Section B Basic Constants II.

No.	Formula	Name	CAS #	DelHf0, kJ/mol	DelGf0, kJ/mol	DelHb, kJ/mol	DelHm, kJ/mol	V liq, cm <sup>3</sup> /mol	T lig, K	Dipole, Debye
1	Ar	argon	7440-37-1	0.00	0.00	6.43		29.10	90.00	0.0
2	Br <sub>2</sub>	bromine	7726-95-6	30.91	3.13	58.80		$\frac{25.10}{51.51}$	298.15	0.0
3	BrD	deuterium bromide	13536-59-9	-37.12	-53.79	20.00		51.51	250.15	0.0
4	BrF <sub>3</sub>	bromine trifluoride	7787-71-5	-255.64	-229.51	47.57		48.84	298.15	1.1
5	BrF <sub>5</sub>	bromine pentafluoride	7789-30-2	$\frac{-428.86}{}$	$\frac{-351.65}{-351.65}$	30.60		71.02	298.15	$\frac{1.1}{1.5}$ $\frac{0.8}{0.8}$
6	BrH	hydrogen bromide	10035-10-6	-36.26	$\frac{-53.30}{-53.30}$	20.00		35.85	193.15	$\frac{1.8}{0.8}$
7	CBrClF <sub>2</sub>	bromochlorodifluoromethane	353-59-3	-435.20	-411.80			91.67	298.15	
8	CBrF <sub>3</sub>	bromotrifluoromethane	75-63-8	-649.80	-622.90			74.42	213.15	0.7
9	CBr <sub>2</sub> F <sub>2</sub>	dibromodifluoromethane	75-61-6	-386.60	-375.70			93.04	298.15	$\begin{array}{c} 0.7 \\ \hline 0.5 \\ \hline 0.5 \\ \hline 0.5 \\ \hline 0.0 \\ \hline 0.0 \\ \hline 0.0 \\ \hline 1.5 \\ \hline 1.4 \\ \hline 1.3 \\ \hline 1.1 \\ \end{array}$
10	CClF <sub>3</sub>	chlorotrifluoromethane	75-72-9	-704.20	-663.60	15.73		68.87	193.15	0.5
11	$CCl_2F_2$	dichlorodifluoromethane (R-12)	75-71-8	-490.80	-451.70	20.08		82.91	243.15	0.5
12	CCl <sub>3</sub> F	trichlorofluoromethane (R-11)	75-69-4	-283.70	-244.40	25.06	6.89	93.82	298.15	$\overline{0.5}$
13	CCl <sub>4</sub>	tetrachloromethane	56-23-5	-95.81	-53.53	29.82	3.28	97.07	298.15	$\overline{0.0}$
14	CF₄ .	tetrafluoromethane	75-73-0	-933.50	-888.80			56.41	153.15	$\overline{0.0}$
15	CHBrF <sub>2</sub>	bromodifluoromethane	1511-62-2	-429.50	-412.90			64.81	253.15	1.5
16	CHClF <sub>2</sub>	chlorodifluoromethane (R-22)	75-45-6	-482.80	-451.70	20.22	4.12	59.08	213.15	1.4
17	CHCl <sub>2</sub> F	dichlorofluoromethane	75-43-4	-284.90	-254.40	25.13		75.96	298.15	1.3
18	CHCl <sub>3</sub>	trichloromethane	67-66-3	-102.93	-70.09	29.24	8.80	80.68	298.15	
19	CHF <sub>3</sub>	trifluoromethane (R-23)	75-46-7	-693.30	-658.80			51.66	213.15	1.6
20	$CH_2Cl_2$	dichloromethane	75-09-2	-95.40	-68.84	28.06	4.60	64.53	298.15	$ \begin{array}{r}     \hline     1.6 \\     \hline     1.8 \\     \hline     2.0 \\     \hline     1.5 \\     \hline     1.9 \\     \hline     1.8 \end{array} $
21	$CH_2F_2$	difluoromethane	75-10-5	-452.30	-424.70			42.91	223.15	2.0
22	$CH_2O_2$	methanoic acid (formic acid)	64-18-6	-378.60	-35.06	22.69	12.72			1.5
23	CH <sub>3</sub> Cl	chloromethane	74-87-3	-81.96	-58.42	21.58	6.43	50.59	253.15	1.9
24	CH <sub>3</sub> F	fluoromethane	593-53-3	-237.70	-213.70			40.60	213.15	1.8
25	$CH_3NO_2$	nitromethane	75-52-5	-74.70	-6.90	33.99		53.96	298.15	3.1
26	$CH_4$	methane	74-82-8	-74.52	-50.45	8.17	0.94	35.54	90.68	$\frac{0.0}{1.7}$
27	$CH_4O$	methanol	67-56-1	-200.94	-162.24	35.21	3.18	40.73	298.15	1.7
28	$CH_4S$	methanethiol (methyl mercaptan)	74-93-1	-22.59	-9.52	_	5.91	55.52	293.15	1.3
29	CH₅N	methanamine (methyl amine)	74-89-5	-22.53	32.73	25.60	6.13	47.34	298.15	$\frac{\overline{1.3}}{\underline{1.3}}$

3		carbon monoxide	630-08-0	-110.53	-137.16	6.04	0.84	34.88	81.00	0.1
3	$CO_2$	carbon dioxide	124-38-9	-393.51	-394.38		9.02	·		0.0
3	$C_2Br_2ClF_3$	1,2-dibromo-2-chloro-1,1,2-trifluroethane	354-51-8			31.17		123.67	298.15	· <u></u>
3	$C_2ClF_5$	1-chloro-1,1,2,2,2-pentafluoroethane	76-15-3	-1123.00	-1042.00	19.22	1.88	98.79	233.15	0.3
3	$C_2Cl_2F_4$	1,1-dichloro-1,2,2,2-tetrafluoroethane	374-07-2	-926.80	-845.40	23.08		116.41	298.15	
3.	$C_2Cl_2F_4$	1,2-dichloro-1,1,2,2-tetrafluoroethane	76-14-2	-900.40	-818.10	22.94	1.51	117.37	298.15	0.5
3	$C_2Cl_3F_3$	1,1,2-trichloro-1,2,2-trifluoroethane	76-13-1	-705.80	-627.30	28.08	2.47	119.78	293.15	_
3	$C_2F_4$	tetrafluoroethene	116-14-3	-659.00	-624.10			65.84	197.00	0.0
3		hexafluoroethane	76-16-4	-1343.00	-1257.00	16.15				$\overline{0.0}$
3	$C_2HBrClF_3$	1-bromo-1-chloro-2,2,2-trifluoroethane	151-67-7			28.08		106.35	298.15	
4	$C_2HBrClF_3$	1-bromo-2-chloro-1,1,2-trifluoroethane	354-06-3			28.31		106.29	298.15	
4	C <sub>2</sub> HClF <sub>4</sub>	1-chloro-1,1,2,2-tetrafluoroethane	354-25-6	-903.30	-830.90	22.30		105.06	273.15	
4	C <sub>2</sub> HClF <sub>4</sub>	1-chloro-1,2,2,2-tetrafluoroethane	2837-89-0	-924.70	-851.80	22.53		99.88	295.00	
		(R-124)								
4	C <sub>2</sub> HCl <sub>2</sub> F <sub>3</sub>	1,1-dichloro-2,2,2-trifluoroethane (R-123)	306-83-2	-743.90	-668.90	25.72		103.90	295.00	
4	$C_2HCl_2F_3$	1,2-dichloro-1,2,2-trifluoroethane	354-23-4					103.72	295	
		(R-123a)								
4	$C_2HF_5$	pentafluoroethane	354-33-6	-1105.00	-1030.00			98.61	293.48	1.5
4	$C_2HF_5O$	pentafluorodimethyl ether (E-125)	3822-68-2					105.66	295.00	
4	$C_2H_2$	ethyne (acetylene)	74-86-2	190.92	201.30		21.28	43.47	203.15	0.0
4	$C_2H_2F_2$	1,1-difluoroethene	75-38-7	-336.81	-313.06			103.78	297.00	1.4
4	$C_2H_2F_4$	1,1,1,2-tetrafluoroethane (R-134a)	811-97-2	-907.10	-838.40			75.38	255.00	
5	$C_2H_2F_4$	1,1,2,2-tetrafluoroethane (R-134)	359-35-3	-892.40	-824.60			74.42	275.00	0.0
5	$C_2H_3ClF_2$	1-chloro-1,1-difluoroethane	75-68-3	-529.70	-465.70	22.38	2.69	90.49	298.15	2.1
5	$C_2H_3Cl_2F$	1,1-dichloro-1-fluoroethane (R-141b)	1717-00-6	-339.70	-276.20	25.94		93.63	290.00	_
5	$C_2H_3F_3$	1,1,1-trifluoroethane (R-143a)	420-46-2	-745.60	-678.70	18.99	6.19	75.38	245.00	2.3
5	$C_2H_3F_3$	1,1,2-trifluoroethane (R-143)	430-66-0	-669.40	-609.40			69.22	300.00	_
5.		ethene (ethylene)	74-85-1	52.50	68.48	13.53	3.35	51.07	183.15	0.0
5	$C_2H_4Br_2$	1,2-dibromoethane	106-93-4	-38.94	-1.06	34.77		86.62	298.15	1.0
5	$C_2H_4Cl_2$	1,1-dichloroethane	75-34-3	-130.12	-73.23	28.85	7.87	84.73	298.15	2.0
5	$C_2H_4Cl_2$	1,2-dichloroethane	107-06-2	-126.78	-70.20	31.98	8.84	79.45	298.15	$ \begin{array}{r}     \hline     \underline{2.0} \\     \hline     \underline{1.8} \\     \underline{2.3} \\     \overline{1.3} \end{array} $
5	$C_2H_4F_2$	1,1-difluoroethane (R-152a)	75-37-6	-500.80	-443.30	21.56		65.76	252.47	2.3
6	$C_2H_4O_2$	ethanoic acid (acetic acid)	64-19-7	-432.25	-374.27	23.70	11.72	57.53	298.15	1.3
6		methyl methanoate (methyl formate)	107-31-3	-352.40	-294.90	27.92		62.14	298.15	1.8
6	$C_2H_5Br$	bromoethane	74-96-4	-63.60	-25.70	27.04		75.12	298.15	$\frac{2.0}{2.0}$
6	$C_2H_5Cl$	chloroethane	75-00-3	-112.26	-60.43	24.53	4.45	72.58	298.15	$\overline{2.0}$

