LPG COLUMN DESIGN AND SIMULATION

(A Report on Project)

Submitted as a part of course work in M. Tech (Gas Engineering)

By

76

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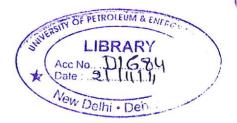
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LPG COLUMN DESIGN AND SIMULATION

A Project report submitted to the University of Petroleum and Energy Studies-Rajahmundry campus in partial fulfillment of the requirement for the degree of

MASTER OF TECHNOLOGY (GAS ENGINEERING)

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APRIL - 2009



DECLARATION

This is to declare that the project report entitled "LPG Column Design and Simulation" has been prepared and submitted by me in all aspects, in partial fulfillment of the requirement for the award of the degree of Master of Technology [Gas Engineering] in University of Petroleum and Energy Studies-Rajahmundry.

The content of this report has not been submitted to any university or institution by me for the award of any degree or diploma.

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CERTIFICATE

This is to certify that the project work entitled "LPG Column Design and Simulation" being submitted by Mr.Saravana Raj.N (R030307011), in partial fulfillment of the requirement for the award of the degree of Master of Technology [Gas Engineering] in University of Petroleum and Energy Studies-Rajahmundry, is a bonafide project work carried out by him under my guidance.

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ACKNOWLEDGEMENTS

I wish to acknowledge my sincere thanks to my mentor and guide **Prof. K.V.Rao**, Academic Head, College of Engineering Studies, UPES - Rajahmundry for his valuable advice and resourceful guidance.

I would like to thank Dr. K.V.Mohana Rao, Centre Director, Rajahmundry Regional Centre, UPES. Mr. J.S Prasad, Course Coordinator - M.Tech (Gas Engineering) for his constant support and guidance.

(N.SARAVANA RAJ)

EXTENDED ABSTRACT

A project report to design LPG Column located in Gas Processing Plant which is processing 3 million cubic meters/day for recovering LPG. All the calculations are performed based upon particular composition which is given for this project.

Gas Processing plant has two UNITs, namely Crude Stabilisation Unit and Gas Processing Unit. The associated natural gas contains sour/rich gas as well unsaturated crude, the unsaturated crude is treated in CSU to get saturated crude and the sour/rich gas is processed in gas processing unit to get LPG, Naphtha, C₂C₃ and Lean gas.

Here our focus is mainly LPG Distillation Column which is part of LPG plant, The plant operation is as follows, Sweetened gas from GSU flows to knock out drum, from where liquid present is separated, then the gas is sent through exchangers and precooled. The precooled gas will be sent to KOD where liquefied hydrocarbon and water are separated out and the gas flows to the molecular sieve drier where the moisture is reduced, the dried gas is passed through filters, to retain any foreign or dust particles. Then the dried gas flows through 1st stage chiller. The process fluid next enters in 1st stage liquid separator where the partially condensed hydrocarbon liquids separated out. Gas vapours from the top of the vessel flows through 2nd stage chiller. Then the H/C fluid enters in 2nd stage liquid separator where 2nd stage liquid is separated. The Second Stage Vapours (SSV) after separation of liquid hydrocarbon (Condensate) are passed through and delivered as feed stock to C₂C₃ Recovery Unit. Also SSV can be sent directly to consumers through bypass line if C₂C₃ Recovery Unit is under shut down.

COLD BOX which serves to chill the feed gas by exchanging heat with various cold streams of the plant. Separator liquid from feed gas separator sent to

LEF column to remove the lighter fractions. The bottom liquid goes to LPG column. If ethane propane plant is under shutdown the LEF gases can be sent to consumer line after compressing through residue gas compressor. Also these gases can be used for regeneration of gas dryers.

The bottom liquid of the LEF column is allowed to enter the LPG column. The top product is Liquefied petroleum gas and bottom product is Natural gas liquids.

This particular LPG column is design using manually (SPREADSHEETS) as well as simulator (CHEMCAD 6.0.1). The various Tray columns (sieve – tray column, value tray column and Bubble cap tray column) and also packed column sizing are designed by CHEMCAD 6.0.1 Simulator. Costing for these different columns are prepared in CHEMCAD 6.0.1 Simulator and finally the cost comparison chart is carried out using SPREADSHEETS.

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1. INTRODUCTION

LPG is a mixture of light hydrocarbons which are gaseous at normal temperatures and pressures, and which liquefy readily at moderate pressures or reduced temperature. It is odourless and so, for safety reasons, a pungent compound, ethyl mercaptan, is added to make any leaks easily detectable.

Motor vehicles run on propane or a mixture of propane and butane. Propane and butane are gases at atmospheric pressure and temperature but can easily be liquefied for storage by an increase in pressure. The most common blend by volume is 60 per cent propane to 40 per cent butane. Industry uses propane or butane and the petrochemical industry may use both as a feedstock. Households use only propane.

1.1 History

Liquefied petroleum Gas is a mixture of $C_2 - C_3$ fractions. It is recovered from refinery Gases and Natural gas streams. It is essentially a twentieth century development long before time. Use has been made of gas compressed containers. Limited amounts of this product were sold in cylinders in England as early as 1810. In 1870, pintsch gas was developed and later used in railway car lighting. Around 1907. blangas formed by craking oil, was liquefied by compressing it to about 1800 psig. Several containers in the United States, despite attendant difficulties such as high transportation cost and complexity of equipment needed for its utilization.

1.2 Scope

Crude oil is a combustible liquid consisting of a mixture of hydrocarbon and compounds such as O₂, S, and N₂. Unstabilised crude oil is received through pipeline from offshore platforms (after two stage separation). A project report to design LPG Column located in Gas Processing Plant, processing 3 million cubic meters/day for recovering LPG. Important scope of this project is design, simulation and cost analysis done by using CHAMCAD 6.0.1 simulator which is capable of simulating any kind of feed composition.

1.3 Limitation

Here design project is for particular composition and feed condition which is mentioned in table 1, this designed column will not be support unless the feed condition and composition ,the basic parameters of distillation column like diameter, height, number of stages, wall thickness etc are vary with feed composition to composition as well as feed conditions. CHEMCAD 6.0.1 simulator is limited upto only 50 number of stages. More than that is not possible in this version.

1.4 Delimitation

Since project is done using CHEMCAD 6.0.1, any kind of feed condition and composition can be specified in simulator to sizing the Distillation column as well as analysis on characteristics of tower, feed stream and product stream. This simulator is capable of designing sieve tray, value tray, bubble cap tray and also packed column, this CHEMCAD 6.0.1 is special to costing of equipment and economics of plant.

1.5 Feed Composition of Gas Processing Plant

Table 1: Feed Composition Gas Processing Plant

| Components | Molecular weight | Mole % | Moles | Kgs |
|----------------------------------|--------------------------|----------|----------|-------------------|
| N ₂ | 28 | 1.42 | 79.3847 | 2222.7716 |
| CO ₂ | 44 | 1.75 | 97.83324 | 4304.6256 |
| CH ₄ | 16 | 69.24 | 3870.842 | 61933.472 |
| C ₂ H ₆ | 30 | 13.05 | 729.5564 | 21886.692 |
| C ₃ H ₈ | 44 | 9.27 | 518.2366 | 22802.4104 |
| i C ₄ H ₁₀ | 58 | 1.66 | 92.8018 | 5382.5044 |
| n C ₄ H ₁₀ | 58 | 2.35 | 131.3761 | 7619.8138 |
| i C ₅ H ₁₂ | 72 | 0.59 | 32.9838 | 2374.8336 |
| n C ₅ H ₁₂ | 72 | 0.67 | 37.4562 | 2696.8464 |
| Average mo | lecular weight = 23.4728 | <u> </u> | t | otal = 131224 Kgs |
| Water in the Fe | eed = 118.46 Kg/ (106Nm | n3/day) | | |
| Associated gas | = 3 x 118.46 Kg = 355.3 | 8 | | |

2. LITERATURE REVIEW

2.1 Growth of Gas Processing Plant [8]

Crude oil is a combustible liquid consisting of a mixture hydrocarbon and compound O_2 , S, and N_2 . Unstabilised crude oil is received through pipeline from offshore platforms (after two stage separation). Stabilisation of crude oil is being done with three stage separation having intermediate dehydration unit. Here Table 2 is explain about growth of Gas Processing Plant.

Table 2: Growth of Gas Processing Plant.

| TECHNOLOGICAL SCHEME PHASES | YEAR INSTALLED |
|--------------------------------|---------------------------------------|
| Technological Scheme Phase I | 1981 (CSU I & LPG I) |
| | CSU I - Crude Stabilization Unit I |
| | LPG I - Liquefied Petroleum Gas I |
| Technological Scheme Phase II | 1984 (CSU II, LPG II & GT I & II) |
| | CSU II - Crude Stabilization Unit I |
| | LPG II - Liquefied Petroleum Gas I |
| | GT I & II - Gas Turbine I & II |
| Technological Scheme Phase III | 1990(GSU12 &13 ,EPRU,CFU I) |
| | GSU12 &13 - Gas Sweetening Unit 12&13 |
| | EPRU - Ethane Propane Recovery Plant |
| | CFU I - Condensate Fraction Unit I |
| Minas Plant & CFU II | 1994 |
| Gas Turbine III | 2001 |

2.2 Overview of Gas Processing Plant [1]

Following figures (figure 1 & 2) are explaining the overview idea about Gas Processing Plant, and also process description have described.

GAS PROCESSING PLANT

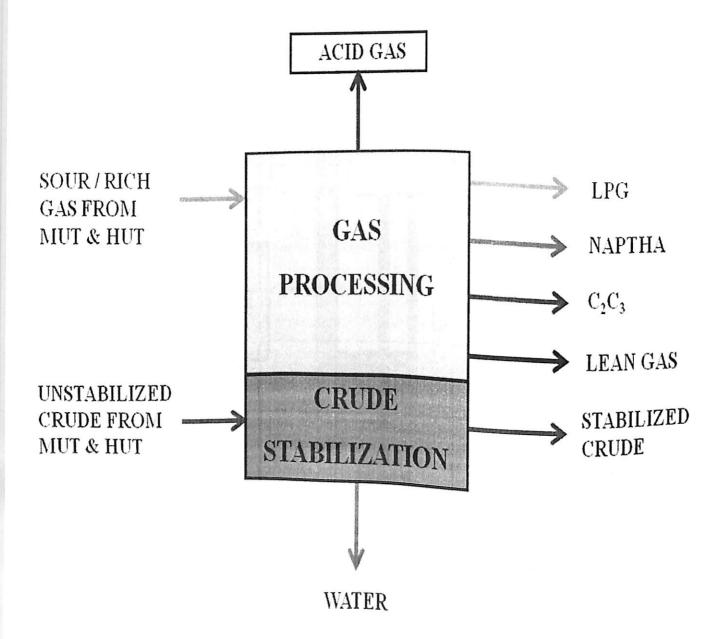


Figure 1: Simple Overview of Gas Processing Plant

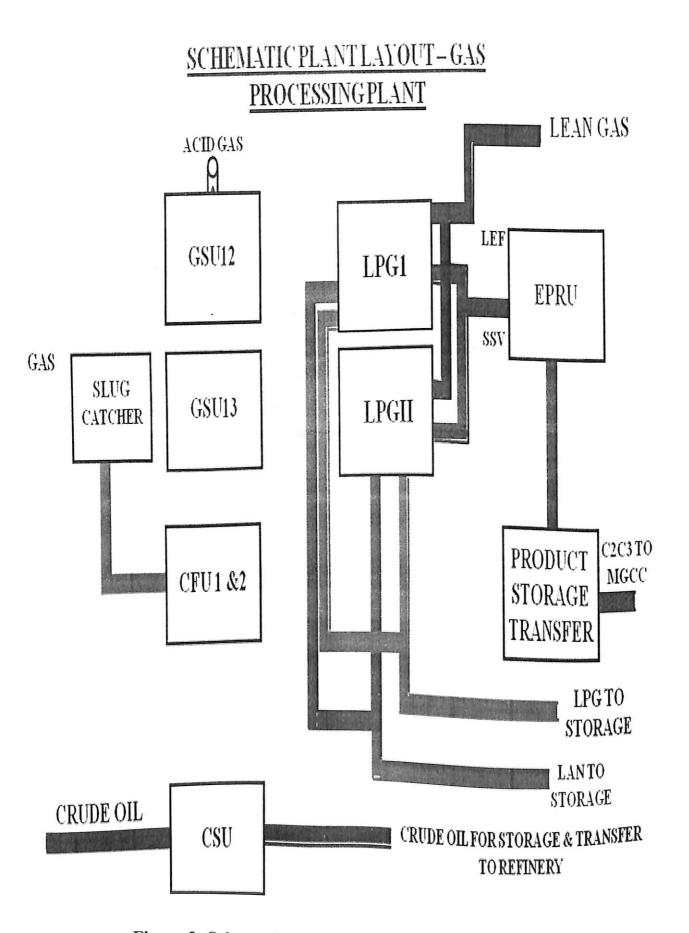


Figure 2: Schematic Plant Layout - Gas Processing Plant

2.2.1 Slug Catcher

It is based on the reducing of the fluid velocity and subsequent gravity separation. To hold the slug liquid reaching URAN at the time of pigging of gas pipeline. Provides room for gas - condensate separation and storage volume of liquid condensate. To continuously send the separated liquid condensate to condensate fractionation unit I & II for further processing.

To partially stabilize the liquid and inject into crude stabilization unit in case of CFU I & II or both in shutdown.

2.2.1 Crude Stabilization Units

Stabilisation of crude oil is required to minimize oil loss through evaporation during storage and transportation for which volatile components are removed. Solids and Salts present in crude oil is removed to minimize corrosive and abrasive wear of equipments. Water content of crude oil is reduced to minimize the cost of pumping and ensure BS& W requirement of refineries. Crude oil is stored in tanks and pumped to refineries as per their requirement.

2.2.3 Crude Sweetening Unit

Removal of undesirable materials such as acid gas includes mainly H₂S and CO₂. Acid gas is highly toxic. It is highly corrosive in presence of water. Forms hydrates and dry ice in cryogenic processes causing serious chocking and corrosion.

Commercial sweetening processes:

- 1. Chemical absorption
- 2. Physical absorption
- 3. Hybrid.

The absorption is at lower temperature and higher pressure and regeneration at high temperature and low pressure. Sulfinol solution is being recycled.

Criteria for process selection is depend on the concentration of impurities in the gas and the degree of removal desired.

GSU plant has two unit GSU 12 and GSU13.the absorption is at lower temperature and higher pressure and regeneration at high temperature and low pressure. Sulfinol solution is being recycled.

2.2.4 Condensate Fraction Unit

Stripping of lighter fraction and separating LPG and LAN with the help of fractionating column from condensate.

2.2.5 LPG Plant

LPG plant is recovering a maximum of butane – propane mixture from gases. The process involves cooling of gases under pressure to condense out LPG & natural gas liquid. The noncondensable gas goes as feed stock of C_2 - C_3 plant propane

2.2.6 Ethane Propane Recovery Plant

 C_2 – C_3 feed streams are cooled to partially condensed them. the refrigeration is provided by passing the high pressure feed stream through expander and by a propane refrigeration system, the condensed feeds are fed in demethaniser to separate methane from top and C_2 - C_3 mix from bottom. C_2 - C_3 stored in spheres and pumped to MGCC NAGOTHANE for their petrochemical use.

2.3 Process Details of LPG – I Plant [1]

MAJOR SECTIONS OF LPG-I PLANT

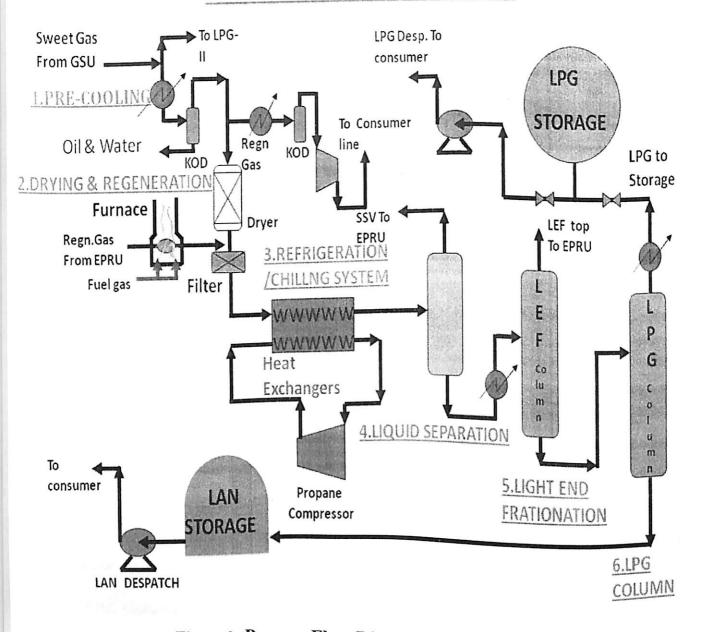


Figure 3: Process Flow Diagram of LPG Plant

2.3.1 Pre cooling & Drying:

Sweetened gas from GSU flows to knock out drum, from where any liquid present is separated, then the gas is sent through exchangers and precooled to 25° C. The precooled gas will be sent to KOD where liquefied hydrocarbon and water are separated out and sent to CSU through . The gas then flows to the molecular sieve drier where the moisture is reduced to less than 5ppm or a due point of -60° C.

2.3.2 Refrigeration:

The dried gas is passed through filters, to retain any foreign or dust particles. Then the dried gas flows through, and cooled to -22^{0} C in the 1st stage chiller. The process fluid next enters in 1st stage liquid separator where the partially condensed hydrocarbon liquids separated out. Gas vapours from the top of the vessel flows through and finally cooled to -37^{0} C in the 2nd stage chiller. Then the H/C fluid enters in 2nd stage liquid separator where 2nd stage liquid is separated. The Second Stage Vapours (SSV) after separation of liquid hydrocarbon (Condensate) are passed through and delivered as feed stock at a temp of 35^oC, to C_2C_3 Recovery Unit. Also SSV can be sent directly to consumers through bypass line if C_2C_3 Recovery Unit is under shut down.

2.3.3 *Cold Box:*

Chill down section and forms COLD BOX which serves to chill the feed gas by exchanging heat with various cold streams of the plant. These exchangers have aluminum brazed plates. The box is filled with insulation material – Pearlite and sealed with Nitrogen.

2.3.4 Light End Fractionator:

Separator liquid from feed gas separator sent to LEF column at around 20°C to remove the lighter fractions. The gas coming out from the top goes to LEF reflux drum and liquid is knocked out by external refrigeration, and the remaining gas is called as LEF top which is sent to ethane propane column after it gets heat from cold box. The bottom liquid goes to LPG column. If ethane propane plant is under shutdown the LEF gases can be sent to consumer line after compressing through residue gas compressor. Also these gases can be used for regeneration of gas dryers.

2.3.5 LPG Column:

The bottom liquid of the LEF column can enter the LPG column. The top product is Liquefied petroleum gas and bottom product is Natural gas liquids.

2.3.6 Propane Column:

The propane is used as a refrigerant in the refrigeration system. The propane losses which occur during refrigeration of feed gas requires make-up. Therefore a propane column is designed in LPG plant to recover propane taking the LPG as a feed to the column.

2.4 Uses and Properties of LP Gases

LP gases are mainly used for cooking purposes and as a fuel. The motor-vehicle sector continues to be the fastest-growing market for LPG. Motorists account for about 42% of World's domestic LPG consumption and automotive use has been growing at a rate of about 17 per cent per year. This growth has been largely due to the tax-free status of LPG which results in a significant pump price differential compared to petrol. LPG vehicle conversion kits are also exempt from salestax. If the production of LPG is abundant in the country, the gases can be used as a petrochemical feed stoke.

The properties of LP Gases are summarized in the following tables (Table 1 & Table 2). The figure for K – Value of several hydrocarbons at various Temperatures and pressures are also presented (De Priester chart – figure 4). The K – Values taken from these figure are used in the design calculations of distillation columns.

2.5 Initial Data for LPG Column [1]

Top temperature $= 60^{\circ}\text{C} = 333 \text{ K}$ Feed temperature $= 108^{\circ}\text{C} = 381 \text{ K}$ Bottom temperature $= 157^{\circ}\text{C} = 430 \text{ K}$

Table 3: Material Balance around LPG Column

| C | - i i id)K os | Vapor (LPG)Kgs | Liquid(Natural gasoline)Kgs |
|----------------------------------|----------------------|----------------|-----------------------------|
| Components | Feed (LEF Liquid)Kgs | | - , 3 |
| C_2H_6 | 163.571 | 163.571 | - |
| C_3H_8 | 12112.806 | 12112.806 | - |
| _ ` | | 5025.787 | - |
| i C ₄ H ₁₀ | 5025.787 | 7342.501 | 1.75/ |
| nC_4H_{10} | 7343.117 | 7342.301 | 1.756 |
| i C ₅ H ₁₂ | 2348.522 | 215.98 | 2132.861 |
| n C ₅ H ₁₂ | 2680.197 | 74.864 | 2604.883 |
| 3-112 | | 24934.5 | 4739.5 |
| | 29674 | | |

Table 4: Physical property of LP gas |5|

(All values at 60 °F and 14.696 psia unless otherwise stated)

| PROPERTIES | PROPANE | ISO – BUTANE | BUTANE |
|--|--------------------------------|--------------------------------|--------------------------------|
| | 44.09 | 58.12 | 58.12 |
| Molecular weight | -43.7 | +10.9 | +31.1 |
| Boiling point °F | -42.1 | -11.7 | -0.5 |
| Boiling point °C | | -255.0 | -216.9 |
| Freezing point °F | -305.8 | -233.0 | -210.7 |
| Density of liquid Specific gravity,(air=1) Degrees, API | 0.508 147.2 4.23 | 0.563 119.8 4.69 | 0.584 110.6 4.87 |
| LB per gallon Density of vapour (ideal gas) Specific gravity.(air=1) | 1.522 116.2 | 2.006 153.1 | 2.006 153.1 |
| Lb gas per 1000cu ft Total heat value (after vaporization) Btu per cu ft Btu per lb | 2,563 21,663 91,740 | 3,269 21,258 99,790 | 3.39 21.008 103,830 |
| Btu per gal of liquid Critical constants Pressure, psia | 617.4 206.2 | 537.0 272.7 | 550.1 306.0 |
| Temperature, °F Specific heat, Btu/lb °F Cp, vapour Cv, vapour Cp/Cv | 0.388 0.343 1.13 0.58 | 0.387 0.348 1.11 0.56 | 0.397 0.361 1.10 0.55 |
| Cp, liquid 60 °F Latent heat of vaporization at boiling point, Blu por the | 183.3 | 157.5 | 165.6 |
| Btu per lb Vapour pressure. psia 0°F 70°F | 37.8 124.3 188.7 | 11.5 45.0 71.8 | 7.3 31.3 51.6 |
| 100°F 100°F(ASTM10), psig max 130°F Properties are for commercial products an | 210 274.5 | 109.5 | 70 80.8 |

Table 4: Combustion Data for LP – gases [5]

(All values at 60 °F and 14.696 psia unless otherwise stated)

| PROPERTIES | PROPANE | ISO - BUTANE | BUTANE |
|--|-----------|-----------------|-----------|
| Flash temperature, °F | -156 | -117 | -101 |
| Ignition temperature, °F | 932 | 950 | 896 |
| Maximum Flame temperature in air, °F | | | |
| Observed | 3497 | 3452 | 3443 |
| · | 3583 | 3583 | 3583 |
| Calculated | | | |
| Flammability limits, % gas in air | 2.37 | 1.30 | 1.36 |
| Lower | 9.50 | 8.44 | 8.41 |
| Higher | 7.50 | | 01.12 |
| Maximum rate flame propagation in 1 in, tube | 32 | 33 | 33 |
| Inches per second | 4.6 – 4.8 | 3.6 - 3.8 | 3.6 – 3.8 |
| Percentage gas in air | | | |
| Required for complete combustion (ideal gas) | 23.9 | 31.1 | 31.1 |
| Air, cu ft per cu ft gas | 15.7 | 15.5 | 15.5 |
| lb per lb gas | 5.0 | 6.5 | 6.5 |
| Oxygen, cu ft per cu ft gas | 3.63 | 3.58 | 3.58 |
| lb per lb gas | | | |
| Product of combustion (ideal gas) | 3.0 | 4.0 | 4.0 |
| Carbon dioxide, cu ft per cu ft gas | 2.99 | 3.03 | 3.03 |
| lb per lb gas | 4.0 | 5.0 | 5.0 |
| Water vapour, cu ft per cu ft gas | 18.9 | 24.6 | 24.6 |
| Nitrogen, cu ft per cu ft gas | | | |

Properties are for commercial products and vary with composition

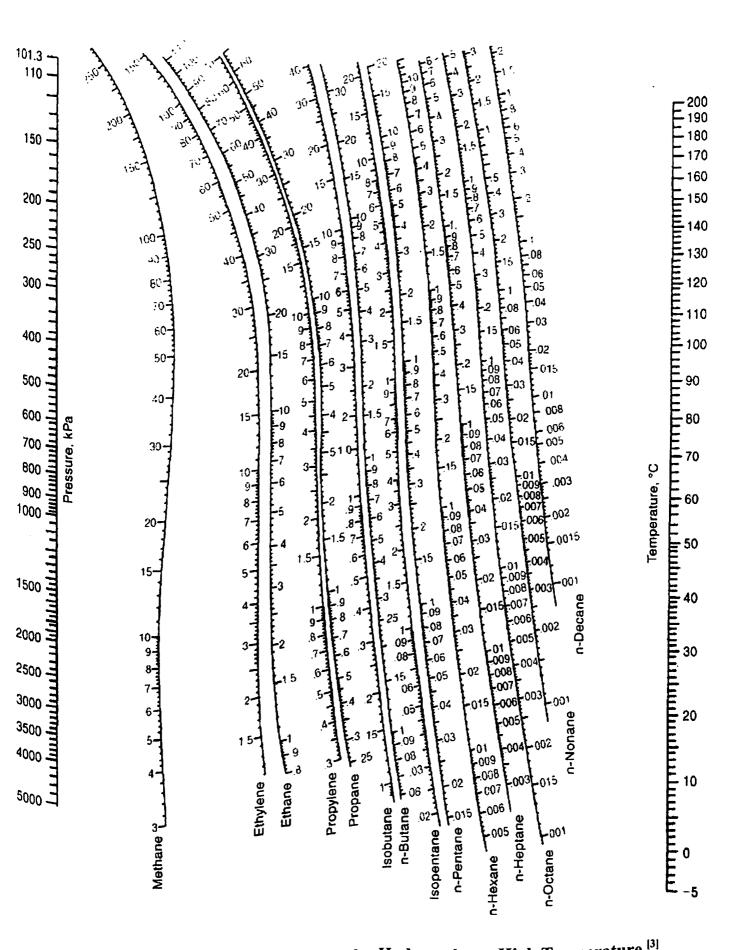


Figure 4: De Priester Chart – k Values for Hydrocarbons, High Temperature [3]

PROCESS DESIGN OF DISTILLATION COLUMN – EXCEL – 2007 3 **PROGRAME**

To estimate the stage, and the condenser and reboiler temperatures, procedures are required for calculating dew and bubble points. then design the sieve tray column. Which involves following three steps.

