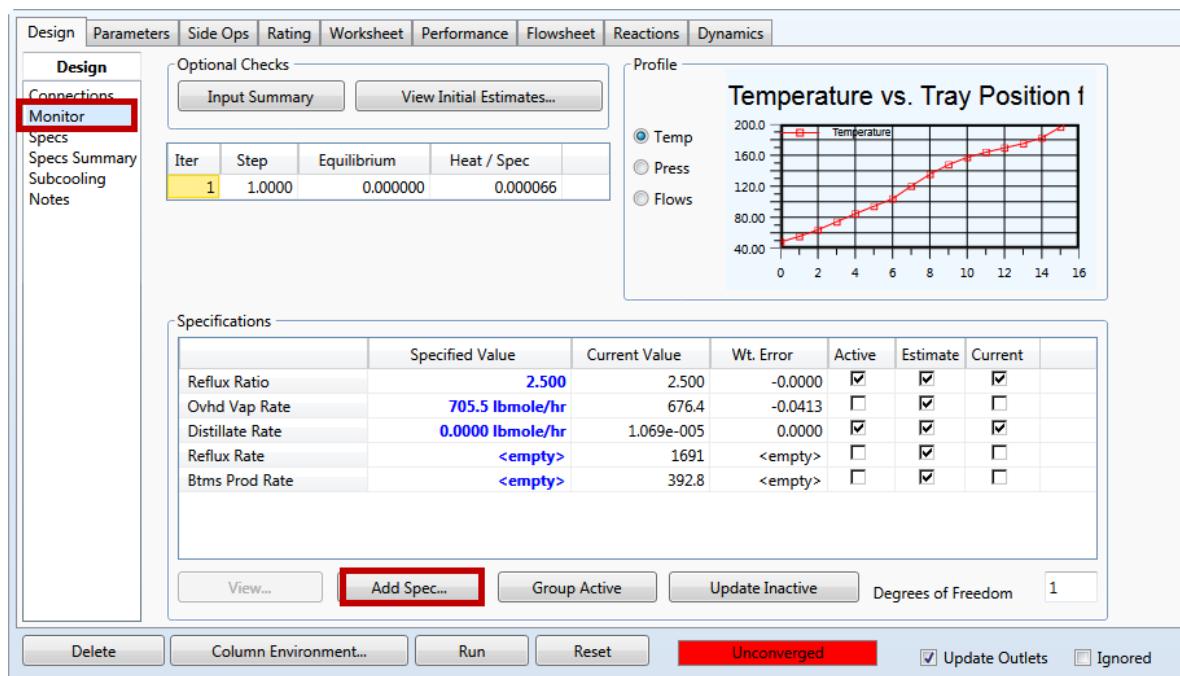
The outlet from the pump is then fed to the de-ethanizer:







Add the Column component ratio from the monitor page as follows:



**Add Specs - DC2 (COL...)**

Column Specification Types

- Column Cold Properties Spec
- Column Component Flow
- Column Component Fraction
- Column Component Ratio**
- Column Component Recovery
- Column Cut Point
- Column Draw Rate
- Column DT (Heater/Cooler) Spec
- Column Dt Spec
- Column Duty
- Column Duty Ratio
- Column Feed Ratio
- Column Gap Cut Point
- Column Liquid Flow
- Column Physical Properties Spec
- Column Pump Around
- Column Reboil Ratio Spec
- Column Recovery
- Column Reflux Feed Ratio Spec
- Column Reflux Fraction Spec

**Add Spec(s)...**

**Comp Ratio Spec: C2/C3**

Name	C2/C3
Stage	Reboiler
Flow Basis	Mole Fraction
Phase	Liquid
Spec Value	1.000e-002
Numerator:	Ethane
Denominator:	Propane

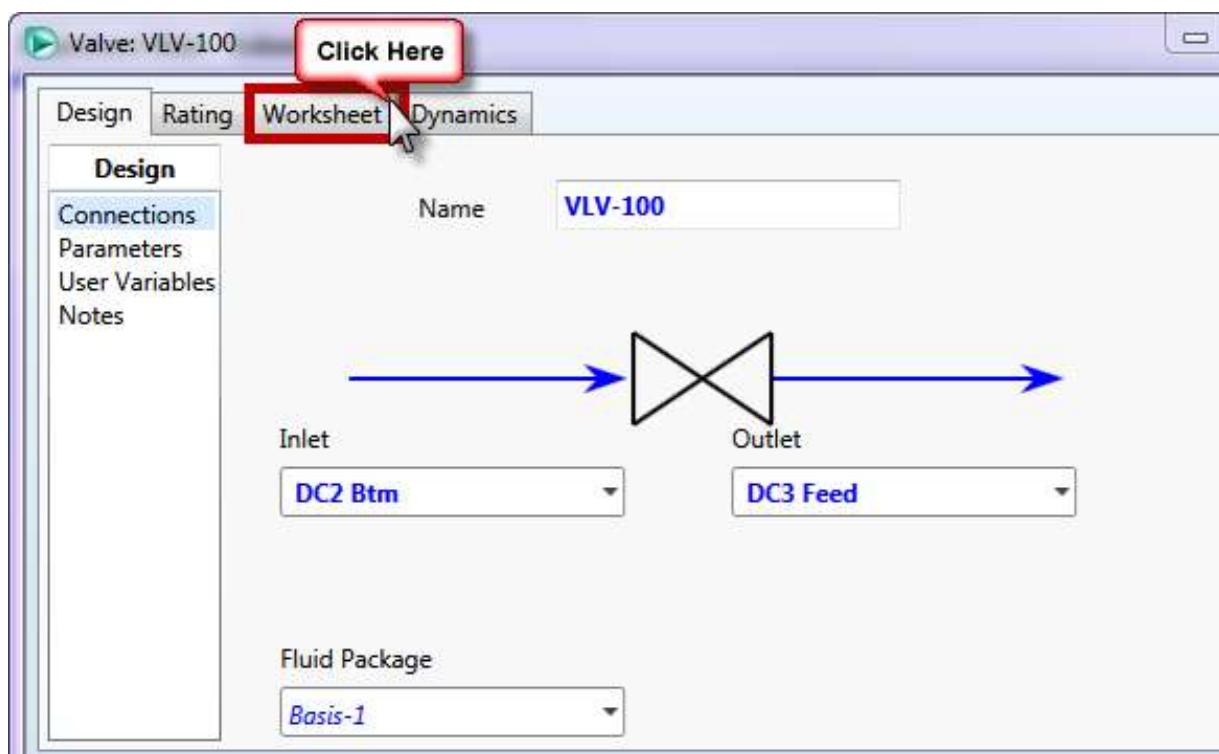
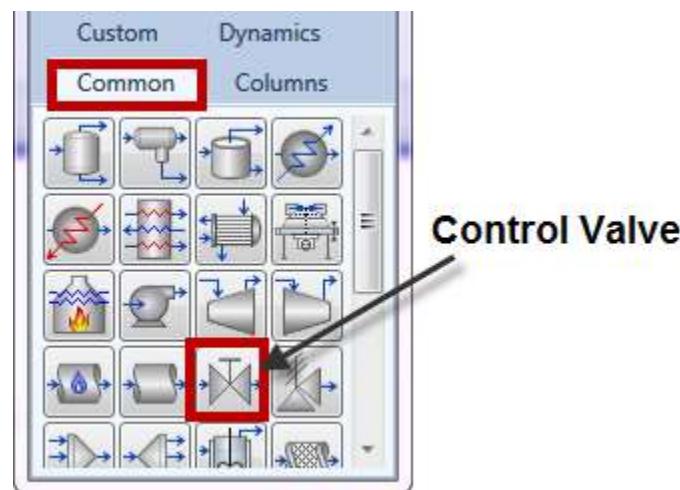
Target Type  Stream  Stage

**Parameters** **Summary** **Spec Type**

Delete



Add a valve on the bottom liquid stream:



Outlet pressure from the valve = 1690 kPa (from the **Worksheet** tab)

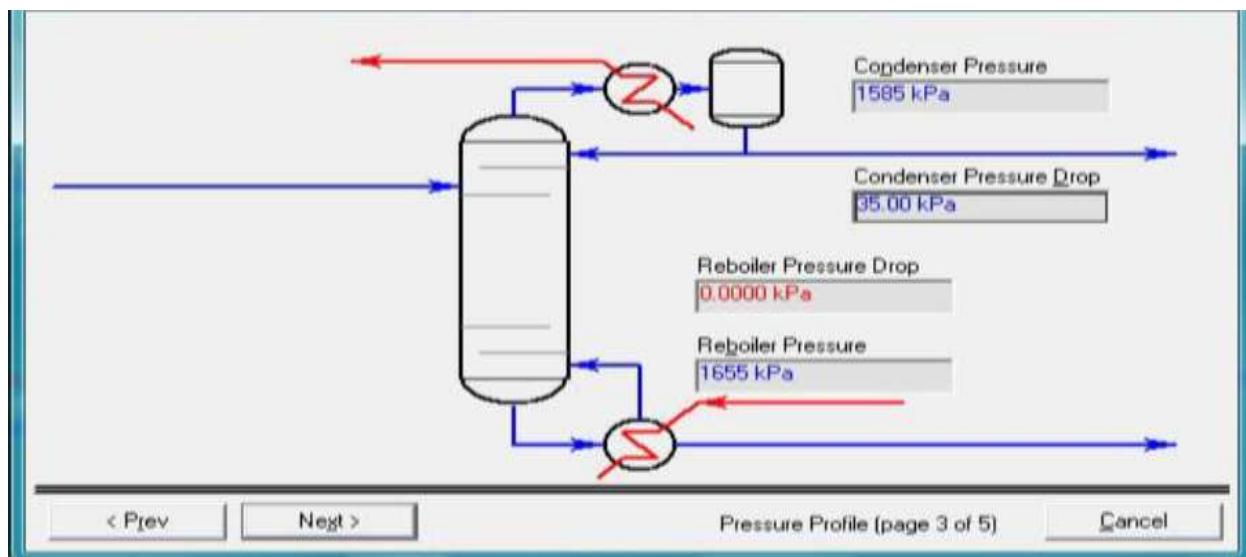
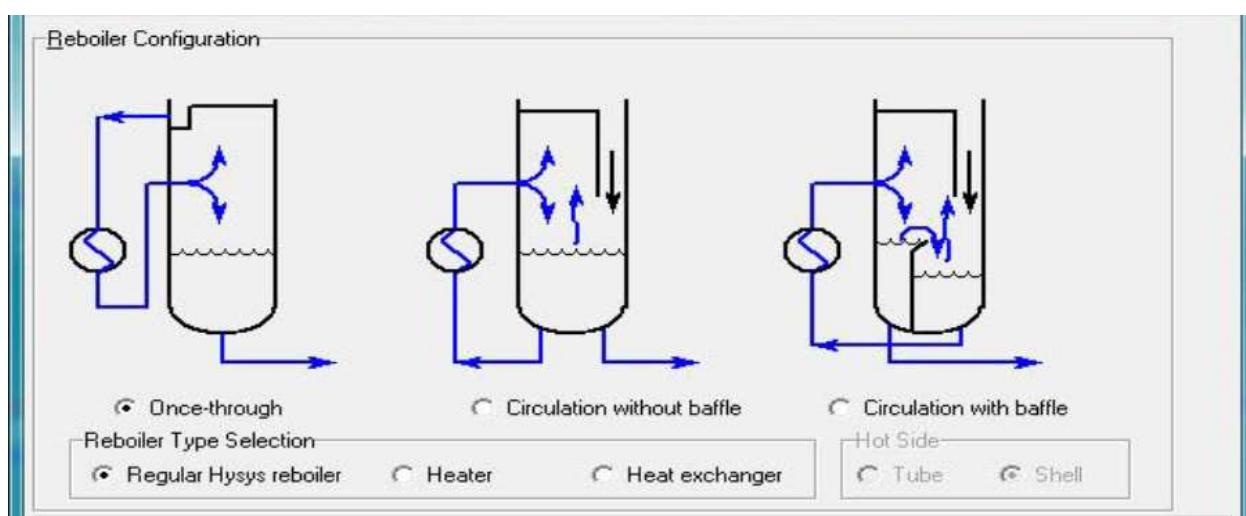
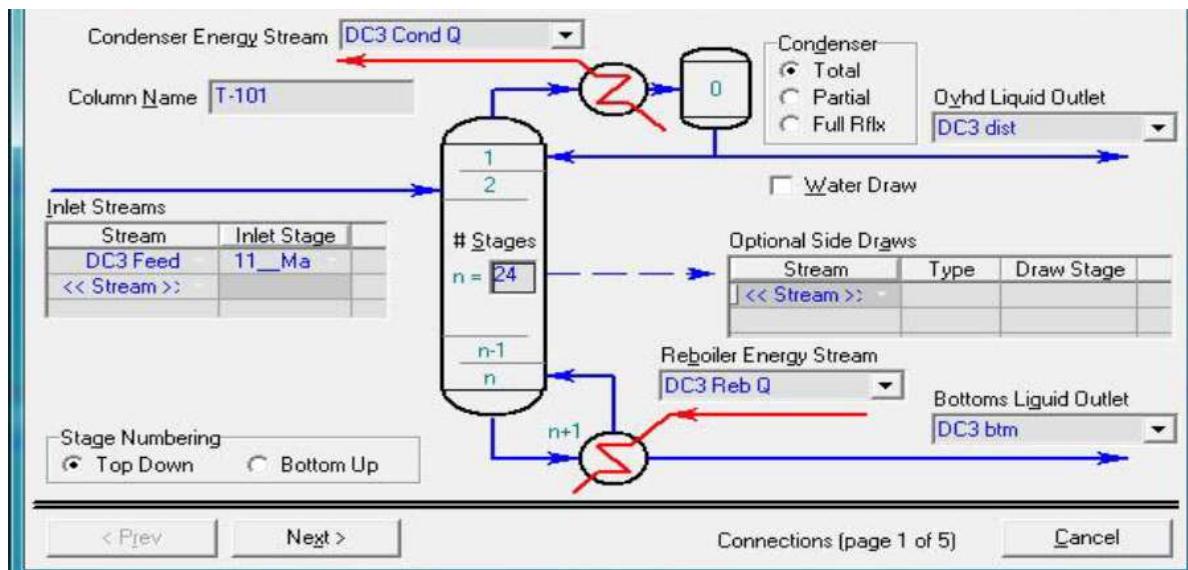
Worksheet	Name	Dc2 liq	DC3 Feed
Conditions	Vapour	0.0000	0.3210
Properties	Temperature [C]	91.16	67.39
Composition	Pressure [kPa]	2792	1690
PF Specs	Molar Flow [kgmole/h]	179.2	179.2
	Mass Flow [kg/h]	9036	9036
	Std Ideal Liq Vol Flow [m <sup>3</sup> /h]	16.78	16.78
	Molar Enthalpy [kJ/kgmole]	-1.233e+005	-1.233e+005
	Molar Entropy [kJ/kgmole-C]	121.5	122.3
	Heat Flow [kJ/h]	-2.211e+007	-2.211e+007

