

## CHAPTER V

### RESULTS AND DISCUSSION

#### 5.1 The Soave-Redlich-Kwong Equation of State

Generally, the SRK equation of state is good for estimating the K values of paraffin in the normal region. The normal region is defined as the area 1 in Figure 5.1. The SRK equation of state gives an average percentage deviation of 5.30 against the experimental data (7).

#### 5.2 The Heavy-Pseudocomponent

The critical behavior of carbon dioxide and paraffin system is shown in Figure 5.2. It shows that the curve of the critical loci of the binary carbon dioxide-paraffin of butane and heavier paraffins are similar. Therefore, any paraffin which is lighter than butane should not be included in the heavy-pseudocomponent.

#### 5.3 Validity of Using Pseudocomponents

The pseudocomponents model is applied in order to scale down the calculation task, both in memory spaces and CPU time of computers. The validity of using pseudocomponents in the model was tested by calculating K values by using the SRK equation of state. The K values were calculated for both cases : with individual components and with a pseudocomponent. Experimental data are from references (42). The results are presented in Table 5.1-5.18.

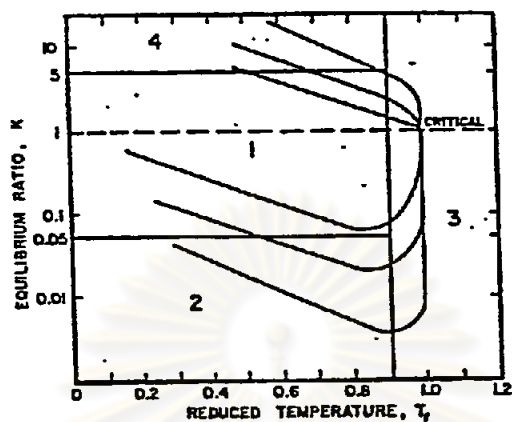


FIGURE 5.1 DEFINITION OF AREAS FOR ERROR ANALYSIS (7)

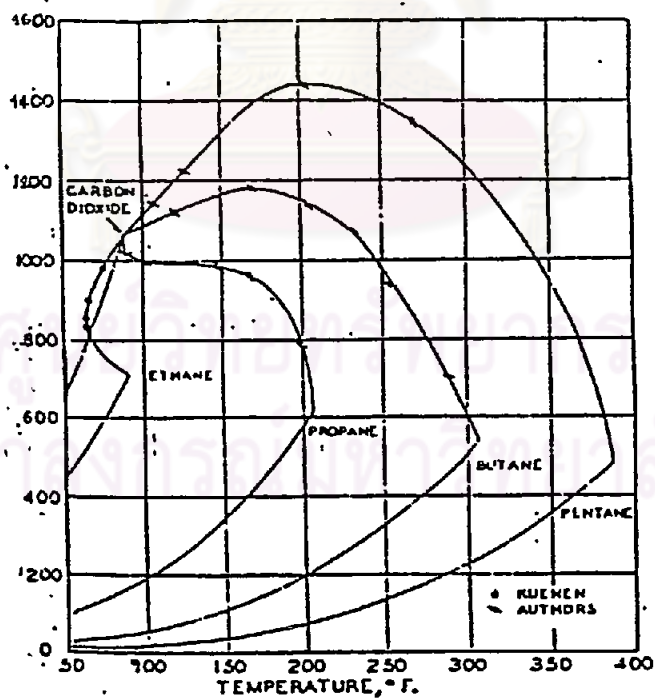


FIGURE 5.2 CRITICAL LOCI OF CARBON DIOXIDE-PARAFFIN SYSTEMS (27)

The average absolute deviation (26) is defined as

$$\%AAD = \frac{\sum_{i=1}^N |d_i|}{N} \times 100$$

where  $d_i$  are error

$N$  is the number of data points.

a) Using Heavy-Pseudocomponent ( $C_{M+}$ )

For  $M = 6$ , (pseudocomponent includes  $i-C_4$  and all components heavier than  $i-C_4$ )

the range of %AAD is 0.7 to 4.7

(see Table 5.1-5.6)

For  $M = 8$ , (pseudocomponent includes  $i-C_5$  and all components heavier than  $i-C_5$ )

the range of %AAD is 1.2 to 11.5

(see Table 5.7-5.12)

b) Using Light-Pseudocomponent ( $C_{M-}$ )

(pseudocomponent includes  $C_1$ ,  $C_2$  and  $C_3$ )

the range of %AAD is 0.5 to 8.0

It may be concluded that the heavy-pseudocomponent that includes  $i$ -butane and all paraffin heavier than  $i$ -butane is the best heavy-pseudocomponent.

#### 5.4 Study of Predicting the Compositions of Pseudocomponents

In the proposed general model, two pseudocomponents are provided, namely, the light pseudocomponent and the heavy pseudocomponent. The problem of flash calculation is that the liquid and vapor phase compositions are not known initially. Only the feed compositions is known. Consequently, the vapor and the liquid properties of the pseudocomponents cannot be evaluated.

However, to perform the flash calculation, the properties of pseudocomponents must be estimated.

A study was done to find a method for estimating the initial values of the vapor and liquid compositions of the components comprising the pseudocomponents. For natural gas under normal practical conditions,  $y_i$ 's are assumed to be slightly different from  $z_i$ 's, so that when normalized, one obtains initial approximations for the light pseudocomponent as follows

$$y_{i,ps} = z_i / \sum_{i=1}^{M-1} z_i \quad (5.1)$$

For the liquid phase composition, applying Equation (4.3), one has

$$x_{i,ps} = (z_i / K_i) / \sum_{i=1}^{M-1} (z_i / K_i) \quad (5.2)$$

when the ideal solution behavior is applied in order to estimate the K values, Equation (5.3) becomes

$$x_{i,ps} = (Pz_i / P_i) / \sum_{i=1}^{M-1} (Pz_i / P_i) \quad (5.3)$$

where  $P_i$  is the vapor pressure of component  $i$

Similarly, for the heavy pseudocomponent, one has

$$y_{i,ps} = z_i / \sum_{i=M}^N z_i \quad (5.4)$$

and

$$x_{i,ps} = (Pz_i / P_i) / \sum_{i=M}^N (Pz_i / P_i) \quad (5.5)$$

A study based on the experimental values in Reference (42) showed that Equation (5.1) to (5.2) gave rather good approximations, according to Table 5.19-5.24, and Figure 5.3-5.6. For the ideal case, it did not good prediction of the initial liquid phase composition. It is thus proposed that the initial estimation of the liquid phase composition be set by

$$x_i = 1.0/N$$

where  $N$  is the total number of components.

### 5.5 Comparison between the Ordinary Vapor-Liquid Equilibrium Calculation Model and the Proposed General Model with Pseudocomponents

In the ordinary model (Figure 4.3), the number of equations to be solved simultaneously equal to the number of components in natural gas. The more the components are in natural gas, the more the equations must be solved. For the proposed model (Figure 4.4), the components which are similar in physical properties are grouped into one pseudocomponent. The critical properties of the pseudocomponent are calculated by Kay's rule. Thus the number of equations to be solved in this model are less than that of the ordinary model. The calculation of the liquid phase composition ( $x_1$ ) and the vapor phase composition ( $y_1$ ) by flash calculations at low and moderate pressure systems were done using both models. The results are presented in Table 5.28-5.55.

Comparison of the results produced by the ordinary model and the proposed model are summarized as follows :

	The Ordinary Model		The Proposed Model	
	L-phase	V-phase	L-phase	V-phase
CPU time(SU)	1.25		0.85	
Memory Spaces(K)	32		36	
%AAD(Comparison with the experimental data)	1.1-3.1	0.0-0.2	0.6-3.1	0.0-0.1
%AAD(Comparison with the ordinary model)	-	-	1.4-2.5	0.0-0.04

It can be concluded that the proposed yields results faster than the ordinary model. The computing time saving is on the average about 32%. But the proposed model requires more spaces to store the initial property values. Compared with the experimental data, the proposed model gives better results than those of the ordinary model. This may be because of propable damping of the fluctuation in property values of the individual components forming the pseudocomponen. Because of good estimation for predicting properties in vapor phase of the SRK equation of state, the percentage average deviation of the results in vapor phase is less than in liquid phase.

It may be concluded that the proposed general model with pseudocomponents is successful in saving computing time by as much as 32%.



คุรุณย์วิทยทรรพยากร  
จุฬาลงกรณ์มหาวิทยาลัย

THE AMOUNT OF COMPONENT, N = 16

T = 41,000 °F      P = 300.00 PSIA      R = 10.73C      PSIA-CUBIC FT/LB-MOLE<sup>OR</sup>

FEED RATE = 1,117,670 MOLE/HR      V/F = 0.500000

TABLE 5.1 COMPARISON OF EQUILIBRIUM K VALUES FROM THE ORDINARY METHOD AND THE PROPOSED GENERAL MODEL FOR HEAVY-PSEUDOCOMPONENT (M=6) AT T = 41.0 °F    P = 300.0 PSIA

ID	COMPONENT	X (%)	Y (%)	K	TC (K)	PC (PSIA)	W
1	C02	0.088	0.395300	4.4920450	548.00	1073.00	0.22500
2	N2	0.015	0.828000	55.1999900	227.20	472.00	0.04000
3	C1	3.564	05.016220	9.9165650	343.30	673.10	0.01400
4	C2	5.574	7.720530	1.3835060	549.77	703.30	0.09900
5	C3	3.238	3.367210	0.4035820	605.95	617.40	0.15200
6	IC4	7.904	1.225400	0.1541760	734.65	529.10	0.13500
7	NC4	9.630	0.919200	0.0941170	703.31	550.70	0.20100
8	IC5	9.678	0.279400	0.0293820	829.80	483.00	0.22200
9	NC5	6.855	0.139200	0.0202860	845.60	499.50	0.25400
10	C6	15.402	0.077600	0.0063310	514.20	439.70	0.30100
11	C7	11.893	0.012490	0.0010910	572.31	396.90	0.35000
12	C8	10.774	0.004340	0.0004020	1024.31	362.10	0.40200
13	C9	4.454	0.001120	0.0002510	1073.00	345.00	0.44600
14	C10	0.869	0.000080	0.0000920	1114.70	306.00	0.48900
15	C11	0.062	0.000000	0.0000000	1153.70	292.00	0.50100
16	C12	0.000	0.000000	0.0000000	1137.70	263.00	0.53900

IPH = 1  
 A = 0.09317      B1 = 0.02932  
 J = 0.05745      RR = 0.00293  
 Z = 0.93092

APH = 0  
 A = 1.06592      B1 = 0.09002  
 J = 0.96779      RR = 0.09596  
 Z = 0.11023

COMPONENT	X	Y	K(EXPI)	K(CAL)	K(MODEL)	TOLERANCE
C02	0.088000	0.395300	4.4920450	3.4344990	3.4696860	0.0351861
N2	0.015000	0.828000	55.1999700	23.3267900	29.0891700	0.7621765
C1	3.564000	5.016220	9.9165650	7.9602000	8.1041060	0.1390367
C2	5.574000	7.720530	1.3835060	1.2731400	1.2870370	0.0108976
C3	3.238000	3.367210	0.4035820	0.3335300	0.3401210	0.0062580
C4+	7.904000	1.225400	0.3066278	0.1316900	0.0071788	-0.1245112
AAO =						4.6362230

THE AMOUNT OF COMPONENT, N = 16

T = 30.000 ° F      P = 300.00 PSIA      R = 10.730 PSIA-CUBIC FT/LB-MOLE ° R

FEED RATE = 0.503350 MOLE/HR      V/F = 0.500000

TABLE 5.2 COMPARISON OF EQUILIBRIUM K VALUES FROM THE ORDINARY METHOD

AND THE PROPOSED GENERAL MODEL FOR HEAVY-PSEUDOCOMPONENT (M=6)

AT T = 30.0 ° F      P = 300.0 PSIA

ID	COMPONENT	X	Y	K	TC	PC	W
		(%)	(%)		(R)	(PSIA)	
1	CO2	0.065	0.250400	3.8523080	543.00	1073.00	0.22500
2	N2	0.019	0.849000	44.6315700	227.20	492.00	0.04000
3	C1	0.522	84.699400	8.8876580	343.30	673.10	0.01400
4	C2	1.764	7.922510	2.1020110	549.77	708.30	0.05900
5	C3	0.770	3.713010	0.3902220	665.95	617.40	0.15200
6	C4	10.308	1.187900	0.1151390	734.65	529.10	0.18500
7	NC4	11.833	0.814000	0.0631030	765.21	550.70	0.20100
8	IC5	12.244	0.310400	0.0253250	829.80	483.00	0.22200
9	NC5	7.713	0.133900	0.0172940	845.60	489.50	0.25400
10	C6	14.212	0.051000	0.0059880	914.20	439.70	0.30100
11	C7	7.321	0.023650	0.0022140	972.21	375.90	0.35000
12	C8	7.332	0.004650	0.0000340	1024.31	362.10	0.40200
13	C9	2.898	0.000960	0.0000000	1073.00	345.00	0.44600
14	C10	0.848	0.000100	0.0000000	1114.70	306.00	0.48900
15	C11	0.149	0.000010	0.0000000	1153.70	232.00	0.56100
16	C12	0.000	0.000000	0.0000000	1187.70	263.00	0.53900

IPH = 1  
 A = 0.10483      B1 = 0.03058  
 J = 0.07332      RR = 0.00321  
 Z = 0.92444

IPH = 0  
 A = 1.01681      B1 = 0.08720  
 Q = 0.92201      RR = 0.08866  
 Z = 0.10720

COMPONENT	X	Y	K (EXPT)	K (CAL)	K (MODEL)	TOLERANCE
CO2	0.065000	0.250400	3.8523080	3.130300	3.1702670	0.0322571
N2	0.019000	0.849000	44.6315700	27.6572500	20.4361700	0.7785154
C1	0.522000	84.699400	8.8876580	7.5251300	7.6611310	0.1360016
C2	1.764000	7.922510	2.1020110	1.1419690	1.1516180	0.0096493
C3	0.770000	3.713010	0.3902220	0.2502300	0.2915384	0.0012584
C4+	10.307990	1.187900	0.2357000	0.1096800	0.0087765	-0.1009035
MAAD =						4.5333860



THE AMOUNT OF COMPONENT, N = 16

T = 30.000 °F      P = 300.00 PSIA      R = 10.730 PSIA-CUBIC FT/LB-MOLE °R

FEED RATE = 0.790030 MOLE/HR      V/F = 0.500000

TABLE 5.3 COMPARISON OF EQUILIBRIUM K VALUES FROM THE ORDINARY METHOD AND THE PROPOSED GENERAL MODEL FOR HEAVY-PSEUDOCOMPONENT (M=6) AT T = 30.0°F P = 300.0 PSIA

ID	COMPONENT	X (%)	Y (%)	K	TC (R)	PC (PSIA)	W
1	CO2	0.056	0.177200	3.1642850	548.00	1073.00	0.22500
2	N2	0.029	0.937020	32.3110300	227.20	492.00	0.04000
3	C1	10.818	04.920380	7.8432230	343.30	673.10	0.01400
4	C2	7.039	7.785640	1.1052010	549.77	703.30	0.09900
5	C3	10.875	3.664360	0.3367250	665.55	617.40	0.15200
6	IC4	9.692	1.150220	0.1185790	734.65	529.10	0.18500
7	NC4	11.420	0.800010	0.0076950	765.31	550.70	0.20100
8	IC5	10.886	0.303310	0.0278670	829.80	493.00	0.22200
9	NC5	7.205	0.127000	0.0176120	845.60	489.50	0.25400
10	C6	13.468	0.091790	0.0068100	914.20	439.70	0.30100
11	C7	8.920	0.027400	0.0030690	972.31	396.90	0.35000
12	C8	6.320	0.003610	0.0008850	1024.31	362.10	0.40200
13	C9	2.411	0.000640	0.0002640	1073.00	345.00	0.44600
14	C10	0.727	0.000110	0.0000000	1114.70	306.00	0.48900
15	C11	0.127	0.000010	0.0000000	1153.70	232.00	0.50100
16	C12	0.000	0.000000	0.0000000	1187.70	253.00	0.53900

IPH = 1  
 A = 0.10425      B1 = 0.03052  
 Q = 0.07280      RR = 0.00318  
 Z = 0.92502

IPH = 0  
 A = 0.93221      B1 = 0.08344  
 Q = 0.84181      RR = 0.07778  
 Z = 0.19375

COMPONENT	X	Y	K (EXP)	K (ICAL)	K (MODEL)	TOLERANCE
CO2	0.056000	0.177200	3.1642850	3.1124650	3.1446050	0.0321360
N2	0.029000	0.937020	32.3110300	26.9741900	27.7234800	0.7452820
C1	10.817990	04.920380	7.8432230	7.4175790	7.5505770	0.1329975
C2	7.039000	7.785640	1.1052010	1.1247590	1.1441510	0.0093913
C3	10.875000	3.664360	0.3367250	0.2893200	0.2904493	0.0011253
C4+	9.692000	1.150220	0.1182707	0.1132700	0.0091656	0.1001044
MAAD =						4.4322500

THE AMOUNT OF EQUIPMENT = 16

T = -8.000 ° F      P = 1115.00 PSIA      R = 10.730 PSIA-CUBIC FT/LB-MOLE<sup>OR</sup>

FEED RATE = 2.630090 MOLE/HR      V/F = 0.500000

TABLE 5.4 COMPARISON OF EQUILIBRIUM K VALUES FROM THE ORDINARY METHOD AND THE PROPOSED GENERAL MODEL FOR HEAVY-PSEUDOCOMPONENT (M=6) AT T = -8.0 ° F      P = 1115.0 PSIA

ID	COMPONENT	X (%)	Y (%)	K	TC (R)	PC (PSIA)	W
1	CJ2	0.589	7.097400	1.2049900	548.00	1073.00	0.22500
2	Y2	0.472	4.210720	0.9210190	227.20	492.00	0.04000
3	C1	37.190	91.002390	2.4210660	343.30	673.10	0.01400
4	C2	4.604	3.046180	0.6548110	349.77	793.30	0.09900
5	C3	3.491	0.582920	0.1652300	665.95	617.40	0.15200
6	IC4	2.137	0.135280	0.0620500	734.65	529.10	0.18500
7	NC4	2.995	0.138320	0.0423870	705.31	350.70	0.20100
8	IC5	3.240	0.059490	0.0181670	829.30	433.00	0.22200
9	NC5	2.597	0.036710	0.0133860	945.60	489.50	0.23400
10	C6	7.432	0.048440	0.0064500	914.20	439.70	0.30100
11	C7	9.005	0.025980	0.0026220	972.31	376.70	0.35000
12	C8	10.591	0.008570	0.0003010	1024.31	362.10	0.40200
13	C9	7.952	0.003420	0.0004300	1073.00	345.00	0.44600
14	C10	4.317	0.000910	0.0002100	1114.70	326.00	0.48300
15	C11	1.696	0.000180	0.0001100	1151.70	282.00	0.50100
16	C12	0.894	0.000040	0.0000400	1187.70	263.00	0.53900

IPH = 1  
 A = 0.43511      BI = 0.11903  
 Q = 0.33191      RR = 0.05179  
 Z = 0.56228

IPH = 0  
 A = 4.12757      BI = 0.33923  
 Q = 3.67326      RR = 1.40019  
 Z = 0.40801

COMPONENT	X	Y	K (EXP)	K (CAL)	K (MODEL)	TOLERANCE
CJ2	0.537000	7.077400	1.2049900	1.0457000	1.0500500	0.0943411
Y2	0.472000	4.210720	0.9210190	0.7024490	0.8722060	0.1697569
C1	37.190000	91.002390	2.4210660	2.1752090	2.2054050	0.0311756
C2	4.604000	3.046180	0.6548110	0.4810800	0.4827543	0.0016745
C3	3.491000	0.582920	0.1652300	0.1710500	0.170968	-0.0001522
C4+	2.137000	0.135280	0.1478521	0.8684000	0.0006044	-0.8657950
%AAD =						1.0367540

TABLE 5.5 COMPARISON OF EQUILIBRIUM K VALUES FROM THE ORDINARY METHOD AND THE PROPOSED GENERAL MODEL FOR HEAVY-PSEUDOCOMPONENT (M=6)

AT T = -10.0°F P = 800.0 PSIA

THE AMOUNT OF COMPONENT, N = 16

T = -10.000 °F P = 800.00 PSIA R = 10.730 PSIA-CUBIC FT/LB-MOLE<sup>OR</sup>

FEED RATE = 2.543590 MOLE/HR V/F = 0.500000

ID	COMPONENT	X (%)	Y (%)	K	TC (R)	PC (PSIA)	W
1	C02	0.143	0.373660	0.2613000	543.00	1073.00	0.22500
2	N2	0.006	0.205190	34.1983300	227.20	492.00	0.04000
3	C1	31.340	97.267850	3.0990310	343.30	673.10	0.01400
4	C2	2.363	2.092600	0.8841960	549.77	703.30	0.09700
5	C3	0.165	0.029010	0.1655930	605.95	517.40	0.15200
6	IC4	0.056	0.005200	0.0923040	734.65	529.10	0.18500
7	NC4	0.177	0.005300	0.0384300	765.31	550.70	0.20100
8	IC5	0.605	0.006700	0.0110600	829.80	493.00	0.22200
9	NC5	0.498	0.003900	0.0071100	845.60	489.50	0.25400
10	C6	1.702	0.003710	0.0022930	914.20	439.70	0.30100
11	C7	3.929	0.003010	0.0007040	972.31	396.90	0.35700
12	C8	5.857	0.001500	0.0002700	1024.21	362.10	0.40200
13	C9	14.467	0.000850	0.0000600	1073.00	345.00	0.44600
14	C10	25.830	0.000450	0.0000230	114.70	306.00	0.48900
15	C11	10.695	0.000110	0.0000100	1133.70	292.00	0.50100
16	C12	1.902	0.000010	0.0000050	1197.70	263.00	0.53900

IPH = 1.1  
 A = 3.26355 BI = 0.07949  
 J = 0.17775 RR = 0.02095  
 Z = 0.81319

IPH = 0  
 A = 6.34436 BI = 0.36397  
 J = 5.34791 RR = 2.30917  
 Z = 3.41194

COMPONENT	X	Y	K (EXP)	K (CAL)	K (MODEL)	TOLERANCE
C02	0.143000	0.373660	0.2613000	1.1140600	1.1180860	0.0048265
C2	0.006000	0.205190	34.1983300	10.7217500	10.9390700	0.2179203
C1	31.337993	97.267350	3.0990310	2.8262500	2.3593040	0.0330544
C2	2.363000	2.092600	0.8841767	0.4337200	0.4355869	0.0016669
C3	0.163000	0.029010	0.1655930	0.1156600	0.1137812	0.0001212
C++	0.055000	0.005200	0.1535367	0.0460500	0.0000077	-0.0468423
%AAD =						1.1019110

THE AMOUNT OF COMPONENT, N = 16

T = 15.000 °F      P = 850.00 PSIA      R = 10.730 PSIA-CUBIC FT/LB-MOLE °R

FEED RATE = 1.357700 MOLE/HR      V/F = 0.500000

TABLE 5.6 COMPARISON OF EQUILIBRIUM K VALUES FROM THE ORDINARY METHOD

AND THE PROPOSED GENERAL MODEL FOR HEAVY-PSEUDOCOMPONENT (M=6)

AT T = 15.0 °F      P = 850.0 PSIA

ID	COMPONENT	X (%)	Y (%)	K	TC (R)	PC (PSIA)	M
1	CO2	0.000	0.000000	0.000000	548.00	1073.00	0.22500
2	N2	0.103	0.944000	9.1650470	227.20	492.00	0.04000
3	C1	27.411	86.800350	3.1633360	343.30	673.10	0.01400
4	C2	15.266	0.197730	0.5364300	549.77	708.30	0.09900
5	C3	13.796	2.557410	0.1851800	665.95	617.40	0.15200
6	IC4	8.911	0.657100	0.0736640	734.65	529.10	0.18500
7	NC4	7.995	0.505800	0.0503530	765.31	550.70	0.20100
8	IC5	7.738	0.186900	0.0241300	829.80	483.00	0.22200
9	NC5	4.285	0.078700	0.0183480	845.60	489.50	0.25400
10	C6	6.677	0.056720	0.0084850	914.20	439.70	0.30100
11	C7	3.533	0.012330	0.0034500	972.31	396.90	0.35000
12	C8	1.737	0.002620	0.0015080	1024.31	362.10	0.40200
13	C9	0.441	0.000280	0.0006350	1073.00	345.00	0.44600
14	C10	0.094	0.000030	0.0003190	1114.70	306.00	0.48900
15	C11	0.012	0.000000	0.0000000	1153.70	282.00	0.50100
16	C12	0.001	0.000000	0.0000000	1187.70	263.00	0.52900

IPH = 1  
 A = 0.30079      3I = 0.08688  
 Q = 0.20635      RR = 0.02613  
 Z = 0.77792

IPH = 0  
 A = 1.53505      0I = 0.17928  
 Q = 1.32362      RR = 0.27521  
 Z = 0.24129

COMPONENT	X	Y	K (EXP)	K (CAL)	K (MODEL)	TOLERANCE
CO2	0.000000	0.000000	0.000000	1.2423250	1.2463130	0.0039835
N2	0.103000	0.944000	9.1650470	7.8123300	7.9326440	0.1203146
C1	27.410790	86.800350	3.1633360	2.6232790	2.6460330	0.0227594
C2	15.265390	0.197730	0.5364300	0.5430300	0.5490063	0.0005763
C3	13.795390	2.557410	0.1851800	0.1781400	0.1773841	-0.0003559
C4+	8.911000	0.657100	0.0736640	0.0711700	0.07152430	-0.0059212
AAD =						0.7516475

THE AMOUNT OF COMPONENT, N = 16

T = 41.000 °F      P = 300.00 PSIA      R = 10.730 PSIA-CUBIC FT/LB-MOLE °R

FEED RATE = 1.319570 MOLE/HR      V/F = 0.500000

ID	COMPONENT	X (%)	Y (%)	K (EXPI)	TC (R)	PC (PSIA)	W
1	CO2	0.088	0.395300	4.4920450	548.00	1073.00	0.22500
2	N2	0.015	0.828000	55.1999900	227.20	492.00	0.04000
3	C1	3.564	85.016220	9.9165650	243.30	673.10	0.01400
4	C2	5.574	7.720530	1.3835060	549.77	708.30	0.09900
5	C3	8.238	3.369210	0.4085820	565.95	617.40	0.15200
6	IC4	7.904	1.225400	0.1548760	734.65	529.10	0.18500
7	NC4	9.630	0.910200	0.0944170	755.31	550.70	0.20100
8	IC5	9.678	0.279800	0.0288820	829.80	483.00	0.22200
9	NC5	6.855	0.139200	0.0202850	845.60	489.50	0.25400
10	C6	15.402	0.097600	0.0063310	914.20	439.70	0.30100
11	C7	11.893	0.012990	0.0010910	972.31	396.90	0.35000
12	C8	10.774	0.004340	0.0004020	1024.31	362.10	0.40200
13	C9	4.454	0.001120	0.0002510	1073.00	345.00	0.44600
14	C10	0.869	0.000080	0.0000920	1114.70	306.00	0.48900
15	C11	0.062	0.000000	0.0000000	1153.70	282.00	0.50100
16	C12	0.000	0.000000	0.0000000	1187.70	263.00	0.53900

IPH = 1.

