

CRITICAL CONSTANTS OF ORGANIC COMPOUNDS

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The parameters of the liquid-gas critical point are important constants in determining the behavior of fluids. This table lists the critical temperature, pressure, and molar volume, as well as the normal boiling point, for over 850 organic substances. The properties and their units are:

T_b : Normal boiling point in K at a pressure of 101.325 kPa (1 atmosphere); an "s" following the value indicates a sublimation point (temperature at which the solid is in equilibrium with the gas at a pressure of 101.325 kPa)

T_c : Critical temperature in K

P_c : Critical pressure in MPa

V_c : Critical molar volume in $\text{cm}^3 \text{mol}^{-1}$

The listed values of the critical constants are critically evaluated using the NIST ThermoData Engine, TDE (Ref. 1), designed to implement the dynamic data evaluation concept (Refs. 2–5). This concept requires large electronic databases capable of storing essentially all relevant experimental data known to date with detailed descriptions of metadata and uncertainties. The combination of these electronic databases with expert-system software, designed to automatically generate recommended property values based on available experimental and predicted data, leads to the ability to

produce critically evaluated data dynamically or "to order." The evaluated data have been generated only for compounds for which experimental data for critical properties are available. Group contribution methods such as Joback-Reed (Ref. 6), Constantinou-Gani (Ref. 7), Marrero-Pardillo (Ref. 8), and Wilson-Jasperson (Ref. 9) as well as quantitative structure-property relationship (QSPR) methods (Ref. 5) were used within the TDE environment to validate available experimental data. Each recommended value in the table is characterized with a combined expanded uncertainty (Ref. 10) (level of confidence, approximately 95%) listed in parentheses. Only references to original experimental data actually used by TDE to generate critically evaluated data are indicated for each compound. Compounds are listed alphabetically by name.

The values of the normal boiling temperatures provided in the table along with the combined expanded uncertainties listed in parentheses have also been critically evaluated using TDE. Additional details on the determination of the normal boiling temperatures using TDE can be found in the Physical Constants of Organic Compounds table in Section 3. The remaining values of the normal boiling temperatures (without uncertainties) are taken from the compilation presented in the 91st Edition of the *CRC Handbook of Chemistry and Physics*.

Name	Mol. Form.	T_b /K	T_c /K	P_c /MPa	$V_c/\text{cm}^3 \text{mol}^{-1}$	Ref.
Acetaldehyde	$\text{C}_2\text{H}_4\text{O}$	293.9(0.6)	462(8)	7.5(1)	154(5)	11–13
Acetic acid	$\text{C}_2\text{H}_4\text{O}_2$	391.0(0.2)	593(2)	5.79(0.03)	171(2)	14–21
Acetic anhydride	$\text{C}_4\text{H}_6\text{O}_3$	412.6(0.3)	606(1)	4.00(0.08)	294(12)	22
Acetone	$\text{C}_3\text{H}_6\text{O}$	329.23(0.07)	508.1(0.2)	4.7(0.1)	221(20)	18, 23–31
Acetonitrile	$\text{C}_2\text{H}_3\text{N}$	354.8(0.2)	545.47(0.07)	4.88(0.01)	173(59)	32–40
Acetophenone	$\text{C}_8\text{H}_8\text{O}$	475.2(0.2)	709.5(0.7)	4.01(0.05)	373(40)	11, 41
Acetylene	C_2H_2	188.45 s	308.4(0.4)	6.24(0.04)	119(11)	42–48
Acrylonitrile	$\text{C}_3\text{H}_3\text{N}$	350.3(0.2)	540(2)	4.6(0.1)	211(10)	49
Allene	C_3H_4	238.3(0.3)	394(4)	6.5(0.7)	167(8)	50
Allyl alcohol	$\text{C}_3\text{H}_6\text{O}$	370.0(0.5)	539.8(0.6)	5.76(0.04)	222(9)	51
Allylamine	$\text{C}_3\text{H}_7\text{N}$	327(2)	540.0(0.7)	4.83(0.03)	217(11)	52
Allyl ethyl ether	$\text{C}_5\text{H}_{10}\text{O}$	338(4)	518(10)	3(2)	320(10)	21
2-Aminobiphenyl	$\text{C}_{12}\text{H}_{11}\text{N}$	571.4(0.2)	838(2)	3.52(0.03)	548(99)	53
2-Aminoethanol	$\text{C}_2\text{H}_7\text{NO}$	443.4(0.4)	671(3)	8.0(0.5)	207(13)	41
2-(2-Aminoethoxy)ethanol	$\text{C}_4\text{H}_{11}\text{NO}_2$	496.2(0.1)	721(4)	4.88(0.1)	333(19)	54
N-(2-Aminoethyl)ethanolamine	$\text{C}_4\text{H}_{12}\text{N}_2\text{O}$	515(5)	739(2)	4.53(0.09)	340(17)	55
Amyl orthosilicate	$\text{C}_{20}\text{H}_{44}\text{O}_4\text{Si}$		714(14)			57
Aniline	$\text{C}_6\text{H}_7\text{N}$	457.2(0.4)	704(7)	5.3(0.1)	291(3)	39, 40, 58–60
Anisole	$\text{C}_7\text{H}_8\text{O}$	426.8(0.2)	646.1(0.2)	4.2(0.1)	355(12)	25, 39, 40, 49
Benzene	C_6H_6	353.23(0.07)	562.0(0.1)	4.90(0.02)	257(11)	19, 27, 28, 34, 61–97
Benzeneacetic acid	$\text{C}_8\text{H}_8\text{O}_2$	541(2)	766(8)	3.9(0.3)	372(16)	98
Benzenebutanoic acid	$\text{C}_{10}\text{H}_{12}\text{O}_2$	569(2)	783(8)	3.2(0.2)	493(18)	98
Benzeneethanol	$\text{C}_8\text{H}_{10}\text{O}$	493(3)	724(4)	4.0(0.2)	390(15)	99
Benzeneheptanoic acid	$\text{C}_{13}\text{H}_{18}\text{O}_2$	585(27)	798(8)	2.5(0.3)	662(21)	98
Benzenehexanoic acid	$\text{C}_{12}\text{H}_{16}\text{O}_2$	574(27)	794(8)	2.6(0.3)	611(20)	98
Benzenepentanoic acid	$\text{C}_{11}\text{H}_{14}\text{O}_2$	583(1)	790(8)	3.1(0.2)	526(19)	98
Benzenepropanoic acid	$\text{C}_9\text{H}_{10}\text{O}_2$	557(2)	776(8)	3.5(0.2)	440(17)	98
Benzo[b]thiophene	$\text{C}_8\text{H}_6\text{S}$	494.0(0.4)	764(2)	4.68(0.04)	359(59)	100
Benzonitrile	$\text{C}_7\text{H}_5\text{N}$	464(1)	691(9)	4.22(0.04)	348(7)	39, 40, 101, 102
Benzophenone	$\text{C}_{13}\text{H}_{10}\text{O}$	579.0(0.2)	830(2)	3.0(0.1)	568(44)	101
Benzyl alcohol	$\text{C}_7\text{H}_8\text{O}$	478.4(0.2)	715(3)	4.3(0.2)	333(13)	103
[1,1'-Bicyclohexyl]-2-one	$\text{C}_{12}\text{H}_{20}\text{O}$	537	787(70)	3(7)	584(20)	104