No.	Formula	Name	CAS #	DelHf0, kJ/mol	DelGf0, kJ/mol	DelHb, kJ/mol	DelHm, kJ/mol	V liq, cm³/mol	T liq, K	Dipole, Debye
64	C <sub>2</sub> H <sub>5</sub> F	fluoroethane	353-36-6	-264.40	-212.40			56.67	213.15	2.0
65	$C_2H_6$	ethane	74-84-0	-83.82	-31.86	14.70	2.86	46.15	90.36	$\frac{0.0}{1.7}$
66	$C_2H_6O$	ethanol	64-17-5	-234.95	-167.73	38.56	5.01	58.68	298.15	1.7
67	$C_2H_6O$	dimethyl ether	115-10-6	-184.11	-112.92	21.51	4.94	69.07	293.15	$ \begin{array}{r} 1.3 \\ 1.6 \\ 1.5 \\ 1.3 \\ 1.0 \\ 0.0 \end{array} $
68	$C_2H_6S$	ethanethiol (ethyl mercaptan)	75-08-1	-46.02	-2.07	26.79	4.98	74.58	298.15	1.6
69	$C_2H_6S$	2-thiapropane (dimethylsulfide)	75-18-3	-37.20	7.41	27.00	7.98	73.77	298.15	1.5
70	$C_2H_7N$	ethanamine (ethyl amine)	75-04-7	-47.47	36.28			66.59	298.15	1.3
71	$C_2H_7N$	<i>N</i> -methylmethanamine (dimethyl amine)	124-40-3	-18.80	68.80	26.40	5.94	68.73	293.15	1.0
72	$C_3F_8$	octafluoropropane (R-218)	76-19-7	-1737.62	-1610.31			140.00	295.00	0.0
73	$C_3HF_7$	1,1,1,2,3,3,3-heptafluoropropane (R-227ea)	431-89-0					121.40	295.00	
74	C <sub>3</sub> H <sub>2</sub> ClF <sub>5</sub> O	2-chloro-1,1,2-trifluoroethyl	13838-16-9			30.09				
	3 2 3	difluoromethyl ether (enflurane)								
78	$C_3H_3F_5$	1,1,1,2,2-pentafluoropropane (R-245cb)	1814-88-6					112.82	295.00	
79	$C_3H_3F_5$	1,1,1,3,3-pentafluoropropane (R-245fa)	460-73-1					99.50	295.00	
81	$C_3H_3F_5O$	2-(difluoromethoxy)-1,1,1-trifluoroethane (E-245)	1885-48-9					107.71	295.00	
82	$C_3H_3NO$	1,2-oxazole(isoxazole)	288-14-2	82.02		37.17		51.408	298.15	
83	$C_3H_4$	1-propyne (methyl acetylene)	74-99-7	$1\overline{10.00}$	150.60	22.15		62.31	273.15	0.7
84	$C_3H_4$	1,2-propadiene	463-49-0	190.92	201.30	20.60	4.40	60.95	243.15	$\overline{0.2}$
85	$C_3H_6$	propene (propylene)	115-07-1	20.00	62.50	18.42	3.00			$\overline{0.4}$
86	$C_3H_6$	cyclopropane	75-19-4	53.30	104.40	20.05	5.44	71.76	233.15	$\begin{array}{c} 0.7 \\ \hline 0.2 \\ \hline 0.4 \\ \hline 0.0 \end{array}$
87	$C_3H_6Cl_2$	1,2-dichloropropane	78-87-5			36.17		100.815	323.14	
88	$C_3H_6O$	2-propen-1-ol (allyl alcohol)	107-18-6	-123.60		47.30				
89	$C_3H_6O$	propanone (acetone)	67-64-1	-217.10	-152.60	29.10	5.69	73.94	298.15	2.9

90	$C_3H_6O_2$	propanoic acid	79-09-4	-452.80	-369.60	31.14	10.66	74.97	298.15	$\frac{1.5}{1.7}$ $\frac{2.0}{2.0}$
91	$C_3H_6O_2$	methyl ethanoate (methyl acetate)	79-20-9	-408.80	-321.40	30.32	7.49	79.89	298.15	1.7
92	$C_3H_6O_2$	ethyl methanoate (ethyl formate)	109-94-4	-388.30	-303.00	29.91		80.93	298.15	<u>2.0</u>
93	$C_3H_6O_3$	dimethylcarbonate	616-38-6	-571		37.7		84.82	298.15	
94	C <sub>3</sub> H <sub>7</sub> Cl	1-chloropropane	540-54-5	-133.18	-52.51	27.18		88.95	298.15	2.0
95	$C_3H_8$	propane	74-98-6	-104.68	-24.29	19.04	3.53	74.87	233.15	$\overline{0.0}$
96	$C_3H_8O$	1-propanol	71-23-8	-255.20	-159.81	41.44	5.20	75.14	298.15	1.7
97	$C_3H_8O$	2-propanol	67-63-0	-272.70	-173.32	39.85	5.38	76.92	298.15	1.7
98	$C_3H_8O$	methyl ethyl ether	540-67-0	-216.50	-117.13	25.69		85.85	293.15	$   \begin{array}{r}     \hline       0.0 \\       \hline       1.7 \\       \hline       1.7 \\       \hline       1.2   \end{array} $
99	$C_3H_8S$	2-thiabutane (methyl ethyl sulfide)	624-89-5	-59.29	11.92	29.53	9.76	91.02	298.15	
100	$C_3H_9N$	1-propanamine (propyl amine)	107-10-8	-70.10	41.90	29.55		83.11	298.15	1.3
101	$C_3H_9N$	2-propanamine (methyl ethyl amine)	75-31-0	-83.70	32.23	27.83		86.33	298.15	
102	$C_3H_9N$	N, N-dimethylmethanamine (trimethyl	75-50-3	-23.60	99.30	22.94	6.55	94.28	298.15	0.6
	3 ,	amine)								_
103	C <sub>3</sub> H <sub>o</sub> NO	2-ethoxymethanamine	109-83-1			<u>57</u>		82.21	323.14	
	3 ,	(2-methylaminoethanol)				_				
104	$C_4F_8$	octafluorocyclobutane	115-25-3					120.94	253.15	0.0
105	$C_4F_{10}$	decafluoro-2-methylpropane	354-92-7					156.91	293.15	
106	$C_4H_4O$	furan	110-00-9	-34.73	0.94	27.10	3.80	73.09	298.15	0.7
107	$C_4H_4S$	thiophene	110-02-1	114.90	126.10	31.48	4.97	79.47	298.15	$\frac{0.7}{0.5}$ $\frac{1.8}{1.8}$
108	$C_4H_5N$	pyrrole	109-97-7	108.20	160.30	38.75	7.91	69.38	294.00	1.8
109	$C_4H_6$	1-butyne	107-00-6	162.26	198.77	24.52	6.03	83.22	298.15	$\overline{0.8}$
110	$C_4H_6$	1,3-butadiene	106-99-0	110.00	150.60	22.47	7.98	88.04	298.15	$\overline{0.0}$
111	$C_4H_6O_3$	acetic anhydride	108-24-7	-575.70	-476.80					$\frac{0.0}{3.0}$
112	$C_4H_8$	cyclobutane	287-23-0	28.40	112.20	24.19	1.09	81.43	298.15	0.0
113	$C_4H_8$	1-butene	106-98-9	-0.54	70.37	22.07	3.96	95.34	298.15	$\overline{0.3}$
114	$C_4H_8$	trans-2-butene	624-64-6	-11.00	63.34	23.34	9.76	93.65	298.15	$\overline{0.0}$
115	$C_4H_8$	cis-2-butene	590-18-1	-7.40	65.46	22.72	7.58	91.01	298.15	$\overline{0.3}$
116	$C_4H_8$	2-methylpropene	115-11-7	-17.10	58.18	21.53	5.93	95.24	298.15	$   \begin{array}{r}     \hline     0.3 \\     \hline     0.5 \\     \hline     3.3 \\     \hline     1.7 \\     \hline     1.5   \end{array} $
117	$C_4^{\stackrel{1}{}}H_8^{\stackrel{2}{}}O$	butanone (methyl ethyl ketone)	78-93-3	-238.60	-146.50	31.30	8.44	90.13	298.15	3.3
118	$C_4H_8O$	tetrahydrofuran	109-99-9	-184.18	-79.57	29.81	8.54	81.71	298.15	1.7
119	$C_4H_8O_2$	butanoic acid	107-92-6	-473.60		40.45	11.08	92.46	298.15	1.5
120	$C_4H_8O_2$	2-methylpropanoic acid	79-31-2	-484.20				93.42	298.15	1.3
121	$C_4H_8O_2$ $C_4H_8O_2$	1,3-dioxane	505-22-6	$\frac{-337.30}{-337.30}$	-203.90	34.37		,	2,0.10	1.0
122	$C_4H_8O_2$ $C_4H_8O_2$	1,4-dioxane	123-91-1	-314.70	-180.20	34.16	12.85	85.29	293.15	0.0
	C4-18O2	1,1 010110110	123 /1 1	211.70	100.20	20	12.00	00.27	2,5.15	0.0