A = 0.09823      B1 = 0.02983  
 Q = 0.06751      RR = 0.00293  
 Z = 0.93086

IPH = 0

A = 1.09831      B1 = 0.09178  
 J = 0.39810      RR = 0.10080  
 Z = 0.11209

COMPONENT	-X-	-Y-	K (EXPI)	K (SRK)	K (MODEL)	-TOLERANCE
CO2	0.088000	0.395300	4.4920450	3.4344990	3.4618270	0.0273275
N2	0.015000	0.828000	55.1999900	28.3263900	-28.8787900	0.5518036
C1	3.564000	85.016220	9.9165650	7.9650200	9.0678470	0.1028271
C2	5.574000	7.720530	-1.3835060	-1.2781400	1.2865630	0.0084238
C3	8.238000	3.369210	0.4085820	0.3333530	0.3398067	0.0059537
IC4	7.904000	1.225400	0.1548760	0.1316900	0.1321259	0.0004358
NC4	9.630000	0.910200	0.0944170	0.0890900	0.0891824	0.0000924
C5+	9.678000	0.279800	0.0573350	-0.0672998	0.0031035	-0.0641963
ZAAD =						1.6819976

TABLE 5.7      COMPARISON OF EQUILIBRIUM K VALUES FROM THE ORDINARY METHOD  
 AND THE PROPOSED GENERAL MODEL FOR HEAVY-PSEUDOCOMPONENT (M=8)  
 AT T = 41.0 °F      P = 300.0 PSIA

THE AMOUNT OF COMPUTATION = 16

T = 30.000 °F      P = 300.00 PSIA      R = 10.733 PSIA-CUBIC FT/LB-MOLE°R

FEED RATE = 0.730030 MOLE/HR      V/F = 0.500000

TABLE 5.8 COMPARISON OF EQUILIBRIUM K VALUES FROM THE ORDINARY METHOD AND THE PROPOSED GENERAL MODEL FOR HEAVY-PSEUDOCOMPONENT (M=8)

AT T = 30.0°F      P = 300.0 PSIA

ID	COMPONENT	X (%)	Y (%)	K (EXP)	TC (R)	PC (PSIA)	W
1	CO2	0.056	0.177200	3.1642850	548.00	1073.00	0.22500
2	N2	0.029	0.937020	32.3110300	227.20	472.00	0.04000
3	C1	10.818	84.920830	7.6432230	343.30	673.10	0.01400
4	C2	7.039	7.785640	1.1052010	543.77	708.30	0.05900
5	C3	10.875	3.664960	0.3367250	665.95	617.40	0.15200
6	IC4	9.692	1.150220	0.1185790	734.65	529.10	0.18500
7	HC4	11.420	0.304010	0.0076950	765.31	550.70	0.20100
8	IC5	10.886	0.303110	0.0273670	823.80	403.00	0.22200
9	HC5	7.205	0.127000	0.0170120	145.60	489.50	0.25400
10	C6	13.468	0.071790	0.0068100	514.20	439.70	0.30100
11	C7	3.520	0.027400	0.0033050	972.31	335.90	0.35000
12	C8	6.328	0.005610	0.0008850	1024.31	362.10	0.40200
13	C9	2.411	0.000640	0.0002640	1073.00	345.00	0.44600
14	C10	0.727	0.000110	0.0000000	1114.70	305.00	0.48700
15	C11	0.127	0.000010	0.0000000	1153.70	292.00	0.50100
16	C12	0.000	0.000000	0.0000000	1187.70	263.00	0.53900

IPH = 1  
 A = 0.10432      BI = 0.03053  
 Y = 0.07286      RR = 0.00319  
 Z = 0.92495

IPH = 0  
 A = 0.95846      BI = 0.08497  
 Y = 0.96627      RR = 0.08144  
 Z = 0.10537

COMPONENT	X	Y	K (EXP)	K (SRK)	K (MODEL)	TOLERANCE
CO2	0.056000	0.177200	3.1642850	3.1124650	3.1332990	0.0258303
N2	0.029000	0.937020	32.3110300	26.9741900	27.5320100	0.5578156
C1	10.817390	84.920830	7.6432230	7.4175790	7.5137150	0.1011362
C2	7.039000	7.785640	1.1052010	1.1347550	1.1422580	0.0074987
C3	10.875000	3.664960	0.3367250	0.2932000	0.2902702	0.0009502
IC4	9.692000	1.150220	0.1185790	0.1092700	0.1095417	0.0002717
HC4	11.420000	0.304010	0.0076950	0.0732500	0.0732526	0.0000026
C5+	10.885990	0.303110	0.0565070	0.0310400	0.03033489	-0.0455911
MAAD =						1.5654200

THE AMOUNT OF COMPONENT, N = 16

T = 30.000 ° F      P = 300.00 PSIA      R = 10.730 PSIA-CUBIC FT/L3-MOLE°R

FEED RATE = 3.53330 MOLE/HR      V/F = 0.500000

ID	COMPONENT	X	Y	K	TC	PC	W
		(#)	(%)	(EXP)	(R)	(PSIA)	
1	C02	0.065	0.253430	3.8523080	548.00	1073.00	0.22500
2	N2	0.019	0.048330	44.6315700	227.20	492.00	0.04000
3	C1	9.522	84.699430	9.8876580	243.30	673.10	0.01400
4	C2	3.764	7.922510	2.1023110	149.77	708.30	0.09900
5	C3	9.770	3.713010	0.3802220	165.55	617.40	0.15200
6	IC4	10.308	1.187700	0.1151390	734.65	529.10	0.18500
7	NC4	11.833	0.818400	0.0691030	705.31	550.70	0.20100
8	IC3	12.244	0.310400	0.0251290	129.80	483.00	0.22200
9	NC5	7.713	0.133500	0.0173940	145.60	439.50	0.25400
10	C6	14.212	0.035130	0.005380	514.20	439.70	0.30100
11	C7	9.321	0.023550	0.0022140	972.31	376.90	0.35000
12	C8	7.332	0.004650	0.0000340	1024.31	362.10	0.40200
13	C9	2.898	0.000860	0.0000000	1073.00	345.00	0.44600
14	C10	0.848	0.000130	0.0000000	1114.70	306.00	0.48900
15	C11	0.149	0.000010	0.0000000	1153.70	282.00	0.50100
16	C12	0.000	0.000000	0.0000000	1187.70	263.00	0.53900

IPH = 1  
 A = 0.10490      BI = 0.03059  
 Q = 0.07339      BR = 0.00321  
 Z = 0.92437

IPH = 0  
 A = 1.04589      BI = 0.08882  
 J = 0.94919      BR = 0.09290  
 Z = 0.10892

COMPONENT	X	Y	K (EXP)	K (SRK)	K (MODEL)	TOLERANCE
C02	0.065000	0.253430	3.8523080	3.1280000	3.1634650	0.0254354
N2	0.019000	0.048330	44.6315700	27.6572500	23.2234500	0.5711975
C1	9.522000	84.699430	9.8876580	7.5251300	7.6269440	0.1018143
C2	3.764000	7.922510	2.1023110	1.1413450	1.1435340	0.0075645
C3	9.770000	3.713010	0.3802220	0.2902800	0.2913132	0.0010332
IC4	10.307390	1.187700	0.1151390	0.1096000	0.1100050	0.0003250
NC4	11.833000	0.818400	0.0691030	0.0732800	0.0733259	0.0000458
C5+	12.244000	0.310400	0.0251290	0.0130300	0.0032767	0.0497532
TAAD =						1.5320120

TABLE 5.9 COMPARISON OF EQUILIBRIUM K VALUES FROM THE ORDINARY METHOD AND THE PROPOSED GENERAL MODEL FOR HEAVY-PSEUDOCOMPONENT (M=8) AT T = 30.0 ° F      P = 300.0 PSIA

THE AMOUNT OF COMPONENT, N = 16

T = -10.000 °F      P = 800.00 PSIA      R = 10.730 PSIA-CUBIC FT/LB-MOLE<sup>0</sup>R

FEED RATE = 2.543590 MOLE/HR      V/F = 0.500000

ID	COMPONENT	X (%)	Y (%)	K (EXP)	TC (R)	PC (PSIA)	W
1	CO2	0.143	0.373660	0.2613000	548.60	1073.00	0.22500
2	N2	0.006	0.205190	34.1983300	227.20	492.00	0.04000
3	C1	31.340	97.267850	3.0990310	143.30	673.10	0.01400
4	C2	2.363	2.092600	0.8841960	549.77	708.30	0.09900
5	C3	0.169	0.028010	0.1655930	665.95	617.40	0.15200
6	IC4	0.056	0.005200	0.0928040	734.65	529.10	0.18500
7	NC4	0.177	0.006800	0.0384300	765.31	550.70	0.20100
8	IC5	0.605	0.006700	0.0110600	829.80	483.00	0.22200
9	NC5	0.498	0.003900	0.0078180	845.60	489.50	0.25400
10	C6	1.702	0.003910	0.0022930	914.20	439.70	0.30100
11	C7	3.929	0.003010	0.0007640	972.31	396.90	0.35000
12	C8	5.857	0.001580	0.0002700	1024.31	362.10	0.40200
13	C9	14.467	0.000850	0.0000600	1073.60	345.00	0.44600
14	C10	25.830	0.000650	0.0000230	1114.70	306.00	0.48900
15	C11	10.695	0.000110	0.0000100	1153.70	282.00	0.50100
16	C12	1.902	0.000010	0.0000050	1187.70	263.00	0.53900

IPH = L  
 A = 0.25356      B1 = 0.07949  
 J = 0.17775      RR = 0.02095  
 Z = 0.81318

IPH = O  
 A = 6.37330      B1 = 0.36497  
 J = 5.87512      RR = 2.32606  
 Z = 0.41296

COMPONENT	X	Y	K (EXP)	K (SRK)	K (MODEL)	TOLERANCE
CO2	0.143000	0.373660	0.2613000	1.1140600	1.1187120	0.0046520
N2	0.006000	0.205190	34.1983300	10.7217500	10.9306900	0.2089434
C1	31.333990	97.267150	3.0990310	2.8262500	2.8579260	0.0316763
C2	2.363000	2.092600	0.8841960	0.4339200	0.4355212	0.0016012
C3	0.169000	0.023010	0.1655930	0.1156600	0.1157777	0.0001177
IC4	0.056000	0.005200	0.0928040	0.0468500	0.0469138	0.0000638
NC4	0.177000	0.006900	0.0384300	0.0306800	0.0306328	-0.0007472
IC5	0.605000	0.006700	0.0223030	0.0240500	0.0000095	-0.0240405
%AAD =						0.4786583

TABLE 5.10 COMPARISON OF EQUILIBRIUM K VA USES FROM THE ORDINARY METHOD AND THE PROPOSED GENERAL MODEL FOR HEAVY-PSEUDOCOMPONENT (M=8) AT T = -10.0 °F P = 800.0 PSIA



THE AMOUNT OF COMPONENT, N = 16

T = 15,000 °F      P = 850.00 PSIA      R = 10,730 PSIA-CUBIC FT/LB-MOLE °R

FEED RATE = 1.357700 MOLE/HR      V/F = 0.500000

ID	COMPONENT	-X (%)	Y (%)	-K (EXP)	TC (R)	-PC (PSIA)	* K
1	CO2	0.000	0.000000	0.000000	548.00	1073.00	0.22500
2	N2	0.103	0.944000	9.1650470	227.20	492.00	0.04000
3	CO	27.411	86.800350	3.1633360	343.30	673.10	0.01400
4	C2	15.266	8.197730	0.5364300	549.77	708.30	0.09900
5	C3	13.796	2.557410	0.1851800	605.95	617.40	0.15200
6	IC4	9.911	0.657100	0.0736640	714.65	529.10	0.18500
7	NC4	9.995	0.505800	0.0505530	765.31	550.70	0.20100
8	IC5	7.738	0.186900	0.0241300	829.80	483.00	0.22200
9	NC5	4.285	0.078700	0.0181480	845.60	489.50	0.25400
10	C6	6.677	0.056720	0.0084850	914.20	439.70	0.30100
11	C7	3.533	0.012330	0.0036900	972.31	396.90	0.35000
12	C8	1.737	0.002620	0.0015080	1024.31	352.10	0.40200
13	C9	0.441	0.000280	0.0003500	1073.00	305.00	0.44600
14	C10	0.094	0.000030	0.0003190	1114.70	306.00	0.48900
15	C11	0.012	0.000000	0.0000000	1153.70	282.00	0.50100
16	C12	0.001	0.000000	0.0000000	1187.70	263.00	0.53900

IPH = .1  
 A = 0.30091      B1 = 0.08690  
 Q = -0.20645      RR = -0.02615  
 Z = 0.77779

IPH = 0  
 A = 1.55570      B1 = 0.18108  
 Q = 1.34183      RR = 0.28171  
 Z = 0.24333

COMPONENT	X	Y	K (EXP)	K (SRK)	K (MODEL)	TOLERANCE
CO2	0.000000	0.000000	0.0000000	1.2423290	1.2449930	0.0026636
N2	0.103000	0.944000	9.1650470	7.8123300	7.8883350	0.0760050
CO	27.410990	86.800350	3.1633360	2.6232790	2.6376110	0.0143319
C2	15.205990	8.197730	0.5364300	0.5480300	0.5486849	0.0006549
C3	13.795390	2.557410	0.1851800	0.1701400	0.1700028	-0.0001372
IC4	9.911000	0.657100	0.0736640	0.0611700	0.0610733	-0.0000967
NC4	9.995000	0.505800	0.0505530	0.0576200	0.0574678	-0.0001522
C5+	7.733000	0.186900	0.0569150	0.0548000	0.0505950	-0.00489220

FAAD = 0.2196110

TABLE 5.11 COMPARISON OF EQUILIBRIUM K VALUES FROM THE ORDINARY METHOD AND THE PROPOSED GENERAL MODEL FOR HEAVY-PSEUDOCOMPONENT (M=8) AT T = 15,000 °F P = 850.0 PSIA

THE AMOUNT OF COMPONENT, N = 16

T = -8.000 ° F      P = 1115.00 PSIA      R = 10.730 PSIA-CUBIC FT/LB-MOLE ° R

FEED RATE = 2.538890 MOLE/HK      V/F = 0.500000

LD	COMPONENT	X (%)	Y (%)	K (EXP)	TC (R)	PC (PSIA)	H
1	C02	0.589	7.097400	1.2049900	548.00	1073.00	0.22500
2	N2	0.472	4.210720	8.9210190	227.20	492.00	0.04000
3	C1	37.190	91.002390	2.4210660	243.30	673.10	0.01400
4	C2	4.604	3.046880	0.6548110	549.77	708.30	0.09900
5	C3	3.491	0.582920	0.1652300	665.95	617.40	0.15200
6	IC4	2.137	0.135280	0.0626500	734.65	529.10	0.18500
7	NC4	2.995	0.128320	0.0423870	765.21	550.70	0.20100
8	IC5	3.240	0.059490	0.0181670	829.80	483.00	0.22200
9	NC5	2.597	0.036710	0.0139860	845.60	489.50	0.25400
10	C6	7.432	0.048440	0.0064500	914.20	439.70	0.30100
11	C7	9.805	-0.025980	0.0026220	972.31	396.90	0.35000
12	C8	10.591	0.008570	0.0008010	1024.31	362.10	0.40200
13	C9	7.952	0.003420	0.0004300	1073.00	345.00	0.44600
14	C10	4.317	0.000910	0.0002100	1114.70	306.00	0.48900
15	C11	1.696	-0.000180	0.0001100	1153.70	282.00	0.50100
15	C12	0.894	0.000040	0.0000400	1187.70	263.00	0.53900

IPH = 1  
 A = 0.43526      B1 = 0.11906  
 Q = 0.33203      RR = 0.05187  
 Z = 0.65210

IPH = 0  
 A = 4.24025      B1 = -0.34505  
 Q = 3.77615      RR = 1.46308  
 Z = 0.41406

COMPONENT	X	Y	K(EXP)	K(SRK)	K(MODEL)	TOLERANCE
C02	0.583300	7.097400	1.2049900	1.0457050	1.0494320	0.0037231
N2	0.472000	4.210720	8.9210190	4.7024490	6.8360180	0.1335688
C1	37.190000	91.002390	2.4210660	2.1752090	2.1997140	0.0245047
C2	4.604000	3.046880	0.6548110	0.4810800	0.4824766	0.0012966
C3	3.491000	0.582920	0.1652300	0.1710500	0.1709797	-0.0000703
IC4	-2.137000	0.135280	0.0626500	0.8664000	0.0866004	-0.7797996
NC4	2.995000	0.128320	0.0423870	0.0607500	0.0605356	-0.0002144
C6+	-3.240000	0.059490	0.0428160	0.0067158	0.0005744	-0.0661454

-3AAD = -11.5140700

TABLE 5.12 COMPARISON OF EQUILIBRIUM K VALUES FROM THE ORDINARY METHOD  
 AND THE PROPOSED GENERAL MODEL FOR HEAVY-PSEUDOCOMPONENT (M=8)  
 AT T = -8.0 ° F      P = 1115.0 PSIA

T = 41.000 ° F      P = 300.00 PSIA      R = 10.730 PSIA-CUBIC FT/LB-MOLE °R

FEED RATE = 1.319670 MOLE/HR      V/F = 0.530000

TABLE 5.13 COMPARISON OF EQUILIBRIUM K VALUES FROM THE ORDINARY METHOD

AND THE PROPOSED GENERAL MODEL FOR LIGHT-PSEUDOCOMPONENT (C<sub>1+</sub>)

AT T = 41.0 ° F      P = 300.0 PSIA

ID	COMPONENT	X (I)	Y (T)	K (EXP)	TC (R)	PC (PSIA)	M
1	CO2	0.088	0.393300	4.4920450	348.00	1073.00	0.22500
2	N2	0.015	0.028000	55.1997900	227.20	492.00	0.04000
3	C1	3.564	85.016220	9.5165650	343.20	673.10	0.01400
4	C2	5.574	7.72530	1.3025060	549.77	708.30	0.09900
5	C3	3.238	3.167210	0.4025620	465.55	617.40	0.15200
6	IC4	7.904	1.225400	0.1540760	734.65	523.10	0.18500
7	NC4	3.630	0.910200	0.0944170	763.31	550.70	0.20100
8	IC5	3.678	0.273800	0.0283820	829.80	483.00	0.22200
9	NC5	6.855	0.137200	0.0202860	845.60	489.50	0.25400
10	C6	15.402	0.097600	0.0063310	914.20	439.70	0.30100
11	C7	11.893	0.012990	0.0010910	972.31	376.90	0.35000
12	C8	10.774	0.004340	0.0004020	1024.21	362.10	0.40200
13	C9	4.454	0.001120	0.0002510	1073.00	345.00	0.44600
14	C10	0.869	0.000080	0.0000920	1114.70	306.00	0.48900
15	C11	0.062	0.000000	0.0000000	1153.70	282.00	0.50100
16	C12	0.000	0.000000	0.0000000	1137.70	263.00	0.53900

IPH = 1  
 A = 0.08640      BI = 0.02819  
 J = 0.05741      RR = 0.00244  
 Z = 0.94187

IPH = 0  
 A = 1.34231      BI = 0.11571  
 J = 1.71321      RR = 0.21317  
 Z = 0.13340

COMPONENT	X	Y	K (EXP)	K (SFK)	K (MODEL)	TOLERANCE
CO2	0.000333	0.003753	4.4920450	2.4144590	3.5546090	0.2301893
N2	0.000150	0.003283	55.1997900	28.3205700	35.7977200	7.4707230
C1+	0.223763	0.961059	11.7086500	7.5650200	0.0232332	-7.9417860
IC4	0.079043	0.012254	0.1540760	0.1116500	0.0306944	-0.1009956
NC4	0.096303	0.005102	0.0944170	0.0390900	0.0214360	-0.0678034
IC5	0.016783	0.002738	0.0283820	0.0337400	0.0058075	-0.0279325
NC5	0.063550	0.001392	0.0202860	0.0240000	0.0014357	-0.0227143
C6	0.154023	0.000975	0.0063310	0.0067600	0.0003891	-0.0063709
C7	0.113933	0.000100	0.0010910	0.0015000	0.0001042	-0.0017958
C8	0.107743	0.000043	0.0004020	0.0005100	0.0000290	-0.0004820
C9	0.044540	0.000011	0.0002510	0.0001100	0.00001264	-0.0000196
C10	0.003693	0.000001	0.0000920	0.0000400	0.00000345	-0.0000055