Name	Mol. Form.	T_b/K	T_c/K	P_c/MPa	$V_c/cm^3 mol^{-1}$	Ref.
1,1'-Bicyclopentyl	$C_{10}H_{18}$	463.61(0.03)	690(2)	3.27(0.03)	497(41)	105
Biphenyl	$C_{12}H_{10}$	528.3(0.3)	773(5)	3.43(0.06)	481(69)	39, 106–109
Bis(2-aminoethyl)amine	$C_4H_{13}N_3$	479.6(0.3)	710(2)	4.43(0.07)	350(21)	55
1,1-Bis(difluoromethoxy)-1,2,2,2-tetrafluoroethane	$C_4H_2F_8O_2$	319.78(0.08)	450(1)	2.40(0.08)	410(4)	110
Bis(difluoromethyl) ether	$C_2H_2F_4O$	278.6(0.4)	420.2(0.1)	4.16(0.06)	223(13)	111
Bis(2-ethylhexyl) phthalate	$C_{24}H_{38}O_4$	657	835(9)	1.1(0.2)	1495(27)	112
Bis(2-hydroxyethyl)methylamine	$C_5H_{13}NO_2$	518(1)	742(4)	4.2(0.4)	404(17)	54
Bis(2,2,2-trifluoroethyl) ether	$C_4H_4F_6O$	336.91	476.31(0.09)	2.78(0.01)	365(1)	113
Bis(trimethylsilyl)methane	$C_7H_{20}Si_2$	406	573.9(0.3)			114
Bromochlorodifluoromethane	$CBrClF_2$	269.2(0.7)	428(12)	4.31(0.01)	229(16)	115
Bromodifluoromethane	$CHBrF_2$	257.5(0.5)	412.0(0.3)	5.2(0.1)	173(18)	116
Bromoethane	C_2H_5Br	311.3(0.6)	503.9(0.4)	6.2(0.1)	214(10)	117, 118
1-Bromo-2-fluorobenzene	C_6H_4BrF	427	669.6(0.6)	4.3(0.6)	342(18)	119
1-Bromo-3-fluorobenzene	C_6H_4BrF	423	652.0(0.4)	4.2(0.6)	337(18)	119
1-Bromo-4-fluorobenzene	C_6H_4BrF	423(2)	654.8(0.4)	4.2(0.2)	338(18)	119
1-Bromopropane	C_3H_7Br	343.9(0.2)	536.9(0.1)	4.33(0.06)	271(6)	120
Bromotrifluoromethane	$CBrF_3$	215.3(0.4)	340.06(0.05)	3.96(0.01)	199(6)	121
1-Bromo-2-(trifluoromethyl)benzene	$C_7H_4BrF_3$	440.7	656.5(0.4)	3.3(0.8)	415(24)	119
1-Bromo-3-(trifluoromethyl)benzene	$C_7H_4BrF_3$	424.7	627.1(0.4)	3.2(0.7)	413(24)	119
1-Bromo-4-(trifluoromethyl)benzene	$C_7H_4BrF_3$	433	629.8(0.4)	3.2(0.8)	413(24)	119
1,3-Butadiene	C_4H_6	268.5(0.2)	425(1)	4.35(0.07)	221(23)	122, 123
Butanal	C_4H_8O	347.9(0.2)	537(2)	4.41(0.1)	258(9)	104, 124
Butane	C_4H_{10}	272.6(0.5)	425.2(0.1)	3.79(0.01)	257(4)	29, 125–139
1,4-Butanediamine	$C_4H_{12}N_2$	429(10)	651(7)	4.5(0.5)	317(14)	140
1,2-Butanediol	$C_4H_{10}O_2$	469.57(0.06)	680(2)	5.4(0.1)	298(12)	141
1,3-Butanediol	$C_4H_{10}O_2$	481.3(0.1)	679(17)	4.7(0.1)	302(67)	54, 141
1,4-Butanediol	$C_4H_{10}O_2$	502.6(0.4)	724(4)	5.5(0.2)	307(14)	49, 55
Butanenitrile	C_4H_7N	390.8(0.4)	585.40(0.07)	3.82(0.05)	265(6)	35, 39, 40, 142
1-Butanethiol	$C_4H_{10}S$	371.5(0.5)	570.1(0.6)	4.01(0.02)	324(12)	143, 144
Butanoic acid	$C_4H_8O_2$	436.8(0.1)	623(6)	4.0(0.3)	292(10)	145–147
1-Butanol	$C_4H_{10}O$	390.8(0.2)	563.0(0.4)	4.43(0.07)	280(14)	34, 80, 148–155
2-Butanol	$C_4H_{10}O$	372.5(0.2)	535(4)	4.2(0.1)	269(4)	148, 149, 153, 156
2-Butanone	C_4H_8O	352.8(0.2)	537(1)	4.18(0.02)	274(30)	25, 30, 33, 34, 38, 157
1-Butene	C_4H_8	276.87(0.08)	419.3(0.1)	4.00(0.05)	236(14)	123, 158–161
<i>cis</i> -2-Butene	C_4H_8	274.03(0.09)	435.7(0.2)	4.23(0.02)	235(4)	86, 123, 158
<i>trans</i> -2-Butene	C_4H_8	266.8(0.2)	428.6(0.1)	4.03(0.02)	238(4)	86, 123, 158
2-Butoxyethanol	$C_6H_{14}O_2$	444(2)	633.9(1)	3.3(0.1)	424(15)	11, 41
1- <i>tert</i> -Butoxy-2-ethoxyethane	$C_8H_{18}O_2$	421.2	585(3)	2.5(0.4)	546(14)	162
2-Butoxyethyl acetate	$C_8H_{16}O_3$	464.2(0.9)	640(2)	2.7(0.2)	551(21)	144, 163
1- <i>tert</i> -Butoxy-2-methoxyethane	$C_7H_{16}O_2$	404(15)	574(1)	2.8(0.7)	480(13)	162
1-Butoxy-2-propanol	$C_7H_{16}O_2$	445(3)	625(1)	2.7(0.1)	479(20)	32
Butyl acetate	$C_6H_{12}O_2$	381(4)	578(10)	3.16(0.06)	403(6)	162, 164–167
<i>sec</i> -Butyl acetate	$C_6H_{12}O_2$	371(1)	571.1(0.5)	3.01(0.1)	398(14)	164, 165
<i>tert</i> -Butyl acetate	$C_6H_{12}O_2$	399.1(0.1)	541(4)	3.0(0.1)	399(7)	54
Butyl acrylate	$C_7H_{12}O_2$	419.8(0.6)	597.4(0.6)	2.76(0.03)	445(7)	51
Butylamine	$C_4H_{11}N$	335.86(0.08)	531.9(0.2)	4.20(0.04)	291(13)	168
<i>sec</i> -Butylamine	$C_4H_{11}N$	317.17(0.07)	514.3(0.2)	4.0(0.2)	284(11)	168
<i>tert</i> -Butylamine	$C_4H_{11}N$	350.1(0.2)	483.7(0.6)	3.85(0.06)	293(23)	75
Butylbenzene	$C_{10}H_{14}$	446.4(0.4)	660.5(0.1)	2.89(0.03)	498(18)	79, 86, 88
<i>sec</i> -Butylbenzene	$C_{10}H_{14}$	442.2(0.3)	652(1)	2.94(0.03)	488(39)	58
<i>tert</i> -Butylbenzene	$C_{10}H_{14}$	456.4(0.3)	648(1)	3.00(0.03)	474(30)	58, 64
Butyl benzoate	$C_{11}H_{14}O_2$	522(3)	725(14)	2.4(0.3)	594(10)	169
Butyl butanoate	$C_8H_{16}O_2$	438.1(0.1)	612(3)	2.4(0.2)	550(9)	162
Butylcyclohexane	$C_{10}H_{20}$	444.8(0.4)	653.1(0.4)	2.57(0.07)	547(14)	170, 171
<i>tert</i> -Butylcyclohexane	$C_{10}H_{20}$	454.0(0.6)	652.0(0.4)	2.82(0.09)	537(15)	170
<i>tert</i> -Butyl ethyl ether	$C_6H_{14}O$	345.8(0.1)	509(2)	3.0(0.2)	394(4)	172
Butyl methyl ether	$C_5H_{12}O$	343.2(0.3)	512.7(0.1)	3.37(0.02)	340(2)	25, 173, 174
Butyl propanoate	$C_7H_{14}O_2$	418.2(0.1)	594(1)	2.8(0.2)	464(10)	162
Butyl vinyl ether	$C_6H_{12}O$	367(1)	540(1)	3.12(0.01)	379(10)	175
γ -Butyrolactone	$C_4H_6O_2$	477.8(0.4)	731(1)	5(1)	246(18)	49

Name	Mol. Form.	T_b/K	T_c/K	P_c/MPa	$V_c/cm^3 mol^{-1}$	Ref.
Chlorobenzene	C_6H_5Cl	404.8(0.2)	632.4(0.1)	4.5(0.1)	303(79)	19, 33, 34
1-Chlorobutane	C_4H_9Cl	351.5(0.2)	539.2(0.6)	4.1(0.2)	303(11)	119
2-Chlorobutane	C_4H_9Cl	341.4	518.6(0.6)	3.4(0.2)	307(14)	119
1-Chloro-2,4-difluorobenzene	$C_6H_3ClF_2$	400	609.6(0.4)	4.0(0.7)	333(16)	119
1-Chloro-2,5-difluorobenzene	$C_6H_3ClF_2$	401	612.5(0.4)	4.0(0.7)	333(16)	119
1-Chloro-3,4-difluorobenzene	$C_6H_3ClF_2$	400	609.2(0.4)	4.0(0.6)	333(16)	119
1-Chloro-3,5-difluorobenzene	$C_6H_3ClF_2$	391.7	592.0(0.4)	3.9(0.7)	327(16)	119
1-Chloro-1,1-difluoroethane	$C_2H_3ClF_2$	264.03(0.07)	410.31(0.05)	4.06(0.03)	230(6)	113, 176–178
1-Chloro-2,2-difluoroethane	$C_2H_2ClF_2$	254.3(0.5)	400.5(0.7)	4.54(0.07)	197(6)	179
Chlorodifluoromethane	$CHClF_2$	232.3(0.5)	369.30(0.05)	4.98(0.01)	165(2)	180–191
2-Chloro-2-(difluoromethoxy)-1,1,1-trifluoroethane	$C_3H_2ClF_5O$	322.4(0.1)	467.8(0.6)	3.05(0.03)	316(24)	192
Chloroethane	C_2H_5Cl	285.4(0.2)	460.3(0.4)	5.24(0.04)	198(11)	193
Chloroethene	C_2H_3Cl	259.3(0.3)	425(5)	5.60(0.03)	171(9)	194
1-Chloro-2-fluorobenzene	C_6H_4ClF	410.8	633.8(0.4)	4.3(0.6)	319(21)	119
1-Chloro-3-fluorobenzene	C_6H_4ClF	401(25)	615.9(0.4)	4.2(0.6)	324(21)	119
1-Chloro-4-fluorobenzene	C_6H_4ClF	403	620.1(0.4)	4.2(0.4)	322(18)	119
1-Chloroheptane	$C_7H_{15}Cl$	432(2)	614(8)	3.1(0.6)	492(14)	119
1-Chlorohexane	$C_6H_{13}Cl$	408.1(0.5)	599(3)	3.3(0.3)	422(12)	119
Chloromethane	CH_3Cl	249.0(0.3)	416.24(0.04)	6.72(0.03)	136(2)	195, 196
2-Chloro-2-methylbutane	$C_5H_{11}Cl$	358(1)	509.1(0.6)	3.2(0.5)	397(15)	119
3-Chloro-3-methylpentane	$C_6H_{13}Cl$	389	528(3)	3(1)	414(14)	119
2-Chloro-2-methylpropane	C_4H_9Cl	324.0(0.5)	497.8(0.1)	3.7(0.4)	308(13)	34
1-Chlorooctane	$C_8H_{17}Cl$	456(3)	643(2)	2.5(0.4)	543(13)	119
Chloropentafluoroacetone	C_5ClF_5O	280.9(0.9)	410.6(0.1)	2.89(0.01)	277(21)	197
Chloropentafluorobenzene	C_6ClF_5	391.11	570(1)	3.2(0.2)	367(25)	198
Chloropentafluoroethane	C_2ClF_5	233.9(0.2)	353.0(0.2)	3.141(0.01)	255(4)	199, 200
1-Chloropentane	$C_5H_{11}Cl$	381.0(0.3)	571.2(0.4)	3.3(0.2)	361(12)	119
1-Chloropropane	C_3H_7Cl	319.3(0.5)	503.3(0.4)	4.56(0.04)	268(24)	118, 193, 201, 202
2-Chloropropane	C_3H_7Cl	308.1(0.6)	482.4(0.4)	4.25(0.04)	245(16)	201, 202
1-Chloro-1,2,2,2-tetrafluoroethane	C_2HClF_4	261.19(0.09)	395.43(0.06)	3.62(0.01)	244(4)	203, 204
4-Chlorotoluene	C_7H_7Cl	435.0(0.2)	615.9(0.5)	2.33(0.09)	377(16)	88
2-Chloro-1,1,1-trifluoroethane	$C_2H_2ClF_3$	279.1(0.6)	425.0(0.2)	4.02(0.02)	232(6)	205
Chlorotrifluoroethene	C_2ClF_3	244.8(0.3)	380.1(0.1)	3.95(0.03)	214(12)	206, 207
2-Chloro-1,1,2-trifluoroethyl difluoromethyl ether	$C_3H_2ClF_5O$	329.9(0.5)	475.0(0.6)	2.98(0.03)	343(25)	192
Chlorotrifluoromethane	$CClF_3$	191.8	301.9(0.2)	3.89(0.01)	180.3(1)	191, 208–216
<i>m</i> -Cresol	C_7H_8O	475.3(0.1)	705.8(0.4)	4.4(0.2)	337(12)	39, 40, 217, 218
<i>o</i> -Cresol	C_7H_8O	464.1(0.1)	697.6(0.2)	4.2(0.2)	336(12)	217, 218
<i>p</i> -Cresol	C_7H_8O	475.0(0.1)	704.6(0.3)	4.1(0.2)	349(13)	217, 218
Cyanogen	C_2N_2	252.1	397(3)	6.2(0.4)	149(8)	219
Cycloheptane	C_7H_{14}	391.9(0.2)	604.2(0.1)	3.85(0.04)	361(12)	34, 220, 221
Cyclohexane	C_6H_{12}	353.8(0.7)	553.4(0.3)	4.07(0.01)	307(12)	19, 34, 78, 82, 84, 88, 90, 144, 163, 170, 171, 220, 222–231
Cyclohexanol	$C_6H_{12}O$	434.0(0.2)	647.1(0.3)	4.3(0.1)	334(33)	49, 232, 233
Cyclohexanone	$C_6H_{10}O$	428.5(0.1)	665(1)	4.61(0.09)	354(12)	162, 166
Cyclohexene	C_6H_{10}	356.0(0.2)	560.45(0.02)	4.43(0.08)	290(20)	86, 201, 202
Cyclohexylamine	$C_6H_{13}N$	406.8(0.5)	626.8(0.9)	3.9(0.6)	349(15)	170
Cyclooctane	C_8H_{16}	424.2(0.1)	647.2(0.4)	3.55(0.06)	417(28)	34, 170, 220, 221
Cyclopentane	C_5H_{10}	322.3(0.1)	511.7(0.2)	4.51(0.07)	264(10)	34, 78, 90, 234–236
Cyclopentanol	$C_5H_{10}O$	413.5(0.2)	619(1)	4.9(0.1)	288(19)	233
Cyclopentanone	C_5H_8O	403.6(0.2)	624(2)	4.59(0.05)	276(17)	233
Cyclopentene	C_5H_8	317.3(0.2)	506.1(0.2)	4.78(0.05)	252(16)	90, 144, 163, 201, 202
Cyclopropane	C_3H_6	242(2)	398.2(0.4)	5.58(0.02)	164(9)	237, 238
Decafluorobiphenyl	$C_{12}F_{10}$	480(2)	640(4)	2.3(0.3)	641(40)	239
1,1,1,2,2,3,4,4,5,5,5-Decafluoro-3,4-bis(trifluoromethyl)-pentane	C_7F_{16}				632(5)	240
1,1,1,2,2,3,4,4,5,5,5-Decafluoro-3-[1,2,2,2-tetrafluoro-1-(trifluoromethyl)ethyl]-4-(trifluoromethyl)pentane	C_9F_{20}	399(24)	581(43)	1(2)	800(7)	240
<i>cis</i> -Decahydronaphthalene	$C_{10}H_{18}$	469.0(0.3)	702(1)	3.2(0.3)	492(19)	241