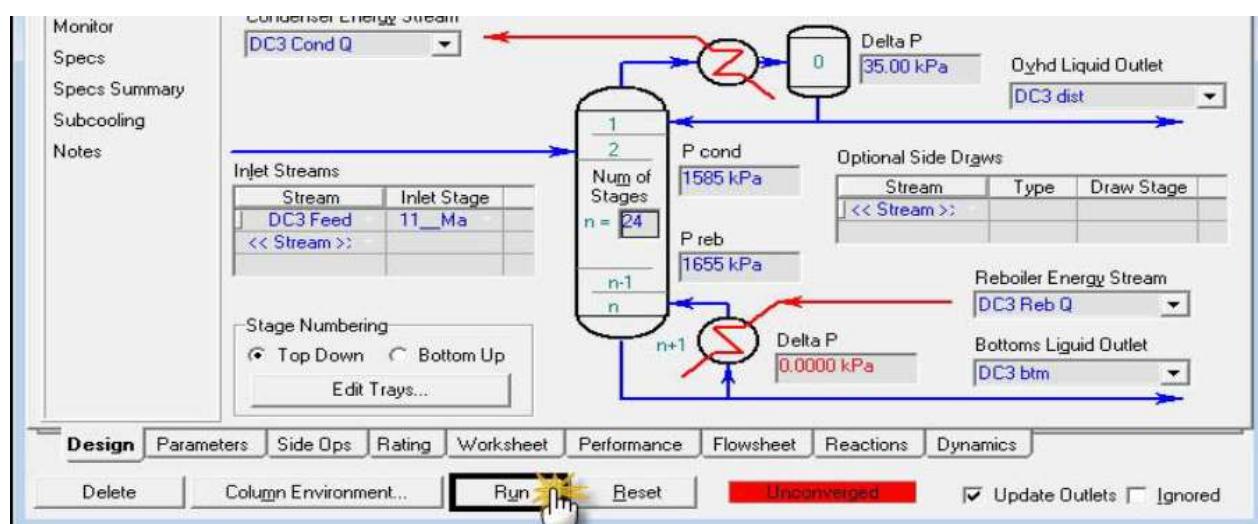
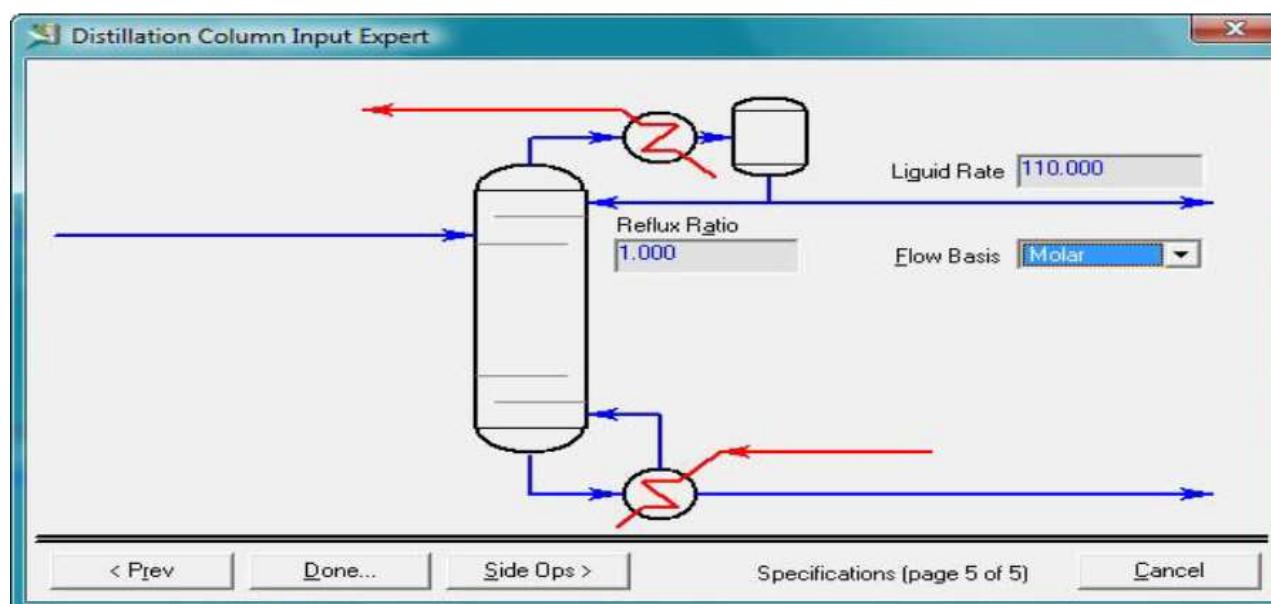
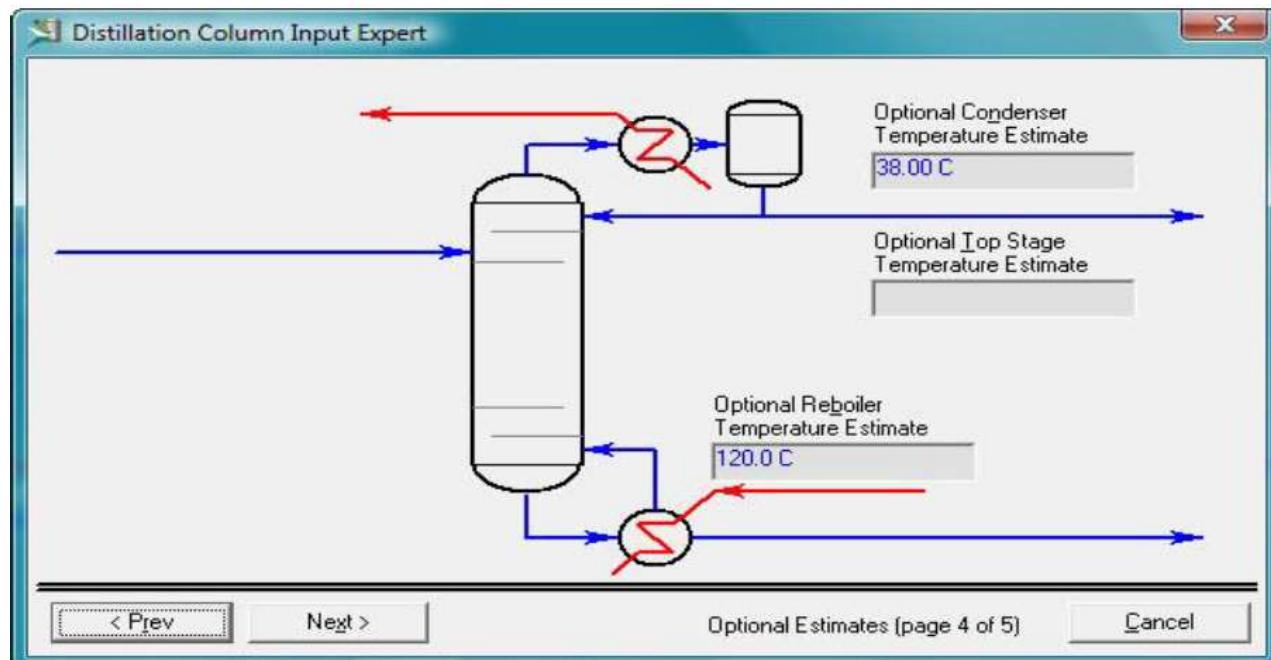
Design Rating **Worksheet** Dynamics

Delete OK  Ignored

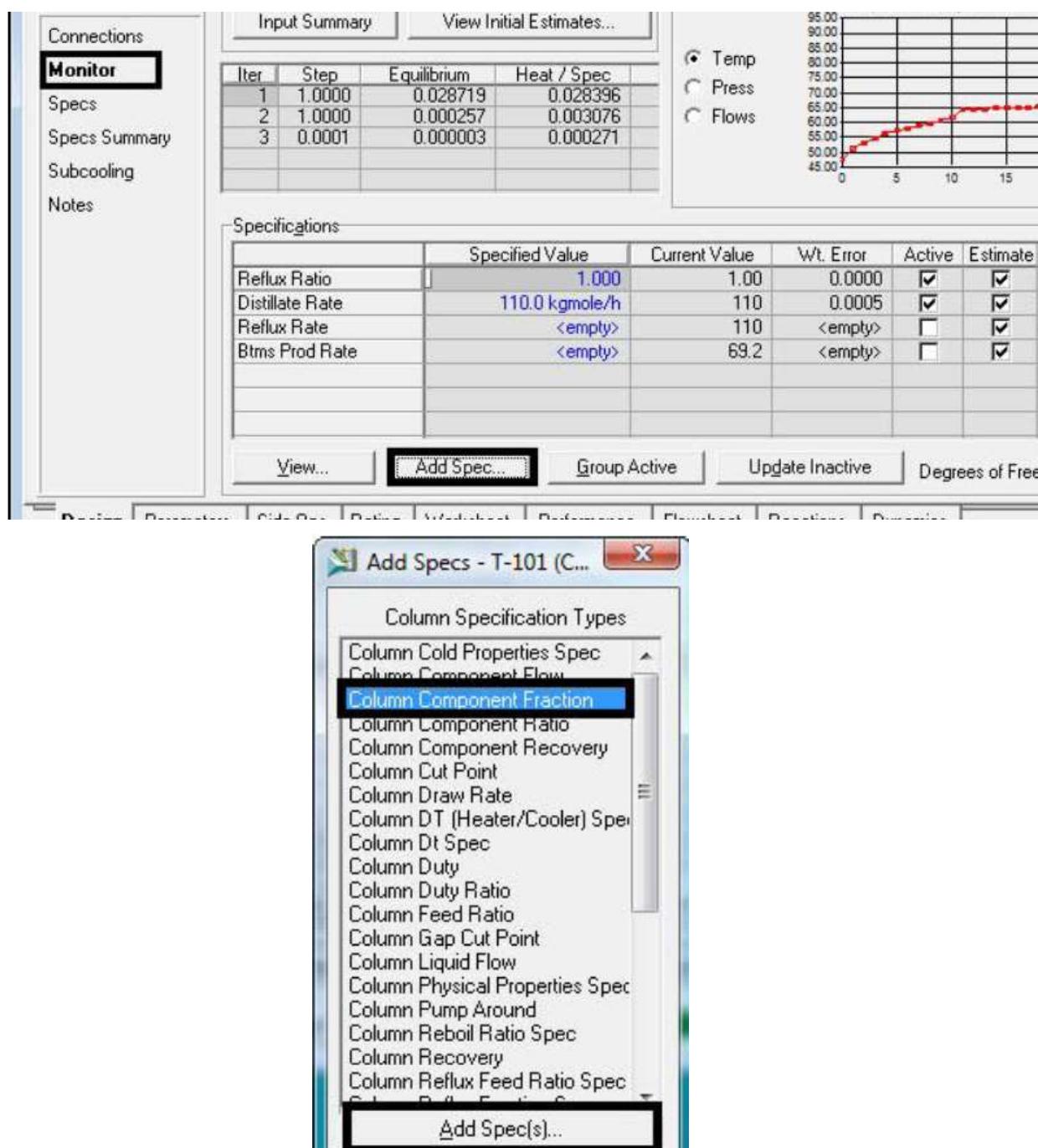
## PART 3: DE- PROPANIZER COLUMN

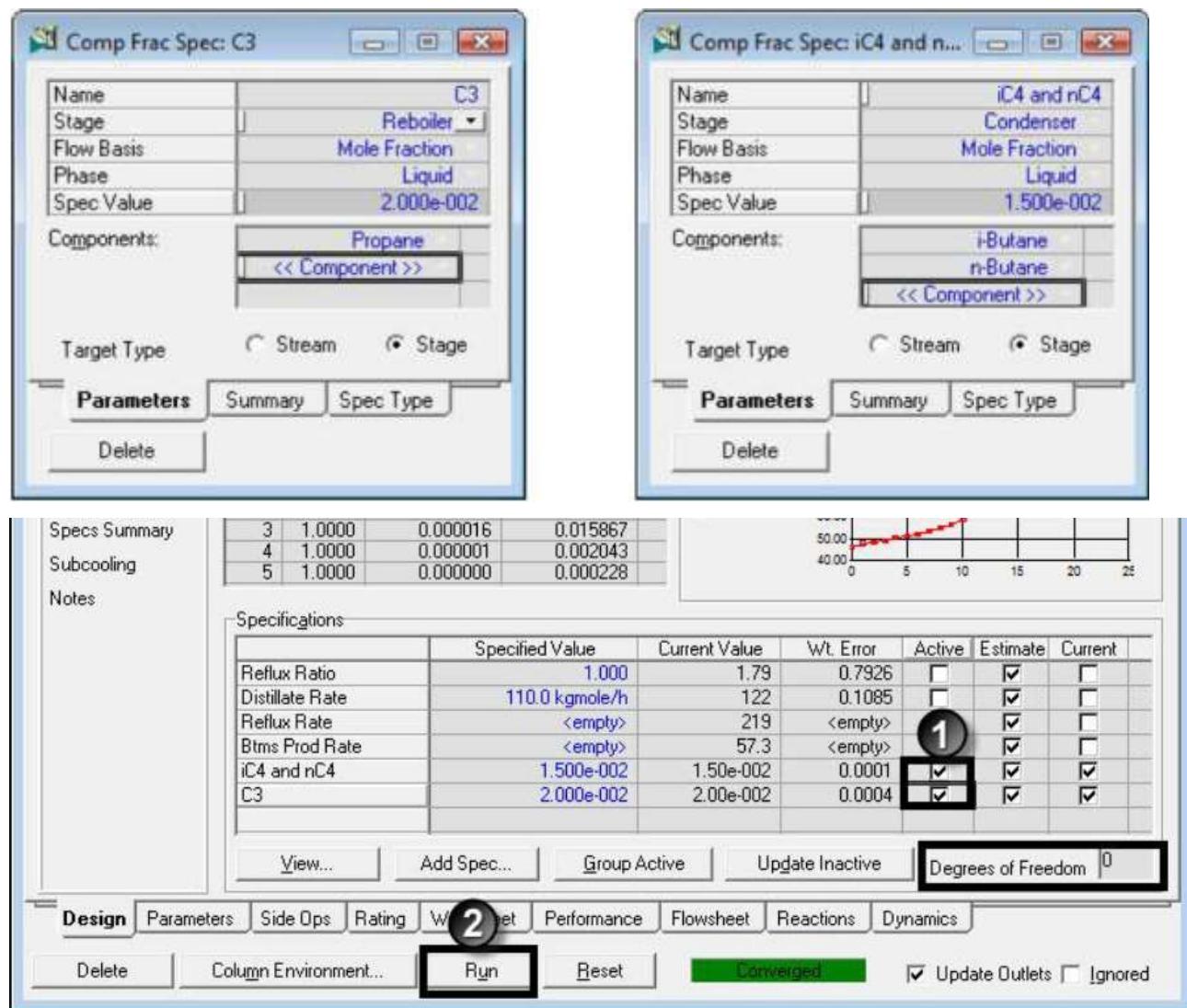
Add a distillation column (De-propanizer):





Now, let's add a new 2 specifications instead of the current:





After running the column, you can view the results from the **Performance** page (for any column)

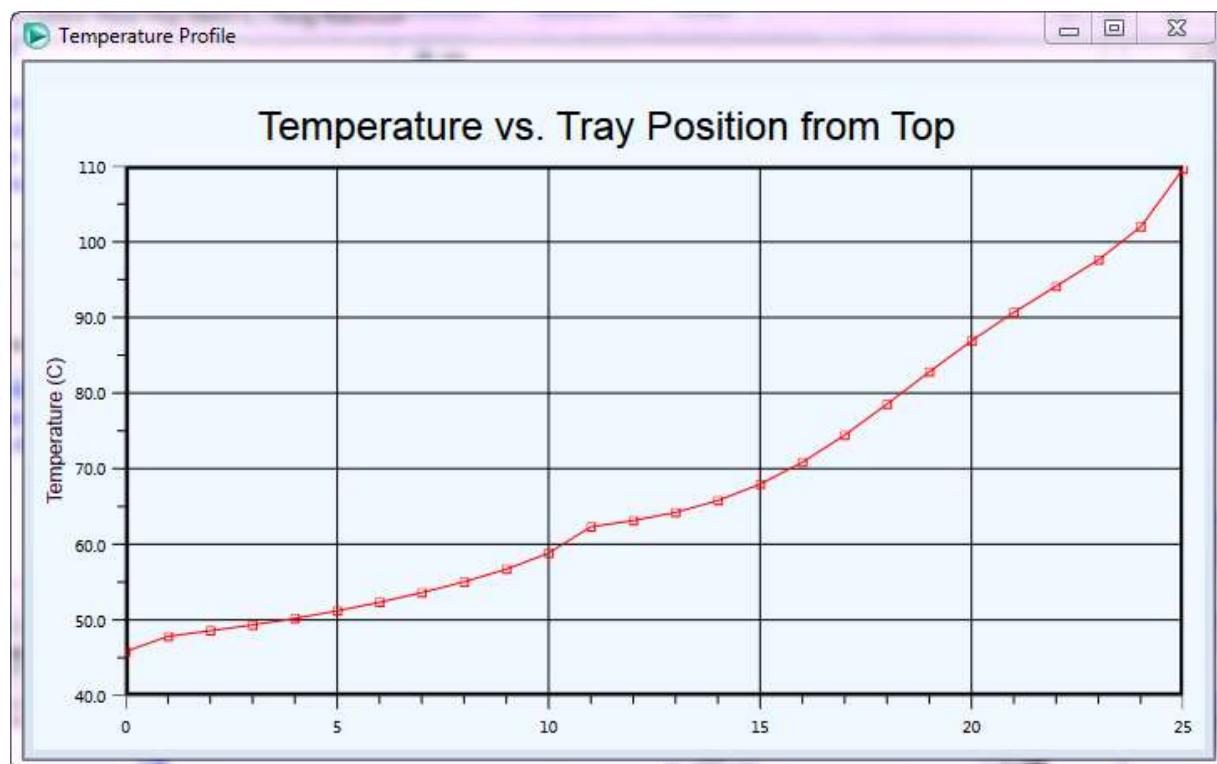
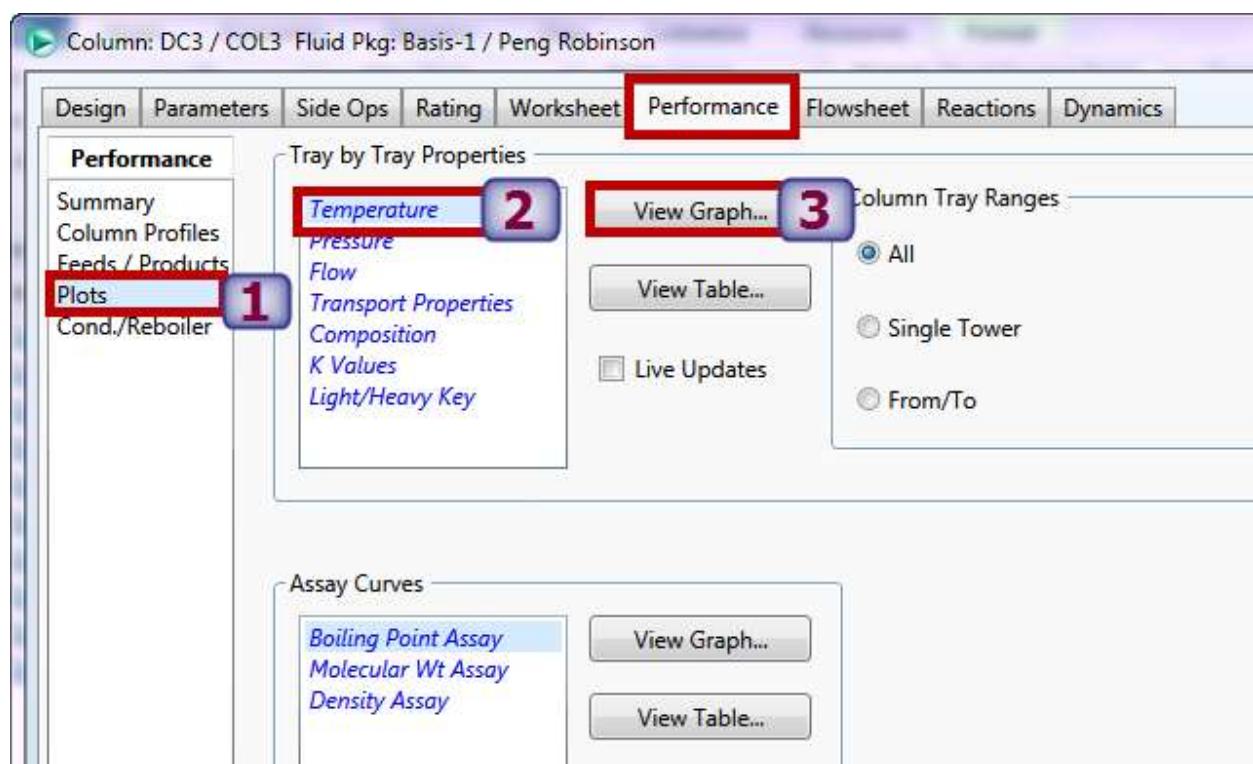
Column: DC3 / COL3 Fluid Pkg: Basis-1 / Peng Robinson