ΣAAD = 1.8992170

THE AMOUNT OF COMPONENT,N = 16

T = 30.000 ° F      P = 300.00 PSIA      R = 10.730 PSIA-CUBIC FT/LB-MOLE R

FEED RATE = 3.503350 MOLE/HR      V/F = 0.500000

TABLE 5.14 COMPARISON OF EQUILIBRIUM K VALUES FROM THE ORDINARY METHOD AND THE PROPOSED GENERAL MODEL FOR LIGHT-PSEUDOCOMPONENT (C<sub>1+</sub>) AT T = 30.0 ° F    P = 300.0 PSIA

ID	COMPONENT	X (%)	Y (%)	K (EXP)	TC (R)	PC (PSIA)	h
1	CO2	0.065	0.250400	3.8523080	548.00	1073.00	0.22500
2	N2	0.019	0.848000	44.6315700	227.20	492.00	0.04000
3	C1	9.522	84.699430	8.8876580	343.20	673.10	0.01400
4	C2	3.764	7.922510	2.1028110	549.17	708.30	0.09900
5	C3	9.770	1.710010	0.3802220	665.55	617.40	0.15200
6	IC4	13.308	1.187900	0.1151350	734.65	529.10	0.18500
7	NC4	11.833	0.818400	0.0691020	765.31	550.70	0.20100
8	IC5	12.244	0.310400	0.0253290	829.80	483.00	0.22200
9	NC5	7.713	0.133500	0.0172940	845.60	489.50	0.25400
10	C6	14.212	0.085180	0.0059880	914.20	439.70	0.30100
11	C7	9.321	0.020650	0.0022140	972.31	396.70	0.35000
12	C8	7.332	0.004650	0.0006340	1024.31	362.10	0.40200
13	C9	2.898	0.000860	0.0000000	1073.00	345.00	0.44600
14	C10	0.848	0.000100	0.0000000	1114.70	306.00	0.48900
15	C11	0.149	0.000010	0.0000000	1153.70	282.00	0.50100
16	C12	0.000	0.000000	0.0000000	1187.70	263.00	0.53900

[PH = 1  
 A = 0.07990      B1 = 0.02754  
 Q = 0.05160      RR = 0.00220  
 Z = 0.94807

[PH = 0  
 A = 2.21589      B1 = 0.11405  
 Q = 2.08879      RR = 0.25281  
 Z = 0.12784

COMPONENT	X	Y	K(EXP)	K(SRK)	K(MODEL)	TOLERANCE
CO2	0.000650	0.002504	3.8523080	2.1380300	3.7700510	0.6320219
N2	0.000190	0.008480	44.6315700	27.6372500	53.5111500	25.8538500
C1+	0.451970	0.983462	11.5549300	7.5251300	0.0002042	-7.5249250
IC5	0.122440	0.003104	0.0253290	0.0269300	0.0002934	-0.0266866
NC5	0.077130	0.001335	0.0172940	0.0191300	0.0000672	-0.0190128
C6	0.142120	0.000852	0.0059880	0.0052000	0.0000193	-0.0051007
C7	0.0933210	0.000206	0.0022140	0.0013800	0.0015532	0.0001732
C8	0.0733200	0.00007	0.0006340	0.0001500	0.0003998	0.0000499
C9	0.0289300	0.000000	0.0000000	0.0000000	0.0000934	0.0000094
C10	0.0084800	0.000001	0.0000000	0.0000000	0.0000289	0.0000089
%AAC =						0.5104037

THE AMOUNT OF COMPONENT, N = 16

T = 30.000 °F P = 300.00 PSIA R = 10.730 PSIA-CUBIC FT/LB-MOLE R

FEED RATE = 0.790030 MOLE/TR V/F = 0.500000

ID	COMPONENT	X (%)	Y (%)	K (EXP)	TC (F)	PC (PSIA)	W
1	CO2	0.056	0.177200	3.1642850	548.00	1073.00	0.22500
2	N2	0.029	0.937020	32.3110300	227.20	492.00	0.04000
3	C1	10.818	84.920880	7.8432220	343.30	673.10	0.01400
4	C2	7.039	7.785640	1.1052010	549.77	709.30	0.09300
5	C3	10.875	3.664860	0.3367250	665.95	617.40	0.15200
6	IC4	9.692	1.150220	0.1185790	734.65	529.10	0.18500
7	NC4	11.420	7.808010	0.0076950	765.31	550.70	0.20100
8	IC5	10.886	0.303610	0.0270670	329.80	483.00	0.22200
9	NC5	7.205	0.127000	0.0170120	845.60	489.50	0.25400
10	C6	13.468	0.091790	0.0063100	914.20	439.70	0.30100
11	C7	8.920	0.027400	0.0033650	972.31	396.90	0.35000
12	C8	6.328	0.005610	0.0003850	1024.21	362.10	0.40200
13	C9	2.411	0.000640	0.0002640	1073.00	345.00	0.44600
14	C10	0.727	0.000110	0.0000000	1114.70	306.00	0.48900
15	C11	0.127	0.000010	0.0000000	1153.70	282.00	0.50100
16	C12	0.000	0.000000	0.0000000	1187.70	263.00	0.53900

IPH = 1  
 A = 0.07386 B1 = 0.02754  
 Q = 0.05156 RR = 0.00220  
 Z = 0.94012

IPH = 0  
 A = 1.93333 B1 = 0.10790  
 Q = 1.81379 RR = 0.20860  
 Z = 0.12222

COMPONENT	X	Y	K (EXP)	K (SRK)	K (MODEL)	TOLERANCE
CO2	0.000560	0.001772	3.1642850	3.1124690	3.6426140	0.5301447
N2	0.000290	0.009370	32.3110300	26.9741900	49.1842100	22.2100200
C1+	0.099440	0.983296	9.4114220	7.4175790	0.0015606	-7.4160190
IC5	0.103360	0.003036	0.0278670	0.0265500	0.0002723	-0.0260777
NC5	0.072050	0.001270	0.0176120	0.0191100	0.0000624	-0.0170476
C6	0.134680	0.000918	0.0063100	0.0051300	0.0000175	-0.0051125
C7	0.039200	0.000274	0.0030690	0.0031800	0.0014556	0.0000756
C8	0.063280	0.000056	0.0003850	0.0003500	0.0003709	0.0000209
C9	0.024110	0.000036	0.0002640	0.0000000	0.0000914	0.0000014
C10	0.007270	0.000011	0.0000000	0.0000200	0.0000261	0.0000011
ΣAOC =						0.5094187

TABLE 5.15 COMPARISON OF EQUILIBRIUM K VALUES FROM THE ORDINARY METHOD AND THE PROPOSED GENERAL MODEL FOR LIGHT-PSEUDOCOMPONENT (C<sub>1+</sub>) AT T = 30.0 °F P = 300.0 PSIA

T = -10.000 ° F      P = 800.00 PSIA      R = 10.730 PSIA-CUBIC FT/L3-MOLE<sup>0</sup>R

FEED RATE = 2.543590 MOLE/HR      V/F = 0.500000

ID	COMPONENT	X (%)	Y (%)	K (EXP)	TC (R)	PC (PSIA)	W
1	CO2	0.143	0.373660	0.2613000	548.00	1073.00	-0.22500
2	N2	0.006	0.205190	34.1983300	227.20	492.00	0.04000
3	C1	31.340	97.267850	3.0990310	243.30	-673.10	0.01400
4	C2	2.363	2.092600	0.8341960	549.77	709.30	0.09900
5	C3	0.169	0.221010	0.1655930	665.95	617.40	0.15200
6	IC4	0.056	0.005200	0.0320640	734.65	529.10	0.18500
7	NC4	0.177	0.205800	0.0384300	765.31	-550.70	0.20100
8	IC5	0.605	0.200700	0.0110600	829.80	493.00	0.22200
9	NC5	0.498	0.003900	0.0078180	845.60	489.50	0.25400
10	C6	1.702	0.003910	0.022930	914.20	439.70	0.30100
11	C7	3.929	0.003010	0.0007640	972.31	396.90	0.35000
12	C8	0.857	0.001580	0.0002700	1024.31	362.10	0.40200
13	C9	14.467	0.003950	0.0000600	1073.00	-345.00	0.44600
14	C10	25.830	0.000650	0.0000230	1114.70	306.00	0.48900
15	C11	10.695	0.000110	0.0000100	-1153.70	-292.00	0.50100
16	C12	1.902	0.000010	0.0000050	1187.70	263.00	0.53900

IPH = 1  
 A = 0.25503      B1 = 0.07862  
 Q = 0.17023      RR = 0.02005  
 Z = 0.32278

IPH = 0  
 A = 4.42351      B1 = 0.30475  
 Q = 4.02588      RR = 1.34807  
 Z = 0.35505

TABLE 5.16 COMPARISON OF EQUILIBRIUM K VALUES FROM THE ORDINARY METHOD  
 AND THE PROPOSED GENERAL MODEL FOR LIGHT-PSEUDOCOMPONENT (C<sub>1+</sub>)  
 AT T = -10.0 ° F    P = 800.0 PSIA

COMPONENT	X	Y	K (EXP)	K (SRK)	K (MODEL)	TOLERANCE
CO2	0.001430	0.003737	0.2613000	1.1140600	1.0910190	-0.0230417
N2	0.000060	0.002052	34.1983300	10.7217500	10.1652300	-0.5565138
C1+	0.333720	0.953815	4.1400100	2.8262500	2.5961400	-0.2301092
IC4	0.000560	0.000052	0.0220640	0.0461500	0.0033645	-0.0304855
NC4	0.001770	0.000068	0.0384300	0.0306800	0.0059204	-0.0248596
IC5	0.006050	0.000067	0.0110600	0.0121500	0.0015496	-0.0107004
NC5	0.004980	0.000039	0.0078180	0.0084400	0.0001316	-0.0081084
C6	0.017020	0.000039	0.0229300	0.0224100	0.0000374	-0.0000326
C7	0.039290	0.000030	0.0007640	0.0006900	0.0000244	-0.0000656
C8	0.008570	0.000016	0.0002700	0.0001900	0.0000065	-0.0000825
C9	0.144670	0.000030	0.0000600	0.0000500	0.0000041	-0.0000019
C10	0.258300	0.000007	0.0000230	0.0000100	0.0000013	0.0000035
SAAC =						0.711051E

THE AMOUNT OF CONDENSATE = 16

T = 15.000 °F      P = 850.00 PSIA      R = 10.733 PSIA-CUBIC FT/LB-MOLE<sup>OR</sup>

FEED RATE = 1.357700 MOLE/HR      V/F = 0.500000

ID	COMPONENT	X (I)	Y (I)	K (EXP)	TC (R)	PC (PSIA)	W
1	C02	0.000	0.000000	0.000000	548.00	1073.00	0.22500
2	H2	0.103	0.944000	9.1650470	227.20	492.00	0.04000
3	C1	27.411	36.000350	0.1633360	343.30	673.10	0.01400
4	C2	15.266	8.197730	0.5364300	549.77	708.30	0.09900
5	C3	13.796	2.557410	0.1851800	665.95	617.40	0.15200
6	IC4	3.911	0.657100	0.0736640	734.65	529.10	0.18500
7	NC4	7.995	0.505800	0.0505530	765.31	550.70	0.26100
8	IC5	7.738	0.180900	0.0241300	829.80	483.00	0.22200
9	NC5	4.285	0.078700	0.0183480	845.60	487.50	0.25400
10	C6	6.677	1.056720	0.0084850	914.20	439.70	0.30100
11	C7	3.533	0.012330	0.0034900	972.31	396.90	0.35000
12	C8	1.737	0.002620	0.0015080	1024.31	362.10	0.40200
13	C9	0.441	0.000280	0.0000350	1173.00	345.00	0.44600
14	C10	0.094	0.000030	0.0000190	1114.70	306.00	0.48900
15	C11	0.012	0.000000	0.0000000	1153.70	282.00	0.50100
16	C12	0.001	0.000000	0.0000000	1187.70	263.00	0.52900

IPH = 1  
 A = 0.25914      B1 = 0.08196  
 Q = 0.17046      RR = 0.02124  
 Z = 0.82456

IPH = 0  
 A = 2.32002      B1 = 0.21171  
 Q = 2.06349      RR = 0.49117  
 Z = 0.26269

COMPONENT	X	Y	K (EXP)	K (SRK)	K (MODEL)	TOLERANCE
C02	0.000000	0.000000	0.000000	1.2423290	1.2739090	0.0365801
H2	0.001030	0.009440	9.1650470	7.8123300	10.7536100	2.9412830
C1	0.274110	0.3600350	0.1633360	2.6232790	0.2475736	-2.3757050
C2	0.152660	0.0819773	0.5364300	0.0311700	0.0149423	-0.0662277
IC4	0.039110	0.0065710	0.0736640	0.0576200	0.0106674	-0.00469526
NC4	0.079950	0.0050580	0.0505530	0.0251900	0.0031917	-0.0225981
IC5	0.077380	0.0018090	0.0241300	0.0191700	0.0027720	-0.00183907
NC5	0.042850	0.0007870	0.0183480	0.0065500	0.002332	-0.0063168
C6	0.066770	0.0005670	0.0084850	0.0023300	0.0000722	-0.0021778
C7	0.035330	0.0001230	0.0034900	0.0007400	0.0000223	-0.00007177
C8	0.017370	0.0000260	0.0015080	0.0001500	0.0000333	-0.00001117
C9	0.004410	0.0000028	0.0000350	0.0000350	0.0000450	-0.0000350
C10	0.000940	0.0000003	0.0000190			
ΣAAD =						1.3627320

TABLE 5.17 COMPARISON OF EQUILIBRIUM K VALUES FROM THE ORDINARY METHOD AND THE PROPOSED GENERAL MODEL FOR LIGHT-PSEUDOCOMPONENT (C<sub>1+</sub>) AT T = 15.0°F P = 850.0 PSIA

THE AMOUNT OF COMPONENT, N = 16

T = -8.000 ° F      P = 1115.00 PSIA      R = 10.730 PSIA-CUBIC FT/LB-MOLE<sup>OR</sup>

FEED RATE = 2.638890 MOLE/HR      V/F = 0.500000

TABLE 5.18 COMPARISON OF EQUILIBRIUM K VALUES FROM THE ORDINARY METHOD AND THE PROPOSED GENERAL MODEL FOR LIGHT-PSEUDOCOMPONENT (C<sub>1+</sub>) AT T = -8.0° F    P = 1115.0 PSIA

ID	COMPONENT	X (%)	Y (%)	K (EXP)	TC (R)	PC (PSIA)	K
-1	CO2	0.589	7.097400	1.2047900	543.00	1073.00	0.22500
2	N2	0.472	4.210720	3.9210190	227.20	492.00	0.04000
3	C1	37.190	91.002370	2.4210600	343.30	673.10	0.01400
4	C2	4.604	3.046880	0.6541110	549.77	708.30	0.09900
5	C3	3.491	0.582920	0.1652300	605.95	617.40	0.15200
6	IC4	2.137	0.133280	0.0626500	734.65	529.10	0.18500
-7	NC4	2.995	0.129320	0.0423870	765.21	550.70	0.20100
8	IC5	3.240	0.050490	0.0181670	829.80	483.00	0.22200
9	NC5	2.597	0.036710	0.0139960	845.60	409.50	0.25400
10	C6	7.432	0.043440	0.0064500	914.20	439.70	0.30100
11	C7	9.809	0.025930	0.0026220	972.31	396.90	0.35000
12	C8	10.591	0.009570	0.0007010	1024.31	362.10	0.40200
13	C9	7.952	0.003420	0.0004300	1073.00	345.00	0.44600
14	C10	4.317	0.000710	0.0002100	1114.70	306.00	0.48900
15	CI1	1.696	0.000180	0.0001100	1153.70	282.00	0.50100
16	CI2	0.894	0.000040	0.0000400	1187.70	263.00	0.53900

IPH = 1  
 A = 0.41413      DI = 0.11685  
 J = 0.28363      RR = 0.04839  
 Z = 0.69087

IPH = 0  
 A = 5.20922      BI = 0.38360  
 J = 4.67848      RR = 1.99024  
 Z = 0.45098

COMPONENT	X	Y	K (EXP)	K (SFK)	K (MJOEL)	TOLERANCE
CO2	0.005390	0.070974	1.2045900	1.0457390	1.0367400	-0.0087693
N2	0.004720	0.042107	3.9210190	6.7024490	7.2626000	0.5601511
C1	0.452850	0.946322	3.2411060	2.1752090	0.9450583	-1.2301510
IC4	0.021370	0.001358	0.0626500	0.8664000	0.0148438	-0.8515512
NC4	0.029950	0.001283	0.0423870	0.0607500	0.0108155	-0.0499245
IC5	0.032400	0.000595	0.0181670	0.0305400	0.0035639	-0.0269761
NC5	0.025970	0.000367	0.0139960	0.0224000	0.0008228	-0.0213772
C6	0.074320	0.000484	0.0064500	0.0085400	0.0002310	-0.0082590
C7	0.098050	0.000260	0.0026220	0.0032900	0.0001007	-0.0031851
C8	0.105910	0.000096	0.0008010	0.0012100	0.0000357	-0.0011743
C9	0.079520	0.000034	0.0004300	0.0004400	0.0000135	-0.0001265
C10	0.043170	0.000007	0.0002100	0.0001700	0.0000121	-0.0000489
ΣAAD =						8.0236350



TABLE 5.19 COMPARISON OF THE INITIAL PHASE COMPOSITION VALUES OF THE EXPERIMENT AND THE ESTIMATION VALUES FOR HEAVY-PSEUDOCOMPONENT AT T = -8.0°F P = 1115.0 PSIA

Component	$x_1$	$y_1$	$z_1$	$K_1$					
CO <sub>2</sub>	0.589	0.70974	0.70973	0.0	T = -8.0°F      P = 1115.0 psia # mole/hr L = 0.00039 V = 2.6385 F = L + V = 2.63889				
N <sub>2</sub>	0.472	4.21072	4.21017	0.0					
C <sub>1</sub>	37.190	91.00240	90.99455	2.421066					
C <sub>2</sub>	4.604	3.04688	3.04710	0.654811					
C <sub>3</sub>	3.491	0.58292	0.58334	0.165230					
					$x_1/\Sigma x_1$	$y_1/\Sigma y_1$	$z_1/\Sigma z_1$	$z_1/K_1$	$\frac{z_1/K_1}{\Sigma(z_1/K_1)}$
i-C <sub>4</sub>	2.137	0.13528	0.13558	0.062650	0.040	0.302	0.298	2.16409	0.052
n-C <sub>4</sub>	2.995	0.12832	0.12874	0.042387	0.056	0.287	0.283	3.03725	0.072
i-C <sub>5</sub>	3.240	0.05949	0.05995	0.018167	0.060	0.133	0.132	3.29994	0.079
n-C <sub>5</sub>	2.597	0.03671	0.03708	0.013986	0.048	0.082	0.082	2.65122	0.063
C <sub>6</sub>	7.432	0.04844	0.04952	0.00645	0.138	0.108	0.109	7.67752	0.183
C <sub>7</sub>	9.805	0.02598	0.02741	0.002622	0.182	0.058	0.060	10.45385	0.249
C <sub>8</sub>	10.591	0.00857	0.01012	0.000801	0.197	0.019	0.022	12.63421	0.301
C <sub>9</sub>	7.592	0.00342	0.00458	0.0					
C <sub>10</sub>	4.317	0.00091	0.00154	0.0					
C <sub>11</sub>	1.696	0.00018	0.00043	0.0					
C <sub>12</sub>	0.894	0.00004	0.00017	0.0					
	= 53.656	= 0.44734	= 0.45512					= 41.91808	

TABLE 5.20 COMPARISON OF THE INITIAL PHASE COMPOSITION VALUES OF THE EXPERIMENT AND THE ESTIMATION VALUES  
FOR HEAVY-PSEUDOCOMPONENT AT T = 15.0°F P = 700.0 PSIA

Component	$x_i$	$y_i$	$z_i$	$K_i$	T = 15 °F      P = 700.0 psia # mole/hr    L = 0.03313 V = 1.3193 F = L + V = 1.35243					
					$x_i/\Sigma x_i$	$y_i/\Sigma y_i$	$z_i/\Sigma z_i$	$z_i/K_i$	$\frac{z_i/K_i}{\Sigma(z_i/K_i)}$	
CO <sub>2</sub>	0.0	0.0	0.0	0.0						
N <sub>2</sub>	0.078	1.01011	0.98729	0.0						
C <sub>1</sub>	22.741	86.63524	85.07018	3.806192						
C <sub>2</sub>	14.343	8.29232	8.44052	0.577630						
C <sub>3</sub>	13.859	2.55278	2.82973	0.184028						
i-C <sub>4</sub>	9.618	0.67387	0.89296	0.069999	0.196	0.446	0.334	12.75675	0.129	
n-C <sub>4</sub>	11.001	0.52806	0.78460	0.047956	0.225	0.350	0.294	16.36083	0.166	
i-C <sub>5</sub>	8.795	0.17622	0.38734	0.020017	0.180	0.117	0.145	19.35055	0.196	
n-C <sub>5</sub>	5.131	0.06521	0.18930	0.012696	0.105	0.043	0.071	14.91021	0.151	
C <sub>6</sub>	7.830	0.05462	0.24508	0.00696	0.160	0.036	0.092	35.16717	0.356	
C <sub>7</sub>	3.903	0.00939	0.10477	0.0		0.006	0.039			
C <sub>8</sub>	2.001	0.00193	0.05090	0.0		0.001	0.019			
C <sub>9</sub>	0.552	0.00021	0.01373	0.0		0.0001	0.005			
C <sub>10</sub>	0.123	0.00002	0.00303	0.0						
C <sub>11</sub>	0.022	0.00000	0.00055	0.0						
C <sub>12</sub>	0.001	0.00000	0.00003	0.0						
= 48.977    = 1.50953    = 2.67229					=98.54551					

TABLE 5.21 COMPARISON OF THE INITIAL PHASE COMPOSITION VALUES OF THE EXPERIMENT AND THE ESTIMATION VALUES FOR HEAVY-PSEUDOCOMPONENT AT  $T = 15.0^\circ\text{F}$   $P = 200.0.0$  PSIA

Component	$x_1$	$y_1$	$z_1$	$K_1$					
$\text{CO}_2$	0.0	0.0	0.0	0.0					
$\text{N}_2$	0.019	0.93738	0.93205	0.0					
$\text{C}_1$	7.577	85.76607	85.31217	11.311396					
$\text{C}_2$	6.028	8.39818	8.38443	1.392303					
$\text{C}_3$	8.828	2.83914	2.87391	0.321401					
					$x_1/\Sigma x_1$	$y_1/\Sigma y_1$	$z_1/\Sigma z_1$	$z_1/K_1$	$\frac{z_1/K_1}{\Sigma(z_1/K_1)}$
i-C <sub>4</sub>	7.698	0.82823	0.86811	0.107526	0.099	0.402	0.348	8.07349	0.100
n-C <sub>4</sub>	10.202	0.67287	0.72819	0.065914	0.132	0.327	0.292	11.04758	0.138
i-C <sub>5</sub>	11.323	0.30302	0.36700	0.026744	0.146	0.147	0.147	13.7227	0.171
n-C <sub>5</sub>	8.028	0.11094	0.15691	0.013811	0.104	0.054	0.063	11.36123	0.141
$\text{C}_6$	18.647	0.11518	0.22277	0.006173	0.240	0.056	0.089	36.08078	0.449
$\text{C}_7$	12.284	0.02390	0.09508	0.0		0.012	0.038		
$\text{C}_8$	6.917	0.00459	0.04472	0.0		0.002	0.018		
$\text{C}_9$	2.018	0.00045	0.01216	0.0		0.0002	0.005		
$\text{C}_{10}$	0.382	0.00003	0.00225	0.0					
$\text{C}_{11}$	0.042	0.00000	0.00025	0.0					
$\text{C}_{12}$	0.006	0.00000	0.00004	0.0					