Name	Mol. Form.	T_b/K	T_c/K	P_c/MPa	$V_c/cm^3 mol^{-1}$	Ref.
<i>trans</i> -Decahydronaphthalene	C ₁₀ H ₁₈	460.4(0.2)	687(1)	3.1(0.1)	499(19)	241
Decamethylcyclopentasiloxane	C ₁₀ H ₃₀ O ₅ Si ₅	486(3)	617.4(0.3)	1.04(0.02)	1201(2)	34, 114
Decanal	C ₁₀ H ₂₀ O	485(3)	674(1)	2.6(0.3)	601(14)	163, 242
Decane	C ₁₀ H ₂₂	447.2(0.1)	618.1(0.9)	2.10(0.03)	621(35)	16, 33, 65, 78, 86, 131, 243–252
1,10-Decanediamine	C ₁₀ H ₂₄ N ₂	535(12)	736(8)	2.4(0.3)	654(28)	140
Decanedioic acid	C ₁₀ H ₁₈ O ₄	647(5)	845(13)	2.5(0.1)	724(21)	253
Decanoic acid	C ₁₀ H ₂₀ O ₂	543(1)	724(5)	1.9(0.8)	638(24)	145, 146
1-Decanol	C ₁₀ H ₂₂ O	502(3)	690(10)	2.3(0.1)	624(87)	149, 152, 254, 255
2-Decanol	C ₁₀ H ₂₂ O	484	668.5(0.3)	2.3(0.5)	646(13)	255
3-Decanol	C ₁₀ H ₂₂ O	490(7)	666.1(0.3)	2.3(0.3)	643(13)	255
4-Decanol	C ₁₀ H ₂₂ O	487(3)	663.7(0.3)	2.3(0.1)	643(13)	255
5-Decanol	C ₁₀ H ₂₂ O	489(5)	663.2(0.4)	2.3(0.4)	646(13)	255
2-Decanone	C ₁₀ H ₂₀ O	484(3)	671.8(0.5)	2.2(0.3)	625(25)	256
3-Decanone	C ₁₀ H ₂₀ O	485(4)	668(1)	2.2(0.2)	628(15)	256
4-Decanone	C ₁₀ H ₂₀ O	479.7	662.9(0.5)	2.2(0.2)	636(18)	256
5-Decanone	C ₁₀ H ₂₀ O	477	661.0(0.4)	2.2(0.2)	628(25)	256
1-Decene	C ₁₀ H ₂₀	444(1)	616.0(0.3)	2.157(0.01)	594(3)	257
Decylbenzene	C ₁₆ H ₂₆	571(1)	752(8)	1.72(0.1)	879(29)	258
Dibenzofuran	C ₁₂ H ₈ O	558.3(0.3)	824(2)	3.37(0.03)	494(32)	259
Dibenzothiophene	C ₁₂ H ₈ S	604.8(0.4)	897(2)	3.9(0.2)	506(108)	260
1,2-Dibromo-1-chloro-1,2,2-trifluoroethane	C ₂ Br ₂ ClF ₃	365.9(0.2)	560.6(0.2)	3.61(0.02)	368(4)	261
1,4-Dibromooctafluorobutane	C ₄ Br ₂ F ₈	371(25)	532(2)	2.4(0.3)	452(29)	198
Dibutylamine	C ₈ H ₁₉ N	435(2)	607.5(0.2)	3.11(0.03)	532(21)	168
1,4-Di- <i>tert</i> -butylbenzene	C ₁₄ H ₂₂	510.4(0.5)	708(2)	2.23(0.01)	732(70)	232
Dibutyl ether	C ₈ H ₁₈ O	414.8(0.3)	584.1(0.2)	2.4(0.2)	521(12)	262
Dibutyl phthalate	C ₁₆ H ₂₂ O ₄	611(9)	797(9)	1.6(0.3)	954(18)	112
<i>m</i> -Dichlorobenzene	C ₆ H ₄ Cl ₂	445(2)	685.7(0.4)	4.2(0.2)	366(22)	119
1,4-Dichlorobenzene	C ₆ H ₄ Cl ₂	447.0(0.2)	669(5)	3.54(0.07)	364(22)	263
Dichlorodiethylsilane	C ₄ H ₁₀ Cl ₂ Si	403(2)	595.7(0.6)	3.06(0.03)	455(4)	264
Dichlorodifluoromethane	CCl ₂ F ₂	243.3(0.1)	384.9(0.2)	4.12(0.01)	218(36)	191, 216, 265
Dichlorodimethylsilane	C ₂ H ₆ Cl ₂ Si	343.6(0.5)	520.3(0.6)	3.49(0.03)	350(5)	266
1,1-Dichloroethane	C ₂ H ₄ Cl ₂	329.4(0.7)	523.4(0.1)	5.1(0.5)	248(12)	267
1,2-Dichloroethane	C ₂ H ₄ Cl ₂	356.5(0.1)	561.5(0.4)	5.4(0.1)	225(8)	33, 34, 268, 269
<i>cis</i> -1,2-Dichloroethene	C ₂ H ₂ Cl ₂	333(2)	535.8(0.4)	5.4(0.3)	220(15)	34
<i>trans</i> -1,2-Dichloroethene	C ₂ H ₂ Cl ₂	320.79(0.08)	515.5(0.2)	5.3(0.2)	216(14)	33, 34
1,1-Dichloro-1-fluoroethane	C ₂ H ₃ Cl ₂ F	305.20(0.09)	477.3(0.1)	4.20(0.02)	253.7(0.6)	178, 270
Dichlorofluoromethane	CHCl ₂ F	282.1	451.6(0.4)	5.20(0.01)	196(1)	271
1,2-Dichlorohexafluoropropane	C ₃ Cl ₂ F ₆	307(2)	451.8(0.1)	2.63(0.07)	365(48)	272
Dichloromethane	CH ₂ Cl ₂	312.9(0.3)	508.0(0.2)	6.35(0.05)	177(13)	273
1,2-Dichloropropane	C ₃ H ₆ Cl ₂	369.6	578(2)	4.63(0.06)	292(6)	119, 232
1,3-Dichloropropane	C ₃ H ₆ Cl ₂	393.9(0.3)	615(3)	4.7(0.5)	299(16)	119
1,1-Dichloro-1,2,2,2-tetrafluoroethane	C ₂ Cl ₂ F ₄	276(1)	418.6(0.8)	3.31(0.03)	294(8)	179
1,2-Dichloro-1,1,2,2-tetrafluoroethane	C ₂ Cl ₂ F ₄	276.8(0.5)	418.74(0.06)	3.25(0.02)	295(3)	191, 274
1,2-Dichloro-1,1,2-trifluoroethane	C ₂ HCl ₂ F ₃	303.1(0.1)	461.6(0.1)	3.77(0.08)	283.2(0.5)	178
2,2-Dichloro-1,1,1-trifluoroethane	C ₂ HCl ₂ F ₃	300.9(0.6)	456.8(0.2)	3.67(0.01)	278(2)	182, 275–280
Didecyl phthalate	C ₂₈ H ₄₆ O ₄	736(4)	870(10)	0.94(0.05)	1807(27)	112
1,1-Diethoxyethane	C ₆ H ₁₄ O ₂	375(2)	539.7(0.4)	3.22(0.08)	426(12)	166
1,2-Diethoxyethane	C ₆ H ₁₄ O ₂	393.8(0.7)	542(3)	2.14(0.02)	432(11)	162
Diethoxymethane	C ₅ H ₁₂ O ₂	359(2)	532(1)	3.4(0.5)	370(10)	162
Diethylamine	C ₄ H ₁₁ N	328.5(0.1)	499.5(0.4)	3.75(0.02)	304(32)	38, 117, 133, 281, 282
<i>p</i> -Diethylbenzene	C ₁₀ H ₁₄	457(1)	657.90(0.03)	2.80(0.08)	494(12)	79, 86
Diethylene glycol	C ₄ H ₁₀ O ₃	518.6(0.2)	753(4)	4.8(0.2)	325(19)	283
Diethylene glycol diethyl ether	C ₈ H ₁₈ O ₃	458(4)	612(10)	2.4(0.7)	587(18)	162
Diethylene glycol dimethyl ether	C ₆ H ₁₄ O ₃	435(2)	617(4)	3.0(0.6)	450(16)	162
Diethylene glycol monobutyl ether	C ₈ H ₁₈ O ₃	505(4)	692(3)	2.8(0.6)	546(26)	41
Diethylene glycol monobutyl ether acetate	C ₁₀ H ₂₀ O ₄	521(2)	694(2)	2.15(0.05)	627(20)	55
Diethylene glycol monoethyl ether	C ₆ H ₁₄ O ₃	475(3)	670(4)	3.2(0.1)	427(23)	49
Diethylene glycol monoethyl ether acetate	C ₈ H ₁₆ O ₄	491(1)	670(12)	2.50(0.06)	524(18)	49, 55
Diethylene glycol monomethyl ether	C ₅ H ₁₂ O ₃	467(2)	672(2)	3.7(0.2)	378(25)	49
Diethylene glycol monopropyl ether	C ₇ H ₁₆ O ₃	488.0(0.4)	680(2)	3.05(0.07)	495(18)	41, 104