Section B Basic Constants II (Continued)

	Б 1	N.	CAS II	DelHf0,	DelGf0,	DelHb,	DelHm,	V liq,	m 1; 17	Dipole,
No.	Formula	Name	CAS #	kJ/mol	kJ/mol	kJ/mol	kJ/mol	cm <sup>3</sup> /mol	T liq, K	Debye
123	$C_4H_8O_2$	methyl propanoate	554-12-1	-427.50	-310.90	32.24		96.93	298.15	1.7
124	$C_4H_8O_2$	ethyl ethanoate (ethyl acetate)	141-78-6	-444.50	-328.00	31.94	10.48	98.55	298.15	1.9
125	$C_4H_8O_2$	propyl methanoate (propyl formate)	110-74-7	-407.60	-293.50	33.61		97.94	298.15	1.9
126	C <sub>4</sub> H <sub>9</sub> Cl	2-chlorobutane	78-86-4	-165.69	-55.06	29.17		106.76	298.15	1.9 1.9 2.1 0.0
127	$C_4H_{10}$	butane	106-97-8	-125.79	-16.57	22.44	4.66	100.48	298.15	0.0
128	$C_4H_{10}$	2-methylpropane	75-28-5	-134.99	-21.44	21.30	4.61	104.36	298.15	0.1
129	$C_4H_{10}N_2$	piperazine	110-85-0	31.6		26.7				
130	$C_4H_{10}O$	1-butanol	71-36-3	$-2\overline{74.60}$	-150.17	43.29	9.28	91.96	298.15	1.8
131	$C_4H_{10}O$	2-methyl-1-propanol (isobutanol)	78-83-1	-282.90	-167.40	41.82		92.91	298.15	1.7
132	$C_4H_{10}O$	2-methyl-2-propanol (tert-butanol)	75-65-0	-325.81	-191.20	39.07	6.79	94.88	298.15	1.7
133	$C_4H_{10}O$	2-butanol (sec-butanol)	78-92-2	-292.75	-167.71	40.75		92.35	298.15	$   \begin{array}{r}     \underline{1.8} \\     \underline{1.7} \\     \underline{1.7} \\     \underline{1.7} \\     \underline{1.7} \\     \underline{1.3}   \end{array} $
134	$C_4H_{10}O$	diethyl ether	60-29-7	-250.80	-120.70	26.52	7.27	104.75	298.15	1.3
135	$C_4H_{10}O_2$	1,2-dimethoxyethane	110-71-4	-342.80		36.69		104.56	298.15	
136	$C_4H_{10}O_2$	1,2-butandiol	26171-83-5			71.55		92.06	323.14	
137	$C_4H_{10}O_2$	1,3-butanediol	107-88-0			74.46		138.54	323.14	
138	$C_4H_{10}S$	3-thiapentane (diethyl sulfide)	352-93-2	-83.22	18.24	31.77	11.90	108.51	298.15	<u>1.6</u>
139	$C_4H_{11}N$	1-butanamine (butyl amine)	109-73-9	-91.76	49.03	31.81		98.97	298.15	1.3
140	$C_4H_{11}N$	<i>N</i> -ethylethanamine (diethyl amine)	109-89-7	-71.70	73.00	29.06		104.24	298.15	<u>1.1</u>
141	$C_4H_{11}N$	2-methyl-1-propanamine (isobutyl amine)	78-81-9	-98.55		30.61		100.45	298.15	1.3 1.1 1.2
142	$C_5F_{12}$	dodecafluoropentane	678-26-2	-2561.86	-2351.67			177.80	293.15	0.0
143	$C_5H_5N$	pyridine	110-86-1	140.37	190.55	35.09	8.28	80.88	298.15	$\frac{2.3}{0.7}$
144	$C_5H_6O$	2-methylfuran	534-22-5	-66.32	-2.36	28.35	8.55	89.73	293.15	0.7
145	$C_5H_8$	1-pentyne	627-19-0	143.10	209.20	26.86		98.91	298.15	$   \begin{array}{r}     \hline       0.9 \\       \hline       0.9 \\       \hline       3.0 \\       \hline       0.0 \\       \hline       0.4 \\   \end{array} $
146	$C_5H_8$	cyclopentene	142-29-0	32.95	110.90	26.76	3.36	91.43	293.15	$\overline{0.9}$
147	$C_5H_8O$	cyclopentanone	120-92-3	$-1\overline{92.80}$	' <u></u> '	36.35		<u></u> -		3.0
148	$C_5H_{10}$	cyclopentane	287-92-3	-77.10	38.92	27.30	0.61	94.73	298.15	$\overline{0.0}$
149	$C_5H_{10}$	1-pentene	109-67-1	-21.30	78.60	25.20	5.81	110.40	298.15	$\overline{0.4}$
150	$C_5H_{10}$	cis-2-pentene	627-20-3	-26.30	73.50	26.06	7.11	107.85	298.15	
151	$C_5H_{10}$	2-methyl-2-butene	513-35-9	-40.80	61.60	26.31	7.60	107.57	298.15	