Step 1: Dew Points And Bubble Points

Step 2: Mccabe-Thiele Design Method

Step 3: Sieve Tray Design

3.1 Dew Points and Bubble Points [3]

To estimate the stage, and the condenser and reboiler temperatures, procedures are required for calculating dew and bubble points. By definition, a saturated liquid is at its bubble point (any rise in temperature will cause a bubble of vapour to form), and a saturated vapour is at its dew point (any drop in temperature will cause a drop of liquid to form). Dew points and bubble points can be calculated from a knowledge of the vapour-liquid equilibrium for the system. In terms of equilibrium constants, the Bubble point is defined by the equation (1) Dew point is calculating by equation (2).

Bubble point:
$$\sum y_i = \sum K_i x_i = 1.0$$
 (1)

Dew point:
$$\sum x_i = \sum (y_i/\kappa_i) = 1.0$$
 (2) [3]

For multicomponent mixtures the temperature that satisfies these equations (3) & (4), at a given system pressure, must be found by trial and error. For binary systems the equations can be Solved more readily because the component compositions are not independent; fixing one fixes the other.

$$y_a = 1 - y_b (3)^{[3]} (4)$$

$$y_a = 1 - x_b \tag{4}$$

$$x_a = 1 - x_b$$

[3]

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| lipboard | | Fant | | Ligaritant | Number | | 21/102 | | |
| C16 | - 2 | | 12.57022E | | | | | | |
| | | fx 0.010947884 | | ======================================= | F | G | Н | | |
| ^ | В | C | D | - | - OTAT | TOA | ICI | TATI | |
|)III | DIE | DOI | NT T | JEW | P() N | | LUU | LAIN | |
| JUL | DDLF | rui | NT - I | | TOIL | hr X _{id} | "Bottom" in K | g "B" in Kg moles | hr X _{i,b} |
| mponent | "Feed" in Kg | | "F" in moles/hr | "Distillate" in K | g "D ₁ "in Kg moles/ | hr -Yid | Dollow in i | • | |
| | reed III Võ | Mol nt | 1 m mores m | 575 Sec. 20 | | 0.010947884 | | 8.50 | |
| athane | 162 571 | | 5.452366667 | 163.571 | 5.452366667 | 0.55276079 | | - | - |
| ropane | 163.571 | 30 | 275.2910455 | 12112.806 | 275.2910455 | 0.173988774 | real er | | - |
| Butane | 12112.806 | 44 | | 5025.787 | 86.6515 | 0.254191582 | 1.756 | 0.030275862 | 0.000459894 |
| Butane | 5025.787 | 58 | 86.6515 | 7342.501 | 126.5948448 | 0.006023185 | 2132.861 | 29.62306944 | 0.449977903 |
| Pentane | 7343.117 | 58 | 126.6054655 | 215.98 | 2.999722222 | 0.002087785 | 2604.883 | 36.17893056 | 0.549562203 |
| Pentane | 2348.522 | 72 | 32.61836111 | 74.864 | 1.039777778 | 1 | | 65.83227586 | 1 |
| entallé | 2680.197 | 72 | 37.22495833 | 24935.509 | 498.0292569 | 1 | | | |
| | | | 563.8436971 | | | | | | |
| | | | | | . Samuel A SUB | 100°C | Trail 02 | 130°C | Trail 03 13 |
| | | | | | | | | | |
| ימקדו | | | | | Trail 01 | 109 | | K. Y.b | K K |
| UBBLE- P | OINT CALE | | | | Trail 01 K _i | K _i X _i ,b | K _i | K _i X _i b | K _i K |
| UBBLE- P | OINT CALE Di"in Kg moles/ | hr X., | 'B" in Kg moles/hr | $X_{i,b}$ | Printed to the second s | 109 | | K ₄ X ₂ b | K, K |
| mponent"] | ONT CALE Di''in Kg moles/ | hr X _{i,d} ' | 'B" in Kg moles/hr | $X_{i,b}$ | Printed to the second s | 109 | | K _i X _i b | K, K |
| mponent"] | D _i "in Kg moles/l | | 'B" in Kg moles/hr | $X_{i,b}$ | Printed to the second s | 109 | | K ₁ X ₂ b |), (*) |
| mponent"] athane | D _i "in Kg moles/1 5.452366667 | 0.010947884 | 'B" in Kg moles/hr | X _{i,b} | Printed to the second s | K _i X _i ,b | K _l | | 16. 16. |
| mponent"] athane ropane Butane | 5.452366667 275.2910455 | 0.010947884 0.55276079 | 'B" in Kg moles/hr - - | | Printed to the second s | K _i X _i ,b | | K ₁ X ₂ b 0.000827809 0.449977903 | - K, K, |
| mponent"] athane ropane Butane | 5.452366667 275.2910455 86.6515 | 0.010947884 0.55276079 0.173988774 | | o.000459894 | - 1.22 0.65 | K _i X _i ,b - 0.000561071 0.292485637 | 1.8 | 0.000827809 | 11 43 44 44 |
| athane ropane Butane Butane Pentane | 5.452366667 275.2910455 86.6515 126.5948448 | 0.010947884 0.55276079 0.173988774 0.254191582 | 0.030275862 29.62306944 | 0.000459894 0.449977903 | 1.22 0.65 0.55 | K _i X _i ,b - 0.000561071 0.292485637 0.302259212 | 1.8 1.09 | 0.000827809 0.449977903 0.494605983 | 11 43 44 44 |
| athane ropane Butane Butane Pentane | 5.452366667 275.2910455 86.6515 126.5948448 2.999722222 | 0.010947884 0.55276079 0.173988774 0.254191582 0.006023185 | 0.030275862 | 0.000459894 0.449977903 0.549562203 | 1.22 0.65 0.55 | K _i X _i ,b - 0.000561071 0.292485637 | 1.8 | 0.000827809 0.449977903 0.494605983 | 11 43 44 44 |
| athane ropane Butane Butane Pentane | 5.452366667 275.2910455 86.6515 126.5948448 | 0.010947884 0.55276079 0.173988774 0.254191582 | 0.030275862 29.62306944 | 0.000459894 0.449977903 | 1.22 0.65 0.55 | K _i X _i ,b - 0.000561071 0.292485637 0.302259212 | 1.8 1.09 | 0.000827809 0.449977903 0.494605983 | 11 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1 |
| athane ropane Butane Butane Pentane Pentane | 5.452366667 275.2910455 86.6515 126.5948448 2.999722222 1.039777778 498.0292569 | 0.010947884 0.55276079 0.173988774 0.254191582 0.006023185 0.002087785 | 0.030275862 29.62306944 36.17893056 | 0.000459894 0.449977903 0.549562203 | K _i - 1.22 0.65 0.55 ΣK _i X _i | K _i X _i ,b 0.000561071 0.292485637 0.302259212 b 0.595305919 | 1.8 1.0.9 ΣK ₄ X | 0.000827809 0.449977903 0.494605983 b 0.945411695 | 14 43 49 4 7 Kg X ₂ th 9 7 |
| mponent "I athane ropane Butane Butane Pentane Pentane | 5.452366667 275.2910455 86.6515 126.5948448 2.999722222 1.039777778 498.0292569 | 0.010947884 0.55276079 0.173988774 0.254191582 0.006023185 0.002087785 | 0.030275862 29.62306944 36.17893056 | 0.000459894 0.449977903 0.549562203 | K _i - 1.22 0.65 0.55 ΣK _i X _i | K _i X _i ,b 0.000561071 0.292485637 0.302259212 b 0.595305919 | 1.3 1.3 1 0.9 <u>YK, X</u> | 0.000827809 0.449977903 0.494605983 .b 0.945411695 | 100-100 (00 100-100 100 |
| mponent "I athane ropane Butane Butane Pentane Pentane | 5.452366667 275.2910455 86.6515 126.5948448 2.99972222 1.039777778 498.0292569 | 0.010947884 0.55276079 0.173988774 0.254191582 0.006023185 0.002087785 | 0.030275862 29.62306944 36.17893056 | 0.000459894 0.449977903 0.549562203 | K _i | K _i X _i ,b 0.000561071 0.292485637 0.302259212 b 0.595305919 | 1.8 1.0.9 ΣK ₄ X | 0.000827809 0.449977903 0.494605983 b 0.945411695 | 100-100 (00 100-100 100 |
| mponent "I athane 'ropane Butane Butane Pentane Pentane | 5.452366667 275.2910455 86.6515 126.5948448 2.99972222 1.039777778 498.0292569 | 0.010947884 0.55276079 0.173988774 0.254191582 0.006023185 0.002087785 | 0.030275862 29.62306944 36.17893056 65.83227586 | 0.000459894 0.449977903 0.549562203 | K _i - 1.22 0.65 0.55 ΣK _i X _i | K _i X _i ,b 0.000561071 0.292485637 0.302259212 b 0.595305919 50°C K _i X _i ,d | 1.8 1.0.9 \(\sum_{K_1} \text{X}\) | 0.000827809 0.449977903 0.494605983 .b 0.945411695 70°C K ₄ X ₂ d | 14 93 44 5 5 6 6 6 6 6 6 6 6 6 6 6 6 6 6 6 6 |
| mponent " athane ropane Butane Butane Pentane Pentane | 5.452366667 275.2910455 86.6515 126.5948448 2.999722222 1.039777778 498.0292569 | 0.010947884 0.55276079 0.173988774 0.254191582 0.006023185 0.002087785 | 0.030275862 29.62306944 36.17893056 | 0.000459894 0.449977903 0.549562203 | K _i | K _i X _i ,b 0.000561071 0.292485637 0.302259212 b 0.595305919 50°C K _i X _i ,d | 13 1 0.9 \(\sum K_4 \) \(\text{X} \) | 0.000827809 0.449977903 0.494605983 b 0.945411695 70°C K _i X _i ,d | Viail 03 44 |
| mponent" athane ropane Butane Butane Pentane Pentane DEW - F | 5.452366667 275.2910455 86.6515 126.5948448 2.999722222 1.039777778 498.0292569 POINT CALE | 0.010947884 0.55276079 0.173988774 0.254191582 0.006023185 0.002087785 1 | 0.030275862 29.62306944 36.17893056 65.83227586 | 0.000459894 0.449977903 0.549562203 | K _i 1.22 0.65 0.55 ΣK _i X _i | K _i X _i ,b 0.000561071 0.292485637 0.302259212 b 0.595305919 50°C K _i X _i ,d 0.045981114 0.829141185 | . K ₁ . 1.8 . 1 . 0.9 . ∑K ₄ X ₂ . Trail 02 . K ₄ . 5.1 . 2.1 | 0.000827809 0.449977903 0.494605983 b 0.945411695 70°C K, X, d | Viail 03 44 |
| mponent "] athane ropane Butane Butane Pentane DEW - F | 5.452366667 275.2910455 86.6515 126.5948448 2.999722222 1.039777778 498.0292569 POINT CALE D _i "in Kg moles/ | 0.010947884 0.55276079 0.173988774 0.254191582 0.006023185 0.002087785 1 | 0.030275862 29.62306944 36.17893056 65.83227586 | 0.000459894 0.449977903 0.549562203 | K _i 1.22 0.65 0.55 ΣK _i X _i Trail 01 K _i | K _i X _i b 0.000561071 0.292485637 0.302259212 b 0.595305919 50°C K _i X _i d 0.045981114 0.829141185 0.121792142 | 1.8 1.0.9 ∑K, X, Trail 02 K, | 0.000827809 0.449977903 0.494605983 .b 0.945411695 70°C K ₁ X ₂ ,d 0.05583421 1.160797659 0.182688213 | Trail 03 44 - X2 - X3 |
| mponent " athane ropane Butane Butane Pentane Pentane DEW - F | 5.452366667 275.2910455 86.6515 126.5948448 2.99972222 1.039777778 498.0292569 POINT CALE D _i "in Kg moles/ 5.452366667 275.2910455 | 0.010947884 0.55276079 0.173988774 0.254191582 0.006023185 0.002087785 1 hr X _{i,d} 0.010947884 0.55276079 | 0.030275862 29.62306944 36.17893056 65.83227586 | 0.000459894 0.449977903 0.549562203 1 | K _i | K _i X _i ,b 0.000561071 0.292485637 0.302259212 b 0.595305919 50°C K _i X _i ,d 0.045981114 0.829141185 0.121792142 0.132179623 | 13 1 0.9 ∑K, X Trall 02 K, | 0.000827809 0.449977903 0.494605983 .b 0.945411695 70°C K, X, d. 0.05583421 1.160797659 0.182688213 0.193185603 | Trail 03 44 - X2 - X3 |
| mponent " athane ropane Butane Butane Pentane DEW - F athane Propane Butane | 5.452366667 275.2910455 86.6515 126.5948448 2.999722222 1.039777778 498.0292569 POINT CALE 5.452366667 275.2910455 86.6515 | 0.010947884 0.55276079 0.173988774 0.254191582 0.006023185 0.002087785 1 hr X _{i,d} 0.010947884 0.55276079 0.173988774 | 0.030275862 29.62306944 36.17893056 65.83227586 | 0.000459894 0.449977903 0.549562203 1 X _{i,b} | K _i 1.22 0.65 0.55 ΣK _i X _i Trail 01 K _i 4.2 1.5 0.7 | K _i X _i ,b 0.000561071 0.292485637 0.302259212 b 0.595305919 50°C K _i X _i ,d 0.045981114 0.829141185 0.121792142 0.132179623 0.01566028 | 1.8 1.0.9 \(\sum K_i \) X. Trall 02 \(K_i \) 5.1 2.1 1.05 0.76 0.37 | 0.000827809 0.449977903 0.494605983 b 0.945411695 70°C K, X,d 0.05583421 1.160797659 0.182688213 0.193185603 0.002128578 | Trail 03 44 - X2 - X3 - 02 - 131 - 02 - 148 - 151 - 519 - 60 |
| mponent " athane ropane Butane Butane Pentane DEW - F Dimponent " Dathane Propane Butane Butane Propane Butane | 5.452366667 275.2910455 86.6515 126.5948448 2.999722222 1.039777778 498.0292569 POINT CALE D _i "in Kg moles/ 5.452366667 275.2910455 86.6515 126.5948448 | 0.010947884 0.55276079 0.173988774 0.254191582 0.006023185 0.002087785 1 hr X _{i,d} 0.010947884 0.55276079 0.173988774 0.254191582 | 0.030275862 29.62306944 36.17893056 65.83227586 | 0.000459894 0.449977903 0.549562203 1 X _{i,b} | K _i 1.22 0.65 0.55 ΣK _i X _i Trail 01 K _i 4.2 1.5 0.7 0.52 | K _i X _i ,b 0.000561071 0.292485637 0.302259212 b 0.595305919 50°C K _i X _i ,d 0.045981114 0.829141185 0.121792142 0.132179623 | 1.8 1 0.9 ∑K, X, Trail 02 K, 5.1 2.1 1.05 0.76 0.37 | 0.000827809 0.449977903 0.494605983 .b 0.945411695 70°C K, X, d 0.05583421 1.160797659 0.182688213 0.193185603 0.002128578 0.900626335 | Trail 03 44 - X2 - X3 - 02 - 131 - 02 - 148 - 151 - 519 - 60 |
| mponent " athane ropane Butane Butane Pentane DEW - F Dimponent " Dathane Propane Butane Butane Propane Butane | 5.452366667 275.2910455 86.6515 126.5948448 2.999722222 1.039777778 498.0292569 POINT CALE 5.452366667 275.2910455 86.6515 126.5948448 2.999722222 | 0.010947884 0.55276079 0.173988774 0.254191582 0.006023185 0.002087785 1 hr X _{i,d} 0.010947884 0.55276079 0.173988774 0.254191582 0.006023185 | 0.030275862 29.62306944 36.17893056 65.83227586 "B" in Kg moles/hr | 0.000459894 0.449977903 0.549562203 1 X _{i,b} | $\begin{array}{c} \textbf{K}_{i} \\ \hline - \\ - \\ 1.22 \\ 0.65 \\ 0.55 \\ \hline \boldsymbol{\Sigma} \textbf{K}_{i} \textbf{X}_{i} \\ \hline \textbf{Trail 01} \\ \textbf{K}_{i} \\ \hline \\ 4.2 \\ 1.5 \\ 0.7 \\ 0.52 \\ 2.6 \\ 0.19 \\ \\ \end{array}$ | K _i X _i ,b 0.000561071 0.292485637 0.302259212 b 0.595305919 50°C K _i X _i ,d 0.045981114 0.829141185 0.121792142 0.132179623 0.01566028 | 1.8 1.0.9 \(\sum K_i \) X. Trall 02 \(K_i \) 5.1 2.1 1.05 0.76 0.37 | 0.000827809 0.449977903 0.494605983 .b 0.945411695 70°C K, X, d 0.05583421 1.160797659 0.182688213 0.193185603 0.002128578 0.900626335 | Trail 03 44 - X2 - X3 - 02 - 131 - 02 - 148 - 151 - 519 - 60 |
| mponent " athane ropane Butane Butane Pentane Pentane DEW - F | 5.452366667 275.2910455 86.6515 126.5948448 2.999722222 1.039777778 498.0292569 POINT CALE D _i "in Kg moles/ 5.452366667 275.2910455 86.6515 126.5948448 | 0.010947884 0.55276079 0.173988774 0.254191582 0.006023185 0.002087785 1 hr X _{i,d} 0.010947884 0.55276079 0.173988774 0.254191582 | 0.030275862 29.62306944 36.17893056 65.83227586 | 0.000459894 0.449977903 0.549562203 1 X _{i,b} | $\begin{array}{c} \textbf{K}_{i} \\ \hline - \\ - \\ 1.22 \\ 0.65 \\ 0.55 \\ \hline \boldsymbol{\Sigma} \textbf{K}_{i} \textbf{X}_{i} \\ \hline \textbf{Trail 01} \\ \textbf{K}_{i} \\ \hline \\ 4.2 \\ 1.5 \\ 0.7 \\ 0.52 \\ 2.6 \\ 0.19 \\ \\ \end{array}$ | K _i X _i b 0.000561071 0.292485637 0.302259212 b 0.595305919 50°C K _i X _i d 0.045981114 0.829141185 0.121792142 0.132179623 0.01566028 0.000396679 | 1.8 1 0.9 ∑K, X, Trail 02 K, 5.1 2.1 1.05 0.76 0.37 | 0.000827809 0.449977903 0.494605983 .b 0.945411695 70°C K, X, d 0.05583421 1.160797659 0.182688213 0.193185603 0.002128578 0.900626335 | 14 43 49 4 7 Kg X ₂ th 9 7 |

3.2 Mccabe-Thiele Design Method 131

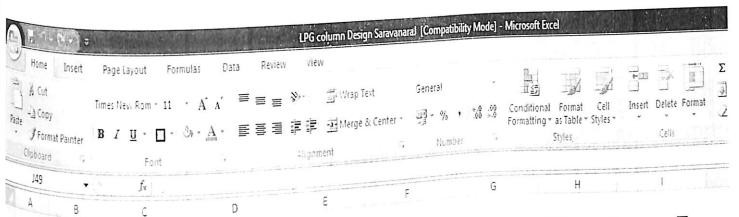
The vapour-liquid equilibrium characteristics (indicated by the shape of the equilibrium curve) of the mixture will determine the number of stages, and hence the number of trays, required for the separation. This is illustrated clearly by applying the McCabe-Thiele method to design a binary column.

The McCabe-Thiele approach is a graphical one, and uses the VLE plot to determine the theoretical number of stages required to effect the separation of a binary mixture. It assumes constant molar overflow and this implies that:

- Molal heats of vaporisation of the components are roughly the same
- Heat effects (heats of solution, heat losses to and from column, etc.) are negligible
- For every mole of vapour condensed, 1 mole of liquid is vaporized.

The design procedure is simple. Given the VLE diagram of the mixture, operating lines are drawn first.