Name	Mol. Form.	T_b/K	T_c/K	P_c/MPa	$V_c/cm^3 mol^{-1}$	Ref.
Diethyl ether	$C_4H_{10}O$	307.5(0.5)	466.8(0.3)	3.64(0.01)	280(5)	19, 77, 95, 155, 196, 284–301
Diethyl oxalate	$C_6H_{10}O_4$	459(1)	618(2)	2.14(0.02)	464(37)	101
Diethyl phthalate	$C_{12}H_{14}O_4$	571(2)	776(9)	2.2(0.2)	687(15)	112
Diethyl succinate	$C_8H_{14}O_4$	490(1)	663(30)	2.26(0.02)	567(44)	101, 302
Diethyl sulfide	$C_4H_{10}S$	365.2(0.2)	557.5(1)	4.0(0.1)	322(8)	32, 303, 304
<i>m</i> -Difluorobenzene	$C_6H_4F_2$	356.1(0.5)	548.4(0.4)	4.20(0.01)	289(21)	119
<i>o</i> -Difluorobenzene	$C_6H_4F_2$	367.0(0.5)	566.0(0.4)	4.28(0.01)	290(21)	119
<i>p</i> -Difluorobenzene	$C_6H_4F_2$	362.0(0.3)	556.9(0.4)	4.28(0.07)	297(22)	119
1,1-Difluoroethane	$C_2H_4F_2$	249.1	386.4(0.1)	4.52(0.01)	178(2)	178, 180, 185, 305–307
1,1-Difluoroethene	$C_2H_2F_2$	187.6(0.8)	302.9(0.6)	4.48(0.05)	155(4)	179, 215
2,2-Difluoroethylbis(trifluoromethyl)amine	$C_4H_3F_8N$	324.5	460.20(0.09)	2.64(0.01)	375(1)	308
Difluoromethane	CH_2F_2	221.50(0.07)	351.28(0.03)	5.79(0.01)	121(4)	306, 309–316
3-Difluoromethoxy-1,1,1,2,2-pentafluoropropane	$C_3H_4F_7O$	319.09	455.1(0.1)	2.77(0.02)	363(1)	113
2-(Difluoromethoxy)-1,1,1-trifluoroethane	$C_3H_3F_5O$	302.3(0.2)	444.9(0.3)	3.43(0.01)	291(19)	113
2,4-Difluorotoluene	$C_7H_6F_2$	390	581.4(0.4)	3.7(0.4)	340(21)	119
2,5-Difluorotoluene	$C_7H_6F_2$	391	587.8(0.4)	3.8(0.5)	341(21)	119
2,6-Difluorotoluene	$C_7H_6F_2$	385	581.8(0.4)	3.7(0.4)	341(21)	119
3,4-Difluorotoluene	$C_7H_6F_2$	385	598.5(0.5)	3.8(0.6)	342(22)	119
Diheptyl phthalate	$C_{22}H_{34}O_4$	633	830(9)	1.24(0.08)	1153(22)	112
Dihexyl phthalate	$C_{20}H_{30}O_4$	652(5)	817(9)	1.3(0.1)	1061(22)	112
3,4-Dihydro-2 <i>H</i> -pyran	C_5H_8O	358.6(0.2)	561(2)	4.63(0.08)	268(34)	75
Diisobutylamine	$C_8H_{19}N$	412.8	584.4(0.2)	3.20(0.06)	518(22)	168
Diisopropylamine	$C_6H_{15}N$	357(3)	523.1(0.2)	3.02(0.02)	407(18)	168
1,4-Diisopropylbenzene	$C_{12}H_{18}$	483.4(0.2)	675(1)	2.30(0.04)	610(65)	317
Diisopropyl ether	$C_6H_{14}O$	341.5(0.2)	500.2(0.7)	2.85(0.04)	386(5)	25, 290, 318, 319
1,2-Dimethoxyethane	$C_4H_{10}O_2$	358.1(0.1)	539(4)	3.91(0.05)	305(9)	162, 166, 175, 290
Dimethoxymethane	$C_3H_8O_2$	315.4(0.2)	488(11)	4.0(0.2)	259(10)	75, 162, 320
1,2-Dimethoxypropane	$C_5H_{12}O_2$	369	543(1)	3.4(0.6)	356(12)	162
2,2-Dimethoxypropane	$C_5H_{12}O_2$	350.5(0.7)	510(3)	4(1)	360(13)	162
Dimethyl adipate	$C_8H_{14}O_4$	504(3)	692(14)	2.5(0.5)	561(13)	321
Dimethylamine	C_2H_7N	280.4(0.4)	437.5(0.4)	5.34(0.05)	188(13)	193, 322, 323
<i>N,N</i> -Dimethylaniline	$C_8H_{11}N$	466(1)	687.7(0.6)	3.63(0.09)	407(15)	39, 40
2,2-Dimethylbutane	C_6H_{14}	322.8(0.2)	489.1(0.5)	3.10(0.01)	363.74(0.02)	86, 324–328
2,3-Dimethylbutane	C_6H_{14}	331.1(0.3)	500.2(0.3)	3.13(0.01)	358(1)	19, 86, 252, 324–330
3,3-Dimethyl-2-butanone	$C_6H_{12}O$	379.2(0.2)	570.9(0.3)	3.67(0.03)	383(6)	164, 165
2,3-Dimethyl-1-butene	C_6H_{12}	328.74(0.04)	497.7(0.9)	3.31(0.01)	346(9)	170
3,3-Dimethyl-1-butene	C_6H_{12}	346.34(0.06)	477.4(0.9)	3.18(0.02)	348(10)	170
2,3-Dimethyl-2-butene	C_6H_{12}	314.39(0.04)	521.0(0.9)	3.4(0.1)	344(7)	170
Dimethyl carbonate	$C_3H_6O_3$	363.26(0.09)	557(1)	4.8(0.2)	251(51)	331, 332
<i>cis</i> -1,3-Dimethylcyclohexane	C_8H_{16}	397.5(0.6)	587.7(0.5)	2.88(0.01)	429(10)	64
<i>cis</i> -1,4-Dimethylcyclohexane	C_8H_{16}	392.4(0.5)	603.2(0.3)	3.44(0.02)	434(7)	164, 165
<i>trans</i> -1,4-Dimethylcyclohexane	C_8H_{16}	397.4(0.7)	588(2)	3.04(0.01)	439(18)	74
Dimethyl disulfide	$C_2H_4S_2$	382.87(0.08)	608(4)	5.1(0.1)	266(8)	54
Dimethyl ether	C_2H_6O	248.3(0.2)	400.1(0.8)	5.31(0.03)	171(3)	125, 174, 186, 333–343
<i>N,N</i> -Dimethylformamide	C_3H_7NO	426.0(0.5)	649.6(0.8)	4.4(0.1)	262(9)	11, 344
Dimethyl glutarate	$C_7H_{12}O_4$	489(4)	682(14)	2.8(0.4)	488(14)	321
2,2-Dimethylheptane	C_9H_{20}	406(1)	576.7(0.5)	2.35(0.07)	546(12)	345
2,2-Dimethylhexane	C_8H_{18}	379.9(0.4)	549.9(0.4)	2.53(0.03)	481(10)	346
2,3-Dimethylhexane	C_8H_{18}	388.8(0.5)	563.5(0.4)	2.63(0.02)	466(16)	346
2,4-Dimethylhexane	C_8H_{18}	382.5(0.4)	553(3)	2.55(0.02)	480(39)	330, 346
2,5-Dimethylhexane	C_8H_{18}	382.2(0.7)	550.0(0.3)	2.49(0.02)	485(20)	19, 346
3,3-Dimethylhexane	C_8H_{18}	385.0(0.6)	562.0(0.4)	2.65(0.02)	450(17)	346
3,4-Dimethylhexane	C_8H_{18}	390.8(0.4)	568.8(0.4)	2.69(0.02)	467(21)	346
Dimethyl malonate	$C_5H_8O_4$	454.2(0.6)	647(1)	3.5(0.1)	368(13)	317, 321
2,7-Dimethylnaphthalene	$C_{12}H_{12}$	535.5(0.3)	775(2)	3.02(0.03)	515(45)	347
Dimethyl octanedioate	$C_{10}H_{18}O_4$	532(8)	723(14)	2.3(0.3)	672(14)	321
Dimethyl oxalate	$C_4H_6O_4$	436.5(0.5)	632(14)	4.0(0.3)	315(12)	321, 348
2,2-Dimethyloxirane	C_4H_8O	324(2)	500(4)	4.4(0.1)	327(31)	54
2,2-Dimethylpentane	C_7H_{16}	352.3(0.3)	520.6(0.6)	2.77(0.04)	409(12)	346, 349
2,3-Dimethylpentane	C_7H_{16}	362.9(0.6)	537.5(0.6)	2.92(0.08)	394(23)	346, 349, 350