152 154 155	$C_5H_{10}$ $C_5H_{10}O$ $C_5H_{10}O$	3-methyl-1-butene 2-pentanone (methyl propyl ketone) 3-pentanone (diethyl ketone)	563-45-1 107-87-9 96-22-0	-27.60 $-259.20$ $-257.90$	76.00 -138.20 -134.30	23.94 33.44 33.45	5.36 10.63 11.59	112.77 107.33 106.41	298.15 298.15 298.15	$\frac{2.5}{2.7}$ $\frac{2.8}{2.8}$
156	$C_5H_{10}O$	3-methyl-2-butanone (methyl isopropyl ketone)	563-80-4	-262.60	-139.30	32.35	9.34	107.91	298.15	2.8
157	$C_5H_{10}O$	2-methyltetrahydrofuran	96-47-9			30.13		100.39	298.15	
158	$C_5H_{10}O_2$	pentanoic acid	109-52-4	-497.00	-351.00	62.40	14.16	109.30	298.15	
159	$C_5H_{10}O_2$	3-methylbutanoic acid	503-74-2	.,,,,,,	221.00	020	110	110.79	298.15	1.0
160	$C_5H_{10}O_2$	methyl butanoate	623-42-7	-450.70	-305.10	33.79		114.42	298.15	1.7
161	$C_5H_{10}O_2$	ethyl propanoate	105-37-3	-463.60	-319.10	33.88		115.54	298.15	$ \begin{array}{r} 1.0 \\ \underline{1.7} \\ \underline{1.8} \\ \underline{2.0} \\ 1.8 \end{array} $
162	$C_5H_{10}O_2$	methyl 2-methylpropanoate	547-63-7	-464.00	-317.00	32.61		114.63	293.15	$\overline{2.0}$
163	$C_5H_{10}O_2$	propyl ethanoate (propyl acetate)	109-60-4	-464.80	-320.20	33.92		115.72	298.15	1.8
164	$C_5H_{10}O_2$	2-methylpropyl methanoate (isobutyl	542-55-2	-436.30	-293.00	33.60		116.96	298.15	1.9
		formate)								
165	$C_5H_{11}Cl$	1-chloropentane	543-59-9	-175.02	-36.68	33.15		121.53	298.15	$\frac{2.2}{0.0}$
166	$C_5H_{12}$	pentane	109-66-0	-146.76	-8.65	25.79	8.40	115.22	298.15	0.0
167	$C_5H_{12}$	2-methylbutane	78-78-4	-153.70	-13.86	24.69	5.16	116.46	298.15	0.1
168	$C_5H_{12}$	2,2-dimethylpropane (neopentane)	463-82-1	-168.11	-15.24	22.74	3.26	122.16	298.15	0.0
169	$C_5H_{12}O$	1-pentanol	71-41-0	-295.60	-142.20	44.40	9.38	108.63	298.15	1.7
170	$C_5H_{12}O$	2-pentanol	6032-29-7	-313.80		41.40				
171	$C_5H_{12}O$	2-methyl-1-butanol	137-32-6	-302.00		55.18				
172	$C_5H_{12}O$	2-methyl-2-butanol	75-85-4	-329.00	-166.00	40.00	4.46	109.50	298.15	1.9 1.8
173	$C_5H_{12}O$	3-methyl-1-butanol	123-51-3	-300.00		44.10		109.22	298.15	1.8
174	$C_5H_{12}O$	3-methyl-2-butanol	598-75-4	-316.40		53.00				
175	$C_5H_{12}O$	ethyl propyl ether	628-32-0	-272.20	-115.00	28.94	8.40	121.35	298.15	1.2
176	$C_5H_{12}S$	3-methyl-1-butanethiol (isopentyl	541-31-1	-114.60	17.70		7.45	125.32	298.15	
		mercaptan)								
177	$C_6ClF_5$	chloropentafluorobenzene	344-07-0			34.76				
178	$C_6F_6$	hexafluorobenzene	392-56-3			31.66				0.0
179	$C_6F_{12}$	dodecafluorocyclohexane	355-68-0							0.0
180	$C_6F_{14}$	tetradecafluorohexane	355-42-0	-2973.99	-2722.34			198.91	293.15	$\frac{0.0}{0.0}$
181	$C_6F_{14}$	tetradecafluoro-2-methylpentane	355-04-4					195.06	293.15	

Section B Basic Constants II (Continued)

No.	Formula	Name	CAS #	DelHf0, kJ/mol	DelGf0, kJ/mol	DelHb, kJ/mol	DelHm, kJ/mol	V liq, cm <sup>3</sup> /mol	T liq, K	Dipole, Debye
184	C <sub>6</sub> HF <sub>5</sub>	pentafluorobenzene	363-72-4			32.15	10.85			
186	C <sub>6</sub> H <sub>5</sub> Cl	chlorobenzene	108-90-7	51.09	98.36	35.19	9.61	102.22	298.15	1.6
187	$C_6H_6$	benzene	71-43-2	82.88	129.75	30.72	9.95	89.41	298.15	$\overline{0.0}$
188	$C_6H_6O$	phenol	108-95-2	-96.40	-32.55	46.18	$1\overline{1.29}$	87.87	298.15	1.6
189	$C_6H_7N$	benzeneamine (aniline)	62-53-3	87.45	167.90	42.44	10.56	91.52	298.15	1.6
190	$C_6H_7N$	2-methylpyridine (2-picoline)	109-06-8	99.16	177.37	36.17	9.72	99.11	298.15	$\frac{1.9}{2.4}$
191	$C_6H_7N$	3-methylpyridine (3-picoline)	108-99-6	106.36	184.62	37.35	14.18	97.81	298.15	$\overline{2.4}$
192	$C_6H_7N$	4-methylpyridine (4-picoline)	108-89-4	102.13	182.08	37.51	12.58	98.02	298.15	
193	$C_6H_8O$	2-cyclohexen-1-one	930-68-7			48.36		99.59	323.14	
194	$C_{6}H_{10}$	cyclohexene	110-83-8	-4.32	106.90	33.42	3.29	$1\overline{01.89}$	298.15	0.6
195	$C_6H_{10}O$	cyclohexanone	108-94-1	$-\overline{230.12}$	-90.87					
196	$C_6H_{10}O$	4-methyl-3-penten-2-one(mesityl oxide)	141-79-7	-178.28		42.7		115.47	298.15	
197	$C_{6}H_{12}$	cyclohexane	110-82-7	-123.10	32.26	29.97	2.63	108.75	298.15	0.3
198	$C_6H_{12}$	methylcyclopentane	96-37-7	-106.00	36.51	29.08	6.93	113.13	298.15	$\overline{0.0}$
199	$C_6H_{12}$	1-hexene	592-41-6	-41.95	86.90	28.28	7.52	125.90	298.15	$\overline{0.4}$
200	$C_{6}H_{12}$	4-methylpent-1-ene	691-37-2	-49.44	85.69	28.57	3.60	129.44	298.15	
201	$C_6H_{12}O$	cyclohexanol	108-93-0	-294.55	$-1\overline{18.05}$		1.76	<u> </u>		<u>1.7</u>
202	$C_6H_{12}O$	2-hexanone (methyl butyl ketone)	591-78-6	-279.80	-129.90	36.35	$1\overline{4.90}$	124.10	298.15	
203	$C_6H_{12}O$	3-hexanone (ethyl propyl ketone)	589-38-8	-278.60	-126.10	35.36	13.49	123.43	298.15	
204	$C_6H_{12}O$	4-methyl-2(methyl isobutyl ketone)	108-10-1	-286.40	-135.10	34.49		125.81	298.15	2.8
205	$C_6H_{12}O$	butylvinylether	111-34-2	-179.20		36.59		130.25	298.15	
206	$C_6H_{12}O_2$	hexanoic acid	142-62-1	·				<u> </u>		
207	$C_6H_{12}O_2$	methyl pentanoate	624-24-8	-471.10	-296.50	35.36		132.75	298.00	
208	$C_6H_{12}O_2$	ethyl butanoate	105-54-4	-485.50	-312.00	35.47		132.95	298.15	1.8
209	$C_6H_{12}O_2$	propyl propanoate	106-36-5	-483.10	-309.70	35.54		132.50	298.15	1.8
210	$C_6H_{12}O_2$	ethyl 2-methylpropanoate	97-62-1	-499.60	-324.00	33.67		133.67	293.15	2.1
211	$C_6H_{12}O_2$	butyl ethanoate (butyl acetate)	123-86-4	-485.30	-312.10	36.28	14.59	132.51	298.15	1.8
212	$C_6H_{12}O_2$	2-methylpropyl ethanoate (isobutyl acetate)	110-19-0	-494.70	-322.00	35.90		133.87	298.15	$   \begin{array}{r}     \underline{1.8} \\     \underline{1.8} \\     \underline{2.1} \\     \underline{1.8} \\     \underline{1.9}   \end{array} $
213	$C_6H_{12}O_2$	pentyl methanoate (pentyl formate)	638-49-3	-448.20	-276.40			131.72	298.15	1.9

