Section A Basic Constants I.

									Vc,	Zc =	
No.	Formula	Name	CAS #	Mol. Wt.	Tfp, K	Tb, K	Tc, K	Pc, bar	cm ³ /mol	PcVc/RTc	Omega
1	Ar	argon	7440-37-1	39.948	83.80	87.27	150.86	48.98	74.57	0.291	-0.002
2	Br_2	bromine	7726-95-6	159.808	265.85	331.90	584.10	103.00	135.00	0.269	0.119
3	BrD	deuterium bromide	13536-59-9	81.918	185.69	206.65	362.00				
4	BrF ₃	bromine trifluoride	7787-71-5	136.899	281.92	398.89	600.00	49.90	114.70	0.115	0.413
5	BrF ₅	bromine pentafluoride	7789-30-2	174.896	212.65	314.31	470.00				
6	BrH	hydrogen bromide	10035-10-6	80.912	186.34	206.46	363.20	85.10			0.069
7	CBrClF ₂	bromochlorodifluoromethane	353-59-3	165.365	113.65	269.20	426.90	42.60	246.00	0.295	0.182
8	CBrF ₃	bromotrifluoromethane	75-63-8	148.910	105.15	215.41	340.15	39.70	200.00	0.275	0.174
9	CBr ₂ F ₂	dibromodifluoromethane	75-61-6	209.816	163.05	295.94	471.30	45.30	250.00	0.286	
10	CClF ₃	chlorotrifluoromethane	75-72-9	104.459	92.00	191.71	301.84	38.73	180.30	0.276	0.175
11	CCl_2F_2	dichlorodifluoromethane (R-12)	75-71-8	120.913	115.19	243.45	385.10	41.30	217.00	0.280	0.179
12	CCl ₃ F	trichlorofluoromethane (R-11)	75-69-4	137.368	162.69	296.81	471.10	44.72	248.00	0.283	0.195
13	CCl ₄	tetrachloromethane	56-23-5	153.822	250.33	349.79	556.30	45.57	276.00	0.271	
14	CF_4	tetrafluoromethane	75-73-0	88.005	89.55	145.11	227.51	37.45	140.70	0.279	0.177
15	CHBrF ₂	bromodifluoromethane	1511-62-2	130.920	258.65	257.68	412.00	47.90	167.00	0.234	0.172
16	CHClF ₂	chlorodifluoromethane (R-22)	75-45-6	86.468	115.73	232.14	369.28	49.86	166.00	0.274	0.221
17	CHCl ₂ F	dichlorofluoromethane	75-43-4	102.923	138.20	281.97	451.52	51.87	196.00	0.271	0.207
18	CHCl ₃	trichloromethane	67-66-3	119.377	209.74	334.33	536.50	55.00	240.00	0.296	
19	CHF ₃	trifluoromethane (R-23)	75-46-7	70.014	117.96	191.11	298.97	48.36	133.00	0.259	0.267
20	CH ₂ Cl ₂	dichloromethane	75-09-2	84.932	176.00	312.79	510.00	61.00			
21	CH_2F_2	difluoromethane	75-10-5	52.024	137.00	221.43	351.26	58.05	121.00	0.245	0.278
22	CH_2O_2	methanoic acid (formic acid)	64-18-6	46.026	281.50	374.04	588.00	58.07			0.316
23	CH ₃ Cl	chloromethane	74-87-3	50.488	175.44	248.95	416.20	66.80	143.00	0.276	0.151
24	CH ₃ F	fluoromethane	593-53-3	34.033	131.35	194.88	315.00	55.48	113.30	0.240	0.204
25	CH ₃ NO ₂	nitromethane	75-52-5	61.040	244.60	374.35	588.00	58.70	173.00	0.208	
26	CH ₄	methane	74-82-8	16.043	90.69	111.66	190.56	45.99	98.60	0.286	0.011
27	CH ₄ O	methanol	67-56-1	32.042	175.49	337.69	512.64	80.97	$1\overline{18.00}$	0.224	0.565
28	CH ₄ S	methanethiol (methyl mercaptan)	74-93-1	48.109	150.18	279.11	470.00	72.30	145.00	0.268	0.150
29	CH ₅ N	methanamine (methyl amine)	74-89-5	31.057	179.69	266.82	430.00	74.20	125.00	0.259	0.283
30	CO	carbon monoxide	630-08-0	28.010	68.15	81.66	132.85	34.94	93.10	0.292	0.045

Section A Basic Constants I (Continued)

No.	Formula	Name	CAS #	Mol. Wt.	Tfn V	Tb, K	Tc, K	Do her	Vc,	Zc = PcVc/RTc	Omaga
NO.	rominia	Name	CAS#	MOI. Wt.	11р, к	10, K	IC, K	rc, bai	CIII / IIIOI	PCVC/KIC	Omega
31	CO_2	carbon dioxide	124-38-9	44.010	216.58		304.12	73.74	94.07	0.274	0.225
32	$C_2Br_2ClF_3$	1,2-dibromo-2-chloro-1,1,2-trifluroethane	354-51-8	276.278	323.15	367.06	561.20	36.10	357.80	0.285	0.251
33	C_2ClF_5	1-chloro-1,1,2,2,2-pentafluoroethane	76-15-3	154.467	173.73	234.08	353.10	31.29	256.00	0.273	0.251
34	$C_2Cl_2F_4$	1,1-dichloro-1,2,2,2-tetrafluoroethane	374-07-2	170.921	216.58	276.59	418.70	32.13	294.00	0.271	0.244
35	$C_2Cl_2F_4$	1,2-dichloro-1,1,2,2-tetrafluoroethane	76-14-2	170.921	180.55	276.58	418.90	32.37	294.00	0.273	0.246
36	$C_2Cl_3F_3$	1,1,2-trichloro-1,2,2-trifluoroethane	76-13-1	187.375	236.95	320.74	487.40	33.78	325.00	0.274	0.249
37	C_2F_4	tetrafluoroethene	116-14-3	100.016	142.00	197.51	306.40	39.44	$\overline{172.00}$	0.266	
38	C_2F_6	hexafluoroethane	76-16-4	138.012	173.05	195.21	293.04	30.39	221.90	0.277	0.257
39	C ₂ HBrClF ₃	1-bromo-1-chloro-2,2,2-trifluoroethane	151-67-7	197.382		323.32	492.20	38.00	302.00	0.280	0.283
40	C ₂ HBrClF ₃	1-bromo-2-chloro-1,1,2-trifluoroethane	354-06-3	197.382		325.70	487.30	34.20	304.00	0.257	0.320
41	C ₂ HClF ₄	1-chloro-1,1,2,2-tetrafluoroethane	354-25-6	136.476	156.15	261.38	400.00	37.60	244.00	0.285	0.260
42	C ₂ HClF ₄	1-chloro-1,2,2,2-tetrafluoroethane (R-124)	2837-89-0	136.476	155.00	261.19	395.60	36.34	243.80	0.269	0.288
43	$C_2HCl_2F_3$	1,1-dichloro-2,2,2-trifluoroethane (R-123)	306-83-2	152.931	166.00	300.81	456.90	36.74	278.05	0.269	0.282
44	$C_2HCl_2F_3$	1,2-dichloro-1,2,2-trifluoroethane (R-123a)	354-23-4	152.931	195	<u>301</u>	461.7		278.00		
45	C_2HF_5	pentafluoroethane	354-33-6	120.022	170.15	225.06	339.17	36.15	211.30	0.271	0.305
46	C_2HF_5O	pentafluorodimethyl ether (E-125)	3822-68-2	136.022	116.00	235.00	354.50	36.31	236.11	0.237	0.326
47	C_2H_2	ethyne (acetylene)	74-86-2	26.038	192.35	188.40	308.30	61.14	112.20	0.268	0.189
48	$C_2H_2F_2$	1,1-difluoroethene	75-38-7	64.035	129.15	187.50	303.20	44.33	153.50	0.270	
49	$C_2H_2F_4$	1,1,1,2-tetrafluoroethane (R-134a)	811-97-2	102.032	172.15	247.04	374.26	40.59	200.80	0.262	0.326
50	$C_2H_2F_4$	1,1,2,2-tetrafluoroethane (R-134)	359-35-3	102.032	184.15	253.10	391.74	46.40	190.40	0.271	0.293
51	$C_2H_3ClF_2$	1-chloro-1,1-difluoroethane	75-68-3	100.495	142.35	264.05	410.30	40.48	231.00	0.267	0.231
52	$C_2H_3Cl_2F$	1,1-dichloro-1-fluoroethane (R-141b)	1717-00-6	116.950	169.60	305.20	477.35	42.50	254.00	0.272	0.225
53	$C_2H_3F_3$	1,1,1-trifluoroethane (R-143a)	420-46-2	84.041	161.85	225.86	346.30	37.92	193.60	0.255	0.259
54	$C_2H_3F_3$	1,1,2-trifluoroethane (R-143)	430-66-0	84.041	189.15	276.85	429.80	52.41	179.20	0.263	0.315
55	C_2H_4	ethene (ethylene)	74-85-1	28.054	103.99	169.42	282.34	50.41	131.10	0.282	0.087
56	$C_2H_4Br_2$	1,2-dibromoethane	106-93-4	187.862	283.05	404.50	582.90	71.50	261.70		
57	$C_2H_4Cl_2$	1,1-dichloroethane	75-34-3	98.959	176.19	330.45	523.00	51.00	236.00	0.277	
58	$C_2H_4Cl_2$	1,2-dichloroethane	107-06-2	98.959	237.46	356.66	561.00	54.00	220.00	0.255	