Name	Mol. Form.	T_b/K	T_c/K	P_c/MPa	$V_c/cm^3 mol^{-1}$	Ref.
2,4-Dimethylpentane	C_7H_{16}	353.5(0.5)	520.0(0.7)	2.74(0.06)	415(19)	346, 349
3,3-Dimethylpentane	C_7H_{16}	359.1(0.6)	536.4(0.4)	2.94(0.02)	413(24)	346
2,3-Dimethyl-1-pentene	C_7H_{14}	357(1)	534(4)	2.9(0.8)	400(11)	170
4,4-Dimethyl-1-pentene	C_7H_{14}	345.6(0.2)	516(4)	2.91(0.01)	406(11)	170
Dimethyl phthalate	$C_{10}H_{10}O_4$	555.8(0.2)	772(9)	2.76(0.08)	557(17)	112
Dimethyl pimelate	$C_9H_{16}O_4$	518(2)	711(14)	2.4(0.1)	608(14)	321
2,3-Dimethylpyridine	C_7H_9N	434.2(0.4)	655.5(0.3)	4.03(0.01)	337(28)	86, 351
2,4-Dimethylpyridine	C_7H_9N	431.5(0.3)	647.1(0.9)	3.83(0.02)	363(34)	90, 352
2,5-Dimethylpyridine	C_7H_9N	430.15(0.05)	644.2(0.3)	4.11(0.04)	371(45)	86
2,6-Dimethylpyridine	C_7H_9N	417.1(0.1)	623.8(0.2)	3.80(0.04)	353(43)	90
3,4-Dimethylpyridine	C_7H_9N	452.2(0.3)	683.8(0.4)	4.06(0.01)	353(13)	86, 351, 352
3,5-Dimethylpyridine	C_7H_9N	445.0(0.1)	667.3(0.3)	3.84(0.02)	369(12)	86, 351
2,6-Dimethylquinoline	$C_{11}H_{11}N$	541.2(0.4)	786(2)	3.27(0.02)	505(24)	529
Dimethyl sebacate	$C_{12}H_{22}O_4$	562(3)	742(14)	2.1(0.2)	695(14)	321
Dimethyl succinate	$C_6H_{10}O_4$	470(1)	662(14)	3.5(0.2)	426(16)	321
Dimethyl sulfide	C_2H_6S	310.47(0.05)	503.0(0.3)	5.40(0.06)	201(9)	52, 117, 304
Dimethyl sulfoxide	C_2H_6OS	465.0(0.9)	707(1)	4.6(0.7)	228(7)	23
1,3-Dimethyl-1,1,3,3-tetraphenyldisiloxane	$C_{26}H_{26}OSi_2$	701(2)	893(9)	1.38(0.1)	1300(89)	353
1,3-Dimethyltricyclo[3.3.1.1 ^{3,7}]decane	$C_{12}H_{20}$	476.53	708(2)	2.86(0.01)	595(115)	141
Dinonyl phthalate	$C_{26}H_{42}O_4$	686	858(9)	1.0(0.3)	1652(26)	112
Diocetyl phthalate	$C_{28}H_{50}O_4$	688(4)	840(9)	1.1(0.1)	1510(22)	112
1,4-Dioxane	$C_4H_8O_2$	374.3(0.3)	587.3(0.1)	5.2(0.2)	251(5)	34, 290, 354
Dipentyl phthalate	$C_{18}H_{26}O_4$	614(40)	811(9)	1.4(0.7)	957(21)	112
Diphenyl ether	$C_{12}H_{10}O$	531.1(0.1)	766.9(0.8)	3.10(0.04)	526(23)	25, 355
Diphenylmethane	$C_{13}H_{12}$	537.3(0.3)	776(9)	3.02(0.07)	546(144)	39, 302, 356, 357
1,3-Diphenyltetramethyldisiloxane	$C_{16}H_{22}OSi_2$		750(8)			353
Dipropylamine	$C_6H_{15}N$	380.6(0.9)	555.8(0.1)	3.6(0.1)	414(15)	168
Dipropylene glycol	$C_6H_{14}O_3$	504(2)	705(4)	3.4(0.1)	444(18)	54
Dipropyl ether	$C_6H_{14}O$	363.2(0.3)	531(2)	2.92(0.05)	402(8)	25, 52
Dipropyl phthalate	$C_{14}H_{18}O_4$	592(2)	784(9)	1.9(0.1)	816(20)	112
Diundecyl phthalate	$C_{30}H_{50}O_4$	711(25)	886(10)	0.89(0.1)	1590(25)	112
Docosane	$C_{22}H_{46}$	642(5)	786(6)	1.0(0.1)	1434(50)	62, 358, 359
Docosanoic acid	$C_{22}H_{44}O_2$	693(3)	837(8)	1.11(0.08)	1485(29)	360
1-Docosanol	$C_{22}H_{46}O$	680(12)	827(8)	1.0(0.6)	1243(20)	361
1,2,2,3,3,4,4,5,5,6,6,7-Dodecafluoro-1-heptanol	$C_7H_4F_{12}O$	444.7(0.7)	589(5)	2.0(0.2)	620(34)	362
Dodecane	$C_{12}H_{26}$	489.4(0.2)	658.8(0.9)	1.80(0.09)	747(18)	34, 86, 118, 243, 244, 363, 364
1,12-Dodecanediamine	$C_{12}H_{26}N_2$	572(13)	767(8)	2.0(0.3)	765(37)	140
Dodecanedioic acid	$C_{12}H_{22}O_4$	621(10)	859(13)	2.1(0.2)	730(19)	253
1-Dodecanethiol	$C_{12}H_{26}S$	550(3)	734(4)	1.81(0.1)	726(25)	99
Dodecanoic acid	$C_{12}H_{24}O_2$	572(1)	743(7)	1.9(0.2)	787(19)	360
1-Dodecanol	$C_{12}H_{26}O$	537.2(0.3)	719.4(0.6)	2.02(0.05)	805(15)	149
2-Dodecanone	$C_{12}H_{24}O$	520(6)	702(4)	1.9(0.9)	742(20)	256
3-Dodecanone	$C_{12}H_{24}O$	523(3)	701(2)	1.9(0.2)	680(19)	256
4-Dodecanone	$C_{12}H_{24}O$	528(5)	697(2)	1.9(0.4)	672(19)	256
5-Dodecanone	$C_{12}H_{24}O$	521(4)	695(5)	1.9(0.4)	678(19)	256
6-Dodecanone	$C_{12}H_{24}O$	522(4)	694(2)	1.9(0.3)	677(19)	256
1-Dodecene	$C_{12}H_{24}$	486.5(0.9)	657.6(0.6)	1.88(0.01)	710(15)	257
Eicosane	$C_{20}H_{42}$	617.2(0.9)	768(6)	1.08(0.05)	1325(48)	358, 359
Eicosanoic acid	$C_{20}H_{40}O_2$	673(6)	820(8)	1.2(0.1)	1346(27)	360
1-Eicosanol	$C_{20}H_{42}O$	629	808(8)	1.1(0.2)	1130(18)	361
1-Eicosene	$C_{20}H_{40}$	620(15)	772(15)	1.1(0.3)	1213(36)	365
Ethane	C_2H_6	184.5(0.4)	305.36(0.04)	4.88(0.01)	146(3)	44, 48, 92, 131, 132, 134, 214, 215, 229, 246, 366–393
1,2-Ethanediamine	$C_2H_8N_2$	390.0(0.5)	613.1(0.3)	6.71(0.04)	204(13)	49
1,2-Ethandiol	$C_2H_6O_2$	470.6(0.1)	719(5)	8.1(0.4)	180(11)	11, 32, 41, 394
1,1-Ethandiol, diacetate	$C_6H_{10}O_4$	441(3)	618(4)	2.9(0.1)	457(13)	54
Ethanethiol	C_2H_6S	308.1(0.1)	498.7(0.3)	5.53(0.08)	208(10)	304