$T = 15.0^\circ\text{F}$   $P = 200.0$  psia

# mole/hr  $L = 0.00770$

$V = 1.3193$

$F = L + V = 1.32700$

$\Sigma x_1 = 77.547$   $\Sigma y_1 = 2.05921$   $\Sigma z_1 = 2.49748$

$\Sigma(z_1/K_1) = 80.2928$

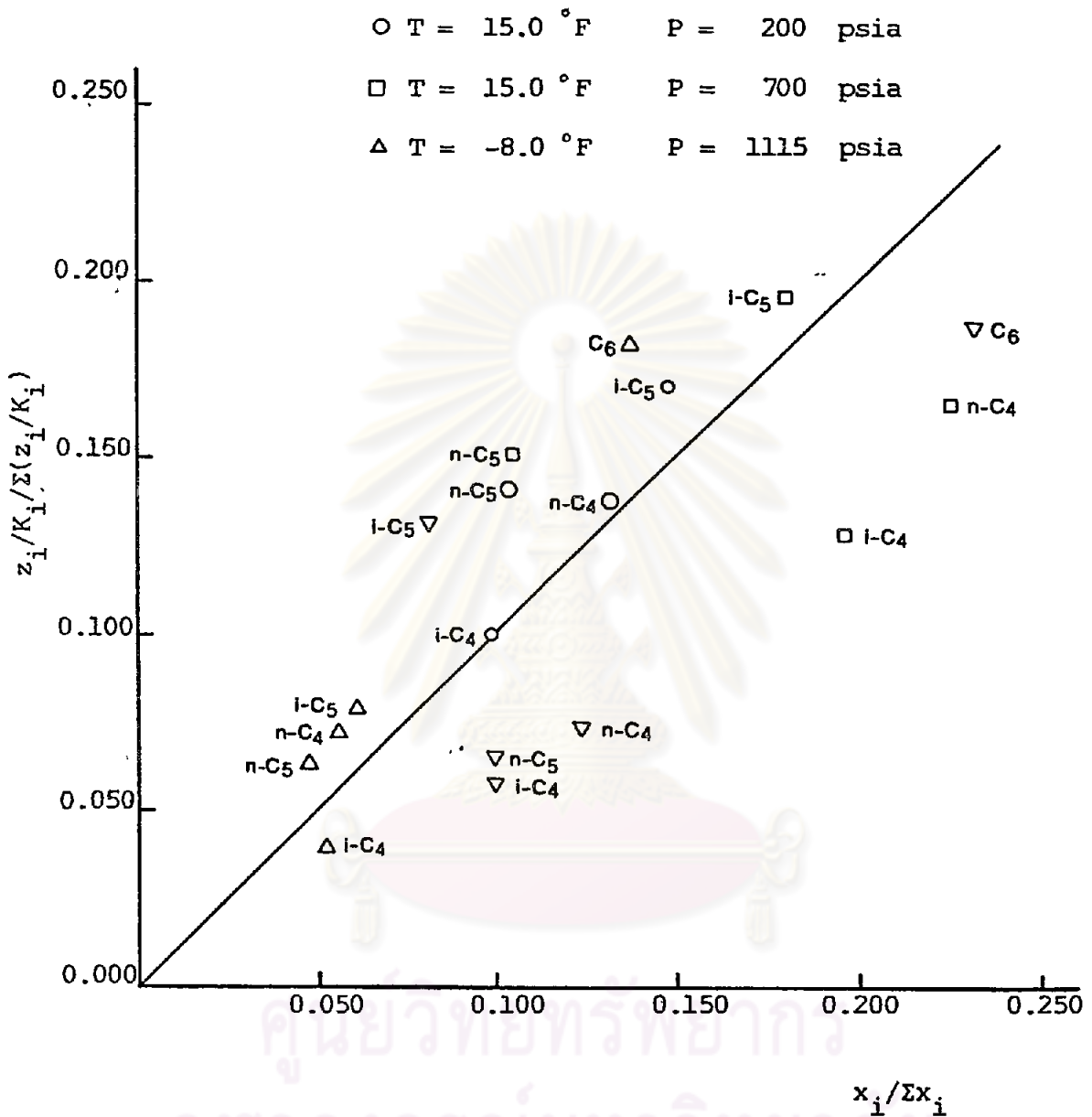


FIGURE 5.3 THE COMPARISON OF EXPERIMENTAL VALUES (ABCISSA) AND ESTIMATED VALUES (ORDINATE) IN THE LIQUID PHASE OF HEAVY-PSEUDOCOMPONENTS

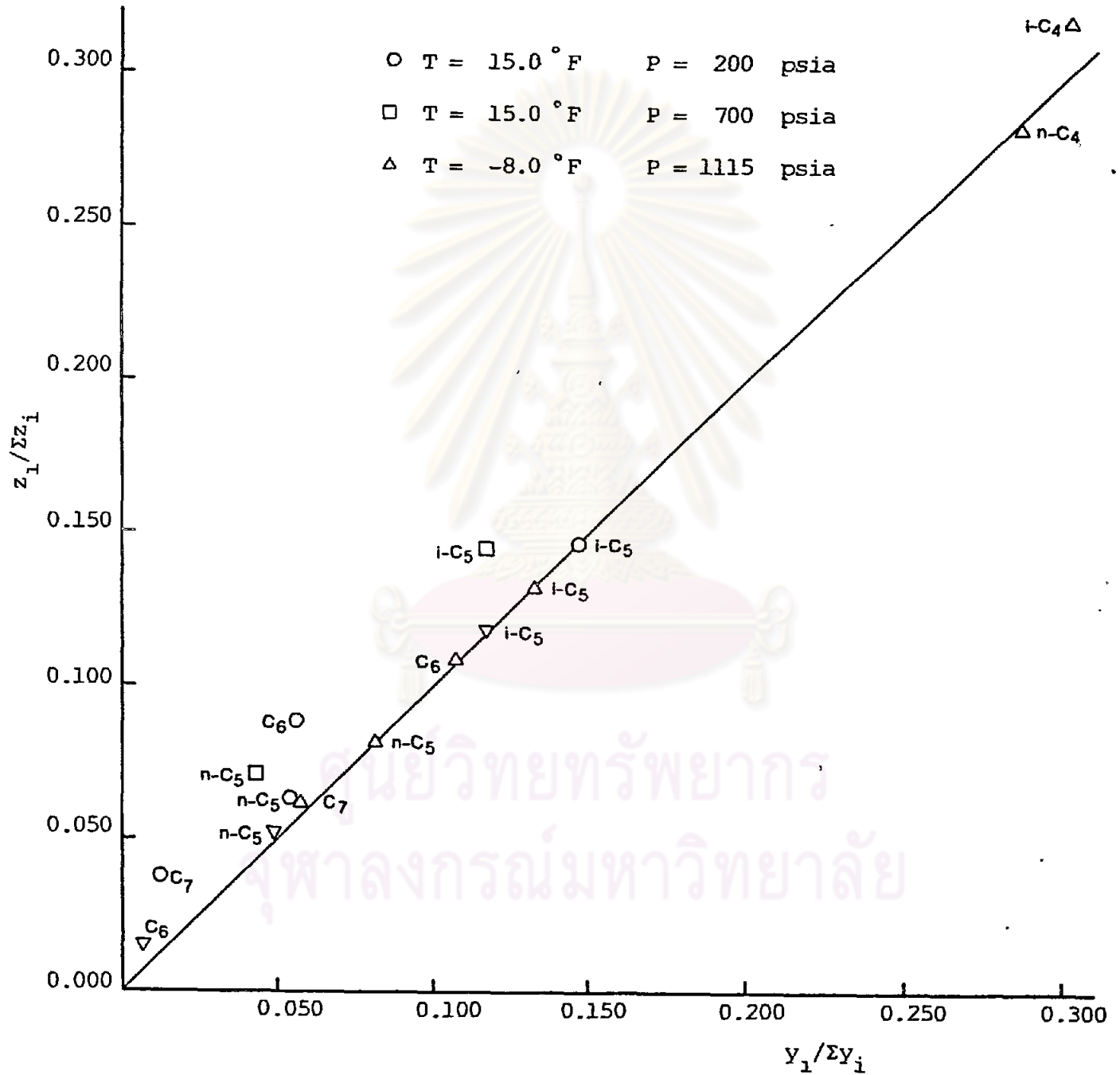


FIGURE 5.4 THE COMPARISON OF EXPERIMENTAL VALUES (ABCISSA) AND ESTIMATED VALUES (ORDINATE) IN THE VAPOR PHASE OF HEAVY-PSEUDOCOMPONENTS

TABLE 5.22 COMPARISON OF THE INITIAL PHASE COMPOSITION VALUES OF THE EXPERIMENT AND THE ESTIMATION VALUES FOR LIGHT-PSEUDOCOMPONENT AT T = 15.0°F P = 200.0 PSIA

Component	$x_1$	$y_1$	$z_1$	$K_1$	$x_1/\Sigma x_1$	$y_1/\Sigma y_1$	$z_1/\Sigma z_1$	$z_1/K_1$	$\frac{z_1/K_1}{\sum (z_1/K_1)}$
CO <sub>2</sub>	0.0	0.0	0.0	0.0	-	-	-	-	-
N <sub>2</sub>	0.019	0.93738	0.93205	0.0	-	-	-	-	-
C <sub>1</sub>	7.577	85.76607	85.31217	11.311396	0.338	0.884	0.883	7.54214	0.335
C <sub>2</sub>	6.028	8.39818	8.38443	1.392303	0.269	0.086	0.087	6.02198	0.268
C <sub>3</sub>	8.828	2.83914	2.87391	0.321401	0.394	0.029	0.030	8.94182	0.397
	22.433	97.0039	96.05705					22.50594	
i-C <sub>4</sub>	7.698	0.82823	0.86811	0.107526	T = 15 °F      P = 200 psia # mole/hr L = 0.0077 V = 1.3193 F = L + V = 1.3270				
n-C <sub>4</sub>	10.202	0.67287	0.72819	0.065914					
i-C <sub>5</sub>	11.323	0.30302	0.36700	0.026744					
n-C <sub>5</sub>	8.028	0.11094	0.15691	0.013811					
C <sub>6</sub>	18.647	0.11518	0.22277	0.006173					
C <sub>7</sub>	12.284	0.02390	0.09508	0.0					
C <sub>8</sub>	6.917	0.00459	0.04472	0.0					
C <sub>9</sub>	2.018	0.00045	0.01216	0.0					
C <sub>10</sub>	0.382	0.00003	0.00225	0.0					
C <sub>11</sub>	0.042	0.00000	0.00025	0.0					
C <sub>12</sub>	0.006	0.00000	0.00004	0.0					

TABLE 5.23 COMPARISON OF THE INITIAL PHASE COMPOSITION VALUES OF THE EXPERIMENT AND THE ESTIMATION VALUES FOR LIGHT-PSEUDOCOMPONENT AT  $T = 15.0^{\circ}\text{F}$   $P = 700.0$  PSIA

Component	$x_1$	$y_1$	$z_1$	$K_1$	$x_1/\Sigma x_1$	$y_1/\Sigma y_1$	$z_1/\Sigma z_1$	$z_1/K_1$	$\frac{z_1/K_1}{\sum (z_1/K_1)}$
CO <sub>2</sub>	0.0	0.0	0.0	0.0	-	-	-	-	-
N <sub>2</sub>	0.078	1.01011	0.98729	0.0	-	-	-	-	-
C <sub>1</sub>	22.741	86.63524	85.07018	3.806192	0.446	0.889	0.883	22.30547	0.429
C <sub>2</sub>	14.343	8.29232	8.44052	0.577630	0.282	0.085	0.088	14.61233	0.279
C <sub>3</sub>	13.859	2.55278	2.82973	0.184028	0.272	0.026	0.029	15.37663	0.294
	50.943	97.48034	96.34043					52.33943	
i-C <sub>4</sub>	9.6184	0.67387	0.89296	0.069999	$T = 15.0^{\circ}\text{F}$ $P = 700.0$ psia # mole/hr $L = 0.03313$ $V = 1.3193$ $F = L + V = 1.35243$				
n-C <sub>4</sub>	11.001	0.52806	0.78460	0.047956					
i-C <sub>5</sub>	8.795	0.17622	0.38734	0.020017					
n-C <sub>5</sub>	5.131	0.06521	0.18930	0.012696					
C <sub>6</sub>	7.830	0.05462	0.024508	0.006969					
C <sub>7</sub>	3.903	0.00939	0.10477	0.0					
C <sub>8</sub>	2.001	0.00193	0.05090	0.0					
C <sub>9</sub>	0.552	0.00021	0.01373	0.0					
C <sub>10</sub>	0.123	0.00002	0.00303	0.0					
C <sub>11</sub>	0.022	0.00000	0.00055	0.0					
C <sub>12</sub>	0.001	0.00000	0.00003	0.0					

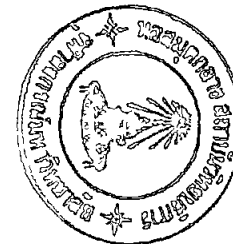


TABLE 5.24 COMPARISON OF THE INITIAL PHASE COMPOSITION VALUES OF THE EXPERIMENT AND THE ESTIMATION VALUES FOR LIGHT-PSEUDOCOMPONENT AT  $T = -8.0^{\circ}\text{F}$   $P = 1115.0$  PSIA

Component	$x_i$	$y_i$	$z_i$	$K_i$	$x_i/\sum x_i$	$y_i/\sum y_i$	$z_i/\sum z_i$	$z_i/K_i$	$\frac{z_i/K_i}{\sum (z_i/K_i)}$
CO <sub>2</sub>	0.589	0.70974	0.70973	0.0	-	-	-	-	-
N <sub>2</sub>	0.472	4.21072	4.21017	0.0	-	-	-	-	-
C <sub>1</sub>	37.190	91.00240	90.99455	2.421066	0.821	0.962	0.962	39.58450	0.821
C <sub>2</sub>	4.604	3.04688	3.04710	0.654811	0.102	0.032	0.032	4.6534	0.102
C <sub>3</sub>	3.491	0.58292	0.58334	0.165230	0.077	0.006	0.006	3.53047	0.077
	45.285	94.6322	94.62499					45.76837	
i-C <sub>4</sub>	2.137	0.13528	0.13558	0.062650					
n-C <sub>4</sub>	2.995	0.12832	0.12874	0.042387					
i-C <sub>5</sub>	3.240	0.05949	0.05995	0.018167					
n-C <sub>5</sub>	2.597	0.03671	0.03708	0.013986					
C <sub>6</sub>	7.432	0.04844	0.04952	0.00645					
C <sub>7</sub>	9.805	0.02598	0.02741	0.002622					
C <sub>8</sub>	10.591	0.00857	0.01012	0.000801					
C <sub>9</sub>	7.952	0.00342	0.00458	0.0					
C <sub>10</sub>	4.317	0.00091	0.00154	0.0					
C <sub>11</sub>	1.696	0.00018	0.00043	0.0					
C <sub>12</sub>	0.894	0.00004	0.00017	0.0					

$T = -8.0^{\circ}\text{F}$        $P = 1115.0$  psia  
 # mole/hr     $L = 0.00039$   
                    $V = 2.6385$   
 $F = L + V = 2.63889$



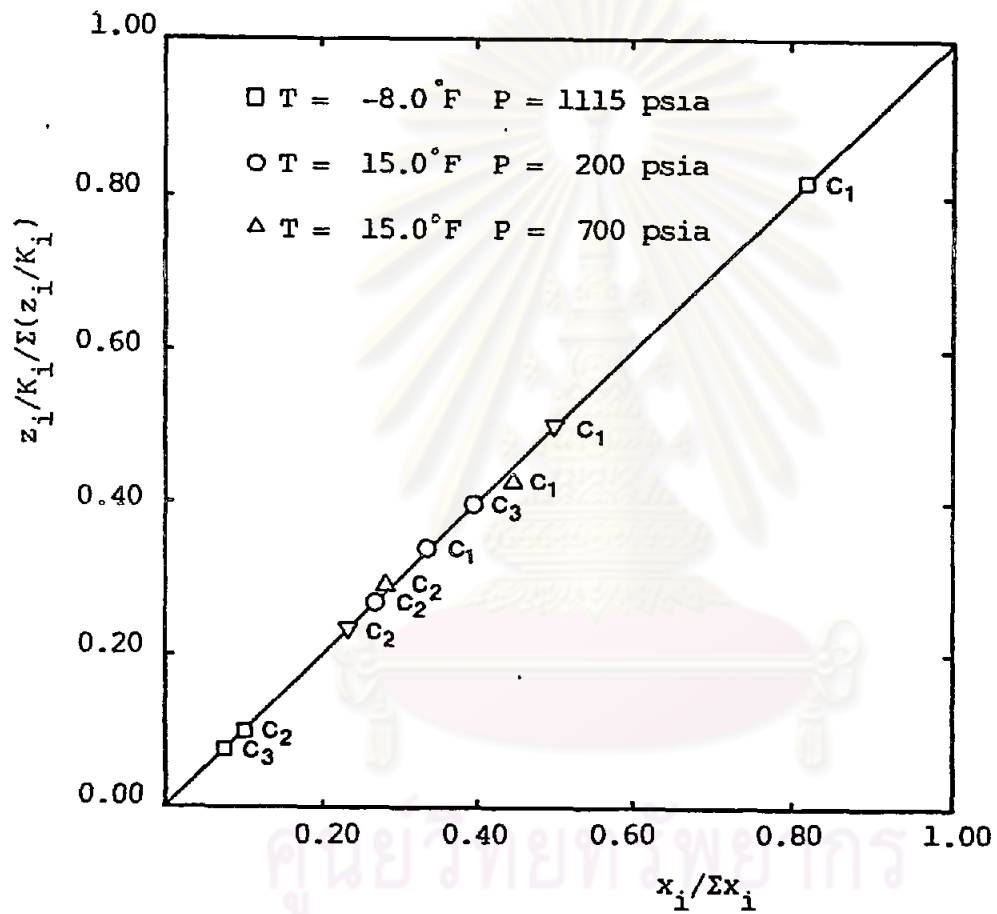


FIGURE 5.5 THE COMPARISON OF EXPERIMENTAL VALUES (ABCISSA) AND ESTIMATED VALUES (ORDIANTE) IN THE LIQUID PHASE OF LIGHT-PSEUDOCOMPONENTS

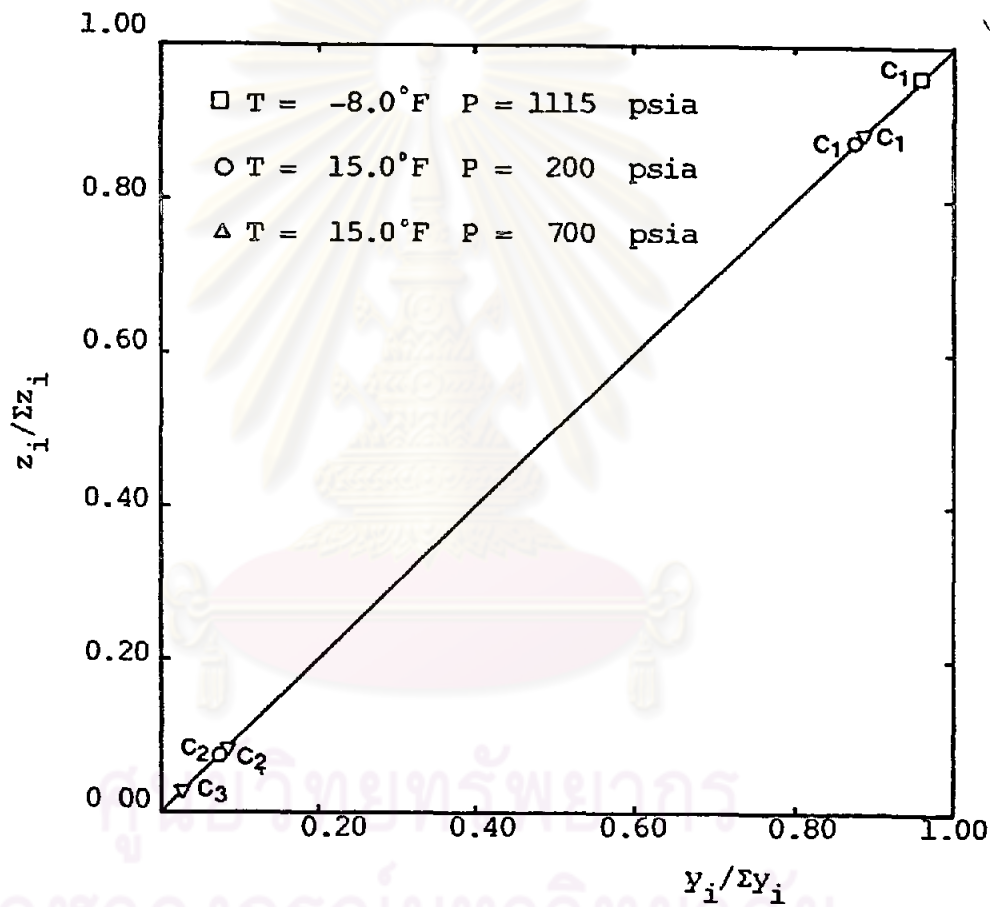


FIGURE 5.6 THE COMPARISON OF EXPERIMENTAL VALUES (ABCISSA) AND ESTIMATED VALUES (ORDINATE) IN THE VAPOR PHASE OF LIGHT-PSEUDOCOMPONENTS

TABLE 5.25 COMPARISON OF THE INITIAL PHASE VALUES OF THE EXPERIMENT AND THE ESTIMATION VALUES FOR IDEAL CASE

AT  $T = 15.0^{\circ}\text{F}$   $P = 200.0$  PSIA

Component	$x_i$	$y_i = z_i$	$P_i$ (atm)	$K_i = P_i/P$	$x_i' = y_i/K_i$		
CO <sub>2</sub>	0.0	0.0	26.53257	1.9495	0.0		
N <sub>2</sub>	0.019	0.93205	-	-	-		
C <sub>1</sub>	7.577	85.31217	506.4543	39.212	2.2926		
C <sub>2</sub>	6.028	8.38443	18.54087	1.3623	6.1546		
C <sub>3</sub>	8.828	2.87391	3.47006	0.255	11.2718		
						$x_i/\Sigma x_i$	$x_i'/\Sigma x_i'$
i-C <sub>4</sub>	7.698	0.86811	1.09927	0.0808	10.748	0.099	0.007
n-C <sub>4</sub>	10.202	0.72819	2.51477	0.185	3.941	0.132	0.002
i-C <sub>5</sub>	11.323	0.36700	0.22640	0.0166	22.062	0.146	0.014
n-C <sub>5</sub>	8.028	0.15691	0.15279	0.0123	13.977	0.104	0.009
C <sub>6</sub>	18.647	0.22277	0.03494	0.0026	86.774	0.240	0.054
C <sub>7</sub>	12.284	0.09508	0.00808	$5.94 \times 10^{-6}$	160.153	0.163	0.099
C <sub>8</sub>	6.917	0.04472	0.00191	$1.41 \times 10^{-4}$	317.661		
C <sub>9</sub>	2.018	0.01246	0.00042	$3.14 \times 10^{-5}$	396.216		
C <sub>10</sub>	0.382	0.00225	0.00009	$7.20 \times 10^{-6}$	312.474		
C <sub>11</sub>	0.042	0.00025	0.00002	$1.62 \times 10^{-6}$	154.659		
C <sub>12</sub>	0.006	0.00004	0.00000	0.0	0.0		