59	$C_2H_4F_2$	1,1-difluoroethane (R-152a)	75-37-6	66.051	156.15	249.10	386.41	45.16	181.00	0.252	0.276
60	$C_2H_4O_2$	ethanoic acid (acetic acid)	64-19-7	60.053	289.77	391.04	594.45	57.90	171.00	0.200	0.445
61	$C_2H_4O_2$	methyl methanoate (methyl formate)	107-31-3	60.053	174.15	304.90	487.20	60.00	172.00	0.255	
62	C_2H_5Br	bromoethane	74-96-4	108.966	154.55	311.50	503.80	62.30	214.90		
63	C ₂ H ₅ Cl	chloroethane	75-00-3	64.514	134.82	285.42	460.30	53.00	199.00	0.276	
64	C_2H_5F	fluoroethane	353-36-6	48.060	129.95	235.43	375.28	50.27	164.00	0.267	0.217
65	C_2H_6	ethane	74-84-0	30.070	90.35	184.55	305.32	48.72	145.50	0.279	0.099
66	C_2H_6O	ethanol	64-17-5	46.069	159.05	351.80	513.92	61.48	167.00	0.240	0.649
67	C_2H_6O	dimethyl ether	115-10-6	46.069	131.65	248.31	400.10	54.00	170.00	0.276	
68	C_2H_6S	ethanethiol (ethyl mercaptan)	75-08-1	62.136	152.15	308.15	499.00	54.90	207.00	0.274	
69	C_2H_6S	2-thiapropane (dimethylsulfide)	75-18-3	62.136	174.85	310.48	503.00	55.30	201.00	0.266	
70	C_2H_7N	ethanamine (ethyl amine)	75-04-7	45.084	192.15	289.75	456.40	56.30	181.80	0.267	0.276
71	C_2H_7N	N-methylmethanamine (dimethyl amine)	124-40-3	45.084	180.96	280.00	437.20	53.40	180.00	0.264	
72	C_3F_8	octafluoropropane (R-218)	76-19-7	188.020	125.60	236.60	345.10	26.80	299.82	0.280	0.325
73	C_3HF_7	1,1,1,2,3,3,3-heptafluoropropane	431-89-0	170.030		293.00	376.00		293.15		
	,	(R-227ea)									
74	C ₃ H ₂ ClF ₅ O	2-chloro-1,1,2-trifluoroethyl	13838-16-9	184.493		329.98	475.03	29.80			0.430
	5 2 5	1.0 4.1.4 (0)									
		difluoromethyl ether (enflurane)									
75	C ₃ H ₂ ClF ₅ O	1-chloro-2,2,2-trifluoroethyl	26675-46-7	184.493		322.42	467.80	30.46			0.402
75	C ₃ H ₂ ClF ₅ O		26675-46-7	184.493		322.42	467.80	<u>30.46</u>			0.402
75 76		1-chloro-2,2,2-trifluoroethyl	26675-46-7 431-63-0	184.493 152.050		322.42 279.00	<u>467.80</u> 412.50	30.46	268.88		0.402 0.382
	$C_3H_2ClF_5O$ $C_3H_2F_6$ $C_3H_2F_6$	1-chloro-2,2,2-trifluoroethyl difluoromethyl ether (isoflurane)			232.65			<u>30.46</u>	268.88 277.21		
76	$C_3H_2F_6$	1-chloro-2,2,2-trifluoroethyl difluoromethyl ether (isoflurane) 1,1,1,2,3,3-hexafluoropropane (R-236ea)	431-63-0	152.050	232.65		412.50	30.46 31.48		0.267	0.382
76 77	$C_3H_2F_6$ $C_3H_2F_6$	1-chloro-2,2,2-trifluoroethyl difluoromethyl ether (isoflurane) 1,1,1,2,3,3-hexafluoropropane (R-236ea) 1,1,1,3,3,3-hexafluoropropane (R-236fa)	431-63-0 690-39-1	152.050 152.050	232.65	279.00	412.50 398.10		277.21	0.267	$\frac{0.382}{0.377}$
76 77 78	$C_{3}H_{2}F_{6}$ $C_{3}H_{2}F_{6}$ $C_{3}H_{3}F_{5}$	1-chloro-2,2,2-trifluoroethyl difluoromethyl ether (isoflurane) 1,1,1,2,3,3-hexafluoropropane (R-236ea) 1,1,1,3,3,3-hexafluoropropane (R-236fa) 1,1,2,2-pentafluoropropane (R-245cb)	431-63-0 690-39-1 1814-88-6	152.050 152.050 134.050	232.65 191.20	279.00	412.50 398.10 380.40		277.21 268.74	0.267	$\frac{0.382}{0.377}$ $\frac{0.382}{0.297}$
76 77 78 79	$C_{3}H_{2}F_{6}$ $C_{3}H_{2}F_{6}$ $C_{3}H_{3}F_{5}$ $C_{3}H_{3}F_{5}$	1-chloro-2,2,2-trifluoroethyl difluoromethyl ether (isoflurane) 1,1,1,2,3,3-hexafluoropropane (R-236ea) 1,1,1,3,3,3-hexafluoropropane (R-245cb) 1,1,1,3,3-pentafluoropropane (R-245fa)	431-63-0 690-39-1 1814-88-6 460-73-1	152.050 152.050 134.050 134.050		279.00 255.10	412.50 398.10 380.40 427.00		277.21 268.74 259.28	0.267	$\begin{array}{c} 0.382 \\ \hline 0.377 \\ \hline 0.297 \\ \hline 0.385 \end{array}$
76 77 78 79 80	$C_3H_2F_6$ $C_3H_2F_6$ $C_3H_3F_5$ $C_3H_3F_5$ $C_3H_3F_5$	1-chloro-2,2,2-trifluoroethyl difluoromethyl ether (isoflurane) 1,1,1,2,3,3-hexafluoropropane (R-236ea) 1,1,1,3,3,3-hexafluoropropane (R-245cb) 1,1,2,2-pentafluoropropane (R-245fa) 1,1,2,2,3-pentafluoropropane (R-245fa) 1,1,2,2,3-pentafluoropropane (R-245ca)	431-63-0 690-39-1 1814-88-6 460-73-1 679-86-7	152.050 152.050 134.050 134.050 134.050		279.00 255.10 301.20	412.50 398.10 380.40 427.00 447.60		277.21 268.74 259.28 256.02	0.267	$\begin{array}{c} 0.382 \\ \hline 0.377 \\ \hline 0.297 \\ \hline 0.385 \end{array}$
76 77 78 79 80	$C_3H_2F_6$ $C_3H_2F_6$ $C_3H_3F_5$ $C_3H_3F_5$ $C_3H_3F_5$	1-chloro-2,2,2-trifluoroethyl difluoromethyl ether (isoflurane) 1,1,1,2,3,3-hexafluoropropane (R-236ea) 1,1,1,3,3,3-hexafluoropropane (R-245cb) 1,1,2,2-pentafluoropropane (R-245cb) 1,1,3,3-pentafluoropropane (R-245fa) 1,1,2,2,3-pentafluoropropane (R-245ca) 2-(difluoromethoxy)-1,1,1-trifluoroethane	431-63-0 690-39-1 1814-88-6 460-73-1 679-86-7	152.050 152.050 134.050 134.050 134.050		279.00 255.10 301.20	412.50 398.10 380.40 427.00 447.60		277.21 268.74 259.28 256.02	0.267 0.237	$\begin{array}{c} 0.382 \\ \hline 0.377 \\ \hline 0.297 \\ \hline 0.385 \end{array}$
76 77 78 79 80 81	C ₃ H ₂ F ₆ C ₃ H ₂ F ₆ C ₃ H ₃ F ₅ C ₃ H ₃ NO	1-chloro-2,2,2-trifluoroethyl difluoromethyl ether (isoflurane) 1,1,1,2,3,3-hexafluoropropane (R-236ea) 1,1,1,3,3,3-hexafluoropropane (R-245cb) 1,1,1,2,2-pentafluoropropane (R-245cb) 1,1,1,3,3-pentafluoropropane (R-245fa) 1,1,2,2,3-pentafluoropropane (R-245ca) 2-(difluoromethoxy)-1,1,1-trifluoroethane (E-245)	431-63-0 690-39-1 1814-88-6 460-73-1 679-86-7 1885-48-9	152.050 152.050 134.050 134.050 134.050 150.050		279.00 255.10 301.20 302.20	412.50 398.10 380.40 427.00 447.60 444.00	31.48	277.21 268.74 259.28 256.02 291.02		0.382 0.377 0.297 0.385 0.353
76 77 78 79 80 81	C ₃ H ₂ F ₆ C ₃ H ₂ F ₆ C ₃ H ₃ F ₅ C ₃ H ₃ F ₅ C ₃ H ₃ F ₅ C ₃ H ₃ F ₅	1-chloro-2,2,2-trifluoroethyl difluoromethyl ether (isoflurane) 1,1,1,2,3,3-hexafluoropropane (R-236ea) 1,1,1,3,3,3-hexafluoropropane (R-245cb) 1,1,1,2,2-pentafluoropropane (R-245cb) 1,1,3,3-pentafluoropropane (R-245fa) 1,1,2,2,3-pentafluoropropane (R-245ca) 2-(difluoromethoxy)-1,1,1-trifluoroethane (E-245) 1,2-oxazole(isoxazole)	431-63-0 690-39-1 1814-88-6 460-73-1 679-86-7 1885-48-9 288-14-2	152.050 152.050 134.050 134.050 134.050 150.050	191.20	279.00 255.10 301.20 302.20 368.61	412.50 398.10 380.40 427.00 447.60 444.00 590.00	31.48 61.00	277.21 268.74 259.28 256.02 291.02 190.94	0.237	0.382 0.377 0.297 0.385 0.353
76 77 78 79 80 81 82 83	C ₃ H ₂ F ₆ C ₃ H ₂ F ₆ C ₃ H ₃ F ₅ C ₃ H ₃ NO C ₃ H ₄ C ₃ H ₄	1-chloro-2,2,2-trifluoroethyl difluoromethyl ether (isoflurane) 1,1,1,2,3,3-hexafluoropropane (R-236ea) 1,1,1,3,3,3-hexafluoropropane (R-245cb) 1,1,1,2,2-pentafluoropropane (R-245fa) 1,1,2,2,3-pentafluoropropane (R-245fa) 1,1,2,2,3-pentafluoropropane (R-245ca) 2-(difluoromethoxy)-1,1,1-trifluoroethane (E-245) 1,2-oxazole(isoxazole) 1-propyne (methyl acetylene)	431-63-0 690-39-1 1814-88-6 460-73-1 679-86-7 1885-48-9 288-14-2 74-99-7	152.050 152.050 134.050 134.050 134.050 150.050 55.058 40.065	191.20 170.50	279.00 255.10 301.20 302.20 368.61 250.12	412.50 398.10 380.40 427.00 447.60 444.00 590.00 402.40	31.48 61.00 56.30	277.21 268.74 259.28 256.02 291.02 190.94 163.50	0.237 0.275	0.382 0.377 0.297 0.385 0.353
76 77 78 79 80 81 82 83 84	C ₃ H ₂ F ₆ C ₃ H ₂ F ₆ C ₃ H ₃ F ₅ C ₃ H ₃ NO C ₃ H ₄	1-chloro-2,2,2-trifluoroethyl difluoromethyl ether (isoflurane) 1,1,1,2,3,3-hexafluoropropane (R-236ea) 1,1,1,3,3,3-hexafluoropropane (R-245cb) 1,1,1,2,2-pentafluoropropane (R-245cb) 1,1,1,3,3-pentafluoropropane (R-245fa) 1,1,2,2,3-pentafluoropropane (R-245ca) 2-(difluoromethoxy)-1,1,1-trifluoroethane (E-245) 1,2-oxazole(isoxazole) 1-propyne (methyl acetylene) 1,2-propadiene	431-63-0 690-39-1 1814-88-6 460-73-1 679-86-7 1885-48-9 288-14-2 74-99-7 463-49-0	152.050 152.050 134.050 134.050 134.050 150.050 55.058 40.065 40.065	191.20 170.50 136.85	279.00 255.10 301.20 302.20 368.61 250.12 238.77	412.50 398.10 380.40 427.00 447.60 444.00 590.00 402.40 394.00	31.48 61.00 56.30 52.50	277.21 268.74 259.28 256.02 291.02 190.94 163.50 173.90	0.237 0.275 0.279	0.382 0.377 0.297 0.385 0.353 0.258
76 77 78 79 80 81 82 83 84 85	C ₃ H ₂ F ₆ C ₃ H ₂ F ₆ C ₃ H ₃ F ₅ C ₃ H ₃ NO C ₃ H ₄ C ₃ H ₄ C ₃ H ₆	1-chloro-2,2,2-trifluoroethyl difluoromethyl ether (isoflurane) 1,1,1,2,3,3-hexafluoropropane (R-236ea) 1,1,1,3,3,3-hexafluoropropane (R-245cb) 1,1,1,2,2-pentafluoropropane (R-245cb) 1,1,1,3,3-pentafluoropropane (R-245fa) 1,1,2,2,3-pentafluoropropane (R-245ca) 2-(difluoromethoxy)-1,1,1-trifluoroethane (E-245) 1,2-oxazole(isoxazole) 1-propyne (methyl acetylene) 1,2-propadiene propene (propylene)	431-63-0 690-39-1 1814-88-6 460-73-1 679-86-7 1885-48-9 288-14-2 74-99-7 463-49-0 115-07-1	152.050 152.050 134.050 134.050 134.050 150.050 55.058 40.065 40.065 42.081	191.20 170.50 136.85 87.89	279.00 255.10 301.20 302.20 368.61 250.12 238.77 225.46	412.50 398.10 380.40 427.00 447.60 444.00 590.00 402.40 394.00 364.90	31.48 61.00 56.30 52.50 46.00	277.21 268.74 259.28 256.02 291.02 190.94 163.50 173.90 184.60	0.237 0.275 0.279 0.280	0.382 0.377 0.297 0.385 0.353 0.258 0.122 0.142
76 77 78 79 80 81 82 83 84 85 86	C ₃ H ₂ F ₆ C ₃ H ₂ F ₆ C ₃ H ₃ F ₅ C ₃ H ₄ C ₃ H ₄ C ₃ H ₆ C ₃ H ₆	1-chloro-2,2,2-trifluoroethyl difluoromethyl ether (isoflurane) 1,1,1,2,3,3-hexafluoropropane (R-236ea) 1,1,1,3,3,3-hexafluoropropane (R-245cb) 1,1,1,3,3-pentafluoropropane (R-245fa) 1,1,2,2,3-pentafluoropropane (R-245fa) 1,1,2,2,3-pentafluoropropane (R-245ca) 2-(difluoromethoxy)-1,1,1-trifluoroethane (E-245) 1,2-oxazole(isoxazole) 1-propyne (methyl acetylene) 1,2-propadiene propene (propylene) cyclopropane	431-63-0 690-39-1 1814-88-6 460-73-1 679-86-7 1885-48-9 288-14-2 74-99-7 463-49-0 115-07-1 75-19-4	152.050 152.050 134.050 134.050 134.050 150.050 55.058 40.065 40.065 42.081 42.081	191.20 170.50 136.85 87.89 145.73	279.00 255.10 301.20 302.20 368.61 250.12 238.77 225.46 240.34	412.50 398.10 380.40 427.00 447.60 444.00 590.00 402.40 394.00 364.90 398.25	31.48 61.00 56.30 52.50 46.00 55.75	277.21 268.74 259.28 256.02 291.02 190.94 163.50 173.90 184.60 162.80	0.237 0.275 0.279 0.280 0.274	0.382 0.377 0.297 0.385 0.353 0.258 0.122 0.142 0.130
76 77 78 79 80 81 82 83 84 85 86 87	C ₃ H ₂ F ₆ C ₃ H ₂ F ₆ C ₃ H ₃ F ₅ C ₃ H ₄ C ₃ H ₄ C ₃ H ₆ C ₃ H ₆ C ₃ H ₆ C ₃ H ₆ Cl ₂	1-chloro-2,2,2-trifluoroethyl difluoromethyl ether (isoflurane) 1,1,1,2,3,3-hexafluoropropane (R-236ea) 1,1,1,3,3,3-hexafluoropropane (R-245cb) 1,1,1,3,3-pentafluoropropane (R-245fa) 1,1,2,2,3-pentafluoropropane (R-245fa) 1,1,2,2,3-pentafluoropropane (R-245ca) 2-(difluoromethoxy)-1,1,1-trifluoroethane (E-245) 1,2-oxazole(isoxazole) 1-propyne (methyl acetylene) 1,2-propadiene propene (propylene) cyclopropane 1,2-dichloropropane	431-63-0 690-39-1 1814-88-6 460-73-1 679-86-7 1885-48-9 288-14-2 74-99-7 463-49-0 115-07-1 75-19-4 78-87-5	152.050 152.050 134.050 134.050 134.050 150.050 55.058 40.065 42.081 42.081 112.187	191.20 170.50 136.85 87.89 145.73	279.00 255.10 301.20 302.20 368.61 250.12 238.77 225.46 240.34 369.43	412.50 398.10 380.40 427.00 447.60 444.00 590.00 402.40 394.00 364.90 398.25 578.00	31.48 61.00 56.30 52.50 46.00 55.75	277.21 268.74 259.28 256.02 291.02 190.94 163.50 173.90 184.60 162.80	0.237 0.275 0.279 0.280 0.274	0.382 0.377 0.297 0.385 0.353 0.258 0.122 0.142 0.130