- Operating lines define the mass balance relationships between the liquid and vapor phases in the
- •There is one operating line for the bottom (stripping) section of the column, and on for the top (rectification or enriching) section of the column.
- •Use of the constant molar overflow assumption also ensures the operating lines are straight lines



McCABE - THIELE Method

R= 1

| | . | | "F" in moles/hr | "Distillate" in Kg | "D _i "in Kg moles/hr | "Bottom" in Kg | "B" in Kg moles |
|--|--|----------------------|--|---|---------------------------------------|-------------------------------|---|
| Component | "Feed" in Kg | Mol wt | "F" in moles.m | 163.571 | 5.452366667 275.2910455 | 77 - 2 mag | |
| Eathane Propane i-Butane n-Butane | 163.571 12112.806 5025.787 7343.117 | 30 44 58 58 | 275.2910455 86.6515 126.6054655 32.61836111 | 12112.806 5025.787 7342.501 215.98 | 86.6515 126.5948448 2.999722222 | 1.756 2132.861 2604.883 | 0.030275862 29.62306944 36.17893056 |
| i-Pentane n-Pentane | 2348.522 | 72 72 | 32.61830111 37.22495833 563.8436971 | 74.864 24935.509 | 1.039777778 498.0292569 | 4739.5 | 65.83227586 |

| Component | <u>.</u> | К воп | K _{avg} | $\sigma_{\rm i}$ |
|-----------|------------------|-------|----------------------|----------------------------|
| Poncil | k _{Top} | | 6.1 | 9.457364341 |
| Eathane | 3.7 | 8.5 | 2.77 | 4.294573643 2.403100775 |
| Propane | 1.34 | 4.2 | 1.55 | 1.88372093 |
| i-Butane | 0.6 | 2.5 | 1.215 | 1 |
| n-Butane | 0.43 | 2 | 0.645 | 0.821705426 |
| i-Pentane | 0.19 | 1.1 | 0.53 | |
| n-Pentane | 0.16 | 0.9 | | |

Cale of non - Key flows

| Cale of non - k | | | $l_i=d_i/(a_i-1)$ | vi=li+di |
|--------------------------------|---|---------------------------------------|---|----------------------------|
| Component | $a_{\mathbf{i}}$ | "D _i "in Kg moles/hr | 0.644688634 | 6.097055301 358.8499745 |
| Eathane Propane i-Butane | 9.457364341 4.294573643 2.403100775 | 5.452366667 275.2910455 86.6515 | 83.55892909 61.75714641 145.9607641 | 149 4096464 |
| | | $\sum l_i$ | • | 11-V'.+b. |

| | | | oles/hr V'i=aipi/(alk- | a_i) $l'_i = V'_i + b_i$ |
|-----------|------------|--------------|--------------------------|-----------------------------|
| Component | a_{i} | "B" in Kg mo | oles/iii · i | 722 64.17138777 |
| - | | 172030 | 27.99245 | 722 64.17138777 |
| n-Pentane | 0.82170542 | 6 36.178930 | $\sum V'_{i} = 27.99245$ | |

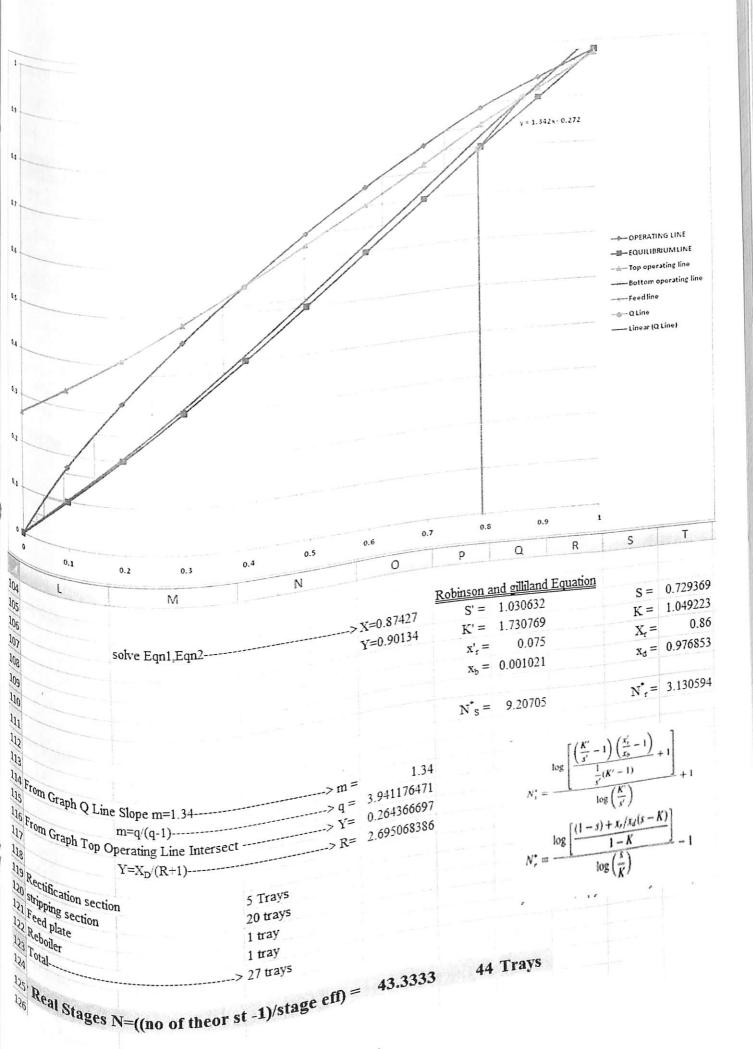
To calculate Maximum Vapour rate (m³/s) from y_i*Mwt

| yi | yi*Mol wt |
|-------------|-------------|
| 0.010947884 | 0.328436528 |
| 0.55276079 | 24.32147475 |
| 0.173988774 | 10.09134891 |
| 0.254191582 | 14.74311177 |
| 0.006023185 | 0.433669302 |
| 0.002087785 | 0.150320486 |

Mwa = 50.06836175

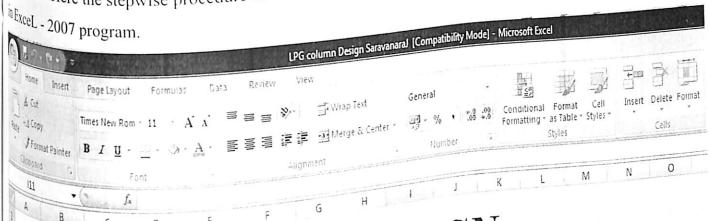
| A B | | | | | G | Н |
|--|----------------------------|----------------------------|---|---|---|-----|
| АВ | C | D _a | Ε | f | | - " |
| Flows of combin | ied kev | | | | | |
| L _a = | 352 0684928 | | | | | |
| V _t = | 482,7028376 | | | | | |
| Γ' _t = | 968.0660567 | | | | | |
| L' _e = | 997.719402 | | | | | |
| | | | | | | |
| slope of top op | erating line | | | | | |
| Slope of top op Slope of bottom Slope of b | $L_{\bullet}V_{\bullet}$ = | 0.729369014 | | | | |
| | | | | | | |
| slope of bottom | | . 222621531 | | | | |
| | Γ, * Λ, *= | 1.030631531 | | | | |
| | 333335 | 0.001020993 | | | | |
| | $X_0 =$ | 0.976853025 | | | | |
| | x ₃ = | 0.795141457 | | | | |
| | X,= | 1.88372093 | | | | |
| <i>X</i> ≈ 0 | α= | 0 | | | | |
| 0.1 | λ= | 0.173076923 | | | | |
| 0.2 | | 0.320158103 | | | | |
| 0.3 | | 0.446691176 | | | | |
| 0.4 | | 0.556701031 0.653225806 | | | | |
| 0.5 | | 0.033223000 | | | | |
| 0.6 | | 0.814655172 | | | | |
| 0.7 | | 0.882833787 | | | | |
| 0.8 | | 0.944300518 | | | | |
| 0.9 | | 1 | | | | |
| 1 | | | | | | |
| line | | | | | | |
| X | | 0 | | | | |
| U | Υ= | 0.1 | | | | |
| 0.1 | | 0.2 | | | | |
| 0.2 | | 0.3 | | | | |
| 0.3 | | 0.4 | | | | |
| 0.4 | | 0.5 | | | | |
| 0.5 | | 0.6 | | | | |
| 0.6 | | 0.7 | | | | |
| 0.7 | | 0.8 | | | | |
| 0.8 | | 0.9 | | | | |
| 0.9 | | 1 | | | | |
| 1 | | | | | | |

| A B | | Č | D | E | F | | G |
|----------------------------------|------------|-----------|-----------------------|---|-----------|----------|------------|
| ⁷⁹ Top operating line | - Fan 1 | >Υ=0.729. | X264 | | | | |
| 80 C= 0.26436 | | | | | | | |
| 81 X= | | <u> </u> | 0.264366697 | | | | |
| 1 | 0 | 1- | 0.337303599 | | | | |
| 3 | 0.1 | | 0.4102405 | | | | |
| 4 | 0.2 | | 0.483177401 | | | | |
| 5 | 0.3 | | 0.556114303 | | | | |
| 6 | 0.4 | | 0.530111333 | | | | |
| | 0.5 | | 0.029031204 | | | | |
| | 0.6 | | 0.774925007 | | | | |
| | 0.7 | | 0.774923007 | | | | |
| | 0.8 | | 0.847801500 | | | | |
| | 0.9 | | | | | | |
| | 1 | | 0.993735711 | | | | |
| 2 | - | | | | Feed line | | |
| Bottom operating lin | ne Eqn | 2>Y=1.031 | X-3.127E-05 | | X | Y | |
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| 8 | 0.2 | | 0.309158185 | | 0.795 | 141457 | 0.4 |
| 9 | 0.3 | | 0.412221338 | | | 141457 | 0.5 |
| 00 | 0.4 | | 0.515284491 | | | | 0.0 |
| | 0.5 | | 0.618347644 | | | 141457 | 0.79514145 |
| DZ | 0.6 | | 0.721410797 | | | 141457 | |
| 03 | 0.7 | | 0.82447395 | | Q line | 1 21 157 | 0.79514145 |
| 04 | 0.8 | | 0.927537103 | | | 141457 | 0.9013 |
| 05 | 0.9 | | 1.030600256 | na principal na minori (il dispositar na minori di madelina na mi | |).87427 | 8.7013 |
| | 1 | | LUJVVV | | | | |



33 Sieve Tray Design 131

Here the stepwise procedure as well as each formulas which involves in that steps, are explain



SIEVE - TRAY DESIGN

STEP 01: Flow Rate Of Enriching Section.

R=L/D

RR= 2.6950684

L=RD

L= 1342.2229 Kmol/hr

G= 1840.2522 Kmol hr

B= 65.832276 Kmol/hr

5 STEP 02: Density.

Liquid Density--->= 634.4994 Kg m³

Vapor Density--->= 4.306843 Kg m³ Molecular Weight->=

STEP 03: Vapor Load:

 $V=D(R+1)*mol.wt/(3600*\rho_V)$

 $V = 303.07447 \text{ m}^3/\text{sec}$

35 STEP 04: Tray Spacing 30 STEP 05: Allowable Vapor Velosity:

V_c=(L/V)*(p

18 inch

Let us assume the tray spacing: 0.457 m

 $V_c = (L/V)^*(\rho_v/\rho_L)$

35 STEP 06: Ctoss section area of column:

Cross section area of column:

" — vapour flow rate

" — vaple vapour flow rate allowable vapor Velocity

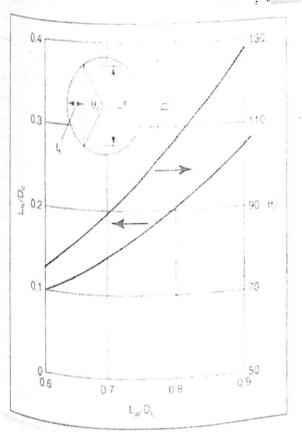
Cross section area of column = 6.12173068 m²

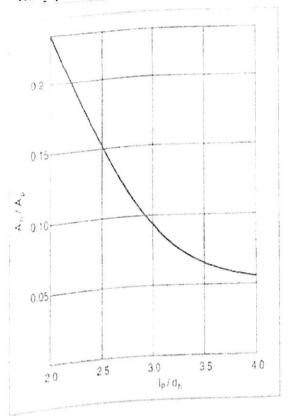
BIROT: Area Total Cohmn Area = Cross Sectional area 1 88 Total Column Area A = 6.956512 m Net Area $A_c = 6.121731 \text{ m}^2$ 12 % Az in A: Down comer area $A_d = 0.834781 \text{ m}^2$ Assume. Active area $A_2 = -5.286949$ Hole area $A_{b} = -0.528695$ Assume. TEP 08: column Diameter. $D_c = ((4*A)\pi)^{-5}$ D_C= 2.9768774 m Column Thinkess is: 0.06 m OD= 3.0368774 m The hole size plate varies from 3mm to 12mm hance we assume as 8 mm 0.008 mm STEP 09: Timkness of sieve plate: Thinkness = 0.1 to 1.2 hole size Let us take 0.6 times Thinkness = 4.8 mm STEP 10: Wire Length (Lw.): 0.0048 m From Graph $L_W/D_C = 0.765$ By using $(Ad Ac)^*100 = 12$ Wire Length $L_W = 2.277311 \text{ m}$ STEP 11: Hight of froath in mm (h.): 0.9 0.8 0.7 I_w/D_c 0.6 $H_{ow} = 43.4 (Q_L/L_W)^{(2.3)}$ $Q_L = flow rate of clear liquid (m³ min)$ Q_L =(Mass flow rate of distillate *R) / (\rho_L * 60) $Q_L = 1.7652473 \text{ m}^3/\text{min}$ $H_{\rm OW} = 36.622264 \text{ mm}$ M STEP 12: Perforated area (A.): 50 mm unperferated strip round pplate edge From graph By using ---> $L_w/D_c = 0.765$ $L_{\rm W}/D_{\rm C} = 0.765$ Assume Angle subtended at plate edge by unperforated strip = 104 ° Mean length, unperforated edge strips = 3.880389 m 50 mm wide claming zones Area of imperforaded edge strips = 0.194019 m² Assume Mean length of clamping zone = 2.887738 m Area of calaming zone = 0.288774 m² Total area for perforations, $A_p = 4.804156 \text{ m}^2$

41

 $A_{\rm b} A_{\rm p} = 0.110049$ 2.82

From graph By using $A_h A_p = 0.110049$





19 STEP 13: Number of holes:

17/ 28

9 in

15

100

197 188

百年

117

118

119

100

DI

120

13 124

135

136

 0.00005024 m^2 Area of one hole =

10523.3862 Number of Holes =

14 STEP 14: Check weeping

Maiximum liquid rate = 1.31652778 Kg s

Maximum lqd rate - 70 per cent turn-down = 0.92156944 Kg/5

mm Weir height 50 Assume From Graph By using h_w - h_{ow} = 86.62226 mm

 h_{W} - $h_{OW} = 86.6222635 \text{ mm}$

 $K_2 = 30.8$ $\tilde{u}_h (min) = (K_2 - 0.90 * (25.4 - d_h))/(\rho_v)^{5.5} m/s$

 $\tilde{\mathbf{u}}_{h}(\min) = 7.29535374 \text{ m/s}$

vapour rate in (moles/sec) = 138.34146 moles/s

Maximum Vapor rate $(m^3/s) = 11.9000289 \text{ m}^3/s$ Actual minimum vapour velosity = $(Min \ vapour \ rate/A_h)$ m/s

Actual minimum vapour velosity 15.7558165 m/s

Actual minimum vapour velosity 13.7074 ####
So minimum operating will be will abow weep point ####

UTSTEP 15: Plate pressure drop 18

129

131

132

134

Maximum vapour velocity through holes \tilde{d}_{E}

ŭ. = 22.50**8**3093 m s

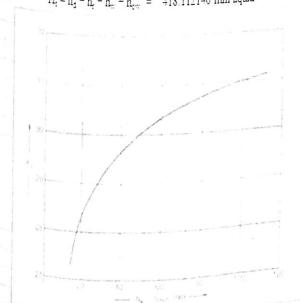
Co = 0.75 From Graph by using (plate thinkness plate dia), (A_k A_p):

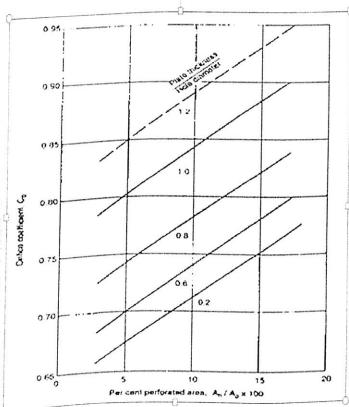
Dry plate drop

 $h_d = 51*((uh \cdot Co)^2)*(\rho_v \cdot \rho_t) = 311.789313 \text{ mm liquid}$ Residual head

 $hr = 12.5*1000 \ \rho_1 = 19.7005702 \ mm \ liquid$

Total plate pressure drop $H_t = h_s - h_t - h_m - h_m = 418.112146 \text{ mm liquid}$





STEP 16: Down comer Equid back-up:

Down comer pressure loss

$$h_{ap} = h_{w} - 10 =$$

40 mm

Area under apron, $A_{2p} = 0.091092 \text{ m}^2$

As this is less than $A_d = 0.834781 \text{ m}^2 \text{ so.} A_d > A_{zp}$

 $h_{tc} = 166(L_{md}/\rho_L A_{pp})^2 = 0.086127 \text{ mm}$

Downcomer backup measured from plate surface, mm

 $H_b = h_a + h_{op} + H_d + H_t = 816.5237 \text{ mm liquid}$

 $H_h = 0.816524 \text{ m}$

Plate spacing = 18 in assume

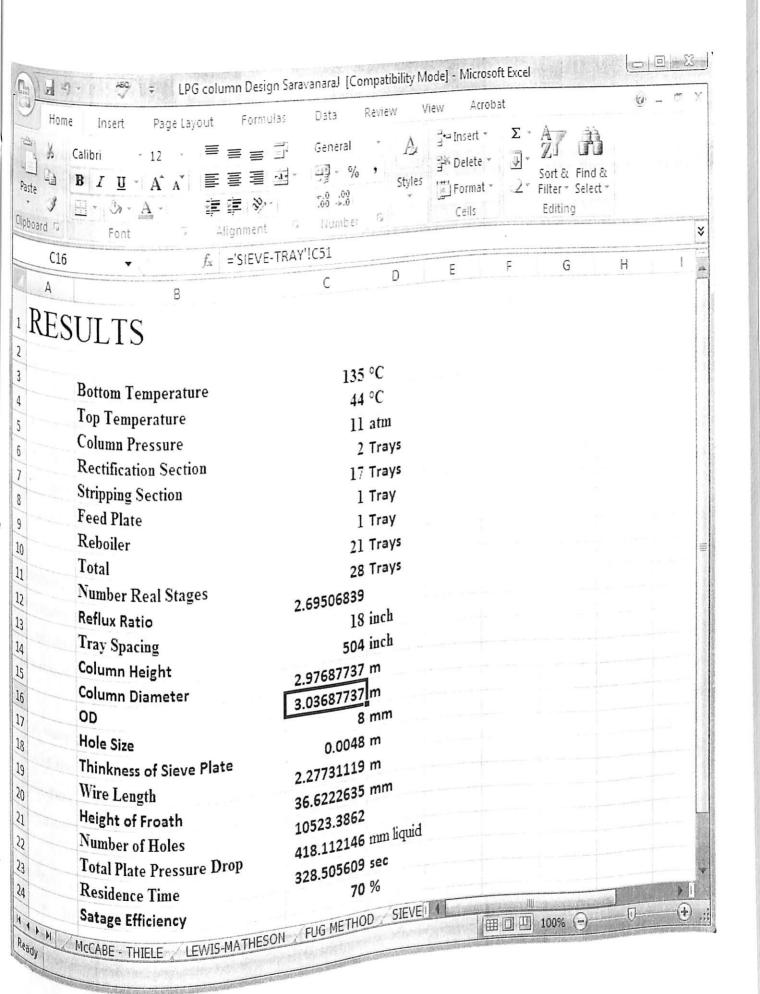
816.523723 <(1/2)*(Plate spacing+Weir length): 1367.2556 mm

So, the diameter and tray spacing – both are satisfactory

STEP 17: Residence time:

 $t_i = (A_d H_{bc} \rho_l) / L_{wd} =$

328.5056



4. PROCESS DESIGN OF DISTILLATION COLUMN - CHEMCAD 6.0.1

4.1 Overview of CHEMCAD 6.0.1 and Its Uses 171

Chemical processing industry (CPI) faces numerous challenges, rising fuel and feedstock costs, reduced engineering staff, shorter product life cycles, increased global competition, and increased regulation. These challenges require that CPI companies seek out and use the best tools to increase productivity and improve engineering decisions.

CHEMCAD is a powerful and flexible chemical process simulation environment, built around three key values of innovation, integration, and open architecture. These values create important advantages for us:

- The latest chemical engineering techniques at our fingertips
- All functionality united in a single software environment
- Seamless connection to the chemical engineering computing environment, with links to tools such as MS Excel and Word and interfaces such as COM, DCOM, OPC, CAPE - OPEN, and

CHEMCAD combines a state - of - the - art graphical user interface (GUI), an extensive XML Chemical component database, a large library of thermodynamic data, and a library of the most Common unit operations to give users the ability to provide significant and measurable returns on their investigations. investment. In addition, the program is customizable to allow custom chemicals, thermodynamics, unit Operations, Calculations, and reporting - all ingredients for a powerful user experience.

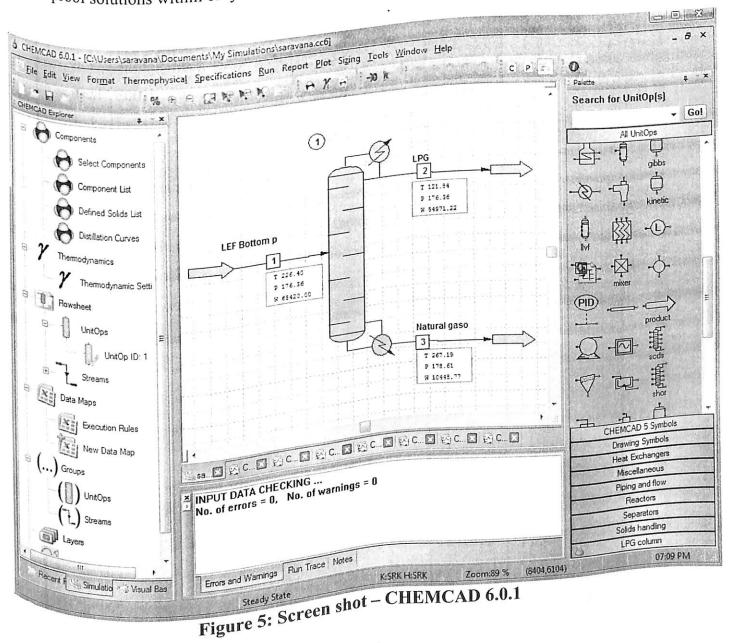
CHEMCAD is capable of modelling continuous, batch, and semi - batch processes, and it can simulate both steady - state and dynamic systems. This program is used extensively around the World for the design, operation, and maintenance of chemical processes in a wide variety of industries, including oil and gas exploration, production, and refining; gas processing; commodity and specialty chemicals; pharmaceuticals; bio fuels; and process equipment manufacturing.

Within all of these industries, chemical engineers work every day with CHEMCAD to address a variety of challenges:

- Initial design of new processes
- Optimization or de bottlenecking of existing processes
- Performance monitoring of processes
- Design and rating of process equipment such as vessels, columns, heatexchangers, piping, valves, and instrumentation

- Evaluation of safety relief devices
- Heat exchanger sizing
- Pressure and flow balancing of complex piping networks
- Reconciliation of plant data
- Economic comparisons of process alternatives
- Advanced process control (APC), including model predictive control (MPC), Real time optimization (RTO), and operator training systems (OTS)
- Binary interaction parameter (BIP) regression from process or lab data

CHEMCAD is capable of delivering the results you need to stay competitive in an increasingly fast and fluid global market. Easy to learn and highly customizable, CHEMCAD can put future - proof solutions within easy reach of our engineering staff



4.2 Methodology for Simulation of LPG Column

4.2.1 Creating a New Simulation

Start by creating a new simulation and giving it a name. To do this, launch CHEMCAD and then Select File > Save to open the Save As dialog box. Navigate to the directory where you want to store the simulation (try My Simulations, located under My Documents) and give your simulation a name, leaving the type as CHEMCAD 6 (*.cc6). Then click Save to create the file and return to the main CHEMCAD window.

4.2.2 Selecting Engineering Units

Select Format > Engineering Units to open the Engineering Unit Selection dialog box. The English units option is the default and is currently highlighted. To change the engineering units system, you would click the Alt SI, SI, or Metric button; you could then change any of the individual Units as well. For this tutorial, you will use English units, so click Cancel to exit this dialog box without making changes.

4.2.3 Drawing the Flow sheet

As described in Chapter 5, creating a flowsheet is a matter of placing UnitOp icons on the Screen, connecting them with streams, and then adding various graphical objects to enhance the drawing. Placing UnitOps. Click the Distillation column tool (see Figure 6). Drag from the Palette to the workspace

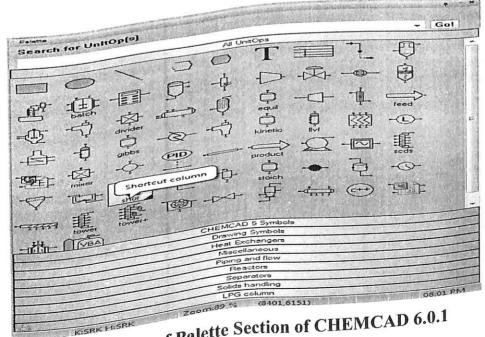


Figure 6: Screen Short view of Palette Section of CHEMCAD 6.0.1

4.2.4 Drawing Streams

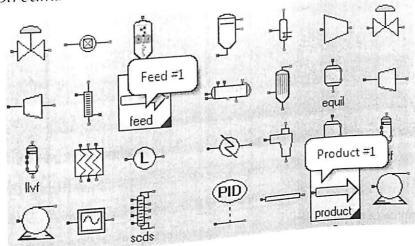
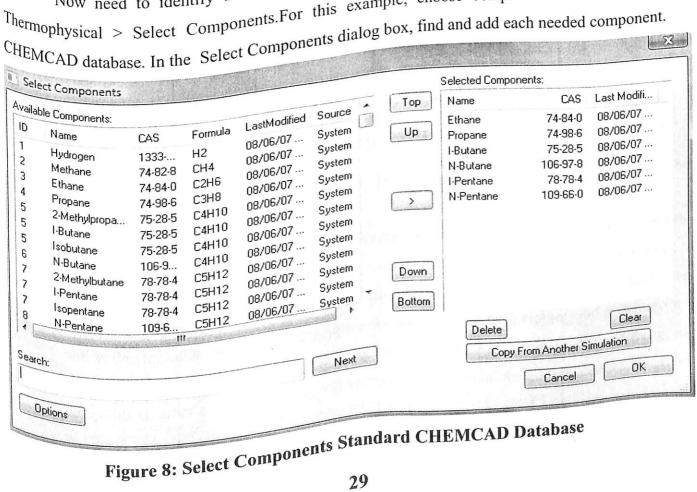


Figure 7: Screen short view of feed and product to distillation column

Click the feed tool (see Figure 7). Drag from the Palette to the workspace place it to feed side of distillation column Then click the product tool (see Figure 7), drag from the palette to the Workspace place it to product side.