$T = 15.0^{\circ}\text{F}$   $P = 200.0$  psia  
 = 13.61 atm

75.547

1614.765

TABLE 5.26 COMPARISON OF THE INITIAL PHASE VALUES OF THE EXPERIMENT AND THE ESTIMATION VALUES FOR IDEAL CASE

AT T = 15.0 °F P = 700.0 PSIA

Component	$x_i$	$y_i = z_i$	$P_i$ (atm)	$K_i = P_i/P$	$x_i' = y_i/K_i$		
CO <sub>2</sub>	0.000	0.000	26.53257	0.55717	0.000	T = 15.0 °F P = 700.0 psia = 47.62 atm	
N <sub>2</sub>	0.078	0.98729	-	-	-		
C <sub>1</sub>	22.741	85.07018	506.4543	10.6353	7.999		
C <sub>2</sub>	14.343	8.44052	18.54087	0.38935	21.678		
C <sub>3</sub>	13.859	2.82973	3.47006	0.07287	38.832		
						$x_i/\Sigma x_i$	$x_i'/\Sigma x_i'$
i-C <sub>4</sub>	9.618	0.89296	1.09927	0.02308	38.690	0.196	0.007
n-C <sub>4</sub>	11.001	0.78460	2.51477	0.05281	14.857	0.225	0.0012
i-C <sub>5</sub>	8.795	0.38734	0.22640	0.00475	81.545	0.180	3.86x10 <sup>-6</sup>
n-C <sub>5</sub>	5.131	0.18930	0.15279	0.00321	58.972	0.005	0.004
C <sub>6</sub>	7.830	0.24508	0.03494	0.00073	335.726	0.160	0.027
C <sub>7</sub>	3.903	0.10477	0.00808	0.000017	6162.941	0.08	0.501
C <sub>8</sub>	2.001	0.05090	0.00191	0.00004	1272.500	0.041	0.103
C <sub>9</sub>	0.552	0.01373	0.00042	0.000009	1525.560	0.011	1.11x10 <sup>-6</sup>
C <sub>10</sub>	0.123	0.00303	0.00009	0.000002	1441.860	0.002	
C <sub>11</sub>	0.022	0.00055	0.00002	0.000000	1375.000		
C <sub>12</sub>	0.001	0.00003	0.00000	0.000000	0.000		

TABLE 5.27 COMPARISON OF THE INITIAL PHASE VALUES OF THE EXPERIMENT AND THE ESTIMATION VALUES FOR IDEAL CASE  
 AT  $T = -8.0^{\circ}\text{F}$   $P = 1115.0$  PSIA

Component	$x_i$	$y_i = z_i$	$P_i$ (atm)	$K_i = P_i/P$	$x'_i = y_i/K_i$	$x_i/\Sigma x_i$	$x'_i/\Sigma x'_i$
$\text{CO}_2$	0.589	4.21072	18.12699	0.239	17.618		
$\text{N}_2$	0.472	0.70973	-	-	-		
$\text{C}_1$	37.190	90.99455	301.6218	3.976	22.886		
$\text{C}_2$	4.604	3.04710	13.15507	0.173	17.613		
$\text{C}_3$	3.491	0.58334	2.22451	0.029	20.115		
i-C <sub>4</sub>	2.137	0.13558	0.65482	0.0086	15.765	0.04	$9.96 \times 10^{-4}$
n-C <sub>4</sub>	2.995	0.12874	1.78649	0.0236	5.455	0.056	$3.44 \times 10^{-4}$
i-C <sub>5</sub>	3.240	0.05995	0.12042	0.0016	59.95	0.060	$3.79 \times 10^{-3}$
n-C <sub>5</sub>	2.597	0.03708	0.07786	0.001	37.08	0.048	$2.34 \times 10^{-3}$
$\text{C}_6$	7.432	0.04952	0.1581	$2.08 \times 10^{-4}$	238.077	0.138	0.015
$\text{C}_7$	9.805	0.02741	0.00323	$4.26 \times 10^{-5}$	643.427	0.183	0.041
$\text{C}_8$	10.591	0.1012	0.00067	$8.83 \times 10^{-6}$	1146.093	0.197	0.072
$\text{C}_9$	7.952	0.00458	0.00013	$1.71 \times 10^{-6}$	2672.112	0.148	0.169
$\text{C}_{10}$	4.317	0.00154	0.000026	$3.43 \times 10^{-7}$	4489.79	0.08	0.284
$\text{C}_{11}$	1.696	0.00043	0.000005	$6.59 \times 10^{-8}$	6525.04	0.03	0.412
$\text{C}_{12}$	0.894	0.00017	0.00000	0.000	0.0		

$T = -8.0^{\circ}\text{F}$   $P = 1115.0$  psia  
 $= 75.85$  atm

53.656

15832.789



THE AMOUNT OF COMPONENT, N = 16

T = -10.00 °F      P = 500.00 PSIA      R = 10.730 PSIA-CUBIC FT/LB-MOLE°R

FEED RATE = 2.638330 MOLE/HR      V/F = 0.500000

INITIAL VALUE OF LIQUID AND VAPOR PHASE COMPOSITION

ID	COMPONENT	X	Y
1	CO2	0.062500	0.005544
2	N2	0.062500	0.002730
3	C1	0.062500	0.966292
4	C2	0.062500	0.022314
5	C3	0.062500	0.002002
6	IC4	0.062500	0.000317
7	NC4	0.062500	0.000294
8	IC5	0.062500	0.000138
9	NC5	0.062500	0.000101
10	C6	0.062500	0.000083
11	C7	0.062500	0.000079
12	C8	0.062500	0.000048
13	C9	0.062500	0.000034
14	C10	0.062500	0.000019
15	C11	0.062500	0.000005
16	C12	0.062500	0.000001
SUM =		1.000000	0.999799

V/F = 0.999755800

V = 2.638345000 MOLE/HR

L = 0.000484467 MOLE/HR

NO. OF ITERATION = 3

THE VAPOR PHASE COMPOSITIONS FROM THE  
 EXPERIMENTAL DATA AND THE ORDINARY  
 VLE CALCULATION MODEL AT T = -10.0°F  
 P = 500.0 PSIA

TABLE 5.28 COMPARISON OF THE LIQUID PHASE AND

ID	COMPONENT	X(EXP)	Y(EXP)	X(MODEL)	Y(MODEL)	K(MODEL)	ERRX	ERRY
1	CO2	0.00228	0.00554	0.00413	0.005544	1.46123	0.00191	-0.000000
2	N2	0.00007	0.00273	0.00019	0.002730	16.07114	0.00012	0.000000
3	C1	0.21054	0.96639	0.27078	0.966448	3.94478	0.06024	0.000061
4	C2	0.03044	0.02231	0.04809	0.022303	0.51266	0.01765	-0.000005
5	C3	0.01310	0.00200	0.01873	0.001999	0.11795	0.00563	-0.000002
6	IC4	0.00612	0.00032	0.00827	0.000315	0.04207	0.00215	-0.000001
7	NC4	0.00797	0.00029	0.01195	0.000291	0.02694	0.00398	-0.000002
9	IC5	0.01355	0.00014	0.01575	0.000134	0.00942	0.00220	-0.000002
9	NC5	0.01318	0.00010	0.01687	0.000097	0.00635	0.00369	-0.000002
10	C6	0.04268	0.00008	0.05149	0.000072	0.00155	0.00881	-0.000006
11	C7	0.09785	0.00007	0.14093	0.000048	0.00038	0.04313	-0.000019
12	C8	0.17314	0.00003	0.15900	0.000012	0.00009	-0.01414	-0.000014
13	C9	0.20957	0.00001	0.14164	0.000003	0.00002	-0.06793	-0.000005
14	C10	0.12938	0.00003	0.08348	0.000000	0.00000	-0.04590	-0.000026
15	C11	0.03810	0.00000	0.02275	0.000000	0.00000	-0.01535	-0.000000
16	C12	0.01032	0.00000	0.00583	0.000000	0.00000	-0.00449	0.000000
SUM =		1.00000	1.00000	1.00000	1.00000	ΣAAD =	1.85838	0.000704

THE AMOUNT OF COMPONENT, N = 16

T = 0.00 °F      P = 500.00 PSIA      R = 10.730 PSIA-CUBIC FT/LB-MOLE<sup>0</sup>R

FEED RATE = 2.633710 MOLE/HR      V/F = 0.500000

INITIAL VALUE OF LIQUID AND VAPOR PHASE COMPOSITION

ID	COMPONENT	X	Y
1	C32	0.062500	0.005940
2	N2	0.062500	0.002810
3	C1	0.062500	0.966192
4	C2	0.062500	0.021952
5	C3	0.062500	0.002042
6	IC4	0.062500	0.000303
7	NC4	0.062500	0.000294
8	IC5	0.062500	0.000136
9	NC5	0.062500	0.000097
10	C6	0.062500	0.000108
11	C7	0.062500	0.000113
12	C8	0.062500	0.000060
13	C9	0.062500	0.000029
14	C10	0.062500	0.000018
15	C11	0.062500	0.000005
16	C12	0.062500	0.000002
SUM =		1.000000	1.000000

V/F = 0.339755800

V = 2.633269000 MOLE/HR

L = 0.000440538 MOLE/HR

NO. OF ITERATION = 3

TABLE 5.29 COMPARISON OF THE LIQUID PHASE AND  
 THE VAPOR PHASE COMPOSITIONS FROM  
 THE EXPERIMENTAL DATA AND THE  
 ORDINARY VLE CALCULATION MODEL AT  
 T = 0.0° F    P = 500.0 PSIA

ID	COMPONENT	X(EXP)	Y(EXP)	X(MODEL)	Y(MODEL)	K(MODEL)	ERRX	ERRY
1	C32	0.00321	0.00594	0.00413	0.005940	1.59715	0.00097	0.000000
2	N2	0.00013	0.00281	0.00020	0.002810	16.14519	0.00007	0.000000
3	C1	0.19913	0.96625	0.26293	0.966347	4.13322	0.06380	0.000094
4	C2	0.03118	0.02185	0.04292	0.021848	0.57249	0.01174	-0.000004
5	C3	0.01306	0.00204	0.01667	0.002039	0.13751	0.00361	-0.000002
6	IC4	0.00630	0.00030	0.00674	0.000302	0.05040	0.00044	-0.000001
7	NC4	0.00980	0.00029	0.01000	0.000292	0.03279	0.00020	-0.000001
8	IC5	0.01213	0.00014	0.01273	0.000133	0.01177	0.00060	-0.000002
9	NC5	0.01204	0.00010	0.01312	0.000094	0.00807	0.00108	-0.000002
10	C6	0.03781	0.00011	0.05304	0.000097	0.00205	0.01523	-0.000008
11	C7	0.08907	0.00011	0.16517	0.000077	0.00052	0.00610	-0.000029
12	C8	0.16319	0.00005	0.18167	0.000020	0.00013	0.01848	-0.000026
13	C9	0.20815	0.00001	0.11964	0.000003	0.00003	-0.08851	-0.000010
14	C10	0.14774	0.00001	0.07936	0.000001	0.00001	-0.06838	-0.000005
15	C11	0.05335	0.00000	0.02387	0.000000	0.00000	0.02948	-0.000001
16	C12	0.02003	0.00000	0.00776	0.000000	0.00000	-0.01227	-0.000000
SUM =		1.00000	1.00000	1.00000	1.000001	ZAAD =	2.44332	0.001165

THE AMOUNT OF COMPONENT, N = 16

T = 50.00 °F      P = 500.00 PSIA      R = 10.730 PSIA-CUBIC FT/LB-MOLE °R

FEED RATE = 1.320530 MOLE/HR      V/F = 0.500000

INITIAL VALUE OF LIQUID AND VAPOR PHASE COMPOSITION

ID	COMPONENT	X	Y
1	C2	0.062500	0.002595
2	V2	0.062500	0.008323
3	C1	0.062500	0.851252
4	C2	0.062500	0.076620
5	C3	0.062500	0.003315
6	IC4	0.062500	0.012398
7	NC4	0.062500	0.008630
8	IC5	0.062500	0.003476
9	NC5	0.062500	0.001519
10	C6	0.062500	0.001224
11	C7	0.062500	0.000387
12	C8	0.062500	0.000137
13	C9	0.062500	0.000034
14	C10	0.062500	0.000011
15	C11	0.062500	0.000001
16	C12	0.062500	0.000000
SUM =		1.000000	0.969920

V/F = 0.778699607  
V = 1.318741000 MOLE/HR  
L = 0.001788139 MOLE/HR

NO. OF ITERATION = 3

THE VAPOR PHASE COMPOSITIONS FROM THE  
 EXPERIMENTAL DATA AND THE ORDINARY  
 VLE CALCULATION MODEL AT T = 50.0 °F  
 P = 500.0 PSIA

Table 5.30 COMPARISON OF THE LIQUID PHASE AND

ID	COMPONENT	X(EXP)	Y(EXP)	X(MODEL)	Y(MODEL)	K(MODEL)	ERRX	ERRY
1	C2	0.00103	0.00260	0.00115	0.002677	2.31748	0.00012	0.000081
2	V2	0.00039	0.00833	0.00054	0.008591	15.79344	0.00015	0.000261
3	C1	0.14339	0.85191	0.17623	0.878570	4.95675	0.03284	0.026658
4	C2	0.07636	0.07662	0.08424	0.078989	0.93225	0.00788	0.002369
5	C3	0.10988	0.00331	0.01225	0.003406	0.27646	-0.09763	0.000098
6	IC4	0.10324	0.01231	0.10823	0.012657	0.11628	0.00499	0.000344
7	NC4	0.11279	0.00853	0.10697	0.008769	0.08151	-0.00582	0.000236
8	IC5	0.10690	0.00338	0.10281	0.003454	0.03341	-0.00409	0.000074
9	NC5	0.06257	0.00146	0.05985	0.001490	0.02475	-0.00272	0.000028
10	C6	0.11327	0.00112	0.13943	0.001081	0.00771	0.02616	0.000038
11	C7	0.07512	0.00003	0.10685	0.000259	0.00241	0.03173	0.000228
12	C8	0.05875	0.00008	0.06958	0.000050	0.00072	0.01083	0.000032
13	C9	0.02627	0.00001	0.02287	0.000005	0.00022	-0.00340	0.000004
14	C10	0.00901	0.00000	0.00817	0.000001	0.00007	-0.00084	0.000002
15	C11	0.00103	0.00000	0.00082	0.000000	0.00003	-0.00021	0.000000
16	C12	0.00000	0.00000	0.00000	0.000000	0.00001	0.00000	0.000000
SUM =		1.00000	1.00000	1.00000	1.00000	FAAD =	1.43374	0.190338



THE AMOUNT OF COMPONENT, N = 16

T = 47.00 °F      P = 815.00 PSIA      R = 10.730 PSIA-CUBIC FT/LB-MOLE °R

FEED RATE = 2.638670 MOLE/HR      V/F = 0.500000

INITIAL VALUE OF LIQUID AND VAPOR PHASE COMPOSITION

ID	COMPONENT	X	Y
1	CO2	0.062500	0.006617
2	N2	0.062500	0.029980
3	C1	0.062500	0.923041
4	C2	0.062500	0.030932
5	C3	0.062500	0.005329
6	IC4	0.062500	0.001090
7	NC4	0.062500	0.001041
8	IC5	0.062500	0.000499
9	NC5	0.062500	0.000368
10	C6	0.062500	0.000423
11	C7	0.062500	0.000315
12	C8	0.062500	0.000198
13	C9	0.062500	0.000098
14	C10	0.062500	0.000045
15	C11	0.062500	0.000017
16	C12	0.062500	0.000007
SUM =		1.000000	0.999999

V/F = 0.799623200  
V = 2.637651000 MOLE/HR  
L = 0.001018524 MOLE/HR

NO. OF ITERATION = 3

THE VAPOR PHASE COMPOSITIONS FROM THE  
 EXPERIMENTAL DATA AND THE ORDINARY  
 VLE CALCULATION MODEL AT T = 47.0°F  
 P = 815.0 PSIA

TABLE 5.31 COMPARISON OF THE LIQUID PHASE AND

ID	COMPONENT	X(EXP)	Y(EXP)	X(MODEL)	Y(MODEL)	K(MODEL)	ERRX	EPRY
1	CO2	0.00405	0.00662	0.00398	0.006618	1.64686	0.00007	0.000001
2	N2	0.00253	0.02998	0.00271	0.029991	10.96481	0.00018	0.000008
3	C1	0.21503	0.92309	0.26213	0.923290	3.48636	0.04710	0.000204
4	C2	0.04723	0.03093	0.04362	0.030927	0.70183	-0.00361	-0.000004
5	C3	0.02556	0.00533	0.02353	0.005322	0.22385	-0.00203	-0.000006
6	IC4	0.01191	0.00109	0.01063	0.001087	0.10122	-0.00128	-0.000003
7	NC4	0.01752	0.00104	0.01441	0.001036	0.07112	-0.00311	-0.000004
8	IC5	0.01744	0.00050	0.01551	0.000493	0.03148	-0.00193	-0.000005
9	NC5	0.01503	0.00037	0.01532	0.000362	0.02339	0.00029	-0.000005
10	C6	0.04752	0.00042	0.05049	0.000404	0.00792	0.00297	-0.000016
11	C7	0.08236	0.00031	0.10140	0.000277	0.00270	0.01904	-0.000033
12	C8	0.12308	0.00019	0.15632	0.000139	0.00088	0.03324	-0.000051
13	C9	0.13932	0.00009	0.14598	0.000043	0.00029	0.00666	-0.000046
14	C10	0.11951	0.00004	0.09513	0.000009	0.00010	-0.02438	-0.000028
15	C11	0.07489	0.00001	0.04021	0.000002	0.00004	-0.03464	-0.000010
16	C12	0.05706	0.00000	0.01864	0.000000	0.00002	-0.03842	-0.000003
SUM =		1.00000	1.000000			FAAD =	1.36850	0.002561

THE AMOUNT OF COMPONENT, N = 16

T = -8.00 °F      ρ = 250.00 PSIA      R = 10.730 PSIA-CUBIC FT/LB-MOLE<sup>o</sup>R

FEED RATE = 2.637800 MOLE/HR      V/F = 0.500000

INITIAL VALUE OF LIQUID AND VAPOR PHASE COMPOSITION

ID	COMPONENT	X	Y
1	C12	0.062500	0.004000
2	N2	0.062500	0.035993
3	C1	0.062500	0.894976
4	C2	0.062500	0.044821
5	C3	0.062500	0.014038
6	IC4	0.062500	0.001311
7	NC4	0.062500	0.002835
8	IC5	0.062500	0.000622
9	NC5	0.062500	0.000632
10	C6	0.062500	0.000294
11	C7	0.062500	0.000264
12	C8	0.062500	0.000144
13	C9	0.062500	0.000045
14	C10	0.062500	0.000018
15	C11	0.062500	0.000006
16	C12	0.062500	0.000002
SUM =		1.000000	1.000000

V/F = 0.799107600

V = 2.637214000 MOLE/HR

L = 0.002585411 MOLE/HR

NO. OF ITERATION = 3

THE VAPOR PHASE COMPOSITIONS FROM THE  
 EXPERIMENTAL DATA AND THE ORDINARY  
 VLE CALCULATION MODEL AT T = -8.0 °F  
 P = 250.0 PSIA

TABLE 5.32 COMPARISON OF THE LIQUID PHASE AND

ID	COMPONENT	X(EXP)	Y(EXP)	X(MODEL)	Y(MODEL)	K(MODEL)	ERRX	ERRY
1	C12	0.00175	0.00400	0.00145	0.004002	2.65062	-0.00030	0.000001
2	N2	0.00110	0.03601	0.00100	0.036026	34.63550	-0.00010	0.000015
3	C1	0.11403	0.89536	0.11202	0.895705	7.67670	-0.00201	0.000345
4	C2	0.06160	0.04481	0.05057	0.044816	0.85083	-0.01103	0.000023
5	C3	0.00821	0.01400	0.007730	0.013979	0.17361	0.06909	-0.000025
6	IC4	0.02246	0.00130	0.02189	0.001292	0.05664	-0.00037	-0.000009
7	NC4	0.07207	0.00280	0.07561	0.002767	0.03514	0.00354	-0.000034
8	IC5	0.04464	0.00060	0.04938	0.000576	0.01121	0.00474	-0.000024
9	NC5	0.06466	0.00060	0.07365	0.000564	0.00735	0.00899	-0.000037
10	C6	0.09214	0.00025	0.11381	0.000189	0.00159	0.02167	-0.000061
11	C7	0.15476	0.00019	0.20462	0.000074	0.00035	0.04986	-0.000114
12	C8	0.16507	0.00006	0.14317	0.000010	0.00007	-0.02190	-0.000052
13	C9	0.07507	0.00001	0.04816	0.000001	0.00001	-0.02691	-0.000008
14	C10	0.03299	0.00000	0.01887	0.000000	0.00000	-0.01412	-0.000001
15	C11	0.01160	0.00000	0.00634	0.000000	0.00000	-0.00526	-0.000000
16	C12	0.00393	0.00000	0.00215	0.000000	0.00000	-0.00178	0.000000
SUM =		1.00000	1.000000			ΣAAD =	1.51173	0.004559

THE AMOUNT OF COMPONENT, N = 16

T = -8.00 ° F      P = 1115.00    PSIA      R = 10.730    PSIA-CUBIC FT/LB-MOLE<sup>OR</sup>  
 FEED RATE = 2.678890    MOLE/HR      V/F = 0.500000

INITIAL VALUE OF LIQUID AND VAPOR PHASE COMPOSITION

ID	COMPONENT	X	Y
1	CO2	0.062500	0.007097
2	N2	0.062500	0.042107
3	C1	0.062500	0.909945
4	C2	0.062500	0.030471
5	C3	0.062500	0.005833
6	IC4	0.062500	0.001356
7	NC4	0.062500	0.001287
8	IC5	0.062500	0.000599
9	NC5	0.062500	0.000371
10	C6	0.062500	0.000495
11	C7	0.062500	0.000274
12	C8	0.062500	0.000101
13	C9	0.062500	0.000046
14	C10	0.062500	0.000015
15	C11	0.062500	0.000004
16	C12	0.062500	0.000002
SUM =		1.000000	1.000004

V/F = 0.999432500  
 V = 2.637163000    MOLE/HR  
 L = 0.001727104    MOLE/HR

NO. OF ITERATION = 3

P ≈ 1115.0 PSIA  
 VLE CALCULATION MODEL AT T = -8.0 ° F  
 THE VAPOR PHASE COMPOSITIONS FROM THE  
 EXPERIMENTAL DATA AND THE ORDINARY

TABLE 5.33 COMPARISON OF THE LIQUID PHASE AND

ID	COMPONENT	X (EXP)	Y (EXP)	X (MODEL)	Y (MODEL)	K (MODEL)	ERRX	ERRY
1	CO2	0.00589	0.00710	0.00722	0.007097	0.96081	0.00133	-0.000000
2	N2	0.00472	0.04211	0.00601	0.042128	6.85053	0.00129	0.000021
3	C1	0.37190	0.91002	0.41306	0.910229	2.15366	0.04116	0.000206
4	C2	0.04604	0.03047	0.06890	0.030448	0.43188	0.02286	-0.000020
5	C3	0.03491	0.00583	0.04043	0.005813	0.14053	0.00552	-0.000016
6	IC4	0.02137	0.00135	0.01994	0.001345	0.06594	-0.00143	-0.000008
7	NC4	0.02995	0.00128	0.02727	0.001272	0.04561	-0.00268	-0.000011
8	IC5	0.03240	0.00059	0.02698	0.000584	0.02116	-0.00542	-0.000011
9	NC5	0.02597	0.00037	0.02282	0.000358	0.01532	-0.00315	-0.000009
10	C6	0.07432	0.00048	0.08227	0.000448	0.00532	0.00795	-0.000037
11	C7	0.09805	0.00026	0.11038	0.000210	0.00186	0.01233	-0.000050
12	C8	0.10591	0.00006	0.08338	0.000053	0.00062	-0.02253	-0.000006
13	C9	0.07952	0.00003	0.05801	0.000012	0.00020	-0.02151	-0.000022
14	C10	0.04317	0.00001	0.02352	0.000002	0.00007	-0.01965	-0.000007
15	C11	0.01696	0.00000	0.00697	0.000000	0.00004	-0.00999	-0.000002
16	C12	0.00894	0.00000	0.00286	0.000000	0.00001	-0.00608	-0.000000
SUM =		1.00000	1.000000			ΣAAD =	1.15552	0.002657