Section A Basic Constants I (Continued)

No.	Formula	Name	CAS #	Mol. Wt.	Tfn K	Tb, K	Tc, K	Pc bar	Vc,	Zc = PcVc/RTc	Omega
90	$C_3H_6O_2$	propanoic acid	79-09-4	74.079	252.31		604.00	45.30	233.00	0.207	0.539
91	$C_3H_6O_2$	methyl ethanoate (methyl acetate)	79-20-9	74.079	175.15	330.09	506.80	46.90	228.00	0.254	
92	$C_3H_6O_2$	ethyl methanoate (ethyl formate)	109-94-4	74.079	193.55	327.47	508.50	47.40	229.00	0.257	0.282
93	$C_3H_6O_3$	dimethylcarbonate	616-38-6	90.084		363.24	557.00	48.00	251.63	0.261	0.336
94	C_3H_7Cl	1-chloropropane	540-54-5	78.541	150.35	319.67	503.10	45.80	254.00	0.278	
95	C_3H_8	propane	74-98-6	44.097	91.45	231.02	369.83	42.48	200.00	0.276	0.152
96	C_3H_8O	1-propanol	71-23-8	60.096	$1\overline{47.00}$	370.93	536.78	51.75	219.00	0.254	0.629
97	C_3H_8O	2-propanol	67-63-0	60.096	183.65	355.39	508.30	47.62	220.00	0.248	0.665
98	C_3H_8O	methyl ethyl ether	540-67-0	60.096	160.00	280.50	437.80	44.00	221.00	0.267	0.236
99	C_3H_8S	2-thiabutane (methyl ethyl sulfide)	624-89-5	76.163	167.24	339.80	533.00	42.60	260.00	0.250	
100	C_3H_9N	1-propanamine (propyl amine)	107-10-8	59.111	188.35	320.38	497.00	48.00	235.00	0.273	0.283
101	C_3H_9N	2-propanamine (methyl ethyl amine)	75-31-0	59.111	177.95	304.93	471.80	45.40	221.00	0.256	0.277
102	C_3H_9N	N,N-dimethylmethanamine (trimethyl amine)	75-50-3	59.111	155.85	276.02	433.30	40.75	254.00	0.266	
103	C ₃ H ₉ NO	2-ethoxymethanamine (2-methylaminoethanol)	109-83-1	75.112		<u>432.39</u>	630.00	53.00	259.01	0.262	0.604
104	C_4F_8	octafluorocyclobutane	115-25-3	200.031	232.96	267.17	388.37	27.78	324.80	0.279	
105	C_4F_{10}	decafluoro-2-methylpropane	354-92-7	238.028		273.15	395.40	24.20	378.00	0.278	
106	C_4H_4O	furan	110-00-9	68.075	187.54	304.44	490.15	55.00	218.00	0.281	
107	C_4H_4S	thiophene	110-02-1	84.142	233.75	357.31	580.00	56.60	219.00	0.252	
108	C_4H_5N	pyrrole	109-97-7	67.090	249.74	403.00	639.70	63.30	200.00	0.238	
109	C_4H_6	1-butyne	107-00-6	54.092	147.29	281.21	440.00	46.00	208.00	0.262	0.245
110	C_4H_6	1,3-butadiene	106-99-0	54.092	164.24	268.62	425.00	43.20	221.00	0.270	0.195
111	$C_4H_6O_3$	acetic anhydride	108-24-7	102.090	199.00	412.69	606.00	40.00			0.456
112	C_4H_8	cyclobutane	287-23-0	56.108	182.42	285.64	460.00	49.90	218.00	0.275	0.185
113	C_4H_8	1-butene	106-98-9	56.108	87.79	266.92	419.50	40.20	240.80	0.278	0.194
114	C_4H_8	trans-2-butene	624-64-6	56.108	167.58	274.03	428.60	41.00	237.70	0.276	0.218
115	C_4H_8	cis-2-butene	590-18-1	56.108	134.25	276.87	435.50	42.10	233.80	0.269	0.203
116	C_4H_8	2-methylpropene	115-11-7	56.108	132.81	266.24	417.90	40.00	238.80	0.275	0.199
117	$C_4^{\dagger}H_8^{\circ}O$	butanone (methyl ethyl ketone)	78-93-3	72.107	186.51	352.71	536.80	42.10	267.00	0.252	0.322