4.2.5 Selecting Components

Now need to identify the components to be used in this simulation. Start by selecting Thermophysical > Select Components. For this example, choose components from the standard Chro-



4.2.6 Defining the Feed Streams

Efficient way to define a single stream is to double - click the stream line. Double - click the line for stream 1, to bring up the Edit Streams dialog box.

| | Cancel | OK |
|------------------------------------|--------------|----|
| Flash | Cano | |
| Stream No. | 1 | |
| Stream Name | LEF Bottom p | |
| | 226.4 | |
| Temp F | 176.3554 | |
| Pres psia Vapor Fraction | 1 | |
| Vapor Fraccion Enthalpy MMBtu/h | -59.97756 | |
| | 29674 | |
| Total flow | kg/h | |
| Total flow unit | kg/h | |
| Comp unit | 163.571 | |
| Ethane | 12112.81 | |
| Propane | 5025.787 | |
| I-Butane | 7343.117 | |
| N-Butane | 2348.522 | |
| I-Pentane | 2680.197 | |
| N-Pentane | | |
| | | |

Figure 9: The Edit Streams Dialog Box

- 1. The Stream Name field can display a stream label of up to 16 alphanumeric characters. This
- 2. The next four fields Temp F, Pres psia, Vapor Fraction, and Enthalpy MMBtu/h—are the thermodynamic properties of the stream. According to the Gibbs Phase Rule, once a mixture's composition is given, specifying any two of these four thermodynamic properties will define the other two. As such, defining the composition, temperature, and pressure for a mixture uniquely defines its vapor fraction and enthalpy. Alternatively, defining the composition, pressure, and enthalpy will uniquely define the mixture's temperature and vapor fraction.

- 3. Since enthalpies are calculated relative to a datum, the calculation of any given stream enthalpy is an involved process which is prone to errors. For this reason, CHEMCAD does not permit to enter stream enthalpy as a constraint.
- 4. In addition to defining the stream's composition, define exactly two of the following properties: temperature, pressure, and vapor fraction. The two variables that specify will display as red text, while the third variable and the value enthalpy will be displayed in black
- 5. An exception to this convention allows to add heat duty with an empty stream. If specify a total component flow rate of zero, specify a temperature, pressure, and enthalpy rate. A stream defined this way is treated as a heat duty, and is added to the heat balance of the unit. The temperature and pressure are arbitrary for this situation.
- 6. The Total flow unit and Comp unit fields work together to provide a variety of ways to define stream compositions. If the selected comp unit is mole, mass, or volume fraction (either globally or locally), then the Total flow unit selection is available. If the selected comp unit is a flow or amount option, then the total flow rate becomes the sum of the component flow rates, and the Total flow unit selection is not available.

4.2.7 Selecting Thermodynamic Options

The Thermodynamics Wizard appears. This tool can suggest thermodynamics options to use With this simulation. CHEMCAD's Thermodynamics Wizard works like as follows.

- First, it looks at the component list and decides what general type of model is required, i.e., equation-of-state, activity model, etc.
- Second, it looks at temperature and pressure ranges that you provide and decides which equation within a given category is best at the limits of those ranges.
- If the method is an activity model, the program then looks at the BIP database to see which model has the most data sets for the current problem. It then calculates the fractional completeness of the BIP matrix. If that fraction is greater than the BIP threshold parameter, it uses the chosen activity method; if not, it uses UNIFAC.

parameter, it uses the crosser of the Thermodynamics Wizard is no replacement for engineering judgment. This tool uses an alphaalgorithm based on general rules, and is therefore fallible. The suggested model might not always be the based

Selecting thermodynamic options basically means selecting a model or method for calculating the best model for the system. Vapor-liquid (or vapor-liquid-liquid) phase equilibrium (called the K- value option) and selecting a method or model for calculating the heat balance (called the enthalpy option). The commands for these selections are located on the Thermophysical menu.

CHEMCAD has a library of dozens of K-value models with a variety of options and about 12 enthalpy models. Making the proper selection from these libraries can sometimes be difficult. For the purposes of this tutorial, assume that you want to use the Peng-Robinson method for both the K-value and enthalpy calculations. Follow these steps to select your thermophysical options:

- 1. Accept the default temperature and pressure ranges in the Thermodynamics Wizard and click
- 2. Click **OK** again to accept the wizard's suggested method of SRK.
- 3. When the Thermodynamic Settings dialog box opens, find the Global KValue Option selection, in the upper left corner of the K - Value Models tab. The current setting is SRK, but for the purposes of the tutorial, you'll need to select the Peng - Robinson model. Click the down arrow at the right end of the selection box to view a long list of K - value choices,

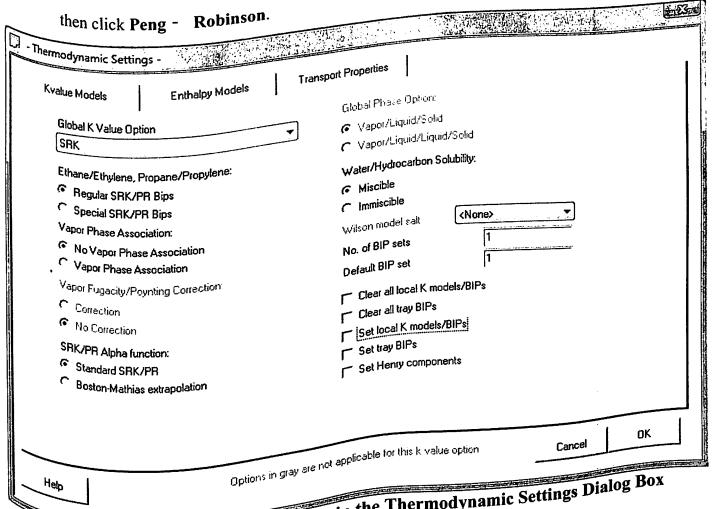


Figure 10: The New K-value Selection in the Thermodynamic Settings Dialog Box

4. Now click the Enthalpy Models tab. The Peng - Robinson method has already been entered as the Global Enthalpy Options selection; this was done automatically because you chose Peng - Robinson as your K - value method. While you do have the option to override this choice, in this case you'll need to keep the Peng - Robinson model; leave all settings as they are and click \mathbf{OK} to return to the main CHEMCAD workspace.

4.2.8 Sizing

After running your simulation, click the UnitOp representing the distillation column and select Sizing > Distillation; choose either Trays or Packing, based on the type of

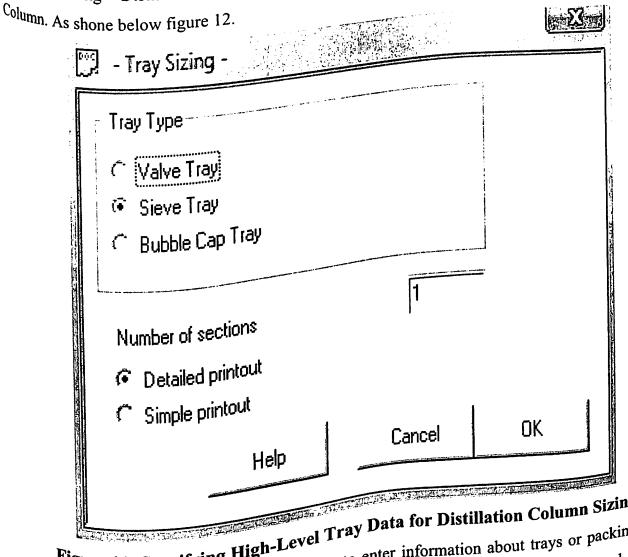


Figure 11: Specifying High-Level Tray Data for Distillation Column Sizing The resulting dialog boxes prompt you to enter information about trays or packing and the The resulting dialog boxes prompt you to chest anisotration about mays of real culation methods. To sizing column, generally need to enter some geometry and hydraulic parameter. parameters to complete column sizing. As shown below figure 13.

| - Sieve Tray - | | | · · · · · · · · · · · · · · · · · · · | | |
|--|-------------------------------------|--|--|--|--|
| Starting Stage Ending Stage Tray diameter Tray spacing No. of passes Hole A / Tot A Weir height Flood correlation Downcomer Clearance Optional flow area Side width Center width Off-center width Off-side width | 1 44 2 1 0.1 0.166667 Fair 0.145833 | ft | Flood percent | Triangular pitch 0.0520833 0.0065 nator <none> <none> oers</none></none> | Section: 1 ft ft ft ft ft ft ft |
| Downcmr A/ Tot A | 0.12 | المعارض والمراجع والمعارضون | The second secon | Cancel | OK |

Figure 12: Specifying Detailed Tray Data

Based on flow sheet values and sizing input, CHEMCAD returns column geometry Based on flow sheet values and SIZINE information such as height and diameter. It also provides hydraulic performance information such as height and diameter. predicted amount of flooding.

To run the simulation, click the Run All button on the toolbar. The program first rechecks the lists 4.2.9 Run the Simulation To run the simulation, click the Run All button. In this case, you should have no errors, although. although you will have warnings about estimates you have not given. You can ignore these warnings about estimates will then proceed. and proceed by clicking Yes. The calculation will then proceed.

When the run finishes, a message box appears: Recycle calculation has converged. To close When the run finishes, a model of dialog box and clear the screen, click OK.

5. SIMULATION REPORTS FOR LPG COLUMN

5.1 Stream C

| .l Stream Composition | | 7.50 | · · · · · · · · · · · · · · · · · · · | |
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| CHEMCAD 6.0.1 Page 1 | . /00 / | | | |
| Job Name: LPG Column | Date: 04/30/2 | | 2 | |
| | 1 | 2 | 3 Natural gaso | |
| Stream No. | LEF Bottom P | LPG | 267.2344 | |
| Stream Name | 1EF 226.4000* | 121.8370 | 178.6497 | |
| Ath B | 176.3554* | 176.3554 | -9.6754 | |
| pres psia | -59.978 | -61.321 | 0.00000 | |
| All para la | 1.0000 | 0.00000 | 144.8203 | |
| | 1.00 | 1095.5776 | 10448.7793 | |
| Total lbmol/h | 1240.3984 | 54971.1758 | 266.9051 | |
| Total lb/h | 65420.0000 | 1634.0311 | 54956.19 | |
| CAL . | 1900.9370 | 415748.28 | 54950.12 | |
| Total and it its/in | 470704.69 | | 0.0000 | |
| Flows and A scin | | 360.6125 | 0.0000 | |
| Ethan Ib/h | 360.6123 | 26704.1680 | 0.0000 | |
| Propaga | -6704.1730 | 11079.9619 | 0.0089 | |
| T)114 | .4079.9055 | 16188.7773 | 0.0009 | |
| | . c109.802' | 474.8668 | 4702.7437 | |
| W-Butane | -477 0000 | 162.7892 | 5746.0269 | |
| 1-pentane | 5908.8232 | 102.1002 | | |
| 1 'Vill's- | 2300° | | | Organica II |
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| or Help, press F1 | | CONTROL OF THE PROPERTY OF THE | Carrie Control | |
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5.2 Feed Stream Property

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| CHEMCAD 6.0.1 Page 1 | | | |
| 20 T | | | |
| Job Name: feed stream I | property Dat | ;e: | |
| 04/27/2009 Time: 21:15 | 5·22 | | |
| rime: 21.1. | | | |
| Stream No. | 1 | | |
| 15 | EF Bottom P | | = |
| · ^- | | | |
| Molar flow lbmol/h | 1240.3984 | | |
| Mass flow lbmol/h Temp F | 65420.0000 226.4000 | | |
| Temp F | 226.4000 176.3554 | | |
| Pres Psia | 1,000 | | |
| Enth Mole fraction | _59.978 | | |
| (C P 11 | 069.3230 | | 1_ |
| βG 2. | 627.7141 | | |
| Std. sp gr. wtr = 1 bec. sp gr. air = 1 | 0.551 | | |
| Std, Sp gr . $wtr = 1Deg_{ree} Sp gr. air = 1$ | 1.821 | | |
| | 125.1802 | | |
| Average mol wt | 52.7411 1.4645 | | |
| WhCAS ID/TES | 44671.3594 | | |
| Actual dens lb/ft3 Std liq ft3/hr Vap 60F safb | | | |
| | 1900.9375 470704.6875 | | |
| ' 16 ' - CLII | | | |
| | 1240.3984 | | ļ |
| Age TOM TDWOT\U | 400 .0000 | | |
| MON TONI | 52.7411 | | |
| JOY A MOT ME | 1.4645 | | |
| JOH, JOHN TES | 44671.3594 | | The state of the s |
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| Std liq ft3/hr Std vap 60F soft | 1900.5375 | | |
| Ch vap 60F sash | 1900.5 470704.6875 470704.8714 | | |
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| P. tgCto. | 0.8628 | | L |
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5.3 Product Stream Property

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| natiew + 15 + | • B 1 0 | 12 · · · 12 · · · · 13 · · · · · 12 · · · · 15 · · · · 16 · · · · 17 · · · · 18 · | |
| 11.1.1.2.1.3.1.4.1.5.1.5.1. | 7 - 1 - 5 - 1 - 9 - 1 - 10 - 1 - | 11 - 1 - 12 - 1 - 13 - 1 - 12 - 1 - 15 | |
| CHEMCAD 6.0.1 | | | |
| Page 1 | | | |
| | | zima. | |
| Job Name: saravana I | 04/27/2 | 009 Time: | |
| 21:24:38 | ace. Var | | |
| | | 3 | |
| Stream No. | 2 | 1 4850 | |
| | LPG | Natural gaso | E |
| Name | | 144.8200 | |
| Name Overall | 1095.5781 | 144.0200 | |
| | 1095.375 54971.2031 | 10448.7607 267.2919 | |
| Mass flow lb/h | 121.8376 | 178.7765 | |
| Temp F | 121.8574 176.3554 | 0.0000 | |
| Pres Psia | 0.0000 | -9.6751 | |
| | -61.321 | 378.2989 | |
| Enth MMBtu/h | 249.5892 | 489.8495 | L |
| Po F | 618.6827 | 489.647 0.627 | |
| Pc Psia Std | 0.539 | 2.491 | |
| * U _ | 1.732 | 94.1471 | |
| Std. Sp gr. wtr = 1 $Degree 3r$ air = 1 | 1.752 | 72.1500 | |
| Degree API | 131.0788 | 30.3757 | |
| Average mol wt Actual der | 50.1755 | 343.9847 | |
| Actual dens lb/ft3 | 30.2482 | 266.9068 | |
| Actual dens lb/ft3 Std lig ft3/hr | 1817.3358 | 266.9000 | 1 |
| 4CV TCO\III | AA 112.79 | 54956.0898 | |
| Std liq ft3/hr Std vap 60F scfh | 1634. 028 415748. 4688 | 144.8200 | |
| A POTII | | 10448.7607 | |
| '401. 474U DD 117 | 1095.5781 | 72.1500 | |
| | - 41 / / / / / | 30.3757 | |
| | | 243.98 4 7 | |
| you dens 15/2+3 | 30.2482 | a66.9068 | |
| Actual dens lb/ft3 Std Vol ft3/hr | 1817.3358 1817.0298 | -4056,0898 | |
| Sta liq ft3/hr | | 56.0904 | |
| ch vap 60F scfb | 1634.4688 415748.4688 35.7113 | 0.0622 | |
| 2 stu/lbmol-E | 0509 | 0.1092 | |
| Viactor | 09573 | 0.0333 | |
| If CD | A 0400 | 3 | |
| Surf tens dyne/cm | 5.5020 | | |
| | | | |

5.4 Topology Report

Page 1 CHEMCAD 6.0.1

Page 1

Page 2

Date: 04/28/2009 Time: 06:20:59 Job Name: LPG Column

FLOWSHEET SUMMARY

Stream Numbers EquipmentLabel

1 -2 -3 1 TOWR

Stream Connections

| Stream | Equip | ment |
|--------|-------|------|
| 1 | From | То |
| 1 | | 1 |
| 2 | 1 | |
| 3 | 1 | |
| | 1 | |

5.5 Mass and Energy Balance

CHEMCAD 6.0.1 Time: 06:32:51

Job Name: saravana Date: 04/28/2009

Calculation mode : Sequential Flash algorithm : Normal

 $e^{quipment}$ Calculation Sequence

 N_0 recycle loops in the flowsheet.

CHEMCAD 6.0.1

Date: 04/28/2009 Time: 06:32:51 Job Name: saravana lb/h Output Input 360.612 $^{0}v_{e_{rall}}$ Mass Balance 1bmol/h 360.612 Output 26704.179 26704.174 11.992 11079.966 Input $\epsilon_{th_{\texttt{a}n_{\texttt{e}}}}$ 11079.964 605.592 16188.813 11.992 16188.803 $\mathfrak{p}_{\mathsf{ropane}}$ 190.630 5177.582 605.⁵⁹² 5177.605 I-Butane 278·527 5908.816 190.630 N_{Butane} 5908.823 71.761 278.527 i.Pentane 65420.000 81.896 71.762 N-Pentane 65420.000 81.896 1240.398 r_{otal} 1240.398

5.6 Thermodynamics Report

CHEMCAD 6.0.1

Page 1

Job Name: LPG Column Date: 04/28/2009 Time: 06:27:49

COMPONENTS

| MILONE | nts | | | Formula |
|--------|-----|---|-----------|---------|
| 1 | ID | # | Name | С2Н6 |
| 2 | | 3 | Ethane | С3Н8 |
| 3 | | 4 | Propane | C4H10 |
| - | | 5 | I-Butane | C4H10 |
| 4 | | 6 | N-Butane | C5H12 |
| 5 6 | | 7 | I-Pentane | C5H12 |
| 0 | | 8 | N-Pentane | |
| | | | | |

THERMODYNAMICS

~value model : SRK Enthalpy model : SRK Liquid density : Li

Library

Std vapor rate reference temperature is 60 F. Atmospher: Atmospheric pressure is

5.7 Net Flow Report

Page 1 CHEMCAD 6.0.1 Time: 22:29:37 Date: 04/27/2009 Job Name: LPG Column 1 $^{\mbox{Unit}}$ type : TOWR Eqp # Unit name: Duties Product Net Flows MMBtu/h Feeds lb/h Vapor -19.01 lb/h Liquid Temp54971.28 Stg lb/h Pres lb/h F psia . 1 79650.22 121.8 134621.36 176.36 2 82165.73 142.8 137137.00 176.41 83566.66 154.8 138537.81 176.46 161.9 83843.00 138814.08 176.51 5 167.0 138450.00 83478.83 65420.00 176.56 6 171.6 82868.21 137839.34 176.62 7 176.2 67658.55 57209.70 176.67 8 70358.48 59909.64 189.4 176.72 g 72185.00 61736.09 198.6 176.77 10 62908.22 205.4 73357.06 11 176.82 63758.44 74207.29 211.3 12 176.88 64564.36 217.3 75013.21 13 176.93 65505.59 75954·45 223.6 14 176.98 66654.29 77103.¹³ 230.2 15 177.04 67971.89 78420.73 236.7 16 177.09 69356.38 79805.21 242.7 17 177.14 70678.48 81127.31 247.9 18 177.20 71841.59 82290.44 ²⁵².1 19 177.25 72798.16 83247.00 255.3 50 177.30 73544.45 83993.29 257.7 2) 74104.77 177.35 84553.60 ²⁵⁹.5 55 74514.27 177.40 84963.10 260.8 53 177.46 74808.13 85257.00 261.7 24 75013.29 177.51 85462.12 262.3 25 75156.92 177.56 85605.76 262.8 56 75257.15 177.61 85706.00 263.1 27 75329.21 85778.03 177.67 263.3 58 75379.72 177.72 85828·⁵⁵ ²⁶³.5 59 75413.58 177.77 85862.41 263.7 30 75426.64 177.83 85875.47 263.8 31 75432.02 177.88 85880.85 264.0 35 75430.48 177.93 85879·³¹ 264.1 33 75423.32 177.98 85872.14 ²⁶⁴.2 75410.73 34 85859.55 178.04 264.4 75399.42 35 85848.24 178.09 ²⁶⁴.5 75383.08 36 85831·⁹⁰ 178.14 264.7 37 75361.45 85810.28 178.19 264.9 38 75333.00 85781.81 178.25 ²⁶⁵.1 75299.43 39 85748.26 178.30 ²⁶⁵.4 75259.49 40 7.99185708.31 178.35 265.6 75213.00 41 10448.82 85661.77 178.41 566.0 75158.57 42 85607.39 178.46 ^{266.3} 75097.77 43 85546.⁵⁹ 178.51 ^{266.7} 267.2 178.56 178.62 82868.211 lb/h. Mass Reflux ratio 171.555 F, 1.449 lotal liquid entering stage 7 at

CHEMCAD 6.0.1 Page 1

| Job Name: LPG Column | Date: 04/27/2009 | Time: 21:51:04 |
|----------------------|------------------|----------------|
| | | 4 |

| i In: | F (| IP # 1 | |
|-------------------------------|--|--|-----------------|
| Unit type : TOWR | Unit name: | | |
| Sta | - 4 5 | 176.36 psia | Y/X |
| Stage # 1 | 121.84 F | 7:7 lb/n | 0.00000 |
| Eth. | Vap 1b/h | 622 5U/01 | 0.00000 |
| Ethane | 0.00000 | -0.00290234 | 0.00000 |
| Propane | 0.00000 | | 0.00000 |
| 1 *~Bii+~ | 0.00000 | - 0 4 E G h D 9 J J | 0.00000 |
| Dit - | 0 0000 | 200 3UL/9 | 0.00000 |
| ' FARL | ი იიიი | 262 5/021 | |
| | U 00000 | 79650.2188 | |
| Total lb/h | 0.0000 | | |
| | - | 176.41 psia | Y/X |
| ^{Stage} # 2 | 142.75 F | | 3.77018 |
| - | 11/11 1/11 | | 1.54981 |
| Ethane | 003 12000 | | 0.80904 |
| , f Up | 07813 | | 0.63694 |
| | 27134.23242 27134.2369 | | 0.32552 |
| N-Butane I-peri | 27134.2369 39645.42969 | | 0.27034 |
| 1-bergue | 39645.423 | | |
| I-pentane | 39645.42 1116.01355 445.48642 | 82165.7344 | |
| N-Pentane Total lb/h | 445.4861 | | |
| TP\P | 134621.3594 | 176.46 psia | Y/X |
| Stage # 3 | | | 4.00734 |
| 4 3 gge # 3 | 154·84 F | | 1.70022 |
| Eth. | 154.0/h Vap 1b/h | | 0.90733 |
| | Vap 107.1 494.92981 494.72656 | 17205.81617 19200.07617 19200.32422 | 0.72112 |
| Propane I-Butane N-But- | | 19200.0762 41348.32422 41348.88452 | 0.37718 |
| N-p ^{out} ane | 50900.72 30311.86133 30312.87891 | 41348.52 3689.88452 3689.58130 | 0.31567 |
| N-Butane I-Pent | 30311.87891 51880.87891 63306 | 3689.58130 2051.58130 | |
| I-pentane N-pentane | 51880 · 63306 2421 · 63875 | 2051.5616 83566.6641 | |
| N-pentane Total | 2421.635 1126.84875 1126.7 0000 | 83500 | |
| Total lb/h | 1126.840 137137.0000 | 176.51 psia | Y/X |
| St. | | 176.31/h Lig 1b/h | 4.14907 |
| Stage # 4 | 161.90 F | Liq 15/11 59.32879 | 1 79092 |
| Et. 4 | 161.90/h Vap 1b/h 59375 | 59.3207 13983.90625 13983.83594 | 0.96736 |
| Ethane Prop | Vap 15/11 431 59375 431 00000 | 13983.90020 17852.83594 17852.87891 | 0 77273 |
| p _{ropane} I-But- | 431.50000 43910.00000 43910.04102 | 17852.83391 42467.87891 42467.14014 | 0.40935 |
| I-Butane | 43910.0002 30280.04102 30287.09375 | 42467.87014 5776.14014 5776.85010 | 0.34402 |
| N-Butane I-pent- | 30280.04175 57537.09375 57537.59668 | 5776.14010 3702.85010 3702.0000 | |
| I-pentane | 57537.095 57537.59668 4145.49097 | 3702.83020 83843.0000 | |
| M-Pentane Total 15 (| 4145.59097 2233.49097 2233.8125 | 830 - | 157 |
| Total lb/h | 2233.4 ⁹⁰ 138537.8125 | 176.56 psia | Y/X |
| 2r TD\P | 1365- | 176.1b/h Lig 65730 | 4.25607 |
| r ∂ge # | 166.97 F | Liq 15/11 55.65730 | 1.85789 |
|) _{Es.} ** 5 | 166.9/h Vap 1b/h Vap 94116 | | 1.01160 |
| Ethane Prop | Vap 15/11 419.94116 419.07422 | 12 ³⁵³ ·34863 16 ¹³³ ·53516 | 0.81067 |
| propane 1-But- | 419·9412 40688·79688 | 16133.3401 16133.53516 40814.53516 | Λ Δ 3300 |
| I-Butane N-Butane | 40688.79688 28932.63672 | 40814.533641 8117.31641 8117.47852 | 0.36495 |
| M-Butane I-pentane | 28932·79672 58656·85254 | 8117.31852 6004.47852 8281 | |
| | 58656.6354 6231.75977 | 6004.478 83478.8281 | |
| | 6231 · 8527 3884 · 75977 3884 · 0781 | | 107 |
| otal ine | 3884 · /5781 138814 · 0781 | 176.62 psia | Y/X |
| Pentane Total lb/h | 13802 | | 4.35621 |
| stage # 6 | 171.56 F | Liq 15712 53.28003 | |
| F. # 6 | 171 · 50/h Vap 16/26962 | J - | |
| Ethane 6 | Vap 15/11 416.26962 | | |
| | 410 | | |

