

COMPONENTS:		ORIGINAL MEASUREMENTS:			
1. Methanamine, (methylanine); CH_5N ; [74-89-5] 2. Butane; C_4H_{10} ; [106-97-8]		Wolff, A.; Höpfner, A.; Höpfner, H.-M <i>Ber. Bunsenges. Phys. Chem.</i> <u>1964</u> , 68, 410-417.			
VARIABLES:		PREPARED BY:			
Composition, temperature		P. G. T. Fogg			
EXPERIMENTAL VALUES:					
Variation of the total vapor pressure/Torr with variation of temperature and of mole fraction of CH_5N in the liquid phase, $x_{\text{CH}_5\text{N}}$					
			T/K		
$x_{\text{CH}_5\text{N}}$	218.15	233.15	253.15	273.15	288.15
0	52.0	125.2	339.5	775.9	1319.4
0.0043	55.2	131.1	349.8	793.3	1348.6
0.0086	58.5	136.8	362.3	814.7	1380.5
0.0334	68.1	160.3	417.8	914.8	1523.8
0.1089	82.3	196.0	513.2	1123.1	1860.3
0.134	83.7	200.5	530.3	1167.9	1934.3
0.208	85.3	207.9	560.9	1257.5	2105.1
0.257	86.2	211.4	575.6	1305.5	2203.1
0.314	86.6	212.0	580.7	1324.4	2225.1
0.365	86.2	212.9	586.4	1342.0	2283.9
0.454	86.6	213.9	590.8	1362.4	2329.1
0.517	86.8	214.1	595.7	1369.4	2349.1
0.532	86.5	213.6	592.4	1369.5	2352.9
0.577	86.6	214.1	592.2	1372.1	2359.7
0.602	86.3	213.4	591.0	1372.0	2362.4
0.654	86.2	212.9	589.5	1370.3	2361.2
0.731	85.3	211.4	585.6	1361.7	2344.0
0.738	85.4	210.3	583.6	1356.7	2343.0
0.774	84.9	209.6	579.3	1347.2	2325.6
Cont.					
AUXILIARY INFORMATION					
METHOD/APPARATUS/PROCEDURE:		SOURCE AND PURITY OF MATERIALS:			
<p>Apparatus described previously was used (1). Liquid mixtures of accurately known composition were introduced into a cell held in a thermostat controlled to $\pm 0.02^\circ\text{C}$. The total vapor pressures were measured by a mercury manometer.</p> <p>The authors calculated activity coefficients of each component from the vapor pressure data by a method described by Barker (2). Constants for Redlich-Kister equations (3) for activity coefficients were evaluated and reported.</p>		<p>1. From commercial <i>reinst</i> methylammonium chloride by reaction with KOH; gas dried with KOH and Li; liquified gas treated with Li and repeatedly fractionated until first and last fractions had vapor pressures between 2191 and 2193 Torr at 20°C.</p> <p>2. Commercial product; dried with P_2O_5 repeatedly fractionated until first and last fractions had the same vapor pressures measured by the manometer as described for</p>			
		ESTIMATED ERROR: Hexane (1). $\delta T/\text{K} = \pm 0.02$ (estimated by authors)			
		REFERENCES: 1. Wolff, H.; Höpfner, A. <i>Z. Elektrochem.</i> <u>1962</u> , 66, 149. 2. Barker, J.A. <i>Aust. J. Chem.</i> <u>1953</u> , 6, 207. 3. Redlich, O.; Kister, A.T. <i>Ind. Eng. Chem.</i> <u>1948</u> , 21, 345.			

COMPONENTS: 1. Methanamine, (methylamine); CH ₅ N; [74-89-5] 2. Butane; C ₄ H ₁₀ ; [106-97-8]	ORIGINAL MEASUREMENTS: Wolff, A.; Höpfner, A.; Höpfner, H.-M. <i>Ber. Bunsenges. Phys. Chem.</i> <u>1964</u> , 68, 410-417.
VARIABLES: Composition, temperature	PREPARED BY: P. G. T. Fogg

EXPERIMENTAL VALUES: Cont.