11	8 C ₄ H ₈ O	tetrahydrofuran	109-99-9	72.107	164.61	339.12	540.20	51.90	224.00	0.259	
11	$9 C_4H_8O_2$	butanoic acid	107-92-6	88.106	267.97	436.87	624.00	40.30	292.00	0.227	0.600
12	$0 C_4H_8O_2$	2-methylpropanoic acid	79-31-2	88.106	227.05	427.57	605.00	37.00	290.00	0.213	0.618
12	$1 C_4H_8O_2$	1,3-dioxane	505-22-6	88.106	228.15	379.20	588.00	45.80	257.00	0.241	
12	$C_4H_8O_2$	1,4-dioxane	123-91-1	88.106	284.15	374.50	587.00	51.70	238.00	0.255	
12	$C_4H_8O_2$	methyl propanoate	554-12-1	88.106	185.65	352.60	530.60	40.00	282.00	0.256	0.349
12	$4 C_4H_8O_2$	ethyl ethanoate (ethyl acetate)	141-78-6	88.106	189.55	350.21	523.20	38.30	286.00	0.252	0.361
12	$5 C_4H_8O_2$	propyl methanoate (propyl formate)	110-74-7	88.106	180.25	354.69	538.00	40.60	285.00	0.259	0.320
12	6 C ₄ H ₉ Cl	2-chlorobutane	78-86-4	92.568	141.85	341.24	520.60	36.80	312.00	0.265	0.267
12	$7 C_4H_{10}$	butane	106-97-8	58.123	134.79	272.66	425.12	37.96	255.00	0.274	0.200
12	$8 C_4H_{10}$	2-methylpropane (isobutane)	75-28-5	58.123	113.54	261.34	407.85	36.40	262.70	0.278	0.186
12	$9 C_4 H_{10} N_2$	piperazine	110-85-0	86.136	384.6	421.772	661.00		267.50		
13		1-butanol	71-36-3	74.123	183.35	390.88	563.05	44.23	275.00	0.260	0.590
13		2-methyl-1-propanol (isobutanol)	78-83-1	74.123		381.04	547.78	43.00	273.00	0.258	0.590
13	$C_{4}H_{10}O$	2-methyl-2-propanol (tert-butanol)	75-65-0	74.123	298.55	355.49	506.21	39.73	275.00	0.260	0.613
13	$3 C_4 H_{10} O$	2-butanol (sec-butanol)	78-92-2	74.123	158.50	372.66	536.05	41.79	269.00	0.252	0.574
13	$4 C_4 H_{10} O$	diethyl ether	60-29-7	74.123	156.86	307.59	466.70	36.40	280.00	0.263	0.281
13	4 10 2	1,2-dimethoxyethane	110-71-4	90.126		<i>358.15</i>	537.00		270.64		
13	$6 C_4 H_{10} O_2$	1,2-butanediol	26171-83-5	90.126		469.58	680.00		303.05		
13	4 10 2	1,3-butanediol	107-88-0	136.154		481.38	676.00		305.00		
13	$8 C_4H_{10}S$	3-thiapentane (diethyl sulfide)	352-93-2	90.189		365.25	557.00	39.60	318.00	0.269	0.295
13	$9 C_4H_{11}N$	1-butanamine (butyl amine)	109-73-9	73.138		349.44	526.80	40.40	290.00	0.267	0.338
14	4 11	N-ethylethanamine (diethyl amine)	109-89-7	73.138		328.60	496.60	37.10	301.00	0.270	
14	$1 C_4H_{11}N$	2-methyl-1-propanamine (isobutyl	78-81-9	73.138	188.55	340.81	528.50	40.20	286.00	0.262	0.220
		amine)									
14	J 12	dodecafluoropentane	678-26-2	288.036		302.35	422.00	20.40	383.10	0.272	0.415
14	2 2	pyridine	110-86-1	79.101		388.37	620.00	56.70	254.00	0.267	0.242
14	5 0	2-methylfuran	534-22-5	82.102		337.87	527.85	47.20	246.40	0.255	0.278
14	2 0	1-pentyne	627-19-0	68.119		313.38	470.00	41.70	277.00	0.296	0.394
14	2 0	cyclopentene	142-29-0	68.119		317.35	506.50	48.00	245.00	0.279	
14		cyclopentanone	120-92-3	84.118		403.72	624.50	46.00	268.00	0.237	0.288
14	- 5 10	cyclopentane	287-92-3	70.134		322.38	511.60	45.08	260.00	0.276	
14	- 3 10	1-pentene	109-67-1	70.134		303.11	464.80	35.60	298.40	0.275	0.237
15		cis-2-pentene	627-20-3	70.134		310.07	475.00	36.90	302.10	0.273	0.253
15	5 10	2-methyl-2-butene	513-35-9	70.134	,	311.70	470.00	38.60	292.00	0.288	0.339
15	C_5H_{10}	3-methyl-1-butene	563-45-1	70.134	104.65	293.21	452.70	35.50	304.90	0.288	0.211

Section A Basic Constants I (Continued)

			G.1.G.11						Vc,	Zc =	
No.	Formula	Name	CAS #	Mol. Wt.	Tfp, K	Tb, K	Tc, K	Pc, bar	cm ³ /mol	PcVc/RTc	Omega
153	$C_5H_{10}O$	cyclopentanol	96-41-3	86.134		413.49	619.50	49.00			0.420
154	$C_5H_{10}O$	2-pentanone (methyl propyl ketone)	107-87-9	86.134	196.34	375.39	561.10	36.90	301.00	0.238	0.346
155	$C_5H_{10}O$	3-pentanone (diethyl ketone)	96-22-0	86.134	234.20	375.14	561.50	37.30	336.00	0.268	0.342
156		3-methyl-2-butanone (methyl isopropyl ketone)	563-80-4	86.134	181.15	367.55	567.70	36.20	310.00	0.238	0.216
157	$C_5H_{10}O$	2-methyltetrahydrofuran	96-47-9	86.134	137.04	353.37	537.00	37.60	267.00	0.225	0.292
158	$C_5H_{10}O_2$	pentanoic acid	109-52-4	102.133	239.45	459.31	643.00	35.80	336.20		0.670
159	$C_5H_{10}O_2$	3-methylbutanoic acid	503-74-2	102.133	243.85	449.68	629.00	34.00			0.651
160	$C_5H_{10}O_2$	methyl butanoate	623-42-7	102.133	187.35	375.90	554.40	34.80	340.00	0.257	0.381
161	$C_5H_{10}O_2$	ethyl propanoate	105-37-3	102.133	199.25	372.18	546.00	33.60	345.00	0.255	0.390
162	$C_5H_{10}O_2$	methyl 2-methylpropanoate	547-63-7	102.133	188.45	365.45	540.80	34.30	339.00	0.259	
163	$C_5H_{10}O_2$	propyl ethanoate (propyl acetate)	109-60-4	102.133	178.15	374.65	549.40	33.30	345.00	0.252	0.389
164	$C_5H_{10}O_2$	2-methylpropyl methanoate (isobutyl	542-55-2	102.133	177.35	371.22	551.00	38.80	352.00	0.298	0.400
		formate)									
165	$C_5H_{11}Cl$	1-chloropentane	543-59-9	106.595	174.15	381.54	552.00				
166	C_5H_{12}	pentane	109-66-0	72.150	143.43	309.22	469.70	33.70	311.00	0.268	0.252
167	C_5H_{12}	2-methylbutane	78-78-4	72.150	113.26	300.99	460.39	33.81	308.30	0.272	0.229
168	C_5H_{12}	2,2-dimethylpropane (neopentane)	463-82-1	72.150	256.58	282.65	433.75	31.99	303.20	0.269	0.197
169	$C_5H_{12}O$	1-pentanol	71-41-0	88.150	194.25	411.16	588.15	39.09	326.00	0.262	0.579
170	$C_5H_{12}O$	2-pentanol	6032-29-7	88.150		392.30	560.30	36.75	329.00	0.260	0.561
171	$C_5H_{12}O$	2-methyl-1-butanol	137-32-6	88.150		403.79	575.40	39.40			0.605
172	$C_5H_{12}O$	2-methyl-2-butanol	75-85-4	88.150	264.40	375.15	545.00	37.90	323.00	0.270	0.478
173	$C_5H_{12}O$	3-methyl-1-butanol	123-51-3	88.150	155.95	403.69	579.40	39.10	325.00	0.264	0.559
174	$C_5H_{12}O$	3-methyl-2-butanol	598-75-4	88.150		385.20	556.10	38.70			
175	$C_5H_{12}O$	ethyl propyl ether	628-32-0	88.150	146.45	337.01	500.60	32.50	339.00	0.265	0.328
176	$C_5H_{12}S$	3-methyl-1-butanethiol (isopentyl mercaptan)	541-31-1	104.216	139.64	391.50	604.00	35.00	364.00	0.254	0.191
177	C_6ClF_5	chloropentafluorobenzene	344-07-0	202.511		391.11	570.81	32.37	376.00	0.256	0.400
178	C_6F_6	hexafluorobenzene	392-56-3	186.056		353.40	516.73	32.75	335.00	0.255	0.396
179	C_6F_{12}	dodecafluorocyclohexane	355-68-0	300.047	335.65		457.29	22.37	459.00	0.270	

