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* XPSIM, Vers. 1.07 *
* Cust/User " " *
* Proj/Problem " " *
* Date MAR 04, 2009 *
* Job " " *
* Page 0001 *
*... extended Process SIMulation ...*
- INPUT -
--- Simulation Input File ---
*-STMT-*
*XPSIM> ...generated by XpsimWin v.1.06 ...
*EXE-OPT> PAGELEN=66
100 RUN ID=PVT05A CUSTOMER=STAFF PROJECT='NORTH AFRICA RESERVOIR'
200 DESC PVT EXPERIMENTS FOR OIL CHARACTERIZATION
300 DESC PRELIMINARY STUDY REQUIRED FOR
400 DESC GATHERING SYSTEM HYDRAULIC STUDY - 24/09/2002
500 DIMENSION INPUT SI PRES=BAR STDVAP=23.69 RTEMP=15.5 RPRES=1
*
*
600 Petroleum Data
*
*
700 HYPO ID=C5+ NAME=C5+ SPGR=.637 MW=71.97 / ID=C6+ NAME=C6+ SPGR=.707 +
MW=84.77 / ID=C7+ NAME=C7+ SPGR=.736 MW=98.27 / ID=C8+ NAME=C8+ +
SPGR=.752 MW=112.09 / ID=C9+ NAME=C9+ SPGR=.767 MW=125.51 / ID=C10+ +
NAME=C10+ SPGR=.761 MW=140.27 / ID=C11+ NAME=C11+ SPGR=.756 +
MW=155.54 TCRIT=370 PCRIT=21.2 / ID=C12+ NAME=C12+ SPGR=.826 +
MW=198.68 TCRIT=425 PCRIT=19.87
800 CALC CRIT=CAVETT
*
*
900 Thermodynamic Data
*
*
1000 METHOD METHOD=SRK
1100 BINARY=C1,C5+ KIJ=0.05,,,,,,,,,,,,,,,,,,,,,
1200 BINARY=C1,C6+ KIJ=0.03,,,,,,,,,,,,,,,,,,,,,
1300 BINARY=C1,C7+ KIJ=0.015,,,,,,,,,,,,,,,,,,,,,
1400 BINARY=C1,C8+ KIJ=0.055,,,,,,,,,,,,,,,,,,,,,
1500 BINARY=C1,C9+ KIJ=0.07,,,,,,,,,,,,,,,,,,,,,
1600 BINARY=C1,C10+ KIJ=0.08,,,,,,,,,,,,,,,,,,,,,
1700 BINARY=C1,C11+ KIJ=0.12,,,,,,,,,,,,,,,,,,,,,
1800 BINARY=C1,C12+ KIJ=0.15,,,,,,,,,,,,,,,,,,,,,
1900 METHOD METHOD=PR
2000 BINARY=C1,C5+ KIJ=0.05,,,,,,,,,,,,,,,,,,,,,
2100 BINARY=C1,C6+ KIJ=0.03,,,,,,,,,,,,,,,,,,,,,
2200 BINARY=C1,C7+ KIJ=0.015,,,,,,,,,,,,,,,,,,,,,
2300 BINARY=C1,C8+ KIJ=0.055,,,,,,,,,,,,,,,,,,,,,
2400 BINARY=C1,C9+ KIJ=0.07,,,,,,,,,,,,,,,,,,,,,
2500 BINARY=C1,C10+ KIJ=0.10,,,,,,,,,,,,,,,,,,,,,
2600 BINARY=C1,C11+ KIJ=0.17,,,,,,,,,,,,,,,,,,,,,
2700 BINARY=C1,C12+ KIJ=0.22,,,,,,,,,,,,,,,,,,,,,
*
*
2800 System Data
*
*
2900 CHEMCOMP H2O / H2S / N2 / CO2 / C1 / C2 / C3 / IC4 / NC4
3000 THERMSET UID=SETF
3100 WOPT K=2
3200 METHODS K=SRK HS=LK CP=IDEAL DV=LK DL=SRK-P IV=LK VIS=IDEAL THC=IDEAL +
SURT=IDEAL
3300 PRINT PETRO METHOD
*
*
3400 PVT Analysis
*
*
3500 STREAM=OIL1
```

* XPSIM, Vers. 1.07 * *... extended Process SIMulation ...*
* Cust/User " " * - INPUT - * Job " " *
* Proj/Problem " " * * Date MAR 04, 2009 *
- Simulation input - Continuation -
*-STMT-
3600 COMP H2O:0 / H2S:0.37 / N2:6.13 / CO2:15.56 / C1:65.85 / C2:4.40 / +
C3:2.18 / IC4:0.56 / NC4:0.95 / C5+:1.03 / C6+:0.66 / C7+:0.65 / +
C8+:0.53 / C9+:0.37 / C10+:0.26 / C11+:0.17 / C12+:0.33
3700 CCE TEMP(CENT)=140 REFSTR=OIL1 DENS=1000 UID=CCE1
3800 DATA PRES=310.3 RV=0.8538 LVF=0.0000 Z=0.980
3900 DATA PRES=303.4 RV=0.8681 LVF=0.0000 Z=0.974
4000 DATA PRES=298.2 RV=0.8792 LVF=0.0000 Z=0.970
4100 DATA PRES=293.5 RV=0.8900 LVF=0.0000 Z=0.966
4200 DATA PRES=288.7 RV=0.9014 LVF=0.0000 Z=0.962
4300 DATA PRES=282.6 RV=0.9167 LVF=0.0000 Z=0.958
4400 DATA PRES=275.9 RV=0.9347 LVF=0.0000 Z=0.954
4500 DATA PRES=270.3 RV=0.9503 LVF=0.0000 Z=0.950
4600 DATA PRES=262.0 RV=0.9755 LVF=0.0000 Z=0.945
4700 DATA PRES=254.4 RV=1.0000 LVF=0.0000 Z=0.941
4800 DATA PRES=248.6 RV=1.0203 LVF=0.00056
4900 DATA PRES=240.6 RV=1.0500 LVF=0.00139
5000 DATA PRES=228.9 RV=1.0979 LVF=0.00268
5100 DATA PRES=215.5 RV=1.1597 LVF=0.00429
5200 DATA PRES=191.0 RV=1.2982 LVF=0.00750
5300 DATA PRES=160.6 RV=1.5404 LVF=0.01079
5400 DATA PRES=123.1 RV=2.0188 LVF=0.01342
5500 DATA PRES=100.8 RV=2.4827 LVF=0.01423
5600 DATA PRES=80.6 RV=3.1345 LVF=0.01439
5700 DATA PRES=59.0 RV=4.3389 LVF=0.01392
5800 CVD TEMP(CENT)=140 REFSTR=OIL1 UID=CVD1
5900 DATA PRES=254.4 RV=0.0000 LVF=0.0000 Z=0.941
6000 DATA PRES=247.9 RV=0.0208 LVF=0.00042 Z=0.919
6100 DATA PRES=237.2 RV=0.0590 LVF=0.00167 Z=0.913
6200 DATA PRES=212.7 RV=0.1483 LVF=0.00432 Z=0.902
6300 DATA PRES=185.2 RV=0.2542 LVF=0.00771 Z=0.893
6400 DATA PRES=153.1 RV=0.3834 LVF=0.01009 Z=0.885
6500 DATA PRES=118.2 RV=0.5285 LVF=0.01167 Z=0.881
6600 DATA PRES=79.8 RV=0.6878 LVF=0.01143 Z=0.882
6700 DATA PRES=38.2 RV=0.8484 LVF=0.01080 Z=0.893
6800 SEP REFSTR=OIL1 STOD=759 GG=1.268 GOR=1000 UID=SEP1
6900 DATA TEMP(K)=307.6 PRES=37.6
7000 LCOMP H2O:0 / H2S:0.20 / N2:0.24 / CO2:4.95 / C1:10.24 / C2:3.15 / +
C3:4.42 / IC4:2.00 / NC4:6.01 / C5+:11.59 / C6+:10.54 / C7+:11.74 / +
C8+:10.63 / C9+:7.90 / C10+:5.60 / C11+:3.67 / C12+:7.13
*
*
7100 Flowsheet Data
*
*
7200 EQUI 3PHASE LLSM=LHLW KEY1=H2O
*
7300 STREAM=WEST TEMP=40.55 PRES=172.8 RATE(M)=100 NORM XBASIS=M
7400 COMP H2O:0 / H2S:0.37 / N2:6.13 / CO2:15.56 / C1:65.85 / C2:4.40 / +
C3:2.18 / IC4:0.56 / NC4:0.95 / C5+:1.03 / C6+:0.66 / C7+:0.65 / +
C8+:0.53 / C9+:0.37 / C10+:0.26 / C11+:0.17 / C12+:0.33
*
7500 PHASENV IN WEST UID=PHASE1
7600 CALC PMAX=400
7700 PRINT PLOT
*
7800 FLASH IN WEST OUT GASA0065(V) LIQA0065(L) UID=CYLIND
7900 CALC ISO TEMP=34.4 PRES=37.6
*
8000 FLASH IN LIQA0065 OUT SGS03(V) STDLIQ(L) UID=STD
8100 CALC ISO TEMP=15.56 PRES=1.013

- Simulation input - Continuation -

-STMT-

8200	FLASH IN GASA0065 SGS03 OUT STDGAS UID=RECOMB
8300	CALC ISO TEMP=15.56 PRES=1.013
8400	FLASH IN WEST OUT DEW UID=DEW
8500	CALC DEW TEMP=140.55 PEST=250
8600	VLE Analysis
8700	VLECURVE STR=WEST
8800	CALC ISO TEMP=140.5 PRES=100,300 POINTS=11
8900	Output Data
9000	PRINT FORMAT=2 MFRAC
9100	END

--- End of Simulation Input File ---

I) * PROBLEM GENERAL DATA *

1) * PROBLEM/PROJECT *
'PVT EXPERIMENTS FOR OIL CHARACTERIZATION'
'PRELIMINARY STUDY REQUIRED FOR'
' GATHERING SYSTEM HYDRAULIC STUDY - 24/09/2002'

2) * UNITS OF MEASURE *
- Input system SI - Output system SI

- INPUT AND OUTPUT UNITS -
Time - HR Weight - KG
Temperature - CENT Pressure - BAR
Energy/Duty - M-KJ Work - KW
Liq volums - CUMT Vap Volume - CUMT
Viscosity - CP Thermal cond - W/MC
Surface tens - DYCM Std Vap Vol - N-CUMT

Standard Vapor Volume is 23.69000 M3/KMOL
Reference Status - Temperature 15.5 - Pressure 1

Enthalpy is 0.0 for ideal gas at 298.15 K
Std entropy is for ideal gas at 298.15 K and 1 Atm

*** HYPOTHETICAL/PETROLEUM COMPONENTS CALCULATION OPTIONS ***

Initial point interval 2.00
End point interval 98.00

Critical props calculation method 'CAVETT '

I.D) * USER'S SUPPLIED THERMODYNAMIC DATA *

1) * METHOD 'SRK' *

No. of COMPONENT parameters 3
No. of BINARY parameters 1

* INTERACTION/BINARY CONSTANTS BETWEEN COMPONENT					
C1	-	'	'	and	C5+
1)	0.500000E-01				
C1	-	'	'	and	C6+
1)	0.300000E-01				
C1	-	'	'	and	C7+
1)	0.150000E-01				
C1	-	'	'	and	C8+
1)	0.550000E-01				
C1	-	'	'	and	C9+
1)	0.700000E-01				
C1	-	'	'	and	C10+
1)	0.800000E-01				
C1	-	'	'	and	C11+
1)	0.120000				
C1	-	'	'	and	C12+
1)	0.150000				

2) * METHOD 'PR' *

No. of COMPONENT parameters 3
No. of BINARY parameters 1

* INTERACTION/BINARY CONSTANTS BETWEEN COMPONENT					
C1	-	'	'	and	C5+
1)	0.500000E-01				
C1	-	'	'	and	C6+
1)	0.300000E-01				
C1	-	'	'	and	C7+
1)	0.150000E-01				
C1	-	'	'	and	C8+
1)	0.550000E-01				
C1	-	'	'	and	C9+
1)	0.700000E-01				
C1	-	'	'	and	C10+
1)	0.100000				
C1	-	'	'	and	C11+
1)	0.170000				

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* XPSIM, Vers. 1.07 *
* Cust/User "STAFF" *
* Proj/Problem "NORTH AFRICA RESERVOIR" *
* Date MAR 04, 2009 *
*... extended Process SIMulation ...*
- INPUT -
* Job "PVT05A" *
C1 - '
1) 0.220000 ' and C12+ - '
'
```

II) * DEFINED COMPONENTS *

* No of Chemical Components 17
No of Hypothetical/Petroleum Components 8

No	1	2	3	4	5
CAS Registry no.	7732-18-5	7783-06-4	7727-37-9	124-38-9	74-82-8
Name	WATER	HYDROGEN SULPHIDE	NITROGEN	CARBON DIOXIDE	METHANE
Component Key	H2O	H2S	N2	CO2	C1
Type	LIBRARY	LIBRARY	LIBRARY	LIBRARY	LIBRARY
Class	INOR	INOR	INOR	INOR	SHYD
Formula	H2O	H2S	N2	CO2	CH4
Molecular weight	18.020	34.080	28.010	44.010	16.040
Boiling point,CENT	100.00	-60.25	-195.80	-78.55	-161.52
Std spec. gravity	1.00100	0.78638	0.64315	0.81906	0.40749
Critical temp,CENT	374.15	100.05	-146.95	31.06	-82.62
Critical pres,BAR	221.200	89.400	34.000	73.825	45.950
Critical vol,CUMT	0.056313	0.098498	0.089489	0.094039	0.099013
Critical Z	0.231459	0.283798	0.289987	0.274489	0.287211
Acentric factor	0.344000	0.104000	0.037000	0.223000	0.011000
Lat. heat, KJ/KMOL	40714.413	18893.595	5552.287	15245.787	8176.826
H form, KJ/KMOL	-241828.891	-20584.362	0.000	-393494.298	-74586.151
G form, KJ/KMOL	-298119.724	-81913.940	-57066.025	-447817.337	-130101.793
Std entr, KJ/KML-C	188.80073	205.70080	191.40074	182.20071	186.20072