	T/K				
x _{CH₅N}	218.15	233.15	253.15	273.15	288.15
0.847	82.5	202.4	561.2	1308.6	2269.0
0.938	70.6	177.4	500.8	1198.0	2102.3
0.957	67.1	169.3	482.2	1158.3	2042.7
1	45.8	126.3	394.2	1001.8	1826.3

$$760 \text{ Torr} = 1 \text{ atm} = 1.013 \times 10^5 \text{ Pa}$$

Constants for calculation of activity coefficients
from the Redlich-Kister equations given below

T/K	A	B	C
218.15	2.261	-0.054	0.412
223.15	2.213	-0.030	0.381
228.15	2.143	-0.017	0.371
233.15	2.095	-0.008	0.353
243.15	1.976	+0.008	0.313
253.15	1.857	+0.032	0.259
263.15	1.741	+0.066	0.214
273.15	1.623	+0.079	0.189
283.15	1.517	+0.087	0.157
288.15	1.461	+0.095	0.150

$$\ln f_1 = A x_2^2 - B x_2^2 (1 - 4 x_1) + C x_2^2 (1 - 8 x_1 + 12 x_1^2)$$

$$\ln f_2 = A x_1^2 + B x_1^2 (1 - 4 x_2) + C x_1^2 (1 - 8 x_2 + 12 x_2^2)$$

where f_1 = activity coefficient of methylamine
 f_2 = activity coefficient of butane
 x_1 = mole fraction of methylamine in the liquid phase
 x_2 = mole fraction of butane in the liquid phase

COMPONENTS:		ORIGINAL MEASUREMENTS:									
1. Methanamine, (methylamine); CH ₅ N; [74-89-5]		Wolff, H.; Höpfner, A. <i>Z. Elektrochem.</i> 1962, 66, 149-159.									
2. Hexane, C ₆ H ₁₄ ; [110-54-3]											
VARIABLES:		PREPARED BY:									
Composition, temperature		P. G. T. Fogg									
EXPERIMENTAL VALUES:											
Variation of the total vapor pressure/Torr with variation of temperature and of mole fraction of CH ₅ N in the liquid phase, $x_{\text{CH}_5\text{N}}$											
T/K											
$x_{\text{CH}_5\text{N}}$	218.15	223.15	228.15	233.15	243.15	253.15	263.15	273.15	283.15	293.15	
0	0.97*	1.52*	2.33*	3.48*	7.34*	14.4	26.5	46.0	76.1	121.7	
0.0068	5.9	7.8	10.1	12.6	20.8	32.9	51.4	78.7	118.5	174.9	
0.0154	11.4	14.3	18.3	23.2	36.0	54.7	80.9	118.4	169.6	239.5	
0.0274	17.3	22.6	28.6	36.1	55.9	83.1	120.3	171.2	238.5	327.6	
0.0415	22.7	30.2	38.4	48.7	75.4	112.8	162.4	228.5	314.3	424.8	
0.0626	29.0	38.5	49.8	63.7	100.1	150.7	217.3	305.9	417.7	560.3	
0.086	32.9	44.0	58.0	74.8	119.9	181.4	264.3	371.0	506.2	672.6	
0.129	37.7	51.1	68.1	89.5	147.5	228.7	338.1	482.3	663.5	887.5	
0.175	39.8	55.1	74.5	99.2	166.2	263.1	395.1	569.5	794.8	1074.1	
0.215	41.1	57.7	77.9	103.9	177.4	284.1	432.8	631.4	887.7	1208.6	
0.282	42.4	59.2	81.1	109.1	189.8	308.3	476.2	706.4	1006.9	1389.4	
0.360	43.6	61.2	83.9	113.6	199.0	327.0	510.9	767.8	1107.1	1545.1	
0.449	43.4	61.5	84.7	115.2	204.0	339.3	536.0	814.4	1188.5	1678.0	
0.543	43.5	61.2	85.4	116.7	207.4	348.6	554.7	849.9	1250.9	1781.9	
0.629	43.7	62.1	86.1	117.9	210.2	355.8	569.7	878.1	1300.4	1863.2	
0.662	43.7	62.0	86.3	117.7	210.6	356.4	574.8	885.3	1314.8	1887.6	
0.735	44.0	62.4	86.7	118.9	214.0	362.3	583.0	905.0	1348.1	1943.3	
0.783	43.8	62.1	86.8	119.2	215.0	365.6	590.1	917.2	1370.0	1982.5	
Cont.											
AUXILIARY INFORMATION											
METHOD/APPARATUS/PROCEDURE:					SOURCE AND PURITY OF MATERIALS:						
Liquid mixtures of accurately known composition were introduced into a cell held in a thermostat controlled to $\pm 0.02^\circ\text{C}$. The total vapor pressures were measured by a mercury manometer.					1. From commercial <i>reinst</i> methylammonium chloride by reaction with KOH; gas dried with KOH and Na; liquified gas treated with Li and repeatedly fractionated until first and last fractions had vapor pressures between 2191 and 2193 Torr at 20°C .						
The authors calculated activity coefficients of each component from the vapor pressure data by a method described by Barker (2). Constants for Redlich-Kister equations (3) for activity coefficients were evaluated and reported. The compositions of the vapor phase were also calculated by the authors.					2. Commercial product; dried over P ₂ O ₅ distilled and repeatedly fractionated until first and last fractions had the same vapor pressures as measured by the manometer.						
					ESTIMATED ERROR:						
					$\delta T/K = \pm 0.02$ (estimated by authors)						
					REFERENCES:						
					1. Wolff, A.; Höpfner, A.; Höpfner, H.-M. <i>Ber. Bunsenges. Phys. Chem.</i> 1964, 68, 410.						
					2. Barker, J.A. <i>Aust. J. Chem.</i> 1953, 6, 207.						
					3. Redlich, O.; Kister, A.T. <i>Ind. Eng. Chem.</i> 1948, 21, 345.						