214	$C_6H_{12}O_2$	3-methylbutyl methanoate (isopentyl formate)	110-45-2	-453.80	-279.80			132.85	298.15	
215	$C_6H_{12}O_3$	2-ethoxyethylacetate	111-15-9			52.61		140.62	323.14	
216	$C_6H_{12}O_3$ $C_6H_{14}$	hexane	110-54-3	-166.92	0.15	28.85	13.07	131.59	298.15	0.0
217	$C_6H_{14}$	2-methylpentane	107-83-5	-174.55	-5.14	27.79	6.27	132.89	298.15	$\frac{0.0}{0.1}$
218	$C_6H_{14}$	3-methylpentane	96-14-0	-172.00	-3.20	28.06	5.30	130.62	298.15	
219	$C_6H_{14}$	2,2-dimethylbutane	75-83-2	-183.97	-9.63	26.31	0.58	133.73	298.15	
220	$C_6H_{14}$	2,3-dimethylbutane	79-29-8	-175.90	-2.05	27.28	0.79	131.17	298.15	0.2
221	$C_6H_{14}O$	1-hexanol	111-27-3	-316.50	-134.13	44.50	15.40	125.19	298.15	$\frac{0.2}{1.8}$
222	$C_6H_{14}O$	2-hexanol	626-93-7	-333.50		41.01		126.07	298.15	
223	$C_6H_{14}O$	3-hexanol	623-37-0	222.20		11101		120.07	2,0.10	
224	$C_6H_{14}O$	2-methyl-1-pentanol	105-30-6							
225	$C_{6}^{14}H_{14}O$	2-methyl-2-pentanol	590-36-3	-349.00		39.59		126.22	298.15	
226	$C_6H_{14}O$	2-methyl-3-pentanol	565-67-3							
227	$C_{6}^{0}H_{14}^{14}O$	4-methyl-1-pentanol	626-89-1	-324.90		44.46		126.22	298.15	
228	$C_6H_{14}O$	4-methyl-2-pentanol	108-11-2	-339.20		44.20		127.20	298.15	
229	$C_{6}^{0}H_{15}^{14}N$	N, N-diethylethanamine (triethyl amine)	121-44-8	-92.70	118.00	31.01		139.96	298.15	0.9
230	$C_{6}^{0}H_{15}^{13}N$	<i>N</i> -propyl-1-propanamine (dipropyl	142-84-7	-113.20		33.47		138.07	298.15	$\frac{0.9}{1.0}$
	0 15	amine)								_
231	$C_{7}F_{14}$	tetradecafluoromethylcyclohexane	355-02-2	-2898.00				195.67	298.15	
232	$C_{7}F_{16}$	hexadecafluoroheptane	335-57-9	-3386.11	-3093.02	36.29		223.92	293.15	0.0
233	$C_7H_3F_5$	pentafluorotoluene	771-56-2			34.75	12.99			
234	$C_7H_8$	toluene	108-88-3	50.17	122.29	33.18	6.85	106.87	298.15	0.4
235	$C_7H_8O$	benzyl alcohol	100-51-6	-72.38	18.20		8.97			1.7
236	$C_7H_8O$	2-methylphenol (o-cresol)	95-48-7	-128.57	$-\overline{34.27}$	45.30	$1\overline{3.94}$	103.64	298.15	1.6
237	$C_7H_8O$	3-methylphenol ( <i>m</i> -cresol)	108-39-4	-132.30	-40.07	47.45	9.41	104.99	298.15	1.8
238	$C_7H_8O$	4-methylphenol ( <i>p</i> -cresol)	106-44-5	-125.35	-31.55	47.45	$1\overline{1.89}$	105.00	298.15	1.6
239	$C_7H_9N$	2,3-dimethylpyridine (2,3-lutidine)	583-61-9	68.28	177.59	39.08	13.48	113.73	298.15	$\overline{2.2}$
240	$C_7H_9N$	2,4-dimethylpyridine (2,4-lutidine)	108-47-4	63.89	172.04	38.53	8.83	115.53	298.15	2.3
241	$C_7H_9N$	2,5-dimethylpyridine (2,5-lutidine)	589-93-5	66.44	174.34	38.68	14.65	115.87	298.15	$\overline{2.2}$
242	$C_7H_9N$	2,6-dimethylpyridine (2,6-lutidine)	108-48-5	58.70	168.40	37.46	13.04	116.52	298.15	1.7
243	$C_7H_9N$	3,4-dimethylpyridine (3,4-lutidine)	583-58-4	70.04	179.34	39.99	14.70	112.32	298.15	0.4 1.7 1.6 1.8 1.6 2.2 2.3 2.2 1.7 1.9 2.6
244	$C_7H_9N$	3,5-dimethylpyridine (3,5-lutidine)	591-22-0	72.80	182.44	39.46	13.11	114.26	298.15	2.6
245	$C_7H_{12}O_2$	butyl-2-propenoate(butylacrylate)	141-32-2	-375.30		47.31				· <u>——</u>

Section B Basic Constants II (Continued)

No.	Formula	Name	CAS #	DelHf0, kJ/mol	DelGf0, kJ/mol	DelHb, kJ/mol	DelHm, kJ/mol	V liq, cm <sup>3</sup> /mol	T liq, K	Dipole, Debye
246	$C_{7}H_{14}$	cycloheptane	291-64-5	-118.10	64.30	33.18	1.88	121.73	298.15	0.0
247	$C_7^{'}H_{14}^{'14}$	methylcyclohexane	108-87-2	-154.70	27.64	31.27	6.75	128.35	298.15	$\overline{0.0}$
248	$C_{7}H_{14}^{14}$	ethylcyclopentane	1640-89-7	-126.90	45.07	31.96	6.87	128.83	298.15	$\overline{0.0}$
249	$C_{7}^{'}H_{14}^{14}$	cis-1,3-dimethylcyclopentane	2532-58-3	-135.90	39.23	30.40	7.37	131.91	298.15	
250	$C_7H_{14}$	trans-1,3-dimethylcyclopentane	1759-58-6	-133.50	41.42	30.80	7.27	132.64	298.15	
251	$C_7H_{14}$	1-heptene	592-76-7	-62.76	95.06	31.09	12.66	141.77	298.15	0.3
252	$C_7H_{14}O_2$	heptanoic acid	111-14-8							
253	$C_7H_{14}O_2$	ethyl pentanoate	539-82-2	-505.90	-303.50	36.96		148.45	293.15	
254	$C_7H_{14}O_2$	ethyl 3-methylbutanoate	108-64-5	-527.90	-324.00	37.00		149.99	293.15	
255	$C_7H_{14}O_2$	propyl butanoate	105-66-8	-505.30	-303.00			149.95	298.15	1.8
256	$C_7H_{14}O_2$	2-methylpropyl propanoate	540-42-1	-512.70	-311.20			150.21	298.15	_
257	$C_7H_{14}O_2$	propyl 2-methylpropanoate	644-49-5	-518.10	-314.00			147.27	273.15	
258	$C_7H_{14}O_2$	3-methylbutyl ethanoate (isopentyl	123-92-2	-511.20	-306.80	37.50		150.26	298.15	1.8
		acetate)								
259	$C_7H_{16}$	heptane	142-82-5	-187.80	8.20	31.77	14.03	147.47	298.15	$\frac{0.0}{0.0}$
260	$C_7H_{16}$	2-methylhexane	591-76-4	-194.60	3.70	30.62	9.19	148.60	298.15	0.0
261	$C_7H_{16}$	3-methylhexane	589-34-4	-191.30	5.30	30.89		146.74	298.15	$\overline{0.0}$
262	$C_7H_{16}$	3-ethylpentane	617-78-7	-189.50	11.40	31.12	9.55	144.40	298.15	$\overline{0.0}$
263	$C_7H_{16}$	2,2-dimethylpentane	590-35-2	-205.81	0.80	29.23	5.82	149.67	298.15	0.0
264	$C_7H_{16}$	2,3-dimethylpentane	565-59-3	-194.10	5.82	30.46		145.05	298.15	$\frac{0.0}{0.0}$
265	$C_7H_{16}$	2,4-dimethylpentane	108-08-7	-201.67	3.51	29.55	6.84	149.95	298.15	0.0
266	$C_7H_{16}$	3,3-dimethylpentane	562-49-2	-201.40	3.50	29.62	6.85	145.40	298.15	0.0
267	$C_7H_{16}$	2,2,3-trimethylbutane	464-06-2	-204.40	5.00	28.90	2.26	146.15	298.15	0.0
268	$C_7H_{16}O$	1-heptanol	111-70-6	-330.90	-119.56	48.10	13.20	141.95	298.15	$\frac{0.0}{0.0}$ $\frac{1.7}{0.0}$
269	$C_7H_{16}O$	2-heptanol	543-49-7							
270	$C_7H_{16}O$	4-heptanol	589-55-9							
271	$C_8F_{18}$	octadecafluorooctane	307-34-6	-3798.24	-3463.70	33.38		253.21	293.15	0.0
272	$C_8H_8O$	methylphenylketone(acetophenone)	98-86-2			55.40		119.94	323.14	
273	$C_8H_{10}$	ethylbenzene	100-41-4	29.92	130.73	35.57	9.18	123.08	298.15	0.4
274	$C_8H_{10}$	1,2-dimethylbenzene ( <i>o</i> -xylene)	95-47-6	19.08	122.05	36.24	13.60	121.25	298.15	0.5