No	6	7	8	9	10
CAS Registry no.	74-84-0	74-98-6	75-28-5	106-97-8	00000-00-0
Name	ETHANE	PROPANE	ISOBUTANE	BUTANE	C5+
Component Key	C2	C3	IC4	NC4	C5+
Type	LIBRARY	LIBRARY	LIBRARY	LIBRARY	PETRO
Class	SHYD	SHYD	SHYD	SHYD	PETR
Formula	C2H6	C3H8	C4H10	C4H10	
Molecular weight	30.070	44.100	58.120	58.120	71.970
Boiling point,CENT	-88.60	-42.10	-11.73	-0.50	37.97
Std spec. gravity	0.35522	0.50430	0.55891	0.58434	0.63700
Critical temp,CENT	32.27	96.67	134.99	152.03	199.93
Critical pres,BAR	48.790	42.490	36.480	37.980	33.488
Critical vol,CUMT	0.148129	0.203227	0.262988	0.254914	0.318220
Critical Z	0.284617	0.280844	0.282728	0.273882	0.270941
Acentric factor	0.099000	0.153000	0.183000	0.200000	0.250739
Lat. heat, KJ/KMOL	14535.497	18848.524	21127.358	21978.088	26237.364
H form, KJ/KMOL	-84015.754	-103855.711	-134606.193	-125771.935	0.000
G form, KJ/KMOL	-152322.057	-184416.002	-222679.880	-218138.991	0.000
Std entr, KJ/KML-C	229.10089	270.20105	295.40115	309.80120	156.48729

No	11	12	13	14	15
CAS Registry no.	00000-00-0	00000-00-0	00000-00-0	00000-00-0	00000-00-0
Name	C6+	C7+	C8+	C9+	C10+
Component Key	C6+	C7+	C8+	C9+	C10+
Type	PETRO	PETRO	PETRO	PETRO	PETRO
Class	PETR	PETR	PETR	PETR	PETR
Formula					
Molecular weight	84.770	98.270	112.090	125.510	140.270
Boiling point,CENT	69.56	97.65	123.67	148.54	167.58
Std spec. gravity	0.70700	0.73600	0.75200	0.76700	0.76100
Critical temp,CENT	238.29	268.39	293.82	317.73	329.93
Critical pres,BAR	33.124	31.327	29.090	27.095	24.260
Critical vol,CUMT	0.340883	0.373737	0.412255	0.451548	0.501459
Critical Z	0.265543	0.260040	0.254413	0.249044	0.242629
Acentric factor	0.318216	0.386998	0.457344	0.524446	0.604638
Lat. heat, KJ/KMOL	29176.403	31810.732	34267.166	36627.102	38443.187
H form, KJ/KMOL	0.000	0.000	0.000	0.000	0.000
G form, KJ/KMOL	0.000	0.000	0.000	0.000	0.000
Std entr, KJ/KML-C	72.90771	57.74207	73.25436	88.16845	168.10503

No	16	17
CAS Registry no.	00000-00-0	00000-00-0
Name	C11+	C12+
Component	Key	Key
	Type	Type
	Class	Class
Formula		
Molecular weight	155.540	198.680
Boiling point,CENT	185.15	248.57
Std spec. gravity	0.75600	0.82600
Critical temp,CENT	370.00	425.00
Critical pres,BAR	21.200	19.870
Critical vol,CUMT	0.595266	0.664265
Critical Z	0.236005	0.227393
Acentric factor	0.687435	0.795084
Lat. heat, KJ/KMOL	40124.206	46239.246
H form, KJ/KMOL	0.000	0.000
G form, KJ/KMOL	0.000	0.000
Std entr, KJ/KML-C	251.88085	157.01906

II-A) * PETROLEUM / HYPOTHETICAL COMPONENTS DATA *

Number	10	11	12	13	14	15
Code	C5+	C6+	C7+	C8+	C9+	C10+
Boiling point, CENT	37.973	69.558	97.649	123.674	148.537	167.582
API density	90.6350	68.6414	60.7554	56.6649	52.9850	54.4396
Specific gravity	0.63700	0.70700	0.73600	0.75200	0.76700	0.76100
K-UOP factor	12.9398	12.0405	11.8739	11.8870	11.8930	12.1646
Molecular weight	71.970	84.770	98.270	112.090	125.510	140.270
Viscosity at 100 F,cst	0.23056	0.31222	0.40318	0.51260	0.65536	0.77579
Viscosity at 210 F,cst	0.27387	0.31444	0.36156	0.41595	0.47991	0.53934
Critical temperature,CENT	199.929	238.294	268.387	293.822	317.727	329.928
Critical pressure,BAR	33.4883	33.1239	31.3267	29.0903	27.0946	24.2601
Acentric factor	0.25074	0.31822	0.38700	0.45734	0.52445	0.60464

Number	16	17
Code	C11+	C12+
Boiling point, CENT	185.148	248.574
API density	55.6693	39.8075
Specific gravity	0.75600	0.82600
K-UOP factor	12.4056	11.8556
Molecular weight	155.540	198.680
Viscosity at 100 F,cst	0.91293	2.1816
Viscosity at 210 F,cst	0.60334	0.96144
Critical temperature,CENT	370.000	425.000
Critical pressure,BAR	21.2000	19.8700
Acentric factor	0.68743	0.79508

III) * PVT/THERMO/TRANSPORT PROPERTIES CALCULATION METHODS *

* CALCULATION SET 1 - SETF *

VLE K-values	SRK	- REDLICH-KWONG-SOAVE
ENTHALPY - Vapor	LK	- LEE-KESLER
ENTHALPY - Liquid	LK	- LEE-KESLER
ENTROPY - Vapor	LK	- LEE-KESLER
ENTROPY - Liquid	LK	- LEE-KESLER
DENSITY - Vapor	LK	- LEE-KESLER
DENSITY - Liquid	SRK-P	- SRK - PENELOUX MOD
VISCOSITY - Vapor	IDEAL	- IDEAL/LIBRARY
VISCOSITY - Liquid	IDEAL	- IDEAL/LIBRARY
HEAT CAPACITY - Vapor	IDEAL	- IDEAL/LIBRARY
HEAT CAPACITY - Liquid	IDEAL	- IDEAL/LIBRARY
CONDUCTIVITY - Vapor	IDEAL	- IDEAL/LIBRARY
CONDUCTIVITY - Liquid	IDEAL	- IDEAL/LIBRARY
ISENTROPIC Vapor Coeff	LK	- LEE-KESLER
SURFACE TENSION	IDEAL	- IDEAL/LIBRARY

- LIMITS AND OPTIONS -
Temperature - Min -223.15 CENT - Max 1200.05 CENT
Pressure - Min 0.100000E-04 BAR - Max 3000.00 BAR

Lowest significant composition 0.10000000E-19

Water K-values are calculated based on
'NGPA vapor pressure chart'
Water thermo-props are calculated based on
'Saturated conditions'
Water solubility in hydr phases calculated based on
'API solubility in HC mixtures'

*
 CONSTANTS/PARAMETERS FOR CALCULATION SET
 1
 SETF

1)
 Parameters of method SRK
 -
 REDLICH-KWONG-SOAVE

No. of component parameters
 3
 No. of binary parameters
 1

Component
 1
 H2O
 -
 'WATER
 '

1)
 1165.14
 2)
 3208.24
 3)
 0.344000

-
 Interaction/Binary parameters

2
 H2S
 1)
 0.000000

3
 N2
 1)
 0.000000

4
 C02
 1)
 0.000000

5
 C1
 1)
 0.500000

6
 C2
 1)
 0.500000

7
 C3
 1)
 0.500000

8
 IC4
 1)
 0.500000

9
 NC4
 1)
 0.500000

10
 C5+
 1)
 0.500000

11
 C6+
 1)
 0.500000

12
 C7+
 1)
 0.500000

13
 C8+
 1)
 0.500000

14
 C9+
 1)
 0.500000

15
 C10+
 1)
 0.500000

16
 C11+
 1)
 0.500000

17
 C12+
 1)
 0.500000

Component
 2
 H2S
 -
 'HYDROGEN SULPHIDE
 '

1)
 671.760
 2)
 1296.64
 3)
 0.104000

-
 Interaction/Binary parameters

3
 N2
 1)
 0.000000

4
 C02
 1)
 0.000000

5
 C1
 1)
 0.900000E-01

6
 C2
 1)
 0.850000E-01

7
 C3
 1)
 0.760000E-01

8
 IC4
 1)
 0.600000E-01

9
 NC4
 1)
 0.600000E-01

10
 C5+
 1)
 0.600000E-01

11
 C6+
 1)
 0.600000E-01

12
 C7+
 1)
 0.600000E-01

13
 C8+
 1)
 0.500000E-01

14
 C9+
 1)
 0.500000E-01

15
 C10+
 1)
 0.500000E-01

16
 C11+
 1)
 0.500000E-01

17
 C12+
 1)
 0.500000E-01

Component
 3
 N2
 -
 'NITROGEN
 '

1)
 227.160
 2)
 493.129
 3)
 0.370000E-01

-
 Interaction/Binary parameters

4
 C02
 1)
 0.000000

5
 C1
 1)
 0.300000E-01

6
 C2
 1)
 0.600000E-01

7
 C3
 1)
 0.900000E-01

8
 IC4
 1)
 0.113000

9
 NC4
 1)
 0.113000

10
 C5+
 1)
 0.140000

11
 C6+
 1)
 0.166000

12
 C7+
 1)
 0.200000

13
 C8+
 1)
 0.200000

```

* XPSIM, Vers. 1.07 *
* Cust/User "STAFF" " *
* Proj/Problem "NORTH AFRICA RESERVOIR " *
14 C9+ 1) 0.200000
15 C10+ 1) 0.200000
16 C11+ 1) 0.200000
17 C12+ 1) 0.200000

Component 4 CO2 - 'CARBON DIOXIDE'
1) 547.578 2) 1070.74 3) 0.223000
- Interaction/Binary parameters
5 C1 1) 0.933000E-01
6 C2 1) 0.136300
7 C3 1) 0.128900
8 IC4 1) 0.128500
9 NC4 1) 0.143000
10 C5+ 1) 0.131000
11 C6+ 1) 0.112000
12 C7+ 1) 0.110000
13 C8+ 1) 0.120000
14 C9+ 1) 0.125000
15 C10+ 1) 0.130000
16 C11+ 1) 0.130000
17 C12+ 1) 0.130000

Component 5 C1 - 'METHANE'
1) 342.954 2) 666.450 3) 0.110000E-01
- Interaction/Binary parameters
6 C2 1) 0.000000
7 C3 1) 0.000000
8 IC4 1) 0.000000
9 NC4 1) 0.000000
10 C5+ 1) 0.500000E-01
11 C6+ 1) 0.300000E-01
12 C7+ 1) 0.150000E-01
13 C8+ 1) 0.550000E-01
14 C9+ 1) 0.700000E-01
15 C10+ 1) 0.800000E-01
16 C11+ 1) 0.120000
17 C12+ 1) 0.150000

Component 6 C2 - 'ETHANE'
1) 549.756 2) 707.640 3) 0.990000E-01
- Interaction/Binary parameters
7 C3 1) 0.000000
8 IC4 1) 0.000000
9 NC4 1) 0.000000
10 C5+ 1) 0.000000
11 C6+ 1) 0.000000
12 C7+ 1) 0.000000
13 C8+ 1) 0.000000
14 C9+ 1) 0.000000
15 C10+ 1) 0.000000
16 C11+ 1) 0.000000
17 C12+ 1) 0.000000

Component 7 C3 - 'PROPANE'
1) 665.676 2) 616.266 3) 0.153000
- Interaction/Binary parameters
8 IC4 1) 0.000000
9 NC4 1) 0.000000
10 C5+ 1) 0.000000
11 C6+ 1) 0.000000
12 C7+ 1) 0.000000
13 C8+ 1) 0.000000

```

```

* XPSIM, Vers. 1.07 *
* Cust/User "STAFF" " *
* Proj/Problem "NORTH AFRICA RESERVOIR " *
14 C9+ 1) 0.00000
15 C10+ 1) 0.00000
16 C11+ 1) 0.00000
17 C12+ 1) 0.00000

Component 8 IC4 - 'ISOBUTANE'
1) 734.652 2) 529.099 3) 0.183000
- Interaction/Binary parameters
9 NC4 1) 0.00000
10 C5+ 1) 0.00000
11 C6+ 1) 0.00000
12 C7+ 1) 0.00000
13 C8+ 1) 0.00000
14 C9+ 1) 0.00000
15 C10+ 1) 0.00000
16 C11+ 1) 0.00000
17 C12+ 1) 0.00000

Component 9 NC4 - 'BUTANE'
1) 765.324 2) 550.854 3) 0.200000
- Interaction/Binary parameters
10 C5+ 1) 0.00000
11 C6+ 1) 0.00000
12 C7+ 1) 0.00000
13 C8+ 1) 0.00000
14 C9+ 1) 0.00000
15 C10+ 1) 0.00000
16 C11+ 1) 0.00000
17 C12+ 1) 0.00000

Component 10 C5+ - 'C5+'
1) 851.541 2) 485.707 3) 0.250739
- Interaction/Binary parameters
11 C6+ 1) 0.00000
12 C7+ 1) 0.00000
13 C8+ 1) 0.00000
14 C9+ 1) 0.00000
15 C10+ 1) 0.00000
16 C11+ 1) 0.00000
17 C12+ 1) 0.00000

Component 11 C6+ - 'C6+'
1) 920.600 2) 480.422 3) 0.318216
- Interaction/Binary parameters
12 C7+ 1) 0.00000
13 C8+ 1) 0.00000
14 C9+ 1) 0.00000
15 C10+ 1) 0.00000
16 C11+ 1) 0.00000
17 C12+ 1) 0.00000

Component 12 C7+ - 'C7+'
1) 974.767 2) 454.356 3) 0.386998
- Interaction/Binary parameters
13 C8+ 1) 0.00000
14 C9+ 1) 0.00000
15 C10+ 1) 0.00000
16 C11+ 1) 0.00000
17 C12+ 1) 0.00000