THE AMOUNT OF COMPONENT, N = 16

T = -10.00 ° F      P = 800.00 PSIA      R = 10.730 PSIA-CUBIC FT/LB-MOLE<sup>0</sup>R

FEED RATE = 2.543530 MOLE/HR      V/F = 0.500000

INITIAL VALUE OF LIQUID AND VAPOR PHASE COMPOSITION

ID	COMPONENT	X	Y
1	CO2	0.062500	0.003736
2	N2	0.062500	0.002052
3	C1	0.062500	0.972654
4	C2	0.062500	0.020926
5	C3	0.062500	0.000280
6	IC4	0.062500	0.000052
7	NC4	0.062500	0.000068
8	IC5	0.062500	0.000067
9	NC5	0.062500	0.000039
10	C6	0.062500	0.000040
11	C7	0.062500	0.000031
12	C8	0.062500	0.000018
13	C9	0.062500	0.000014
14	C10	0.062500	0.000015
15	C11	0.062500	0.000005
16	C12	0.062500	0.000001
SUM =		1.000000	0.999999

V/F = 0.999992100  
V = 2.543536000 MOLE/HR  
L = 0.000053406 MOLE/HR

NO. OF ITERATION = 5

P = 800.0 PSIA  
 VLE CALCULATION MODEL AT T = 10.0°F  
 THE VAPOR PHASE COMPOSITIONS FROM THE  
 EXPERIMENTAL DATA AND THE ORDINARY  
 VLE CALCULATION MODEL AT T = 10.0°F

TABLE 5.34 COMPARISON OF THE LIQUID PHASE AND

ID	COMPONENT	X(EXP)	Y(EXP)	X(MODEL)	Y(MODEL)	K(MODEL)	ERRX	ERRY
1	CO2	0.00143	0.00374	0.00167	0.003737	1.17160	0.00024	-0.000000
2	N2	0.00006	0.00205	0.00003	0.002052	12.78584	0.00002	-0.000000
3	C1	0.31340	0.97268	0.16267	0.972667	3.12449	-0.15073	-0.000011
4	C2	0.02363	0.02093	0.02420	0.020926	0.45188	0.00057	0.000000
5	C3	0.00169	0.00028	0.00124	0.000280	0.11813	-0.00045	0.000000
6	IC4	0.00056	0.00005	0.00056	0.000052	0.04819	0.00000	-0.000000
7	NC4	0.00177	0.00007	0.00116	0.000068	0.03078	-0.00061	0.000000
8	IC5	0.00605	0.00007	0.00233	0.000067	0.01241	-0.00322	0.000000
9	NC5	0.00498	0.00004	0.00244	0.000039	0.00838	-0.00254	0.000000
10	C6	0.01702	0.00004	0.00870	0.000040	0.00236	-0.00832	0.000000
11	C7	0.03929	0.00003	0.02390	0.000031	0.00068	-0.01539	0.000001
12	C8	0.05857	0.00002	0.04912	0.000017	0.00018	-0.00945	0.000001
13	C9	0.14467	0.00001	0.12762	0.000012	0.00005	-0.01705	0.000003
14	C10	0.25830	0.00001	0.36529	0.000010	0.00001	0.10699	0.000004
15	C11	0.10695	0.00000	0.18574	0.000002	0.00001	0.07879	0.000001
16	C12	0.01902	0.00000	0.04277	0.000000	0.00000	0.02375	0.000000
SUM =		1.00000	1.000000	1.000000	1.000000	%AAD =	2.61326	0.000148

THE AMOUNT OF COMPONENT, N = 16

T = 8.00 °F P = 300.00 PSIA R = 10.730 PSIA-CUBIC FT/LB-MOLE °R

FEED RATE = 1.153050 MOLE/HR V/F = 0.500000

INITIAL VALUE OF LIQUID AND VAPOR PHASE COMPOSITION

ID	COMPONENT	X	Y
1	CO2	0.062500	0.010069
2	N2	0.062500	0.004437
3	C1	0.062500	0.951993
4	C2	0.062500	0.031772
5	C3	0.062500	0.001574
6	IC4	0.062500	0.000015
7	NC4	0.062500	0.000014
8	IC5	0.062500	0.000018
9	NC5	0.062500	0.000017
10	C6	0.062500	0.000053
11	C7	0.062500	0.000056
12	C8	0.062500	0.000034
13	C9	0.062500	0.000013
14	C10	0.062500	0.000027
15	C11	0.062500	0.000006
16	C12	0.062500	0.000000
SUM =		1.000000	1.000000

V/F = 0.500000  
V = 1.152992030 MOLE/HR  
L = 0.000058174 MOLE/HR

NO. OF ITERATION = 3

THE VAPOR PHASE COMPOSITIONS FROM  
 THE EXPERIMENTAL DATA AND THE  
 ORDINARY VLE CALCULATION MODEL AT  
 T = 8.0°F P = 300.0 PSIA

TABLE 5.35 COMPARISON OF THE LIQUID PHASE AND

ID	COMPONENT	X(EXP)	Y(EXP)	X(MODEL)	Y(MODEL)	K(MODEL)	ERRX	ERRY
1	CO2	0.00262	0.01007	0.00513	0.010069	2.55399	0.00251	0.000000
2	N2	0.00007	0.00444	0.00022	0.004437	26.08089	0.00015	0.000000
3	C1	0.07704	0.95193	0.18779	0.951963	6.59248	0.11075	0.000040
4	C2	0.03243	0.03177	0.04734	0.031771	0.87280	0.01491	-0.000001
5	C3	0.01646	0.00157	0.01026	0.001573	0.19946	-0.00620	-0.000000
6	IC4	0.00008	0.00002	0.00028	0.000015	0.06951	0.00020	-0.000000
7	NC4	0.00021	0.00001	0.00040	0.000014	0.04521	0.00019	-0.000000
8	IC5	0.00130	0.00002	0.00154	0.000018	0.01536	0.00024	-0.000000
9	NC5	0.00173	0.00002	0.00210	0.000017	0.01054	0.00037	-0.000000
10	C6	0.02294	0.00005	0.02617	0.000051	0.00253	0.00323	-0.000002
11	C7	0.06476	0.00005	0.09964	0.000047	0.00061	0.03488	-0.000007
12	C8	0.09969	0.00003	0.16972	0.000018	0.00014	0.07003	-0.000012
13	C9	0.11754	0.00001	0.16686	0.000003	0.00003	-0.01068	-0.000005
14	C10	0.42584	0.00001	0.27146	0.000002	0.00001	-0.15438	-0.000009
15	C11	0.12610	0.00000	0.06580	0.000000	0.00000	0.06030	0.000000
16	C12	0.01119	0.00000	0.00530	0.000000	0.00000	-0.00589	0.000000
SUM =		1.00000	1.000002			%AAD =	2.96814	0.000484

THE AMOUNT OF COMPONENT, N = 16

T = 8.00 ° F P = 700.00 PSIA R = 10.730 PSIA-CUBIC FT/LB-MOLE ° R

FEED RATE = 2.453830 MOLE/HR V/F = 0.500000

INITIAL VALUE OF LIQUID AND VAPOR PHASE COMPOSITION

ID	COMPONENT	X	Y
1	CO2	0.062500	0.010155
2	N2	0.062500	0.004651
3	C1	0.062500	0.951288
4	C2	0.062500	0.031921
5	C3	0.062500	0.001755
6	IC4	0.062500	0.000015
7	NC4	0.062500	0.000014
8	IC5	0.062500	0.000013
9	NC5	0.062500	0.000015
10	C6	0.062500	0.000048
11	C7	0.062500	0.000051
12	C8	0.062500	0.000031
13	C9	0.062500	0.000013
14	C10	0.062500	0.000024
15	C11	0.062500	0.000004
16	C12	0.062500	0.000001
SUM =		1.000000	0.999999

V/F = 0.999877900  
 V = 2.453800000 MOLE/HR  
 L = 0.090030109 MOLE/HR

NO. OF ITERATION = 3

ORDINARY VLE CALCULATION MODEL AT  
 T = 8.0 ° F P = 700.0 PSIA  
 THE VAPOR PHASE COMPOSITIONS FROM  
 THE EXPERIMENTAL DATA AND THE

TABLE 5.36 COMPARISON OF THE LIQUID PHASE AND

ID	COMPONENT	X(EXP)	Y(EXP)	X(MODEL)	Y(MODEL)	K(MODEL)	ERRX	ERRY
1	CO2	0.00285	0.01016	0.00947	0.010155	1.31993	0.00662	-0.000000
2	N2	0.00009	0.00465	0.00055	0.004652	10.42605	0.00046	0.000000
3	C1	0.25105	0.95131	0.38169	0.951346	3.06853	0.13064	0.000036
4	C2	0.02279	0.03192	0.07540	0.031917	0.52112	0.05261	-0.000005
5	C3	0.00885	0.00176	0.01493	0.001754	0.14461	0.00608	-0.000001
6	IC4	0.00005	0.00001	0.00031	0.000015	0.05851	0.00026	-0.000000
7	NC4	0.00015	0.00001	0.00043	0.000014	0.03987	0.00028	-0.000000
8	IC5	0.00092	0.00001	0.00100	0.000013	0.01580	0.00009	-0.000000
9	NC5	0.00128	0.00001	0.00162	0.000015	0.01131	0.00034	-0.000000
10	C6	0.01516	0.00005	0.01713	0.000046	0.00330	0.00197	-0.000001
11	C7	0.04357	0.00005	0.05805	0.000046	0.00097	0.01449	-0.000004
12	C8	0.06792	0.00003	0.09724	0.000021	0.00027	0.02932	-0.000008
13	C9	0.10162	0.00001	0.08130	0.000005	0.00008	-0.02032	-0.000005
14	C10	0.36321	0.00001	0.20955	0.000004	0.00002	-0.15366	-0.000009
15	C11	0.09597	0.00000	0.04242	0.000000	0.00001	-0.05355	-0.000001
16	C12	0.02452	0.00000	0.00889	0.000000	0.00000	-0.01563	-0.000000
SUM =		1.00000	1.00000	1.00000	1.00000	SAAD =	3.03951	0.000445

THE AMOUNT OF COMPONENT, N = 16

T = 0.00 ° F      P = 300.00 PSIA      R = 10.730 PSIA-CUBIC FT/LB-MOLE ° R

FEED RATE = 2.250860 MOLE/HR      V/F = 0.500000

INITIAL VALUE OF LIQUID AND VAPOR PHASE COMPOSITION

ID	COMPONENT	X	Y
1	CO2	0.062500	0.006002
2	N2	0.062500	0.003101
3	C1	0.062500	0.964610
4	C2	0.062500	0.022708
5	C3	0.062500	0.002301
6	IC4	0.062500	0.000330
7	NC4	0.062500	0.000311
8	IC5	0.062500	0.000151
9	NC5	0.062500	0.000111
10	C6	0.062500	0.000123
11	C7	0.062500	0.000112
12	C8	0.062500	0.000075
13	C9	0.062500	0.000040
14	C10	0.062500	0.000019
15	C11	0.062500	0.000004
16	C12	0.062500	0.000001
SUM =		1.000000	0.979999

V/F = 0.779755800  
 V = 2.250412000 MOLE/HR  
 L = 0.009447273 MOLE/HR

NO. OF ITERATION = 3

THE VAPOR PHASE COMPOSITIONS FROM THE  
 EXPERIMENTAL DATA AND THE ORDINARY  
 VLE CALCULATION MODEL AT T = 0.0 ° F  
 P = 300.0 PSIA

ID	COMPONENT	X(EXP)	Y(EXP)	X(MODEL)	Y(MODEL)	K(MODEL)	FRRX	ERRY
1	CO2	0.00172	0.00600	0.00270	0.006003	2.43330	0.00098	0.000000
2	N2	0.00006	0.00310	0.00012	0.003102	28.44411	0.00006	0.000000
3	C1	0.12591	0.96467	0.15784	0.964792	6.68532	0.03193	0.000120
4	C2	0.01849	0.02271	0.03065	0.022707	0.81029	0.01216	0.000002
5	C3	0.00894	0.00230	0.01429	0.002299	0.17595	0.00535	0.000002
6	IC4	0.00425	0.00033	0.00602	0.000329	0.05983	0.00177	0.000001
7	NC4	0.00706	0.00031	0.00891	0.000309	0.03790	0.00185	0.000001
8	IC5	0.01056	0.00015	0.01285	0.000148	0.01259	0.00229	0.000002
9	NC5	0.01112	0.00011	0.01397	0.000108	0.00843	0.00285	0.000002
10	C6	0.04042	0.00012	0.06171	0.000109	0.00194	0.02129	0.000011
11	C7	0.09897	0.00010	0.17672	0.000072	0.00045	0.07775	0.000032
12	C8	0.18355	0.00006	0.23975	0.000021	0.00010	0.05620	0.000040
13	C9	0.24508	0.00002	0.16423	0.000003	0.00002	0.08085	0.000010
14	C10	0.17514	0.00001	0.08440	0.000000	0.00000	-0.09074	0.000006
15	C11	0.05016	0.00000	0.01959	0.000000	0.00000	-0.03057	0.000001
16	C12	0.01859	0.00000	0.00625	0.000000	0.00000	-0.01233	0.000000
SUM =		1.00000	1.000000			ΣAAD =	2.68101	0.001507

THE AMOUNT OF COMPONENT, N = 16

T = -10.00 °F    P = 500.00 PSIA    R = 10.720    FSTA-CU3IC FT/LB-MOLE<sup>R</sup>

FEED RATE = 2.638830 MOLE/HR    V/F = 0.500000

INITIAL VALUE OF LIQUID AND VAPOR PHASE COMPOSITION

ID	COMPONENT	X	Y
1	CO2	0.062500	0.005544
2	N2	0.062500	0.002730
3	C1	0.062500	0.966292
4	C2	0.062500	0.022314
5	C3	0.062500	0.002002
6	IC4	0.062500	0.000317
7	NC4	0.062500	0.000274
8	IC5	0.062500	0.000138
9	NC5	0.062500	0.000101
10	C6	0.062500	0.000093
11	C7	0.062500	0.000079
12	C8	0.062500	0.000048
13	C9	0.062500	0.000034
14	C10	0.062500	0.000019
15	C11	0.062500	0.000005
16	C12	0.062500	0.000001
SUM =		1.000000	0.997779

V/F = 0.778513200  
V = 2.634835000 MOLE/HR  
L = 0.003994942 MOLE/HR

NO. OF ITERATION = 2

THE VAPOR PHASE COMPOSITIONS FROM  
THE EXPERIMENTAL DATA AND THE  
PROPOSED MODEL AT T = -10.0°F  
P = 500.0 PSIA

TABLE 5.38 COMPARISON OF THE LIQUID PHASE AND

ID	COMPONENT	X(EXP)	Y(EXP)	X(MODEL)	Y(MODEL)	K(MODEL)	ERRX	ERRY
1	CO2	0.00228	0.00554	0.00331	0.005541	1.51901	0.00137	0.000002
2	N2	0.00007	0.00273	0.00013	0.002731	13.55620	0.00008	0.000004
3	C1	0.21054	0.96639	0.20444	0.966471	4.28653	0.01514	0.001003
4	C2	0.03044	0.02231	0.03812	0.022263	0.52955	0.01164	-0.000023
5	C3	0.01310	0.00200	0.01493	0.001777	0.11976	0.00344	-0.000020
6	IC4	0.00612	0.00032	0.00590	0.000316	0.04849	0.00040	-0.000000
7	NC4	0.00797	0.00029	0.00847	0.000293	0.03132	0.00138	-0.000000
8	IC5	0.01355	0.00014	0.01080	0.000136	0.01143	-0.00163	0.000000
9	NC5	0.01318	0.00010	0.01155	0.000099	0.00777	-0.00043	0.000000
10	C6	0.04268	0.00008	0.03573	0.000079	0.00200	-0.00324	0.000000
11	C7	0.09785	0.00007	0.11303	0.000064	0.00051	0.02698	-0.000003
12	C9	0.17314	0.00003	0.17493	0.000024	0.00012	0.01979	-0.000002
13	C9	0.20957	0.00001	0.20102	0.000007	0.00003	0.01234	-0.000001
14	C10	0.12938	0.00003	0.12131	0.000001	0.00001	0.01557	-0.0000025
15	C11	0.03810	0.00000	0.03667	0.000000	0.00000	0.00240	-0.000000
16	C12	0.01032	0.00000	0.00951	0.000000	0.00000	0.00018	0.000000
SUM =		1.00000	1.00000	1.00000	1.00000	KAAD =	0.72633	0.006319



THE AMOUNT OF COMPONENT A = 16

T = 0.00 ° F      P = 500.00 PSIA      R = 10.730 PSIA-CL3IC FT/LB-MOLE<sup>OR</sup>

FEED RATE = 2.633710 MOLE/HR.      V/F = 0.50000

INITIAL VALUE OF LIQUID AND VAPOR PHASE COMPOSITION

ID	COMPONENT	X	Y
1	CO2	0.062500	0.005940
2	N2	0.062500	0.002810
3	C1	0.062500	0.966192
4	C2	0.062500	0.021852
5	C3	0.062500	0.002042
6	IC4	0.062500	0.000303
7	NC4	0.062500	0.000294
8	IC5	0.062500	0.000136
9	NC5	0.062500	0.000097
10	C6	0.062500	0.000108
11	C7	0.062500	0.000113
12	C8	0.062500	0.000060
13	C9	0.062500	0.000029
14	C10	0.062500	0.000018
15	C11	0.062500	0.000005
16	C12	0.062500	0.000002
SUM =		1.000000	1.000000

V/F = 0.998506300  
 V = 2.634722000 MOLE/HR  
 L = 0.003987312 MOLE/HR

NO. OF ITERATION = 2

TABLE 5.39 COMPARISON OF THE LIQUID PHASE AND  
 THE VAPOR PHASE COMPOSITIONS FROM THE  
 EXPERIMENTAL DATA AND THE PROPOSED  
 GENERAL MODEL AT T = 0.0 ° F  
 P = 500.0 PSIA

ID	COMPONENT	X(EXP)	Y(EXP)	X(MODEL)	Y(MODEL)	X(MODEL)	EPPX	ERRY
1	CO2	0.00321	0.00594	0.00339	0.005937	1.65729	0.00038	0.000003
2	N2	0.00013	0.00281	0.00014	0.002811	19.42580	0.00002	0.000004
3	C1	0.19913	0.96625	0.20489	0.966350	4.46485	0.01752	0.001060
4	C2	0.03118	0.02185	0.03493	0.021808	0.59027	0.00580	0.000023
5	C3	0.01306	0.00204	0.01372	0.002021	0.13948	0.00145	0.000019
6	IC4	0.00630	0.00030	0.00505	0.000303	0.05674	0.00096	0.000000
7	NC4	0.00980	0.00029	0.00745	0.000293	0.03719	0.00193	0.000000
8	IC5	0.01213	0.00014	0.00922	0.000135	0.01383	0.00239	0.000000
9	NC5	0.01204	0.00010	0.00948	0.000096	0.00755	0.00201	0.000000
10	C6	0.03781	0.00011	0.03850	0.000103	0.00253	0.00296	0.000002
11	C7	0.08907	0.00011	0.13373	0.000095	0.00067	0.05234	0.000010
12	C8	0.16319	0.00005	0.19383	0.000035	0.00017	0.04174	0.000012
13	C9	0.20915	0.00001	0.16724	0.000008	0.00004	0.03131	0.000006
14	C10	0.14774	0.00001	0.12614	0.000001	0.00001	0.01436	0.000005
15	C11	0.05335	0.00000	0.03524	0.000000	0.00000	0.01186	0.000001
16	C12	0.02003	0.00000	0.01297	0.000000	0.00000	0.00632	0.000000
SUM =		1.00000	1.000002	1.00000	1.000002	1.2037	1.2037	0.007143

THE AMOUNT OF COMPONENT, N = 16

T = 50.00 ° F      P = 500.00 PSIA      R = 10.730 PSIA-CLBIC FT/LB-MOLE<sup>0</sup>R

FEED RATE = 1.320500 MOLE/HR      V/F = 0.500000

TABLE 5.40 COMPARISON OF THE LIQUID PHASE AND

THE VAPOR PHASE COMPOSITIONS FROM

THE EXPERIMENTAL DATA AND THE

PROPOSED GENERAL MODEL AT T = 50.0° F

P = 500.0 PSIA

INITIAL VALUE OF LIQUID AND VAPOR PHASE COMPOSITION

ID	COMPONENT	X	Y
1	CO2	0.062500	0.002595
2	N2	0.062500	0.008323
3	C1	0.062500	0.851252
4	C2	0.062500	0.076620
5	C3	0.062500	0.003315
6	IC4	0.062500	0.012398
7	NC4	0.062500	0.008630
8	IC5	0.062500	0.003476
9	NC5	0.062500	0.001519
10	C6	0.062500	0.001224
11	C7	0.062500	0.000387
12	C8	0.062500	0.000137
13	C9	0.062500	0.000034
14	C10	0.062500	0.000011
15	C11	0.062500	0.000001
16	C12	0.062500	0.000000
SUM =		1.000000	0.963920

V/F = 0.373352700

V = 1.285310000 MOLE/HR

L = 0.035219100 MOLE/HR

NO. OF ITERATION = 2

ID	COMPONENT	XI (EXP)	YI (EXP)	XI (MODEL)	YI (MODEL)	KI (MODEL)	ERRX	ERRY
1	CO2	0.00103	0.00260	0.00125	0.002667	2.28385	0.00012	0.000039
2	N2	0.00039	0.00833	0.00062	0.008642	14.99790	0.00018	0.000205
3	C1	0.14335	0.85191	0.19602	0.830545	4.81664	0.03715	0.017702
4	C2	0.07636	0.07662	0.08945	0.077423	0.92806	0.00603	-0.000158
5	C3	0.10988	0.00331	0.01207	0.003141	0.27959	-0.09876	-0.000206
6	IC4	0.10324	0.01231	0.10160	0.012477	0.13168	-0.00367	0.000007
7	NC4	0.11279	0.00853	0.10023	0.008660	0.09264	-0.02047	0.000019
8	IC5	0.10690	0.00338	0.09354	0.003442	0.03946	-0.02075	0.000019
9	NC5	0.06257	0.00146	0.05452	0.001472	0.02934	-0.01235	0.000011
10	C6	0.11327	0.00112	0.12694	0.001130	0.00954	0.00365	-0.000004
11	C7	0.07512	0.00003	0.10353	0.000302	0.00312	0.02025	0.000264
12	C8	0.05875	0.00008	0.07793	0.000071	0.00093	0.01305	-0.000012
13	C9	0.02627	0.00001	0.02961	0.000009	0.00031	0.00100	-0.000001
14	C10	0.00901	0.00000	0.01140	0.000001	0.00010	0.00155	-0.000001
15	C11	0.00103	0.00000	0.00113	0.000000	0.00004	0.00004	-0.000000
16	C12	0.00000	0.00000	0.00000	0.000000	0.00001	0.00000	0.000000
SUM =		1.00000	1.00000			KAAD =	1.93154	0.116773

THE AMOUNT OF COMPONENT, N = 16

T = 47.00 ° F      P = 815.00 PSIA      R = 10.730 PSIA-CUBIC FT/LB-MOLE<sup>OR</sup>

FEED RATE = 2.533670 MOLE/HR      V/F = 0.500000

INITIAL VALUE OF LIQUID AND VAPOR PHASE COMPOSITION

ID	COMPONENT	X	Y
1	CO2	0.062500	0.006617
2	N2	0.062500	0.029980
3	C1	0.062500	0.923041
4	C2	0.062500	0.030932
5	C3	0.062500	0.005329
6	IC4	0.062500	0.001090
7	NC4	0.062500	0.001041
8	IC5	0.062500	0.000499
9	NC5	0.062500	0.000368
10	C6	0.062500	0.000423
11	C7	0.062500	0.000315
12	C8	0.062500	0.000198
13	C9	0.062500	0.000098
14	C10	0.062500	0.000045
15	C11	0.062500	0.000017
16	C12	0.062500	0.000007
SUM =		1.000000	0.999999

V/F = 0.794394600  
V = 2.623856000 MOLE/HR  
L = 0.014313420 MOLE/HR

NO. OF ITERATION = 2

TABLE 5.41 COMPARISON OF THE LIQUID PHASE AND  
 THE VAPOR PHASE COMPOSITIONS FROM  
 THE EXPERIMENTAL DATA AND THE  
 PROPOSED GENERAL MODEL AT T = 47.0 ° F  
 P = 815.0 PSIA

