COMPONENT (I)		V(I) (ML/MOL)	R(I)	0(1)
DIBUTYL ETHER	C8H18O	170.44	6.0925	5.1760
2-ETHYL-1-HEXANOL	C8H180	156.38	6.1511	5.2080
QUINOLINE	CSH7N	118.18	4.7923	3.1530
ISOPROPYLBENZENE	C9H12	140.17	5.2708	4. 6440
NICCTINE	C18H14N2	168.73	6.4898	4.6210
RIPHENYL	C12H10	155.77	6.0434	4.2408

CCMPONENT (1)		V(1)	*(1)	0111	
				2.6120	_
PROPIONIC ACID	C 3H6O2	74.98	2.0769		ð
N.N-DIMETHYLFORMANIDE	C3H7NO	17.43	1.0856	2.7360	b
N-METHYL ACETAMIDE	C 3H740	76.96	3.1061	2.7320	3
1-PREPANOL	C 3H0I3	75.14	2.7709	2.5120	₽.
2-PRCPANITL	C 3H90	76.92	2.7791	2.5080	Appendix A
1.2-PROPANEDIOL	C 3H50 S	73.64	3.2824	2.7840	
GLYCEROL	C3H903	73.19	3-5857	3.0600	
LSOPROPALAMINE	Carr	46.66	2.9439	2.6700	
PETHYL VINYL RETONE	C4H69	#L.16	3.017#	2.6640	
METHACRYLIC ACID	CAHADZ	84.70	1.3197	3.0600	
VINYL ACETATE	C4H672	93.10	3.2485	2.9040	
2-BUTANONE	CAHBO	40-17	3.2479	2.8760	
BUTTRALDEHYDE	CAHAT	PR-27	1.2479	2-8760	
TETRAHYDRCFURAN	CAHRO	81.55	2.9415	2.7200	
HUTTRIC ACID	C4H802	92.43	3.5512	3.1520	
1.4-DIDXANE	C4H872	45-71	3.1854	2.6400	
ETHYL ACETATE	C4H802	98.49	3.4746	3-1100	
1-HYDROXY-2-BUTANONE	C4H802	87.57	3.5504	3.1480	
A.N-DIMETHYLACFTAMIDE	C4H94D	93.02	3.5332	2.9680	
*CRPHOL I NE	C4H9ND	87.52	3.4740	2.7960	
1-BUTANOL	C4H100 .	41.47	3.4543	3.0520	
2-BUTANOL	C4H100	92.35	3.4535	3.0480	
2-METHYL-L-PPOPANOL	C4H100	92.91	3.4535	3.0480	
TERT-BUTANOL	CAHLOD	94-88	3.4528	3-1240	
2.3-BUTANEDI OL	C4HEDO2	62.99	3.7560	3.3200	
1.2-DIMETHOXYETHANE	CAHEODZ	104.44	3.6388	3.2560	
2-ETHOXYETHANOL	C4H1002	97.41	3.5952	3.2920	

COMPONENT (1)		V(I) (ML/MOL) N(1)	0(1)
TETRAHYDROFURAN	C41180	81.55	2.9415	2.7288
BUTYRIC ACID	C4H802	92.43	3.5512	3.1528
1,4-DIOXANE	C4H802	85.71	3.1854	2.6488
ETHYL ACETATE	C4H8O2	98.49	3.4786	3.1160
3-HYDROXY-2-BUTANONE	C4H802	87.57	3.5584	3.148#
ISOBUTYRIC ACID	C4H8O2	93.44	3.5584	3.1488
ISOPROPYL FORMATE	C4H8O2	100.95	3.4911	3.1128
SULFOLANE	C4H8025	95.27	4.8358	3. 2808
N, N-DIMETHYLACETAMIDE	C4H9NO	93.02	3.5332	2.9688
MORPHOLINE	C4119NO	87.52	3.4748	2.7960
1-BUTANOL	C4H18O	91.97	3.4543	3.8528
2-BUTANGL	C4H18O	92.35	3.4535	3.0488
TERT-BUTANOL	C4H180	94.88	3.4528	3.128#
DIETHYL ETHER	C4H1#0	184.75	3.3949	3.0160
2-METHYL-1-PROPANOL	C4H18O	92.91	3.4535	3.040#
1,4-BUTANEDIOL	C4H1802	88.35	3.7576	3.328#
2,3-BUTANEDIOL	C4H1802	62.99	3.756#	3.3200
DIETHYLENE GLYCOL	C4H1003	94.78	4.8813	3.5680