```

Component 13 C8+	- 'C8+	'
1) 1020.55	2) 421.919	3) 0.457344
- Interaction/Binary parameters		
14 C9+	1) 0.00000	
15 C10+	1) 0.00000	
16 C11+	1) 0.00000	
17 C12+	1) 0.00000	
Component 14 C9+	- 'C9+	'
1) 1063.58	2) 392.975	3) 0.524446
- Interaction/Binary parameters		
15 C10+	1) 0.00000	
16 C11+	1) 0.00000	
17 C12+	1) 0.00000	
Component 15 C10+	- 'C10+	'
1) 1085.54	2) 351.864	3) 0.604638
- Interaction/Binary parameters		
16 C11+	1) 0.00000	
17 C12+	1) 0.00000	
Component 16 C11+	- 'C11+	'
1) 1157.67	2) 307.481	3) 0.687435
- Interaction/Binary parameters		
17 C12+	1) 0.00000	
Component 17 C12+	- 'C12+	'
1) 1256.67	2) 288.191	3) 0.795084

2) Parameters of method LK - LEE-KESLER

General component parameters are used

3) Parameters of method SRK-P - SRK - PENELOUX MOD

No. of component parameters 3
 No. of binary parameters 1

Component 1 H2O	- 'WATER	'
1) 0.00000	2) 0.00000	3) 0.00000
- Interaction/Binary parameters		
2 H2S	1) 0.00000	
3 N2	1) 0.00000	
4 CO2	1) 0.00000	
5 C1	1) 0.00000	
6 C2	1) 0.00000	
7 C3	1) 0.00000	
8 IC4	1) 0.00000	
9 NC4	1) 0.00000	
10 C5+	1) 0.00000	
11 C6+	1) 0.00000	
12 C7+	1) 0.00000	
13 C8+	1) 0.00000	
14 C9+	1) 0.00000	
15 C10+	1) 0.00000	
16 C11+	1) 0.00000	
17 C12+	1) 0.00000	

Component 2 H2S

- 'HYDROGEN SULPHIDE'

1) 0.00000 2) 0.00000 3) 0.00000

- Interaction/Binary parameters

3 N2 1) 0.00000

4 CO2 1) 0.00000

5 C1 1) 0.00000

6 C2 1) 0.00000

7 C3 1) 0.00000

8 IC4 1) 0.00000

9 NC4 1) 0.00000

10 C5+ 1) 0.00000

11 C6+ 1) 0.00000

12 C7+ 1) 0.00000

13 C8+ 1) 0.00000

14 C9+ 1) 0.00000

15 C10+ 1) 0.00000

16 C11+ 1) 0.00000

17 C12+ 1) 0.00000

Component 3 N2

- 'NITROGEN'

1) 0.00000 2) 0.00000 3) 0.00000

- Interaction/Binary parameters

4 CO2 1) 0.00000

5 C1 1) 0.00000

6 C2 1) 0.00000

7 C3 1) 0.00000

8 IC4 1) 0.00000

9 NC4 1) 0.00000

10 C5+ 1) 0.00000

11 C6+ 1) 0.00000

12 C7+ 1) 0.00000

13 C8+ 1) 0.00000

14 C9+ 1) 0.00000

15 C10+ 1) 0.00000

16 C11+ 1) 0.00000

17 C12+ 1) 0.00000

Component 4 CO2

- 'CARBON DIOXIDE'

1) 0.00000 2) 0.00000 3) 0.00000

- Interaction/Binary parameters

5 C1 1) 0.00000

6 C2 1) 0.00000

7 C3 1) 0.00000

8 IC4 1) 0.00000

9 NC4 1) 0.00000

10 C5+ 1) 0.00000

11 C6+ 1) 0.00000

12 C7+ 1) 0.00000

13 C8+ 1) 0.00000

14 C9+ 1) 0.00000

15 C10+ 1) 0.00000

16 C11+ 1) 0.00000

17 C12+ 1) 0.00000

Component 5 C1

- 'METHANE'

1) -116.630 2) 667.800 3) 0.287600

- Interaction/Binary parameters

6 C2 1) 0.00000

7 C3 1) 0.00000

8 IC4 1) 0.00000

9 NC4 1) 0.00000

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* XPSIM, Vers. 1.07 *
* Cust/User "STAFF" " *
* Proj/Problem "NORTH AFRICA RESERVOIR " *
* Date MAR 04, 2009 *

*... extended Process SIMulation ...*
* Job "PVT05A" " *

10 C5+ 1) 0.00000
11 C6+ 1) 0.00000
12 C7+ 1) 0.00000
13 C8+ 1) 0.00000
14 C9+ 1) 0.00000
15 C10+ 1) 0.00000
16 C11+ 1) 0.00000
17 C12+ 1) 0.00000

Component 6 C2 - 'ETHANE'
1) 90.0900 2) 707.800 3) 0.278900
- Interaction/Binary parameters
7 C3 1) 0.00000
8 IC4 1) 0.00000
9 NC4 1) 0.00000
10 C5+ 1) 0.00000
11 C6+ 1) 0.00000
12 C7+ 1) 0.00000
13 C8+ 1) 0.00000
14 C9+ 1) 0.00000
15 C10+ 1) 0.00000
16 C11+ 1) 0.00000
17 C12+ 1) 0.00000

Component 7 C3 - 'PROPANE'
1) 206.010 2) 616.300 3) 0.276300
- Interaction/Binary parameters
8 IC4 1) 0.00000
9 NC4 1) 0.00000
10 C5+ 1) 0.00000
11 C6+ 1) 0.00000
12 C7+ 1) 0.00000
13 C8+ 1) 0.00000
14 C9+ 1) 0.00000
15 C10+ 1) 0.00000
16 C11+ 1) 0.00000
17 C12+ 1) 0.00000

Component 8 IC4 - 'ISOBUTANE'
1) 274.980 2) 529.100 3) 0.275000
- Interaction/Binary parameters
9 NC4 1) 0.00000
10 C5+ 1) 0.00000
11 C6+ 1) 0.00000
12 C7+ 1) 0.00000
13 C8+ 1) 0.00000
14 C9+ 1) 0.00000
15 C10+ 1) 0.00000
16 C11+ 1) 0.00000
17 C12+ 1) 0.00000

Component 9 NC4 - 'BUTANE'
1) 305.650 2) 550.700 3) 0.272800
- Interaction/Binary parameters
10 C5+ 1) 0.00000
11 C6+ 1) 0.00000
12 C7+ 1) 0.00000
13 C8+ 1) 0.00000
14 C9+ 1) 0.00000
15 C10+ 1) 0.00000
16 C11+ 1) 0.00000
17 C12+ 1) 0.00000

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* XPSIM, Vers. 1.07 *
* Cust/User "STAFF" " *
* Proj/Problem "NORTH AFRICA RESERVOIR " *
* Date MAR 04, 2009 *
* Page 0017 *
* Job "PVT05A" " *

... extended Process - SIMULATION ...
- INPUT -

Component 10 C5+ - 'C5+'
1) 0.00000 2) 0.00000 3) 0.00000
- Interaction/Binary parameters
11 C6+ 1) 0.00000
12 C7+ 1) 0.00000
13 C8+ 1) 0.00000
14 C9+ 1) 0.00000
15 C10+ 1) 0.00000
16 C11+ 1) 0.00000
17 C12+ 1) 0.00000

Component 11 C6+ - 'C6+'
1) 0.00000 2) 0.00000 3) 0.00000
- Interaction/Binary parameters
12 C7+ 1) 0.00000
13 C8+ 1) 0.00000
14 C9+ 1) 0.00000
15 C10+ 1) 0.00000
16 C11+ 1) 0.00000
17 C12+ 1) 0.00000

Component 12 C7+ - 'C7+'
1) 0.00000 2) 0.00000 3) 0.00000
- Interaction/Binary parameters
13 C8+ 1) 0.00000
14 C9+ 1) 0.00000
15 C10+ 1) 0.00000
16 C11+ 1) 0.00000
17 C12+ 1) 0.00000

Component 13 C8+ - 'C8+'
1) 0.00000 2) 0.00000 3) 0.00000
- Interaction/Binary parameters
14 C9+ 1) 0.00000
15 C10+ 1) 0.00000
16 C11+ 1) 0.00000
17 C12+ 1) 0.00000

Component 14 C9+ - 'C9+'
1) 0.00000 2) 0.00000 3) 0.00000
- Interaction/Binary parameters
15 C10+ 1) 0.00000
16 C11+ 1) 0.00000
17 C12+ 1) 0.00000

Component 15 C10+ - 'C10+'
1) 0.00000 2) 0.00000 3) 0.00000
- Interaction/Binary parameters
16 C11+ 1) 0.00000
17 C12+ 1) 0.00000

Component 16 C11+ - 'C11+'
1) 0.00000 2) 0.00000 3) 0.00000
- Interaction/Binary parameters
17 C12+ 1) 0.00000