| Propane | CHEMCAD 6.0.1 | | 04 | |
|--|--|--------------|---------------------|--------------|
| Propane | | 54/27/200 |) Time: 21:51:04 | |
| 1-9utane | Name: LPG Column | Dite: U4/2/ | 27703 | 1.91991 |
| 1. | Propano | 2057 66016 | 11342.87793 | 1.05256 |
| Tentane | 1-Butana | 37213 31055 | 14415.60211 | 0.84578 |
| Stage # 7 | ""Bilt > = - | 57003.30078 | 37578.7103 | 0.45507 |
| Tentane | L'Yen+- | 9573.02734 | 10504.0007 | 0.38439 |
| Stage # 7 | · Fenta- | 6186.38770 | 9973.37 | |
| Stage # 7 | Total lb/h | 138450.0000 | | |
| Ethane 413.89233 8705.40039 1.094 i-Butane 25495.62500 27665.23828 0.477 i-Butane 10959.79883 10269.07227 i-Pentane 1055.50293 67658.5469 i-Pentane 21659.36035 1214 116.05371 i-Pentane 10569.36035 32386.89648 0.550 i-Pentane 27665.19531 11983.77148 0.470 i-Pentane 27665.9533 i-Pentane 27665.9533 1216.05371 i-Pentane 3688.07520 8890.53320 1.065 i-Pentane 4688.07520 8890.53320 1.065 i-Pentane 10.44367 2374.8944 1.306 i-Pentane 32386.85742 13848.28516 i-Pentane 32386.85742 13848.28516 i-Pentane 32386.85742 13848.28516 i-Pentane 32386.85742 13848.28516 i-Pentane 5489.13818 72185.0000 i-Pentane 34871.75781 1216.05332 1.376 i-Pentane 34871.75781 11675.0366 i-Pentane 34871.75781 1310.30566 i-Pentane 35307.50781 19118.39063 0.58 | | | 176.67 psia | Y/X |
| Richane | ^{Sta} ge # 7 | 176.17 F | - · - I D/II | 4.45814 |
| Alane | | tran lb/h | | 1.98343 |
| Table 1945 1965 | p. clugue | 113 89233 | -AC 10033 | 1.09471 |
| Notation 25495.6230 27665.2305 0.407 1.Pentane 10959.79883 10269.07227 0.404 1.Pentane 10959.79883 10269.07227 107658.5469 10763.2469 10763.2469 10.44367 1.Pentane 10.44367 1.Pentane 10.44367 1.Pentane 10.44367 1.Pentane 10.44367 1.Pentane 10.569.36035 32386.89648 0.550 1.Pentane 10.569.36035 32386.89648 0.550 1.Pentane 10.569.36035 32386.89648 0.550 1.Pentane 10.569.36035 1.Pentane 10.569.36035 1.Pentane 1.Pentane 10.568.45020 1.Pentane 1.Pentane 4542.15674 70358.4844 1.Pentane 10.44367 1.Pentane 10.4368 1.Pentane 10.44367 1.Pentane 10.4368 1.Pentane 10.43688 1.Pentane 10.4 | ropane | 22047 050/8 | 20 ANU 20 | 0.88200 |
| Stage # 9 198.59 F 198.1647 70358.4844 Total 1b/h 1978.0959.79883 176.77 psia 177.77 psi | Natione | 05/05 62500 | | 0.47791 |
| Tentane | DII + _ | r 2767 41000 | | 0.40461 |
| Stage # 8 | | 10050 19000 | 10269.072 | |
| Stage # 8 189.43 F Liq lb/h 4.682 Ethane 42.13249 4688.07568 1.217 Propane 8705.40039 10073.24609 0.991 LButane 10569.36035 32386.89648 0.550 M-Butane 27665.19531 1198.377148 0.470 L-Pentane 5685.45020 11216.05371 11216.05371 M-Pentane 4542.15674 70358.4844 1207 Total lb/h 57209.6953 176.77 Psia Y/ Stage # 9 198.59 F Liq lb/h 4.825 Liq lb/h 2.49081 2.273 Sthane 10.44367 2.49081 2.273 Ethane 10.4367 2374.89941 1.30 L-Butane 4688.07520 8890.53320 1.065 L-Butane 10.4367 2374.89941 1.30 L-Pentane 7261.88742 13848.2856 0.518 L-Pentane 7261.8818 72185.0000 72185.0000 R-Pentane 7249081 149.52332 1 | Total | 9155.50233 | | |
| Stage # 8 | at Ip/µ | 137839.545 | 176 72 Psia | Y/X |
| Ethane Vap 1b/h 10.47568 1.217 Propane 8705.40039 4688.07568 1.217 Propane 8705.40039 10073.24609 0.991 1Butane 10569.36035 32386.89648 0.550 N-Butane 27665.19531 11983.77148 0.470 N-Pentane 5685.45020 11216.05371 70358.4844 Total lb/h 57209.6953 176.77 psia 176.77 psia Stage # 9 198.59 F Liq 1b/h 2.49081 2.273 Lindane 10.44367 2.49081 2.273 2.49081 2.273 Rthane 10.44367 2374.89941 1.305 2.49081 1.305 Perbane 4688.07520 8890.53320 1.068 0.516 N-Butane 10073.24414 34471.79688 0.604 N-Pentane 32386.85742 138448.28516 0.516 N-Pentane 32386.85742 13848.28516 0.516 N-Pentane 5489.13818 72185.0000 72185.0000 Rtha | Stage " | -2 13 F | | 4.68214 |
| Sthane 42.13249 4688.07300 1.217 Ptopane 8705.40039 10073.24469 0.991 1.Butane 10569.36035 32386.89648 0.550 N-Butane 27665.19531 11983.77148 0.470 1.Pentane 5685.45020 11216.05371 11216.05371 N-Pentane 4542.15674 70358.4844 70358.4844 Potal lb/h 57209.6953 176.77 psia Y/ Stage # 9 198.59 F Liq lb/h 2.49081 2.273 Stage # 9 198.59 F Liq lb/h 2.49081 2.273 Stage # 9 198.59 F Liq lb/h 2.49081 2.273 Stage # 10 10.44367 2374.89941 1.305 N-Butane 10073.24414 34871.79688 0.604 N-Pentane 7261.87842 13848.28516 0.518 N-Pentane 7261.87842 13848.28516 0.518 N-Pentane 7261.87842 12196.9316 72185.0000 Stage # 10 205.40 F Liq lb/h | | - 1n/u | | 2.15514 |
| Copane | Ethane | | | 1.21776 |
| N-Butane | . FUM . | -ar 1000 | | 0.99139 |
| T-Pentane | | A 1000 | 32386.89648 | 0.55062 |
| N-Pentane | n-Butane | | 11983.771 | 0.4/001 |
| Total lb/h 57209.6953 Stage # 9 198.59 F Liq 1b/h 4.825 Ethane 10.44367 2.49081 2.273 2.49081 2.273 1.305 Ptopane 4688.07520 8890.53320 1.063 1.064 1.068 1.068 1.Butane 10073.24414 34871.79688 0.604 0.516 1.3848.28516 0.516 0.516 N-Butane 32386.85742 13848.28516 0.516 1.76.82 psia 72185.0000 7.76.82 psia 72185.0000 | N. Pentane | | 11216.03344 | |
| Stage # 9 | Totale | | 70350. | |
| Stage # 9 198.59 F Liq 1b/m 2.49081 2.273 Ethane 10.44367 2374.89941 1.305 Propane 4688.07520 8890.53320 1.065 1.Butane 10073.24414 34871.79688 0.604 M.Butane 32386.85742 13848.28516 0.518 1.Pentane 7261.87842 12196.93164 0.518 N.Pentane 7261.87842 12196.93164 12196.93164 N.Pentane 5489.13818 72185.0000 72185.0000 70tal lb/h 59909.6367 176.82 psia Y. Stage # 10 205.40 F Liq lb/h 4.93 Ethane 2.49081 1149.52332 1.37 Propane 2.49081 1149.52332 1.37 I.Putane 34871.75781 16175.00488 0.55 I.Putane 34871.75781 16175.00488 0.55 I.Pentane 9126.31709 73357.0625 73357.0625 Total lb/h 61736.0898 176.88 psia Y Stage # 11 211.33 F 0.13046 2.43 Vap lb/h | lb/h | 57209.6953 | .76 77 psia | Y/X |
| Ethane | Stan | | 1/0. 1b/h | 4.82948 |
| Ethane 10.44367 2374.89941 1.305 1-popane 4688.07520 8890.53320 1.065 1-Butane 10073.24414 34871.79688 0.604 N-Butane 32386.85742 13848.28516 0.518 N-pentane 5489.13818 72185.0000 Stage # 10 205.40 F Liq 1b/h 2.368 Propane 2374.89966 7414.10498 1.12 1-butane 34871.75781 16175.30566 1-pentane 34871.75781 16175.30566 N-Butane 34871.75781 16175.30566 N-butane 34871.75781 16175.30566 N-pentane 9126.39453 13310.30566 N-pentane 6470.01709 Total lb/h 61736.0898 N-Butane 1b/h 61736.0898 N-Butane 34871.75781 16175.30566 N-pentane 6470.01709 Total lb/h 61736.0898 N-Butane 35307.50781 176.88 psia | _{-d} ge # 9 | 198.59 F | 2.49081 | 2.273/2 |
| Copane | Eth- | | | 1.30500 |
| ## ## ## ## ## ## ## ## ## ## ## ## ## | propa- | 10.415 | 8890.53320 | 1.00970 |
| 10 10 10 10 10 10 10 10 | 1-Bnt | 4688.01 | 34871.72516 | 0.51837 |
| Tentane 7261.87842 72185.0000 Total lb/h 59909.6367 Stage # 10 205.40 F Liq 0.57677 2.366 Ethane 2.49081 1149.52332 1.37 Ethane 2374.89966 7414.10498 1.12 I.Butane 34871.75781 16175.00488 0.55 I.Pentane 9126.39453 13310.30566 I.Pentane 9126.39453 13310.30566 I.Pentane 6470.01709 73357.0625 Total lb/h 61736.0898 Stage # 11 211.33 F 16.88 psia Y I. Stage # 11 21.33 F 16.88 psia Y I. Stage # 11 21.33 F 16.88 psia Y I. Stage # 11 21.33 F 16.88 psia Y I. Stage # 11 21.33 F 16.88 psia Y I. Stage # 11 21.33 F 16.88 psia Y I. Stage # 11 21.33 F 16.88 psia Y I. Stage # 11 21.33 F 16.88 psia Y I. Stage # 11 21.33 F 16.88 psia Y I. Stage # 11 21.33 F 16.88 psia Y I. Stage # 18 Psi | N-Butan | 10073.2 | 13840.02164 | 0.32 |
| Total lb/h Stage # 10 205.40 F Vap lb/h Vap lb/h Vap lb/h Vap lb/h Propane 1.Butane 1.Pentane Total lb/h Stage # 11 Stage # 11 Stage # 11 Stage # 12 2.49081 1.149.52332 1.37. 1.249081 1.149.52332 1.37. 1.25332 1.37. 1.26888 1.29 1.290488 1.205.40 F 1.29081 1.29081 1.29081 1.29081 1.2909.66 1.29081 1.29081 1.29081 1.29081 1.29081 1.29081 1.29081 1.29081 1.29081 1.29081 1.29081 1.29081 1.29081 1.20088 1.20081 1.20 | % | 32360.87842 | 12196. 0000 | |
| Stage # 10 Stage # 11 Stage | To entanc | 5489.13818 | 72100 | |
| Stage # 10 205.40 F L140.57677 2.363 Ethane 2.49081 1149.52332 1.37 Propane 2.374.89966 7414.10498 1.12 I-Butane 8890.53223 35307.50488 0.64 N-Butane 34871.75781 16175.00488 0.55 I-Pentane 9126.39453 13310.30566 0.55 N-Pentane 6470.01709 73357.0625 73357.0625 Total lb/h 61736.0898 176.88 psia Y Stage # 11 211.33 F 0.13046 2.43 Propane 0.57677 536.30585 1.42 Propane 1149.52332 5904.14648 1.18 N-Butane 7414.10498 34001.51172 0.68 N-Butane 7414.10498 34001.51172 0.68 N-Pentane 35307.50781 1918.39063 0.58 N-Pentane 35307.50781 194646.80371 14646.80371 N-Pentane 12583.38965 74207.2891 | otal lh/h | 59909.6367 | 176.82 Psia | Y/X |
| Ethane | Sta | | 1/0 1b/h | 4.93790 |
| Ethane Propane 1.1249081 1.149.52498 1.125 1.126 1.127 1.127 1.128 1.128 1.129 | ************************************** | 205·40 F | 0.57677 | 1.37114 |
| Propage 2374.89966 7414.54688 0.64 | € _{t k} | Vap 10,0081 | 1149.52332 | 1 12933 |
| Stage # 11 211.33 F 16.175.00488 1.1 | brous | 2.0966 | 74141688 | 0 64510 |
| Sutane 8890.75781 16175.30566 13310.30566 13310.30566 13310.30566 13310.30566 13310.30566 13310.30566 13310.30566 13310.30566 13310.30566 13310.30566 13310.30566 13310.30566 13310.30565 13310.30565 13310.30566 13310.30565 13310.30566 13310.30565 13310.30566 13310.30565 13310.30566 13310.30565 13310.30565 13310.30565 13310.30565 13310.30566 13310.30565 13310.30565 13310.30565 13310.30565 13310.30566 13310.30565 13310.30565 13310.30565 13310.30565 13310.30566 13310.30566 13310.30565 13310.30566 13310.30565 13310.30565 13310.30565 13310.30565 13310.30565 13310.30565 13310.30566 13310.30565 13010.30565 13010. | l'But- | 2374.63223 | 35307.00488 | 0.55582 |
| N-pentane 9126.39453 133157.0625 Total lb/h 61736.0898 Stage # 11 211.33 F 1149/h 5.03 Ethane 0.57677 536.30585 1.42 Propane 149.52332 5904.14648 1.18 N-pentane 35307.50781 19118.39063 N-pentane 35307.50781 19118.39063 N-pentane 11453.11328 14646.80371 N-pentane 11453.11328 14646.80371 | N-Buta- | 8890.75781 | 161/3,20566 | - |
| Stage # 11 | h ent- | 348/1.39453 | 13310.0625 | |
| Stage # 11 211.33 F 114 0.13046 2.43 Pthane 0.57677 536.30585 1.42 Propane 1149.52332 5904.14648 1.18 N.Butane 7414.10498 34001.51172 0.68 N.Butane 7414.10498 34001.51172 0.68 N.Butane 35307.50781 19118.39063 0.58 N.Pentane 35307.50781 14646.80371 N.Pentane 11453.11328 14646.80371 N.Pentane 11453.38965 | To bentar | 6470.01709 | 7355 | - /17 |
| Rethane 0.57677 536.30585 1.42 Perhane 0.57677 5904.14648 1.18 I.Butane 149.52332 5904.14648 1.18 I.Butane 7414.10498 34001.51172 0.68 I.Putane 35307.50781 19118.39063 0.58 I.Pentane 35307.50781 14646.80371 14646.80371 N.Pentane 11453.11328 14646.80371 14646.80371 | of al lb (| 64736.0898 | 176.88 psla | Y/X |
| # 11 | St. | 01, | Lig 1b/h | 5.03240 |
| Vap 157677 530.14648 1.18 1.18 1.18 1.18 1.18 1.18 1.18 1. | _a ge # 11 | 211.33^{+} | 0.13040 | 1.42944 |
| Frond 1.49.52332 5904.51172 0.68 N.Butane 149.52332 34001.51172 0.68 N.Butane 7414.10498 34001.51172 0.68 1.putane 35307.50781 19118.80371 N.Pentane 11453.11328 14646.80371 N.Pentane 11453.38965 24207.2891 | \$h1 | Vap 15,7677 | 536.30565 | 1 18203 |
| N. Butane 1149.328 34001.31063 0.58 N. Butane 7414.10498 34001.31063 0.58 N. Putane 35307.50781 19118.80371 N. Pentane 11453.11328 14646.80371 1453.31328 14646.80371 1583.38965 | prope | 0.53332 | | v (8137 |
| 7414.50781 19116.80371 1. Pentane 35307.50781 14646.80371 1. Pentane 11453.11328 14646.80371 1. Pentane 11453.38965 | (~, rdpc | 1149.52 | 34001.312 | 0.58937 |
| N-Pentane 35307.11328 1403 11453.11328 1403 11453.38965 24207.2891 | We can- | 7414.50781 | 19118.80371 | 0 • • |
| Pentane 11433 38965 24207.2891 | | 3530 / 1 328 | | |
| Total 1b/h 62908.2188 7583.2 7420.7 7420.7 | "`Pentane | 11453.38965 | 74207.2891 | |
| ^{otal} lb/h 62908.2188 | J | 7503. | 1420 | |
| 42 | etal 1r. | 62908.2188 | 40 | |
| | ±n\ν | 02- | 42 | |

| GREAD 6.0.1 | | | |
|-------------------------------|--|--|--------------------|
| Job Name: LPG Column | Date: 04/2//20 | | |
| ^{Stage} # 12 | 2:7.27 F | 176.93 psia | Y/X |
| - | Vap lb/h | Liq lb/h | 5.12536 |
| Ethane | 0.13046 | 0.02893 | 2.51757 |
| rronana | 536.30591 | 242.12886 | 1.48845 |
| "Blift a | 536.30371 | 4508 57324 | 1.23564 |
| " Dilt > " | 5904.14648 | 21276 759// | 0.71969 |
| · FBnt - | 34001.47656 | 22736.66602 | 0.62395 |
| | 14396.49707 | 16249.05664 | |
| Total lb/h | 8919.88770 | 75013.2109 | |
| | 63758.4414 | : 3 | |
| ^{Stage} # 13 | 2 E 7 F | 176.98 psia | Y/X |
| | 223.57 F | tia 1b/n | 5.22074 |
| Ethane Pro | Vap lb/h | V 0005+ | 2.59961 |
| , f UP - | 0.02893 | 105.97120 | 1.55162 |
| | 242.12888 | 3306.00000 | 1.29336 |
| N-Butane I-po- | 4508.57275 | 27513.86914 | 0.76086 |
| I-bent | 31276.72070 | -020 POID | 0.66178 |
| I-pentane N-pent | 10014 7/555 | 0 1111000 | |
| N-pentane Total | 10500 1300 | 75954.4453 | |
| 17 TP/P | 64564.3633 | an psia | |
| ^{Stage} # 14 | - T | 177.04 psia | Y/X |
| ³⁶ # 14 | 230.17 F | Liq 1b/h 0.00135 | 5.31363 |
| ^६ ۲ ኤ - | $-10/\Pi$ | 45.04261 | 2.68383 |
| p _r opane I-Bu- | V (1002- | 1 / 0 0 0 | 1.61783 |
| I-Butane | 105.97121 | 2331.0525 23174.35156 23174.35164 | 1.35436 0.80507 |
| N-But- | - a c 10000 | | 0.80307 |
| ('P_rque | 10 X 10/- | | 0.70201 |
| N-P-ucaue | - ~ 1 () () () | 20072.430 77103.1328 | |
| lotal ane | | | |
| , 1D\P | 12363.075 | 177.09 psia | Y/X |
| Stage # 15 | | . 10/11 | 5.40477 |
| , 30 # 15 | 236.69 F | Liq 10,0029 | 2.76719 |
| | 10/11 | - 623/4 | 1.68393 |
| propane I-But- | 0.00135 | - 0 81.732 | 1.41546 |
| I-Butane | 45.04261 45.0082 | 1583.010 18731.77148 18736.55859 | 0.84978 |
| M-Butane I-butane | 45.042 2331.09082 23174.31445 23174.29102 | 18731.7711 18731.7711 36026.55859 36026.96680 | 0.74402 |
| h ent- | 23174.329102 | 36026.33000 22059.96680 22059.7344 | |
| 1/1 'Lan- | 26/30. 55176 | 22059.900 | |
| W_pentane Total lb/h | 26758.29176 14345.55176 66654.2891 | | |
| | 66654.20 | 177.14 psia | Y/X |
| stage # 16 | 70 F | | 5.48478 |
| Ft. # 16 | 242.70 F | | 2.84302 |
| Ethane | Vap 10,0029 | 7.51175 | 1.74500 |
| propane Lauta | 0.00 | 1040.81995 | 1.47226 |
| I Butane | | | 0.89187 |
| 1. Butar | 1583.813 | 14590.1201 40250.59375 40250.16016 | 0.78315 |
| W. ent. | 18731. | 40250.33016 23916.16016 23916.2109 | |
| Topentane | 31304.04980 | 23916.1009 79805.2109 | |
| egl rie | 31304.609 16333.04980 67971.8906 | | /st |
| Stage # 17 | 67971.05 | 177.20 psia | Y/X |
| " <i>96</i> " | | | 5.55148 |
| | 247.88 F | 7 00002 | 2.90770 1.79767 |
| $b_{c}^{\mu}q^{\mu}$ | Vap 10/0006 | 2.96535 | 1./9/5 |
| propane I But 2 | 0.51175 | 664.58356 | |
| l'opane Butane | 7.5119 | | |
| A116 | 1040.01 | | |
| | | | |

| 0.0.1 | | -1.51.04 | |
|---|--|---|--------------------|
| Job Name: LPG Column | Date: 04/27/20 | 09 Time: 21:51:04 | |
| | Date. V. D. | | 1.52147 |
| N-Butane | 14590.08789 | 11007.28418 | 0.92870 |
| 'Tenta- | 35528.71094 | 43912.32031 | 0.81748 |
| " rent | 18189.24609 | 25540.16211 | |
| lotal lb/h | 69356.3750 | 81127.3125 | |
| | 69356.5756 | | |
| ^{Stage} # 18 | 052 07 F | 177.25 psia | Y/X |
| | 252.07 F | Liq 1b/h | 5.60428 |
| Ethane | Vap 1b/h 0.00001 | 0.00000 1.14975 | 2.95974 |
| \$ \$0n | 2.96535 | 1.14575 | 1.84041 |
| | 664.58356 | 414.39511 | 1.56152 |
| | 11007.24512 | 8089.31299 | 0.95893 0.84571 |
| | 39190.43750 | 46900.29297 26885.28320 | 0.845/1 |
| N-Pentane Total | 19813.24805 | 26885.2032 82290.4375 | |
| <pre>lotal lb/h</pre> | 70678.4766 | | |
| - TD/U | 706/8.4/0- | 177.30 psia | Y/X |
| ^{Stage} # 19 | 25 22 F | | 5.64442 |
| ³⁶ # 19 | 255.32 F | | 2.99971 |
| Ethane Pro- | Vap 1b/h 0.00000 | A 13941 | 1.87345 |
| | 1 149/5 | - 2 67932 | 1,59255 |
| I-Butane N-But | 11 4 39511 | ~ 1 n40° | 0.98247 |
| N-Butane I-Pent | 8089.27441 | | 0.86774 |
| I-pentane | 8089.21 42178.41016 | 49216.33 | |
| R-Pentane Pentane | | 27953.200 83247.0000 | |
| Total lb/h | 71841.5938 | os. | Y/X |
| St. | /1841. | 177.35 psia | 5.67410 |
| Stage # 20 | 257.73 F | 177.36/h Liq 1b/h | 3.02938 |
| Et L | 10/11 | 0.00000 0.16607 | 1.89807 |
| Ethane Propane | | - Ob /4± | 1 61570 |
| 1 ^p dgu6 | a 13941 | | 1 00011 |
| Mrp_cgDe | - 2 57933 | 4126.335 50936.29688 50937.53516 | 0.88427 |
| 1/2 , dDe | | 50936.2900 50936.2901 28777.53516 28777.2891 | - |
| 4/2 - 1/2 n - | 5823.123 44494.66797 44494.35352 | 28777.5331 83993.2891 | |
| Pentane Pentane Potal lb (| 44494.867 22226.35352 22226.1641 | 83995 | 151 |
| Pentane lotal lb/h | 22226.3334 72798.1641 | 177.40 psia | Y/X |
| Stage # 21 | | 177.40/h Liq 1b/h | 5.69561 3.05089 |
| ». * 21 | 259.50 F | Liq 15/11 0.00000 | 1.91595 |
| GP1 | 259.30/h Vap 1b/h 0.00000 | | 1 63254 |
| eropane 1-bane Mut- | 0.06607 | | 1 01290 |
| I.Butane | | 2890.075 | 0.89631 |
| N'Butane 1'Pentane | 152.93, | 62160° a2438 | O s |
| N_Pentale | 4126.23406 | 52166·44328 52166·73438 29405·736016 | |
| The street and street | 46214.61914 | Q455- | |
| (9) 116 | 4621 ⁴ ·41914 23050·61914 2544·4531 | 46 psia | Y/X |
| Pentane lotal lb/h stage # 22 | 23050·615 73544·4531 | 177.46 Psia 177.46 Psia 1b/h Lig 0.00000 | 5.71095 |
| .46 h | 76 F | Ti9 20000 | 3.06620 |
| Ethane Props | 260.76 F | υ΄ ₂ 2318 | 1.92870 1.64451 |
| $\delta r_{A}^{O\nu}$ | 260.7h Vap 1b/h 0.00000 | | 1.02214 |
| 1 TO IN | 0.06224 | 54.03095 2007.19495 2007.18750 | 0.90491 |
| Nautane 1.pentane | 0.08256 91.28356 91.03467 | 2007.19450 53015.18750 5306.64258 | 0.50 |
| h perane | 91·28367 2890·03467 2890·57031 | 53015.18758 29886.64258 29886.1016 | |
| Thutane Thentane Totaltane | 2890.03431 47444.57031 | 29886.6423 29886.1016 84963.1016 | |
| .dl .16 | 4744 ⁴ ·82031 23678·82034 | ia | |
| Totane Total lb/h Stage | 23678·820 74104·7734 | 177.51 psia | |
| stage # 53 | 191" | ٠. | |
| 23 | 261.65 F | | |
| -3 | 20- | 4.1 | |