COMPONENTS:		ORIGINAL MEASUREMENTS:									
1. Methanamine, (methylamine); CH ₅ N; [74-89-5]		Wolff, H; Höpfner, A. Z. <i>Elektrochem.</i>									
2. Hexane, C ₆ H ₁₄ ; [110-54-3]		1962, 66, 149-159.									
VARIABLES:		PREPARED BY:									
Composition, temperature		P. G. T. Fogg									
EXPERIMENTAL VALUES: Cont.											
T/K											
x _{CH₅N}	218.15	223.15	228.15	233.15	243.15	253.15	263.15	273.15	283.15	293.15	
0.827	44.0	62.6	87.4	120.1	216.1	369.1	597.0	929.0	1392.6	2017.1	
0.877	44.3	63.2	88.7	121.6	219.9	375.3	607.6	948.8	1422.4	2062.9	
0.925	44.6	63.6	89.1	122.6	222.1	380.1	617.5	964.4	1448.3	2104.5	
0.969	44.9 [‡]	64.9	90.8	125.3	227.4	389.0	631.5	985.9	1483.7	2152.8	
1	44.9 [‡]	65.4	91.9	126.3	229.8	394.2	640.5	1001.8	1506.6	2191.7	
760 Torr = 1 atm = 1.013 x 10 ⁵ Pa											
*by extrapolation											
‡ corrected value of 45.8 given in later ref. (1)											
Constants* for calculation of activity coefficients from the Redlich-Kister equations given below											
	T/K	A	B	C							
	218.15	2.392	0.122	0.527							
	223.15	2.338	0.130	0.472							
	228.15	2.266	0.138	0.395							
	233.15	2.213	0.150	0.384							
	243.15	2.087	0.163	0.300							
	253.15	1.956	0.180	0.250							
	263.15	1.833	0.198	0.212							
	273.15	1.709	0.214	0.185							
	283.15	1.588	0.224	0.164							
	293.15	1.470	0.231	0.148							
$\ln f_1 = A x_2^2 - B x_2^2(1 - 4 x_1) + C x_2^2(1 - 8 x_1 + 12 x_1^2)$ $\ln f_2 = A x_1^2 + B x_1^2(1 - 4 x_2) + C x_1^2(1 - 8 x_2 + 12 x_2^2)$											
where f_1 = activity coefficient of methylamine f_2 = activity coefficient of hexane x_1 = mole fraction of methylamine in the liquid phase x_2 = mole fraction of hexane in the liquid phase.											
* revised values given by the authors in a later paper (1).											