Component 17 C12+ - 'C12+'
1) 0.00000 2) 0.00000 3) 0.00000

```

*** PVT ANALYSIS ***

1) *** STREAM 'OIL1' ***

* No *	Component	* - Mols -
2	HYDROGEN SULPHIDE	0.003700
3	NITROGEN	0.061300
4	CARBON DIOXIDE	0.155600
5	METHANE	0.658500
6	ETHANE	0.044000
7	PROPANE	0.021800
8	ISOBUTANE	0.005600
9	BUTANE	0.009500
10	C5+	0.010300
11	C6+	0.006600
12	C7+	0.006500
13	C8+	0.005300
14	C9+	0.003700
15	C10+	0.002600
16	C11+	0.001700
17	C12+	0.003300

2) *** EXPERIMENT NO. 1 CCE1 ***
COSTANT COMPOSITION EXPANSION

Temperature 140.000 CENT
Saturation density 1000.0000 KG/CUMT
Reference stream OIL1

* EXPERIMENTAL POINTS *

POINT NO. 1
Pressure 310.300 BAR
Relative volume 0.854
Liquid volume fraction 0.000
Gas Z factor 0.980

POINT NO. 2
Pressure 303.400 BAR
Relative volume 0.868
Liquid volume fraction 0.000
Gas Z factor 0.974

POINT NO. 3
Pressure 298.200 BAR
Relative volume 0.879
Liquid volume fraction 0.000
Gas Z factor 0.970

POINT NO. 4
Pressure 293.500 BAR
Relative volume 0.890
Liquid volume fraction 0.000
Gas Z factor 0.966

POINT NO.	5		
Pressure	288.700	BAR	
Relative volume	0.901		
Liquid volume fraction			0.000
Gas Z factor	0.962		
POINT NO.	6		
Pressure	282.600	BAR	
Relative volume	0.917		
Liquid volume fraction			0.000
Gas Z factor	0.958		
POINT NO.	7		
Pressure	275.900	BAR	
Relative volume	0.935		
Liquid volume fraction			0.000
Gas Z factor	0.954		
POINT NO.	8		
Pressure	270.300	BAR	
Relative volume	0.950		
Liquid volume fraction			0.000
Gas Z factor	0.950		
POINT NO.	9		
Pressure	262.000	BAR	
Relative volume	0.976		
Liquid volume fraction			0.000
Gas Z factor	0.945		
POINT NO.	10		
Pressure	254.400	BAR	
Relative volume	1.000		
Liquid volume fraction			0.000
Gas Z factor	0.941		
POINT NO.	11		
Pressure	248.600	BAR	
Relative volume	1.020		
Liquid volume fraction			0.001
POINT NO.	12		
Pressure	240.600	BAR	
Relative volume	1.050		
Liquid volume fraction			0.001
POINT NO.	13		
Pressure	228.900	BAR	
Relative volume	1.098		
Liquid volume fraction			0.003
POINT NO.	14		
Pressure	215.500	BAR	
Relative volume	1.160		
Liquid volume fraction			0.004

POINT NO.	15
Pressure	191.000 BAR
Relative volume	1.298
Liquid volume fraction	0.007
POINT NO.	16
Pressure	160.600 BAR
Relative volume	1.540
Liquid volume fraction	0.011
POINT NO.	17
Pressure	123.100 BAR
Relative volume	2.019
Liquid volume fraction	0.013
POINT NO.	18
Pressure	100.800 BAR
Relative volume	2.483
Liquid volume fraction	0.014
POINT NO.	19
Pressure	80.600 BAR
Relative volume	3.135
Liquid volume fraction	0.014
POINT NO.	20
Pressure	59.000 BAR
Relative volume	4.339
Liquid volume fraction	0.014

3) *** EXPERIMENT NO. 2 CVD1 ***
COSTANT VOLUME DEPLETION

Temperature 140.000 CENT
Reference stream OIL1

* EXPERIMENTAL POINTS *

POINT NO.	1		
Pressure	254.400	BAR	
Relative volume	0.000		
Liquid volume fraction	0.000		0.000
Gas Z factor	0.941		
POINT NO.	2		
Pressure	247.900	BAR	
Relative volume	0.021		
Liquid volume fraction	0.000		0.000
Gas Z factor	0.919		
POINT NO.	3		
Pressure	237.200	BAR	
Relative volume	0.059		
Liquid volume fraction	0.002		0.002
Gas Z factor	0.913		
POINT NO.	4		
Pressure	212.700	BAR	
Relative volume	0.148		
Liquid volume fraction	0.004		0.004
Gas Z factor	0.902		
POINT NO.	5		
Pressure	185.200	BAR	
Relative volume	0.254		
Liquid volume fraction	0.008		0.008
Gas Z factor	0.893		
POINT NO.	6		
Pressure	153.100	BAR	
Relative volume	0.383		
Liquid volume fraction	0.010		0.010
Gas Z factor	0.885		
POINT NO.	7		
Pressure	118.200	BAR	
Relative volume	0.528		
Liquid volume fraction	0.012		0.012
Gas Z factor	0.881		
POINT NO.	8		
Pressure	79.800	BAR	
Relative volume	0.688		
Liquid volume fraction	0.011		0.011
Gas Z factor	0.882		

POINT NO. 9
Pressure 38.200 BAR
Relative volume 0.848
Liquid volume fraction 0.011
Gas Z factor 0.893

4) *** EXPERIMENT NO. 3 SEP1 ***
TEST SEPARATORS

Reference stream OIL1
Oil density 759.000 KG/CUMT
Gas gravity 1.2680
Gas/oil ratio 1000.000 NM3/M3

* EXPERIMENTAL POINTS *

POINT NO. 1
Temperature 307.600 K
Pressure 37.600 BAR

* LIQUID COMPOSITION *

* No *	Component	* - Mols -
2	HYDROGEN SULPHIDE	0.002000
3	NITROGEN	0.002400
4	CARBON DIOXIDE	0.049495
5	METHANE	0.102390
6	ETHANE	0.031497
7	PROPANE	0.044196
8	ISOBUTANE	0.019998
9	BUTANE	0.060094
10	C5+	0.115888
11	C6+	0.105389
12	C7+	0.117388
13	C8+	0.106289
14	C9+	0.078992
15	C10+	0.055994
16	C11+	0.036696
17	C12+	0.071293

*** STREAM 'WEST ' ***

- Temperature 40.550 CENT - Pressure 172.8000 BAR

GLOBAL STREAM

* No *	Component	* - Mols	- Mol fr-	- Weight	- Wt fr-
2	HYDROGEN SULPHIDE	0.370	0.003700	12.610	0.004723
3	NITROGEN	6.130	0.061300	171.701	0.064312
4	CARBON DIOXIDE	15.560	0.155600	684.796	0.256495
5	METHANE	65.850	0.658500	1056.234	0.395619
6	ETHANE	4.400	0.044000	132.308	0.049557
7	PROPANE	2.180	0.021800	96.138	0.036009
8	ISOBUTANE	0.560	0.005600	32.547	0.012191
9	BUTANE	0.950	0.009500	55.214	0.020681
10	C5+	1.030	0.010300	74.129	0.027766
11	C6+	0.660	0.006600	55.948	0.020956
12	C7+	0.650	0.006500	63.875	0.023925
13	C8+	0.530	0.005300	59.408	0.022252
14	C9+	0.370	0.003700	46.439	0.017394
15	C10+	0.260	0.002600	36.470	0.013660
16	C11+	0.170	0.001700	26.442	0.009904
17	C12+	0.330	0.003300	65.564	0.024558

* TOTAL *		100.000	1.000000	2669.823	1.000000
		KMOL		KG	

* XPSIM, Vers. 1.07 *
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* Proj/Problem "NORTH AFRICA RESERVOIR" *

*** UNIT 1 - 'PHASE1' - 'PHASE ENVELOPE'
--- Feed Streams ---
'WEST'
- 3 Phase VLE (LHLW separation) equilibrium applied

1) * CALCULATION PARAMETERS *

Maximum envelope pressure 400.0000 BAR
Maximum bubble point pressure 400.0000 BAR

Calculation starts from DEW POINT curve
Initial pressure 9.8067 BAR
Maximum temperature stepsize 5.0000 CENT
Maximum pressure stepsize 4.8344 BAR

1) * PRINTOUT OPTION *

Envelope plot requested

*** UNIT 2 - 'CYLIND' - 'FLASH'
--- Feed Streams ---
'WEST'
- 3 Phase VLE (LHLW separation) equilibrium applied

1) * CALCULATION TYPE *
Isothermal/Isobaric flash
Temperature is fixed at 34.400 CENT
Pressure is fixed at 37.6000 BAR

*** UNIT 3 - 'STD' - 'FLASH'
--- Feed Streams ---
from unit 2 - CYLIND 'LIQA0065'
- 3 Phase VLE (LHLW separation) equilibrium applied

1) * CALCULATION TYPE *
Isothermal/Isobaric flash
Temperature is fixed at 15.560 CENT
Pressure is fixed at 1.0130 BAR

extended Process SIMulation ...*
- INPUT
- UNIT 1 - PHASE1 -

Job "PVT05A"
Date MAR 04, 2009

*** UNIT 4 - 'RECOMB' - 'FLASH' ***
--- Feed Streams --- - Product Streams -
from unit 2 - CYLIND 'GASA0065' 'STDGAS'
from unit 3 - STD 'SGS03'
- 3 Phase VLE (LHLW separation) equilibrium applied

- 1) * CALCULATION TYPE *
Isothermal/Isobaric flash
Temperature is fixed at 15.560 CENT
Pressure is fixed at 1.0130 BAR

*** UNIT 5 - 'DEW' - 'FLASH' ***
--- Feed Streams --- - Product Streams -
'WEST' 'DEW'
- 3 Phase VLE (LHLW separation) equilibrium applied

- 1) * CALCULATION TYPE *
Dew-point
Temperature is fixed at 140.550 CENT
Pressure is variable, estimate 250.0000 BAR

V) **** STREAM / UNITS DICTIONARY ****

1) STREAMS-UNITS CROSS-REFERENCE TABLE

No	Stream Id	Description/Service	Product of Unit	Feed of Unit(s)		
1	WEST			1 PHASE1	2 CYLIND	5 DEW
2	GASA0065		2 CYLIND	4 RECOMB		
3	LIQA0065		2 CYLIND	3 STD		
4	SGS03		3 STD	4 RECOMB		
5	STDLIQ		3 STD			
6	STDGAS		4 RECOMB			
7	DEW		5 DEW			

2) UNITS-STREAMS CROSS-REFERENCE TABLE

No	Unit	Unit Description	Feed Streams		Product Streams	
1	PHASE1		1	WEST		
2	CYLIND		1	WEST	2 GASA0065 3 LIQA0065	
3	STD		3	LIQA0065	4 SGS03 5 STDLIQ	
4	RECOMB		2 4	GASA0065 SGS03	6 STDGAS	
5	DEW		1	WEST	7 DEW	

VI) **** RUN STATISTICS ****

* No of streams	7
process operations	5
topping structures	0
control operations	0
set operations	0

The calculation is NOT a RECYCLE PROBLEM
Calculation sequence defined by input sequence

*** VAPOR/LIQUID EQUILIBRIUM CURVES - SET NO. 1 ***

VLE calculation for stream 'WEST'
3 PHASE flash calculation used
Key component 1 1 H2O WATER
Calculation model 1

No. of VLE curves 1

1) - FLASH AT GIVEN T/P -
Initial temperature 140.500 CENT - End temperature 140.500 CENT
Initial pressure 100.0000 BAR - End pressure 300.0000 BAR
No. of points 11

*** PVT ANALYSIS TRACE FOLLOWS ***

EXPERIMENT NO.	1	CCE1	SIMULATED
* 'BUBBLE ' POINT CALCULATION FAILED *			
EXPERIMENT NO.	2	CVD1	SIMULATED
EXPERIMENT NO.	3	SEP1	SIMULATED

*** PVT ANALYSIS ***

1) *** EXPERIMENT NO. 1 CCE1 ***
COSTANT COMPOSITION EXPANSION

Temperature 140.000 CENT

* FLUID COMPOSITION *		
* No	* Component	* Mols
2	HYDROGEN SULPHIDE	0.003700
3	NITROGEN	0.061300
4	CARBON DIOXIDE	0.155600
5	METHANE	0.658500
6	ETHANE	0.044000
7	PROPANE	0.021800
8	ISOBUTANE	0.005600
9	BUTANE	0.009500
10	C5+	0.010300
11	C6+	0.006600
12	C7+	0.006500
13	C8+	0.005300
14	C9+	0.003700
15	C10+	0.002600
16	C11+	0.001700
17	C12+	0.003300

Saturation pressure, BAR -Exp 254.4000 -Calc 252.9194
Saturation pressure error(abs) 1.4806 (rel) 0.0058
Saturation density 1000.0000 KG/CUMT

* EXPERIMENTAL AND CALCULATED POINTS *

POINT NO. 1					
Pressure 310.300 BAR					
	- Experimental	-	- Calculated	-	- Err(abs) -
Relative volume	0.8538		0.8528		0.001002
Liquid volume frac.	0.000000		0.000000		0.000000
Gas Z factor	0.980000		0.965978		0.014022
					0.014308
POINT NO. 2					
Pressure 303.400 BAR					
	- Experimental	-	- Calculated	-	- Err(abs) -
Relative volume	0.8681		0.8674		0.000657
Liquid volume frac.	0.000000		0.000000		0.000000
Gas Z factor	0.974000		0.960718		0.013282
					0.013636
POINT NO. 3					
Pressure 298.200 BAR					
	- Experimental	-	- Calculated	-	- Err(abs) -
Relative volume	0.8792		0.8790		0.000171
Liquid volume frac.	0.000000		0.000000		0.000000
Gas Z factor	0.970000		0.956864		0.013136
					0.013542

* XPSIM, Vers. 1.07 *		*... extended Process SIMulation ...*		* Page 0032 *	
* Cust/User "STAFF " *		PVT ANALYSIS -		* Job "PVT05A " *	
* Proj/Problem "NORTH AFRICA RESERVOIR " *		- EXPERIMENT 1 - CCE1 -		* Date MAR 04, 2009 *	
POINT NO.	4				
Pressure	293.500 BAR				
	- Experimental -	- Calculated -	- Err(abs) -	- Err(rel) -	
Relative volume	0.8900	0.8899	0.000068	0.000077	
Liquid volume frac.	0.000000	0.000000	0.000000	0.000000	
Gas Z factor	0.966000	0.953464	0.012536	0.012978	
POINT NO.	5				
Pressure	288.700 BAR				
	- Experimental -	- Calculated -	- Err(abs) -	- Err(rel) -	
Relative volume	0.9014	0.9015	0.000113	0.000126	
Liquid volume frac.	0.000000	0.000000	0.000000	0.000000	