| CHEMCAD 6.0.1 | | . 04 | |
|-------------------------------------|--|---|---------------------|
| | | 9 Time: 21:51:04 | |
| Job Name: LPG Column | r:t-: 04/27/2009 | | Y/X |
| 1 | 1 h / h | Liq lb/h | 5.72167 |
| Ethane | Tap lb/h 0.00000 | 0.00000 | 3.07697 1.93768 |
| *Tonan | 0.02318 | 31.83530 | 1.93766 |
| Dist - | 54.05692 | - 20E 66U4U | 1.02862 |
| 1 1117 - | 2007.15527 | 76 1070 | 0.91099 |
| | 48293.30859 | - 0 7 3 3 3 3 3 3 3 | 0.5- |
| | 24159.72852 | 85257.0000 | |
| lotal lb/h | 74514.2734 | | |
| \$ | | 177.56 psia | Y/X |
| ^{Stage} # 24 | 262.29 F | 10/11 | 5.72913 |
| Eth | $n = \frac{1}{2}$ | | 3.08449 1.94399 |
| Ethane Pro- | U ()UUU | 0.00318 0.00318 18.67446 | 1.65890 |
| propane I-Rus | 0.00860 | | 1.03318 |
| N-B | 31.83526 | | 0.91527 |
| 1-p-caue | 1385.62085 | 53921·01727 30569·94727 30562·1172 | • |
| N-D-cutane | . AAEA 7303' | 30569·34 85462·1172 | |
| N-Pentane | - 1506 4327 | 85402 | |
| 1 7p/h | 74808.1328 | 177.61 psia | Y/X |
| app ^f ^c | 75 F | 17/.02/h Liq 1b/h | 0.00000 3.08991 |
| 25 | 262.75 F | Liq 10/10 0.00000 | 1.94848 |
| Ethane | Vap 1b/h 0.00000 | $0.00117 \\ 0.00117 \\ 0.2378$ | 1 6631 ⁰ |
| propane | 2 00310 | $ \begin{array}{r} 0.00278 \\ 10.92378 \\ 10.74176 \\ 652.72656 \end{array} $ | 1 03640 |
| li ^{-p} , rgue | . ~ 67442 | $652 \cdot 72656$ | 0.91828 |
| 1 10 caub | | 652.74156 54106.72656 54106.35938 30835.7578 | |
| I "" "" GDG | | 30835.35936 85605.7578 | |
| H.Pentane Total lb/h | 49199.1320 24843.03320 24843.2891 | 8500 | Y/X |
| 1 40/2 | 24843.032 75013.2891 | 177.67 psia | 0.0000 |
| gde " | | 177.67 h Liq 1b/h 0.00000 | a 09354 |
| 26 | 263.09 F | 0,0043 | a a5157 |
| Ethane | Vap 1b/h 0.00000 | 0.0007 6.37756 6.37421 | 4 666UJ |
| ["hqUb | 0.0001 0.00117 0.02375 | 6.37421 446.37421 | 1.03869 |
| I Was saba | | 446.37424 54172.27734 54172.93750 | 0.92044 |
| | $ \begin{array}{r} 10.925 \\ 52.70221 \\ 652.95547 \end{array} $ | 54172·27750 31080·93750 310706·0000 | |
| | 652 · 7024 49384 · 85547 49384 · 44141 | 25/5 | |
| Total lb/h | 49384.831 25108.44141 25108.9219 | 177.72 psia 177.1b/h | Y/X |
| 1p/h | 25108·441 75156.9219 | 177.72 p | 0.00000 |
| Stage # 27 | | 7 10 - A000 | ~ UAD42 |
| * 27 | 263.34 F | | 1.95357 |
| Ethane | vap 15,000 | 0.00027 3.71877 3.01308 | 1.66802 1.04035 |
| 1 1. Pdp - | 0.50043 | 3.71878 304.81308 304.66016 | 0.92203 |
| | 0.0003 6.37752 6.33463 | 304.81306 66016 54145.83984 | 0.52 |
| 1 Putane | 6.37463 446.33463 41016 | 54145.66013 54145.83984 31323.83983 | |
| I Butane I Pentane N Pentane Totale | 446.3340 49450.41016 49450.02344 | | *** |
| thentane thentane total lb. | 49450.4102 25354.02344 25357.1484 | 77 ps1a | Y/X |
| Total lb/h | 25354.023 75257.1484 | 177.77 psia 177.77 psia 1b/h Liq 0.00000 | 0.00000 3.09720 |
| 1 '36 " | - 2 F | 1.19 00000 | . ለካካኒሀ |
| Ethane | 263.53 F | 0,000 | |
| Propane | vap -0000 | | 1.04161 |
| 1 1645 | 0.50016 | 2·16025 207·90465 207·21875 | 1.0 |
| lopane Mutane Lobutar | 0.0001 3.71873 3.7350 | 207.90405 54041.21875 | |
| Rutane 1. Pentane | $\begin{array}{r} 3.71870 \\ 77350 \\ 304.78906 \end{array}$ | 24. | |
| tane tane entane | 304·77396 49423·78906 | 45 | |
| | 423- | u J | |

| CHEMCAD 6.0.1 | | -1 0/ | |
|---|---|--|-------------------------------|
| | .ca (200 | 9 Time: 21:51:04 | |
| Job Name: LPG Column | Date: 04/27/200 | | 0.92323 |
| II-Da. | | 22577 25911 | |
| ll-Pentane | 25596.92773 | 85828.5469 | |
| lotal lb/h | 75329.2109 | 177.83 psia | Y/X |
| ^{Stage} # 29 | 263.70 F | . 1h/H | 0.00000 |
| | Vap 1b/h | | 3.09897 |
| ^{Eth} ane | 0 00000 | a 00004 | 1.95663 |
| (0na. | ი იიიინ | 1 26014 | 1.67092 |
| | o 16619 | .1 66333 | 1.04262 |
| | 207 86508 | 78120 | 0.92417 |
| I-pentane | 10010 3390° | | |
| N-pentane | 25050 3470- | 85862.30 | |
| Tp/h | 75379.7188 | 177.88 psia | Y/X |
| Stage # 30 | 0.4 F | | 0.00000 |
| 30 ge # 30 | 263.84 F | Liq 10,00000 0.00001 | 3.10231 |
| Ethane | Vap 1b/h 0.00000 | 2 VUUA7 | 1.95861 |
| | 0 0000Z | - 73/02 | 1.67247 1.04351 |
| I-Butane | 2 261100 | 96.40909 96.375 | 0.92492 |
| N-Butane | . 4 67 390 | 96.40575 53619.09375 53619.22656 | 0.52 |
| I pentane | . A UIIZ- | 53619.095 32159.22656 32159.4688 85875.4688 | |
| N-pentane Total | 49144.30102 26125.79102 | 858737 | /3/ |
| 1 70/2 | 26125.75 75413.5781 | 177.93 psia | y/X 0.00000 |
| Stage # 31 | 263.97 F | 177.55/h Liq 1b/h 0.00000 | 2 102/0 |
| , [#] 31 | | 0.0000 | 1 95933 |
| Ethane | Vap 15/10000 0.00000 | | 1 67321 |
| propane I-but | ~ VUUV - | | 1.04429 |
| I N'V _ CQDE | 71/10 | 65.567 53306.71094 53308.14844 | 0.92568 |
| I (" dDo | 96.36951 96.3266 | 53306·7104 32508·14844 32508·8516 | |
| 1 4/V CQDC | 96.3076 48897.22266 48897.31445 | ando | |
| Pentane Total 1b/h | 48897.222445 26432.31445 75426.6406 | og psia | Y/X |
| | 75420 | 177.98 h Liq 1b/h | 0.00000 3.10311 |
| ode # € | 264·10 F | Liq 1b/11 0.00000 | 4 9600' |
| \(\xi_1 \) | 264.10/h Vap 00000 | 0.000727 | 4 67400 |
| brane | 0.5000 | 0.2472 0.56070 44.34609 | 4 04504 |
| propane Inbutane | 0.0005 0.42574 0.52815 | 44.56070 24609 52925.24609 52925.25781 | 0.92642 |
| Mutane I-pentane | 0.42315 65.52815 63.203 | 52925.24081 32909.25781 32909.3125 | |
| | 65.5203 48584.83203 48584.24023 | 32909·2370 32909·3125 85879·3125 | |
| Rentane Pentane Potal la | 48584.8320 26781.24023 26781.20234 | , nsia | Y/X |
| Total lb/h | 26781·2402 75432·0234 | 178.04 psia | 00000 |
| $\begin{cases} st_{ade} & # 33 \end{cases}$ | , J | 1.19 -0000 | 2 10300 |
| Pr | 264·24 F | 0.000 | 4 96002 |
| 1 10 | Vap 12000 | | 1.93490 1.67490 1.04583 |
| 1 1 1 2 2 2 2 2 2 2 2 2 2 2 2 2 2 2 2 2 | 0.5000 | | 0.92719 |
| | 0.00023 0.24723 0.2113 | 62468 - 460 - 1 | 0.72 |
| 1, pe ane | 44.52109 | 23313 1406 | |
| Butane Pentane Pentane | 18203. 1766 | 850 ng psia | y/X |
| Pentane Total lh | 271000 4845 | 178.09 F Liq 1b/h | 0.00000 |
| 1 de 10 D | 7543 20 F | Liq 10,10000 | |
| ξ _{th} 34 | 264.5 h | | |
| Ethane | Vap 15/11 0.00000 | 46 | |
| i e | | 40 | |

| CHEMCAD 6.0.1 | | . 01 | |
|--|---|---|------------------------------|
| | | 9 Time: 21:51:04 | |
| Job Name: LPG Column | Tate: 04.27/200 | | 3.10430 |
| · | | 0.0000 | 1.96177 |
| Propane | 0.0000 | ი იგვვს | 1.67584 |
| 1 ''011+ | 0.14349 | an 52693 | 1.04668 |
| 1 " DII + 5 | 30.21972 | 35930 | 0.92802 |
| · Pant | 47746.60547 | | |
| | · 16.16.35324 | 85859.5547 | |
| lotal lb/h | 75423.3203 | | /٧ |
| | | 178.14 psia | Y/X |
| Stage # 35 | 264.54 F | . 16/11 | 0.00000 3.10423 |
| | $\frac{1}{2}$ | | 1.96243 |
| Ethane Pro- | 0.0000 | 0.00000 | 1.67669 |
| propane | 0.00000 | 0.04828 | 1.04761 |
| I-Butane | n n8323 | 13.90978 | 0.92896 |
| li-Butane i-Par | 20 48730 | 13.905 51295.76563 51295.51172 | 0.52 |
| li-pantane | os 48430 | 51295.703 34538.51172 345848.2422 | |
| Il-Pentane Total | - 0.3 0 / h / J 0 0 | 85840. | |
| 1 7p/h | 75410.7266 | 178·19 psia | Y/X |
| Stan | | 178.19/h Liq 1b/h | 0.00000 |
| ^{Stage} # 36 | 264.71 F | Lig 15/11 0.00000 | 3.10578 |
| l Eth | 1 h/U | ∆11110 ¥ | 1.96385 |
| Ptopane | | | 1.67803 |
| I.Butane | 0.00000 | | 1.04869 0.92998 |
| N-Butane I-Pent | 0.04824 | | 0.92950 |
| (" _dDe | 13.87020 | 50555.64010 35266.81250 35266.8984 | |
| I-Pentane I-Pentane Total It | 13.87453 46573.89453 | 35266.8123 85831.8984 | |
| l otal in | 46573.60938 28811.60938 75399.4219 | - nsia | Y/X |
| 1 10/1 | 75399.32 | 178·25 psia | 0.00000 |
| cg∂e # | . 00 F | 1.19 -0000 | 2 10700 |
| * 37 | 264.90 F | 0.0000 | a 06551 |
| Ethane | Vap 1b/h 0.00000 | 0.00017 0.01617 | 1.67960 |
| 1 Lydne | - 2000 | 0.01018 6.35008 | 1.04993 1.049115 |
| 1 1/2 dha | | 6.35094 49692.08594 49692.82422 | 0.9522 |
| | | 49692·08534 36111·82422 361810·2813 | |
| | | 36111.82423 85810.2813 | |
| Trentane Total lb/h | 45833.776 29539.90625 29539.0781 | goo | Y/X |
| Jp/p | 29539.9002 75383.0781 | 178.30 psia | 0.00000 |
| stage # 38 | | + 10 - 0000 | 2 1 1 0 0 7 |
| ^{8r′} | 265.12 F | • | 1.96752 |
| | Vap 12000 | $0.00034 \\ 0.00934 \\ 0.27190$ | 1.68146 1.05136 1.0249 |
| topane l'But- | 0000 | 0.00950 4.27190 | 0.93249 |
| | 0.00061 0.01614 0.31050 | 4.27194 4.52344 48687.52344 48687.00781 | 0.952 |
| Nautane Lentane Lentane | 0.01010 6.31050 6.31094 | 48687.00781 | |
| Hent's | 6.31094 | 48687.52341 37090.00781 37091.8125 85781.8125 | |
| Totale Total Total Total Total Total Total Total Total | $ \begin{array}{r} 6.31094 \\ 4.21094 \\ 44970.21094 \\ 91406 \\ 30384.914453 \\ 30384.4453 \end{array} $ | gor | Y/X |
| 91, 16 | 3038 ⁴ ·91453 75361·4453 | 178.35 psia 178.1b/h | 0.00000 |
| Total lb/h Stage # 39 | 7530- | 4 i a ~ ~ ~ ~ ~ ~ ~ ~ ~ ~ ~ ~ ~ ~ ~ ~ ~ ~ | |
| 30 | 265.37 F | | |
| kthane 39 | 265.37/h Vap 00000 | 0.00537 | 1.68330 |
| Propane 1, Buta | Vap 15/11 0.00000 | 0.00334 2.86146 2.8766 | 0.93403 |
| | 0.00000 | 2.86146 2.09766 47525.29688 | 0.95 |
| 1 gnrave | $0.00_{-2.32}$ | 47525.29688 | |
| h 60 116 | 0.00932 4.23232 | 47525.09788 38220.29688 382748.2578 85748.2578 | |
| bengane | 43965.63261 | 85740 | |
| N.pentane Potal lb/h | 43965·652 31363·09961 31363·0000 | | |
| | 31363.0990 75333.0000 | _ | |
| | • - | 47 | |

| CHEMCAD 6.0.1 | | • | |
|---|--|---|--------------------|
| 9 | Date: 0: 27/200 | 9 Time: 21:51:0 | 4 |
| Job Name: LPG Column | Date: " 2//200 | • - | |
| ti . | | 178.41 PS14 | Y/X |
| Stage # 40 | 205.55 F | 7 i 7 1 h/n | 0.00000 |
| Fra. | var lo h | 0.0000 | 3.11585 |
| Ethane Pro- | 0.0000 | 0.00000 | 1.97241 |
| Propane I-Butane | 0.00000 | 0.00308 | 1.68601 1.05489 |
| II-Butane | 0.00534 | 1.90505 | 0.93581 |
| | 2.82100 | 46184.59375 | 0.9550- |
| | 42803.21875 | 39521.81641 85708.3125 | |
| lotal lb/h | 32493.39063 75299.4297 | | |
| -1 1p/h | 75299.32 | 178.46 psia | Y/X |
| ^{Stage} # 41 | - cr 07 F | · in/u | 0.00000 |
| | 265.97 F | | 3,11955 |
| Ethane Pro | Vap 1b/h 0.00000 | ~ UUUUU | 1.97548 |
| | 0.0000 | 2 001/2 | 1.68885 |
| I-Butane | a nn305 | 25/23 | 1.05707 0.93785 |
| N-Butane | 1 26540 | - 60000 | 0.93703 |
| I-pentane | 718/3 | 44645.01172 41015.01172 | |
| N-pentane | | 41015.017 85661.7734 | |
| lotal lb/h | 75259.4922 | 178.51 psia | Y/X |
| St. | | 178·31 F | 0.00000 |
| Stage # 42 | 266.32 F | 178.32/h Liq 1b/h 0.00000 | ი იიიიი |
| 1 thi | | ∆000° | 1.97905 |
| browe and | | | 1.69214 |
| Propane I-Butane | 0.00000 | | 1.05957 0.94018 |
| 11/2 rgUe | 0.00172 0.00178 1.21768 | | 0.94010 |
| 1'V cdUb | - 67500 | 42886·2402 42720·32422 42720·3906 | |
| 1 (1/2 1/2 D c | 39923.62500 39923.62507 35288.10547 35288.10000 | 850 ⁰ | |
| Pentane Otal lb/h | 35288.103 75213.0000 | 178.56 Psia | Y/X |
| 1 10/1 | 75213. | 178.30/h Liq 1b/h | 0.00000 |
| , dge ₦ | 266.73 F | Ti9 20000 | 0.0000 |
| 1 Orl | 266 /h Vap 1b/h | 0,000 | 1.98312 |
| topane | vap 16/11 0.00000 | | 1.69589 |
| 1 ("hqDe | 0.00000 | | 1.06242 0.94284 |
| 1 1/2 Cdn | 0.00095 0.00095 | 0.52550 40887.06250 40807.00000 | 0.9420 |
| | 0.00034 0.77948 0.7500 | 40887.00000 44659.00000 4546.5859 | |
| | 38164.3797 | a 5 5 7 | |
| Pentane Potal lb/h | 38164.37797 36993.41797 75158.5703 | ca psia | Y/X |
| 1b/h | 75158. | 178.62 F 1b/h Liq 0.00000 | 0.00000 |
| \$\\ \frac{\stage}{\text{stage} \pm 44} | 10 F | Tig _00000 | ~ nnuuu |
| En. # 44 | 26 ⁷ ·19 F | ° ~~000 | 1 987/9 |
| Ethane | Vap -0000 | | 1.70018 |
| 1692 | 0. 2000 | | 1.06565 0.94586 |
| | 0.00051 0.00051 | | 0.94500 |
| 1 outane | $0.00038 \\ 0.48358 \\ 0.18531$ | 4721.87025 5726.90625 5726.8164 | |
| Butane I pentane I pentane Total lb (| $0.48531 \\ 36165.19531 \\ 36165.08984 \\ 232.08534$ | 5726.90020 10448.8164 | |
| orantane | 36165·19984 38932·08984 | * | |
| Pentane Potal 1b/h | 38932·0896 75097·7734 | | |
| 7.11 | 13- | | |

CHEMCAD 6.0.1

Mame: LPG COLUMN | Late: 0: 27-2009 Time: 22:31:24

| nit tyr | Pe : TOWF | | | Eqp # 1 | | Thermal | Surface |
|----------------------------------|----------------|----------------|-------------------------------|-------------------------|--------------------------|---|---------------------|
| Lion | C: TOWF | e Unit | nam d : | | | N | tension |
| rīQuid | | | Actual | Actual density | viscosity | Btu/hr-ft. | -F dyne/cm 5.502 |
| Stg | I | Average | vol rate | 1b/ft3 | cP 0.0957 | 0.049 | 5.391 |
| i | lb/h | mol wt | ft3/hr | 30.25 | a 0958 | 0.048 0.047 | 5.247 |
| 2 | 79650 | 50.18 | 2633.22 | 30.41 | a 0957 | 0.047 | 5.171 |
| 3 | 82166 | 53.41 | 2702.20 2747.20 | 30.42 | 0 0950 | 0.047 | 5.126 |
| 4 | 83567 83843 | 55.20 | 2754.63 | 30.44 | 0 0950 | 0.046 | 5.083 5.030 |
| 5 6 | 83479 | 56.34 | 2739.84 | 30.47 30.50 | 0.0962 | 0.046 | 4.837 |
| 7 | 82868 | 57.27 58.21 | 2716.83 | 30.52 | 0 0301 | 0.045 0.045 | 4.751 |
| 8 | 67659 | 59.18 | 2216.51 | 30.42 | a 0950 | 0.043 | 4.650 |
| 9 | 70358 | 60.72 | 2313.03 | 30.38 | ~ U321 | 0.044 | 4.660 4.746 |
| 10 | 72185 | 61.81 | 2375.91 | 30.34 | a 1960 | 0.044 | 4.753 |
| lı la | 73357 | 62.71 | 2418.22 | 30.38 | 0.0965 | 0.044 | 4.721 |
| 12 13 | 74207 | 63.60 | 2442.45 | 30.48 30.52 | ~ UAR2 | $\begin{array}{c} 0.044 \\ 0.044 \end{array}$ | 4.674 |
| 13 14 | 75013 75054 | 64.58 | 2188.32 | 20.53 | a 0999 | 0.044 | 4.626 |
| 15 | 75954 77103 | 65.66 | 2525.33 | 30.54 | $^{\circ}$ 1015 | 0.044 | 4.586 4.558 |
| 16 | 78421 | 66.79 67.89 | 2569.20 | 20.51 | a 1030 | 0.044 | 4.539 |
| 17 | 79805 | 68.89 | 2616.04 | 20.49 | 0.1044 0.1056 | 0.044 | 4.527 |
| 18 | 81127 | 69.73 | 2660.70 | $30 \cdot 48$ | Λ 106 ⁵ | 0.044 | 4.520 4.515 |
| 50 1 <i>0</i> | 82290 | 70.39 | 2700.00 2732.12 | 30.47 30.46 | 0 10/4 | 0.045 | 4.512 |
| 51 | 83247 | 70.90 | 2757.01 | 20.40 | 0 10// | 0.045 | 4.509 |
| 55 | 83993 | 71.27 | 2775.13 | 20.40 | ~ 108 [⊥] | 0.045 | 4.507 |
| 53 | 84554 84963 | 71.54 | 2200.37 | 20.40 | 0.1084 0.1086 0.1086 | 0.045 0.045 | 4.505 4.504 |
| 24 | 85257 | 71.73 71.86 | ~~04.4- | 20 40 | Δ 108' | 0.045 | 4.502 |
| ζ <u>ς</u> | 85462 | 71.80 | | 30.45 30.45 30.45 | 2 1 () B O | 0.045 | 4.501 |
| २ _६ २७ | 85606 | 72.02 | 2810.99 2814.45 2814.01 | 20 43 | A 1002 | 0.045 | 4.499 |
| 58 | 85706 | 72.06 | | 20 42 | 2 1 U 0 2 | 0.045 | 4.498 |
| Śθ | 85778 | 72.09 | | 20 42 | 0.1089 | 0.045 | 4.495 |
| 30 | 85829 85862 | 72.11 | | 20.4^{3} | 2 1000 | 0 045 | 1 494 |
| 31 | 85875 | 72.12 72.13 | | 30·44 30·44 30·44 | - 11190 | 0 045 | n 493 |
| 32 25 | 85881 | 72.13 | | ~ ~ 4 ~ | ~ 1112V | 0.045 | 1 49Z |
| 32 33 34 35 36 38 | 85879 | 72.14 | 2821.46 | 20 42 | 0.1090 0.1090 0.1090 | 0.045 | 4.489 |
| 35 | 85872 | 72.14 | | -0 47 | ~ 1 <i>U</i> 2× | 0.045 | α ΔΒ΄ |
| 36 | 85860 85860 | 72.15 | | - 1 4 7 | ∧ 1U2 [±] | 0.045 | Λ ΔΒΟ |
| 38 | 85848 85832 | 72.15 | | 30.42 30.41 30.41 | A 1025 | 0.045 | 4 400 |
| 39 | 85782 | 72.15 | | - 0 4 - | A 1026 | 0.045 | 1 4b2 |
| 90 91 | 85748 | 72.15 72.15 | | A 40 | 0.1092 0.1092 0.1092 | $0.044 \\ 0.044$ | 4.484 |
| 42 | 8570a | 72 15 | 2818.67 | | 0.1093 | 0.03. | |
| 44 43 41 | 85662 | 72.15 | | . ^ _ / | 0.20 | | |
| 44 | 85607 | 72.15 | | 30.38 | | | |
| S. | 85547 10449 | 72.15 | 2813.93 | | | | |
| $e^{j\mathcal{E}}$ | -0449 | 72.15 | | | | | |
| l | | Liq H | | | | | |
| Ś | | MMBtu. | /h | | | | |
| 3 | | 00 85 | | | | | |

T, C,. D.T. 149

49

-88.85 -89.13 -89.213

60.35

61.38

65⁹⁰⁸

-59.4 -60.153 -60.785

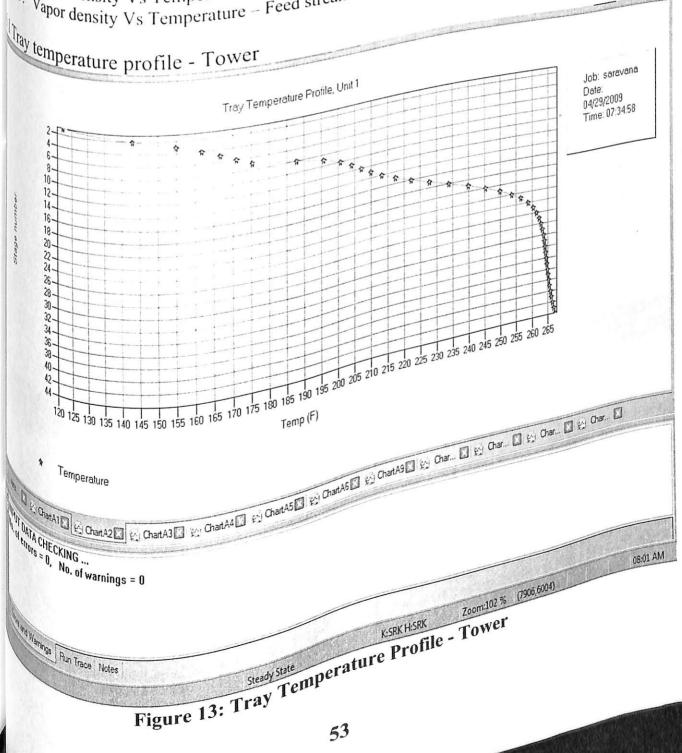
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Page 4
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 21
 ŹĮ
           -61.285
 23
           -61.663
 24
            -61.94
 25
           -62.149
 26
           -62.275
 27
           -62.369
 58
           -62.431
 29
           -62.473
 30
           -62.499
 31
           -62.511
 32
           -62.505
 33
           -62.492
 34
           -62.471
 35
           -62.442
 36
           -62.406
 37
           -62.367
 38
           -62.319
 39
           -62.262
 10
           -62.193
 41
           -62.113
 ٩Į
           -62.019
43
           -61.912
 44
           -61.788
           -61.649
                                                                Page 1
Column Diagnosis Report
CHEMCAD 6.0.1
                                     Time: 06:30:07
, Mame: LPG COLUMN
                   Date: 04/28/2009
has convereged
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6. ANALYSIS ON CHARACTERISTICS OF TOWER AND FEED STREAM - CHEMICAD 6.0.1

Following charts are generated by CHEMCAD 6.0.1 which are explain about characteristics of well as stream in varies conditions.