Name	Mol. Form.	T_b/K	T_c/K	P_c/MPa	$V_c/cm^3 mol^{-1}$	Ref.
Ethanol	C_2H_6O	351.39(0.09)	515(1)	6.25(0.04)	169(4)	19, 25, 34, 80, 85, 128, 148–150, 152, 222, 395–409
Ethoxybenzene	$C_8H_{10}O$	443.0(0.2)	647(2)	3.45(0.05)	407(16)	39, 40
2-Ethoxyethyl acetate	$C_6H_{12}O_3$	429.8(0.4)	609(2)	3.07(0.03)	443(38)	141, 144, 163
2-Ethoxy-2-methylbutane	$C_7H_{16}O$	374.6(0.4)	546(2)	2.83(0.09)	448(29)	410
1-Ethoxy-1,1,2,2,3,3,4,4,4-nonafluorobutane	$C_6H_5F_9O$	350.04	482.0(0.1)	1.98(0.01)	518(2)	113
Ethyl acetate	$C_4H_8O_2$	350.2(0.2)	523.27(0.07)	3.88(0.02)	288(19)	19, 23, 400, 411–413
Ethylamine	C_2H_7N	289.8(0.2)	456.5(0.9)	5.6(0.1)	183(22)	193, 414
Ethylbenzene	C_8H_{10}	409.3(0.4)	617.1(0.1)	3.61(0.01)	365(61)	68, 78, 79, 86, 88, 172, 345
Ethyl benzoate	$C_9H_{10}O_2$	485.6(0.2)	700(14)	3.01(0.05)	470(12)	169
Ethyl butanoate	$C_6H_{12}O_2$	394.2(0.4)	566.1(0.1)	3.2(0.3)	421(30)	415, 416
Ethyl <i>trans</i> -2-butenolate	$C_6H_{10}O_2$	413(5)	599(10)	3(2)	382(7)	21
Ethylcyclohexane	C_8H_{16}	404.9(0.4)	606.9(0.4)	3.27(0.04)	431(12)	170, 171
Ethylcyclopentane	C_7H_{14}	376.6(0.6)	569.48(0.05)	3.40(0.07)	377(3)	236
Ethyl 2,2-dimethylpropanoate	$C_7H_{14}O_2$	391.4(0.4)	566(2)	2.88(0.02)	461(8)	417
Ethylene	C_2H_4	169.3(0.3)	282.35(0.03)	5.06(0.01)	130.9(0.2)	61, 246, 371, 418–436
Ethyl 3-ethoxypropanoate	$C_7H_{14}O_3$	441(2)	621(3)	2.7(0.1)	478(19)	11, 41
Ethyl formate	$C_3H_6O_2$	327.2(0.1)	508.5(0.5)	4.78(0.03)	223(47)	413, 415
Ethyl heptanoate	$C_9H_{18}O_2$	461(2)	634(1)	2.3(0.2)	587(8)	162
3-Ethylhexane	C_8H_{18}	391.6(0.5)	565.5(0.4)	2.61(0.02)	450(15)	346
Ethyl hexanoate	$C_8H_{16}O_2$	438(1)	615(1)	2.6(0.2)	528(8)	162
2-Ethylhexanoic acid	$C_8H_{16}O_2$	500.6(0.1)	674(1)	2.75(0.06)	543(8)	144, 163
2-Ethyl-1-hexanol	$C_8H_{18}O$	459.3(0.2)	640.2(0.3)	3.0(0.2)	508(29)	437
2-Ethylhexyl acetate	$C_{10}H_{20}O_2$	473(1)	642(2)	2.02(0.01)	644(10)	410
Ethyl 3-methylbutanoate	$C_7H_{14}O_2$	408(3)	584(6)	3(1)	463(9)	147, 162
Ethyl methyl ether	C_3H_8O	279(2)	437.8(0.2)	4.39(0.06)	219(5)	174, 304, 438
3-Ethyl-2-methylpentane	C_8H_{18}	388.8(0.6)	567.1(0.4)	2.70(0.02)	442(22)	346
3-Ethyl-3-methylpentane	C_8H_{18}	391.3(0.9)	576.5(0.4)	2.77(0.01)	463(13)	346
Ethyl 2-methylpropanoate	$C_6H_{12}O_2$	384(2)	554(4)	3.1(0.3)	421(76)	415
Ethyl methyl sulfide	C_3H_8S	339.8(0.3)	533(10)	4.62(0.03)	260(6)	303
Ethyl nonanoate	$C_{11}H_{22}O_2$	497(5)	664(1)	2.0(0.4)	715(8)	162
Ethyl octanoate	$C_{10}H_{20}O_2$	479(1)	652(12)	2(2)	657(8)	147, 162
3-Ethylpentane	C_7H_{16}	366.5(0.4)	540.7(0.4)	2.90(0.03)	412(14)	346, 349
Ethyl pentanoate	$C_7H_{14}O_2$	415(3)	593(1)	2.8(0.4)	466(10)	162
2-Ethylphenol	$C_8H_{10}O$	477.6(0.1)	703(1)	3.7(0.3)	388(15)	218
3-Ethylphenol	$C_8H_{10}O$	491.5(0.1)	716(1)	3.8(0.3)	393(15)	218
4-Ethylphenol	$C_8H_{10}O$	491.12(0.06)	716(1)	3.1(0.6)	395(15)	218
Ethyl propanoate	$C_5H_{10}O_2$	372.0(0.2)	547(1)	3.37(0.05)	343(23)	19, 64, 162, 413, 415
Ethyl propyl ether	$C_5H_{12}O$	336(3)	500.2(0.4)	3.37(0.01)	343(44)	25, 304
S-Ethyl thioacetate	C_4H_8OS	387(3)	590.5(0.2)	4.1(0.1)	320(10)	49
4-Ethyltoluene	C_9H_{12}	435.1(0.6)	640.2(0.5)	3.23(0.04)	446(12)	439, 440
Ethyl vinyl ether	C_4H_8O	309(2)	475(2)	4.06(0.04)	262(8)	290
Fluorobenzene	C_6H_5F	357.8(0.3)	560.10(0.07)	4.55(0.01)	272(9)	86, 441
Fluoroethane	C_2H_5F	235.4(0.3)	375.2(0.2)	5.02(0.01)	164(3)	206, 442, 443
Fluoromethane	CH_3F	194.8	317.42(0.01)	5.88(0.01)	112.41(0.01)	444, 445
2-Fluorotoluene	C_7H_7F	387(2)	591.2(0.4)	3.9(0.3)	323(47)	119
3-Fluorotoluene	C_7H_7F	389(2)	591.8(0.4)	3.9(0.3)	332(17)	119
4-Fluorotoluene	C_7H_7F	389.8(0.4)	592.1(0.8)	3.85(0.01)	332(17)	119
Formic acid	CH_2O_2	374	588(10)		115.88(0.08)	146
Furan	C_4H_4O	304.4(0.2)	490.2(0.2)	5.43(0.08)	218(3)	241, 290
Glycerol	$C_3H_8O_3$	562(3)	850(9)	7.6(0.8)	251(15)	394
Heneicosane	$C_{21}H_{44}$	632(6)	778(8)	1.0(0.1)	1366(48)	359
Heptadecane	$C_{17}H_{36}$	576(2)	736(1)	1.33(0.07)	1081(28)	359
Heptadecanoic acid	$C_{17}H_{34}O_2$	635(4)	792(8)	1.4(0.1)	1130(24)	360
1-Heptadecanol	$C_{17}H_{36}O$	597	780(8)	1.4(0.1)	1097(18)	361
1-Heptadecene	$C_{17}H_{34}$	574(3)	734(7)	1.34(0.08)	1053(35)	365
2,2,3,3,5,5,6-Heptafluoro-1,4-dioxane	$C_4HF_7O_2$	312.5(0.1)	453(1)	2.86(0.06)	359(4)	110
1,1,1,2,2,3,3-Heptafluoropentan-4-one	$C_5H_3F_7O$	337.4	476.55(0.08)	2.57(0.01)	394.2(0.7)	308

Name	Mol. Form.	T_b/K	T_c/K	P_c/MPa	$V_c/cm^3 mol^{-1}$	Ref.
1,1,1,2,3,3,3-Heptafluoropropane	$C_3H_7F_7$	257.65	375.0(0.1)	2.93(0.01)	299(7)	446–449
1,1,1,2,4,4,4-Heptafluoro-2-trifluoromethoxybutane	$C_5H_2F_{10}O$	322.73(0.09)	447(1)	2.15(0.06)	465(5)	110
1,1,1,2,2,3,3-Heptafluoro-3-(trifluoromethoxy)propane	$C_4F_{10}O$	280.0(0.4)	391.7(0.7)	1.89(0.05)	431(31)	450
2,2,4,4,6,8,8-Heptamethylnonane	$C_{16}H_{34}$	519(1)	692(4)	1.53(0.01)	957(27)	451
1,1,1,3,5,5,5-Heptamethyltrisiloxane	$C_7H_{22}O_2Si_3$	416(1)	553.4(0.6)	1.48(0.02)	828(3)	452
Heptanal	$C_7H_{14}O$	426(3)	616.8(0.4)	3.2(0.2)	434(7)	242
2-Heptanamine	$C_7H_{17}N$	414(4)	598(2)	2.9(0.3)	455(18)	170
Heptane	C_7H_{16}	371.53(0.07)	540.1(0.2)	2.74(0.01)	428(15)	34, 36, 62, 75, 78, 86, 131, 133–135, 157, 243, 244, 254, 286, 324, 346, 359, 394, 432, 453–463
Heptanedioic acid	$C_7H_{12}O_4$	615.1	842(13)	3.3(0.2)	463(15)	253
Heptanoic acid	$C_7H_{14}O_2$	495(2)	678(2)	3.0(0.3)	476(5)	145, 146, 464
1-Heptanol	$C_7H_{16}O$	451(1)	632.4(0.6)	3.1(0.2)	430(9)	149, 254, 465
2-Heptanol	$C_7H_{16}O$	432	608.4(0.6)	3.0(0.1)	442(2)	149, 465
3-Heptanol	$C_7H_{16}O$	436(2)	605.4(0.3)	3.1(0.4)	451(3)	465
4-Heptanol	$C_7H_{16}O$	434(2)	602.6(0.3)	3.1(0.6)	455(4)	465
2-Heptanone	$C_7H_{14}O$	424.1(0.3)	611.4(0.2)	2.98(0.04)	436(4)	25, 172, 466
3-Heptanone	$C_7H_{14}O$	419(2)	606.6(0.2)	3.0(0.1)	433(5)	466
4-Heptanone	$C_7H_{14}O$	417(1)	602.0(0.2)	3.0(0.3)	434(5)	466
1-Heptene	C_7H_{14}	370(2)	537.3(0.3)	2.85(0.02)	409(2)	51, 52, 86, 124, 257, 467
<i>cis</i> -2-Heptene	C_7H_{14}	371(2)	548.5(0.6)	3.0(0.3)	410(7)	170
<i>trans</i> -2-Heptene	C_7H_{14}	369(2)	542.8(0.4)	3.0(0.2)	410(7)	170
<i>trans</i> -3-Heptene	C_7H_{14}	367(1)	538.6(0.7)	3.0(0.2)	411(7)	170
Heptylbenzene	$C_{13}H_{20}$	515(4)	708(7)	2.1(0.2)	680(16)	258
Heptyl orthosilicate	$C_{28}H_{60}O_4Si$		778(16)			57
Hexacosane	$C_{26}H_{54}$	688(11)	816(8)	0.8(0.2)	1740(59)	358
Hexadecane	$C_{16}H_{34}$	560.0(0.7)	722.2(0.8)	1.4(0.2)	1009(53)	34, 244, 245
Hexadecanoic acid	$C_{16}H_{32}O_2$	624(6)	785(8)	1.5(0.2)	1059(23)	360
1-Hexadecanol	$C_{16}H_{34}O$	598(2)	770(8)	1.47(0.1)	1019(17)	361
1-Hexadecene	$C_{16}H_{32}$	558(1)	718(7)	1.4(0.2)	986(33)	365
Hexaethyldisiloxane	$C_{12}H_{30}OSi_2$	525(7)	692.9(0.1)	1.7(0.7)	955(2)	34
Hexafluoroacetylacetone	$C_5H_2F_6O_2$	342(2)	485.1(0.5)	2.9(0.2)	313(49)	468
Hexafluorobenzene	C_6F_6	353.3(0.2)	516.4(0.5)	3.28(0.01)	337(4)	74, 198, 286, 469–473
2,2,4,4,5,5-Hexafluoro-1,3-dioxolane	$C_3F_6O_2$	251.0(0.2)	368.1(0.7)	2.72(0.04)	293(30)	116
Hexafluoroethane	C_2F_6	195.0(0.1)	292.9(0.2)	3.03(0.01)	223(3)	172, 474–477
1,1,1,3,3,3-Hexafluoro-2-methoxy-2-(trifluoromethyl)propane	$C_5H_3F_9O$	327(1)	463(1)	2.37(0.07)	448(5)	110
1,1,1,2,3,3-Hexafluoro-3-(2,2,3,3,3-pentafluoropropoxy)propane	$C_6H_3F_{11}O$	360.64	486.48(0.07)	1.95(0.01)	529(2)	113
1,1,1,2,3,3-Hexafluoropropane	$C_3H_2F_6$	277.65	412.40(0.06)	3.42(0.01)	270(5)	203, 447, 448, 478, 479
1,1,1,3,3,3-Hexafluoropropane	$C_3H_2F_6$	271.8(0.2)	398.07(0.06)	3.18(0.01)	262(18)	203
1,1,1,2,3,3-Hexafluoro-3-(2,2,3,3-tetrafluoropropoxy)propane	$C_6H_4F_{10}O$	379.07	516.2(0.3)	2.2(0.2)	543(34)	113
1,1,1,2,3,3-Hexafluoro-3-(2,2,2-trifluoroethoxy)propane	$C_5H_3F_9O$	345.87	475.74(0.09)	2.23(0.02)	455(2)	113
Hexamethylbenzene	$C_{12}H_{18}$	541(3)	758(2)	2.6(0.4)	581(15)	25
1,1,1,5,5,5-Hexamethyl-3,3-bis[(trimethylsilyl)oxy]trisiloxane	$C_{12}H_{36}O_4Si_5$	494.6(0.2)	622.6(0.2)	1.03(0.02)	1323(90)	34
Hexamethyldisiloxane	$C_6H_{18}OSi_2$	373.6(0.3)	518.7(0.6)	1.95(0.02)	629(15)	480
2,6,10,15,19,23-Hexamethyltetracosane	$C_{30}H_{62}$	693(6)	796(2)	0.60(0.04)	2060(70)	481
Hexanal	$C_6H_{12}O$	402.8(0.4)	592(3)	3.4(0.2)	378(7)	163, 242
Hexane	C_6H_{14}	341.87(0.06)	507.5(0.1)	3.03(0.01)	366.0(0.8)	19, 29, 33, 34, 38, 52, 77, 78, 84, 86, 96, 106, 124, 131, 133, 134, 243, 244, 254, 281, 286, 318, 324–328, 401, 456, 458, 470, 482–488
1,6-Hexanediamine	$C_6H_{16}N_2$	470(2)	685(7)	3.6(0.5)	446(17)	140