ISOBUTYN ALDENYDE

BUTYLAMINE

FURFURAL

PYRIDINE

ISOPRENE

METHYL CYCLOPROPYL KETONE

METHYL METHACRYLATE

N-METHYLPYRROLIDONE

DIETHYLAMINE

C4HBO

C4H11N

C4H11H

C5H402

CSHSN

C5H8

C5H80

C 511 9NO

3.2471

91.39

99.56

184.25

83.23

88.86

100.04

186.78

3.6191

3.6836

3.1688

2.9993

3.3638

3.4681

3.9215

3.9818

3.1648

3.1720

2.4848

2.1130

3.0120

2.7960

3.564#

3.2000

2.8728

BUTYL ACETATE	7777				
	C6H12G2	124.82	4.8985	4.2080	
DIACETONE ALCOHOL	C6H12O2	83.11	5.5010	4.7320	
ISOPENTYL FORMATE	esu1202	131.85	4.8274	4.1960	
PROPYL PROPIONATE	C6H12O2	131.03	3.00011	,,,,,,,	
					9111
and the second second		V(1) (ML/	MOL) R(.)	4111
M PONENT (I)		2012-03:00 ct 0 2:0			
	C6H14	131.6	4.4	4998	3.8560
CANE		142.2	R 4.	7421	4.0880
SOPROPYL ETHER	C6H14D	141.1			
	C6H14O	125.2	3 4.	8031	4.1320

	C6H14	131.61	4.4998	3.8560
HEXANE	C6H14O	142.28	4.7421	4.0880
DIISOPROPYL ETHER		125.23	4.8031	4.1320
1-HEXANOL	C6H14O	125.24	4.8023	4.1280
2-HEXANDL	C6H14O			4.3720
2-BUTOXY-ETHANOL	C6H14O2	131.84	5.0470	
DIISOPROPYLAMINE	C6H15N	142.53	5.0308	4.2440
ETHYL BUTYL AMINE	C6H15N	130.40	5.0324	4.2520
	C6H15N	136.82	5.0316	4.2480
N-ETHYL-SEC-BUTYLAMINE	C6H15N	139.95	5.0118	4.2560
TRIETHYLAMINE	C7H602	113.60	4.3230	3.3440
BENZOIC ACID	C7H8	106.85	3,9228	2.9688
TOLUENE	C7H80	103.55	4.2260	3.2560
BENZYL ALCOHOL	C7H80	95.28	4.2867	3.2480
2-METHYLPHENOL	C7H80	104.97	4.2867	3.2480
3-WETHATBHENOF	C7HEO	93.71	4.2567	3.2480
4-METHYLPHENOL	C7H802	109.99	4.5306	3.4880
GUAIACOL	C7H9N	115.59	4.4693	3.249€
2,4-DIMETHYLPYRIDINE	C7H9N	116.48	4.4693	3. 2490
2,6-DIMETHYLPYRIDINE	C7H1402	150.26	5.5010	4.7320
ISOPENTYL ACETATE	C7H16	147.47	5.1742	4.3960
HEFTANE	СВИБ	115.57	4.3671	3.2960
STYRENE	свиво	117.36	4.6941	3.6080
ACETOPHENONE	свиза	123.07	4.5972	3.5080
ETHYLBENZENE	CSHIO	123.93	4.6578	3.5360
P-XYLENE	CEHILN	131.95	5.1437	3.7890
2-METHYL-S-ETHYLPYRIDINE	C#H16	155.97	5.6156	4.9200
2,4,4-TRIMETHYL-1-PENTENE	1			

COMPONENT (1)		V(I) (ML/MOL)	R(I)	Q(I)
2-METHYL-2-BUTENE	C5H1#	105.90	3.5919	3.2200
2-METHYL-3-BUTEN-2-OL	C5H180	86.24	3.8978	3.1760
3-PENTANONE	CSH1SO	106.41	3.9223	3.4168
ISOPROPYL ACETATE	C5H1#02	117.83	4.1522	3.6520
PENTANE	C5H12	116.11	3.8254	3.3168
3-METHYLBUTANOL	C5H12O	109.22	4.1279	3,5888
1-PENTANOL	C5H12O	100.63	4.1287	3.5920
TERT-PENTANOL	C5H12O	109.38	4.1272	3.6688
DIMETHYLISOPROPYL AMINE	C5H13N	121.96	4.3367	3.7128
ETHYL ISOPROPYL AMINE	CSH13N	124.87	4.3572	3.7060
BENZENE	C6H6	89.41	3.1875	2.4000
PHENOL	CEHEO	53.14	3.5517	2.6800
ANILINE	C6H7H	91.53	3.7165	2.8160
2-METHYLPYRIDINE	C6H7N	99.03	3.7343	2.6818
3-METHYLPYRIDINE	CEHTN	97.88	3.7343	2.6810
4-METHYLPYRIDINE	CEHTN	98.01	3.7343	2.6818
PHENYLHYDRAZINE	C 6H EN 2	98.46	4.2492	3.2240
CYCLOHEXANONE	CEH180	104.18	4.1433	3.2400
METHYLDIHYDROPYRAN	C6H18O	89.90	3.0711	2.3840
MESITYLOXIDE	C6H18O	114.23	4.3632	3.8606
CYCLOHEXANE	C6H12	108.75	4.0464	3.240
CYCLOHEXANOL	C6H12O	183.43	4.3489	3.512
BUTYL ACETATE	C6H15O5	132.55	4.8274	4.196
DIACETONE ALCOHOL	C6H12G2	124.82	4.8985	4.308
ISOPENTYL FORMATE	C6H1202	83.11	5.5010	4.732
PROPYL PROPIONATE	C6H12O2	131.85	4.8274	4.196