180	C_6F_{14}	tetradecafluorohexane	355-42-0	338.044	186.05	329.75	448.70	18.70	573.20	0.274	0.513
181	C_6F_{14}	tetradecafluoro-2-methylpentane	355-04-4	338.044		330.75	453.00	18.20	550.00	0.266	
182	$C_{6}F_{14}$	tetradecafluoro-3-methylpentane	865-71-4	338.044	158.15	331.55	450.00	16.90	531.00	0.240	
183	C_6F_{14}	tetradecafluoro-2,3-dimethyl butane	354-96-1	338.044	258.15	332.95	463.00	18.70	528.00	0.255	
184	C_6HF_5	pentafluorobenzene	363-72-4	168.066	225.85	358.89	530.97	35.37	324.00	0.260	0.374
185	$C_6H_2F_4$	1,2,4,5-tetrafluorobenzene	327-54-8	150.076		363.50	543.35	37.99	<u> </u>		0.357
186	C_6H_5Cl	chlorobenzene	108-90-7	112.558	227.90	404.91	632.40	45.20	308.00	0.265	0.251
187	C_6H_6	benzene	71-43-2	78.114	278.68	353.24	562.05	48.95	256.00	0.268	0.210
188	C_6H_6O	phenol	108-95-2	94.113	314.05	455.04	694.25	61.30	229.00	0.243	0.442
189	C_6H_7N	benzeneamine (aniline)	62-53-3	93.128	266.85	457.17	699.00	53.10	273.90	0.256	0.380
190	C_6H_7N	2-methylpyridine (2-picoline)	109-06-8	93.128	206.45	402.50	621.00	46.00	335.00	0.260	0.299
191	C_6H_7N	3-methylpyridine (3-picoline)	108-99-6	93.128	254.96	417.28	645.00	44.80	288.00	0.241	0.279
192	C_6H_7N	4-methylpyridine (4-picoline)	108-89-4	93.128	276.80	418.49	645.70	46.60	325.62	0.253	0.305
193	C_6H_8O	2-cyclohexen-1-one	930-68-7	96.131		445.49	685.00	45.30	305.18	0.243	0.308
194	C_6H_{10}	cyclohexene	110-83-8	82.145	169.67	356.12	560.40		296.88		
195	$C_6H_{10}O$	cyclohexanone	108-94-1	98.144		428.59	653.00	40.00			0.299
196	$C_6H_{10}O$	4-methyl-3-penten-2-one (mesityl oxide)	141-79-7	98.147		402.86	605.00		353.43		
197	C_6H_{12}	cyclohexane	110-82-7	84.161	279.69	353.93	553.50	40.73	308.00	0.273	0.211
198	C_6H_{12}	methylcyclopentane	96-37-7	84.161	130.72	344.98	532.79	37.84	319.00	0.272	0.227
199	C_6H_{12}	1-hexene	592-41-6	84.161	133.34	336.63	504.00	31.43	355.10	0.266	0.281
200	C_6H_{12}	4-methylpent-1-ene	691-37-2	84.161	119.5	326.82	495.00	32.90	357.9	0.286	0.257
201	$C_6H_{12}O$	cyclohexanol	108-93-0	100.161	297.65	433.94	648.00	40.75	333.88	0.253	0.366
202	$C_6H_{12}O$	2-hexanone (methyl butyl ketone)	591-78-6	100.161	217.40	400.70	587.00	33.20	378.00	0.254	0.393
203	$C_6H_{12}O$	3-hexanone (ethyl propyl ketone)	589-38-8	100.161	217.50	396.65	582.80	33.20	378.00	0.256	0.380
204	$C_6H_{12}O$	4-methyl-2-pentanone (methyl isobutyl	108-10-1	100.161	189.15	389.15	574.60	32.70	340.60	0.256	0.351
		ketone)									
205	$C_6H_{12}O$	butylvinylether	111-34-2	100.163		367.13	540.50	<u>32.00</u>	383.62	0.273	0.358
206	$C_6H_{12}O_2$	hexanoic acid	142-62-1	116.160		478.38	<u>662.00</u>	<u>32.00</u>	377.20		0.694
207	$C_6H_{12}O_2$	methyl pentanoate	624-24-8	116.160		400.55	567.00	31.90	416.30	0.267	
208	$C_6H_{12}O_2$	ethyl butanoate	105-54-4	116.160	175.15	394.69	566.00	30.60	421.00	0.274	0.463
209	$C_6H_{12}O_2$	propyl propanoate	106-36-5	116.160	197.25	395.64	578.00	30.90	394.00	0.253	0.373
210	$C_6H_{12}O_2$	ethyl 2-methylpropanoate	97-62-1	116.160	185.00	383.00	553.00	30.70	421.00	0.281	
211	$C_6H_{12}O_2$	butyl ethanoate (butyl acetate)	123-86-4	116.160	199.65	399.12	579.00	30.90	412.80	0.253	0.407
212	$C_6H_{12}O_2$	2-methylpropyl ethanoate (isobutyl acetate)	110-19-0	116.160	174.25	389.72	561.00	31.60	413.00	0.266	0.456

Section A Basic Constants I (Continued)

No.	Formula	Name	CAS #	Mol. Wt.	Tfn K	Tb, K	Tc, K	Pc har	Vc,	Zc = PcVc/RTc	Omega
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213	$C_6H_{12}O_2$	pentyl methanoate (pentyl formate)	638-49-3	116.160	199.65	403.55	576.00	30.90	412.00	0.254	
214	$C_6H_{12}O_2$	3-methylbutyl methanoate (isopentyl formate)	110-45-2	116.160	179.65	397.28	578.00	31.30	411.40	0.252	0.400
215	$C_6H_{12}O_3$	2-ethoxyethylacetate	111-15-9	132.165		429.74	610.60	31.66	443.5	0.277	0.523
216	C_6H_{14}	hexane	110-54-3	86.177	177.84	341.88	507.60	30.25	368.00	0.264	0.300
217	C_6H_{14}	2-methylpentane	107-83-5	86.177	119.48	333.40	497.50	30.10	366.70	0.267	0.278
218	$C_{6}H_{14}$	3-methylpentane	96-14-0	86.177	110.26	336.40	504.40	31.20	366.70	0.273	0.273
219	C_6H_{14}	2,2-dimethylbutane	75-83-2	86.177	173.33	322.87	488.70	30.80	359.10	0.272	0.233
220	C_6H_{14}	2,3-dimethylbutane	79-29-8	86.177	144.35	331.12	499.90	31.30	357.60	0.269	0.248
221	$C_6H_{14}O$	1-hexanol	111-27-3	102.177	229.20	430.44	611.40	35.10	381.00	0.263	0.573
222	$C_6H_{14}O$	2-hexanol	626-93-7	102.177	223.00	413.02	586.20	33.80	383.00	0.266	0.562
223	$C_6H_{14}O$	3-hexanol	623-37-0	102.177		408.94	582.40	33.60	383.00	0.266	0.539
224	$C_6H_{14}O$	2-methyl-1-pentanol	105-30-6	102.177		420.76	604.40	34.50			0.498
225	$C_6H_{14}O$	2-methyl-2-pentanol	590-36-3	102.177	171.15	394.51	559.50	34.70	380.00	0.283	0.573
226	$C_6H_{14}O$	2-methyl-3-pentanol	565-67-3	102.177		401.20	576.00	34.60			
227	$C_6H_{14}O$	4-methyl-1-pentanol	626-89-1	102.177		424.93	603.50	36.30	380.00	0.275	0.588
228	$C_6H_{14}O$	4-methyl-2-pentanol	108-11-2	102.177		404.86	574.50	33.10	380.00	0.263	0.552
229	$C_6H_{15}N$	N,N-diethylethanamine (triethyl amine)	121-44-8	101.192	158.45	362.90	535.00	30.00	390.00	0.263	
230	$C_6H_{15}N$	N-propyl-1-propanamine (dipropyl amine)	142-84-7	101.192	210.15	382.30	550.00	31.40	402.00	0.276	
231	$C_{7}F_{14}$	tetradecafluoromethylcyclohexane	355-02-2	350.055	228.45	349.45	485.91	20.19	570.00	0.285	
232	C_7F_{16}	hexadecafluoroheptane	335-57-9	388.052	195.15	355.59	475.00	16.50	664.30	0.248	0.561
233	$C_7H_3F_5$	pentafluorotoluene	771-56-2	182.093	243.35	390.65	566.52	31.24	384.00	0.255	0.415
234	C_7H_8	toluene	108-88-3	92.141	178.16	383.79	591.75	41.08	316.00	0.264	0.264
235	C_7H_8O	benzyl alcohol	100-51-6	108.140	257.80	478.46	715.00	43.00			0.390
236	C_7H_8O	2-methylphenol (o-cresol)	95-48-7	108.140	302.95	464.17	697.60	50.00	282.00	0.243	0.436
237	C_7H_8O	3-methylphenol (m-cresol)	108-39-4	108.140	284.95	475.38	705.70	45.60	312.00	0.241	0.452
238	C_7H_8O	4-methylphenol (p-cresol)	106-44-5	108.140	307.89	475.12	704.50	51.50	277.00	0.244	0.510
239	C_7H_9N	2,3-dimethylpyridine (2,3 lutidine)	583-61-9	107.155	249.55	434.30	655.40	40.90	367.00	0.275	0.345
240	C_7H_9N	2,4-dimethylpyridine (2,4 lutidine)	108-47-4	107.155	209.23	431.55	647.00	38.70	367.00	0.264	0.351