**Click Here**

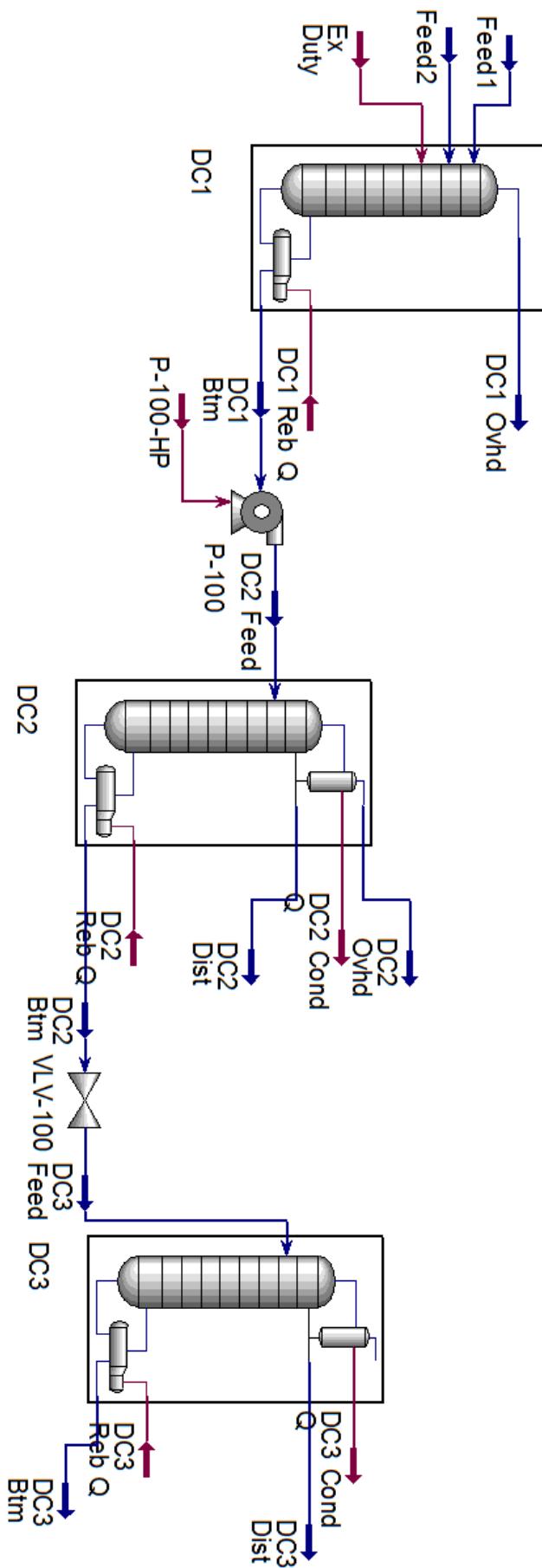
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<b>Feeds</b> <table border="1"> <thead> <tr> <th colspan="2">DC3 Feed</th> </tr> </thead> <tbody> <tr> <td>Flow Rate (kgmole/h)</td> <td>178.1833</td> </tr> <tr> <td>Nitrogen</td> <td>0.0000</td> </tr> <tr> <td>CO<sub>2</sub></td> <td>0.0000</td> </tr> <tr> <td>Methane</td> <td>0.0000</td> </tr> <tr> <td>Ethane</td> <td>0.0067</td> </tr> <tr> <td>Propane</td> <td>0.6666</td> </tr> <tr> <td>i-Butane</td> <td>0.1260</td> </tr> </tbody> </table> <b>Products</b> <table border="1"> <thead> <tr> <th colspan="3">DC3 Dist</th> <th>DC3 Btm</th> </tr> </thead> <tbody> <tr> <td>Flow Rate (kgmole/h)</td> <td>120.6237</td> <td>57.5596</td> <td></td> </tr> <tr> <td>Nitrogen</td> <td>0.0000</td> <td>0.0000</td> <td></td> </tr> <tr> <td>CO<sub>2</sub></td> <td>0.0000</td> <td>0.0000</td> <td></td> </tr> <tr> <td>Methane</td> <td>0.0000</td> <td>0.0000</td> <td></td> </tr> <tr> <td>Ethane</td> <td>0.0098</td> <td>0.0000</td> <td></td> </tr> <tr> <td>Propane</td> <td>0.9752</td> <td>0.0200</td> <td></td> </tr> <tr> <td>i-Butane</td> <td>0.0138</td> <td>0.3613</td> <td></td> </tr> </tbody> </table>									DC3 Feed		Flow Rate (kgmole/h)	178.1833	Nitrogen	0.0000	CO <sub>2</sub>	0.0000	Methane	0.0000	Ethane	0.0067	Propane	0.6666	i-Butane	0.1260	DC3 Dist			DC3 Btm	Flow Rate (kgmole/h)	120.6237	57.5596		Nitrogen	0.0000	0.0000		CO <sub>2</sub>	0.0000	0.0000		Methane	0.0000	0.0000		Ethane	0.0098	0.0000		Propane	0.9752	0.0200		i-Butane	0.0138	0.3613	
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<b>Reflux Ratio</b> 1.817 <b>Boilup Ratio</b> 4.204 <div style="display: flex; justify-content: space-around; margin-top: 10px;"> <span><input checked="" type="radio"/> Flows</span> <span><input type="radio"/> Energy</span> <span>Basis</span> <div style="display: flex; justify-content: space-around;"> <span><input checked="" type="radio"/> Molar</span> <span><input type="radio"/> Mass</span> <span><input type="radio"/> Ideal</span> </div> <div style="display: flex; justify-content: space-around; margin-top: 5px;"> <span><input type="radio"/> Liq Vol @Std Cond</span> <span><input type="radio"/> Act. V</span> </div> </div> <table border="1"> <thead> <tr> <th></th> <th>Temperature [C]</th> <th>Pressure [kPa]</th> <th>Net Liquid [kgmole/h]</th> <th>Net Vapour [kgmole/h]</th> <th>Net Feed [kgmole/h]</th> <th>Net Draw [kgmole/h]</th> </tr> </thead> <tbody> <tr> <td>Condenser</td> <td>45.83</td> <td>1585</td> <td>219.135</td> <td></td> <td></td> <td></td> </tr> <tr> <td>1_Main TS</td> <td>47.83</td> <td>1620</td> <td>220.901</td> <td>339.759</td> <td></td> <td></td> </tr> <tr> <td>2_Main TS</td> <td>48.58</td> <td>1622</td> <td>219.835</td> <td>341.525</td> <td></td> <td></td> </tr> <tr> <td>3_Main TS</td> <td>49.34</td> <td>1623</td> <td>218.431</td> <td>340.458</td> <td></td> <td></td> </tr> <tr> <td>4_Main TS</td> <td>50.20</td> <td>1625</td> <td>216.760</td> <td>339.055</td> <td></td> <td></td> </tr> <tr> <td>5_Main TS</td> <td>51.20</td> <td>1626</td> <td>214.845</td> <td>337.384</td> <td></td> <td></td> </tr> <tr> <td>6_Main TS</td> <td>52.33</td> <td>1628</td> <td>212.688</td> <td>335.469</td> <td></td> <td></td> </tr> <tr> <td>7_Main TS</td> <td>53.61</td> <td>1629</td> <td>210.236</td> <td>333.311</td> <td></td> <td></td> </tr> <tr> <td>8_Main TS</td> <td>55.04</td> <td>1631</td> <td>207.290</td> <td>330.860</td> <td></td> <td></td> </tr> <tr> <td>9_Main TS</td> <td>56.71</td> <td>1632</td> <td>203.256</td> <td>327.914</td> <td></td> <td></td> </tr> <tr> <td>10_Main TS</td> <td>58.85</td> <td>1634</td> <td>196.071</td> <td>323.879</td> <td></td> <td></td> </tr> <tr> <td>11_Main TS</td> <td>62.32</td> <td>1635</td> <td>311.932</td> <td>316.695</td> <td>178.18</td> <td></td> </tr> <tr> <td>12_Main TS</td> <td>63.13</td> <td>1637</td> <td>211.458</td> <td>251.372</td> <td></td> <td></td> </tr> </tbody> </table>										Temperature [C]	Pressure [kPa]	Net Liquid [kgmole/h]	Net Vapour [kgmole/h]	Net Feed [kgmole/h]	Net Draw [kgmole/h]	Condenser	45.83	1585	219.135				1_Main TS	47.83	1620	220.901	339.759			2_Main TS	48.58	1622	219.835	341.525			3_Main TS	49.34	1623	218.431	340.458			4_Main TS	50.20	1625	216.760	339.055			5_Main TS	51.20	1626	214.845	337.384			6_Main TS	52.33	1628	212.688	335.469			7_Main TS	53.61	1629	210.236	333.311			8_Main TS	55.04	1631	207.290	330.860			9_Main TS	56.71	1632	203.256	327.914			10_Main TS	58.85	1634	196.071	323.879			11_Main TS	62.32	1635	311.932	316.695	178.18		12_Main TS	63.13	1637	211.458	251.372		
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# Save Your Case!



*Oil*

*Stabilization*

*Optimization*

## **Workshop**

A poor-boy stabilization scheme is used to separate an oil and gas mixture into a stabilized oil and a saleable gas. A simple three-stage separation with heating between each stage is used and the object of the exercise is to select the let-down pressure and temperatures such that the products revenue less the utilities cost is maximized. A special tool in HYSYS, the Optimizer, will be used to find the optimum operating conditions. HYSYS includes additional modelling and decision support tools that can be used to enhance the usability of your models. In this module, you will use the HYSYS optimization tool available in HYSYS to investigate the debottlenecking and optimization of a crude column.

## **Learning Objectives**

Once you have completed this section, you will be able to:

- Use the Optimizer tool in HYSYS to optimize flowsheets
- Use the Spreadsheet to perform calculations

## Oil Stabilization

A feed stream @  $10^{\circ}\text{C}$ , 4125 kPa with a flowrate of 1 MMSCFD is fed to a heater (duty= $4.25 \times 10^5$  kJ/hr) before entering the first separator where the separated liquid is heated in a second heater (duty= $3.15 \times 10^5$  kJ/hr). The outlet from the heater is then sent to a letdown valve in order to decrease the pressure to 2050 kPa before entering the second separator where the separated liquid is heated through a third heater (duty= $1.13 \times 10^5$  kJ/hr). The outlet from the third heater is then throttled through a valve (outlet pressure = 350 kPa) and then fed to a third separator to obtain the final liquid oil product.

Each gas stream from the 2<sup>nd</sup> & 3<sup>rd</sup> separators is fed to a separate compressor to raise the pressure to 4125 kPa and then mixed (using a mixer) with the gas stream from the 1<sup>st</sup> separator to get the final gas product stream.

Notes:

- Pressure drop across all heaters and separators are 0.0.

Comp	Mol frac	Comp	Mol frac
C1	0.316	n-C5	0.053
C2	0.158	C6	0.027
C3	0.105	C7	0.026
i-C4	0.105	C8	0.026
n-C4	0.105	C9	0.026
i-C5	0.053		

Calculate:

- The total liquid product = ..... barrel/hr
- The total gas product = .....  $\text{m}^3$ \_gas/hr

## Oil Stabilization Optimization

In this case, we want to maximize the total operating profit while achieving an RVP of Liquid Product less than 96.5 kPa. The incomes from the Plant are both the Gas and Liquid Products. The operating costs are the Steam Costs for each Heater plus the Power Cost for each Compressor.

Profit = Income - Cost

Profit = (Gas Product + Liquid Product) - (Steam Costs + Compression Cost)

### **Prices & costs:**

Oil Price = 15 \$/bbl

Gas Price = 0.106 \$/m<sup>3</sup>\_gas

Steam Cost = 0.682 \$/kW-h

Compression Cost = 0.1 \$/kW-h

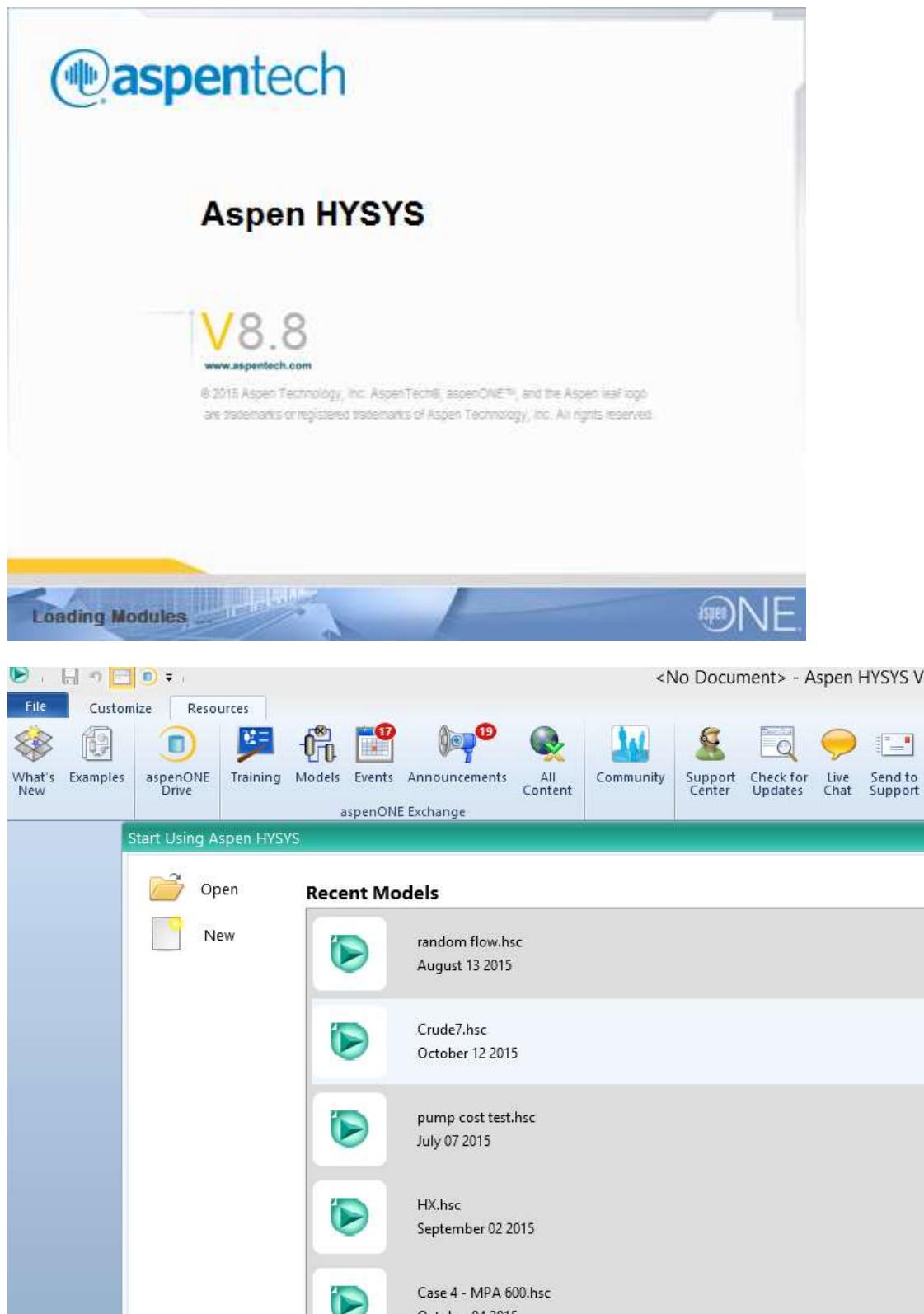
### **The variables to be adjusted:**

- Heater duties (for the 3 heaters). Use range of 0 – 1e6 kJ/hr
- Valves outlet pressures.
  - Use range of 650 – 3500 kPa for the first valve
  - Use range of 70 – 1000 kPa for the second valve

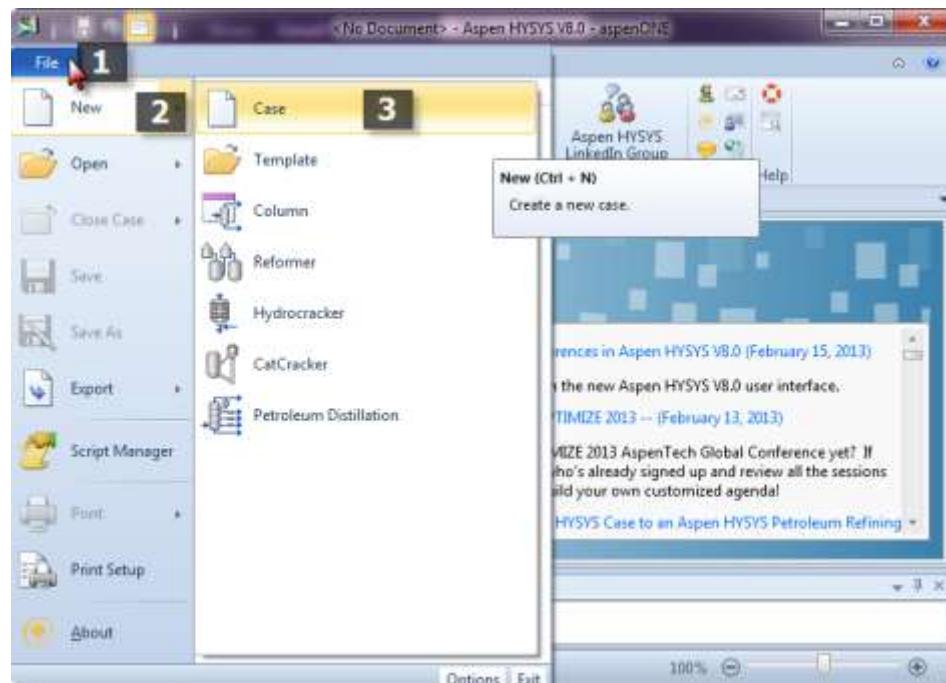
Calculate:

- The optimum values for the adjusted variables
- The maximum profit ..... \$/hr

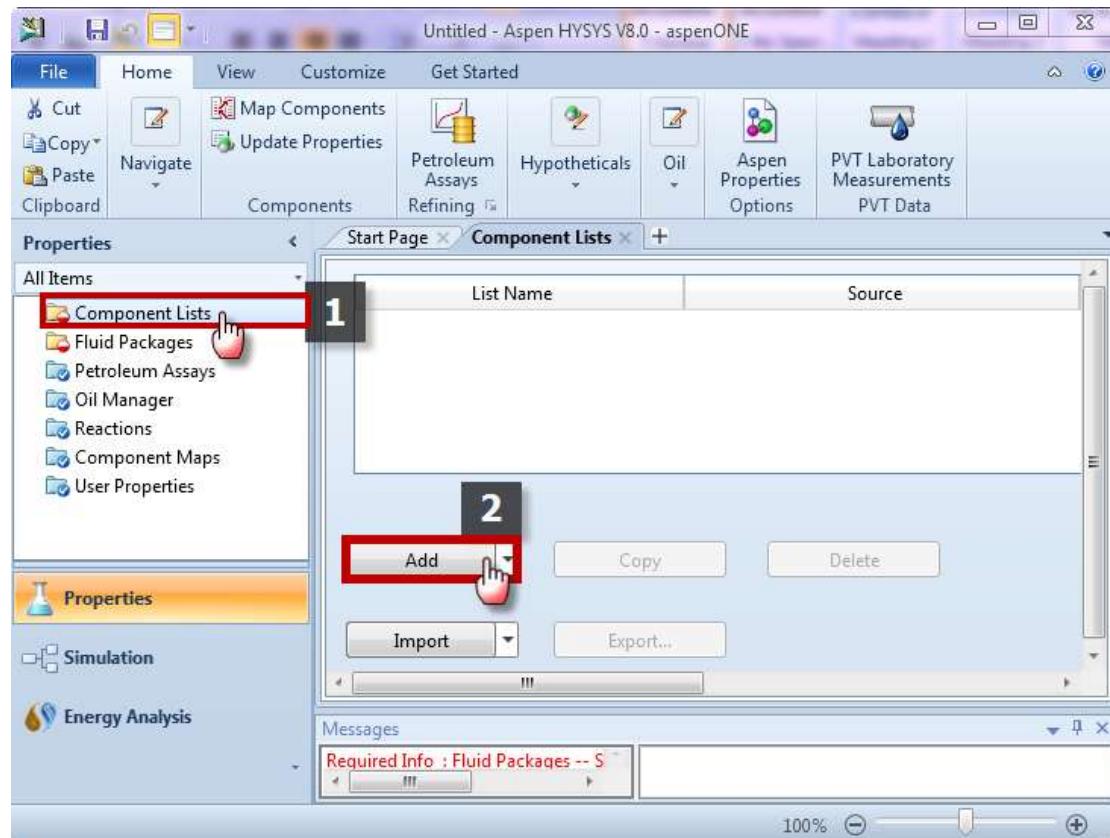
To start the program, From Start Menu, Select All Programs >>  
Aspen Tech >> Process Modeling V8 >>> Aspen HYSYS >>  
Aspen HYSYS



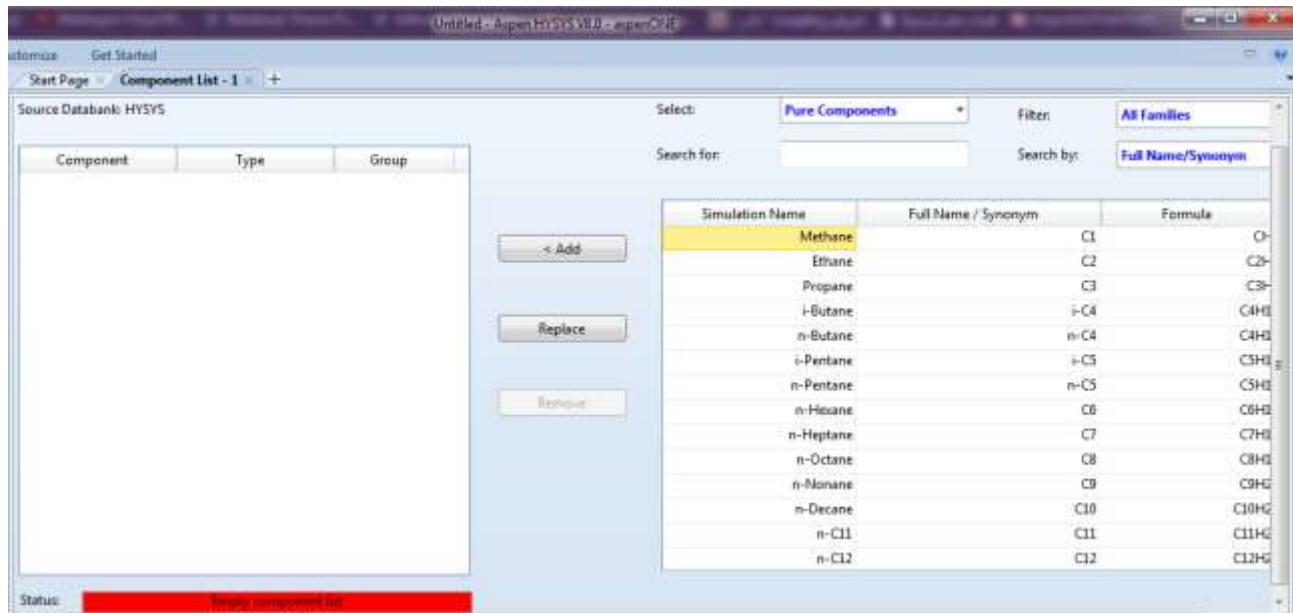
10- First, Start a new case



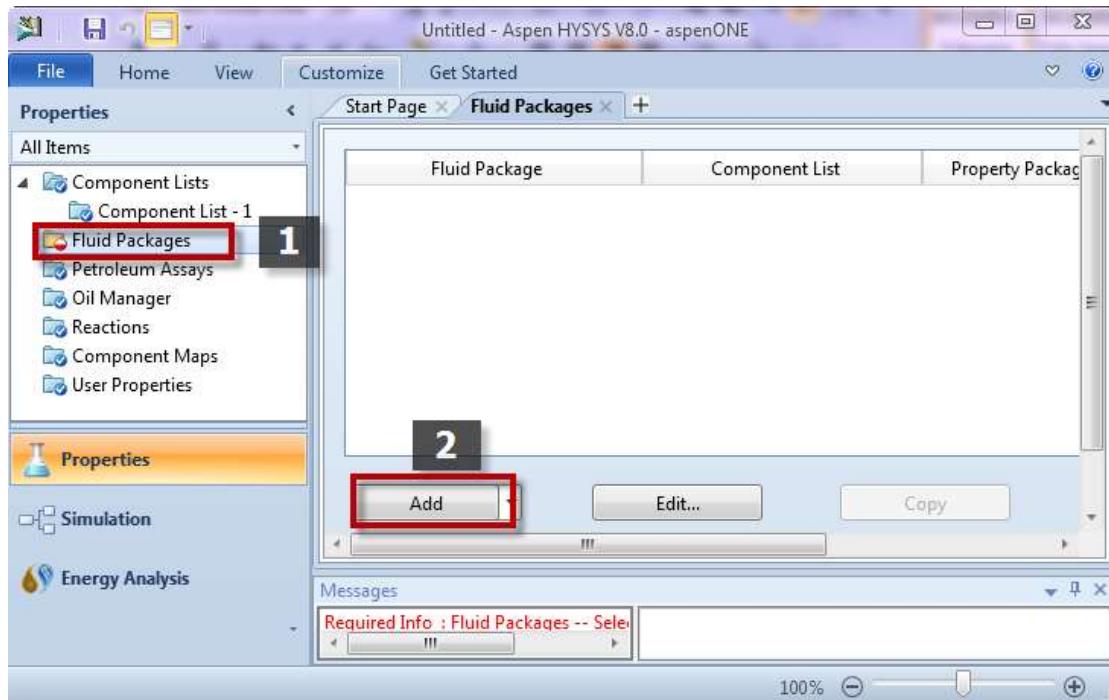
11- Add the Components



12- Choose the system components from the databank:

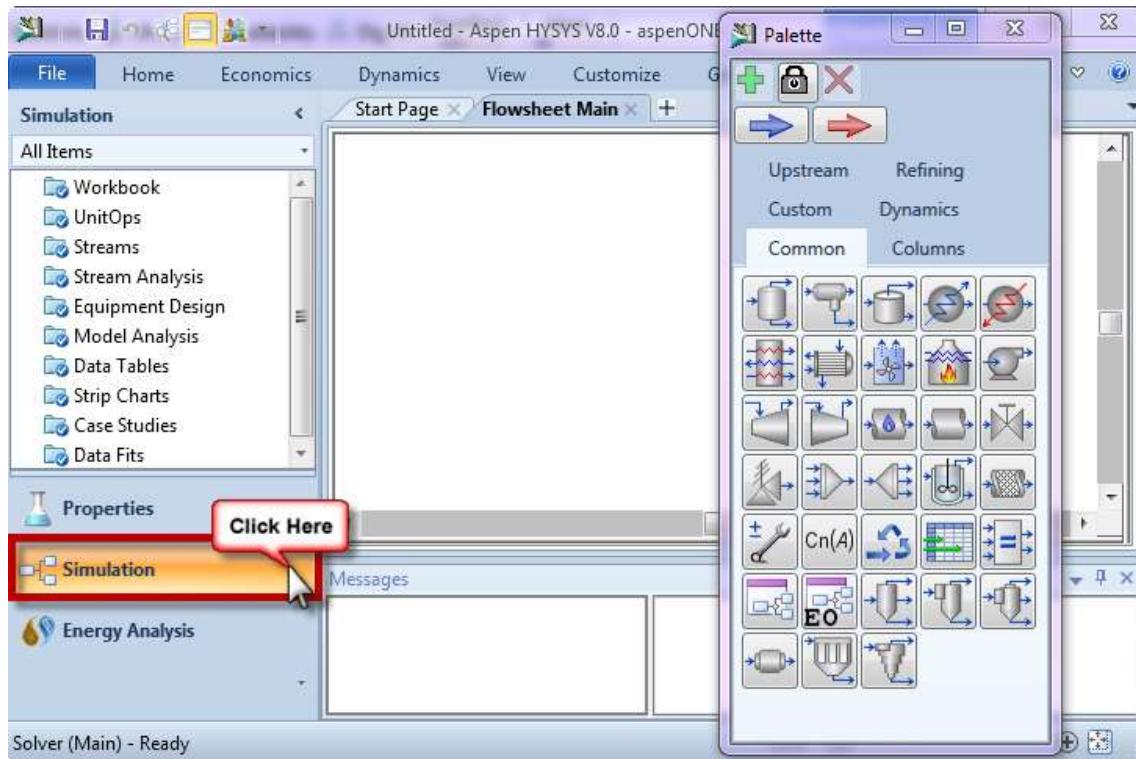


Now, select the suitable fluid package



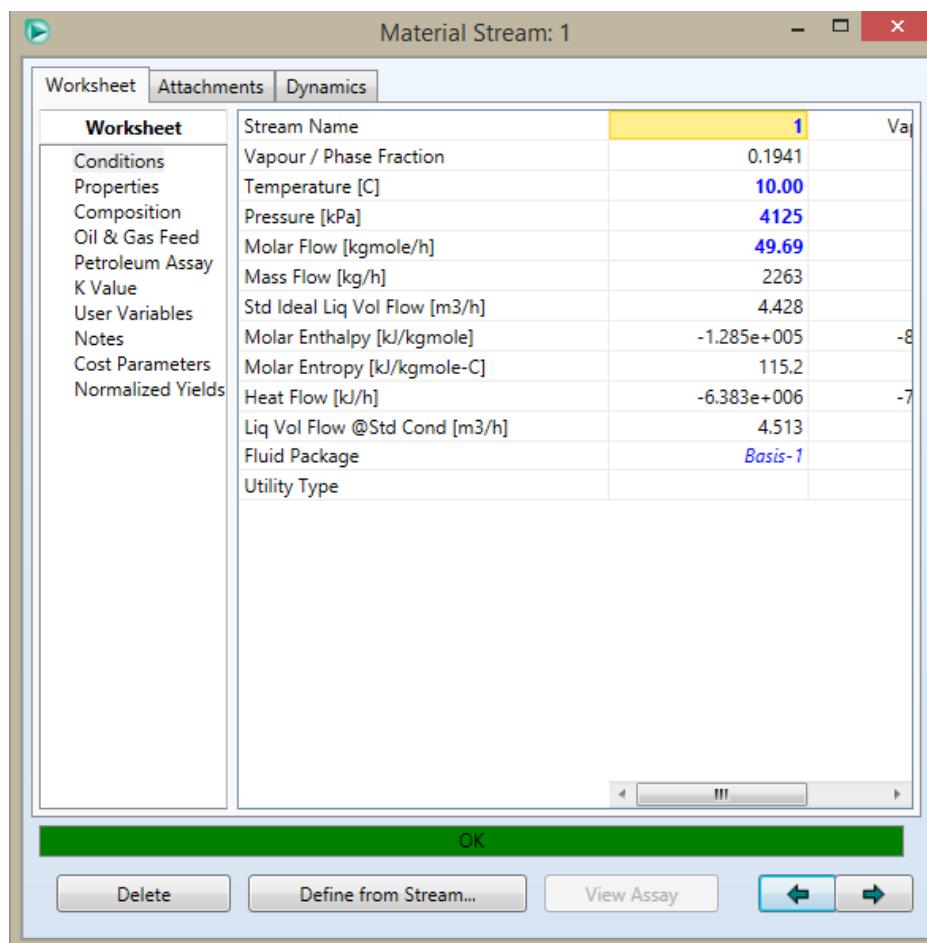
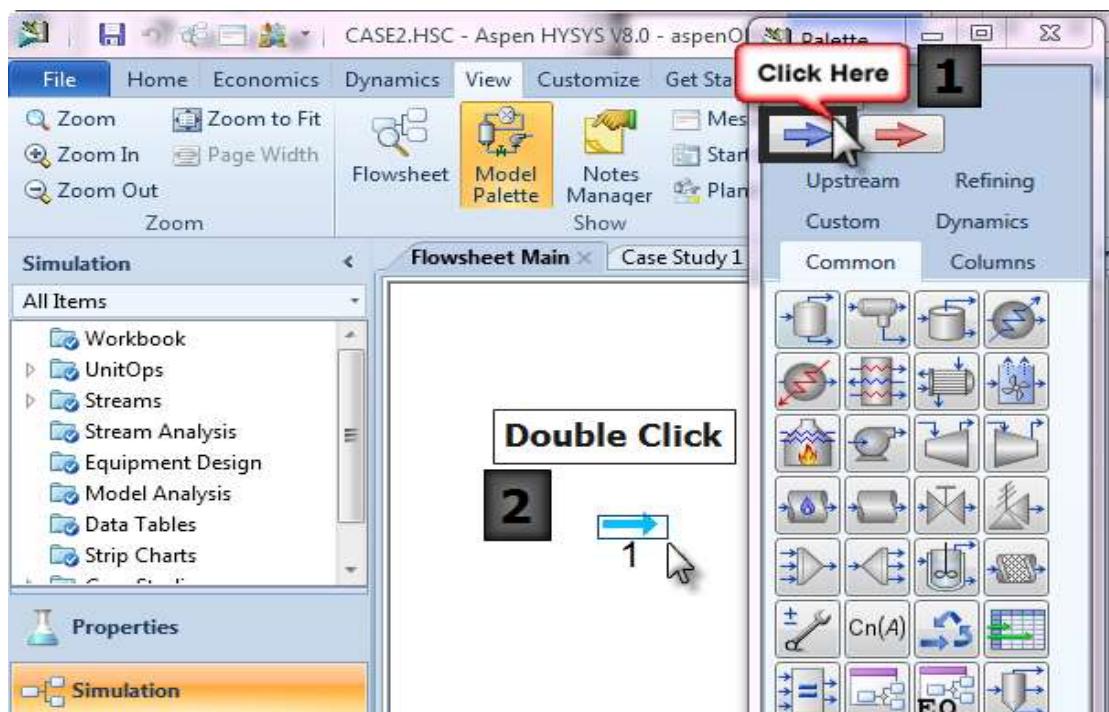
In this case, select Peng-Robinson

Now you can start drawing the flow sheet for the process by clicking the Simulation button:

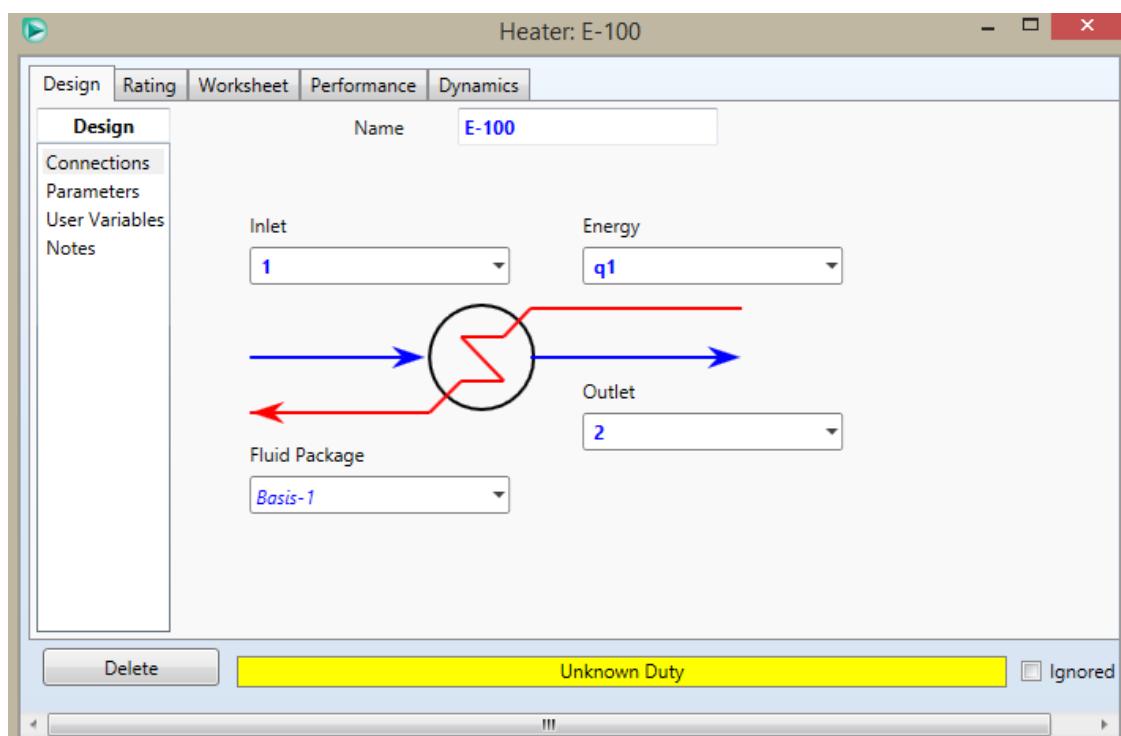
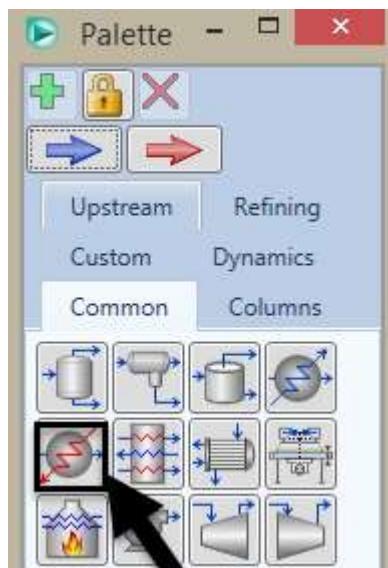