Name	Mol. Form.	T_b/K	T_c/K	P_c/MPa	$V_c/cm^3 mol^{-1}$	Ref.
1,6-Hexanedioic acid	$C_6H_{10}O_4$	610.5	841(13)	3.8(0.3)	449(17)	253
1,6-Hexanediol	$C_6H_{14}O_2$	481	741(10)	4.1(0.1)	404(13)	489
Hexanenitrile	$C_6H_{11}N$	436.6(0.3)	633.8(0.2)	2.99(0.06)	378(8)	35
Hexanoic acid	$C_6H_{12}O_2$	478.0(0.6)	661(7)		413(15)	145, 146, 360, 464
1-Hexanol	$C_6H_{14}O$	430.0(0.7)	611.0(0.4)	3.40(0.09)	381(30)	34, 149, 151, 152, 254, 255, 456
2-Hexanol	$C_6H_{14}O$	413	585.9(0.5)	3.3(0.3)	406(8)	149, 255, 437
3-Hexanol	$C_6H_{14}O$	416(2)	582.4(0.4)	3.3(0.1)	378(14)	64, 254, 255
2-Hexanone	$C_6H_{12}O$	400.8(0.1)	586.7(0.5)	3.31(0.04)	377(4)	25, 172, 466
3-Hexanone	$C_6H_{12}O$	396.6(0.3)	583.1(0.5)	3.32(0.01)	378(4)	25, 466
Hexatriacontane	$C_{36}H_{74}$	777(7)	872(9)	0.47(0.07)	2711(2)	358
1-Hexene	C_6H_{12}	336.5(0.1)	504.1(0.9)	3.20(0.03)	381(9)	86, 201, 202, 252, 257, 490
<i>cis</i> -2-Hexene	C_6H_{12}	342.0(0.5)	513.4(0.9)	3.34(0.06)	347(7)	170
<i>trans</i> -2-Hexene	C_6H_{12}	341.00(0.09)	509.0(0.7)	3.16(0.01)	353(7)	170
<i>cis</i> -3-Hexene	C_6H_{12}	339.5(0.5)	510(1)	3.29(0.01)	351(7)	170
<i>trans</i> -3-Hexene	C_6H_{12}	340.21(0.09)	507(2)	3.18(0.01)	352(7)	170
5-Hexen-2-one	$C_6H_{10}O$	402.2(0.5)	593.5(0.6)	3.51(0.04)	359(10)	51
Hexyl acetate	$C_8H_{16}O_2$	444.2(0.7)	618(1)	2.5(0.1)	526(8)	162
Hexylamine	$C_6H_{15}N$	405(1)	592.3(0.7)	3.4(0.3)	402(15)	170
Hexylbenzene	$C_{12}H_{18}$	499(2)	695(7)	2.4(0.2)	620(14)	258
Hexyl benzoate	$C_{13}H_{18}O_2$	550(26)	748(14)	2.0(0.3)	658(14)	169
Indan	C_9H_{10}	451.0(0.4)	684.8(0.4)	3.95(0.03)	385(22)	25
Isobutanal	C_4H_8O	337.2(0.2)	543.6(0.6)	5.12(0.08)	283(10)	164, 165
Isobutane	C_4H_{10}	261.4(0.5)	407.84(0.07)	3.64(0.02)	256(7)	125, 132, 491, 492
Isobutene	C_4H_8	266.1(0.2)	418.0(0.3)	4.00(0.04)	240(3)	123, 158, 493, 494
Isobutyl acetate	$C_6H_{12}O_2$	390.0(0.6)	562(2)	2.97(0.06)	369(37)	166, 201, 202, 415
Isobutylbenzene	$C_{10}H_{14}$	445.8(0.4)	650(3)	3.0(0.2)	493(15)	96
Isobutyl butanoate	$C_8H_{16}O_2$	430(1)	611(6)	2.5(0.3)	524(9)	147
Isobutylcyclohexane	$C_{10}H_{20}$	444.5	642.1(0.6)	2.61(0.07)	550(14)	170
Isobutyl formate	$C_5H_{10}O_2$	371.5(0.3)	551(4)	3.9(0.4)	359(25)	415
Isobutyl isobutanoate	$C_8H_{16}O_2$	421(3)	602(6)	2.5(0.8)	530(9)	147
Isobutyl 3-methylbutanoate	$C_9H_{18}O_2$	442(3)	621(6)	2.3(0.7)	581(9)	147
Isobutyl propanoate	$C_7H_{14}O_2$	409(2)	586(8)	3(1)	462(9)	162, 167
Isopentane	C_5H_{12}	300.98(0.06)	460.37(0.09)	3.35(0.06)	313(17)	19, 86, 96, 127, 495
Isopentyl acetate	$C_7H_{14}O_2$	414.8(0.7)	586.1(0.4)	2.76(0.07)	464(9)	166
Isopentyl butanoate	$C_9H_{18}O_2$	458.0(0.3)	619(6)	3(1)	595(10)	147
Isopentyl nitrite	$C_5H_{11}NO_2$	372(3)	626(16)	5.07(0.04)	386.2(0.1)	52
Isopentyl propanoate	$C_8H_{16}O_2$	446(4)	611(6)	2.5(0.9)	523(9)	147
Isopropyl acetate	$C_5H_{10}O_2$	361.8(0.2)	531.1(0.6)	3.31(0.04)	343(4)	64, 166, 172, 411, 490
Isopropylamine	C_3H_9N	304.9(0.2)	472.2(0.9)	4.55(0.07)	231(5)	75, 170
Isopropylbenzene	C_9H_{12}	425.5(0.2)	631(1)	3.2(0.1)	423(5)	79, 90, 96, 172
Isopropylcyclohexane	C_9H_{18}	427.5(0.4)	632.2(0.4)	3.1(0.2)	484(14)	170
Isopropyl formate	$C_4H_8O_2$	341(2)	534.6(0.5)	3.95(0.03)	294(11)	164, 165
1-Isopropyl-4-methylbenzene	$C_{10}H_{14}$	450(2)	654(8)	2.8(0.1)	495(16)	96, 147
(1 <i>S</i> ,2 <i>R</i> ,5 <i>S</i>)-2-Isopropyl-5-methylcyclohexanol	$C_{10}H_{20}O$	489(3)	694(5)	2.7(0.6)	539(14)	302
Isopropyl methyl ether	$C_4H_{10}O$	303.9(0.5)	464.4(0.2)	3.76(0.01)	287(11)	25
Isoquinoline	C_8H_7N	516.3(0.6)	803(8)	5.07(0.03)	380(17)	218
<i>d</i> -Limonene	$C_{10}H_{16}$	450.8(0.5)	653(2)	2.81(0.02)	498(11)	496
Mesityl oxide	$C_6H_{10}O$	402.8(0.4)	605(2)	3.85(0.02)	353(27)	497
Methane	CH_4	111.6(0.2)	190.56(0.02)	4.60(0.01)	99(3)	132, 134, 498–508
Methane- <i>d</i> ₄	CD_4		189.2(0.6)		98(3)	507
Methanethiol	CH_4S	279.1(0.1)	469.9(0.3)	7.24(0.09)	148(4)	304
Methanol	CH_4O	337.6(0.7)	512.7(0.6)	8.01(0.03)	117(4)	19, 38, 67, 80, 85, 87, 131, 148, 152, 155, 196, 405, 458, 482, 509–516
1-Methoxy-2,4-dimethylbenzene	$C_9H_{12}O$	465	682(4)	3.2(0.7)	451(18)	162
2-Methoxy-1,4-dimethylbenzene	$C_9H_{12}O$	467	677(1)	3.2(0.7)	451(15)	162
2-Methoxyethanol	$C_3H_8O_2$	397.4(0.1)	598(1)	5.28(0.08)	263(6)	49
2-Methoxyethyl acetate	$C_5H_{10}O_3$	415(3)	603(3)	3.6(0.5)	368(15)	162