241	C_7H_9N	2,5-dimethylpyridine (2,5 lutidine)	589-93-5	107.155	257.65	430.14	644.20	39.80	367.00	0.273	0.369
242	C_7H_9N	2,6-dimethylpyridine (2,6 lutidine)	108-48-5	107.155	267.03	416.91	623.80	39.80	367.00	0.282	0.373
243	C_7H_9N	3,4-dimethylpyridine (3,4 lutidine)	583-58-4	107.155	262.15	452.29	683.80	40.90	367.00	0.264	0.337
244	C_7H_9N	3,5-dimethylpyridine (3,5 lutidine)	591-22-0	107.155	266.65	445.06	667.20	38.70	367.00	0.256	0.351
245	$C_7H_{12}O_2$	butyl-2-propenoate(butylacrylate)	141-32-2	128.175		419.77	644.00		427.54		0.312
246	C_7H_{14}	cycloheptane	291-64-5	98.188	265.15	391.95	604.30	38.40	359.00	0.274	0.242
247	C_7H_{14}	methylcyclohexane	108-87-2	98.188	146.56	374.09	572.19	34.71	368.00	0.268	0.235
248	C_7H_{14}	ethylcyclopentane	1640-89-7	98.188	134.70	376.59	569.50	33.97	375.00	0.269	0.270
249	$C_{7}H_{14}$	cis-1,3-dimethylcyclopentane	2532-58-3	98.188	139.45	364.71	551.00	34.00	363.30	0.277	0.276
250	C_7H_{14}	trans-1,3-dimethylcyclopentane	1759-58-6	98.188	139.18	363.90	553.00	34.00	363.30	0.276	0.253
251	C_7H_{14}	1-heptene	592-76-7	98.188	153.45	366.79	537.30	29.20	409.00	0.267	0.343
252	$C_7H_{14}O_2$	heptanoic acid	111-14-8	130.187		495.35	679.00	29.00	429.70		0.712
253	$C_7H_{14}O_2$	ethyl pentanoate	539-82-2	130.187	181.95	419.25	570.00	27.80	449.00	0.263	
254	$C_7H_{14}O_2$	ethyl 3-methylbutanoate	108-64-5	130.187	173.85	407.45	588.00	27.30	447.00	0.250	
255	$C_7H_{14}O_2$	propyl butanoate	105-66-8	130.187	177.95	416.50	600.00	27.80	449.00	0.250	0.399
256	$C_7H_{14}O_2$	2-methylpropyl propanoate	540-42-1	130.187	201.75	409.75	592.00	27.30	447.00	0.248	0.375
257	$C_7H_{14}O_2$	propyl 2-methylpropanoate	644-49-5	130.187		408.55	589.00	27.30	447.00	0.249	
258	$C_7H_{14}O_2$	3-methylbutyl ethanoate (isopentyl	123-92-2	130.187	194.65	415.20	599.00	28.10	442.00	0.249	
		acetate)									
259	C_7H_{16}	heptane	142-82-5	100.204	182.59	371.57	540.20	27.40	428.00	0.261	0.350
260	C_7H_{16}	2-methylhexane	591-76-4	100.204	154.89	363.18	530.10	27.30	421.00	0.261	0.331
261	C_7H_{16}	3-methylhexane	589-34-4	100.204	149.35	365.00	535.20	28.10	404.00	0.255	0.323
262	C_7H_{16}	3-ethylpentane	617-78-7	100.204	154.57	366.63	540.50	28.90	415.80	0.267	0.311
263	C_7H_{16}	2,2-dimethylpentane	590-35-2	100.204	149.37	352.32	520.40	27.70	415.80	0.266	0.287
264	C_7H_{16}	2,3-dimethylpentane	565-59-3	100.204	82.60	362.91	537.30	29.10	393.00	0.256	0.297
265	C_7H_{16}	2,4-dimethylpentane	108-08-7	100.204	153.94	353.62	519.70	27.40	417.50	0.265	0.304
266	C_7H_{16}	3,3-dimethylpentane	562-49-2	100.204	138.25	359.19	536.30	29.50	414.10	0.274	0.269
267	C_7H_{16}	2,2,3-trimethylbutane	464-06-2	100.204	248.15	354.01	531.10	29.50	397.60	0.266	0.250
268	$C_7H_{16}O$	1-heptanol	111-70-6	116.203	239.20	449.81	631.90	31.50	435.00	0.261	0.588
269	$C_7H_{16}O$	2-heptanol	543-49-7	116.203		434.20	608.30	30.21	442.00		
270	$C_7H_{16}O$	4-heptanol	589-55-9	116.203			602.60		432.00		
271	C_8F_{18}	octadecafluorooctane	307-34-6	438.059		379.05	502.20	16.60	534.40	0.281	0.619
272	C_8H_8O	methylphenylketone(acetophenone)	98-86-2	120.153		475.26	713.00	40.30	380.23	0.258	0.361
273	C_8H_{10}	ethylbenzene	100-41-4	106.167	178.18	409.36	617.15	36.09	374.00	0.263	0.304
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Section A Basic Constants I (Continued)

No.	Formula	Name	CAS #	Mol. Wt.	Tfn V	Tb, K	Tc, K	Do hor	Vc,	Zc = PcVc/RTc	Omega
274	C_8H_{10}	1,2-dimethylbenzene (o-xylene)	95-47-6	106.167		417.59	630.30	37.32	370.00	0.263	0.312
275	C_8H_{10}	1,3-dimethylbenzene (m-xylene)	108-38-3	106.167	225.28		617.00	35.41	375.00	0.259	0.327
276	C_8H_{10}	1,4-dimethylbenzene (p-xylene)	106-42-3	106.167	286.41		616.20	35.11	378.00	0.259	0.322
277	$C_8H_{10}O$	2-ethylphenol	90-00-6	122.167	269.84		703.00	43.00	342.00	0.252	0.475
278	$C_8H_{10}O$	3-ethylphenol	620-17-7	122.167	269.15	491.58	718.80	41.50	342.00	0.237	0.489
279	$C_8H_{10}O$	4-ethyl-phenol	123-07-9	122.167	318.18	491.15	716.40	40.50	342.00	0.233	0.491
280	$C_8H_{10}O$	2,3-dimethylphenol (2,3-xylenol)	526-75-0	122.167	345.95	490.03	722.80	46.30	470.00	0.263	0.496
281	$C_8H_{10}O$	2,4-dimethylphenol (2,4-xylenol)	105-67-9	122.167	297.68	484.09	707.60	42.80	510.00	0.249	0.506
282	$C_8H_{10}O$	2,5-dimethylphenol (2,5-xylenol)	95-87-4	122.167	347.97	484.29	706.90	42.80	470.00	0.249	0.514
283	$C_8H_{10}O$	2,6-dimethylphenol (2,6-xylenol)	576-26-1	122.167	318.75	474.18	701.00	46.30	520.00	0.272	0.489
284	$C_8H_{10}O$	3,4-dimethylphenol (3,4-xylenol)	95-65-8	122.167	333.95	500.11	729.80	42.80	460.00	0.241	0.512
285	$C_8H_{10}O$	3,5-dimethylphenol (3,5-xylenol)	108-68-9	122.167	336.75	494.85	715.60	41.30	610.00	0.237	0.544
286	C_8H_{16}	cyclooctane	292-64-8	112.215	287.95	424.31	647.20	35.70	410.00	0.271	0.254
287	C_8H_{16}	t-1,4-dimethylcyclohexane	2207-04-7	112.215	239.85	429.75	587.70				
288	C_8H_{16}	1-octene	111-66-0	112.215	171.46	394.44	567.00	26.80	468.00	0.266	0.393
289	$C_8H_{16}O_2$	octanoic acid	124-07-2	112.215	289.45	512.01	695.00	26.40			0.734
290	$C_8H_{16}O_2$	2-ethylhexanoic acid	149-57-5	144.218		500.66	675.00		554.68		
291	$C_8H_{16}O_2$	propyl pentanoate	141-06-0	144.214	202.45	440.65	613.00	25.20	504.00	0.249	
292	$C_8H_{16}O_2$	2-methylpropyl butanoate	539-90-2	144.214		430.05	611.00	24.90	502.00	0.246	
293	$C_8H_{16}O_2$	propyl 3-methylbutanoate	557-00-6	144.214		429.05	609.00	24.90	502.00	0.247	
294	$C_8H_{16}O_2$	3-methylbutyl propanoate	105-68-0	144.214		433.35	611.00	25.50	497.00	0.249	
295	$C_8H_{16}O_2$	2-methylpropyl-2-methylpropanoate	97-85-8	144.214	192.45	421.75	602.00	24.50	501.00	0.245	
296	C_8H_{18}	octane	111-65-9	114.231	216.39	398.82	568.70	24.90	492.00	0.259	0.399
297	C_8H_{18}	2-methylheptane	592-27-8	114.231	164.13	390.80	559.60	24.80	488.20	0.260	0.378
298	$C_8^{0}H_{18}^{10}$	3-methylheptane	589-81-1	114.231	152.63	392.08	563.60	25.50	471.10	0.253	0.371
299	C_8^{10}	4-methylheptane	589-53-7	114.231	152.21	390.86	561.70	25.40	476.00	0.259	0.371
300	$C_8^{0}H_{18}^{10}$	3-ethylhexane	619-99-8	114.231		391.69	565.40	26.10	460.50	0.253	0.362
301	$C_8^{0}H_{18}^{10}$	2,2-dimethylhexane	590-73-8	114.231	151.98	379.99	549.80	25.30	478.00	0.265	0.339
302	C_8^{10}	2,3-dimethylhexane	584-94-1	114.231		388.76	563.40	26.30	468.20	0.263	0.347
303	$C_8^{0}H_{18}^{10}$	2,4-dimethylhexane	589-43-5	114.231		382.58	553.50	25.60	472.00	0.263	0.344

304	C_8H_{18}	2,5-dimethylhexane	592-13-2	114.231	181.99	382.26	550.00	24.90	482.00	0.262	0.357
305	C_8H_{18}	3,3-dimethylhexane	563-16-6	114.231	147.06	385.12	562.00	26.50	442.80	0.251	0.320
306	C_8H_{18}	3,4-dimethylhexane	583-48-2	114.231		390.88	568.80	26.90	458.80	0.265	0.338
307	C_8H_{18}	3-ethyl-2-methylpentane	609-26-7	114.231	158.18	388.81	567.00	27.00	445.30	0.254	0.331
308	$C_8^{0}H_{18}^{0}$	3-ethyl-3-methylpentane	1067-08-9	114.231	182.31	391.42	576.50	28.10	455.10	0.267	0.305
309	$C_8^{0}H_{18}^{10}$	2,2,3-trimethylpentane	564-02-3	114.231	160.90	382.99	563.40	27.30	436.00	0.254	0.298
310	$C_8^{0}H_{18}^{10}$	2,2,4-trimethylpentane (isooctane)	540-84-1	114.231	165.80	372.39	543.90	25.70	469.70	0.266	0.304
311	$C_8^{0}H_{18}^{0}$	2,3,3-trimethylpentane	560-21-4	114.231	172.48	387.92	573.50	28.20	455.10	0.269	0.291
312	$C_8^{0}H_{18}^{10}$	2,3,4-trimethylpentane	565-75-3	114.231	163.97	386.62	566.30	27.30	456.20	0.267	0.316
313	C_8H_{18}	2,2,3,3-tetramethylbutane	594-82-1	114.231	373.94	379.44	567.80	28.70	482.00	0.280	0.248
314	$C_8H_{18}O$	1-octanol	111-87-5	130.230	257.65	468.33	652.50	28.60	490.00	0.258	0.594
315	$C_8H_{18}O$	2-octanol	123-96-6	130.230	241.15	453.03	637.00	28.10	493.00	0.262	0.531
316	$C_8H_{18}O$	3-octanol	589-98-0	130.230			628.50		515.00		
317	$C_8H_{18}O$	4-octanol	589-62-8	130.230		440.72	625.10		515.00		
318	$C_8H_{18}O$	2-ethyl-1-hexanol	104-76-7	130.230	203.20	457.77	640.50	27.99			0.558
319	$C_8H_{19}N$	n-octylamine	111-86-4	129.246		451.70	641.00	26.17	517.00	0.254	0.446
320	$C_8H_{19}N$	N-butyl-1-butanamine (dibutyl amine)	111-92-2	129.246	211.15	432.80	602.30	25.70	512.00	0.263	
321	C_9F_{20}	eicosafluorononane	375-96-2	488.067		398.45	523.90	15.60	590.60	0.282	0.635
322	C_9H_7N	quinoline	91-22-5	129.161	257.55	510.30	782.00	45.00	402.00	0.278	0.315
323	C_9H_7N	isoquinoline	119-65-3	129.161	299.62	516.35	803.00	47.00	402.00	0.283	
324	C_9H_{10}	indan	496-11-7	118.178	221.74	451.00	684.90	39.50	393.00	0.273	
325	C_9H_{12}	propylbenzene	103-65-1	120.194	173.65	432.35	638.35	32.00	440.00	0.265	0.345
326	C_9H_{12}	1-methylethylbenzene (cumene)	98-82-8	120.194	177.12	425.52	631.00	32.09	434.70	0.261	0.326
327	C_9H_{12}	1-ethyl-4-methylbenzene	622-96-8	120.194	210.81	435.13	640.20	32.30	440.00	0.259	0.364
328	C_9H_{12}	1,2,3-trimethylbenzene	526-73-8	120.194	247.77	449.23	664.50	34.54	435.00	0.267	0.367
329	C_9H_{12}	1,2,4-trimethylbenzene	95-63-6	120.194	229.35	442.49	649.10	32.32	435.00	0.256	0.377
330	C_9H_{12}	1,3,5-trimethylbenzene (mesitylene)	108-67-8	120.194	228.43	437.90	637.30	31.27	430.00	0.252	0.399
331	C_9H_{18}	1-nonene	124-11-8	126.242	191.80	420.03	594.00	23.30	526.00	0.248	0.411
332	$C_9H_{18}O_2$	nonanoic acid	112-05-0	158.241		527.74	711.00	24.30			0.748
333	$C_9H_{18}O_2$	3-methylbutyl butanoate	106-27-4	158.241		452.09	619.00	23.30	500.00	0.226	0.583
334	C_9H_{20}	nonane	111-84-2	128.258	219.66	423.97	594.60	22.90	555.00	0.257	0.445
335	C_9H_{20}	2-methyloctane	3221-61-2	128.258	192.79	416.44	587.00	23.10	529.00	0.250	0.423
336	C_9H_{20}	2,2-dimethylheptane	1071-26-7	128.258	160.16	405.97	577.80	23.50	525.00	0.257	0.383
337	C_9H_{20}	2,2,5-trimethylhexane	3522-94-9	128.258	167.37	397.24	569.80		519.00		
338	C_9H_{20}	2,2,3,3-tetramethylpentane	7154-79-2	128.258	263.26	413.44	607.60	27.40	478.00	0.269	0.304