ID	COMPONENT	X(EXP)	Y(EXP)	X(MODEL)	Y(MODEL)	K(MODEL)	ERRX	ERRY
1	CO2	0.00405	0.00662	0.00362	0.006608	1.69211	-0.00013	0.000015
2	N2	0.00253	0.02998	0.00235	0.030024	11.81847	0.00002	0.000153
3	C1	0.21503	0.92309	0.23417	0.923417	3.65476	0.03856	0.003728
4	C2	0.04723	0.03093	0.03957	0.030751	0.71995	-0.00436	-0.000066
5	C3	0.02556	0.00533	0.02116	0.005211	0.22825	-0.00265	-0.000091
6	IC4	0.01191	0.00109	0.00853	0.001085	0.11789	-0.00267	-0.000000
7	NC4	0.01752	0.00104	0.01141	0.001036	0.03414	-0.00517	-0.000000
8	IC5	0.01744	0.00050	0.01177	0.000496	0.03903	-0.00469	-0.000000
9	NC5	0.01503	0.00037	0.01151	0.000365	0.02940	-0.00257	-0.000001
10	C6	0.04752	0.00042	0.02654	0.000416	0.01056	-0.00795	-0.000002
11	C7	0.08236	0.00031	0.07362	0.000304	0.00383	-0.00264	-0.000004
12	C8	0.12308	0.00019	0.12553	0.000181	0.00133	0.01133	0.000000
13	C9	0.13932	0.00009	0.15376	0.000077	0.00046	0.02719	0.000011
14	C10	0.11951	0.00004	0.14436	0.000026	0.00017	0.03632	0.000011
15	C11	0.07485	0.00001	0.07746	0.000007	0.00003	0.00904	0.000004
16	C12	0.05706	0.00000	0.04421	0.000001	0.00003	-0.00910	-0.000002
SUM =		1.00000	1.000004			FAAC =	1.04356	0.025677

THE AMOUNT OF COMPONENT, N = 16

T = -8.00 ° F.      P = 250.00    PSIA      R = 10.730    PSIA-CUBIC FT/LB-MOLE<sup>0.8</sup>  
 FEED RATE = 2.639800    MOLE/HR      V/F = 0.500000

INITIAL VALUE OF LIQUID AND VAPOR PHASE COMPOSITION

ID	COMPONENT	X	Y
1	CO2	0.062500	0.004000
2	N2	0.062500	0.035993
3	C1	0.062500	0.894976
4	C2	0.062500	0.044821
5	C3	0.062500	0.014038
6	IC4	0.062500	0.001311
7	NC4	0.062500	0.002835
8	IC5	0.062500	0.000622
9	NC5	0.062500	0.000632
10	C6	0.062500	0.000294
11	C7	0.062500	0.000264
12	C8	0.062500	0.000144
13	C9	0.062500	0.000045
14	C10	0.062500	0.000018
15	C11	0.062500	0.000006
16	C12	0.062500	0.000002
SUM =		1.000000	1.000000

V/F = 0.993023200  
 V = 2.621387000    MOLE/HR  
 L = 0.018410630    MOLE/HR

NO. OF ITERATION = 2

THE VAPOR PHASE COMPOSITIONS FROM  
 THE EXPERIMENTAL DATA AND THE  
 PROPOSED GENERAL MODEL AT T = -8.0 ° F  
 P = 250.0 PSIA

ID	COMPONENT	X(EXP)	Y(EXP)	X(MODEL)	Y(MODEL)	K(MODEL)	EPRX	TRAY
1	CO2	0.00175	0.00400	0.00220	0.004000	2.46668	-0.00012	0.000016
2	N2	0.00110	0.03601	0.00180	0.036085	26.02831	0.00029	0.000274
3	C1	0.11403	0.89536	0.18672	0.896543	6.52544	0.02394	0.004931
4	C2	0.06160	0.04481	0.07451	0.044561	0.80840	-0.00625	-0.000066
5	C3	0.00821	0.01400	0.10653	0.013527	0.17257	0.07050	-0.000421
6	IC4	0.02246	0.00130	0.02563	0.001272	0.06748	-0.00353	-0.000023
7	NC4	0.07207	0.00280	0.08581	0.002709	0.04291	-0.00866	-0.000050
8	IC5	0.04464	0.00060	0.05135	0.000549	0.01453	-0.00669	-0.000040
9	NC5	0.06466	0.00060	0.07363	0.000528	0.00974	-0.01022	-0.000070
10	C6	0.09214	0.00025	0.09572	0.000161	0.00229	-0.02142	-0.000083
11	C7	0.15476	0.00019	0.14774	0.000053	0.00054	-0.04560	-0.000129
12	C8	0.16507	0.00006	0.09720	0.000008	0.00012	-0.09325	-0.000354
13	C9	0.07507	0.00001	0.03230	0.000001	0.00003	-0.05121	-0.000003
14	C10	0.03295	0.00000	0.01263	0.000000	0.00001	-0.02366	-0.000001
15	C11	0.01160	0.00000	0.00424	0.000000	0.00000	-0.00346	-0.000000
16	C12	0.00393	0.00000	0.00144	0.000000	0.00000	-0.00287	0.000000
SUM =		1.00000	1.00000	1.00000	1.000000	1.00000	2.35406	0.03956

THE AMOUNT OF COMPONENT,N = 16

T = -8.00 °F      P = 1115.00 PSIA      R = 10.730 PSIA-CUBIC FT/LB-MOLE°R

FEED RATE = 2.638390 MOLE/HR      V/F = 0.500000

INITIAL VALUE OF LIQUID AND VAPOR PHASE COMPOSITION

ID	COMPONENT	X	Y
1	CO2	0.062500	0.007097
2	N2	0.062500	0.042107
3	C1	0.062500	0.909745
4	C2	0.062500	0.030471
5	C3	0.062500	0.005833
6	IC4	0.062500	0.001356
7	NC4	0.062500	0.001287
8	IC5	0.062500	0.000599
9	NC5	0.062500	0.000371
10	C6	0.062500	0.000495
11	C7	0.062500	0.000274
12	C8	0.062500	0.000101
13	C9	0.062500	0.000046
14	C10	0.062500	0.000015
15	C11	0.062500	0.000004
16	C12	0.062500	0.000002
SUM =		1.000000	1.000004

V/F = 0.709463500  
V = 2.611078000 MOLE/HR  
L = 0.027812000 MOLE/HR

NO. OF ITERATION = 2

TABLE 5.43 COMPARISON OF THE LIQUID PHASE AND THE VAPOR PHASE COMPOSITIONS FROM THE EXPERIMENTAL DATA AND THE PROPOSED GENERAL MODEL AT T = -8.0°F  
P = 1115.0 PSIA

ID	COMPONENT	X(EXP)	Y(EXP)	X(MODEL)	Y(MODEL)	K(MODEL)	ERRX	ERY
1	CO2	0.00589	0.00710	0.0085	0.007062	0.89118	0.00206	-0.000007
2	N2	0.00472	0.04211	0.01081	0.042299	4.43717	0.00485	0.000346
3	C1	0.37190	0.91002	0.59602	0.910691	1.73038	0.15631	0.003984
4	C2	0.04604	0.03047	0.07881	0.029944	0.43086	0.02371	-0.000416
5	C3	0.03491	0.00583	0.02983	0.005500	0.15653	0.00035	-0.000309
6	IC4	0.02137	0.00135	0.01257	0.001350	0.12177	-0.01025	0.000007
7	NC4	0.02995	0.00128	0.01673	0.001281	0.03653	-0.01509	0.000003
8	IC5	0.03240	0.00059	0.01407	0.000595	0.04803	-0.01995	0.000003
9	NC5	0.02597	0.00037	0.01168	0.000368	0.03576	-0.01564	0.000002
10	C6	0.07432	0.00048	0.02626	0.000490	0.01531	-0.04223	0.000007
11	C7	0.09805	0.00026	0.04573	0.000268	0.02665	-0.05756	0.000007
12	C8	0.10591	0.00006	0.03962	0.000097	0.00276	-0.07005	0.000033
13	C9	0.07952	0.00003	0.04232	0.000041	0.00110	-0.04206	0.000007
14	C10	0.04317	0.00001	0.02733	0.000012	0.00051	-0.01894	0.000003
15	C11	0.01696	0.00000	0.01117	0.000003	0.00031	-0.00708	0.000001
16	C12	0.00834	0.00000	0.00713	0.000001	0.00015	-0.00263	0.000001
SUM =		1.00000	1.000002			RAAD =	3.05958	0.032132

THE AMOUNT OF COMPONENT, N = 16

T = 0.00 ° F      P = 300.00 PSIA      R = 10.730 PSIA-CUBIC FT/LB-MOLE<sup>0</sup>R

FEED RATE = 2.250830 MOLE/HR      V/F = 0.500000

INITIAL VALUE OF LIQUID AND VAPOR PHASE COMPOSITION

ID	COMPONENT	X	Y
1	CO2	0.062500	0.006002
2	N2	0.062500	0.003101
3	C1	0.062500	0.964610
4	C2	0.062500	0.022708
5	C3	0.062500	0.002301
6	IC4	0.062500	0.000330
7	NC4	0.062500	0.000311
8	IC5	0.062500	0.000151
9	NC5	0.062500	0.000111
10	C6	0.062500	0.000123
11	C7	0.062500	0.000112
12	C8	0.062500	0.000075
13	C9	0.062500	0.000040
14	C10	0.062500	0.000019
15	C11	0.062500	0.000004
16	C12	0.062500	0.000001
SUM =		1.000000	0.999999

V/F = 0.99543700

V = 2.247332000 MOLE/HR

L = 0.003327370 MOLE/HR

NO. OF ITERATION = 2

TABLE 5.44 COMPARISON OF THE LIQUID PHASE AND  
 THE VAPOR PHASE COMPOSITIONS FROM  
 THE EXPERIMENTAL DATA AND THE  
 PROPOSED GENERAL MODEL AT T = 0.0 ° F  
 P = 300.0 PSIA

ID	COMPONENT	X(EXP)	Y(EXP)	X(MODEL)	Y(MODEL)	K(MODEL)	ERRX	ERRY
1	C02	0.00172	0.00600	0.00197	0.006001	2.53359	0.00065	0.000005
2	N2	0.00006	0.00310	0.00008	0.003102	32.63286	0.00004	0.000004
3	C1	0.12591	0.96467	0.11054	0.964791	7.25635	0.00719	0.001151
4	C2	0.01849	0.02271	0.02244	0.022678	0.84015	0.00853	-0.000007
5	C3	0.00894	0.00230	0.01050	0.002284	0.17978	0.00378	-0.000015
6	IC4	0.00425	0.00033	0.00429	0.000329	0.06390	0.00091	-0.000000
7	NC4	0.00706	0.00031	0.00625	0.000307	0.04117	0.00046	-0.000000
8	IC5	0.01056	0.00015	0.00850	0.000149	0.01395	0.00015	-0.000001
9	NC5	0.01112	0.00011	0.00951	0.000109	0.00949	0.00042	-0.000001
10	C6	0.04042	0.00012	0.04310	0.000117	0.00225	0.01140	-0.000003
11	C7	0.09897	0.00010	0.14106	0.000091	0.00054	0.07089	-0.000014
12	C8	0.18355	0.00006	0.25673	0.000037	0.00012	0.12559	-0.000024
13	C9	0.24508	0.00002	0.22122	0.000007	0.00003	0.02130	-0.000015
14	C10	0.17514	0.00001	0.12437	0.000001	0.00001	0.02537	-0.000005
15	C11	0.05016	0.00000	0.02943	0.000000	0.00000	0.01472	-0.000001
16	C12	0.01859	0.00000	0.00543	0.000000	0.00000	0.00717	-0.000000
SUM =		1.00000	1.00000	1.00004	1.00000	ERRAD =	1.86663	0.00778

THE AMOUNT OF COMPONENT, N = 16

T = 8.00 °F      P = 300.00 PSIA      R = 10.730 PSIA-CUBIC FT/LB-MOLE °R

FEED RATE = 1.153030 MOLE/HR      V/F = 0.500000

INITIAL VALUE OF LIQUID AND VAPOR PHASE COMPOSITION

ID	COMPONENT	X	Y
1	C22	0.062500	0.010069
2	N2	0.062500	0.004437
3	C1	0.062500	0.951893
4	C2	0.062500	0.031772
5	C3	0.062500	0.001574
6	IC4	0.062500	0.000015
7	NC4	0.062500	0.000014
8	IC5	0.062500	0.000018
9	NC5	0.062500	0.000017
10	C6	0.062500	0.000053
11	C7	0.062500	0.000056
12	C8	0.062500	0.000034
13	C9	0.062500	0.000013
14	C10	0.062500	0.000027
15	C11	0.062500	0.000006
16	C12	0.062500	0.000000
SUM =		1.000000	1.000000

V/F = 0.500000

V = 1.152317000 MOLE/HR

L = 0.000232677 MOLE/HR

NO. OF ITERATION = 2

ID	COMPONENT	X(EXP)	Y(EXP)	X(MODEL)	Y(MODEL)	K(MODEL)	ERRX	EPY
1	C02	0.00262	0.01007	0.00534	0.010069	2.54509	0.00134	0.000000
2	N2	0.00007	0.00444	0.00023	0.004438	25.83601	0.00010	0.000000
3	C1	0.07704	0.95193	0.19602	0.951963	6.55877	0.06811	0.000064
4	C2	0.03243	0.03177	0.04925	0.031771	0.87122	0.00404	-0.000000
5	C3	0.01646	0.00157	0.01065	0.001573	0.19938	-0.00857	-0.000000
6	IC4	0.00008	0.00002	0.00027	0.000015	0.07435	0.00012	-0.000000
7	NC4	0.00021	0.00001	0.00033	0.000014	0.04851	0.00008	-0.000000
8	IC5	0.00130	0.00002	0.00146	0.000019	0.01682	-0.00022	-0.000000
9	NC5	0.00173	0.00002	0.00193	0.000017	0.01158	-0.00026	-0.000000
10	C6	0.02294	0.00005	0.02426	0.000051	0.00285	-0.00497	-0.000001
11	C7	0.00676	0.00005	0.05173	0.000048	0.00070	0.00314	-0.000006
12	C8	0.00969	0.00003	0.10053	0.000019	0.00016	0.01922	-0.000011
13	C9	0.011754	0.00001	0.10633	0.000023	0.00004	-0.03993	-0.000005
14	C10	0.042584	0.00001	0.27007	0.000022	0.00001	-0.01994	-0.000009
15	C11	0.012610	0.00000	0.06797	0.000020	0.00000	-0.07577	0.000000
16	C12	0.01119	0.00000	0.00553	0.000000	0.00000	-0.00712	0.000000
SUM =		1.00000	1.000001			RAAD =	2.92394	0.000009

TABLE 5.45 COMPARISON OF THE LIQUID PHASE AND  
 THE VAPOR PHASE COMPOSITIONS FROM  
 THE EXPERIMENTAL DATA AND THE  
 PROPOSED GENERAL MODEL AT T = 8.0 °F  
 P = 300.0 PSIA

THE AMOUNT OF COMPONENT, N = 16

T = 8.00 °F      P = 700.00 PSIA      R = 10.730 PSIA-CUBIC FT/LB-MOLE °R

FEED RATE = 2.453330 MOLE/HR      V/F = 0.500000

INITIAL VALUE OF LIQUID AND VAPOR PHASE COMPOSITION

ID	COMPONENT	X	Y
1	C02	0.062500	0.010155
2	I2	0.062500	0.004651
3	C1	0.062500	0.031288
4	C2	0.062500	0.031921
5	C3	0.062500	0.031755
6	IC4	0.062500	0.030015
7	NC4	0.062500	0.030014
8	IC5	0.062500	0.000013
9	NC5	0.062500	0.030015
10	C6	0.062500	0.030043
11	C7	0.062500	0.030051
12	C8	0.062500	0.000031
13	C9	0.062500	0.000013
14	C10	0.062500	0.000024
15	C11	0.062500	0.000004
16	C12	0.062500	0.000001
SUM =		1.000000	0.991999

V/F = 0.193877900

V = 2.453342030 MOLE/HR

L = 0.003337872 MOLE/HR

NO. OF ITERATION = 2

THE VAPOR PHASE COMPOSITIONS FROM  
 THE EXPERIMENTAL DATA AND THE  
 PROPOSED GENERAL MODEL AT T = 8.0 °F  
 P = 700.0 PSIA

ID	COMPONENT	X(EXP)	Y(EXP)	X(MODEL)	Y(MODEL)	K(MODEL)	ERRX	ERRY
1	C02	0.00265	0.01016	0.01027	0.010155	1.28749	0.00503	0.000000
2	I2	0.00009	0.00465	0.00064	0.004651	9.54257	0.00040	0.000000
3	C1	0.25105	0.03131	0.42403	0.031340	2.92411	0.07430	0.000055
4	C2	0.02279	0.03192	0.08073	0.031917	0.51531	0.03515	0.000004
5	C3	0.00885	0.00176	0.01572	0.001754	0.14545	0.00321	0.000001
6	IC4	0.00005	0.03001	0.00023	0.030015	0.06959	0.00017	0.000000
7	NC4	0.00015	0.03001	0.00037	0.030014	0.04701	0.00015	0.000000
8	IC5	0.00092	0.00001	0.00060	0.000013	0.01963	0.00026	0.000000
9	NC5	0.00128	0.00001	0.00137	0.000015	0.01397	-0.00022	0.000000
10	C6	0.01516	0.00005	0.01403	0.000046	0.03429	0.00436	0.000001
11	C7	0.04357	0.00005	0.04622	0.000047	0.00132	0.00011	0.000003
12	C8	0.06752	0.00003	0.07667	0.000023	0.00039	0.000757	0.000005
13	C9	0.10162	0.00001	0.07203	0.000006	0.00011	0.004623	0.000004
14	C10	0.06321	0.00001	0.20250	0.000005	0.00004	0.20785	0.000001
15	C11	0.03597	0.00000	0.04275	0.000000	0.00002	0.06317	0.000001
16	C12	0.02452	0.00000	0.00924	0.000000	0.00000	0.01743	0.000000
SUM =		1.00000	1.00000	1.00000	1.00000	1.00000	2.52500	0.000000

TABLE 5.46 COMPARISON OF THE LIQUID PHASE AND



THE AMOUNT OF COMPONENT, N = 16

T = -10.00 ° F      P = 500.00 PSIA      R = 10.730      PSIA-CUBIC FT/LB-MOLE ° R

FEED RATE = 2.639830 MOLE/HR      V/F = 0.500000

INITIAL VALUE OF LIQUID AND VAPOR PHASE COMPOSITION

ID	COMPONENT	X	Y
1	C12	0.062500	0.005544
2	N2	0.062500	0.002730
3	C1	0.062500	0.966292
4	C2	0.062500	0.022314
5	C3	0.062500	0.002002
6	IC4	0.062500	0.000317
7	NC4	0.062500	0.000294
8	IC5	0.062500	0.000138
9	NC5	0.062500	0.000101
10	C6	0.062500	0.000083
11	C7	0.062500	0.000079
12	C8	0.062500	0.000048
13	C9	0.062500	0.000034
14	C10	0.062500	0.000019
15	C11	0.062500	0.000005
16	C12	0.062500	0.000001
SUM =		1.000000	0.999999

V/F = 0.993513200  
 V = 2.63983000 MOLE/HR  
 L = 0.003934942 MOLE/HR

NO. OF ITERATION = 2

TABLE 5.47 COMPARISON OF THE LIQUID PHASE AND  
 THE VAPOR PHASE COMPOSITIONS FROM THE  
 ORDINARY VLE CALCULATION MODEL AND THE  
 PROPOSED GENERAL MODEL AT T = -10.0 ° F  
 P = 500.0 PSIA

ID	COMPONENT	X(ORDINARY)	Y(ORDINARY)	X(MODEL)	Y(MODEL)	K(MODEL)	ERRX	ERRY
1	C02	0.00419	0.00554	0.00331	0.005541	1.51901	-0.00054	0.000002
2	N2	0.00019	0.00273	0.00013	0.002731	18.55620	-0.00004	0.000004
3	C1	0.27078	0.96645	0.20444	0.966471	4.23653	-0.04510	0.000943
4	C2	0.04809	0.02231	0.03812	0.022263	0.52955	-0.00601	-0.000024
5	C3	0.01973	0.00200	0.01493	0.001979	0.11976	-0.00219	-0.000018
6	IC4	0.00827	0.00031	0.00593	0.000316	0.04848	-0.00175	0.000001
7	NC4	0.01195	0.00029	0.00847	0.000293	0.03132	-0.00260	0.000007
8	IC5	0.01575	0.00001	0.01080	0.000136	0.01143	-0.00383	0.0000123
9	NC5	0.01687	0.00010	0.01155	0.000099	0.00777	-0.00412	0.000002
10	C6	0.05145	0.00007	0.02573	0.000079	0.00200	-0.01205	0.000007
11	C7	0.14098	0.00005	0.11303	0.000064	0.00051	-0.01615	0.000016
12	C8	0.15900	0.00001	0.17495	0.000024	0.00012	0.03413	0.000012
13	C9	0.14164	0.00000	0.20102	0.000007	0.00003	0.08027	0.000004
14	C10	0.08348	0.00000	0.13131	0.000001	0.00001	0.06147	0.000001
15	C11	0.02275	0.00000	0.03667	0.000000	0.00000	0.01775	0.000000
16	C12	0.00583	0.00000	0.00951	0.000000	0.00000	0.00467	0.000000
SUM =		1.00000	1.00000	1.00000	1.00000	9.440 =	1.92915	0.007239

THE AMOUNT OF COMPONENT, N = 16

T = 0.00 ° F      P = 500.00 PSIA      R = 10.730      FSIA-CUDIC FT/LB-MOLE<sup>OR</sup>

FEED RATE = 2.638710 MOLE/HR      V/F = 0.500000

INITIAL VALUE OF LIQUID AND VAPOR PHASE COMPOSITION

ID	COMPONENT	X	Y
1	CO2	0.062500	0.005740
2	N2	0.062500	0.002810
3	C1	0.062500	0.966192
4	C2	0.062500	0.021852
5	C3	0.062500	0.002042
6	IC4	0.062500	0.000303
7	NC4	0.062500	0.000294
8	IC5	0.062500	0.000136
9	NC5	0.062500	0.000097
10	C6	0.062500	0.000108
11	C7	0.062500	0.000113
12	C8	0.062500	0.000060
13	C9	0.062500	0.000029
14	C10	0.062500	0.000018
15	C11	0.062500	0.000005
16	C12	0.062500	0.000002
SUM =		1.000000	1.000000

V/F = 0.113506300  
 V = 2.634722000 MOLE/HR  
 L = 0.003987312 MOLE/HR

NO. OF ITERATION = 2

TABLE 5.48 COMPARISON OF THE LIQUID PHASE AND THE VAPOR PHASE COMPOSITIONS FROM THE ORDINARY VLE CALCULATION MODEL AND THE PROPOSED GENERAL MODEL AT T = 0.0° F P = 500.0 PSIA

ID	COMPONENT	X(ORDINARY)	Y(ORDINARY)	X(MODEL)	Y(MODEL)	K(MODEL)	ERRX	ERRY
1	CO2	0.00418	0.00594	0.00337	0.005937	1.65729	-0.00059	0.000003
2	N2	0.00020	0.00281	0.00014	0.002811	18.42580	-0.00005	0.000004
3	C1	0.26293	0.96635	0.20487	0.966358	4.46485	-0.04628	0.000966
4	C2	0.04292	0.02185	0.03491	0.021808	0.59027	-0.00574	-0.000018
5	C3	0.01667	0.00204	0.01372	0.002021	0.13948	-0.00216	-0.000016
6	IC4	0.00674	0.00030	0.00563	0.000303	0.03674	-0.00140	0.000001
7	NC4	0.01000	0.00029	0.00745	0.000273	0.03719	-0.00213	0.000001
8	IC5	0.01273	0.00013	0.00922	0.000135	0.01383	-0.00278	0.000007
9	NC5	0.01312	0.00007	0.00948	0.000096	0.00955	-0.00309	0.000002
10	C6	0.05304	0.00010	0.03856	0.000103	0.00253	-0.01227	0.000006
11	C7	0.16517	0.00008	0.13373	0.000095	0.00067	-0.02376	0.000013
12	C8	0.18167	0.00002	0.19380	0.000035	0.00017	0.02376	0.000015
13	C9	0.11964	0.00000	0.16724	0.000008	0.00004	0.05720	0.000005
14	C10	0.07936	0.00000	0.12614	0.000001	0.00001	0.05402	0.000000
15	C11	0.02387	0.00000	0.02924	0.000000	0.00000	0.01762	0.000000
16	C12	0.00776	0.00000	0.01297	0.000000	0.00000	0.00575	0.000000
SUM =		1.00000	1.00000	1.00000	1.00000	ZAAD =	1.61634	0.00000