Gas Z factor	0.962000	0.950076	0.011924	0.012395	
POINT NO.	6				
Pressure	282.600 BAR				
	- Experimental -	- Calculated -	- Err(abs) -	- Err(rel) -	
Relative volume	0.9167	0.9169	0.000225	0.000245	
Liquid volume frac.	0.000000	0.000000	0.000000	0.000000	
Gas Z factor	0.958000	0.945900	0.012100	0.012631	
POINT NO.	7				
Pressure	275.900 BAR				
	- Experimental -	- Calculated -	- Err(abs) -	- Err(rel) -	
Relative volume	0.9347	0.9348	0.000109	0.000117	
Liquid volume frac.	0.000000	0.000000	0.000000	0.000000	
Gas Z factor	0.954000	0.941487	0.012513	0.013117	
POINT NO.	8				
Pressure	270.300 BAR				
	- Experimental -	- Calculated -	- Err(abs) -	- Err(rel) -	
Relative volume	0.9503	0.9506	0.000286	0.000301	
Liquid volume frac.	0.000000	0.000000	0.000000	0.000000	
Gas Z factor	0.950000	0.937944	0.012056	0.012691	
POINT NO.	9				
Pressure	262.000 BAR				
	- Experimental -	- Calculated -	- Err(abs) -	- Err(rel) -	
Relative volume	0.9755	0.9755	0.000022	0.000023	
Liquid volume frac.	0.000000	0.000000	0.000000	0.000000	
Gas Z factor	0.945000	0.932950	0.012050	0.012752	
POINT NO.	10				
Pressure	254.400 BAR				
	- Experimental -	- Calculated -	- Err(abs) -	- Err(rel) -	
Relative volume	1.0000	1.0000	0.000000	0.000000	
Liquid volume frac.	0.000000	0.000000	0.000000	0.000000	
Gas Z factor	0.941000	0.928660	0.012340	0.013114	
POINT NO.	11				
Pressure	248.600 BAR				
	- Experimental -	- Calculated -	- Err(abs) -	- Err(rel) -	
Relative volume	1.0203	1.0204	0.000144	0.000141	
Liquid volume frac.	0.000560	0.000660	0.000100	0.178767	
POINT NO.	12				
Pressure	240.600 BAR				
	- Experimental -	- Calculated -	- Err(abs) -	- Err(rel) -	
Relative volume	1.0500	1.0509	0.000945	0.000900	
Liquid volume frac.	0.001390	0.001941	0.000551	0.396074	

* XPSIM, Vers. 1.07 *					* Page 0033 *				
* Cust/User "STAFF " *					* Job "PVT05A " *				
* Proj/Problem "NORTH AFRICA RESERVOIR " *					* Date MAR 04, 2009 *				
POINT NO. 13									
Pressure	228.900 BAR								
	- Experimental	-	-	Calculated	-	-	Err(abs)	-	Err(rel)
Relative volume	1.0979			1.0999			0.002039		0.001857
Liquid volume frac.	0.002680			0.003473			0.000793		0.295908
POINT NO. 14									
Pressure	215.500 BAR								
	- Experimental	-	-	Calculated	-	-	Err(abs)	-	Err(rel)
Relative volume	1.1597			1.1636			0.003926		0.003385
Liquid volume frac.	0.004290			0.004883			0.000593		0.138288
POINT NO. 15									
Pressure	191.000 BAR								
	- Experimental	-	-	Calculated	-	-	Err(abs)	-	Err(rel)
Relative volume	1.2982			1.3067			0.008474		0.006528
Liquid volume frac.	0.007500			0.006595			0.000905		0.120692
POINT NO. 16									
Pressure	160.600 BAR								
	- Experimental	-	-	Calculated	-	-	Err(abs)	-	Err(rel)
Relative volume	1.5404			1.5529			0.012527		0.008133
Liquid volume frac.	0.010790			0.007366			0.003424		0.317350
POINT NO. 17									
Pressure	123.100 BAR								
	- Experimental	-	-	Calculated	-	-	Err(abs)	-	Err(rel)
Relative volume	2.0188			2.0410			0.022248		0.011020
Liquid volume frac.	0.013420			0.006665			0.006755		0.503331
POINT NO. 18									
Pressure	100.800 BAR								
	- Experimental	-	-	Calculated	-	-	Err(abs)	-	Err(rel)
Relative volume	2.4827			2.5141			0.031423		0.012657
Liquid volume frac.	0.014230			0.005608			0.008622		0.605882
POINT NO. 19									
Pressure	80.600 BAR								
	- Experimental	-	-	Calculated	-	-	Err(abs)	-	Err(rel)
Relative volume	3.1345			3.1769			0.042374		0.013519
Liquid volume frac.	0.014390			0.004385			0.010005		0.695280
POINT NO. 20									
Pressure	59.000 BAR								
	- Experimental	-	-	Calculated	-	-	Err(abs)	-	Err(rel)
Relative volume	4.3389			4.3995			0.060574		0.013961
Liquid volume frac.	0.013920			0.002942			0.010978		0.788648

2) *** EXPERIMENT NO. 2 CVD1 ***
COSTANT VOLUME DEPLETION

Temperature 140.000 CENT

* FLUID COMPOSITION *		
* No	* Component	* - Mols -
2	HYDROGEN SULPHIDE	0.003700
3	NITROGEN	0.061300
4	CARBON DIOXIDE	0.155600
5	METHANE	0.658500
6	ETHANE	0.044000
7	PROPANE	0.021800
8	ISOBUTANE	0.005600
9	BUTANE	0.009500
10	C5+	0.010300
11	C6+	0.006600
12	C7+	0.006500
13	C8+	0.005300
14	C9+	0.003700
15	C10+	0.002600
16	C11+	0.001700
17	C12+	0.003300

Saturation pressure, BAR - Exp 254.4000 - Calc 35.9841
Saturation pressure error(abs) 218.4159 (rel) 0.8586

* EXPERIMENTAL AND CALCULATED POINTS *

POINT NO.	1				
Pressure	254.400 BAR				
	- Experimental -	-	Calculated -	- Err(abs) -	- Err(rel) -
Removed vap. frac.	0.0000		0.0000	0.000000	0.000000
Liquid volume frac.	0.000000		0.000000	0.000000	0.000000
Gas Z factor	0.941000		0.928660	0.012340	0.013114

* XPSIM, Vers. 1.07 *

* Cust/User "STAFF" *
* Proj/Problem "NORTH AFRICA RESERVOIR" *
* CALCULATED VAPOR COMPOSITION *

* No * Component * - Mols -
2 HYDROGEN SULPHIDE 0.003700
3 NITROGEN 0.061300
4 CARBON DIOXIDE 0.155600
5 METHANE 0.658500
6 ETHANE 0.044000
7 PROPANE 0.021800
8 ISOBUTANE 0.005600
9 BUTANE 0.009500
10 C5+ 0.010300
11 C6+ 0.006600
12 C7+ 0.006500
13 C8+ 0.005300
14 C9+ 0.003700
15 C10+ 0.002600
16 C11+ 0.001700
17 C12+ 0.003300

POINT NO. 2
Pressure 247.900 BAR
- Experimental - - Calculated - - Err(abs) - - Err(rel) -
Removed vap. frac. 0.0208 0.0225 0.001746 0.083953
Liquid volume frac. 0.000420 0.000824 0.000404 0.961758
Gas Z factor 0.919000 0.925786 0.006786 0.007384
* CALCULATED VAPOR COMPOSITION *
* No * Component * - Mols -
2 HYDROGEN SULPHIDE 0.003700
3 NITROGEN 0.061333
4 CARBON DIOXIDE 0.155637
5 METHANE 0.658776
6 ETHANE 0.044002
7 PROPANE 0.021794
8 ISOBUTANE 0.005597
9 BUTANE 0.009493
10 C5+ 0.010285
11 C6+ 0.006585
12 C7+ 0.006478
13 C8+ 0.005273
14 C9+ 0.003673
15 C10+ 0.002575
16 C11+ 0.001668
17 C12+ 0.003132

POINT NO. 3
Pressure 237.200 BAR
- Experimental - - Calculated - - Err(abs) - - Err(rel) -
Removed vap. frac. 0.0590 0.0606 0.001633 0.027677
Liquid volume frac. 0.001670 0.002429 0.000759 0.454588
Gas Z factor 0.913000 0.921814 0.008814 0.009654

extended Process SIMulation ...*
- PVT ANALYSIS -
- EXPERIMENT 2 - CVD1 -
* Page 0035 *
* Job "PVT05A" *
* Date MAR 04, 2009 *

* XPSIM, Vers. 1.07 *

* Cust/User "STAFF" *
* Proj/Problem "NORTH AFRICA RESERVOIR" *
* CALCULATED VAPOR COMPOSITION *

* No * Component * - Mols -
2 HYDROGEN SULPHIDE 0.003699
3 NITROGEN 0.061401
4 CARBON DIOXIDE 0.155715
5 METHANE 0.659349
6 ETHANE 0.044006
7 PROPANE 0.021781
8 ISOBUTANE 0.005590
9 BUTANE 0.009478
10 C5+ 0.010254
11 C6+ 0.006553
12 C7+ 0.006432
13 C8+ 0.005216
14 C9+ 0.003614
15 C10+ 0.002522
16 C11+ 0.001600
17 C12+ 0.002790

POINT NO. 4
Pressure 212.700 BAR
- Experimental - - Calculated - - Err(abs) - - Err(rel) -
Removed vap. frac. 0.1483 0.1508 0.002495 0.016827
Liquid volume frac. 0.004320 0.005234 0.000914 0.211680
Gas Z factor 0.902000 0.915054 0.013054 0.014473
* CALCULATED VAPOR COMPOSITION *
* No * Component * - Mols -
2 HYDROGEN SULPHIDE 0.003698
3 NITROGEN 0.061541
4 CARBON DIOXIDE 0.155894
5 METHANE 0.660563
6 ETHANE 0.044021
7 PROPANE 0.021758
8 ISOBUTANE 0.005577
9 BUTANE 0.009448
10 C5+ 0.010185
11 C6+ 0.006481
12 C7+ 0.006326
13 C8+ 0.005083
14 C9+ 0.003477
15 C10+ 0.002395
16 C11+ 0.001439
17 C12+ 0.002114

POINT NO. 5
Pressure 185.200 BAR
- Experimental - - Calculated - - Err(abs) - - Err(rel) -
Removed vap. frac. 0.2542 0.2560 0.001778 0.006994
Liquid volume frac. 0.007710 0.007423 0.000287 0.037256
Gas Z factor 0.893000 0.911643 0.018643 0.020877

extended Process SIMulation ...*
- PVT ANALYSIS -
- EXPERIMENT 2 - CVD1 -
* Page 0036 *
* Job "PVT05A" *
* Date MAR 04, 2009 *

* XPSIM, Vers. 1.07 *

* Cust/User "STAFF" *
* Proj/Problem "NORTH AFRICA RESERVOIR" *
* CALCULATED VAPOR COMPOSITION *

* No * Component * - Mols -
2 HYDROGEN SULPHIDE 0.003698
3 NITROGEN 0.061673
4 CARBON DIOXIDE 0.156097
5 METHANE 0.661763
6 ETHANE 0.044048
7 PROPANE 0.021742
8 ISOBUTANE 0.005565
9 BUTANE 0.009420
10 C5+ 0.010117
11 C6+ 0.006404
12 C7+ 0.006207
13 C8+ 0.004930
14 C9+ 0.003316
15 C10+ 0.002244
16 C11+ 0.001256
17 C12+ 0.001519

POINT NO. 6
Pressure 153.100 BAR
- Experimental - - Calculated - - Err(abs) - - Err(rel) -
Removed vap. frac. 0.3834 0.3829 0.000499 0.001301
Liquid volume frac. 0.010090 0.008687 0.001403 0.139058
Gas Z factor 0.885000 0.913564 0.028564 0.032275

* CALCULATED VAPOR COMPOSITION *
* No * Component * - Mols -
2 HYDROGEN SULPHIDE 0.003701
3 NITROGEN 0.061789
4 CARBON DIOXIDE 0.156330
5 METHANE 0.662916
6 ETHANE 0.044094
7 PROPANE 0.021740
8 ISOBUTANE 0.005557
9 BUTANE 0.009399
10 C5+ 0.010054
11 C6+ 0.006326
12 C7+ 0.006076
13 C8+ 0.004756
14 C9+ 0.003126
15 C10+ 0.002063
16 C11+ 0.001050
17 C12+ 0.001021

POINT NO. 7
Pressure 118.200 BAR
- Experimental - - Calculated - - Err(abs) - - Err(rel) -
Removed vap. frac. 0.5285 0.5231 0.005353 0.010128
Liquid volume frac. 0.011670 0.008799 0.002871 0.246036
Gas Z factor 0.881000 0.922751 0.041751 0.047391

extended Process SIMulation ...*
- PVT ANALYSIS -
- EXPERIMENT 2 - CVD1 -
* Page 0037 *
* Job "PVT05A" *
* Date MAR 04, 2009 *

* XPSIM, Vers. 1.07 *

* Cust/User "STAFF" *
* Proj/Problem "NORTH AFRICA RESERVOIR" *
* CALCULATED VAPOR COMPOSITION *

* No * Component * - Mols -
2 HYDROGEN SULPHIDE 0.003708
3 NITROGEN 0.061852
4 CARBON DIOXIDE 0.156562
5 METHANE 0.663733
6 ETHANE 0.044164
7 PROPANE 0.021769
8 ISOBUTANE 0.005560
9 BUTANE 0.009401
10 C5+ 0.010028
11 C6+ 0.006277
12 C7+ 0.005972
13 C8+ 0.004603
14 C9+ 0.002947
15 C10+ 0.001886
16 C11+ 0.000860
17 C12+ 0.000677

POINT NO. 8
Pressure 79.800 BAR
- Experimental - - Calculated - - Err(abs) - - Err(rel) -
Removed vap. frac. 0.6878 0.6766 0.011168 0.016237
Liquid volume frac. 0.011430 0.007437 0.003993 0.349329
Gas Z factor 0.882000 0.940470 0.058470 0.066293

* CALCULATED VAPOR COMPOSITION *
* No * Component * - Mols -
2 HYDROGEN SULPHIDE 0.003719
3 NITROGEN 0.061806
4 CARBON DIOXIDE 0.156717
5 METHANE 0.663706
6 ETHANE 0.044258
7 PROPANE 0.021853
8 ISOBUTANE 0.005588
9 BUTANE 0.009456
10 C5+ 0.010106
11 C6+ 0.006325
12 C7+ 0.005996
13 C8+ 0.004584
14 C9+ 0.002877
15 C10+ 0.001790
16 C11+ 0.000738
17 C12+ 0.000481

POINT NO. 9
Pressure 38.200 BAR
- Experimental - - Calculated - - Err(abs) - - Err(rel) -
Removed vap. frac. 0.8484 0.8384 0.010002 0.011789
Liquid volume frac. 0.010800 0.004324 0.006476 0.599659
Gas Z factor 0.893000 0.967196 0.074196 0.083086

extended Process SIMulation ...*
- PVT ANALYSIS -
- EXPERIMENT 2 - CVD1 -
* Job "PVT05A" *
* Date MAR 04, 2009 *