Section A Basic Constants I (Continued)

			CAS #	M 1 37/2	TDC IZ	TDI 17	TD 17	D 1	Vc,	Zc =	
No.	Formula	Name	CAS #	Mol. Wt.	11p, K	Tb, K	Tc, K	Pc, bar	cm ³ /mol	PcVc/RTc	Omega
339	C_9H_{20}	2,2,3,4-tetramethylpentane	1186-53-4	128.258	152.06	406.18	592.70	25.30	490.00	0.258	0.301
340	C_9H_{20}	2,2,4,4-tetramethylpentane	1070-87-7	128.258	206.64	395.44	574.60	24.90	504.00		0.314
341	C_9H_{20}	2,3,3,4-tetramethylpentane	16747-38-9	128.258	171.05	414.71	607.10	26.70	493.00	0.261	0.309
342	$C_9H_{20}O$	1-nonanol	143-08-8	144.257	268.15	486.52	668.90	26.30	544.00	0.261	0.633
343	$C_9H_{20}O$	2-nonanol	628-99-9	144.257		466.70	649.60	25.30	575.00	0.269	
344	$C_{10}F_{22}$	docosafluorodecane	307-45-9	538.075		417.35	542.30	14.50	624.20	0.279	
345	$C_{10}H_{8}$	naphthalene	91-20-3	128.174	351.35	491.16	748.40	40.50	407.00	0.265	0.304
346	$C_{10}H_{12}$	1,2,3,4-tetrahydronaphthalene	119-64-2	132.205	237.35	480.75	720.00	36.50	408.00	0.249	
347	$C_{10}H_{12}$	1-methylindan	767-58-8	132.205		463.80	694.10	35.30	448.00	0.274	
348	$C_{10}H_{12}$	2-methylindan	824-63-5	132.205		464.50	695.30	35.30	448.00	0.274	
349	$C_{10}H_{12}$	4-methylindan	824-22-6	132.205		478.60	716.40	35.30	448.00	0.266	
350	$C_{10}H_{12}$	5-methylindan	874-35-1	132.205		475.10	711.20	35.30	448.00	0.267	
351	$C_{10}H_{14}$	butylbenzene	104-51-8	134.221	185.19	456.42	660.50	28.90	497.00	0.262	0.393
352	$C_{10}H_{14}$	2-methylpropylbenzene (isobutylbenzene)	538-93-2	134.221	221.67	445.90	650.00	30.50	480.00	0.271	0.383
353	$C_{10}H_{14}$	1,4-diethylbenzene	105-05-5	134.221	230.30	456.90	657.90	28.03	480.50	0.247	0.403
354	$C_{10}H_{14}$	1-(1-methylethyl)-4-methylbenzene	99-87-6	134.221	205.22	450.26	652.00	28.00	497.00	0.248	0.376
		(p-cymene)									
355	$C_{10}H_{14}$	1,2,4,5-tetramethylbenzene (durene)	95-93-2	134.221	352.45	470.00	676.00	29.00	482.00	0.249	0.423
356	$C_{10}H_{18}$	cis-bicyclo[4.4.0]decane (cis-decalin)	493-01-6	138.253	230.14	468.92	703.60	32.00	480.00	0.265	0.276
357	$C_{10}H_{18}$	trans-bicyclo[4.4.0]decane (trans-decalin)	493-02-7	138.253	242.75	460.42	687.00	32.00	480.00	0.272	0.303
358	$C_{10}H_{20}$	1-decene	872-05-9	140.269	206.89	443.75	617.00	22.20	584.00	0.253	
359	$C_{10}H_{20}O_2$	decanoic acid	334-48-5	172.268	305.15	541.92	<u>726.00</u>	22.30			0.749
360	$C_{10}H_{22}$	decane	124-18-5	142.285	243.49	447.30	617.70	21.10	624.00	0.256	0.490
361	$C_{10}H_{22}$	2,2,5-trimethylheptane	20291-95-6	142.285		423.90	598.90	22.40	569.00	0.256	0.398
362	$C_{10}H_{22}$	3,3,5-trimethylheptane	7154-80-5	142.285		428.83	609.60	23.20	564.00	0.258	0.383
363	$C_{10}H_{22}$	2,2,3,3-tetramethylhexane	13475-81-5	142.285	219.19	433.46	623.00	25.10			0.366
364	$C_{10}H_{22}$	2,2,5,5-tetramethylhexane	1071-81-4	142.285	260.60	410.61	581.40	21.90			0.377
365	$C_{10}H_{22}O$	1-decanol	112-30-1	158.284	280.05	504.25	684.40	23.70	600.00	0.252	0.661
366	$C_{10}H_{24}N_4$	octamethylethenetetramine	996-70-3	200.326		<i>477.44</i>	680.00		646.21		
367	$C_{11}H_{10}$	1-methylnaphthalene	90-12-0	142.200	242.69	517.84	772.00	36.00	462.00	0.259	0.348

368	$C_{11}H_{10}$	2-methylnaphthalene	91-57-6	142.200	307.71	514.20	761.00	35.40	462.00	0.258	0.374
369	$C_{11}H_{24}$	undecane	1120-21-4	156.312	247.57	469.08	639.00	19.80	689.00	0.257	0.537
370	$C_{11}H_{24}O$	1-undecanol	112-42-5	172.311		521.24	705.00	22.40			0.656
371	$C_{12}H_{10}$	1,1'-biphenyl	92-52-4	154.211	342.35	528.23	773.00	33.80	497.00	0.261	0.404
372	$C_{12}H_{12}$	1,6-dimethylnaphthalene	575-43-9	156.227	257.00	539.50	784.00	31.00	517.00	0.246	
373	$C_{12}^{12}H_{12}^{12}$	2,7-dimethylnaphthalene	582-16-1	156.227	368.85	535.00	775.00	32.30			
374		1,3,5-triethylbenzene	102-25-0	162.276		488.93	679.00	24.35	624.14	0.269	0.527
375	$C_{12}H_{20}$	1,3-dimethyltricyclo[3.3.1.1 ^{3,7}]decane	702-79-4	164.292		476.44	708.00		571.45		
	-1220	(1,3-dimethyladamantane)									
376	$C_{12}H_{24}$	1-dodecene	112-41-4	168.323	237.95	486.95	658.00	19.30			
377	$C_{12}H_{26}$	dodecane	112-40-3	170.338	263.57	489.48	658.00	18.20	754.00	0.251	0.576
378	$C_{12}^{12}H_{26}^{20}O$	1-dodecanol	112-53-8	186.338	297.10	537.79	720.00	20.80	718.00	0.249	0.684
379	$C_{13}H_{12}$	diphenylmethane	101-81-5	168.238	298.39	537.65	760.00	27.10	563.00	0.241	0.481
380	$C_{13}H_{28}$	tridecane	629-50-5	184.365	267.76	508.63	675.00	16.80	823.00	0.246	0.618
381	$C_{13}^{13}H_{28}^{28}O$	1-tridecanol	112-70-9	200.365		553.72	734.00	19.35			0.712
382	$C_{14}^{13}H_{10}^{28}$	phenanthrene	85-01-8	178.233	372.35	611.55	869.00	28.70	554.00	0.220	0.479
383	$C_{14}^{14}H_{10}^{10}$	anthracene	120-12-7	178.233	492.65	614.39	869.30	28.70	554.00	0.220	0.501
384	$C_{14}^{14}H_{22}^{10}$	1,4-di(trimethylmethyl)benzene	1012-72-2	190.330	350.80	510.43	708.00	23.00	732.0	0.286	0.506
	- 14 22	(p-ditertbutylbenzene)									
385	$C_{14}H_{30}$	tetradecane	629-59-4	198.392	279.01	526.76	693.00	15.70	894.00	0.244	0.644
386	$C_{14}H_{30}O$	1-tetradecanol	112-72-1	214.392		569.04	747.00	18.10			0.744
387	$C_{15}H_{32}$	pentadecane	629-62-9	212.419	283.08	543.83	708.00	14.80	966.00	0.243	0.685
388	$C_{15}H_{32}O$	1-pentadecanol	629-76-5	228.419		583.68	759.00	17.00			0.778
389	$C_{16}H_{34}$	hexadecane	544-76-3	226.446	291.32	559.98	723.00	14.00	1034.00	0.241	0.718
390	$C_{16}H_{34}$	2,2,4,4,6,8,8-heptamethylnonane	4390-04-9	226.446		520.25	693.00	15.70			0.548
391	$C_{16}H_{34}O$	1-hexadecanol	4485-13-6	242.446	322.45	597.53	770.00	16.10			0.818
392	$C_{17}H_{36}$	heptadecane	629-78-7	240.473	295.13	574.56	736.00	13.40	1103.00	0.242	0.753
393	$C_{17}H_{36}O$	1-heptadecanol	1454-85-9	256.472	327.00	611.12	780.00	15.00			0.853
394	$C_{18}^{17}H_{14}^{30}$	1,2-diphenylbenzene	84-15-1	230.309	329.35	605.15	857.00	29.90			
395	$C_{18}^{10}H_{14}$	1,3-diphenylbenzene	92-06-8	230.309	360.15	636.15	883.00	24.80	724.00	0.245	
396	$C_{18}H_{14}$	1,4-diphenylbenzene	92-94-4	230.309	483.25	649.15	908.00	29.90	729.00	0.289	
397	$C_{18}^{13}H_{38}^{14}$	octadecane	593-45-3	254.500	301.32	588.30	747.00	12.90	$1\overline{189.00}$	0.247	0.800
398	$C_{18}^{10}H_{38}^{30}O$	1-octadecanol	112-92-5	270.499	331.00	623.57	790.00	$\overline{14.40}$			0.892
399	$C_{19}^{10}H_{40}^{30}$	nonadecane	629-92-5	268.527	305.25	602.34	755.00	11.60			0.845
	17 70										