THE AMOUNT OF COMPONENT A = 15

T = 50.00 °F      P = 500.00 PSIA      K = 10.730      PSIA-CUBIC FT/LB-MOLE<sup>0</sup>

FEED RATE = 1.111310 MOLE/HR      V/F = 0.500000

INITIAL VALUE OF LIQUID AND VAPOR PHASE COMPOSITION

ID	COMPONENT	X	Y
1	CO2	0.062500	0.002395
2	N2	0.062500	0.003323
3	C1	0.062500	0.031252
4	C2	0.062500	0.076620
5	C3	0.062500	0.003315
6	IC4	0.062500	0.023980
7	NC4	0.062500	0.008630
8	IC5	0.062500	0.003476
9	NC5	0.062500	0.001519
10	C6	0.062500	0.001224
11	C7	0.062500	0.000387
12	C8	0.062500	0.000137
13	C9	0.062500	0.000034
14	C10	0.062500	0.000011
15	C11	0.062500	0.000001
16	C12	0.062500	0.000000
SUM =		1.000000	0.969720

V/F = 0.500000  
 V = 1.28531000 MOLE/HR  
 L = 0.03321710 MOLE/HR

NO. OF ITERATION = 2

THE VAPOR PHASE COMPOSITIONS FROM  
 THE ORDINARY VLE CALCULATION MODEL  
 AND THE PROPOSED GENERAL MODEL AT  
 T = 50.0 °F    P = 500.0 PSIA

TABLE 5.49 COMPARISON OF THE LIQUID PHASE AND

ID	COMPONENT	X(ORDINARY)	Y(ORDINARY)	X(MODEL)	Y(MODEL)	K(MODEL)	ERRX	ERRY
1	CO2	0.00115	0.00263	0.00125	0.00267	2.28365	0.00000	-0.000043
2	I2	0.00054	0.00859	0.00062	0.00864	14.99790	0.00003	-0.000056
3	C1	0.17623	0.07857	0.19602	0.083545	4.81666	0.00431	0.003956
4	C2	0.00842	0.07899	0.00845	0.077423	0.92806	0.07397	-0.002527
5	C3	0.01225	0.00034	0.01207	0.001141	0.27895	-0.00113	0.002761
6	IC4	0.10823	0.01256	0.10163	0.013477	0.13168	-0.01466	-0.000335
7	NC4	0.10657	0.00877	0.10023	0.003650	0.09264	-0.01465	-0.000217
8	IC5	0.10281	0.00345	0.09354	0.003442	0.03946	-0.01666	-0.000055
9	NC5	0.05985	0.00119	0.05452	0.001492	0.02324	-0.00063	-0.000017
10	C6	0.13943	0.00108	0.12654	0.001130	0.00954	-0.00251	0.000035
11	C7	0.10685	0.00003	0.10033	0.003302	0.00312	-0.01148	0.000272
12	C8	0.00698	0.00005	0.00751	0.000071	0.00058	0.06482	0.000020
13	C9	0.02287	0.00001	0.02561	0.000007	0.00021	0.00440	-0.000004
14	C10	0.00817	0.00000	0.00814	0.000001	0.00010	0.00239	0.000000
15	C11	0.00032	0.00000	0.00031	0.000000	0.00004	0.00027	0.000000
16	C12	0.00000	0.00000	0.00000	0.000000	0.00001	0.00000	0.000000
SUM =		1.00000	1.00000	1.00000	1.00000	2.440	1.50579	0.095610



THE AMOUNT OF COMPONENT, N = 16

T = 47.00 ° F      P = 815.00 PSIA      R = 10.730 FSIA-CUBIC FT/LB-MOLE°R

FEED RATE = 2.638670 MOLE/HR      V/F = 0.500000

INITIAL VALUE OF LIQUID AND VAPOR PHASE COMPOSITION

ID	COMPONENT	X	Y
1	CO2	0.062500	0.006617
2	N2	0.062500	0.029980
3	C1	0.062500	0.923041
4	C2	0.062500	0.030932
5	C3	0.062500	0.005329
6	IC4	0.062500	0.001090
7	NC4	0.062500	0.001041
8	IC5	0.062500	0.000499
9	NC5	0.062500	0.000368
10	C6	0.062500	0.000423
11	C7	0.062500	0.000315
12	C8	0.062500	0.000198
13	C9	0.062500	0.000098
14	C10	0.062500	0.000045
15	C11	0.062500	0.000017
16	C12	0.062500	0.000007
SUM =		1.000000	0.999999

V/F = 0.394394600  
 V = 2.623356000 MOLE/HR  
 L = 0.014013420 MOLE/HR

NO. OF ITERATION = 2

TABLE 5.50 COMPARISON OF THE LIQUID PHASE AND THE VAPOR PHASE COMPOSITIONS FROM THE ORDINARY VLE CALCULATION MODEL AND THE PROPOSED GENERAL MODEL AT T = 47.0° F P = 815.0 PSIA

ID	COMPONENT	X(ORDINARY)	Y(ORDINARY)	X(MODEL)	Y(MODEL)	K(MODEL)	ERRX	EPRY
1	CO2	0.00398	0.00662	0.00362	0.00668	1.69211	-0.00006	0.000014
2	N2	0.00271	0.02997	0.00233	0.030024	11.81847	-0.00016	0.000144
3	C1	0.26213	0.92329	0.23417	0.923417	3.65476	-0.00854	0.003525
4	C2	0.04362	0.03093	0.03553	0.030751	0.71995	-0.00075	-0.000062
5	C3	0.02353	0.00532	0.02116	0.005211	0.22825	-0.00062	-0.000062
6	IC4	0.01063	0.00109	0.00853	0.001085	0.11789	-0.00139	0.000032
7	NC4	0.01441	0.00104	0.01141	0.001036	0.08414	-0.00206	0.000093
8	IC5	0.01551	0.00049	0.01177	0.000496	0.03903	-0.00276	0.000024
9	NC5	0.01532	0.00036	0.01151	0.000365	0.02940	-0.00286	0.000024
10	C6	0.05045	0.00042	0.03653	0.000416	0.01056	-0.01072	0.000014
11	C7	0.01040	0.00028	0.07362	0.000304	0.00383	-0.02169	0.000028
12	C8	0.01562	0.00014	0.12593	0.000181	0.00133	-0.01991	0.000042
13	C9	0.014598	0.00004	0.15375	0.000077	0.00046	0.02053	0.000024
14	C10	0.09513	0.00001	0.14433	0.000026	0.00017	0.06120	0.000017
15	C11	0.04021	0.00000	0.07746	0.000007	0.00009	0.00009	0.000005
16	C12	0.01864	0.00000	0.04421	0.000001	0.00003	0.02924	0.000001
SUM =		1.00000	1.000004			1.11111	1.41477	0.024763

THE AMOUNT OF COMPONENT N = 16

T = -8.00 ° F      P = 250.00 PSIA      R = 10.730      FSIA-CUBIC FT/LB-MOLE<sup>0</sup>R

FEED RATE = 2.639800 MOLE/HR      V/F = 0.500000

INITIAL VALUE OF LIQUID AND VAPOR PHASE COMPOSITION

ID	COMPONENT	X	Y
1	CO2	0.062500	0.004000
2	N2	0.062500	0.035993
3	C1	0.062500	0.894976
4	C2	0.062500	0.044821
5	C3	0.062500	0.014038
6	IC4	0.062500	0.001311
7	NC4	0.062500	0.002835
8	IC5	0.062500	0.000622
9	NC5	0.062500	0.000632
10	C6	0.062500	0.000294
11	C7	0.062500	0.000264
12	C8	0.062500	0.000144
13	C9	0.062500	0.000045
14	C10	0.062500	0.000018
15	C11	0.062500	0.000006
16	C12	0.062500	0.000002
SUM =		1.000000	1.000000

V/F = 0.33028200  
V = 2.621389000 MOLE/HR  
L = 0.018410630 MOLE/HR

NO. OF ITERATION = 2

TABLE 5.51 COMPARISON OF THE LIQUID PHASE AND  
 AND THE VAPOR PHASE COMPOSITIONS  
 FROM THE ORDINARY VLE CALCULATION  
 MODEL AND THE PROPOSED GENERAL  
 MODEL AT T = -8.0 ° F P = 250.0 PSIA

ID	COMPONENT	X(ORDINARY)	Y(ORDINARY)	X(MODEL)	Y(MODEL)	K(MODEL)	ERRX	ERRY
1	CO2	0.00145	0.00400	0.00220	0.004000	2.46668	0.00018	0.000015
2	N2	0.00100	0.03603	0.00183	0.036085	26.02831	0.00039	0.000210
3	C1	0.11202	0.89570	0.18672	0.896543	6.52544	0.02595	0.004586
4	C2	0.05057	0.04482	0.07491	0.044561	0.80840	0.00478	-0.000067
5	C3	0.07730	0.01398	0.10653	0.013527	0.17257	-0.00141	-0.000396
6	IC4	0.02189	0.00129	0.02563	0.001272	0.06748	-0.00276	-0.000014
7	NC4	0.07561	0.00277	0.08581	0.002709	0.04291	-0.01220	-0.000066
8	IC5	0.04938	0.00058	0.05133	0.000549	0.01453	-0.01143	-0.000024
9	NC5	0.07365	0.00056	0.07363	0.000528	0.00974	-0.01721	-0.000034
10	C6	0.11381	0.00019	0.09572	0.000161	0.00229	-0.04309	-0.000027
11	C7	0.20462	0.00007	0.14774	0.000053	0.00054	-0.09546	-0.000015
12	C8	0.14317	0.00001	0.09720	0.000003	0.00012	-0.07135	-0.000001
13	C9	0.04816	0.00000	0.03233	0.000001	0.00003	-0.02430	0.000001
14	C10	0.01887	0.00000	0.01263	0.000000	0.00001	-0.00954	0.000000
15	C11	0.00634	0.00000	0.00424	0.000000	0.00000	-0.00320	0.000000
16	C12	0.00215	0.00000	0.00144	0.000000	0.00000	-0.00109	0.000000
SUM =		1.00000	1.00000	1.00000	1.00000	KAAD =	2.04076	0.033992

THE AMOUNT OF COMPONENT, N = 16

T = -8.00 ° F      P = 1115.00    PSIA      R = 10.730    PSIA-CUBIC FT/LB-MOLE<sup>OR</sup>

FEED RATE = 2.6309 JO    MOLE/HR      V/F = 0.500000

INITIAL VALUE OF LIQUID AND VAPOR PHASE COMPOSITION

ID	COMPONENT	X	Y
1	C2	0.062500	0.007097
2	N2	0.062500	0.042107
3	C1	0.062500	0.909945
4	C2	0.062500	0.030471
5	C3	0.062500	0.005833
6	IC4	0.062500	0.001356
7	NC4	0.062500	0.001287
8	IC5	0.062500	0.000599
9	NC5	0.062500	0.000371
10	C6	0.062500	0.000495
11	C7	0.062500	0.000274
12	C8	0.062500	0.000101
13	C9	0.062500	0.000046
14	C10	0.062500	0.000015
15	C11	0.062500	0.000004
16	C12	0.062500	0.000002
SUM =		1.000000	1.000004

V/F = 0.939463500  
 V = 2.6110780 JO    MOLE/HR  
 L = 0.0278120 JO    MOLE/HR

NO. OF ITERATION = 2

TABLE 5.52 COMPARISON OF THE LIQUID PHASE AND THE VAPOR PHASE COMPOSITIONS FROM THE ORDINARY VLE CALCULATION MODEL AND THE PROPOSED GENERAL MODEL AT T = -8.0 ° F    P = 1115.0 PSIA

ID	COMPONENT	X(ORDINARY)	Y(ORDINARY)	X(MODEL)	Y(MODEL)	K(MODEL)	ERRX	ERRY
1	C02	0.00722	0.00710	0.00899	0.007062	0.89113	0.00073	-0.000009
2	N2	0.00601	0.04213	0.01081	0.042299	4.43717	0.00356	0.000326
3	C1	0.41306	0.91023	0.59682	0.910691	1.73038	0.11515	0.003773
4	C2	0.06890	0.03045	0.07881	0.029944	0.43086	0.00005	-0.000395
5	C3	0.04043	0.00581	0.03983	0.005500	0.15658	0.00517	-0.000293
6	IC4	0.01994	0.00135	0.01257	0.001350	0.12177	-0.00882	0.000010
7	NC4	0.02727	0.00127	0.01673	0.001231	0.09653	-0.01241	0.000014
8	IC5	0.02698	0.00058	0.01407	0.000576	0.04803	-0.01453	0.000014
9	NC5	0.02282	0.00036	0.01163	0.000363	0.03576	-0.01249	0.000012
10	C6	0.08227	0.00045	0.02623	0.000490	0.01531	-0.05018	0.000043
11	C7	0.11038	0.00021	0.04575	0.000268	0.00665	-0.05989	0.000053
12	C8	0.08338	0.00015	0.03962	0.000097	0.00275	-0.04832	0.000044
13	C9	0.05801	0.00001	0.04232	0.000041	0.00110	-0.02055	0.000027
14	C10	0.02352	0.00000	0.02739	0.000012	0.00051	0.00071	0.000010
15	C11	0.00697	0.00000	0.01117	0.000003	0.00031	0.00291	0.000003
16	C12	0.00286	0.00000	0.00713	0.000001	0.00015	0.00345	0.000001
SUM =		1.00000	1.000002	ΣAAD =		2.21075	0.031502	

THE AMOUNT OF COMPONENT, N = 16

T = 0.00 °F P = 300.00 PSIA R = 16.720 FSIA-CUBIC FT/LB-MOLE<sup>0</sup>R

FEED RATE = 2.250850 MOLE/HR V/F = 0.500000

INITIAL VALUE OF LIQUID AND VAPOR PHASE COMPOSITION

ID	COMPONENT	X	Y
1	CO2	0.062500	0.006002
2	N2	0.062500	0.003101
3	C1	0.062500	0.764610
4	C2	0.062500	0.022708
5	C3	0.062500	0.002301
6	IC4	0.062500	0.000330
7	NC4	0.062500	0.000311
8	IC5	0.062500	0.000151
9	NC5	0.062500	0.000111
10	C6	0.062500	0.000123
11	C7	0.062500	0.000112
12	C8	0.062500	0.000075
13	C9	0.062500	0.000040
14	C10	0.062500	0.000019
15	C11	0.062500	0.000004
16	C12	0.062500	0.000001
SUM =		1.000000	0.999999

V/F = 0.998543700  
 V = 2.247532000 MOLE/HR  
 L = 0.003327370 MOLE/HR

NO. OF ITERATION = 2

TABLE 5.53 COMPARISON OF THE LIQUID PHASE AND THE VAPOR PHASE COMPOSITIONS FROM THE ORDINARY VLE CALCULATION MODEL AND THE PROPOSED GENERAL MODEL AT T = 0.0°F P = 300.0 PSIA

ID	COMPONENT	X(ORDINARY)	Y(ORDINARY)	X(MODEL)	Y(MODEL)	K(MODEL)	ERRX	EPY
1	CO2	0.062500	0.006000	0.006197	0.006001	2.53359	-0.00033	0.000004
2	N2	0.062500	0.003100	0.003008	0.003102	32.63265	-0.00002	0.000003
3	C1	0.15784	0.96479	0.11054	0.964791	7.25635	-0.02474	0.001030
4	C2	0.03065	0.02271	0.02244	0.022678	0.04015	-0.00363	-0.000005
5	C3	0.01429	0.00230	0.01055	0.002234	0.17978	-0.00157	-0.000013
6	IC4	0.00602	0.00033	0.00427	0.000329	0.06390	-0.00086	-0.000001
7	NC4	0.00891	0.00031	0.00625	0.000309	0.04117	-0.00139	0.000001
8	IC5	0.01285	0.00015	0.00890	0.000147	0.01395	-0.00214	0.000002
9	NC5	0.01397	0.00011	0.00957	0.000109	0.00948	-0.00243	0.000001
10	C6	0.06171	0.00011	0.04310	0.000117	0.00225	-0.00781	0.000003
11	C7	0.17672	0.00007	0.14106	0.000091	0.00054	-0.00686	0.000019
12	C8	0.23975	0.00002	0.25673	0.000037	0.00012	0.06939	0.000014
13	C9	0.16423	0.00000	0.00000	0.000007	0.00003	0.10215	0.000004
14	C10	0.08440	0.00000	0.12437	0.000001	0.00001	0.06535	0.000001
15	C11	0.01959	0.00000	0.02543	0.000000	0.00000	0.01595	0.000000
16	C12	0.00625	0.00000	0.00943	0.000000	0.00000	0.00516	0.000000
SUM =		1.000000	1.000000	1.000000	1.000000	MAAD =	1.94905	0.005926

THE AMOUNT OF COMPONENT, A = 16

T = 8.00 °F P = 300.00 PSIA R = 10.730 PSIA-CUBIC FT/LB-MOLE°R

FEED RATE = 1.153030 MOLE/HR V/F = 0.500000

INITIAL VALUE OF LIQUID AND VAPOR PHASE COMPOSITION

ID	COMPONENT	X	Y
1	C2	0.062500	0.010069
2	V2	0.062500	0.004437
3	C1	0.062500	0.951893
4	C2	0.062500	0.031772
5	C3	0.062500	0.001574
6	IC4	0.062500	0.000015
7	NC4	0.062500	0.000014
8	IC5	0.062500	0.000018
9	NC5	0.062500	0.000017
10	C6	0.062500	0.000053
11	C7	0.062500	0.000056
12	C8	0.062500	0.000034
13	C9	0.062500	0.000013
14	C10	0.062500	0.000027
15	C11	0.062500	0.000006
16	C12	0.062500	0.000000
SUM =		1.000000	1.000000

V/F = 0.33377900  
 V = 1.152317030 MOLE/HR  
 L = 0.032232637 MOLE/HR

NO. OF ITERATION = 2

TABLE 5.54 COMPARISON OF THE LIQUID AND THE VAPOR PHASE COMPOSITIONS FROM THE ORDINARY VLE CALCULATION MODEL AND THE PROPOSED GENERAL MODEL AT T = 8.0°F P = 300.0 PSIA

ID	COMPONENT	X(ORDINARY)	Y(ORDINARY)	X(MODEL)	Y(MODEL)	K(MODEL)	ERRX	ERRY
1	C2	0.00513	0.01007	0.00534	0.010069	2.54509	-0.00117	0.000001
2	N2	0.00022	0.00444	0.00023	0.004438	25.83601	-0.00005	0.000000
3	C1	0.18779	0.95197	0.19602	0.951963	6.55877	-0.04264	0.000024
4	C2	0.04734	0.03177	0.04925	0.031771	0.87122	-0.01087	0.000001
5	C3	0.01026	0.00157	0.01063	0.001573	0.19938	-0.00237	0.000000
6	IC4	0.00028	0.00001	0.00027	0.000015	0.07435	-0.00008	0.000000
7	NC4	0.00040	0.00001	0.00039	0.000014	0.04951	-0.00011	0.000000
9	IC5	0.00154	0.00002	0.00146	0.000018	0.01682	-0.00046	0.000000
9	NC5	0.00210	0.00002	0.00199	0.000017	0.01158	-0.00063	0.000000
10	C6	0.02617	0.00005	0.02426	0.000051	0.00285	-0.00010	0.000000
11	C7	0.09964	0.00005	0.09173	0.000048	0.00070	-0.003174	0.000001
12	C8	0.16972	0.00002	0.16053	0.000019	0.00016	-0.05081	0.000001
13	C9	0.10686	0.00000	0.10633	0.000003	0.00004	-0.02815	0.000000
14	C10	0.27146	0.00000	0.27107	0.000002	0.00001	-0.00556	-0.000000
15	C11	0.06580	0.00000	0.06737	0.000000	0.00000	-0.01947	0.000000
16	C12	0.00530	0.00000	0.00553	0.000000	0.00000	-0.00123	0.000000
SUM =		1.00000	1.00000	1.00000	1.00000	2AAD =	1.62211	0.000176



THE AMOUNT OF COMPONENT A = 16

T = 8.00 °F      P = 700.00 PSIA      R = 10.730 PSIA-LOGIC FT/LB-MOLE<sup>2</sup>R

FEED RATE = 2.453880 MOLE/HR      V/F = 0.500000

INITIAL VALUE OF LIQUID AND VAPOR PHASE COMPOSITION

ID	COMPONENT	X	Y
1	C12	0.062500	0.010155
2	C12	0.062500	0.004651
3	C1	0.062500	0.951288
4	C2	0.062500	0.031921
5	C3	0.062500	0.001755
6	IC4	0.062500	0.000015
7	NC4	0.062500	0.000014
8	IC5	0.062500	0.000013
9	NC5	0.062500	0.000015
10	C6	0.062500	0.000043
11	C7	0.062500	0.000051
12	C9	0.062500	0.000031
13	C9	0.062500	0.000013
14	C10	0.062500	0.000024
15	C11	0.062500	0.000004
15	C12	0.062500	0.000001
SUM =		1.000000	0.999999

V/F = 0.500000  
V = 2.453880 MOLE/HR  
L = 0.000000 MOLE/HR

NO. OF ITERATION = 2

TABLE 5.55 COMPARISON OF THE LIQUID PHASE AND THE VAPOR PHASE COMPOSITIONS FROM THE ORDINARY VLE CALCULATION MODEL AND THE PROPOSED GENERAL MODEL AT T = 8.0 °F P = 700.0 PSIA

ID	COMPONENT	X(ORDINARY)	Y(ORDINARY)	X(MODEL)	Y(MODEL)	K(MODEL)	ERRX	ERRY
1	C12	0.06947	0.01015	0.01027	0.010155	1.28949	-0.00159	0.000000
2	C12	0.00655	0.00465	0.00664	0.004651	9.54257	-0.00006	0.000000
3	C1	0.98169	0.95135	0.94248	0.951390	2.92411	-0.05634	0.000019
4	C2	0.07540	0.03192	0.06673	0.031917	0.51531	-0.01346	0.000000
5	C3	0.01493	0.00175	0.01572	0.001754	0.14545	-0.00287	0.000000
6	IC4	0.00031	0.00001	0.00023	0.000015	0.06959	-0.00009	-0.000000
7	NC4	0.00043	0.00001	0.00039	0.000014	0.04701	-0.00013	-0.000000
8	IC5	0.00100	0.00001	0.00089	0.000013	0.01961	-0.00004	-0.000000
9	NC5	0.00162	0.00001	0.0013	0.000015	0.01397	-0.00006	-0.000000
10	C6	0.01713	0.00004	0.01403	0.000046	0.00423	-0.00063	-0.000000
11	C7	0.05806	0.00005	0.04622	0.000047	0.00132	-0.02260	0.000001
12	C9	0.09724	0.00002	0.07867	0.000023	0.00029	-0.01689	-0.000002
13	C9	0.08130	0.00001	0.07223	0.000006	0.00011	-0.00251	-0.000001
14	C10	0.00555	0.00000	0.00253	0.000005	0.00004	-0.00547	0.000000
15	C11	0.00442	0.00000	0.00217	0.000000	0.00002	-0.00092	0.000000
15	C12	0.00087	0.00000	0.00052	0.000000	0.00000	-0.00180	0.000000
SUM =		1.00000	1.00000	1.00000	1.00000	1.45501	-0.000167	