275	$C_8H_{10}$	1,3-dimethylbenzene ( <i>m</i> -xylene)	108-38-3	17.32	118.89	35.66	11.57	123.47	298.15	$\frac{0.3}{0.1}$
276	$C_8H_{10}$	1,4-dimethylbenzene ( <i>p</i> -xylene)	106-42-3	18.03	121.48	35.67	16.81	123.93	298.15	0.1
277	$C_8H_{10}O$	2-ethylphenol	90-00-6	-145.23	-23.15	46.20	13.94	120.41	298.15	
278	$C_8H_{10}O$	3-ethylphenol	620-17-7	-146.06	-25.01	48.90	9.41	121.25	298.15	
279	$C_8H_{10}O$	4-ethyl-phenol	123-07-9	-144.05	-21.43	48.90	11.89	120.84	298.15	
280	$C_8H_{10}O$	2,3-dimethylphenol (2,3 xylenol)	526-75-0	-157.19	-33.20	47.60	21.02			
281	$C_8H_{10}O$	2,4-dimethylphenol (2,4 xylenol)	105-67-9	-162.88	-41.07	47.40		119.75	293.15	$\frac{2.0}{1.5}$
282	$C_8H_{10}O$	2,5-dimethylphenol (2,5 xylenol)	95-87-4	-161.63	-39.52	47.10	23.38	119.30	293.15	1.5
283	$C_8H_{10}O$	2,6-dimethylphenol (2,6 xylenol)	576-26-1	-161.74	-38.89	44.70	18.90			
284	$C_8H_{10}O$	3,4-dimethylphenol (3,4 xylenol)	95-65-8	-156.56	-34.13	50.10	18.13	118.04	293.15	$\frac{1.7}{1.8}$
285	$C_8H_{10}O$	3,5-dimethylphenol (3,5 xylenol)	108-68-9	-161.54	-39.26	49.70	18.00	119.54	293.15	1.8
286	$C_8H_{16}$	cyclooctane	292-64-8	-124.40	91.38	35.90	2.41	134.87	298.15	0.0
287	$C_8H_{16}$	t-1,4-dimethylcyclohexane	2207-04-7	-168.300	62.43	37.88	12.33	147.97	298.15	
288	$C_8H_{16}$	1-octene	111-66-0	-83.59	103.20	34.07	15.57	157.85	298.15	0.3
289	$C_8H_{16}O_2$	octanoic acid	124-07-2				21.36			_
290	$C_8H_{16}O_2$	2-ethylhexanoic acid	149-57-5			76.31		161.81	323.15	
291	$C_8H_{16}O_2$	propyl pentanoate	141-06-0	-525.70	-294.40					
292	$C_8^{10}C_2^{2}$	2-methylpropyl butanoate	539-90-2	-533.00	-302.50			168.24	293.15	
293	$C_8H_{16}O_2$	propyl 3-methylbutanoate	557-00-6	-547.10	-314.00			167.11	293.15	
294	$C_8^{10}C_2^{2}$	3-methylbutyl propanoate	105-68-0	-529.20	-295.80			166.72	298.15	
295	$C_8^{10}C_2^{2}$	2-methylpropyl 2-methylpropanoate	97-85-8	-546.20	-313.00			164.82	273.15	
296	C <sub>8</sub> H <sub>18</sub>	octane	111-65-9	-208.75	16.27	34.41	20.65	163.53	298.15	0.0
297	$C_8^{18}$	2-methylheptane	592-27-8	-215.35	11.96	33.26	11.92	164.63	298.15	
298	$C_8^{"}H_{18}^{"}$	3-methylheptane	589-81-1	-212.50	13.00	33.66	11.38	162.78	298.15	
299	$C_8H_{18}$	4-methylheptane	589-53-7	-211.96	15.98	33.35	10.84	163.06	298.15	
300	$C_8H_{18}$	3-ethylhexane	619-99-8	-210.71	17.11	33.59		161.01	298.15	
301	$C_8H_{18}$	2,2-dimethylhexane	590-73-8	-224.60	10.68	32.07	6.78	165.29	298.15	
302	$C_8H_{18}$	2,3-dimethylhexane	584-94-1	-213.80	15.65	33.17		161.31	298.15	
303	$C_8H_{18}$	2,4-dimethylhexane	589-43-5	-219.24	11.51	32.51		164.08	298.15	
304	$C_8H_{18}$	2,5-dimethylhexane	592-13-2	-222.51	9.95	32.54	12.95	165.70	298.15	
305	$C_8H_{18}$	3,3-dimethylhexane	563-16-6	-219.99	13.68	32.31	7.11	161.81	298.15	
306	$C_8H_{18}$	3,4-dimethylhexane	583-48-2	-212.67	16.99	33.24	7.11	159.73	298.15	
307	$C_8H_{18}$	3-ethyl-2-methylpentane	609-26-7	-212.80	19.23	32.93	11.34	159.72	298.15	
308	$C_8H_{18}$	3-ethyl-3-methylpentane	1067-08-9	-214.85	22.88	32.78	10.84	157.88	298.15	
309	$C_8H_{18}$	2,2,3-trimethylpentane	564-02-3	-214.85 $-219.95$	17.97	31.94	8.62	160.43	298.15	
507	C811 <sub>18</sub>	2,2,5 dimenty pentane	JUT-UZ=J	217.73	17.77	31.77	0.02	100.73	270.13	

Section B Basic Constants II (Continued)

No.	Formula	Name	CAS #	DelHf0, kJ/mol	DelGf0, kJ/mol	DelHb, kJ/mol	DelHm, kJ/mol	V liq, cm <sup>3</sup> /mol	T liq, K	Dipole, Debye
310	C <sub>8</sub> H <sub>18</sub>	2,2,4-trimethylpentane (isooctane)	540-84-1	-224.01	14.21	30.79	9.04	166.07	298.15	
311	$C_8H_{18}$	2,3,3-trimethylpentane	560-21-4	-218.45	18.56	32.12	$\frac{0.84}{0.86}$	158.15	298.15	
312	$C_8H_{18}$	2,3,4-trimethylpentane	565-75-3	-217.32	19.28	32.36	9.26	159.74	298.15	
313	$C_8H_{18}$	2,2,3,3-tetramethylbutane	594-82-1	-225.73	10.31	31.42	7.54	139.02	298.15	
314	$C_8H_{18}O$	1-octanol	111-87-5	-356.90	$-1\overline{16.59}$	46.90	7.0	158.37	298.15	2.0
315	$C_8H_{18}O$	2-octanol	123-96-6	220.50	110.00	44.40		159.38	298.15	1.6
318	$C_8H_{18}O$	2-ethyl-1-hexanol	104-76-7							$\frac{2.0}{1.6}$
319	$C_8H_{19}N$	<i>n</i> -octylamine	111-86-4	-173.46		54.63		162.33	298.15	
320	$C_8H_{19}N$	<i>N</i> -butyl-1-butanamine (dibutyl amine)	111-92-2			38.44		170.58	298.15	1.1
321	$C_{9}F_{20}$	eicosafluorononane	375-96-2	-4210.36	-3834.38			271.15	298.15	$\frac{0.0}{2.3}$ $\frac{2.7}{2.7}$
322	$C_0H_7N$	quinoline	91-22-5			49.70	10.80	118.50	298.15	$\overline{2.3}$
323	$C_9H_7N$	isoquinoline	119-65-3			49.00		118.39	293.15	$\overline{2.7}$
324	$C_9H_{10}$	indan	496-11-7	60.75	166.61	39.63		123.15	298.15	
325	$C_9H_{12}$	propylbenzene	103-65-1	7.91	137.58	38.20	9.27	140.20	298.15	
326	$C_9H_{12}$	1-methylethylbenzene (cumene)	98-82-8	4.00	139.05	37.50	7.79	140.17	298.15	0.8
327	$C_9H_{12}$	1-ethyl-4-methylbenzene	622-96-8	-2.05	130.28	38.40	13.36	139.60	293.15	
328	$C_9H_{12}$	1,2,3-trimethylbenzene	526-73-8	-9.50	124.96	40.00	8.18	133.22	278.70	
329	$C_9H_{12}$	1,2,4-trimethylbenzene	95-63-6	-13.81	117.50	39.20	13.19	136.00	278.70	
330	$C_9H_{12}$	1,3,5-trimethylbenzene (mesitylene)	108-67-8	-15.90	118.26	39.00	9.51	142.99	298.15	0.1
331	$C_9H_{18}$	1-nonene	124-11-8	-104.00	111.80	36.31	18.08	174.05	298.15	
332	$C_9H_{18}O_2$	nonanoic acid	112-05-0				20.28			
333	$C_9H_{18}O_2$	3-methylbutyl butanoate	106-27-4	-551.50	-289.20			184.02	298.15	
334	$C_9H_{20}$	nonane	111-84-2	-228.86	25.00	36.91	15.50	179.70	298.15	0.0
335	$C_9H_{20}$	2-methyloctane	3221-61-2	-235.85	20.30	36.10	18.00	180.75	298.15	
336	$C_9H_{20}$	2,2-dimethylheptane	1071-26-7	-246.10	18.10	34.60	8.90	181.51	298.15	
337	$C_9H_{20}$	2,2,5-trimethylhexane	3522-94-9	-253.26	14.00	40.30	6.20	182.39	298.15	
338	$C_9H_{20}$	2,2,3,3-tetramethylpentane	7154-79-2	-237.11	37.60	34.30	2.33	170.34	298.15	
339	$C_9H_{20}$	2,2,3,4-tetramethylpentane	1186-53-4	-234.97	35.40	33.70	0.50	174.45	298.15	
340	$C_9H_{20}$	2,2,4,4-tetramethylpentane	1070-87-7	-242.25	34.20	38.30	9.75	179.23	298.15	
341	$C_9H_{20}$	2,3,3,4-tetramethylpentane	16747-38-9	-236.31	38.30	34.50	9.00	170.76	298.15	