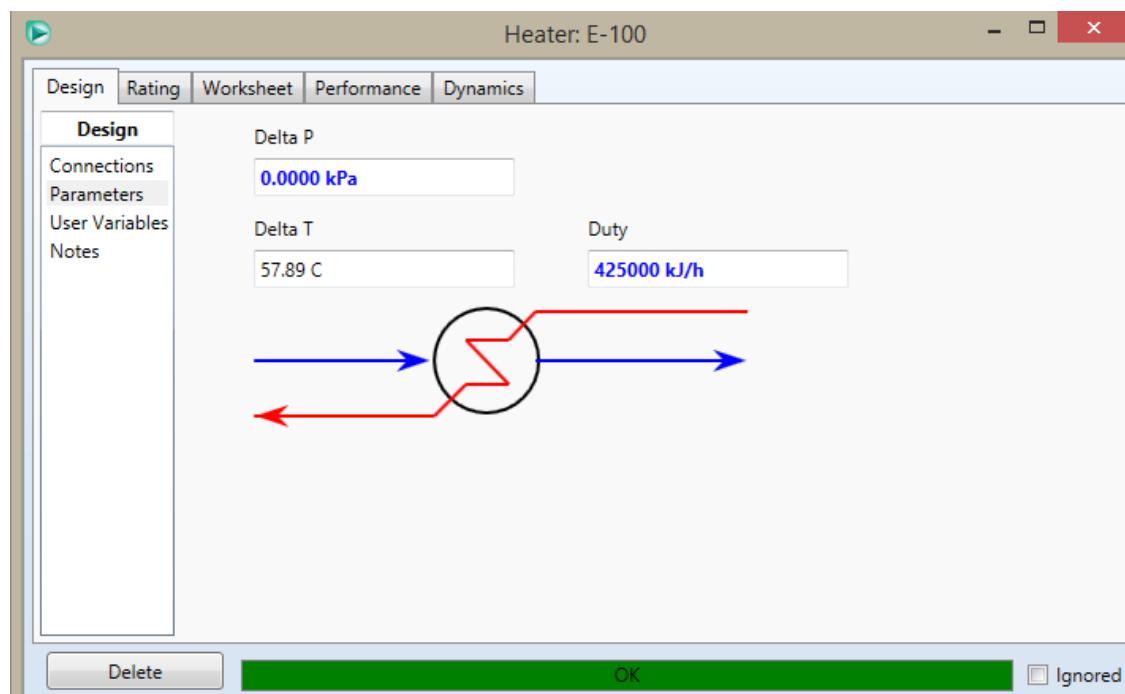
Now add a material stream to define the feed stream composition and conditions

From the palette:

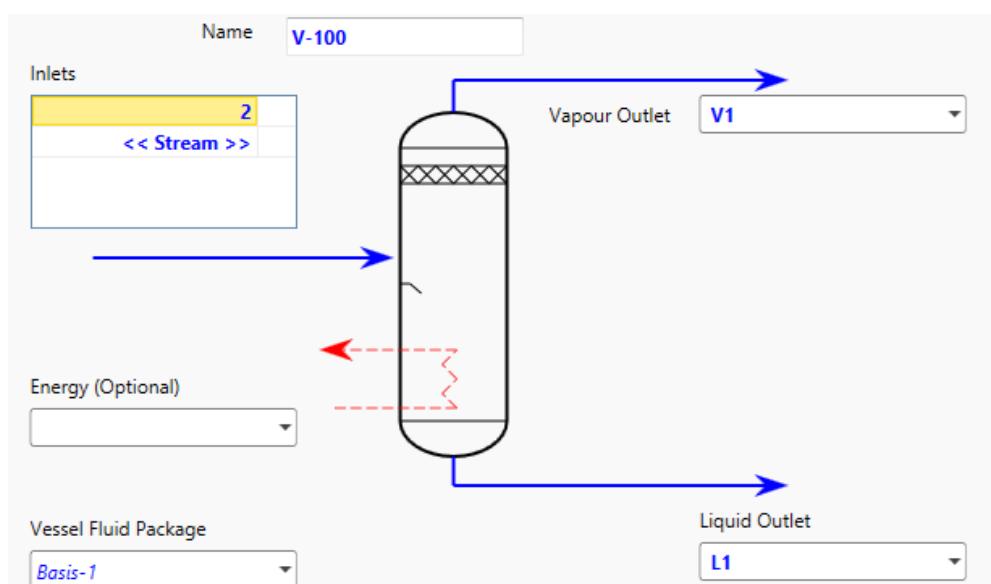
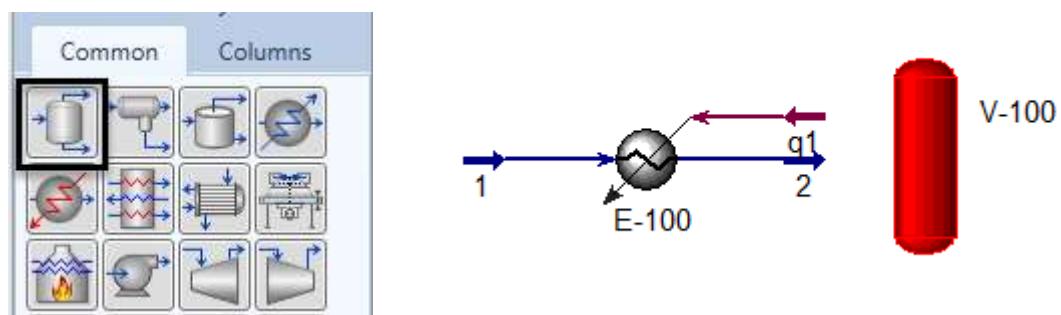


Add a heater with a duty of  $4.25 \times 10^5$  kJ/hr and pressure drop of 0.0

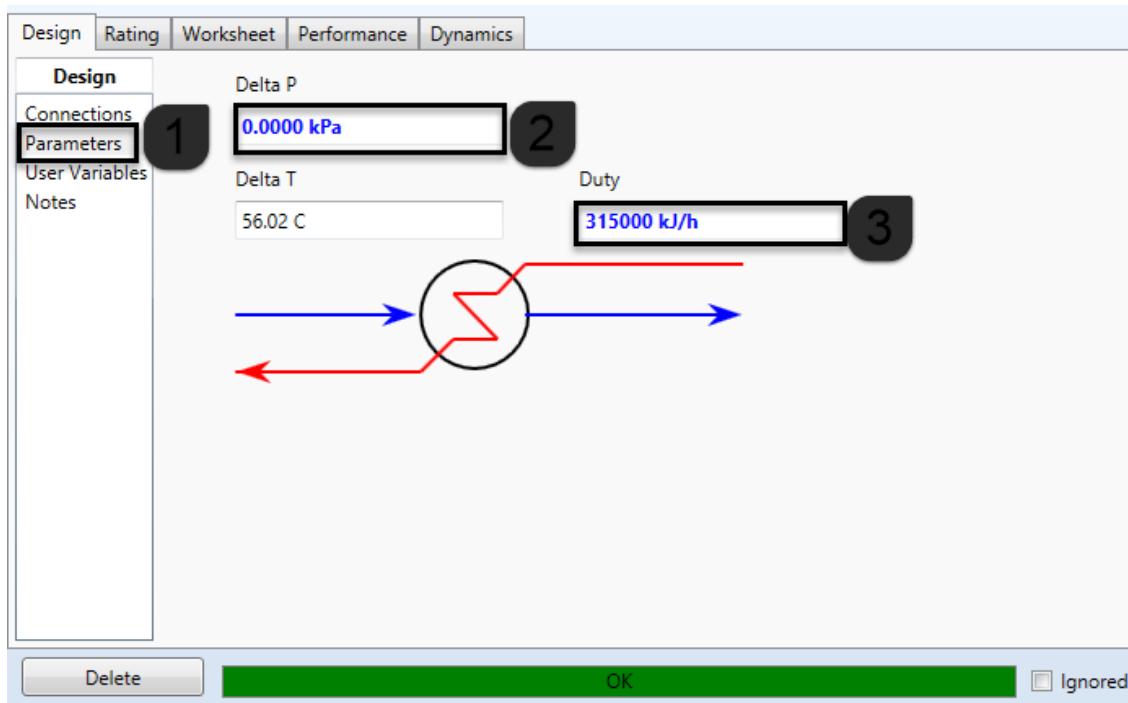




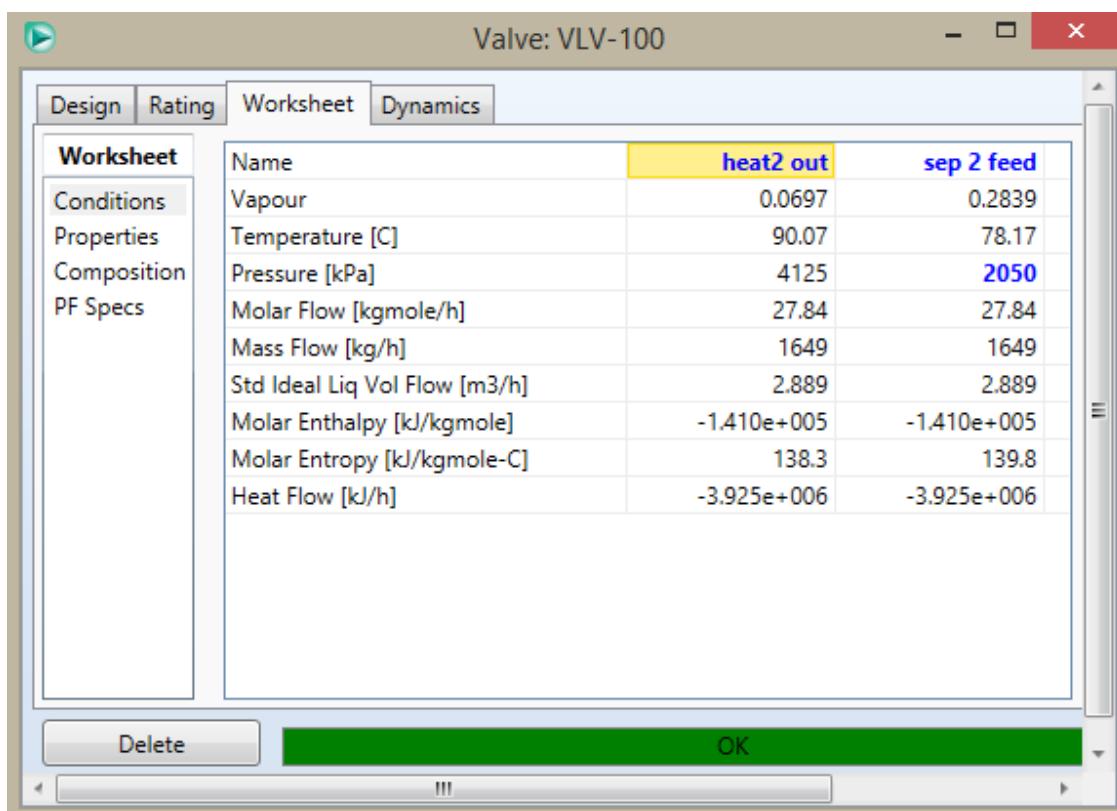
Add the first separator



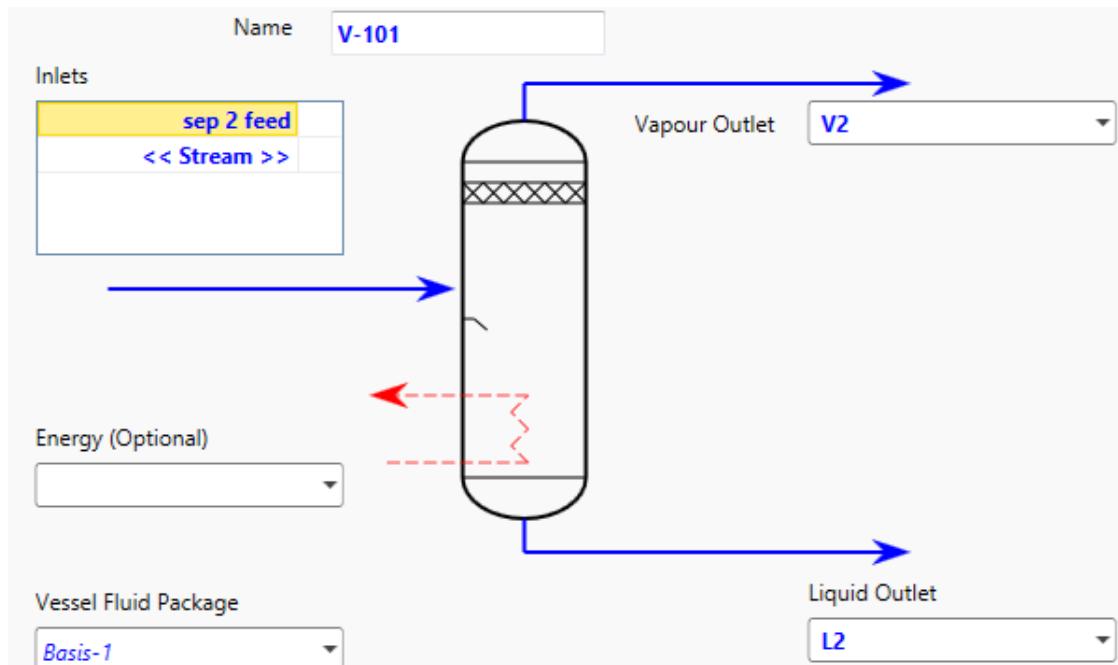
The **liquid** stream is then heated, add a second heater with a pressure drop of 0.0 & duty of  $3.15 \times 10^5$



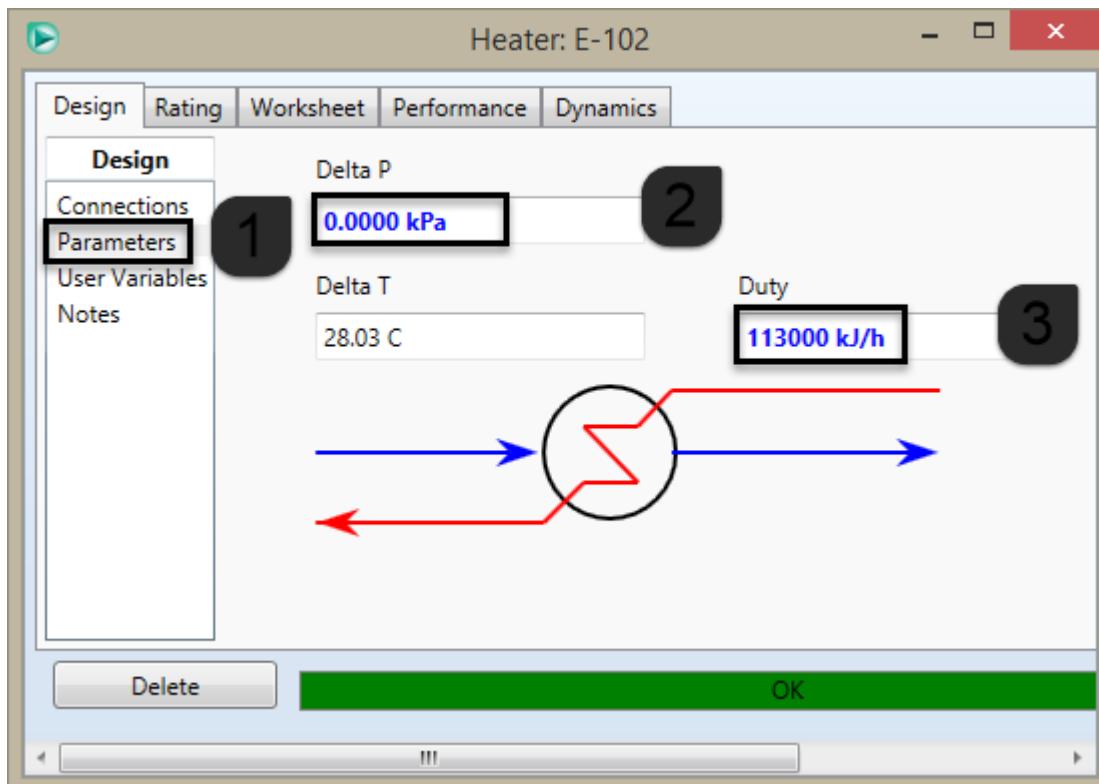
Add a valve with outlet pressure of 2050 kPa