<p>COMPONENTS:</p> <p>1. Methanamine, (methylamine); CH_5N; [74-89-5]</p> <p>2. Nonane; C_9H_{20}; [111-84-2]</p>	<p>ORIGINAL MEASUREMENTS:</p> <p>Wolff, H.; Höpfner, A.; Höpfner, H.-M <i>Ber. Bunsenges. Phys. Chem.</i> <u>1964</u>, 68, 410-417.</p>																																																																																															
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<p>EXPERIMENTAL VALUES:</p> <p>Variation of the total vapor pressure/Torr with variation of temperature and of mole fraction of CH_5N in the liquid phase, $x_{\text{CH}_5\text{N}}$</p> <table border="1" data-bbox="131 554 1210 1098"> <thead> <tr> <th>$x_{\text{CH}_5\text{N}}$</th> <th colspan="4">T/K</th> </tr> <tr> <th></th> <th>233.15</th> <th>253.15</th> <th>273.15</th> <th>293.15</th> </tr> </thead> <tbody> <tr><td>0</td><td>0.0</td><td>0.2</td><td>0.9</td><td>3.4</td></tr> <tr><td>0.0083</td><td>10.2</td><td>20.6</td><td>37.0</td><td>61.5</td></tr> <tr><td>0.0165</td><td>19.3</td><td>39.5</td><td>71.3</td><td>116.7</td></tr> <tr><td>0.0245</td><td>27.6</td><td>56.8</td><td>103.4</td><td>169.8</td></tr> <tr><td>0.0324</td><td>35.1</td><td>73.7</td><td>134.0</td><td>221.2</td></tr> <tr><td>0.0403</td><td>41.9</td><td>89.1</td><td>163.3</td><td>270.8</td></tr> <tr><td>0.0452</td><td>45.7</td><td>97.0</td><td>176.0</td><td>286.3</td></tr> <tr><td>0.0862</td><td>72.2</td><td>162.2</td><td>308.0</td><td>515.1</td></tr> <tr><td>0.1278</td><td>88.6</td><td>212.5</td><td>420.0</td><td>725.1</td></tr> <tr><td>0.188</td><td>103.1</td><td>263.2</td><td>546.9</td><td>977.5</td></tr> <tr><td>0.251</td><td>111.3</td><td>297.9</td><td>644.4</td><td>1193.2</td></tr> <tr><td>0.311</td><td>115.9</td><td>319.7</td><td>714.2</td><td>1359.2</td></tr> <tr><td>0.381</td><td>118.9</td><td>337.1</td><td>774.3</td><td>1518.4</td></tr> <tr><td>0.440</td><td>120.2</td><td>347.0</td><td>809.9</td><td>1613.1</td></tr> <tr><td>0.533</td><td>121.2</td><td>357.1</td><td>853.4</td><td>1745.7</td></tr> <tr><td>0.588</td><td>121.4</td><td>361.5</td><td>872.4</td><td>1807.1</td></tr> <tr><td>0.692</td><td>121.6</td><td>366.3</td><td>899.5</td><td>1898.2</td></tr> </tbody> </table> <p style="text-align: right;">Cont.</p>		$x_{\text{CH}_5\text{N}}$	T/K					233.15	253.15	273.15	293.15	0	0.0	0.2	0.9	3.4	0.0083	10.2	20.6	37.0	61.5	0.0165	19.3	39.5	71.3	116.7	0.0245	27.6	56.8	103.4	169.8	0.0324	35.1	73.7	134.0	221.2	0.0403	41.9	89.1	163.3	270.8	0.0452	45.7	97.0	176.0	286.3	0.0862	72.2	162.2	308.0	515.1	0.1278	88.6	212.5	420.0	725.1	0.188	103.1	263.2	546.9	977.5	0.251	111.3	297.9	644.4	1193.2	0.311	115.9	319.7	714.2	1359.2	0.381	118.9	337.1	774.3	1518.4	0.440	120.2	347.0	809.9	1613.1	0.533	121.2	357.1	853.4	1745.7	0.588	121.4	361.5	872.4	1807.1	0.692	121.6	366.3	899.5	1898.2
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COMPONENTS: 1. Methanamine, (methylamine); CH ₅ N; [74-89-5] 2. Nonane; C ₉ H ₂₀ ; [111-84-2]	ORIGINAL MEASUREMENTS: Wolff, H.; Höpfner, A.; Höpfner, H.-M <i>Ber. Bunsenges. Phys. Chem.</i> <u>1964</u> , 68, 410-417.
VARIABLES: Composition, temperature	PREPARED BY: P. G. T. Fogg