342	$C_9H_{20}O$	1-nonanol	143-08-8	-381.20	-111.92	54.40		174.92	298.15	1.7
343	$C_9H_{20}O$	2-nonanol	628-99-9							
344	$C_{10}F_{22}$	docosafluorodecane		-4622.48	-4205.05					0.0
345	$C_{10}H_{8}$	naphthalene	91-20-3	150.30	223.50	43.40	19.12	129.13	333.15	0.0
346	$C_{10}H_{12}$	1,2,3,4-tetrahydronaphthalene	119-64-2	26.61	166.89			136.27	293.15	
347	$C_{10}H_{12}$	1-methylindan	767-58-8							
348	$C_{10}H_{12}$	2-methylindan	824-63-5							
349	$C_{10}H_{12}$	4-methylindan	824-22-6							
350	$C_{10}H_{12}$	5-methylindan	874-35-1							
351	$C_{10}H_{14}$	butylbenzene	104-51-8	-13.14	145.39	38.87	11.22	156.78	298.15	$\frac{0.4}{0.3}$
352	$C_{10}H_{14}$	2-methylpropylbenzene	538-93-2	-20.34	140.20	37.80	12.50	158.08	298.15	0.3
353	$C_{10}H_{14}$	1,4-diethylbenzene	105-05-5	-20.37	140.30	39.40	10.60	156.45	298.15	0.1
354	$C_{10}H_{14}$	1-(1-methylethyl)-4-methylbenzene	99-87-6	-27.74	136.70	38.20	9.66	157.49	298.15	0.0
355	$C_{10}H_{14}$	1,2,4,5-tetramethylbenzene	95-93-2	-44.56	120.50	45.52	21.00	152.28	298.15	0.0
356	$C_{10}H_{18}$	cis-bicyclo[4.4.0]decane (cis-decalin)	493-01-6	-169.20	85.60	41.00		154.83	298.15	0.0
357	$C_{10}H_{18}$	trans-bicyclo[4.4.0]decane (trans-decalin)	493-02-7	-182.10	74.20	40.20		159.66	298.15	0.0
358	$C_{10}H_{20}$	1-decene	872-05-9	-124.20	121.10			189.30	293.15	
359	$C_{10}H_{20}O_2$	decanoic acid	334-48-5				28.01			
360	$C_{10}H_{22}$	decane	124-18-5	-249.53	33.30	38.75	28.78	195.95	298.15	0.0
361	$C_{10}H_{22}$	2,2,5-trimethylheptane	20291-95-6	-272.21	22.20	36.20	12.00	196.44	298.15	
362	$C_{10}H_{22}$	3,3,5-trimethylheptane	7154-80-5	-259.87	32.60	36.40	14.00	192.48	298.15	
363	$C_{10}H_{22}$	2,2,3,3-tetramethylhexane	13475-81-5	-257.99	47.20	36.20	12.40	187.02	298.15	
364	$C_{10}H_{22}$	2,2,5,5-tetramethylhexane	1071-81-4	-285.89	19.50	42.40	9.80	199.06	298.15	
365	$C_{10}H_{22}O$	1-decanol	112-30-1	-396.70	-98.45	49.80	38.00	191.51	298.15	1.8
366	$C_{10}H_{24}N_4$	octamethylethenetetramine	996-70-3	132.9		53.85		232.40	298.15	
367	$C_{11}H_{10}$	1-methylnaphthalene	90-12-0	115.20	216.40	46.00	6.94	139.37	293.15	$\frac{0.5}{0.4}$
368	$C_{11}H_{10}$	2-methylnaphthalene	91-57-6	114.90	215.00	46.50	12.13	145.77	333.15	0.4
369	$C_{11}H_{24}$	undecane	1120-21-4	-270.16	41.25	41.20	22.32	212.24	298.15	0.0
370	$C_{11}H_{24}O$	1-undecanol	112-42-5							
371	$C_{12}H_{10}$	1,1'-biphenyl	92-52-4	182.42	281.08	48.30	18.57	155.77	347.00	0.0
372	$C_{12}H_{12}$	1,6-dimethylnaphthalene	575-43-9	79.80	210.90			155.93	293.15	
373	$C_{12}H_{12}$	2,7-dimethylnaphthalene	582-16-1	79.00	210.70	59.50	23.35			
374	$C_{12}H_{18}$	1,3,5-triethylbenzene	102-25-0			59.22		193.39	323.14	
375	$C_{12}H_{20}$	1,3-dimethyltricyclo[3.3.1.1 <sup>3,7</sup> ]decane	702-79-4			49.37		186.61	323.14	
		(1,3-dimethyladamantane)						<del></del>		

Section B Basic Constants II (Continued)

No.	Formula	Name	CAS #	DelHf0, kJ/mol	DelGf0, kJ/mol	DelHb, kJ/mol	DelHm, kJ/mol	V liq, cm <sup>3</sup> /mol	T lig, K	Dipole, Debye
376	$C_{12}H_{24}$	1-dodecene	112-41-4	-165.50	138.00			222.06	293.15	
377		dodecane	112-41-4	$\frac{-103.30}{-290.79}$	49.53	43.40	36.58	$\frac{222.00}{228.59}$	293.13	0.0
378	$C_{12}H_{26}$	1-dodecanol	112-40-3	-290.79 $-443.10$		43.40	30.36	220.39	290.13	0.0
378 379	$C_{12}H_{26}O$		101-81-5	$\frac{-443.10}{165.20}$	-87.13			167.23	293.15	0.4
	$C_{13}H_{12}$	diphenylmethane	629-50-5		57.01	15 65	20.50			$\frac{0.4}{0.0}$
380	$C_{13}H_{28}$	tridecane	629-30-3 112-70-9	-311.42	57.81	45.65	28.50	244.94	298.15	0.0
381	$C_{13}H_{28}O$	1-tridecanol		207.50	200.20		16.47			0.0
382	$C_{14}H_{10}$	phenanthrene	85-01-8	207.50	308.20		16.47			$\frac{0.0}{0.0}$
383	$C_{14}H_{10}$	anthracene	120-12-7	230.90	333.70		28.83	226.15	252 15	0.0
384	$C_{14}H_{22}$	1,4-di(trimethylmethyl)benzene (p-ditertbutylbenzene)	1012-72-2				22.48	236.17	373.15	
385	$C_{14}H_{30}$	tetradecane	629-59-4	-332.05	66.09	47.61	45.61	261.32	298.15	0.0
386	$C_{14}^{14}H_{30}^{30}O$	1-tetradecanol	112-72-1							
387	$C_{15}^{14}H_{32}^{30}$	pentadecane	629-62-9	-352.68	74.37	49.45	34.80	277.71	298.15	0.0
388	$C_{15}H_{32}O$	1-pentadecanol	629-76-5							
389	$C_{16}^{15}H_{34}^{32}$	hexadecane	544-76-3	-373.31	82.65	51.21	51.84	294.11	298.15	0.0
390	$C_{16}^{10}H_{34}^{34}$	2,2,4,4,6,8,8-heptamethylnonane	4390-04-9							
391	$C_{16}H_{34}O$	1-hexadecanol	4485-13-6				34.29			
392	$C_{17}^{10}H_{36}^{34}$	heptadecane	629-78-7	-393.94	90.93	52.89	40.50	310.45	298.15	0.0
393	$C_{17}^{17}H_{36}^{30}O$	1-heptadecanol	1454-85-9	-546.30	-44.67					
394	$C_{18}^{17}H_{14}^{30}$	1,2-diphenylbenzene	84-15-1							
395	$C_{18}^{13}H_{14}^{14}$	1,3-diphenylbenzene	92-06-8							
396	$C_{18}^{13}H_{14}^{14}$	1,4-diphenylbenzene	92-94-4				35.50			0.7
397	$C_{18}^{13}H_{38}^{14}$	octadecane	593-45-3	-414.57	99.21	54.46	61.39	326.66	298.15	$\overline{0.0}$
398	$C_{18}^{18}H_{38}^{38}O$	1-octadecanol	112-92-5	-566.90	-36.22					1.7
399	$C_{19}^{18}H_{40}^{38}$	nonadecane	629-92-5	-435.20	107.49	56.02	45.82	343.25	298.15	$\begin{array}{c} 0.7 \\ \hline 0.0 \\ \hline 1.7 \\ \hline 0.0 \end{array}$
400	$C_{19}^{19}H_{40}^{40}O$	1-nonadecanol	1454-84-8							
401	$C_{20}^{19}H_{42}^{40}$	eicosane	112-95-8	-455.83	115.77	57.49	69.87	361.18	298.15	0.0
402	$C_{20}^{20}H_{42}^{42}O$	1-eicosanol	629-96-9	-608.10	-19.43					_
403	$C_{21}^{20}H_{44}^{42}$	heneicosane	629-94-7				47.70	381.11	313.15	0.0
404	$C_{22}^{21}H_{46}^{44}$	docosane	629-97-0				49.96	399.14	318.15	0.0