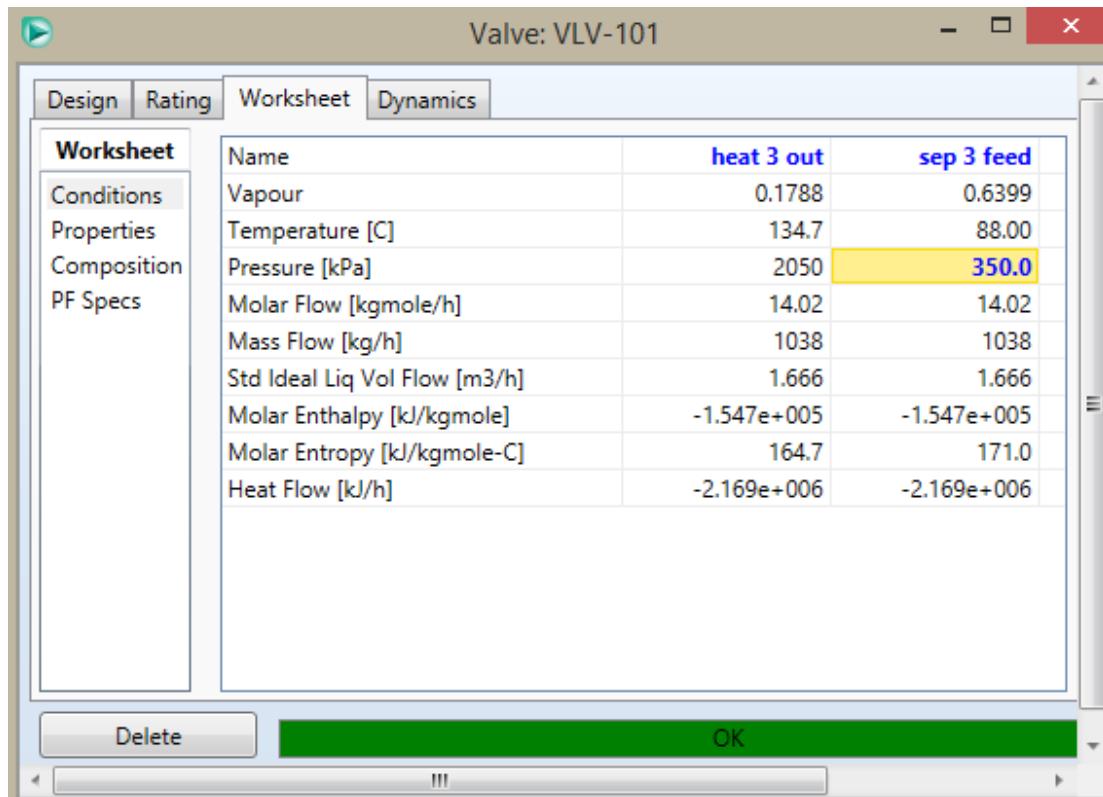
Add the second separator



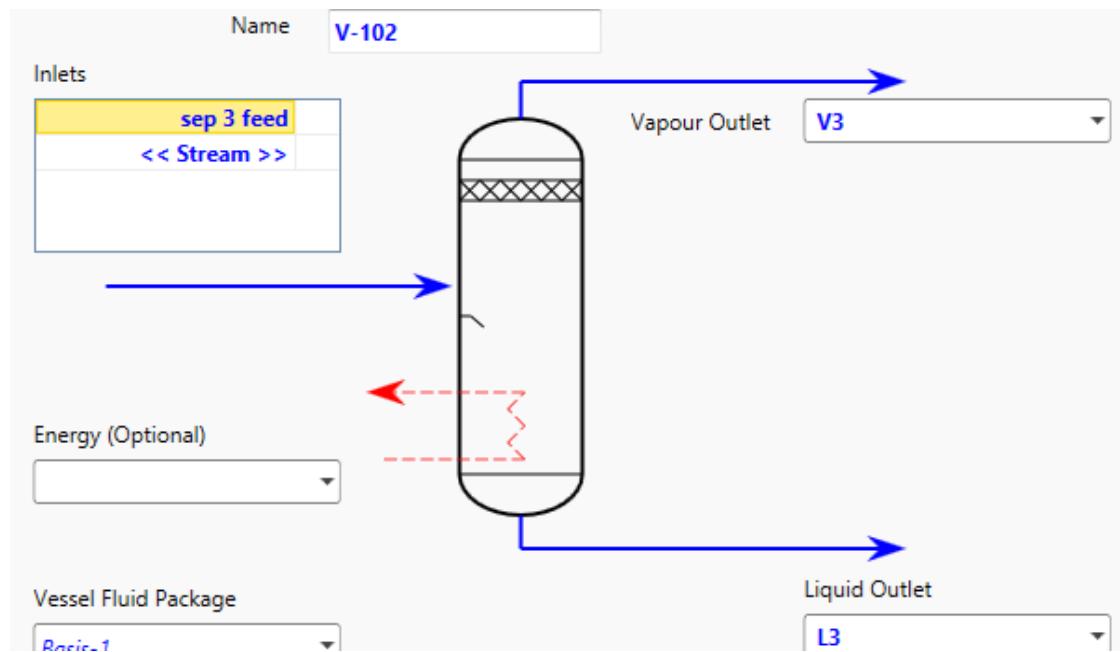
The liquid from the second separator is now fed to a third heater with a pressure drop of 0.0 & duty of  $1.13 * 10^5$



Add a second valve with an outlet pressure of 350 kPa

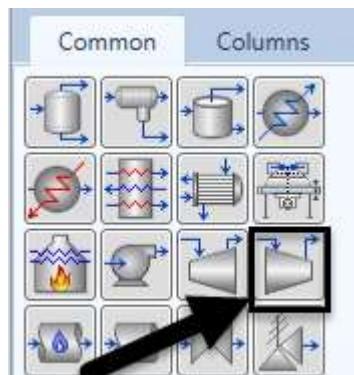


Add the third separator



The **vapor** from the **second separator** is fed to a compressor to raise the pressure to 4125 kPa

Add a compressor from the palette



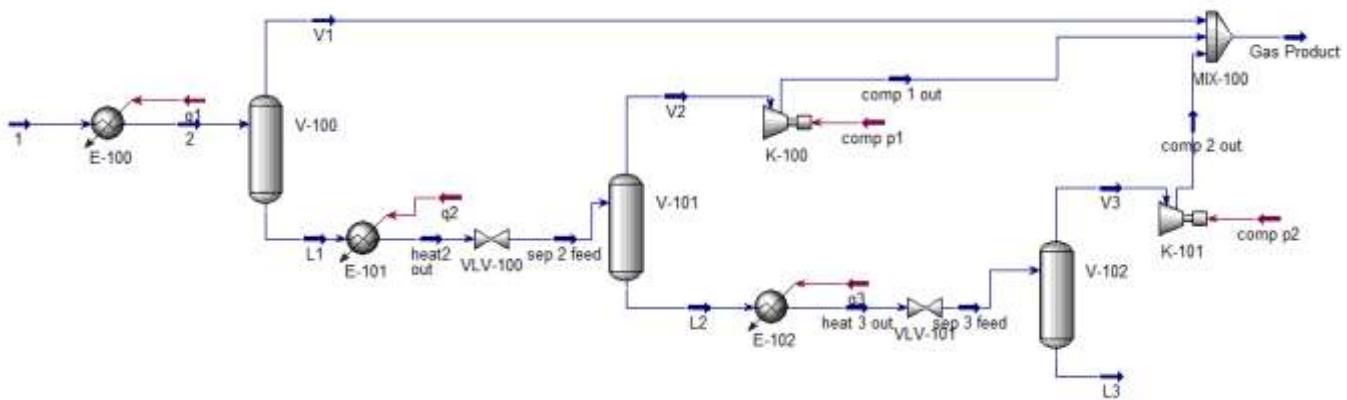
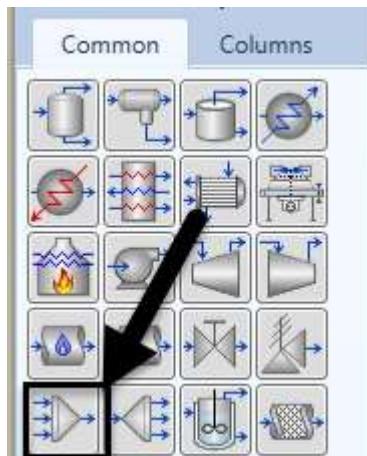
	Name	V2	comp 1 out	comp p1
Conditions	Vapour	1.0000	1.0000	<empty>
Properties	Temperature [C]	106.7	145.9	<empty>
Composition	Pressure [kPa]	2050	4125	<empty>
PF Specs	Molar Flow [kgmole/h]	13.82	13.82	<empty>
	Mass Flow [kg/h]	611.4	611.4	<empty>
	LiqVol Flow [m3/h]	1.223	1.223	<empty>
	Molar Enthalpy [kJ/kgmole]	-1.043e+005	-1.019e+005	<empty>
	Molar Entropy [kJ/kgmole-C]	175.2	176.6	<empty>
	Heat Flow [kJ/h]	-1.441e+006	-1.408e+006	3.313e+004

The **vapor** from the **third separator** is fed to a second compressor to raise the pressure to 4125 kPa

Add a **second** compressor with an outlet pressure of 4125kPa

	Name	V3	comp 2 out	comp p2
Conditions	Vapour	1.0000	1.0000	<empty>
Properties	Temperature [C]	88.00	191.0	<empty>
Composition	Pressure [kPa]	350.0	4125	<empty>
PF Specs	Molar Flow [kgmole/h]	8.970	8.970	<empty>
	Mass Flow [kg/h]	549.9	549.9	<empty>
	LiqVol Flow [m3/h]	0.9460	0.9460	<empty>
	Molar Enthalpy [kJ/kgmole]	-1.270e+005	-1.183e+005	<empty>
	Molar Entropy [kJ/kgmole-C]	183.1	187.8	<empty>
	Heat Flow [kJ/h]	-1.139e+006	-1.061e+006	7.753e+004

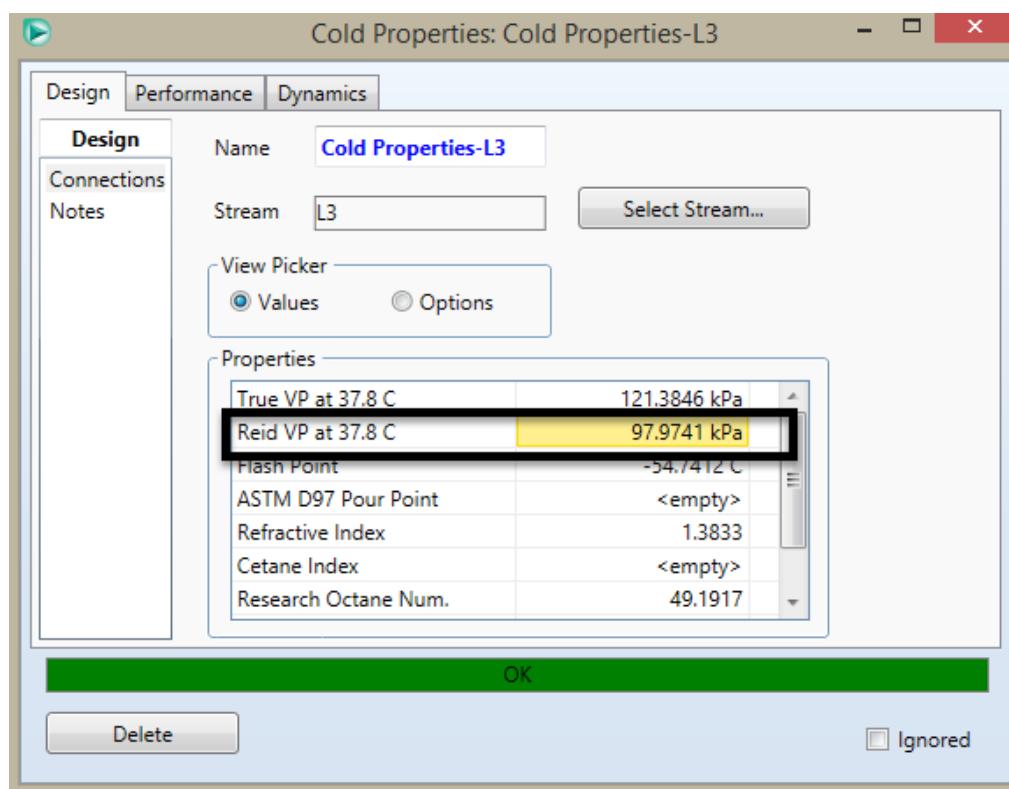
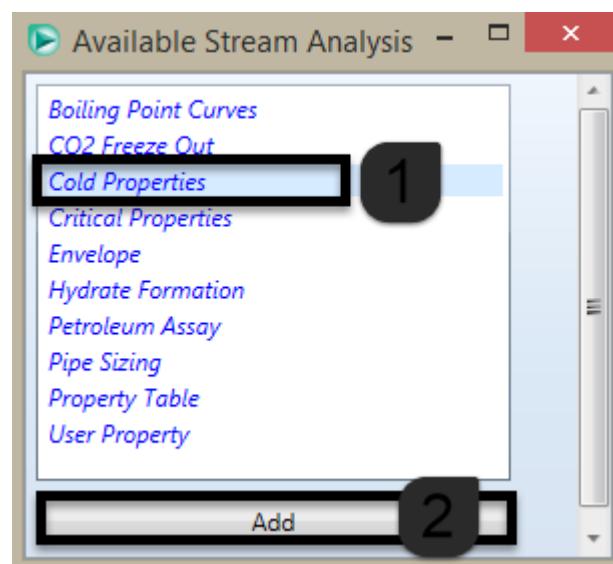
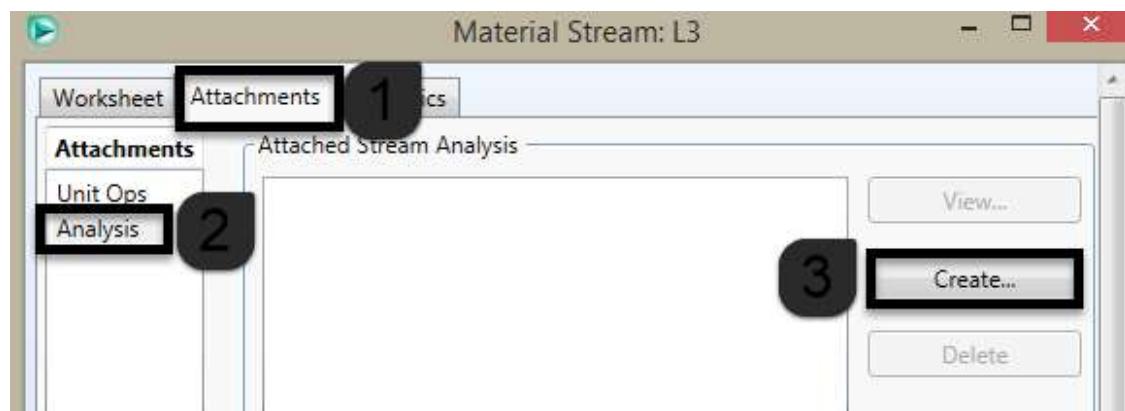
The three vapor streams will be mixed using a mixer



The RVP of the Liquid Product stream should be about 96.5 kPa to satisfy the pipeline criterion.

Use **cold properties** analysis to see the current Reid Vapor Pressure for the liquid product from the third separator:

From the attachments tab, select Analysis and then Create:



## Changing the Units

We need to change the default unit set to fit this case

Change the units as follows:

Molar flow: m<sup>3</sup>\_gas/hr

Liq. Vol. Flow: barrel/hr

Std. Vol. Flow: barrel/hr



Set defaults options for units of measurement

**Display Units**

	Unit	
Molar Concentr.	ppmmol	<b>m3/h_(gas)</b>
Molar Density	kgmole/m3	
Molar Electrical	S-m2/kgmole	
Molar Enthalpy	kJ/kgmole	
<b>Molar Flow</b>	<b>m3/h_(gas)</b>	
Molar Ratio Cor.	mol/mol	
Molar Volume	m3/kmol	

Set defaults options for units of measurement

**Display Units**

	Unit	
Lead Level in Ga	kgPb/m3	
Length	m	
<b>Liq. Vol. Flow</b>	<b>barrel/hr</b>	
Liquid Specific F	sm3/s/Pa	
Mass	kg	
Mass Concentra	wt %	

## Optimization

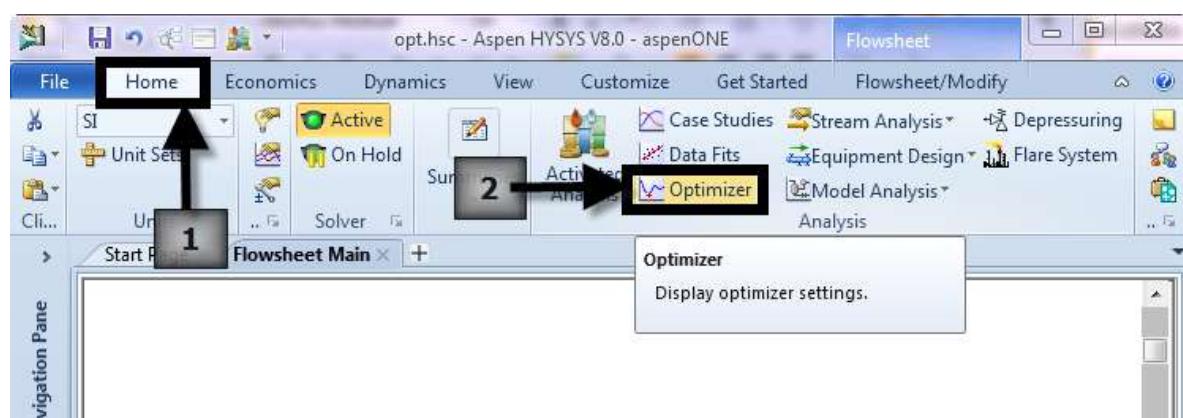
HYSYS contains a multi-variable Steady State Optimizer. Once your flowsheet has been built and converged, you can use the Optimizer tool to find the operating conditions which minimize or maximize an Objective Function. The Optimizer owns its own Spreadsheet for defining the Objective Functions as well as any constraint expressions to be used. This allows you to construct Objective Functions which maximize profit, minimize utilities or minimize exchanger UA.