EXPERIMENTAL VALUES: Cont.

	T/K			
$x_{\text{CH}_5\text{N}}$	233.15	253.15	273.15	293.15
0.735	122.4	368.9	909.1	1932.4
0.767	122.2	369.3	915.7	1954.2
0.938	122.3	379.2	958.7	2100.7
0.968	123.4	383.6	974.4	2129.8
1	126.3	394.2	1001.8	2191.7

$$760 \text{ Torr} = 1 \text{ atm} = 1.013 \times 10^5 \text{ Pa}$$

Constants for calculation of activity coefficients
from the Redlich-Kister equations given below

T/K	A	B	C
233.15	2.323	0.338	0.393
243.15	2.174	0.362	0.336
253.15	2.024	0.379	0.286
263.15	1.881	0.392	0.255
273.15	1.738	0.399	0.229
283.15	1.597	0.401	0.214
293.15	1.463	0.400	0.204

$$\ln f_1 = A x_2^2 - B x_2^2 (1 - 4 x_1) + C x_2^2 (1 - 8 x_1 + 12 x_1^2)$$

$$\ln f_2 = A x_1^2 + B x_1^2 (1 - 4 x_2) + C x_1^2 (1 - 8 x_2 + 12 x_2^2)$$

where f_1 = activity coefficient of methylamine

f_2 = activity coefficient of nonane

x_1 = mole fraction of methylamine in the liquid phase

x_2 = mole fraction of nonane in the liquid phase.

COMPONENTS:			ORIGINAL MEASUREMENTS:
1. Methanamine, (methylanine); CH_5N ; [74-89-5] 2. Decane; $\text{C}_{10}\text{H}_{22}$; [124-18-5]			Gerrard, W. <i>Solubility of Gases and Liquids,</i> <i>Plenum, 1976, Chapter 10.</i>
VARIABLES:			PREPARED BY:
Temperature, pressure			C. L. Young
EXPERIMENTAL VALUES:			
T/K	P/mmHg	P/ 10^5 Pa	Mole fraction of methylanine in liquid, $x_{\text{CH}_3\text{NH}_2}$
273.15	100	0.133	0.028
	200	0.267	0.061
	300	0.400	0.097
	400	0.533	0.141
	500	0.667	0.193
	600	0.800	0.262
	700	0.933	0.350
278.15	760	1.013	0.413
	100	0.133	0.022
	200	0.267	0.047
	300	0.400	0.075
	400	0.533	0.109
	500	0.667	0.144
	600	0.800	0.190
283.15	700	0.933	0.256
	760	1.013	0.318
	100	0.133	0.020
	200	0.267	0.044
	300	0.400	0.070
	400	0.533	0.100
	500	0.667	0.132
	600	0.800	0.168
	700	0.933	0.206
	760	1.013	0.253
AUXILIARY INFORMATION			
METHOD/APPARATUS/PROCEDURE:		SOURCE AND PURITY OF MATERIALS:	
<p>Amine was passed into a known weight pf pure liquid in a bubbler tube at a total pressure measured by a manometer assembly. The amount of absorbed gas was estimated by weighing. The temperature was manually controlled to within 0.2K. The apparatus and procedure are described by Gerrard [1,2].</p>		<p>1. British Drug Houses or Cambrian Gases sample.</p> <p>2. Purified and attested by conventional procedures.</p>	
		ESTIMATED ERROR:	
		$\delta T/K = \pm 0.1$; $\delta x/x = \pm 3\%$ (estimated by compiler)	
		REFERENCES:	
		<p>1. Gerrard, W. <i>J. Appl. Chem. Biotechnol.</i> <u>1972</u>, 22 623-650.</p> <p>2. Gerrard, W. <i>Solubility of Gases and Liquids.</i> Plenum Press, New York. <u>1976</u>. Chapter 1.</p>	