```
* XPSIM, Vers. 1.07 *
* Cust/User "STAFF" " *
* Proj/Problem "NORTH AFRICA RESERVOIR " *
* CALCULATED VAPOR COMPOSITION *
* No * Component * - Moles -
2 HYDROGEN SULPHIDE 0.003732
3 NITROGEN 0.061485
4 CARBON DIOXIDE 0.156465
5 METHANE 0.661173
6 ETHANE 0.044319
7 PROPANE 0.022014
8 ISOBUTANE 0.005665
9 BUTANE 0.009620
10 C5+ 0.010451
11 C6+ 0.006679
12 C7+ 0.006493
13 C8+ 0.005130
14 C9+ 0.003326
15 C10+ 0.002099
16 C11+ 0.000857
17 C12+ 0.000493

extended Process SIMulation ...*
- PVT ANALYSIS -
- EXPERIMENT 2 - CVD1 -

* Job "PVT05A" " *
* Date MAR 04, 2009 *
```

3) *** EXPERIMENT NO. 3 SEP1 ***
TEST SEPARATORS

* FLUID COMPOSITION *		
* No	* Component	* - Mols -
2	HYDROGEN SULPHIDE	0.003700
3	NITROGEN	0.061300
4	CARBON DIOXIDE	0.155600
5	METHANE	0.658500
6	ETHANE	0.044000
7	PROPANE	0.021800
8	ISOBUTANE	0.005600
9	BUTANE	0.009500
10	C5+	0.010300
11	C6+	0.006600
12	C7+	0.006500
13	C8+	0.005300
14	C9+	0.003700
15	C10+	0.002600
16	C11+	0.001700
17	C12+	0.003300

Oil density,KG/CUMT	759.000	719.747	39.252878	0.051717
Gas gravity	1.26800	0.81570	0.452303	0.356706
Gas/oil ratio,NM3/M3	1000.0000	3668.7276	2668.727553	2.668728

* EXPERIMENTAL AND CALCULATED POINTS *

POINT NO.	1
Temperature, K	307.600
Pressure, BAR	37.600
Oil density, KG/CUMT	719.747
Gas gravity	0.815697
Gas/oil ratio,NM3/M3	3668.7276

* CALCULATED VAPOR COMPOSITION *		
* No	* Component	* - Mols -
2	HYDROGEN SULPHIDE	0.003694
3	NITROGEN	0.064104
4	CARBON DIOXIDE	0.160289
5	METHANE	0.685785
6	ETHANE	0.044524
7	PROPANE	0.020659
8	ISOBUTANE	0.004797
9	BUTANE	0.007582
10	C5+	0.005527
11	C6+	0.001872
12	C7+	0.000822
13	C8+	0.000260
14	C9+	0.000065
15	C10+	0.000019
16	C11+	0.000002
17	C12+	0.000000

* XPSIM, Vers. 1.07 *			*... extended Process SIMulation ...*			* Page 0041 *		
* Cust/User "STAFF			- PVT ANALYSIS -			* Job "PVT05A		
* Proj/Problem "NORTH AFRICA RESERVOIR			- EXPERIMENT 3 - SEP1 -			* Date MAR 04, 2009 *		
			* LIQUID Composition *					
			Experimental			Calculated		
			Err(abs)					
			Mols			Mols		
			-			-		
* No	* Component							
2	HYDROGEN SULPHIDE	0.002000	0.003827	0.001827				
3	NITROGEN	0.002400	0.002712	0.000312				
4	CARBON DIOXIDE	0.049495	0.057640	0.008145				
5	METHANE	0.102390	0.088430	0.013960				
6	ETHANE	0.031497	0.033051	0.001554				
7	PROPANE	0.044196	0.045634	0.001438				
8	ISOBUTANE	0.019998	0.022381	0.002383				
9	BUTANE	0.060094	0.049583	0.010511				
10	C5+	0.115888	0.110032	0.005857				
11	C6+	0.105389	0.105383	0.000006				
12	C7+	0.117388	0.125123	0.007735				
13	C8+	0.106289	0.110598	0.004309				
14	C9+	0.078992	0.079645	0.000653				
15	C10+	0.055994	0.056530	0.000536				
16	C11+	0.036696	0.037181	0.000485				
17	C12+	0.071293	0.072243	0.000950				

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* XPSIM, Vers. 1.07 *
* Cust/User "STAFF" *
* Proj/Problem "NORTH AFRICA RESERVOIR" *
* Date MAR 04, 2009 *

*... extended Process SIMulation ...*
* Job "PVT05A" *
* PVT ANALYSIS -
* EXPERIMENT 3 - SEP1 -

*** SIMULATION TRACE FOLLOWS ***

... New/output binary file 'PVT05A' initialized ...

. Stream 'WEST' flashed .
. Feed/Recycle streams processed .
*** CALCULATION TERMINATED, PRESSURE LIMIT REACHED ***
*** CALCULATION TERMINATED, TOO MANY INDEPENDENT VARIABLES CHANGES ***
Unit op 1 - 'PHASE1' calculated
Unit op 2 - 'CYLIND' calculated
Unit op 3 - 'STD' calculated
Unit op 4 - 'RECOMB' calculated
Unit op 5 - 'DEW' calculated

*** SOLUTION OBTAINED ***

... Base case results loaded on binary file 'PVT05A' ...
```

```
* XPSIM, Vers. 1.07 *
* Cust/User "STAFF"
* Proj/Problem "NORTH AFRICA RESERVOIR"

*... extended Process SIMulation ...*
* - VLE/BVLE CALCULATIONS -
* - STREAM WEST - VLE CURVE -

* Job "PVT05A"
* Date MAR 04, 2009

* ISOTHERMAL/ISOBARIC FLASH CURVE - STREAM WEST
Initial temperature 140.500 °C - End temperature 140.500 °C
Initial pressure 100.0000 bar - End pressure 300.0000 bar
No. of points 11
1) Temp 140.500 °C Pres 100.0000 bar V/F ratio 0.98950
2) Temp 140.500 °C Pres 120.0000 bar V/F ratio 0.98937
3) Temp 140.500 °C Pres 140.0000 bar V/F ratio 0.98971
4) Temp 140.500 °C Pres 160.0000 bar V/F ratio 0.99048
5) Temp 140.500 °C Pres 180.0000 bar V/F ratio 0.99170
6) Temp 140.500 °C Pres 200.0000 bar V/F ratio 0.99337
7) Temp 140.500 °C Pres 220.0000 bar V/F ratio 0.99551
8) Temp 140.500 °C Pres 240.0000 bar V/F ratio 0.99817
9) * PHASE CHANGE * Temp 140.500 °C Pres 251.8041 bar V/F ratio 1.00000
10) Temp 140.500 °C Pres 260.0000 bar V/F ratio 1.00000
11) Temp 140.500 °C Pres 280.0000 bar V/F ratio 1.00000
12) Temp 140.500 °C Pres 300.0000 bar V/F ratio 1.00000
```

```

* XPSIM, Vers. 1.07 *
* Cust/User "STAFF" " *
* Proj/Problem "NORTH AFRICA RESERVOIR " *
* Date MAR 04, 2009 *

*** UNIT 1 - 'PHASE1 ' - ' PHASE ENVELOPE ' ***
--- Feed Streams --- - Product Streams -
'WEST '

1) * FEEDS *

Stream WEST
Temperature, °C 40.55
Pressure, bar 172.800
Total rate, kmol/h 100.00
Vapor, kmol/h 94.411
Liquid, kmol/h 5.5893

2) * METHOD *
K-values calculation method 'SRK '

3) *** DEW POINT EQUILIBRIUM CURVE ***

Cricondentherm Point - Temp 184.764 °C - Pres 78.7096 bar
Cricondenbar Point - *** NOT DETERMINED ***
Critical Point - *** NOT DETERMINED ***
Max Temp/Pres Point - Temp 184.764 °C - Pres 399.3893 bar

1) - DEW PT - PRES 9.8067 bar - TEMP 153.265 °C
2) - DEW PT - PRES 10.7873 bar - TEMP 155.203 °C
3) - DEW PT - PRES 11.8661 bar - TEMP 157.136 °C
4) - DEW PT - PRES 13.0527 bar - TEMP 159.061 °C
5) - DEW PT - PRES 14.3579 bar - TEMP 160.974 °C
6) - DEW PT - PRES 15.7937 bar - TEMP 162.872 °C
7) - DEW PT - PRES 17.3731 bar - TEMP 164.749 °C
8) - DEW PT - PRES 19.1104 bar - TEMP 166.600 °C
9) - DEW PT - PRES 21.0215 bar - TEMP 168.419 °C
10) - DEW PT - PRES 23.1236 bar - TEMP 170.201 °C
11) - DEW PT - PRES 25.4360 bar - TEMP 171.936 °C
12) - DEW PT - PRES 27.9796 bar - TEMP 173.618 °C
13) - DEW PT - PRES 30.7775 bar - TEMP 175.237 °C
14) - DEW PT - PRES 33.8553 bar - TEMP 176.781 °C
15) - DEW PT - PRES 37.2408 bar - TEMP 178.240 °C
16) - DEW PT - PRES 40.9649 bar - TEMP 179.599 °C
17) - DEW PT - PRES 45.0614 bar - TEMP 180.844 °C
18) - DEW PT - PRES 49.5675 bar - TEMP 181.956 °C
19) - DEW PT - PRES 51.9285 bar - TEMP 182.446 °C
20) - DEW PT - PRES 54.4631 bar - TEMP 182.907 °C
21) - DEW PT - PRES 57.1213 bar - TEMP 183.326 °C
22) - DEW PT - PRES 59.9094 bar - TEMP 183.696 °C
23) - DEW PT - PRES 62.8335 bar - TEMP 184.019 °C
24) - DEW PT - PRES 65.2059 bar - TEMP 184.233 °C
25) - DEW PT - PRES 67.6679 bar - TEMP 184.413 °C
26) - DEW PT - PRES 70.2228 bar - TEMP 184.557 °C
27) - DEW PT - PRES 72.8742 bar - TEMP 184.664 °C
28) - DEW PT - PRES 75.6257 bar - TEMP 184.729 °C
29) - DEW PT - PRES 78.4811 bar - TEMP 184.764 °C
30) - DEW PT - PRES 78.7096 bar - TEMP MAX 184.764 °C
31) - DEW PT - PRES 81.4443 bar - TEMP 184.737 °C
32) - DEW PT - PRES 83.8267 bar - TEMP 184.690 °C
33) - DEW PT - PRES 86.2787 bar - TEMP 184.613 °C
34) - DEW PT - PRES 88.8025 bar - TEMP 184.506 °C
35) - DEW PT - PRES 91.4001 bar - TEMP 184.367 °C
36) - DEW PT - PRES 94.0736 bar - TEMP 184.195 °C
37) - DEW PT - PRES 96.8254 bar - TEMP 183.986 °C

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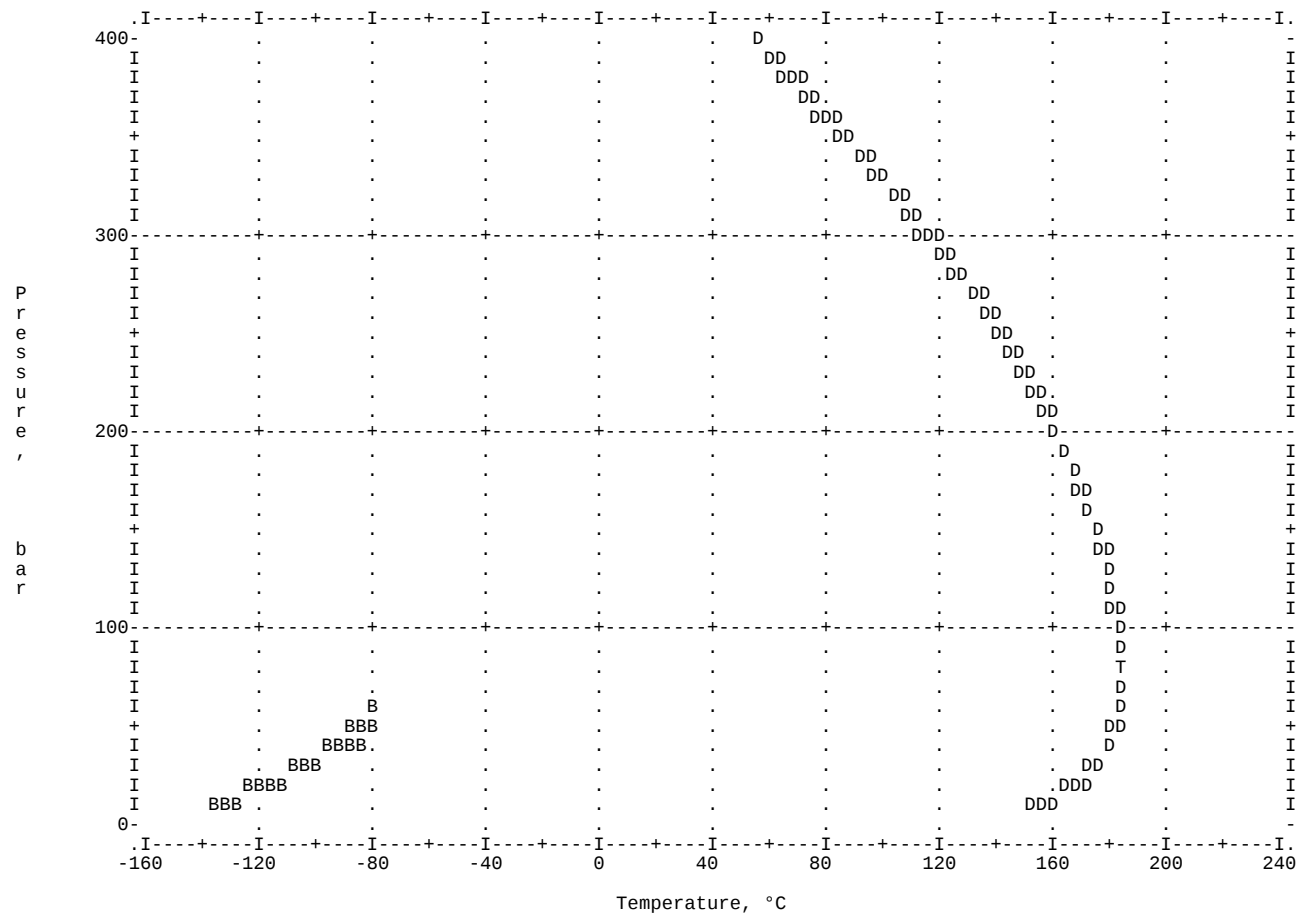
38)	-	DEW	PT - PRES	99.6577	bar	-	TEMP	183.744	°C
39)	-	DEW	PT - PRES	104.4921	bar	-	TEMP	183.261	°C
40)	-	DEW	PT - PRES	106.8819	bar	-	TEMP	182.992	°C
41)	-	DEW	PT - PRES	109.4437	bar	-	TEMP	182.682	°C
42)	-	DEW	PT - PRES	112.0668	bar	-	TEMP	182.344	°C
43)	-	DEW	PT - PRES	114.7528	bar	-	TEMP	181.974	°C
44)	-	DEW	PT - PRES	117.1450	bar	-	TEMP	181.626	°C
45)	-	DEW	PT - PRES	119.5872	bar	-	TEMP	181.254	°C
46)	-	DEW	PT - PRES	122.0802	bar	-	TEMP	180.856	°C
47)	-	DEW	PT - PRES	124.6252	bar	-	TEMP	180.431	°C
48)	-	DEW	PT - PRES	127.2233	bar	-	TEMP	179.979	°C
49)	-	DEW	PT - PRES	129.8755	bar	-	TEMP	179.502	°C
50)	-	DEW	PT - PRES	132.5830	bar	-	TEMP	178.974	°C
51)	-	DEW	PT - PRES	135.3470	bar	-	TEMP	178.450	°C
52)	-	DEW	PT - PRES	140.1814	bar	-	TEMP	177.460	°C
53)	-	DEW	PT - PRES	145.0158	bar	-	TEMP	176.412	°C
54)	-	DEW	PT - PRES	149.8502	bar	-	TEMP	175.309	°C
55)	-	DEW	PT - PRES	154.6846	bar	-	TEMP	174.152	°C