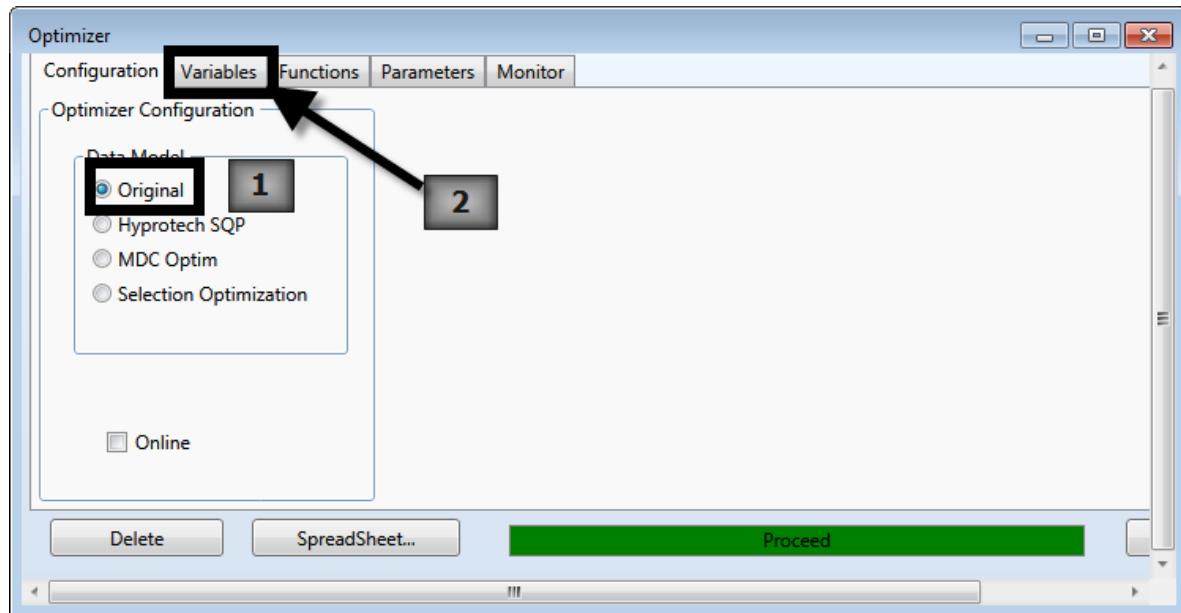
In this case, we want to maximize the total operating profit while achieving an RVP of Liquid Product less than 96.5 kPa. The incomes from the Plant are both the Gas and Liquid Products. The operating costs are the Steam Costs for each Heater plus the Power Cost for each Compressor.

Profit = Income - Cost

Profit= (Gas Product + Liquid Product) – (Steam Costs + Compression Cost)

Use the Optimizer tool:

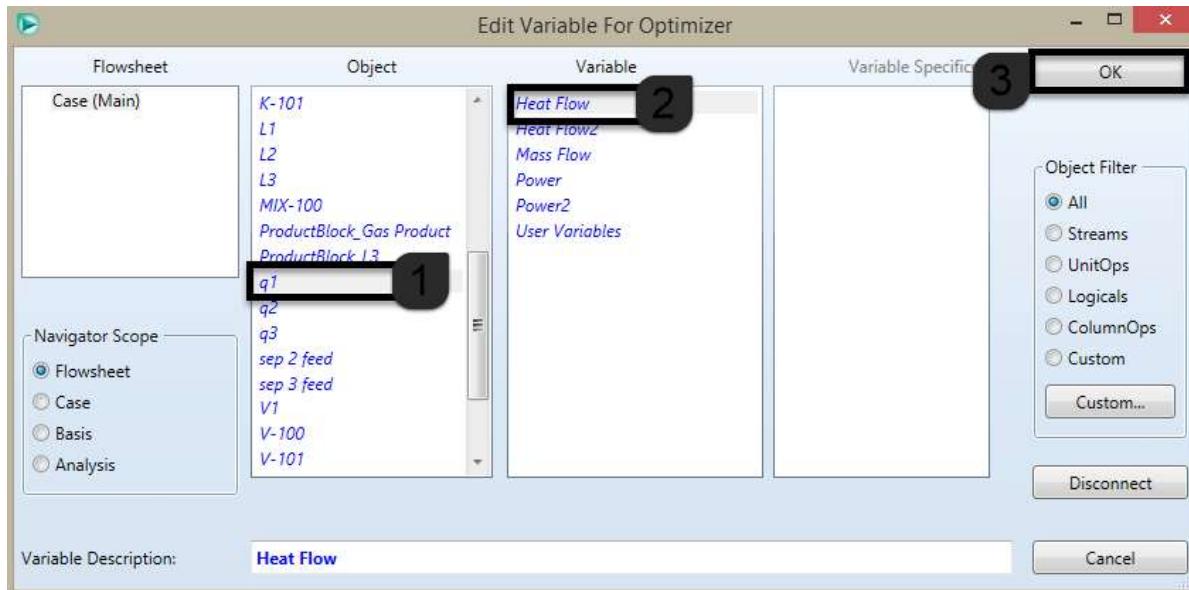




The variables to be optimized in order to maximize the profit should be added now, these variables are:

The 3 Heaters' duties and Valves outlet pressures



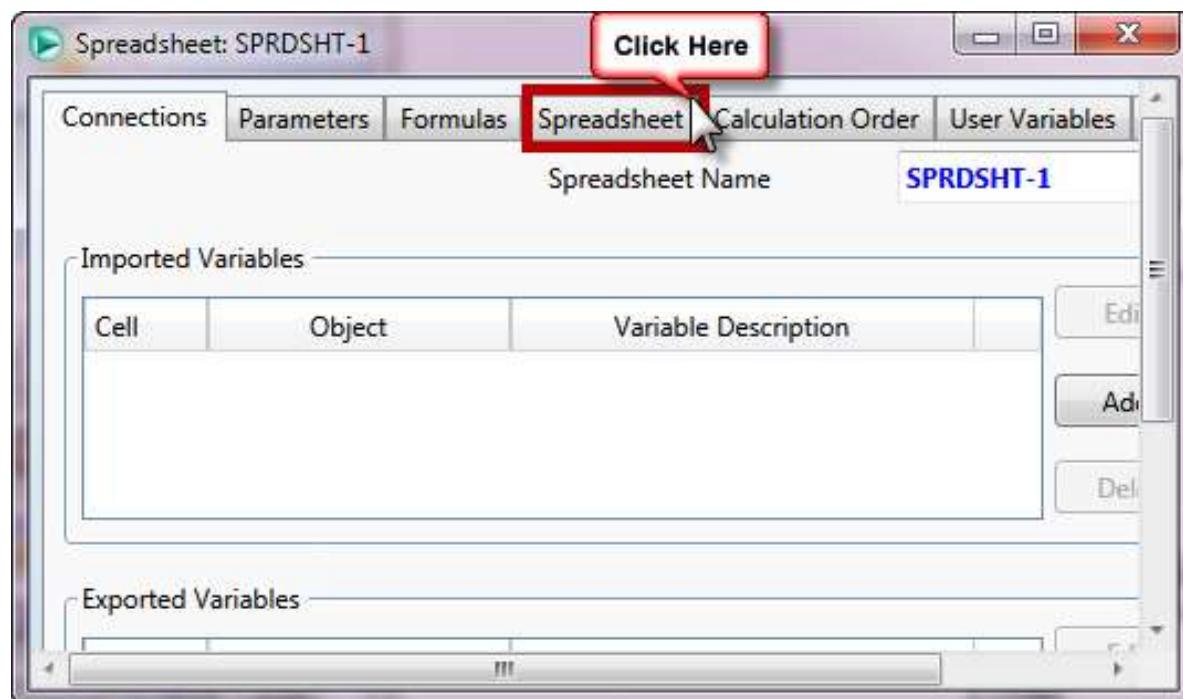
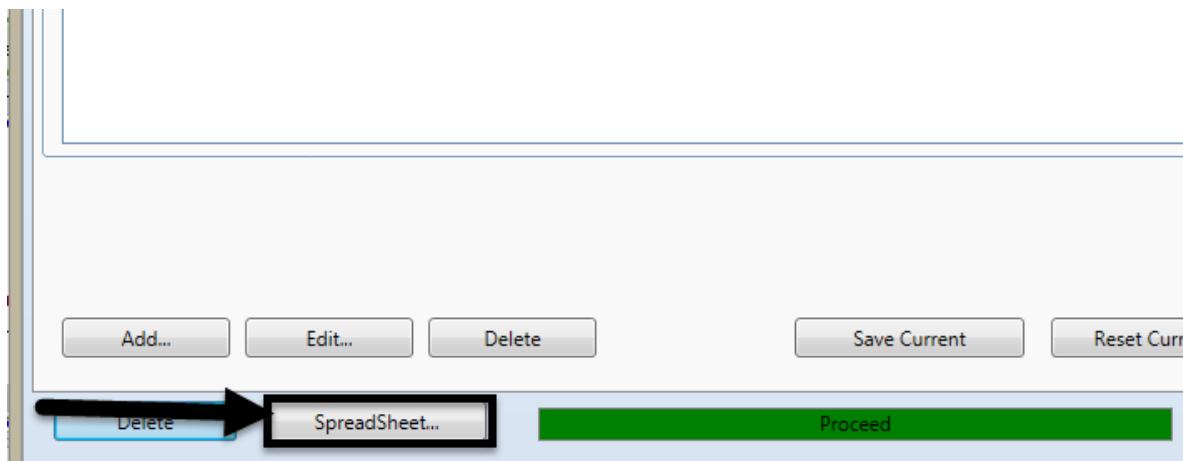


After adding all the 5 variables, set the upper and lower ranges for each variable as follows:

Adjusted (Primary) Variables							
Object	Variable Description	Low Bound	Current Value	High Bound	Reset Value	Enabled	
q1	Heat Flow	0.0000	4.250e+005	1.000e+006	<empty>	<input checked="" type="checkbox"/>	
q2	Heat Flow	0.0000	3.150e+005	1.000e+006	<empty>	<input checked="" type="checkbox"/>	
q3	Heat Flow	0.0000	1.130e+005	1.000e+006	<empty>	<input checked="" type="checkbox"/>	
sep 2 feed	Pressure	650.0	2050	3500	<empty>	<input checked="" type="checkbox"/>	
sep 3 feed	Pressure	70.00	350.0	1000	<empty>	<input checked="" type="checkbox"/>	

The Optimizer has its own Spreadsheet for defining the Objective and Constraint functions.

Now we have to start building the profit module using the spreadsheet operation:



Profit= Income - Cost

Profit= (Gas Product + Liquid Product) – (Steam Costs + Compression Cost)

Prices & costs:

Oil Price= 15 \$/bbl

Gas Price =0.106 \$/m<sup>3</sup>\_gas

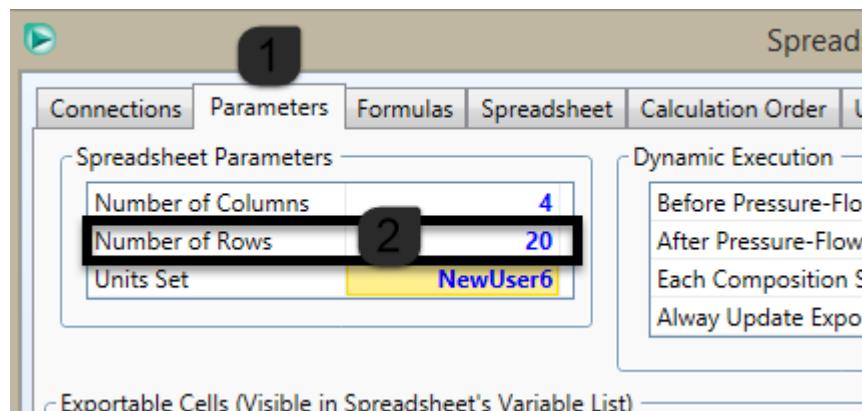
Steam Cost= 0.682 \$/kW-h

Compression Cost=0.1 \$/kW-h

The RVP spec for the liquid should be added in the spread sheet in order to use it as a constraint.

RVP spec= 96.5 kPa

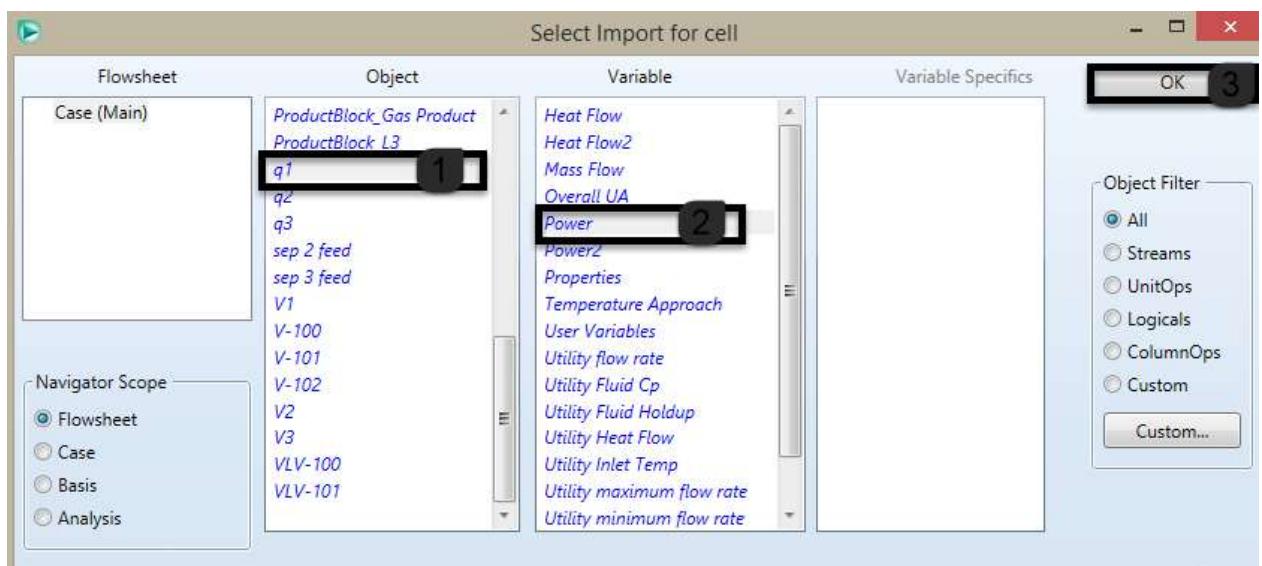
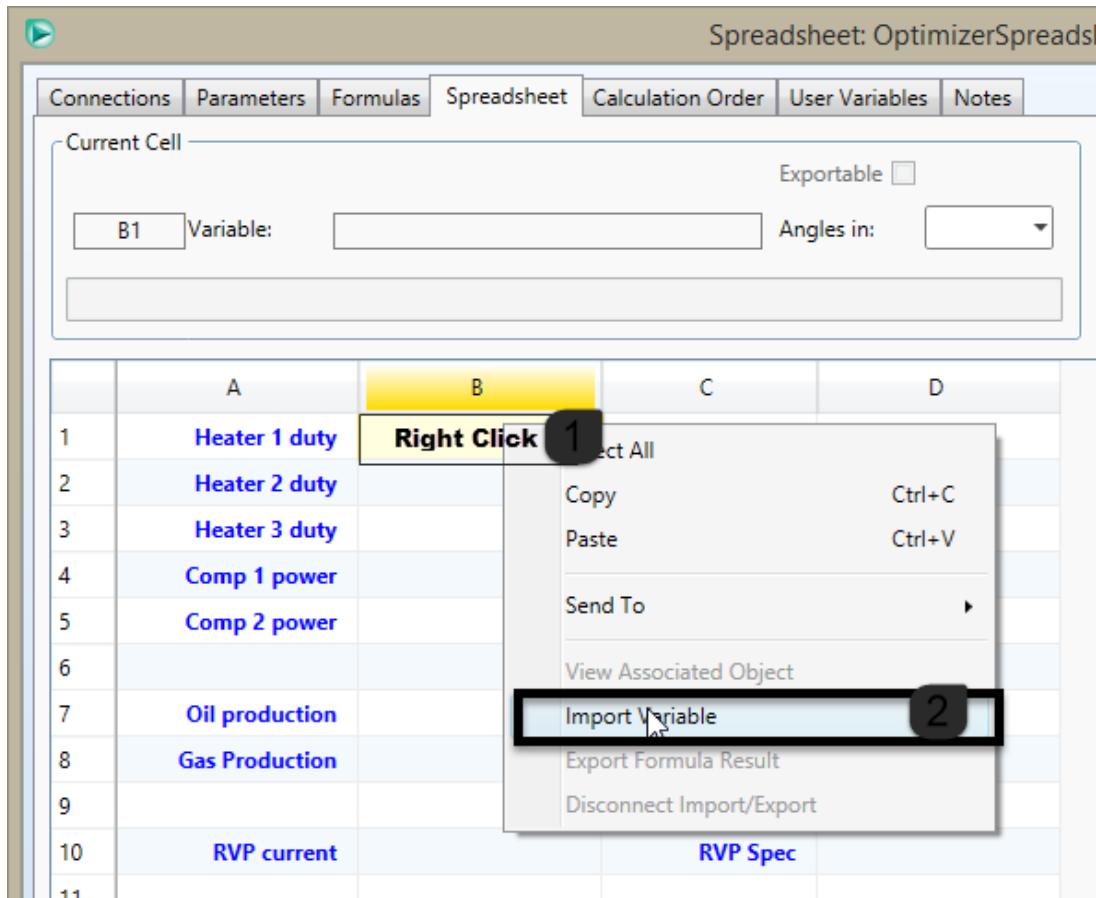
First we need to increase the number of rows in the spread sheet to be 20 from the parameters tab:



Now return to the spreadsheet tab again and add all the profit equation parameters & variables.