405	$C_{23}H_{48}$	tricosane	638-67-5				41.76			$\frac{0.0}{0.0}$
406	$C_{24}H_{50}$	tetracosane	646-31-1				54.89	434.96	324.25	0.0
407	ClD	deuterium chloride	7698-05-7	-93.33	-95.93			31.02	193.15	
408	ClFO <sub>3</sub>	perchloryl fluoride	7616-94-6	-21.44	50.62	19.33		62.70	243.15	0.0
409	ClF <sub>5</sub>	chlorine pentafluoride	13637-63-3	-238.49	-147.11			73.41	298.15	
410	ClH	hydrogen chloride	7647-01-0	-92.31	-95.19	16.15		30.28	183.15	1.1
411	$ClH_4N$	ammonium chloride	12125-02-9							
412	CINO	nitrogen oxychloride	2696-92-6	51.71	65.97	25.78		46.10	261.00	$\frac{1.8}{0.0}$
413	$Cl_2$	chlorine	7782-50-5	0.00	0.00	20.41		45.36	239.00	0.0
414	DH	deuterium hydride	13983-20-5	0.32	-1.46	1.08	0.16			0.0
415	DI	deuterium iodide	14104-45-1	26.23	1.84			46.04	237.00	
416	$D_2$	deuterium	7782-39-0	0.00	0.00	1.23	0.20	24.41	22.70	0.0
417	$D_2$	deuterium, normal	800000-54-8	0.00	0.00	1.23	0.20			$\frac{0.0}{0.0}$
418	$D_2O$	deuterium oxide	7789-20-0	-249.20	-234.53	41.46	6.38	18.13	298.15	1.9
419	$D_2S$	deuterium sulfide	13536-94-2	-23.89	-35.39	18.85	2.37			
420	$D_3N$	trideuteroammonia	13550-49-7	-64.28	-31.69					
421	$D_3P$	trideuterophosphine	13537-03-6							
422	FH	hydrogen fluoride	7664-39-3	-273.30	-275.40			20.69	293.15	1.8
423	$FNO_2$	nitrogen dioxyfluoride	10022-50-1	-108.78	-66.55	18.05				$\begin{array}{c} 1.8 \\ \hline 0.5 \\ \hline 0.0 \\ \hline 1.9 \\ \hline 0.2 \\ \hline 0.0 \\ \hline 0.2 \\ \hline \end{array}$
424	$F_2$	fluorine	7782-41-4	0.00	0.00	6.62		25.16	85.00	$\overline{0.0}$
425	$F_2HN$	difluoroamine	10405-27-3							1.9
426	$F_2N_2$	cis-difluorodiazene	13812-43-6	74.89	114.89					$\overline{0.2}$
427	$F_2N_2$	trans-difluorodiazene	13776-62-0	81.17	120.35					$\overline{0.0}$
428	$F_2O$	oxygen difluoride	7783-41-7	24.70	42.01	11.09		34.93	123.15	$\overline{0.2}$
429	$F_2Xe$	xenon difluoride	13709-36-9					54.09	298.15	$\overline{0.0}$
430	$\tilde{F_3}N$	nitrogen trifluoride	7783-54-2	-129.70	-88.29	11.56		46.11	144.00	$\overline{0.2}$
431	$F_3NO$	trifluoroamine oxide	13847-65-9	-163.30	-96.46					
432	$F_4N_2$	tetrafluorohydrazine	10036-47-2	-8.37	79.66	13.27		69.34	163.00	0.3
433	$\vec{F_4}$ S	sulfur tetrafluoride	7783-60-0	-774.04	$-7\overline{31.01}$	26.44		55.82	195.00	1.0
434	$F_4$ Xe	xenon tetrafluoride	13709-61-0	-187.60				68.41	298.15	$\begin{array}{r} 0.3 \\ \hline 1.0 \\ \hline 0.0 \\ \hline 0.0 \end{array}$
435	$F_6S$	sulfur hexafluoride	2551-62-4	-1220.89	-1117.09			79.81	223.15	$\overline{0.0}$
436	$F_6^0U$	uranium hexafluoride	7783-81-5	-2139.00	-2060.00					0.0
437	нÏ	hydrogen iodide	10034-85-2	26.50	1.70	19.76		46.06	243.15	0.5
438	$H_2$	hydrogen	1333-74-0	0.00	0.00	0.89	0.12	28.39	20.00	$\frac{0.5}{0.0}$
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Section B Basic Constants II (Continued)

No.	Formula	Name	CAS #	DelHf0, kJ/mol	DelGf0, kJ/mol	DelHb, kJ/mol	DelHm, kJ/mol	V liq, cm <sup>3</sup> /mol	T liq, K	Dipole, Debye
								CIII / IIIOI	r nq, ix	
439	$H_2$	hydrogen, normal	800000-51-5	0.00	0.00	0.90	0.12	40.05	200.45	$\frac{0.0}{1.0}$
440	H <sub>2</sub> O	water	7732-18-5	-241.81	-228.42	40.66	6.01	18.07	298.15	$\frac{1.8}{0.9}$
441	$H_2S$	hydrogen sulfide	7783-06-4	-20.63	-33.43	18.68	2.38	34.32	214.00	0.9
442	$H_2S_2$	dihydrogen disulfide	13465-07-1	15.52	-1.67	35.10	7.53	49.85	298.15	
443	$H_2S_3$	dihydrogen trisulfide	13845-23-3					66.23	298.15	
444	$H_2S_4$	dihydrogen tetrasulfide	13845-25-5					82.66	298.15	
445	$H_2S_5$	dihydrogen pentasulfide	13845-24-4							
446	$H_2Se$	hydrogen selenide	7783-07-5			19.70		38.20	231.00	
447	$H_3N$	ammonia	7664-41-7	-45.94	-16.41	23.35	5.66	24.96	239.15	1.5
448	$H_3P$	phosphine	7803-51-2	5.44	13.44	14.60		45.57	183.15	0.6
449	$H_4N_2$	hydrazine	302-01-2	$9\overline{5.40}$	159.38	44.77	12.66	31.79	293.15	3.0
450	He	helium	7440-59-7	0.00	0.00	0.08		32.54	4.30	$\begin{array}{r} 1.5 \\ \underline{0.6} \\ \underline{3.0} \\ \underline{0.0} \end{array}$
451	He	helium-3	14762-55-1	0.00	0.00					$\overline{0.0}$
452	$I_2$	iodine	7553-56-2	62.42	19.33	41.57		67.86	453.00	$\begin{array}{c} \underline{0.0} \\ \underline{1.3} \\ \underline{0.0} \\ \underline{0.2} \\ \underline{0.0} \\ 0.2 \\ \end{array}$
453	Kr	krypton	7439-90-9	0.00	0.00	9.08		34.63	120.00	$\overline{0.0}$
454	NO	nitrogen monoxide (nitric oxide)	10102-43-9	90.25	86.58	$1\overline{3.78}$	2.30	23.44	121.00	$\overline{0.2}$
455	$N_2$	nitrogen	7727-37-9	0.00	0.00	5.58	0.72	34.84	78.00	$\overline{0.0}$
456	$N_2O$	dinitrogen oxide (nitrous oxide)	10024-97-2	82.05	104.18	16.55	6.54	35.90	184.00	$\overline{0.2}$
457	$N_2^2O_4$	dinitrogen tetroxide (nitrogen dioxide)	10544-72-6	9.16	97.85	29.00	14.65	63.59	293.15	
458	Ne	neon	7440-01-9	0.00	0.00	1.71		16.76	27.00	0.0
459	$OT_2$	tritium oxide	14940-65-9	-249.37	-226.24			18.15	298.15	_
460	O, -	oxygen	7782-44-7	0.00	0.00	6.82	0.44	27.85	90.00	0.0
461	$O_2S$	sulfur dioxide	7446-09-5	-296.81	-300.14	24.94	7.40	44.03	263.15	1.6
462	$O_3$	ozone	10028-15-6	142.70	163.10	14.20		35.40	161.00	0.6
463	$O_3$ S	sulfur trioxide	7446-11-9	-395.72	-370.93	40.69	7.53	42.10	298.15	_
464	Rn	radon	10043-92-2	0.00	0.00			50.46	211.15	0.0
465	S	sulfur	7704-34-9	276.98	236.50	10.46	1.61	-		$\overline{0.0}$
466	Se	selenium	7782-49-2	0.00	0.00	95.48				$\frac{0.0}{0.0}$
467	$T_2$	tritium	10028-17-8	0.00	0.00					0.0
468	Xe	xenon	7440-63-3	0.00	0.00	12.57		42.91	165.00	$\begin{array}{c} 0.0 \\ 0.0 \\ 0.0 \\ \hline 0.0 \\ 0.0 \\ \end{array}$