56)	-	DEW	PT - PRES	159.5189	bar	-	TEMP	172.942	°C
57)	-	DEW	PT - PRES	164.3533	bar	-	TEMP	171.680	°C
58)	-	DEW	PT - PRES	166.7530	bar	-	TEMP	171.035	°C
59)	-	DEW	PT - PRES	169.7050	bar	-	TEMP	170.225	°C
60)	-	DEW	PT - PRES	172.7092	bar	-	TEMP	169.381	°C
61)	-	DEW	PT - PRES	175.7666	bar	-	TEMP	168.503	°C
62)	-	DEW	PT - PRES	178.8781	bar	-	TEMP	167.589	°C
63)	-	DEW	PT - PRES	182.0447	bar	-	TEMP	166.638	°C
64)	-	DEW	PT - PRES	185.2673	bar	-	TEMP	165.649	°C
65)	-	DEW	PT - PRES	187.6690	bar	-	TEMP	164.895	°C
66)	-	DEW	PT - PRES	190.1017	bar	-	TEMP	164.126	°C
67)	-	DEW	PT - PRES	194.9361	bar	-	TEMP	162.555	°C
68)	-	DEW	PT - PRES	199.7705	bar	-	TEMP	160.937	°C
69)	-	DEW	PT - PRES	202.1732	bar	-	TEMP	160.111	°C
70)	-	DEW	PT - PRES	204.7278	bar	-	TEMP	159.228	°C
71)	-	DEW	PT - PRES	207.1309	bar	-	TEMP	158.378	°C
72)	-	DEW	PT - PRES	209.5622	bar	-	TEMP	157.514	°C
73)	-	DEW	PT - PRES	214.3966	bar	-	TEMP	155.752	°C
74)	-	DEW	PT - PRES	216.8003	bar	-	TEMP	154.856	°C
75)	-	DEW	PT - PRES	219.2867	bar	-	TEMP	153.923	°C
76)	-	DEW	PT - PRES	221.6907	bar	-	TEMP	153.002	°C
77)	-	DEW	PT - PRES	224.1211	bar	-	TEMP	152.066	°C
78)	-	DEW	PT - PRES	228.9555	bar	-	TEMP	150.161	°C
79)	-	DEW	PT - PRES	233.7899	bar	-	TEMP	148.208	°C
80)	-	DEW	PT - PRES	238.6243	bar	-	TEMP	146.151	°C
81)	-	DEW	PT - PRES	243.4586	bar	-	TEMP	144.157	°C
82)	-	DEW	PT - PRES	248.2930	bar	-	TEMP	142.058	°C
83)	-	DEW	PT - PRES	253.1274	bar	-	TEMP	139.907	°C
84)	-	DEW	PT - PRES	255.5332	bar	-	TEMP	138.816	°C
85)	-	DEW	PT - PRES	258.2745	bar	-	TEMP	137.563	°C
86)	-	DEW	PT - PRES	260.6804	bar	-	TEMP	136.444	°C
87)	-	DEW	PT - PRES	263.1088	bar	-	TEMP	135.309	°C
88)	-	DEW	PT - PRES	265.5150	bar	-	TEMP	134.164	°C
89)	-	DEW	PT - PRES	267.9432	bar	-	TEMP	133.003	°C
90)	-	DEW	PT - PRES	272.7776	bar	-	TEMP	130.644	°C
91)	-	DEW	PT - PRES	275.1842	bar	-	TEMP	129.449	°C
92)	-	DEW	PT - PRES	277.6556	bar	-	TEMP	128.213	°C
93)	-	DEW	PT - PRES	282.4900	bar	-	TEMP	125.749	°C
94)	-	DEW	PT - PRES	287.3244	bar	-	TEMP	123.178	°C
95)	-	DEW	PT - PRES	292.1588	bar	-	TEMP	120.662	°C
96)	-	DEW	PT - PRES	296.9932	bar	-	TEMP	118.040	°C
97)	-	DEW	PT - PRES	299.4006	bar	-	TEMP	116.710	°C
98)	-	DEW	PT - PRES	301.9959	bar	-	TEMP	115.290	°C
99)	-	DEW	PT - PRES	304.6136	bar	-	TEMP	113.796	°C

100)	-	DEW	PT - PRES	307.0213 bar	- TEMP	112.427 °C
101)	-	DEW	PT - PRES	309.4480 bar	- TEMP	111.041 °C
102)	-	DEW	PT - PRES	311.8939 bar	- TEMP	109.622 °C
103)	-	DEW	PT - PRES	314.3018 bar	- TEMP	108.215 °C
104)	-	DEW	PT - PRES	316.7283 bar	- TEMP	106.794 °C
105)	-	DEW	PT - PRES	319.1735 bar	- TEMP	105.336 °C
106)	-	DEW	PT - PRES	324.0079 bar	- TEMP	102.429 °C
107)	-	DEW	PT - PRES	328.8423 bar	- TEMP	99.482 °C
108)	-	DEW	PT - PRES	331.2506 bar	- TEMP	97.996 °C
109)	-	DEW	PT - PRES	333.7499 bar	- TEMP	96.455 °C
110)	-	DEW	PT - PRES	336.2680 bar	- TEMP	94.877 °C
111)	-	DEW	PT - PRES	338.6765 bar	- TEMP	93.366 °C
112)	-	DEW	PT - PRES	341.1023 bar	- TEMP	91.841 °C
113)	-	DEW	PT - PRES	343.5455 bar	- TEMP	90.292 °C
114)	-	DEW	PT - PRES	348.3799 bar	- TEMP	87.216 °C
115)	-	DEW	PT - PRES	353.2143 bar	- TEMP	84.094 °C
116)	-	DEW	PT - PRES	355.6233 bar	- TEMP	82.584 °C
117)	-	DEW	PT - PRES	358.1167 bar	- TEMP	80.990 °C
118)	-	DEW	PT - PRES	360.5258 bar	- TEMP	79.448 °C
119)	-	DEW	PT - PRES	362.9511 bar	- TEMP	77.906 °C
120)	-	DEW	PT - PRES	365.3927 bar	- TEMP	76.346 °C
121)	-	DEW	PT - PRES	367.8020 bar	- TEMP	74.816 °C
122)	-	DEW	PT - PRES	370.2271 bar	- TEMP	73.288 °C
123)	-	DEW	PT - PRES	372.6682 bar	- TEMP	71.735 °C
124)	-	DEW	PT - PRES	375.1255 bar	- TEMP	70.216 °C
125)	-	DEW	PT - PRES	379.9599 bar	- TEMP	67.227 °C
126)	-	DEW	PT - PRES	382.3694 bar	- TEMP	65.763 °C
127)	-	DEW	PT - PRES	384.8254 bar	- TEMP	64.283 °C
128)	-	DEW	PT - PRES	387.2350 bar	- TEMP	62.845 °C
129)	-	DEW	PT - PRES	389.6598 bar	- TEMP	61.420 °C
130)	-	DEW	PT - PRES	392.0997 bar	- TEMP	59.986 °C
131)	-	DEW	PT - PRES	394.5549 bar	- TEMP	58.591 °C
132)	-	DEW	PT - PRES	399.3893 bar	- TEMP	55.879 °C

4) *** BUBBLE POINT EQUILIBRIUM CURVE ***

Cricondentherm	Point - *** NOT DETERMINED ***	
Cricondenbar	Point - *** NOT DETERMINED ***	
Critical	Point - *** NOT DETERMINED ***	
Max Temp/Pres	Point - Temp -80.187 °C - Pres	55.3378 bar
1)	- BUBBLE PT - PRES	9.8067 bar - TEMP -137.830 °C
2)	- BUBBLE PT - PRES	10.7873 bar - TEMP -135.317 °C
3)	- BUBBLE PT - PRES	11.8660 bar - TEMP -132.733 °C
4)	- BUBBLE PT - PRES	13.0513 bar - TEMP -130.080 °C
5)	- BUBBLE PT - PRES	14.3580 bar - TEMP -127.346 °C
6)	- BUBBLE PT - PRES	15.7965 bar - TEMP -124.534 °C
7)	- BUBBLE PT - PRES	17.3803 bar - TEMP -121.640 °C
8)	- BUBBLE PT - PRES	19.1243 bar - TEMP -118.662 °C
9)	- BUBBLE PT - PRES	21.0453 bar - TEMP -115.596 °C
10)	- BUBBLE PT - PRES	23.1618 bar - TEMP -112.439 °C
11)	- BUBBLE PT - PRES	25.4943 bar - TEMP -109.189 °C
12)	- BUBBLE PT - PRES	28.0663 bar - TEMP -105.842 °C
13)	- BUBBLE PT - PRES	30.9004 bar - TEMP -102.402 °C
14)	- BUBBLE PT - PRES	34.0223 bar - TEMP -98.869 °C
15)	- BUBBLE PT - PRES	37.4601 bar - TEMP -95.251 °C
16)	- BUBBLE PT - PRES	41.2431 bar - TEMP -91.558 °C
17)	- BUBBLE PT - PRES	45.3960 bar - TEMP -87.818 °C
18)	- BUBBLE PT - PRES	49.9100 bar - TEMP -84.107 °C
19)	- BUBBLE PT - PRES	53.4347 bar - TEMP -81.478 °C
20)	- BUBBLE PT - PRES	55.3378 bar - TEMP -80.187 °C
21)	- BUBBLE PT - PRES	54.3884 bar - TEMP -80.736 °C

22)	-	BUBBLE	PT	-	PRES	52.5506	bar	-	TEMP	-81.745	°C
23)	-	BUBBLE	PT	-	PRES	48.4686	bar	-	TEMP	-83.843	°C
24)	-	BUBBLE	PT	-	PRES	45.4884	bar	-	TEMP	-85.293	°C
25)	-	BUBBLE	PT	-	PRES	44.6960	bar	-	TEMP	-85.673	°C
26)	-	BUBBLE	PT	-	PRES	44.1714	bar	-	TEMP	-85.916	°C
27)	-	BUBBLE	PT	-	PRES	43.9134	bar	-	TEMP	-86.037	°C
28)	-	BUBBLE	PT	-	PRES	43.7702	bar	-	TEMP	-86.104	°C
29)	-	BUBBLE	PT	-	PRES	43.6790	bar	-	TEMP	-86.145	°C



*** UNIT 2 - 'CYLIND ' - ' FLASH

--- Feed Streams ---

'WEST ' ' ***

- Product Streams -

'GASA0065' VAPOR to unit 4 - RECOMB

'LIQA0065' LIQUID to unit 3 - STD

1) * OPERATING CONDITIONS *

Outlet temperature	34.40 °C
Outlet pressure	37.6000 bar
Heat exchanged	236.630 MJ

2) * FEEDS *

Stream	WEST
Temperature, °C	40.55
Pressure, bar	172.800
Total rate, kmol/h	100.00
Vapor, kmol/h	94.411
Liquid, kmol/h	5.5893

3) * PRODUCTS *

Stream	GASA0065	LIQA0065
Temperature, °C	34.40	34.40
Pressure, bar	37.600	37.600
Total rate, kmol/h	95.430	4.5704
Vapor, kmol/h	95.430	0.0000
Liquid, kmol/h	0.0000	4.5704

*** UNIT 3 - 'STD' - 'FLASH' ***

--- Feed Streams ---

from unit 2 - CYLIND 'LIQA0065'

- Product Streams -

'SGS03' VAPOR to unit 4 - RECOMB

'STDLIQ' LIQUID

1) * OPERATING CONDITIONS *

Outlet temperature	15.56 °C
Outlet pressure	1.0130 bar
Heat exchanged	-5.834 MJ

2) * FEEDS *

Stream	LIQA0065
Temperature, °C	34.40
Pressure, bar	37.600
Total rate, kmol/h	4.5704
Vapor, kmol/h	0.0000
Liquid, kmol/h	4.5704

3) * PRODUCTS *

Stream	SGS03	STDLIQ
Temperature, °C	15.56	15.56
Pressure, bar	1.013	1.013
Total rate, kmol/h	1.2077	3.3627
Vapor, kmol/h	1.2077	0.0000
Liquid, kmol/h	0.0000	3.3627

*** UNIT 4 - 'RECOMB' - 'FLASH' ***

--- Feed Streams ---

from unit 2 - CYLIND 'GASA0065'

from unit 3 - STD 'SGS03'

- Product Streams -

'STDGAS'

1) * OPERATING CONDITIONS *

Outlet temperature	15.56 °C
Outlet pressure	1.0130 bar
Heat exchanged	5.716 MJ

2) * FEEDS *

Stream	GASA0065	SGS03
Temperature, °C	34.40	15.56
Pressure, bar	37.600	1.013
Total rate, kmol/h	95.430	1.2077
Vapor, kmol/h	95.430	1.2077
Liquid, kmol/h	0.0000	0.0000

3) * PRODUCTS *

Stream	STDGAS
Temperature, °C	15.56
Pressure, bar	1.013
Total rate, kmol/h	96.637
Vapor, kmol/h	96.637
Liquid, kmol/h	0.0000

*** UNIT 5 - 'DEW' - 'FLASH' ***

--- Feed Streams ---

'WEST'

- Product Streams -

'DEW'

1) * OPERATING CONDITIONS *

Outlet temperature	140.55 °C
Outlet pressure	251.6927 bar
Heat exchanged	649.707 MJ

2) * FEEDS *

Stream	WEST
Temperature, °C	40.55
Pressure, bar	172.8000
Total rate, kmol/h	100.00
Vapor, kmol/h	94.411
Liquid, kmol/h	5.5893

3) * PRODUCTS *

Stream	DEW
Temperature, °C	140.55
Pressure, bar	251.693
Total rate, kmol/h	100.00
Vapor, kmol/h	100.00
Liquid, kmol/h	0.0000

*** STREAM COMPONENT MOLAR FLOW RATES - kmol/h ***						
Stream Id	WEST	GASA0065	LIQA0065	SGS03	STDLIQ	STDGAS