<p>COMPONENTS:</p> <p>1. Methanamine (methylamine); CH₅N; [74-89-5]</p> <p>2. Decane; C₁₀H₂₂; [124-18-5]</p>	<p>ORIGINAL MEASUREMENTS:</p> <p>Gerrard, W. <i>Solubility of Gases and Liquids</i> Plenum, <u>1976</u>, Chapter 10.</p>
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EXPERIMENTAL VALUES:

T/K	P/mmHg	P/10 ⁵ Pa	Mole fraction of methylamine in liquid, CH ₃ NH ₂
293.15	100	0.133	0.011
	200	0.267	0.023
	300	0.400	0.037
	400	0.533	0.055
	500	0.667	0.078
	600	0.800	0.103
	700	0.933	0.134
	760	1.013	0.156

COMPONENTS:			ORIGINAL MEASUREMENTS:
1. Methanamine; (methylanine); CH_5N ; [74-89-5] 2. Aromatic hydrocarbons			Gerrard, W. <i>Solubility of Gases and Liquids</i> , Plenum, <u>1976</u> , Chapter 10.
VARIABLES:			PREPARED BY:
Pressure			C. L. Young
EXPERIMENTAL VALUES:			Mole fraction of methylanine in liquid, $x_{\text{CH}_3\text{NH}_2}$
T/K	P/mmHg	P/ 10^5Pa	
		Benzene; C_6H_6 ; [71-43-2]	
283.15	760	1.013	0.408
		Methylbenzene (toluene); C_7H_8 ; [108-88-3]	
283.15	760	1.013	0.393
		1,3-Dimethylbenzene (<i>m</i> -xylene); C_8H_{10} ; [108-38-3]	
283.15	100	0.133	0.027
	200	0.267	0.064
	300	0.400	0.100
	400	0.533	0.145
	500	0.667	0.196
	600	0.800	0.275
	700	0.933	0.320
	760	1.013	0.358
		1,3,5-Trimethylbenzene (mesitylene); C_9H_{12} ; [108-67-8]	
283.15	100	0.133	0.032
	200	0.267	0.068
	300	0.400	0.109
	400	0.533	0.155
	500	0.667	0.200
	600	0.800	0.253
	700	0.933	0.310
	760	1.013	0.326
AUXILIARY INFORMATION			
METHOD/APPARATUS/PROCEDURE:		SOURCE AND PURITY OF MATERIALS:	
<p>Amine was passed into a known weight of pure liquid in a bubbler tube at a total pressure measured by a manometer assembly. The amount of absorbed gas was estimated by weighing. The temperature was manually controlled to within 0.2K. The apparatus and procedure are described by Gerrard [1,2].</p>		<p>1. British Drug Houses or Cambrian Gases sample.</p> <p>2. Purified and attested by conventional procedures.</p>	
		ESTIMATED ERROR: $\delta T/K = \pm 0.1$; $\delta x/x = \pm 3\%$ (estimated by compiler)	
		REFERENCES: 1. Gerrard, W. <i>J. Appl. Chem. Biotechnol.</i> <u>1972</u> , 22 623-650. 2. Gerrard, W. <i>Solubility of Gases and Liquids</i> . Plenum Press, New York. <u>1976</u> . Chapter 1.	