- 1. Tray temperature profile Tower
- Pseudo component curve Tower
- 3. Heat curve Tower
- Phase envelope Feed stream
- Liquid viscosity Vs Temperature Feed stream

 Van
- Napor viscosity Vs Temperature Feed stream
 Napor viscosity Vs Temperature Feed stream
 Feed stream
- Napor viscosity Vs Temperature Feed stream
 Napor pressure Vs Temperature Feed stream
- Liquid density Vs Temperature Feed stream
 Vans density Vs Temperature Feed stream
 Feed stream
- 9. Vapor density Vs Temperature Feed stream



6.2 Pseudocomponent Curve - Tower

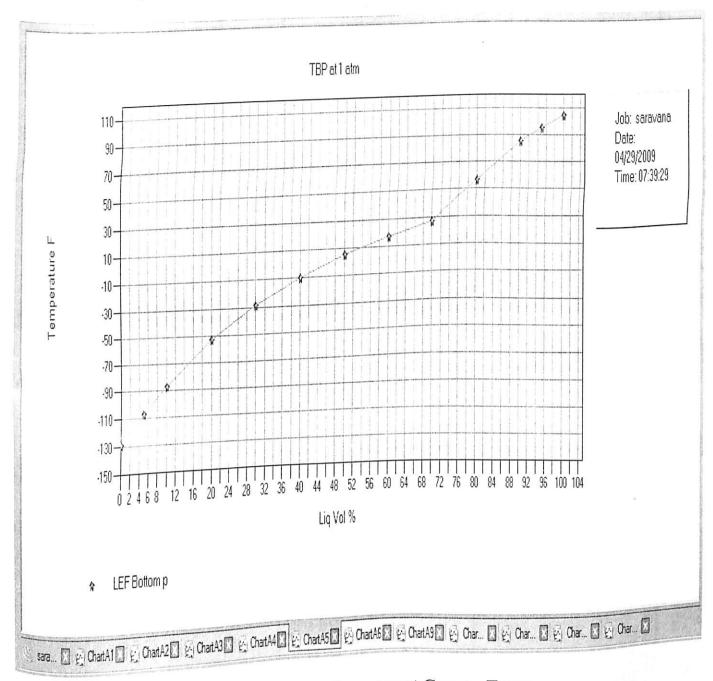


Figure 14: Pseudo Component Curve - Tower

6.3 Heat Curve - Tower

CHEMCAD 6.0.1

Job Name: LPG COLUMN Date: 04/29/2009 Time: 16:09:09

Eqp # 1 Unit type : TOWR Unit name:

Condenser

| NP 1 2 3 4 5 6 7 8 9 10 | Temp F 142.8 140.6 138.5 136.4 134.3 132.3 130.2 128.1 126.1 | Pres psia 176.4 176.4 176.4 176.4 176.4 176.4 176.4 | Del H MMBtu/h 19.1 17.2 15.3 13.4 11.5 9.55 7.64 5.73 3.82 1.91 0.000 | Vapor 1b/h 135286 121386 107556 93808 80143 66562 53063 39647 26319 13091 | Liquid 1b/h 0 13900 27730 41478 55143 68725 82223 95639 108968 122195 135285 | Vap mole frac. 1.0000 0.9029 0.8051 0.7066 0.6075 0.5078 0.4074 0.3064 0.2047 0.1025 0.0000 | Vap mass frac. 1.0000 Dew 0.8973 0.7950 0.6934 0.5924 0.4920 0.3922 0.2931 0.1945 0.0968 0.0000 Bub |
|-------------------------|--|---|---|--|--|---|---|
| 10 11 | 124.0 | 176.4 | 0.000 | 1 | 135285 | 0.0000 | 0.0000 Bub |

Page 1

Reboiler

| NP 1 2 3 4 5 6 7 8 9 10 11 | Temp F 266.7 266.8 266.9 266.9 267.0 267.1 267.1 267.1 | Pres psia 178.6 178.6 178.6 178.6 178.6 178.6 178.6 178.6 178.6 | Del H MMBtu/h 0.000 0.808 1.62 2.43 3.23 4.04 4.85 5.66 6.47 7.28 8.08 | Vapor 1b/h 0 7603 15204 22805 30404 38002 45600 53194 60788 68381 75972 | Liquid 1b/h 86422 78818 71217 63617 56018 48420 40822 33227 25634 18041 10449 | Vap mole frac. 0.0000 0.0880 0.1759 0.2639 0.3518 0.4397 0.5276 0.6155 0.7034 0.7912 0.8791 | Vap mass frac. 0.0000 Bub 0.0880 0.1759 0.2639 0.3518 0.4397 0.5276 0.6155 0.7034 0.7912 0.8791 | |
|----------------------------|---|---|--|---|---|---|---|--|
|----------------------------|---|---|--|---|---|---|---|--|

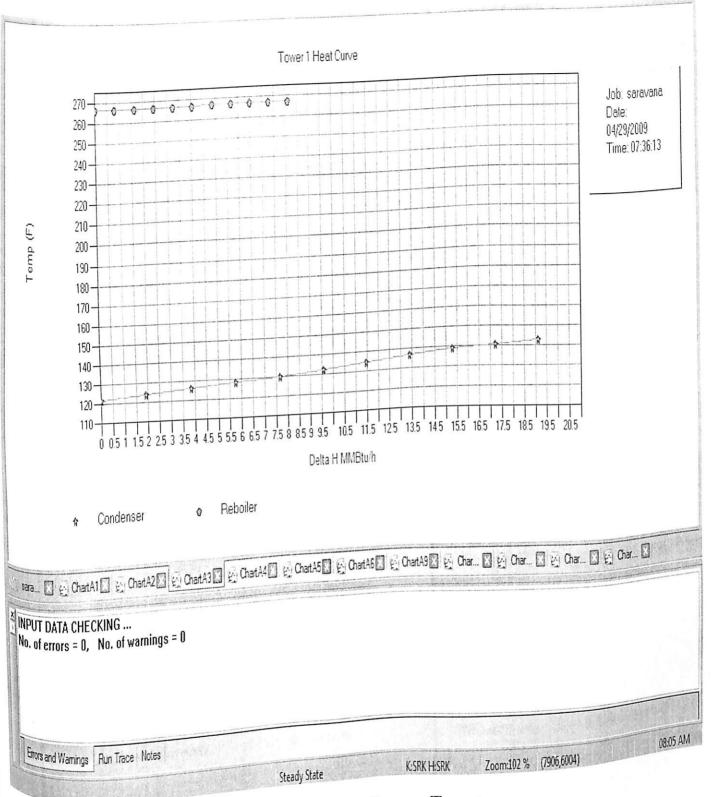


Figure 15: Heat Curve - Tower

6.4 Phase Envelope – Feed Stream

Page 1
CHEMCAD 6.0.1

Job Name: LPG COLUMN Date: 04/29/2009 Time: 08:17:35

Phase Envelope for Stream 1

| | | | | 7 | Zl |
|-----|----------|---------|---------|----------|-------|
| | | Press | Vfrac | Zv | 21 1 |
| No. | Temp | | | | |
| | F | psia | 0.00000 | 0.911 | 0.019 |
| 1 | 56.866 | 64.339 | | 0.900 | 0.022 |
| 2 | 66.866 | 74.866 | 0.00000 | 0.889 | 0.025 |
| | | 86.602 | 0.00000 | | |
| 3 | 76.866 | 99.622 | 0.00000 | 0.876 | 0.029 |
| 4 | 86.866 | | 0.00000 | 0.863 | 0.033 |
| 5 | 96.866 | 113.999 | 0.00000 | 0.849 | 0.038 |
| 6 | 106.866 | 129.807 | 0.00000 | 0.835 | 0.043 |
| | 116.866 | 147.116 | 0.00000 | 0.819 | 0.049 |
| 7 | | 165.994 | 0.00000 | | 0.055 |
| 8 | 126.866 | 186.509 | 0.00000 | 0.802 | |
| 9 | 136.866 | 186.509 | 0.00000 | 0.785 | 0.062 |
| 10 | 146.866 | 208.721 | 0.00000 | 0.766 | 0.070 |
| 11 | 156.866 | 232.687 | 0.00000 | 0.746 | 0.078 |
| 12 | 166.866 | 258.458 | 0.00000 | 0.725 | 0.088 |
| | 176.866 | 286.075 | 0.00000 | 0.702 | 0.099 |
| 13 | 186.866 | 315.570 | 0.00000 | 0.702 | 0.111 |
| 14 | | 346.958 | 0.00000 | 0.678 | |
| 15 | 196.866 | 380.233 | 0.00000 | 0.652 | 0.124 |
| 16 | 206.866 | 300.233 | 0.00000 | 0.625 | 0.140 |
| 17 | 216.866 | 415.354 | 0.00000 | 0.594 | 0.158 |
| 18 | 226.866 | 452.228 | | 0.578 | 0.168 |
| | 231.866 | 471.271 | 0.00000 | 0.569 | 0.173 |
| 19 | 231.000 | 480.930 | 0.00000 | | |
| 20 | 234.366 | 485.791 | 0.00000 | 0.565 | 0.176 |
| 21 | 235.616 | 489.694 | 0.00000 | 0.561 | 0.178 |
| 22 | 236.616 | 489.034 | 0.00000 | 0.557 | 0.181 |
| | 237.616 | 493.610 | 0.00000 | 0.554 | 0.183 |
| 23 | 237.61 | 497.538 | | 0.550 | 0.186 |
| 24 | 238.616 | 501.476 | 0.00000 | 0.546 | 0.188 |
| 25 | 239.616 | 505.426 | 0.00000 | | |
| 26 | 240.616 | 505.420 | 0.00000 | 0.542 | 0.191 |
| 27 | 241.616 | 509.386 | 0.00000 | 0.539 | 0.193 |
| | 242.616 | 513.356 | 0.00000 | 0.535 | 0.196 |
| 28 | 242.010 | 517.334 | 0.00000 | 0.531 | 0.199 |
| 29 | 243.616 | 521.320 | 0.00000 | 0.527 | 0.202 |
| 30 | 244.616 | 525.313 | 0.00000 | | |
| 31 | 245.616 | 525.313 | 0.00000 | 0.523 | 0.204 |
| 32 | 246.616 | 529.313 | 0.00000 | 0.518 | 0.207 |
| 33 | 247.616 | 533.317 | 0.00000 | 0.514 | 0.210 |
| | 247.010 | 537.325 | 0.00000 | 0.510 | 0.213 |
| 34 | 248.616 | 541.335 | 0.00000 | 0.505 | 0.217 |
| 35 | 249.616 | 545.347 | 0.0000 | 0.501 | 0.220 |
| 36 | 250.616 | 545.54 | 0.00000 | | 0.220 |
| 37 | 251.616 | 549.358 | 0.00000 | 0.496 | 0.223 |
| 38 | 251.616 | 553.366 | 0.00000 | 0.491 | 0.227 |
| | 252.616 | 557.369 | 0.00000 | 0.487 | 0.231 |
| 39 | 253.616 | 561.367 | 0.00000 | 0.482 | 0.234 |
| 40 | 254.616 | 501.5 | 0.00000 | 0.476 | 0.238 |
| 41 | 255.616 | 565.355 | 0.00000 | | 0.242 |
| 42 | 256.616 | 569.330 | 0.00000 | 0.471 | 0.242 |
| 43 | | 573.290 | 0.0000 | 0.460 | 0.251 |
| | 257.616 | 581.146 | 0.00000 | 0.454 | 0.256 |
| 45 | 259.616 | 201.21 | 0.00000 | 0.448 | 0.260 |
| 46 | 260.616 | 585.031 | a aaau | | 0.266 |
| 47 | 261.616 | 588.881 | a 00000 | 0.442 | 0.271 |
| 48 | | 592.685 | 0.00000 | 0.436 | 0.277 |
| 49 | 262.616 | 596.432 | 0.0000 | 0.429 | 0.27 |
| 50 | 263.616 | 600.111 | 0.00000 | 0.421 | 0.283 |
| | 264.616 | 600.111 | 0.00000 | J | |
| 51 | 265.616 | 603.702 | = | | |
| | = 00,020 | | | | |

CHEMCAD 6.0.1

| Job Name: | : LPG COLUM | M Date: | 04/29/2009 | Time: 08:17:35 | |
|-----------|--------------------|--------------------|--------------------|----------------|----------------|
| 52 | 266 616 | 607.183 | 0.00000 | 0.414 | 0.290 |
| 53 | 266.616 267.616 | 610.524 | 0.00000 | 0.405 | 0.297 |
| 54 | 268.616 | 613.666 | 0.00000 | 0.397 | 0.305 |
| 55 | 269.616 | 616.524 | 0.00000 | 0.386 | 0.315 |
| 56 | 270.616 | 618.916 | 0.00000 | 0.378 | 0.326 |
| 57 | 271.616 | 614.160 | 0.00000 | 0.382 | 0.366 |
| 1 | 98.617 | 62.771 | 1.00000 | 0.912 | 0.020 |
| 2 | 121.585 | 91.019 | 1.00000 | 0.884 | 0.028 |
| 3 | 139.635 | 119.266 | 1.00000 | 0.858 | 0.037 |
| 4 | 154.685 | 147.513 | 1.00000 | 0.833 | 0.045 |
| 5 | 167.685 | 175.760 | 1.00000 | 0.809 | 0.054 |
| 6 | 179.179 | 204.007 | 1.00000 | 0.786 | 0.063 |
| 7 | 189.513 | 232.254 | 1.00000 | 0.764 | 0.073 |
| 8 | 198.919 | 260.502 | 1.00000 | 0.742 | 0.082 |
| 9 | 207.560 | 288.749 | 1.00000 | 0.720 | 0.092 |
| 10 | 215.557 | 316.996 | 1.00000 | 0.698 | 0.103 |
| 11 | 222.999 | 345.243 | 1.00000 | 0.676 | 0.114 |
| 12 | 229.956 | 373.490 | 1.00000 | 0.654 | 0.126 |
| 13 | 236.479 | 401.737 | 1.00000 | 0.632 | 0.138 |
| 14 | 242.609 | 429.985 | 1.00000 | 0.608 0.584 | 0.151 |
| 15 | 248.373 | 458.232 | 1.00000 | 0.559 | 0.166 |
| 16 | 253.770 | 486.479 | 1.00000 | 0.532 | 0.182 |
| 17 | 258.859 | 514.726 | 1.00000 1.00000 | 0.503 | 0.200 |
| 18 | 263.564 | 542.973 557.097 | 1.00000 | 0.486 | 0.220 0.232 |
| 19 | 265.762 | 564.159 | 1.00000 | 0.478 | 0.232 |
| 20 | 266.817 | 567.689 | 1.00000 | 0.473 | 0.242 |
| 21 | 267.329 | 570.689 | 1.00000 | 0.469 | 0.242 |
| 22 | 267.760 | | 1.00000 | 0.465 | 0.243 |
| 23 | 268.182 | 573.689 | 1.00000 | 0.461 | 0.252 |
| 24 | 268.598 | 576.689 | 1.00000 | 0.457 | |
| 25 | 269.006 | 579.689 | 1.00000 | 0.452 | 0.255 |
| 26 | 269.405 | 582.689 | 1.00000 | 0.448 | 0.258 |
| 27 | 269.795 | 585.689 | 1.00000 | 0.443 | 0.262 |
| 28 | 270.174 | 588.689 | 1.00000 | 0.438 | 0.266 |
| 29 | 270.542 | 591.689 | 1.00000 | 0.433 | 0.270 |
| 30 | 270.896 | 594.689 | 1.00000 | 0.428 | 0.274 |
| 31 | 271.235 | 597.689 | 1.00000 | 0.422 | 0.278 |
| 32 | 271.555 | 600.689 | 1.00000 | 0.416 | 0.283 |
| 33 | 271.852 | 603.689 | 1.00000 | 0.409 | 0.288 0.293 |
| 34 | 272.119 | 606.689 | 1.00000 | 0.401 | |
| 35 36 | 272.348 | 609.689 | 1.00000 | 0.393 | 0.299 |
| 36 | 272.526 | 612.689 | 1.00000 | 0.384 | 0.306 0.320 |
| 37 30 | 272.729 | 615.689 | 1.00000 | 0.359 | 0.320 |
| 38 | 272.116 | 618.689 | 1.00000 | | 0.010 |
| | | | | | |

271.116 F 3c = 616.538 psia 0.344

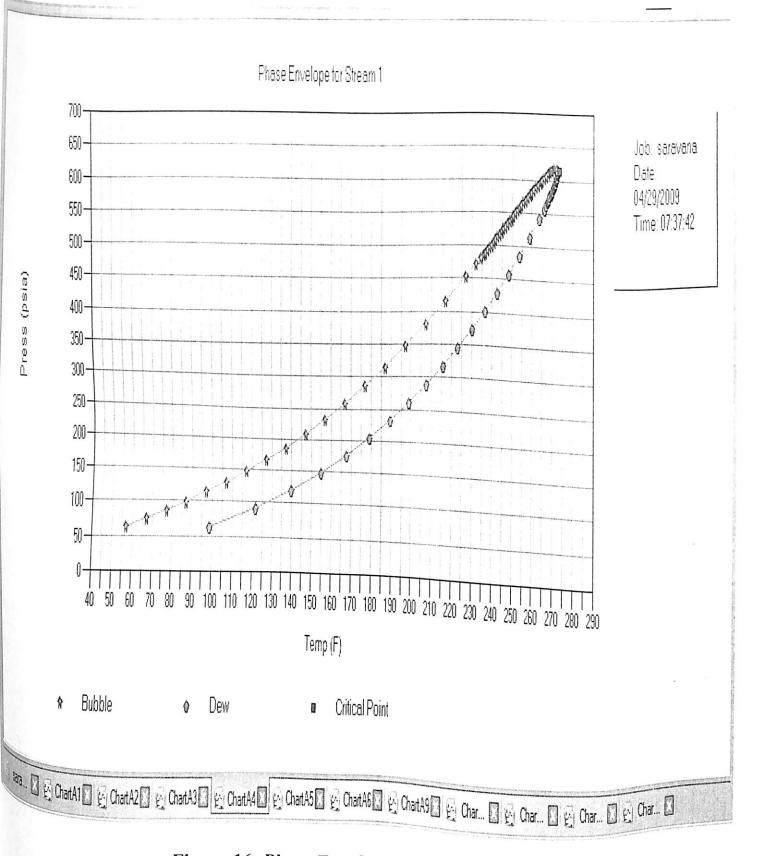


Figure 16: Phase Envelope – Feed Stream

6.5 Liquid Viscosity Vs Temperature – Feed Stream

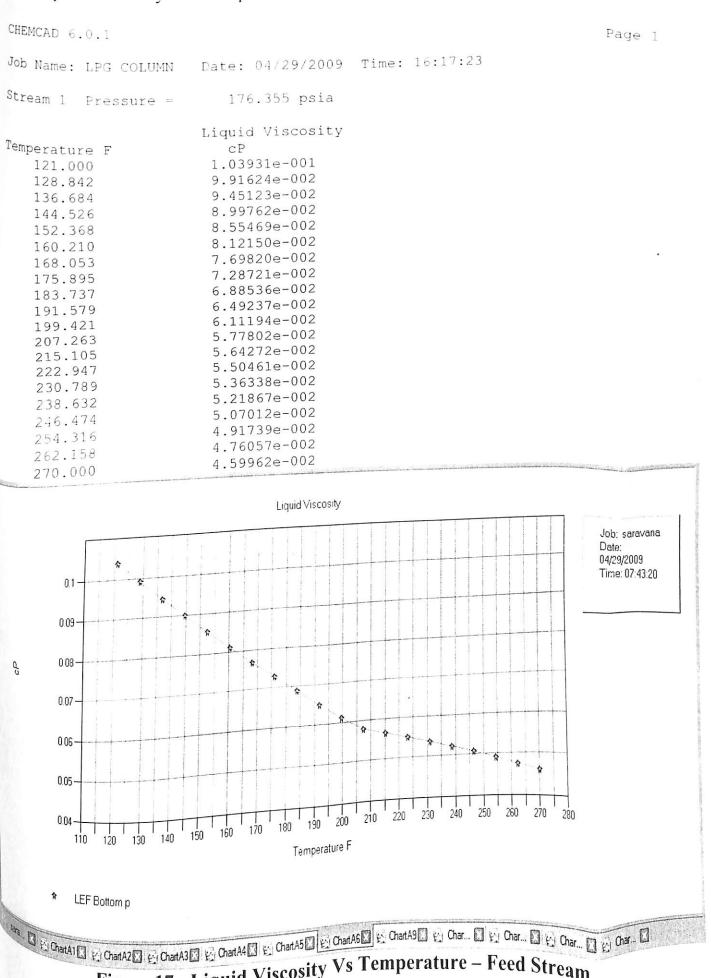
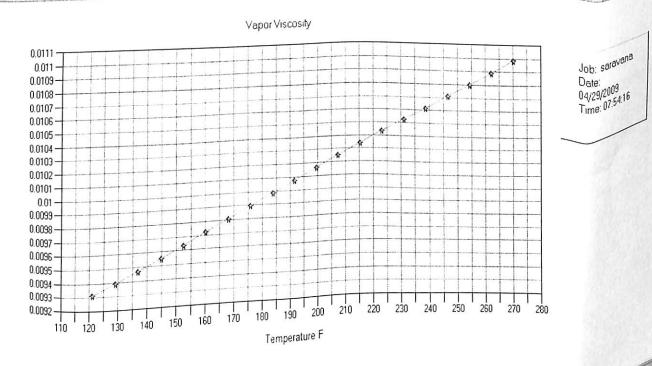