COMPONENT (1)		V(I) (ML/MCL)	R(1)	0(1)
ACRYLIC ACID	C3H4O2	68.56	2.6467	2.4000
3-CHLORO-1-PROPENE	C3H5CL	81.62	2.8108	2.4400
PROFIONITRILE	C 3H 5N	70.90	2.5445	2.2640
ACCITONE	C3H60	74.05	2.5735	2.3360
ALLYL ALCOHOL	C3H60	68.62	2.5498	2.3000
PROFICNIC ALDEHYDE	C3H6O	73.41	2.5735	2.3360
DIDXDLAME	C3H6O2	69.29	2.5110	2.1000
METRYL ACETATE	C3H602	79.84	2.6042	2.5760
PROFICACID	C3H6O2	74.98	2.8768	2.6120
1, 3, 5-TRICKANE	C3H6O3	76.99	2.7549	2.3488
N. N-TIMETHILFORMAMIDE	C3H7NO	77.43	3.0856	2.7360
M-METHALIZETANIDE	C3H7NO	76.96	3.1061	2.7328
1-PROFFAMOL	C 3H 8O	75.14	2.7799	2.5120
0-PROFAMOL	C 3H 60	76.92	2.7791	2.5080
DIMETHOUS ATTRANE	C3H802	89.13	2.9644	2.7160
1.1-PCFAMILIOL	C3H802	71.68	3.2524	2.7840
CLYCEROL	C3H8O3	73.19	3.5857	3.0600
ISOPROPILATIVE	C 3H 9N	\$6.66	2.9439	2.6200
HONLANDE	CIESN	83.01	2.5447	2.6240
SEEL SEELE	C4H 602	90.27	3.2485	2.9040
NINIL ACCESSIO	C48602	93.10	3.2485	2,9040
ACCOUNT ASSETS THE	C48603	94.96	3.5755	3.2160
BEEN-METER PROPRIESE ACTS STREET	C4H7NO	95.32	3.4628	3.5448
I - E LT LACAZ	C4HEC	90.17	3.2479	2.8760
EUT GALLERYES	C4HEO	88.27	3.2479	2.8760
INDERTO ALLEGADE	CEHRO	91.39	3.2471	2.8720

FURE COMPONENT PARAMETERS:				
COMPONENT (1)		VII) (ML/MOL)	R(I)	0(1)
DEUTERIUM OXIDE	D20	18.13	g.920B	1.4888
WATER CHLOROFORM	CHCL3	18.07 80.67 39.31	9.9288 2.8788 1.1957	2.4108
HYDROGEN CYANIDE DICHLOROMETHANE FORMALDEHYDE	CH2CL2 CH2O	64.5Ø 36.83	2.2564	1.9888
FORMIC ACID NITROMETHANE METHANOL	CH2O2 CH3NO2 CH4O	37.91 53.96 48.73	1.5288 2.8886 1.4311	1.5328 1.6680 1.4320
1.1.2.2-TETRACHLORO ETHANE ACETORITRILE 1.2-DICHLORGETHANE	C2H2CL4 C2H3N C2H4CL2	105.68 52.86 78.87	4.1212 1.8781 2.9388	3.3688 1.7248 2.5288
ACETALDEHYDE ETHYLENE OXIDE ACETIC ACID	C 2H 4O C 2H 4O C 2H 4O 2	56.62 49.92 57.54	1.8991 1.5927 2.2024	1.7968 1.3288 2.8728
2-CHLOROETHANOL N-HETHYLFGRMAMIDE ETHANOL	C2H5C LO C2H5NO C2H6O	67.29 59.14 58.69	2.6698 2.4317 2.1055	2.3920 2.1920 1.9720
DIMETHYLSULFOXIDE 1, 2-ETHANEDIOL DIMETHYLAMINE	C 2H 6D 5 C 2H 6D 2 C 2H 7N	71.38 55.92 66.27	2.8266 2.4886 2.3348	2.4728 2.2488 2.8928
ETHYLAMINE ACHYLAMINE	C 2H 7N C 2H 8N 2 C 3H 3N	68.75 67.84 66.30	2.2783 2.7384 2.3144	2. 0840 2. 4720 2. 0520
THIAZGLE	C 3H 3N 5 C 3H 4O	70.95 66.83	2.6688	1.853g 2.124g