Section A Basic Constants I (Continued)

No.	Formula	Name	CAS #	Mol. Wt.	Tfp, K	Tb, K	Tc, K	Pc, bar	Vc, cm ³ /mol	Zc = PcVc/RTc	Omega
400	C ₁₉ H ₄₀ O	1-nonadecanol	1454-84-8	284.526		635.41	799.00	13.80			0.934
401	$C_{20}H_{42}$	eicosane	112-95-8	282.554	309.95	616.84	768.00	10.70			0.865
402	$C_{20}H_{42}O$	1-eicosanol	629-96-9	298.553	339.00	647.69	809.00	13.00			0.954
403	$C_{21}^{20}H_{44}$	heneicosane	629-94-7	296.580	313.65	629.65	778.00	10.30			
404	$C_{22}H_{46}$	docosane	629-97-0	310.607	317.55	641.75	786.00	9.80			
405	$C_{23}H_{48}$	tricosane	638-67-5	324.634	320.65	653.35	790.00	9.20			
406	$C_{24}^{23}H_{50}$	tetracosane	646-31-1	338.661	324.05	664.45	800.00	8.70			
407	CID	deuterium chloride	7698-05-7	37.467	158.51	188.43	323.50				
408	ClFO ₃	perchloryl fluoride	7616-94-6	102.449	125.41	226.49	368.40	53.70	161.00	0.282	
409	ClF ₅	chlorine pentafluoride	13637-63-3	130.445	171.15	259.28	416.00	53.00	230.40		
410	ClH	hydrogen chloride	7647-01-0	36.461	158.97	188.15	324.69	83.10	81.00	0.249	
411	ClH₄N	ammonium chloride	12125-02-9	53.491	793.20	613.16	882.00		165.00		
412	CINO	nitrogen oxychloride	2696-92-6	65.459	213.55	267.77	440.60	91.20			
413	Cl ₂	chlorine	7782-50-5	70.905	172.19	239.12	417.00	77.00	124.00	0.275	
414	DĤ	deuterium hydride	13983-20-5	3.022	16.59	22.13	35.90	14.80	62.70	0.311	-0.176
415	DI	deuterium iodide	14104-45-1	128.919	221.28	237.52	421.80				
416	D_2	deuterium	7782-39-0	4.028	18.63	22.13	38.25	16.50	60.20	0.312	
417	$\overline{D_2}$	deuterium, normal	800000-54-8	4.028	18.72	23.65	38.35	16.65	60.20	0.314	-0.143
418	$D_2^{2}O$	deuterium oxide	7789-20-0	20.028	276.96	374.55	643.89	216.71	56.26	0.228	
419	$\overline{D_2S}$	deuterium sulfide	13536-94-2	36.094	187.15		372.30	89.00	96.00	0.276	
420	$\overline{D_3}N$	trideuteroammonia	13550-49-7	20.049	198.82	242.10	405.50	113.00	72.00	0.241	
421	D_3P	trideuterophosphine	13537-03-6	37.016			323.60	65.00			
422	FH	hydrogen fluoride	7664-39-3	20.006	189.58	292.68	461.00	65.00	69.00	0.117	
423	FNO_2	nitrogen dioxyfluoride	10022-50-1	65.004	107.15	200.75	349.30	82.00			
424	F_2	fluorine	7782-41-4	37.997	53.48	84.95	144.30	52.15	66.20	0.288	0.051
425	F_2HN	difluoroamine	10405-27-3	53.012	157.15	250.15	403.00	63.00	97.00	0.182	
426	F_2N_2	cis-difluorodiazene	13812-43-6	66.010		167.40	272.00	68.00	113.30		
427	F_2N_2	trans-difluorodiazene	13776-62-0	66.010	101.15	161.70	260.00	64.00	113.30		
428	F_2O	oxygen difluoride	7783-41-7	53.996	49.35	128.38	215.20	49.50	97.60	0.270	
429	F_2Xe	xenon difluoride	13709-36-9	169.287	402.18		631.00	93.00	149.00	0.264	
430	F_3N	nitrogen trifluoride	7783-54-2	71.002	66.37	144.11	234.00	45.30	118.75	0.256	0.124
431	F ₃ NO	trifluoroamine oxide	13847-65-9	87.001	112.15	185.65	302.60	64.00	150.10		

432	F_4N_2	tetrafluorohydrazine	10036-47-2	104.007	111.65	198.95	309.20	37.00	181.00	0.222	
433	F_4S	sulfur tetrafluoride	7783-60-0	108.060	148.15	233.15	364.00	43.30	146.00	0.209	0.258
434	F_4Xe	xenon tetrafluoride	13709-61-0	207.284	390.25		612.00	70.40	189.00	0.261	
435	F_6S	sulfur hexafluoride	2551-62-4	146.056	222.45	209.25	318.72	37.60	198.40	0.282	0.208
436	F_6U	uranium hexafluoride	7783-81-5	352.070	337.00	324.96	503.35	45.31	250.00	0.271	0.277
437	ΗĬ	hydrogen iodide	10034-85-2	127.912	222.38	237.57	423.90	90.00	132.70		0.038
438	H_2	hydrogen	1333-74-0	2.016	13.83	20.27	32.98	12.93	64.20	0.303	-0.217
439	H ₂	hydrogen, normal	800000-51-5	2.016	13.56	20.38	33.25	12.97	65.00	0.305	-0.216
440	H ₂ O	water	7732-18-5	18.015	273.15	373.15	647.14	220.64	55.95	0.229	0.344
441	H_2S	hydrogen sulfide	7783-06-4	34.082	187.62	212.84	373.40	89.63	98.00	0.283	0.090
442	H_2S_2	dihydrogen disulfide	13465-07-1	66.148	183.35	344.25	572.00	59.10	150.00	0.186	
443	H_2S_3	dihydrogen trisulfide	13845-23-3	98.214		443.13	738.00	51.30	205.00	0.171	
444	H_2S_4	dihydrogen tetrasulfide	13845-25-5	130.280		513.14	855.00	43.70	260.00	0.160	
445	H_2S_5	dihydrogen pentasulfide	13845-24-4	162.346		558.10	930.00	38.90	315.00	0.158	
446	H ₂ Se	hydrogen selenide	7783-07-5	80.976	207.42	228.25	411.00	89.20			
447	H ₃ N	ammonia	7664-41-7	17.031	195.41	239.82	405.40	113.53	72.47	0.255	0.257
448	H ₃ P	phosphine	7803-51-2	33.998	139.37	185.42	324.50	65.40			
449	H_4N_2	hydrazine	302-01-2	32.045	274.68	386.65	653.01	147.00	101.10	0.282	
450	He	helium	7440-59-7	4.003	2.15	4.30	5.19	2.27	57.30	0.301	-0.390
451	He	helium-3	14762-55-1	3.017	1.01	3.33	3.31	1.14	72.50	0.300	-0.480
452	I_2	iodine	7553-56-2	253.809	386.76	457.56	819.00		155.00		
453	Kr	krypton	7439-90-9	83.800	115.77	119.74	209.40	55.00	91.20	0.288	
454	NO	nitrogen monoxide (nitric oxide)	10102-43-9	30.006	109.51	121.38	180.00	64.80	58.00	0.251	0.582
455	N_2	nitrogen	7727-37-9	28.014	63.15	77.35	126.20	33.98	90.10	0.289	0.037
456	N_2O	dinitrogen oxide (nitrous oxide)	10024-97-2	44.013	182.33	184.67	309.60	72.55	97.00	0.273	
457	N_2O_4	dinitrogen tetroxide (nitrogen dioxide)	10544-72-6	92.011	261.95	302.22	431.01	101.00			1.007
458	Ne	neon	7440-01-9	20.180	24.56	27.07	44.40	27.60	41.70	0.312	-0.016
459	OT_2	tritium oxide	14940-65-9	22.032	277.64		641.72	214.10	56.00	0.225	
460	O, -	oxygen	7782-44-7	31.999	54.36	90.17	154.58	50.43	73.37	0.288	
461	$O_2^{-}S$	sulfur dioxide	7446-09-5	64.065	197.67	263.13	430.80	78.84	122.00	0.269	
462	O_3	ozone	10028-15-6	47.998	80.65	161.80	261.05	55.70	89.40	0.229	0.224
463	O_3S	sulfur trioxide	7446-11-9	80.064	289.95	317.90	490.90	82.10	126.50	0.254	
464	Rn	radon	10043-92-2	222.018	202.15	209.80	377.00	63.00	140.00	0.281	
465	S	sulfur	7704-34-9	32.066	392.75	717.75	1313.01	182.00	158.00	0.263	
466	Se	selenium	7782-49-2	78.960	494.00	957.95	1766.00	271.60	62.30	0.118	
467	T_2	tritium	10028-17-8	6.032			40.00	18.50	55.30	0.308	
468	Xe	xenon	7440-63-3	131.290	161.25	165.01	289.74	58.40	118.00	0.286	