Figure 17: Liquid Viscosity Vs Temperature – Feed Stream

6.6 Vapor Viscosity Vs Temperature – Feed Stream

Page 1 CHEMCAD 6.0.1 Time: 16:18:23 Job Name: LPG COLUMN Date: 04/29/2009 Stream 1 176.355 psia Pressure = Vapor Viscosity Temperature F 9.31369e-003 121.000 9.39012e-003 128.842 9.46959e-003 136.684 9.55145e-003 144.526 9.63523e-003 152.368 9.72057e-003 160.210 9.80718e-003 168.053 9.89484e-003 175.895 9.98336e-003 183.737 1.00726e-002 191.579 1.01624e-002 199.421 1.02527e-002 207.263 1.03434e-002 215.105 1.04344e-002 222.947 1.05256e-002 230.789 1.06171e-002 238.632 1.07087e-002 246.474 1.08003e-002 254.316 1.08921e-002 262.158 1.09839e-002 270.000



BachartAl 函ChartA2 函ChartA3 函ChartA4 函ChartA5 函ChartA5 函ChartA5 函ChartA5 函ChartA5 函ChartA5 函数ChartA5 函数C Figure 18: Vapor Viscosity Vs Temperature – Feed Stream

6.7 Vapor Pressure Vs Temperature – Feed Stream

CHEMCAD 6.0.1 Page 1 Job Name: LPG COLUMN Date: 04/29/2009 Time: 08:20:38 Stream 1 Pressure = 176.355 psia Vapor Pressure Temperature F psia 121.000 6.79130e+005 128.842 5.88018e+007 136.684 5.09130e+009 144.526 4.40825e+011 152.368 1.50150e+005 160.210 1.30006e+007 168.053 1.12564e+009 175.895 9.74629e+010 183.737 8.43873e+012 Warning: K values = 1. 1.00000e+005 191.579 Warning: K values = 1. 199.421 1.00000e+005 Warning: K values = 1. 207.263 1.00000e+005 215.105 8.65841e+006 Warning: K values = 1. 222.947 1.00000e+005 Warning: K values = 1. 1.00000e+005 230.789 Warning: K values = 1. 1.00000e+005 238.632 Warning: K values = 1. 1.00000e+005 246.474 8.65841e+006 254.316 7.49680e+008 262.158 1.00100e+005 270.000 Vapor Pressure 9e12 8.5e12 Job: saravana Date: 04/29/2009 Time: 07:48:12 8e12 -7.5e12-6.5e12-6e12-5.5e12-5e12-4.5e12-3.5e12-3e12-2.5e12-2e12-1.5e12 1~ 1 1 210 220 230 240 250 260 270 5e11-190 160 150 Temperature F LEF Bottom p BE ChartA2回 ChartA3回 ChartA4回 ChartA5回 E ChartA6回 ChartA6U ChartA

Figure 19: Vapor Pressure Vs Temperature – Feed Stream

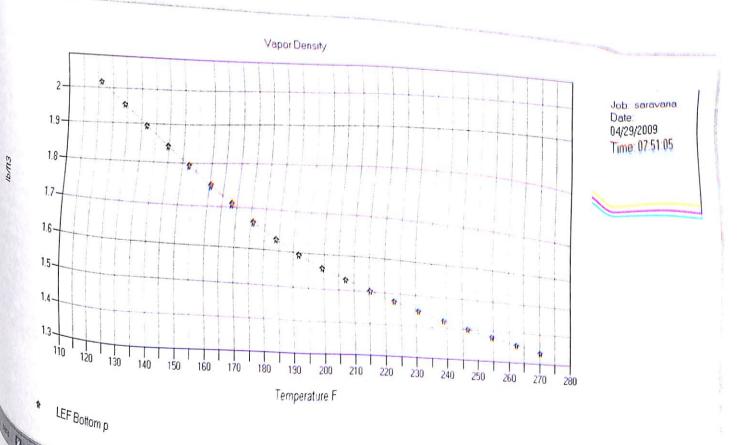
6.8 Liquid Density Vs Temperature – Feed Stream

CHEMCAD 6.0.1 Page -Job Name: LPG COLUMN Date: 04/29/2009 Time: 16:20:25 Stream 1 Pressure = 176.355 psia Liquid Density Temperature F lb/ft3 121.000 3.12027e+001 128.842 3.07572e+001 136.684 3.02916e+001 144.526 2.98021e+001 152.368 2.92837e+001 160.210 2.87293e+001 168.053 2.81285e+001 175.895 2.74641e+001 183.737 2.67057e+001 191.579 2.57880e+001 199.421 2.45099e+001 207.263 2.01451e+001 215.105 2.00079e+001 222.947 1.98598e+001 230.789 1.96983e+001 238.632 1.95198e+001 246.474 1.93186e+001 254.316 1.90845e+001 1.87956e+001 262.158 1.83774e+001 270.000 Liquid Density 31.5 30.5 Job: saravana 29.5 Date: 04/29/2009 28.5 Time: 07:51:35 27.5 26.5 1b/ft3 25.5 24.5 23.5 22.5 21.5 20.5-19.5 18.5 220 17.5-200 180 170 150 140 Temperature F LEF Bottom p

Figure 20: Liquid Density Vs Temperature — Feed Stream

6.9 Vapor Density Vs Temperature – Feed Stream

```
CHEMCAD 6.0.1
Job Name: LPG COLUMN
                     Date: 04/29/2009 Time: 05:21:55
$tream 1 Pressure =
                         176.355 psia
Temperature F
                      Vapor Density
                         lb/ft3
  121.000
                       2.02028e+000
  128.842
                       1.95529e+000
  136.684
                       1.89665e+000
  144.526
                       1.84325e+000
  152.368
                       1.79424e+000
   160.210
                       1.74898e+000
   168.053
                       1.70695e+000
   175.895
                       1.66774e+000
   183.737
                        1.63102e+000
   191.579
                       1.59649e+000
   199.421
                        1.56394e+000
   207.263
                        1.53316e+000
   215.105
                        1.50398e+000
   222.947
                        1.47625e+000
   230.789
                        1.44985e+000
   238.632
                       1.42467e+000
   246.474
                       1.40060e+000
   254.316
                       1.37757e+000
  262.158
                       1.35549e+000
  270.000
                       1.33430e+000
```



Time :

7. LPG COLUMN SIZING

Column Sizing Sieve Tray

Page 1

Name: LPG COLUMN: Late: 04/28/2009 Time: 19:51:51

por load is defined as the vapor from the tray below.

Quid load is defined as the liquid on the tray.

lood correlation: Fair

| 1 | |
|---|--|
| Equip. 1 Tray No. 1 Vapor 104119.000 1b/h 104119.000 ft3/hr 61186.832 ft3/hr 61186.832 1b/ft3 1.702 1b/ft3 | Liquid 49147.703 lb/h 1622.310 ft3/hr 30.295 lb/ft3 6.000 2.000 |
| | |
| Tower internal diameter, ft | 1 |
| Transit Internal diameter, Internal diameter, | Area ft2 |
| Tray spacing, ft No. of tray liquid passes Width ft 4.616 | 3.476 |
| The of travelianid passes width 10 4.610 | 4.616 |
| Tray spacing, ft Length ft bowncomer dimension 1.083 | 0.167 |
| | 3.833 |
| ®: •Nr • • • • • • • • • • • • • • • • • • | 5.562 |
| Avg. Side | |
| | 28.274 |
| Flow Path length ft | 21.322 |
| | 68.868 |
| | 0.015 |
| 1 LCZ | 0.646 |
| flood are gred res | 102956.078 |
| ≥ | 0.230 |
| Aerational entrainment Minimum factor Tray (Weeping) vapor flow | 0.048 |
| Min. Con factor 1b/1 | 0.146 |
| Training (Weening) vapor II | 0.140 |
| Tra Press loca ft | 0.516 |
| | 3.982 |
| | 15.428 |
| h come clearance | 0.130 |
| Downcomer clearance ft Downcomer backup ft Downcomer residence time sec sec | 5.432 |
| Downcomer clearance ft Downcomer backup ft Downcomer residence time Downcomer apparent residence Light of the sec | 164.570 |
| Nowncomer apparent residence liquid holdup ft3 | 176.35 ⁵ |
| Liquid holdup ft3 Design holdup 1b | 0.850 |
| giand hold | 13700.000 |
| heart hold | 0.003 |
| Julan Dra | 0.003 |
| Design pressure psia Allowable stress psia | 0.047 |
| Cowah Cliciency | 0.049 |
| Allowable stress psia Column thickness ft | |
| bolumnian allowance ft | |
| Column thickness ft tom thickness ft | |
| thickness ft | |
| Allego 2- | |

1.2 Costing Sieve Tray

CHEMCAD 6.0.1

Date: 04/28/2009 Time: 20:00:45 Job Name: LPG COLUMN

Page 1

Preliminary Distillation Towers Cost Estimation

Distillation Tower Cost for Equip. 1 Column material = Carbon steel Sieve tray Tray material = Carpon steel Base cost index = 347.5Current cost index = 610.4

| Calculated cost: | = \$ | 118284 |
|---|------|--------|
| Shell cost | = \$ | 45320 |
| Tray cost | = \$ | 24500 |
| Platform & ladders | - ÷ | 330414 |
| Column cost (nurchase) | = \$ | 991241 |
| Column (installed) | = \$ | 0 |
| Con 1 (MI) ECITOR - 1 | • | 0 |
| condenser cost (installed) | = \$ | 0 |
| Condenser cost (installed) Condenser cost (installed) | = \$ | 0 |
| | = \$ | 330414 |
| Rebailer cost (111500000 | = \$ | 991241 |
| Total (niirchass) | = \$ | 9710 |
| Total cost (installed) | | |
| | | |

66

7.3 Column Sizing – Value Tray

Page 1
CHEMCAD 6.0.1

Job Name: LPG COLUMN Date: 04/28/2009 Time: 19:55:01

Vapor load is defined as the vapor from the tray below. Liquid load is defined as the liquid on the tray.

Section: 1 Flood correlation: Glistch

| " Colletation. Gra- | | |
|--|---|--------------------------|
| Equip. 1 | Tray No. 2 | Liquid 50185.875 lb/h |
| _ | Vapor | 50185.675 16711 |
| Tray Loadings | - 4 A 1 I D / 11 | 1644.608 ft3/hr |
| 1 3 2 2 2 2 2 2 2 2 2 2 2 2 2 2 2 2 2 2 | 105157.141 123/hr 60794.498 ft3/hr 730 lb/ft3 | 30.515 lb/ft3 |
| | 60794.498 160/ft3 1.730 lb/ft3 | 1.000 |
| Densi | 1.75 | |
| Density | 1.730 127 | |
| System factor | | 12.000 |
| ATAG TADO . ALT | | _ |
| Valve material : S.S. | | 14.000 |
| Valve thickness gauge | | 5.000 |
| Deck | ft | 2.000 |
| Deck thickness gauge | ft | 1 |
| Tower internal diameter, | Length ft | Area ft2 |
| Itay spacing ft | Length It | 1.922 |
| of travelianta pass- | width 10 3.611 | 3.611 |
| Downcomer dimension | <i>U</i> • • • • • • • • • • • • • • • • • • • | 0.167 |
| side | 0.771 | 3.458 |
| Avg. weir length ft | | 4.566 |
| Weir bei | | 19.635 |
| Weir height, ft | | 15.792 |
| TOW Dath lawath LU | | |
| TOW Dath alth Tt. | | 70.723 |
| -4V 3ros £F3 | | 3.000 |
| | | 235 |
| F 1004 | | 0.386 |
| 4016 3x02 EFO | | 0.082 |
| | | 0.225 |
| Tray press loss, ft | | 0.146 |
| Trau press loss, | | 0.041 |
| | | 0.713 |
| Dry press drop, ft | | 3.000 |
| h. "comer clearance | | 8.413 |
| Downcomer head loss ft | | 0.238 |
| Downcomer head loss ft Downcomer backup ft Downcomer residence time Downcomer apparent reside Downcomer velocity ft/se | sec sec | 3.907 |
| Nowncomer residence time | nce time | 3.907 |
| Downcomes reside | | 119.217 |
| Downcomer apparent restriction ft/se | | 176.409 |
| 7400.2 | c | 0.850 |
| | | 13700.000 |
| Desid holdup 1b | | 0.003 |
| Design pressure psia | | 0.039 |
| Joint efficiency | | 0.052 |
| Allowable stress psia | | |
| Corrosion allowance ft | | |
| Column thickness ft | ••• | |
| Bottom thickness ft | | |
| - cutckiless - | | |

7.4 Costing - Value Tray

Page 1
CHEMCAD 6.0.1

Job Name: LPG COLUMN Date: 04/28/2009 Time: 20:01:55

Preliminary Distillation Towers Cost Estimation

Distillation Tower Cost for Equip. 1
Column material = Carbon steel
Valve tray
Tray material = Carbon steel
Base cost index = 347.5
Current cost index = 610.4

7.6 Costing - Bubble Cap Tray

CHEMCAD 6.0.1

Job Name: LPG COLUMN Date: 04/28/2009 Time: 19:59:33

Preliminary Distillation Towers Cost Estimation

Distillation Tower Cost for Equip. 1
Column material = Carbon steel
Bubble cap tray
Tray material = Carbon steel
Base cost index = 347.5
Current cost index = 610.4

| Calculated cost: Shell cost Tray cost Platform & ladders Column cost (purchase) Column cost (installed) Condenser cost (purchase) Condenser cost (purchase) Reboiler cost (purchase) Reboiler cost (installed) Total cost (purchase) Total cost (purchase) | | \$ \$ \$ \$ \$ \$ \$ \$ \$ \$ \$ | 108430 77716 23377 368035 1104104 0 0 0 368035 1104104 |
|--|---|----------------------------------|---|
| Total cost (purchase) Total cost (installed) | = | \$ | 1104103 |

7.7 Packed Column Sizing

3.

CHEMCAD 6.0.1

Page 1

HTUOV

Job Name: LPG COLUMN Date: 04/29/2009 Time: 00:06:06

Billet-Schultes Correlation

Packing Parameters

50.00000 mm

105.00000 VoidFrac. 0.96000

| Packed | Tower | Design | for | Tower | 1 |
|--------|-------|--------|-----|-------|---|
|--------|-------|--------|-----|-------|---|

| racked | Tower Design | an for Tow | ier 1 | , | Diam | HTUOV |
|----------|--------------|------------|--------------|--------------|-------|-------|
| | rough Dept. | 9 | | LiqLoad | ft | ft |
| Stg | DD | %Flood | VapLoad | 1h/(ft2*5ec/ | 4.772 | 1.536 |
| - 3 | PDrop | %F1000 | ib/(ft2*sec) | 0.780 | 4.772 | 1.593 |
| 2 | psi | | 1.634 | 0.779 | 4.791 | 1.628 |
| 2 | 0.011 | 75.00000 | 1.627 | 0.777 | 4.792 | 1.654 |
| 3 | 0.011 | 75.00000 | 1.624 | 0.774 | 4.787 | 1.673 |
| 4 | 0.011 | 75.00000 | 1.623 | 0.772 | 4.783 | 0.747 |
| 5 | 0.011 | 75.00000 | 1.623 | 1.906 | 2.570 | 0.744 |
| 6 | 0.011 | 75.00000 | 1.346 | 1.900 | 2.634 | 0.744 |
| 7 | | 75.00000 | 1.355 | 1.888 | 2.682 | 0.744 |
| 8 | 0.012 | 75.00000 | 1.361 | 1.875 | 2.714 | |
| 9 | 0.012 | 75.00000 | 1.301 | 1.867 | 2.734 | 0.748 |
| 10 | 0.012 | 75.00000 | 1.365 | 1.862 | 2.746 | 0.746 |
| | 0.012 | 75.00000 | 1.367 | 1.859 | 2.751 | 0.761 |
| 11 | 0.012 | 75.00000 | 1.368 | 1.861 | 2.753 | 0.770 |
| 12 | 0.012 | 75 00000 | 1.372 | 1.862 | 2.755 | 0.776 |
| 13 | 0.012 | 75 NOUUU | 1.374 | 1.862 | 2.755 | 0.779 |
| 14 | 0.012 | 75 NOUUU | 1.375 | 1.863 | 2.756 | 0.782 |
| 15 | | 75 00000 | 1,376 | 1.863 | 2.757 | 0.783 |
| 16 | 0.012 | 75.00000 | 1.376 | 1.863 | 2.757 | 0.785 |
| 17 | 0.012 | 75.00000 | 1.377 | 1.864 | 2.757 | 0.786 |
| 18 | 0.012 | 75.000 | 1.377 | 1.864 | 2.757 | 0.787 |
| 19 | 0.012 | 75.00000 | 1.377 | 1.864 | 2.757 | 0.788 |
| 20 | 0.012 | 75.00000 | 1.377 | 1.004 | 2.758 | 0.789 |
| | 0.012 | 75.00000 | 1.378 | 1.864 | 2.758 | 0.789 |
| 51 | 0.012 | 75.00000 | 1.378 | 1.864 | 2.758 | 0.790 |
| 55 | 0.012 | 75 00000 | 1.378 | 1.864 | 2.758 | 0.791 |
| 53 | 0.012 | 75 00000 | 1.378 | 1.865 | 2.759 | 0.791 |
| 24 | 0.012 | 75 00000 | 1.379 | 1.865 | 2.759 | 0.791 |
| 25 | | 75 00000 | 1.379 | 1.865 | 2.759 | 0.792 |
| 56 | 0.012 | 75 NOUUU | 1.379 | 1.865 | 2.759 | 0.792 |
| 27 | 0.012 | 75.00000 | 1.379 | 1.865 | 2.759 | 0.792 |
| 58 | 0.012 | 75.00000 | 1 380 | 1.866 | 2.759 | 0.792 |
| 29 | 0.012 | 75.00000 | 1 380 | 1.866 | 2.759 | 0.792 |
| 30 | 0.012 | 75.000 | 1.380 | 1.867 | 2.760 | 0.792 |
| 31 | 0.012 | 75.00000 | 1.381 | 1.867 | 2.760 | 0.792 |
| 35 21 | 0.012 | 75.00000 | 1.382 | 1 868 | 2.760 | 0.792 |
| | 0.012 | 75 NOUV | 1.383 | 1.869 | 2.760 | 0.792 |
| 33 | 0.012 | 25 UUUUU | 1.303 | 1.871 | 2.761 | 0.792 |
| 34 | 0.012 | VUUV | 1.384 | 1.874 | 2.762 | 0.791 |
| 35 | | 75 NOUVO | 1.386 | 1.877 | 2.762 | 0.790 |
| 36 | 0.012 | ac 00000 | 1.389 | 1.881 | 2.765 | 0.790 |
| 37 | 0.012 | 2 E UUUUU | 1 394 | 1.00- | 2.770 | 0.790 |
| 86 | 0.012 | 25 UUUU | 1 390 | 1.886 | 2.779 | 0.790 |
| 96 | 0.012 | VIIIO | 1 404 | 1.891 | 2.794 | 0.792 |
| 40 | 0.012 | 75.00000 | 1 413 | 1.896 | 2.815 | |
| 41 | 0.012 | 75.00000 | 1 422 | 1.898 | | |
| 42 | 0.012 | 75.0000 | 1.431 | | | |
| 43 | 0.012 | 75.00000 | J * * | | | |
| ٠,3 | 0.012 | 75.00000 | | | | |
| | | | | 41 | | |

Page 2 CHEMCAD 6.0.1 Job Name: LPG COLUMN Date: 04/29/2009 Time: 00:06:07 88.000 Overall: 2.095 Height ft 0.496 HETP ft 1.266 Vapor Load at Loading 1b/(ft2*sec) Pressure drop psi Liquid Load at Flooding lb/(ft2*sec)
Liquid Load at Flooding lb/(ft2*sec) 1.880 1.562 2.321 Liquid Load at Flooding 1b/(ft2*sec) 3.157 Diameter at Loading ft 2.590 178.580 Diameter at Flooding ft 0.850 Design pressure psia 13700.000 Joint efficiency 0.003 Allowable stress psia 0.036 Corrosion allowance ft 0.089 Column thickness ft Bottom thickness ft 7.8 Costing Packed Tower Page 1 Date: 04/29/2009 Time: 00:08:35 CHEMCAD 6.0.1 Job Name: LPG COLUMN Preliminary Distillation Towers Cost Estimation Distillation Tower Cost for Equip. 1 Column material = Carbon steel Packed tower Base cost index = 347.5Current cost index = 610.4 127057 Calculated cost: = \$ 0 = \$ Shell cost 19272 = \$ Packing cost Platform & ladders 257034 = \$ Column cost (purchase) 771103 = \$ Column cost (installed) 0 Condenser cost (purchase) = \$
Condenser cost (installed) = \$ 0 0 Reboiler cost (purchase) 0 **=** \$ Reboiler cost (installed)

257034

771103

= \$

= \$

Total cost (purchase)

Total cost (installed)

8. DISCUSSION ON RESULTS

From calculation which are done by manually as well as simulator seams the packing tower is a best in cases of purchasing, installation, place occupation and also efficiency. Following Figures 23 & 24 are screen shots of Excel 2007, here cost comparison is worked out by using Excel 2007.

8.1 Purchase Cost Comparison

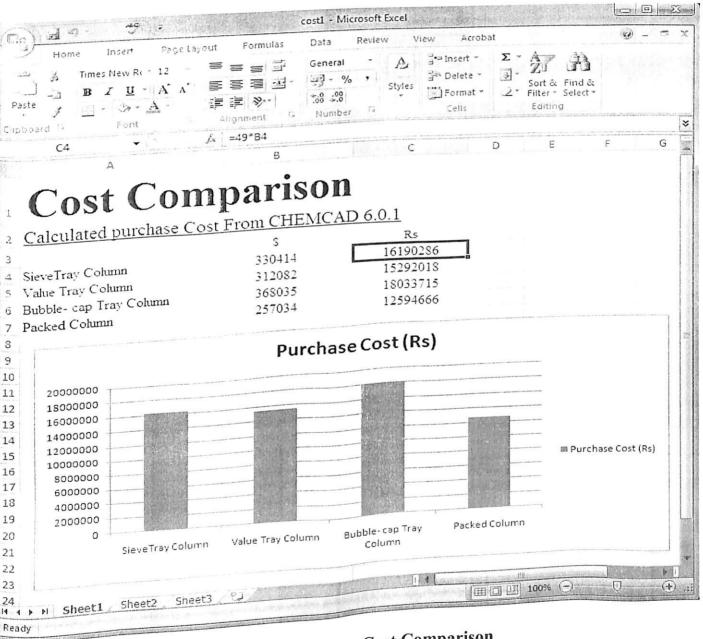


Figure 21: Purchase Cost Comparison

3.2 Installed Cost Comparison

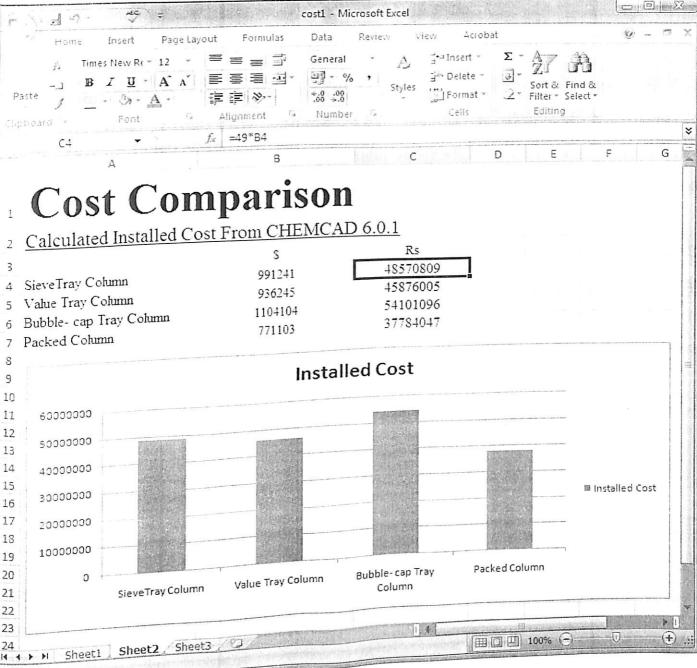


Figure 23: Installed Cost Comparison

Ready

9. CONCLUSIONS

- The project has been successfully focused on Design & Simulation of Gas Processing plant LPG column. LPG column is designed by manually (SPREADSHEATS) as well as simulator (CHEMCAD 6.0.1). The various Tray columns (sieve tray column, value tray column and Bubble cap tray column) and also packed column sizing are designed by CHEMCAD 6.0.1 Simulator.
- Before sizing the tower all the stream composition and property are specified then analysis on characteristics of feed and product streams and also tower are worked out in CHEMCAD 6.0.1. Costing for these different columns are prepared in CHEMCAD 6.0.1 Simulator and finally the cost comparison chart is carried out using SPREADSHEET.
- The project work has enhanced my fundamental knowledge based on "Gas Processing & Handling System Design" as well as "Simulation & Modeling" of M.Tech Gas Engineering.

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