Temperature, °C	40.550	34.400	34.400	15.560	15.560	15.560
Pressure, bar	172.8000	37.6000	37.6000	1.0130	1.0130	1.0130
Phase	MIXED	VAPOR	LIQUID	VAPOR	LIQUID	VAPOR
2 HYDROGEN SULPHIDE	0.370000	0.352498	0.017502	0.015327	0.002174	0.367826
3 NITROGEN	6.130000	6.117604	0.012396	0.012362	0.000034	6.129966
4 CARBON DIOXIDE	15.560000	15.296415	0.263585	0.253437	0.010148	15.549852
5 METHANE	65.850000	65.445719	0.404281	0.399944	0.004337	65.845663
6 ETHANE	4.400000	4.248849	0.151151	0.137657	0.013493	4.386507
7 PROPANE	2.180000	1.971272	0.208728	0.150209	0.058519	2.121481
8 ISOBUTANE	0.560000	0.457629	0.102371	0.050899	0.051472	0.508528
9 BUTANE	0.950000	0.723214	0.226786	0.089329	0.137457	0.812543
10 C5+	1.030000	0.526886	0.503114	0.069797	0.433317	0.596683
11 C6+	0.660000	0.178343	0.481657	0.019856	0.461800	0.198200
12 C7+	0.650000	0.078302	0.571698	0.006792	0.564906	0.085094
13 C8+	0.530000	0.024758	0.505242	0.001659	0.503583	0.026417
14 C9+	0.370000	0.006189	0.363811	0.000312	0.363498	0.006502
15 C10+	0.260000	0.001784	0.258216	0.000069	0.258147	0.001853
16 C11+	0.170000	0.000167	0.169833	0.000004	0.169829	0.000171
17 C12+	0.330000	0.000018	0.329982	0.000000	0.329982	0.000018
TOTALS, kmol/h	100.000000	95.429614	4.570386	1.207662	3.362724	96.637276
VAPOR fraction	0.944107	1.000000	0.000000	1.000000	0.000000	1.000000
LIQUID 1/H fraction	0.055893	0.000000	1.000000	0.000000	1.000000	0.000000
Enthalpy MJ/h	-423.509059	-42.558292	-144.321056	-0.732025	-149.422795	-37.574618

Stream Id	DEW
Temperature, °C	140.550
Pressure, bar	251.6927
Phase	VAPOR
2 HYDROGEN SULPHIDE	0.370000
3 NITROGEN	6.130000
4 CARBON DIOXIDE	15.560000
5 METHANE	65.850000
6 ETHANE	4.400000
7 PROPANE	2.180000
8 ISOBUTANE	0.560000
9 BUTANE	0.950000
10 C5+	1.030000
11 C6+	0.660000
12 C7+	0.650000
13 C8+	0.530000
14 C9+	0.370000
15 C10+	0.260000
16 C11+	0.170000
17 C12+	0.330000
TOTALS, kmol/h	100.000000
VAPOR fraction	1.000000
LIQUID 1/H fraction	0.000000
Enthalpy MJ/h	226.198279

*** STREAM MOLAR COMPOSITIONS ***						
Stream Id	WEST	GASA0065	LIQA0065	SGS03	STDLIQ	STDGAS
Temperature, °C	40.550	34.400	34.400	15.560	15.560	15.560
Pressure, bar	172.8000	37.6000	37.6000	1.0130	1.0130	1.0130
Phase	MIXED	VAPOR	LIQUID	VAPOR	LIQUID	VAPOR
2 HYDROGEN SULPHIDE	0.00370000	0.00369380	0.00382937	0.01269165	0.00064663	0.00380625
3 NITROGEN	0.06130000	0.06410593	0.00271221	0.01023598	0.00001018	0.06343273
4 CARBON DIOXIDE	0.15560000	0.16029002	0.05767242	0.20985761	0.00301782	0.16090946
5 METHANE	0.65850000	0.68580094	0.08845659	0.33117164	0.00128981	0.68136919
6 ETHANE	0.04400000	0.04452338	0.03307173	0.11398674	0.00401255	0.04539146
7 PROPANE	0.02180000	0.02065681	0.04566975	0.12437991	0.01740241	0.02195302
8 ISOBUTANE	0.00560000	0.00479547	0.02239866	0.04214635	0.01530663	0.00526223
9 BUTANE	0.00950000	0.00757851	0.04962077	0.07396841	0.04087675	0.00840817
10 C5+	0.01030000	0.00552120	0.11008126	0.05779494	0.12885895	0.00617446
11 C6+	0.00660000	0.00186885	0.10538643	0.01644203	0.13732922	0.00205097
12 C7+	0.00650000	0.00082053	0.12508735	0.00562410	0.16799045	0.00088056
13 C8+	0.00530000	0.00025943	0.11054697	0.00137395	0.14975451	0.00027336
14 C9+	0.00370000	0.00006486	0.07960177	0.00025873	0.10809640	0.00006728
15 C10+	0.00260000	0.00001870	0.05649758	0.00005688	0.07676724	0.00001917
16 C11+	0.00170000	0.00000175	0.03715940	0.00000322	0.05050338	0.00000177
17 C12+	0.00330000	0.00000019	0.07220005	0.00000017	0.09812935	0.00000019
TOTALS, /h	1.000000	1.000000	1.000000	1.000000	1.000000	1.000000
VAPOR fraction	0.944107	1.000000	0.000000	1.000000	0.000000	1.000000
LIQUID 1/H fraction	0.055893	0.000000	1.000000	0.000000	1.000000	0.000000
Spec Ent, kJ/kmol	-4235.859805	-446.046258	-31583.165628	-606.260734	-44443.116199	-388.891781

Stream Id	DEW
Temperature, °C	140.550
Pressure, bar	251.6927
Phase	VAPOR
2 HYDROGEN SULPHIDE	0.00370000
3 NITROGEN	0.06130000
4 CARBON DIOXIDE	0.15560000
5 METHANE	0.65850000
6 ETHANE	0.04400000
7 PROPANE	0.02180000
8 ISOBUTANE	0.00560000
9 BUTANE	0.00950000
10 C5+	0.01030000
11 C6+	0.00660000
12 C7+	0.00650000
13 C8+	0.00530000
14 C9+	0.00370000
15 C10+	0.00260000
16 C11+	0.00170000
17 C12+	0.00330000
TOTALS, /h	1.000000
VAPOR fraction	1.000000
LIQUID 1/H fraction	0.000000
Spec Ent, kJ/kmol	2262.393632

STREAM SUMMARY						
Stream Id		WEST	GASA0065	LIQA0065	SGS03	STDLIQ
Temperature, °C		40.550	34.400	34.400	15.560	15.560
Pressure, bar		172.8000	37.6000	37.6000	1.0130	1.0130
Phase		MIXED	VAPOR	LIQUID	VAPOR	LIQUID
Flow rate	kmol/h	100.000	95.430	4.570	1.208	3.363
	kg/h	2669.823	2254.203	415.620	44.960	370.660
Vapor fract.		0.944107	1.000000	0.000000	1.000000	0.000000
LIQUID 1/H FRACT		0.055893	0.000000	1.000000	0.000000	1.000000
LIQUID 2/W FRACT		0.000000	0.000000	0.000000	0.000000	0.000000
Mol. weight		26.6982	23.621625	90.937664	37.229354	110.226044
Act. Volume rate	m3/h	10.896660	59.332435	0.577453	28.350302	0.484853
Enthalpy	MJ/h	-423.509059	-42.558292	-144.321056	-0.732025	-149.422795
Spec. enthalpy	kJ/kmol	-4235.860	-446.046	-31583.166	-606.261	-44443.116
	kJ/kg	-158.657	-18.883	-347.306	-16.284	-403.200
*** VAPOUR ***						
Flow rate	kmol/h	94.411	95.430		1.208	
	kg/h	2302.316	2254.203		44.960	
Wt. fraction		0.862348	1.000000		1.000000	
Mol. weight		24.3862	23.6216		37.2294	
Std. Volume rate	N-m3/h	2236.589921	2260.727548		28.609520	
Act. Volume rate	m3/h	10.3438	59.3324		28.3503	
Enthalpy	MJ/h	-314.405646	-42.558292		-0.732025	
Spec. enthalpy	kJ/kmol	-3330.795	-446.046		-606.261	
	kJ/kg	-136.585	-18.883		-16.284	
Spec. Heat cap.	kJ/kmol °C	40.738091	39.142166		56.061979	
	kJ/kg °C	1.670540	1.657048		1.505854	
Spec. Entropy	kJ/kmol °C	148.0693	168.4872		224.0845	
	kJ/kg °C	6.0719	7.1328		6.0190	
Compressibility		0.725752	0.914078		0.990517	
Density	kg/m3	222.5802	37.9928		1.5859	
Isentr. Exponent		1.249446	1.300322		1.176492	
Viscosity	cP	0.2386E-01	0.1339E-01		0.9765E-02	
Thermal Conduct.	W/m°C	0.057321	0.032739		0.019299	
Reference Gas Status - Temperature	15.5		- Pressure 1			
*** LIQUID 1/H ***						
Flow rate	kmol/h	5.589		4.570		3.363
	kg/h	367.507		415.620		370.660
Wt. fraction		0.137652		1.000000		1.000000
Mol. weight		65.7520		90.9377		110.2260
Std. Volume rate	m3/h	0.541464		0.583287		0.502275
Act. Volume rate	m3/h	0.5529		0.5775		0.4849
Enthalpy	MJ/h	-109.103413		-144.321056		-149.422795
Spec. enthalpy	kJ/kmol	-19523.659		-31583.166		-44443.116
	kJ/kg	-296.929		-347.306		-403.200
Spec. Heat cap.	kJ/kmol °C	305.836280		268.390437		253.739510
	kJ/kg °C	4.651360		2.951367		2.301992
Spec. Entropy	kJ/kmol °C	95.1214		55.4700		26.0484
	kJ/kg °C	1.4467		0.6100		0.2363
Density	kg/m3	664.6911		719.7472		764.4794
Sp. Grav. 60/60		0.6787		0.7125		0.7379
Viscosity	cP	0.1033		0.2411		0.4553
Thermal Conduct.	W/m°C	0.121679		0.124252		0.128122
Surface Tension	dyn/cm	8.9238		15.4556		21.7673

STREAM SUMMARY

Stream Id		STDGAS	DEW
Temperature, °C		15.560	140.550
Pressure, bar		1.0130	251.6927
Phase		VAPOR	VAPOR
Flow rate	kmol/h	96.637	100.000
	kg/h	2299.163	2669.823
Vapor fract.		1.000000	1.000000
LIQUID 1/H FRACT		0.000000	0.000000
LIQUID 2/W FRACT		0.000000	0.000000
Mol. weight		23.7917	26.698228
Act. Volume rate	m3/h	2283.212862	12.683557
Enthalpy	MJ/h	-37.574618	226.198279
Spec. enthalpy	kJ/kmol	-388.892	2262.394
	kJ/kg	-16.346	84.739
*** VAPOUR ***			
Flow rate	kmol/h	96.637	100.000
	kg/h	2299.163	2669.823
Wt. fraction		1.000000	1.000000
Mol. weight		23.7917	26.6982
Std. Volume rate	N-m3/h	2289.337068	2369.000000
Act. Volume rate	m3/h	2283.2129	12.6836
Enthalpy	MJ/h	-37.574618	226.198279
Spec. enthalpy	kJ/kmol	-388.892	2262.394
	kJ/kg	-16.346	84.739
Spec. Heat cap.	kJ/kmol °C	38.451944	51.503550
	kJ/kg °C	1.616193	1.929100
Spec. Entropy	kJ/kmol °C	198.6212	160.4206
	kJ/kg °C	8.3483	6.0087
Compressibility		0.996900	0.927958
Density	kg/m3	1.0070	210.4948
Isentr. Exponent		1.276956	1.219736
Viscosity	cP	0.1176E-01	0.2540E-01
Thermal Conduct.	W/m°C	0.027266	0.064209
Reference Gas Status - Temperature	15.5		- Pressure 1

- Calculation time 0.41 sec

- Simulation no. 3 - Ended at 16.06.29 on MAR 04, 2009 -