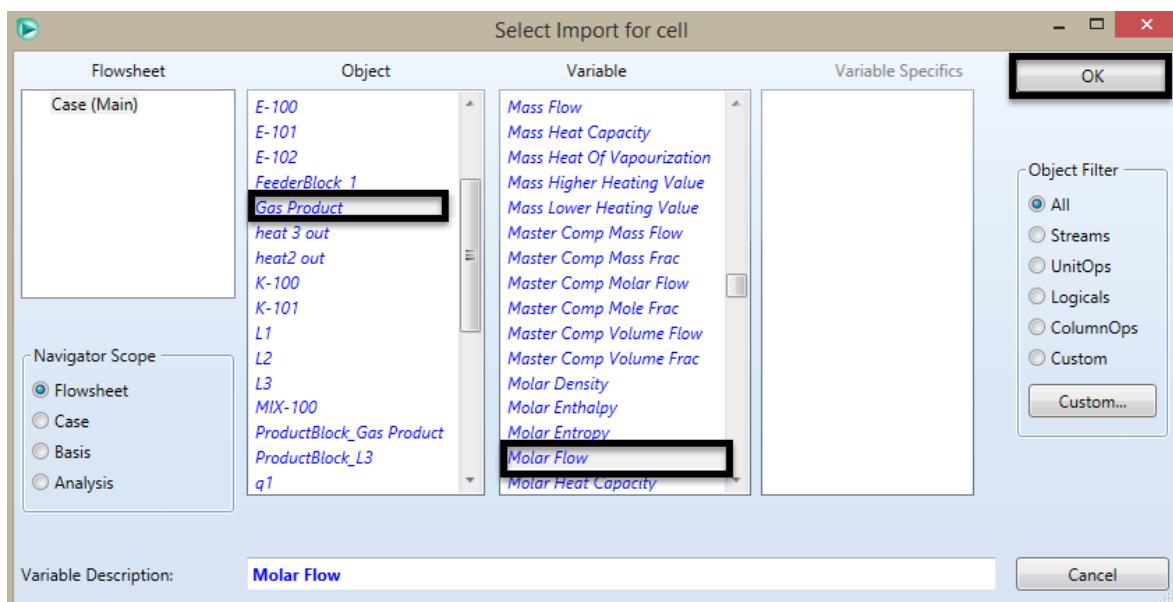
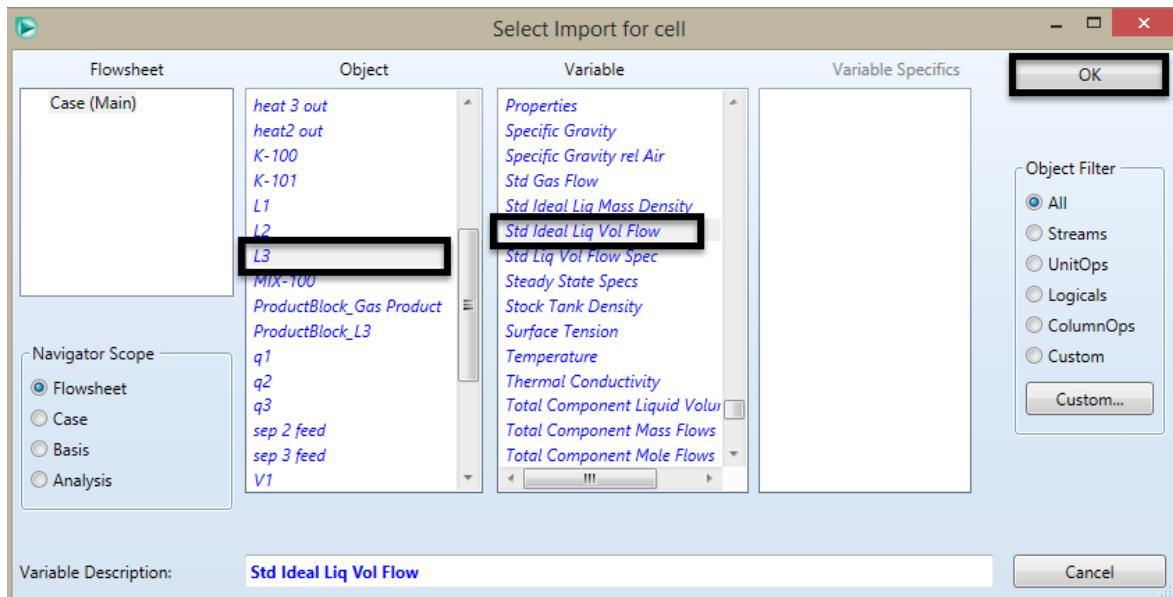
	A	B	C	D
1	Heater 1 duty		Steam Cost	
2	Heater 2 duty			
3	Heater 3 duty			
4	Comp 1 power		Compression Cost	
5	Comp 2 power			
6				
7	Oil production		Oil Price	
8	Gas Production		Gas Price	
9				
10	RVP current		RVP Spec	
11				
12			Income	
13			Cost	
14			Profit	

After adding the labels, import the variables in each labeled cell:

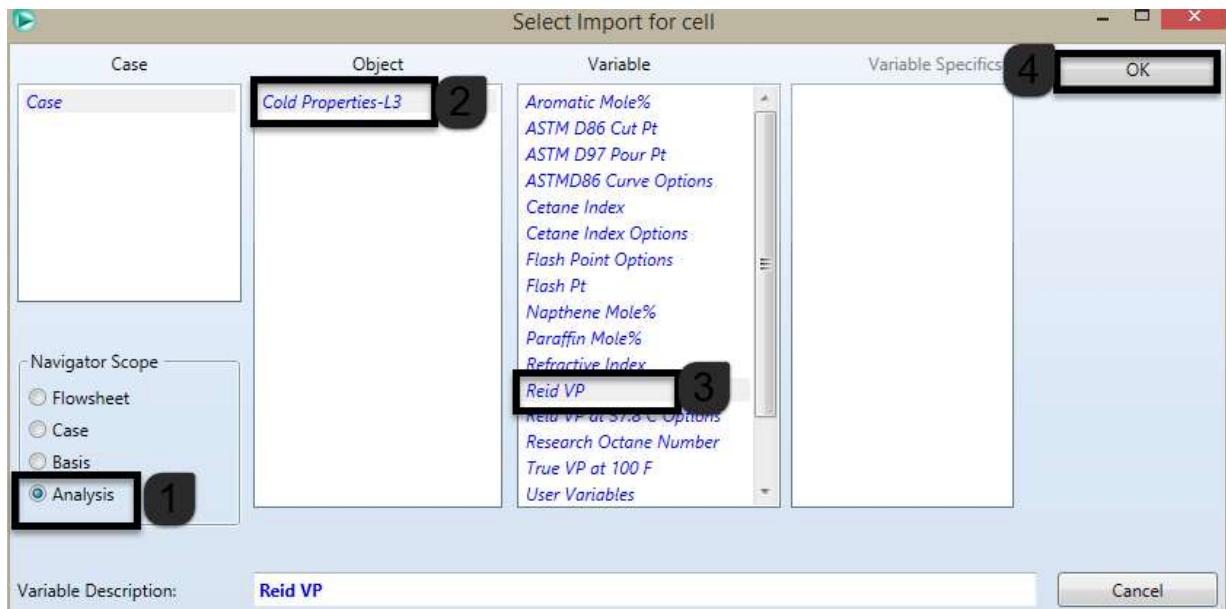


Add all duties for heaters & compressors the same way as we did in the previous step.

Do the same for Oil & gas production flow rates



The RVP current value should be imported from the Analysis:



The Prices should now added manually without importing it

	A	B	C	D
1	Heater 1 duty	118.1 kW	Steam Cost (\$/kW-h)	0.6820
2	Heater 2 duty	87.50 kW		
3	Heater 3 duty	31.39 kW		
4	Comp 1 power	9.203 kW	Compression Cost (\$/kW-h)	0.1000
5	Comp 2 power	21.54 kW		
6				
7	Oil production	4.530 barrel/hr	Oil Price (\$/bbl)	15.00
8	Gas Production	1056 m3/h_(gas)	Gas Price (\$/m3_gas)	0.1060
9				
10	RVP current	97.97 kPa	RVP Spec (kPa)	96.50
11				
12			Income	
13			Cost	
14			Profit	

Calculate the Income, Cost & Profit using the current formulas:

Income: =d7\*b7+d8\*b8

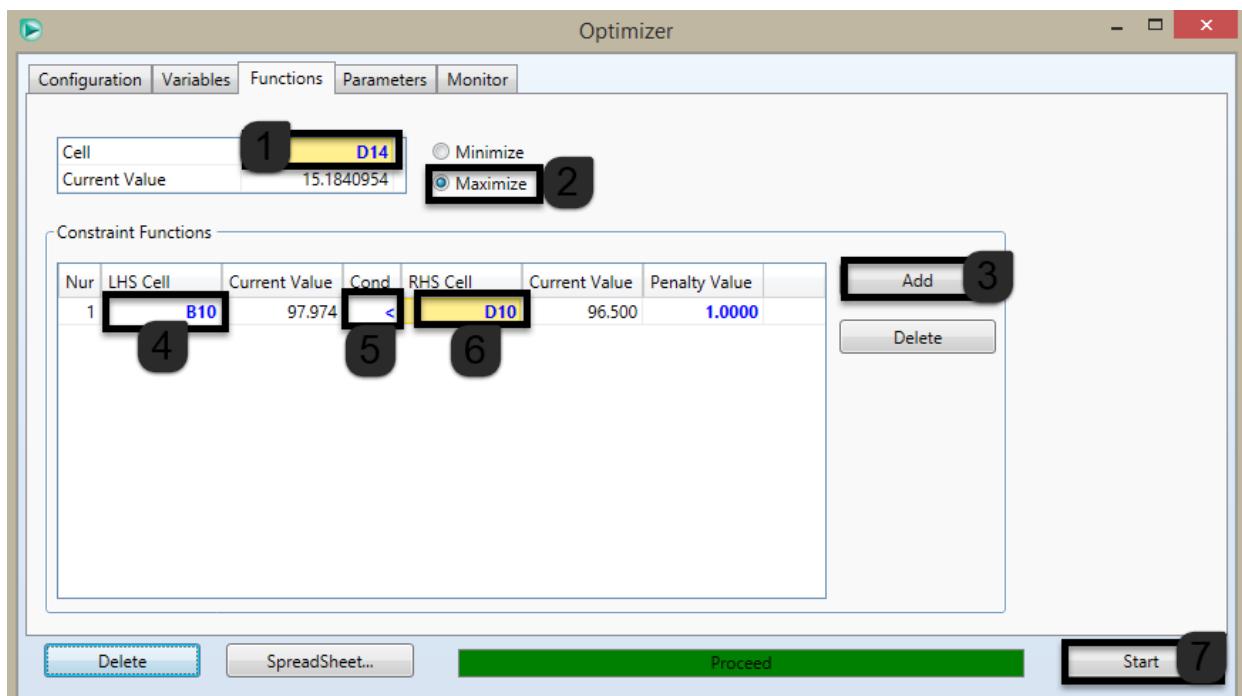
Cost: =(b1+b2+b3)\*d1+(b4+b5)\*d4

Profit: =d12-d13

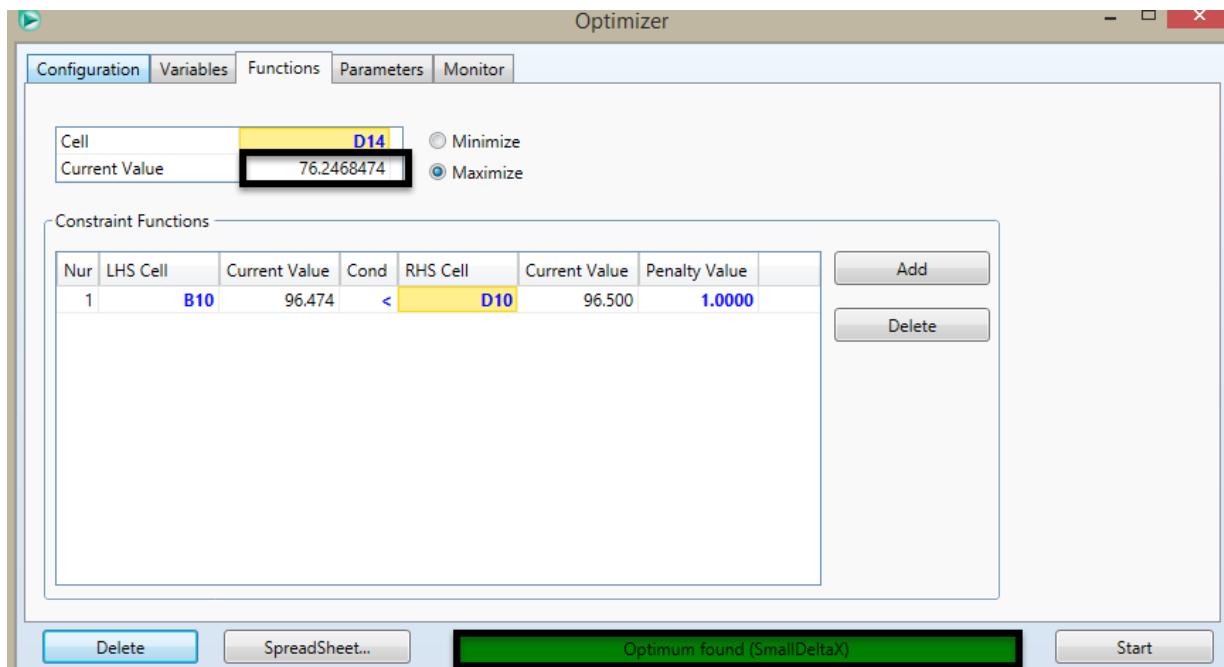
	A	B	C	D
1	Heater 1 duty	118.1 kW	Steam Cost (\$/kW-h)	0.6820
2	Heater 2 duty	87.50 kW		
3	Heater 3 duty	31.39 kW		
4	Comp 1 power	9.203 kW	Compression Cost (\$/kW-h)	0.1000
5	Comp 2 power	21.54 kW		
6				
7	Oil production	4.530 barrel/hr	Oil Price (\$/bbl)	15.00
8	Gas Production	1056 m <sup>3</sup> /h_(gas)	Gas Price (\$/m <sup>3</sup> _gas)	0.1060
9				
10	RVP current	97.97 kPa	RVP Spec (kPa)	96.50
11				
12			Income (\$/hr)	179.9
13			Cost (\$/hr)	164.7 kW
14			Profit (\$/hr)	15.18 kW

After calculating the operating profit @ current conditions, use the optimizer to maximize the profit by changing the 5 variables which we added before

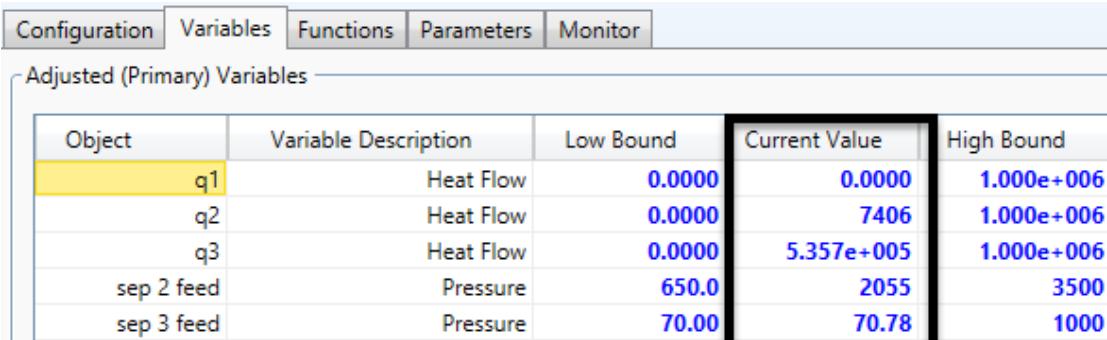
Open the optimizer and go to Functions tab:



You may need to press the Start button 2 or 3 times to ensure reaching the optimum solution as follows:



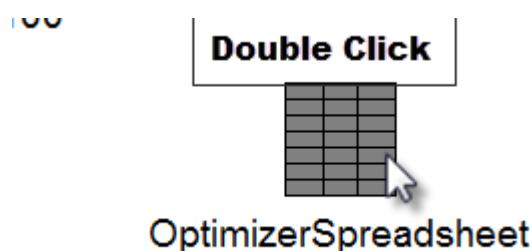
You may see the new values for the variables from the Variables tab:



The screenshot shows the Variables tab. It displays a table of 'Adjusted (Primary) Variables' with columns: Object, Variable Description, Low Bound, Current Value, and High Bound. The table rows are:

Object	Variable Description	Low Bound	Current Value	High Bound
q1	Heat Flow	0.0000	0.0000	1.000e+006
q2	Heat Flow	0.0000	7406	1.000e+006
q3	Heat Flow	0.0000	5.357e+005	1.000e+006
sep 2 feed	Pressure	650.0	2055	3500
sep 3 feed	Pressure	70.00	70.78	1000

Now you can return back to the spreadsheet to observe the results:



	A	B	C	D
1	Heater 1 duty	0.0000 kW	Steam Cost (\$/kW...)	0.6820
2	Heater 2 duty	2.057 kW		
3	Heater 3 duty	148.8 kW		
4	Comp 1 power	3.797 kW	Compression Cost...	0.1000
5	Comp 2 power	109.4 kW		
6				
7	Oil production	5.403 barrel/hr	Oil Price (\$/bbl)	15.00
8	Gas Production	1032 m3/h_(gas)	Gas Price (\$/m3_g...	0.1060
9				
10	RVP current	96.47 kPa	RVP Spec (kPa)	96.50
11				
12			Income (\$/hr)	190.5
13			Cost (\$/hr)	114.2 kW
14			Profit (\$/hr)	76.25 kW

**Save Your Case!**

## Exercise:

One thing you may notice with the Optimized solution is that the Pressure of V3 has been decreased to 70 kPa (10 psia) which is less than atmospheric. This is not a desired condition for the inlet of a compressor. The inlet of the second compressor, cannot be less than 125 kPa (19 psia). What is the maximum profit if you adhere to this guideline?

You may check the website for more data &  
courses

<http://www.adeyab.com>

If you have any inquiry:

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[adeyab@adeyab.com](mailto:adeyab@adeyab.com)

**002 01227549943**

**00966544390045**