Name	Mol. Form.	T_b/K	T_c/K	P_c/MPa	$V_c/cm^3 mol^{-1}$	Ref.
4-Methoxy-1,1,1,2,2,3,3,3-heptafluorobutane	$C_5H_5F_7O$	344.13	481.5(0.2)	2.38(0.01)	431(2)	113
1-Methoxy-1,1,2,2,3,3,3-hexafluoropropane	$C_4H_4F_6O$	341.02	487.0(0.3)	2.9(0.1)	370(25)	113
2-Methoxy-2-methylbutane	$C_6H_{14}O$	359.5(0.1)	536(2)	3.23(0.09)	372(19)	172, 410, 517
5-Methoxy-1,1,2,2,3,3,4,4-octafluoropentane	$C_6H_6F_8O$	395.83	546.1(0.3)	2.40(0.07)	493(30)	113
1-Methoxy-2-propanol	$C_4H_{10}O_2$	393.1(0.6)	579.8(0.3)	4.11(0.04)	304(12)	49
2-Methoxypropene	C_4H_8O	308.8(0.3)	478.5(0.6)	4.2(0.3)	257(13)	518
Methyl acetate	$C_3H_6O_2$	329.8(0.2)	506.7(0.4)	4.73(0.07)	227(22)	19, 411–413
Methylamine	CH_5N	266.8(0.3)	430.6(0.6)	7.61(0.09)	139(1)	130, 193, 322, 323
<i>N</i> -Methylaniline	C_7H_9N	470(1)	702(5)	5.2(0.6)	347(16)	117
2-Methylaniline	C_7H_9N	473.1(0.4)	710(1)	3.6(0.1)	377(52)	519
3-Methylaniline	C_7H_9N	476.4(0.5)	709(10)	4.6(0.5)	346(17)	106
4-Methylaniline	C_7H_9N	474(1)	667(10)	3.3(0.7)	334(16)	106
2-Methylanisole	$C_8H_{10}O$	446(2)	662(1)	3.6(0.3)	397(18)	162
3-Methylanisole	$C_8H_{10}O$	450(2)	665(1)	3.6(0.3)	449(22)	162
4-Methylanisole	$C_8H_{10}O$	448(2)	667(1)	3.6(0.4)	396(15)	162
α -Methylbenzenemethanol	$C_8H_{10}O$	478(4)	699(5)	3.8(0.7)	399(16)	32
Methyl benzoate	$C_8H_8O_2$	472(2)	702(1)	3.8(0.1)	408(11)	517
2-Methylbutanal	$C_5H_{10}O$	363(2)	531.6(0.1)	4.04(0.03)	318(10)	124
Methyl butanoate	$C_5H_{10}O_2$	375.0(0.1)	554.4(0.1)	3.49(0.08)	341(20)	19, 413, 416
3-Methylbutanoic acid	$C_5H_{10}O_2$	449.6(0.2)	629(1)	3.4(0.2)	355(10)	146
2-Methyl-1-butanol	$C_5H_{12}O$	402.1(0.4)	575.4(0.5)	3.9(0.1)	342(8)	254
2-Methyl-2-butanol	$C_5H_{12}O$	375.6	544(1)	3.71(0.05)	326(9)	147, 254
3-Methyl-1-butanol	$C_5H_{12}O$	403.9(0.3)	579(2)	3.9(0.3)	335(7)	21, 91, 222, 254, 511, 520
3-Methyl-2-butanol	$C_5H_{12}O$	386.8(0.4)	556.1(0.5)	3.9(0.4)	336(9)	254
3-Methyl-2-butanone	$C_5H_{10}O$	367.3(0.2)	553.1(0.3)	3.83(0.1)	321(33)	157, 166
2-Methyl-2-butene	C_5H_{10}	311.6(0.4)	470(1)	3.4(0.1)	299(7)	521
3-Methyl-1-butene	C_5H_{10}	293.2(0.2)	452.7(0.5)	3.51(0.04)	305(8)	490
Methyl <i>tert</i> -butyl ether	$C_5H_{12}O$	328.2(0.1)	497.0(0.6)	3.41(0.05)	335(10)	25, 440
Methylcyclohexane	C_7H_{14}	374.0(0.1)	572.3(0.2)	3.48(0.09)	368(3)	34, 74, 78, 86, 88, 171, 235, 236
Methylcyclopentane	C_6H_{12}	344.9(0.2)	532.78(0.05)	3.79(0.05)	322(2)	78, 235, 236
2-Methylcyclopentanone	$C_6H_{10}O$	413(3)	631(2)	4.0(0.6)	328(17)	162
2-Methyl- <i>N,N</i> -dimethylaniline	$C_9H_{13}N$	458(2)	668.0(0.7)	3.12(0.08)	466(15)	39, 40
Methyl dodecanoate	$C_{13}H_{26}O_2$	540(2)	712(5)	1.4(0.4)	842(9)	218
1,1'-Methylenebis[(1-methylethyl)benzene]	$C_{19}H_{24}$	592(36)	795(8)	1.6(0.1)	871(30)	394
Methyl formate	$C_2H_4O_2$	304.8(0.3)	487.16(0.1)	6.01(0.01)	172(6)	19, 412, 413
2-Methylfuran	C_5H_6O	337.0(0.2)	528(3)	4.77(0.08)	252(3)	290
2-Methylheptane	C_8H_{18}	390.8(0.9)	559.6(0.1)	2.50(0.02)	487(12)	86, 346, 522
3-Methylheptane	C_8H_{18}	392.0(0.6)	563.7(0.4)	2.54(0.02)	463(12)	346
4-Methylheptane	C_8H_{18}	390.8(0.5)	561.7(0.4)	2.54(0.02)	480(14)	346
Methyl heptanoate	$C_8H_{16}O_2$	442.8(0.4)	628(2)	2.6(0.4)	521(8)	162
4-Methyl-3-heptanol	$C_8H_{18}O$	430(2)	623.5(0.7)	2.8(0.4)	505(13)	437
5-Methyl-3-heptanol	$C_8H_{18}O$	427(2)	621.2(0.3)	2.8(0.3)	493(13)	437
2-Methyl-3-heptanone	$C_8H_{16}O$	431	615(1)	2.7(0.3)	487(11)	162
5-Methyl-3-heptanone	$C_8H_{16}O$	432(4)	619(4)	2.7(0.7)	484(11)	162
2-Methyl-1-heptene	C_8H_{16}	392(2)	567.5(0.9)	2.6(0.2)	466(10)	170
2-Methyl-2-heptene	C_8H_{16}	395(2)	569(1)	2.6(0.4)	465(11)	170
2-Methylhexane	C_7H_{16}	363.1(0.8)	530.4(0.1)	2.73(0.03)	420(15)	86, 346, 349, 522
3-Methylhexane	C_7H_{16}	365.0(0.1)	535.4(0.5)	2.82(0.06)	405(19)	346, 349
2-Methyl-3-hexanone	$C_7H_{14}O$	407(3)	593(1)	2.9(0.4)	428(13)	162
5-Methyl-2-hexanone	$C_7H_{14}O$	412(2)	604(1)	2.9(0.3)	434(13)	162
2-Methyl-1-hexene	C_7H_{14}	365(2)	542(1)	2.9(0.3)	407(8)	170
5-Methyl-1-hexene	C_7H_{14}	358(1)	528.7(0.4)	2.9(0.2)	410(9)	170
<i>N</i> -Methylhexylamine	$C_7H_{17}N$	418(8)	592(1)	2.8(0.8)	458(20)	170
Methyl isobutanoate	$C_5H_{10}O_2$	365(1)	540.7(0.5)	3.43(0.01)	341(42)	19, 413
Methyl methacrylate	$C_5H_8O_2$	373.8(0.2)	540.3(0.6)	2.97(0.06)	320(6)	51
1-Methylnaphthalene	$C_{11}H_{10}$	517.5(0.9)	771(5)	3.56(0.07)	479(22)	86, 218, 523
2-Methylnaphthalene	$C_{11}H_{10}$	514.2(0.3)	761(3)	3.37(0.06)	464(20)	218
2-Methyloctane	C_9H_{20}	416(1)	582.8(0.2)	2.30(0.02)	547(17)	345, 522
Methyloxirane	C_3H_6O	308	488.11(0.08)	5.44(0.02)	197(42)	290, 524

Name	Mol. Form.	T_b/K	T_c/K	P_c/MPa	$V_c/cm^3 mol^{-1}$	Ref.
Methyl pentafluoroethyl ether	$C_3H_3F_5O$	278.8(0.9)	406.81(0.05)	2.89(0.01)	301(5)	176, 448, 525
2-Methylpentane	C_6H_{14}	333.36(0.09)	497.9(0.2)	3.03(0.01)	371(2)	86, 127, 252, 325–328, 487, 522
3-Methylpentane	C_6H_{14}	336.4(0.5)	504.6(0.2)	3.12(0.01)	368.7(0.3)	252, 324–328, 526
Methyl pentanoate	$C_6H_{12}O_2$	400.51(0.06)	588.9(0.3)	3.20(0.05)	398(6)	164, 165
2-Methyl-1-pentanol	$C_6H_{14}O$	430(6)	604.4(0.5)	3.4(0.2)	410(8)	254
2-Methyl-2-pentanol	$C_6H_{14}O$	394(1)	559.5(0.7)	3.6(0.4)	410(11)	437
2-Methyl-3-pentanol	$C_6H_{14}O$	401.0(0.2)	576(1)	3.5(0.1)	380(9)	254
3-Methyl-3-pentanol	$C_6H_{14}O$	402(4)	575.6(0.6)	3.5(0.2)	376(10)	254
4-Methyl-1-pentanol	$C_6H_{14}O$	424(2)	603.5(0.7)	3.4(0.4)	406(7)	437
4-Methyl-2-pentanol	$C_6H_{14}O$	405.1(0.5)	574.4(0.5)	4(2)	389(9)	437
4-Methyl-2-pentanone	$C_6H_{12}O$	388.8(0.2)	575.4(1)	3.4(0.1)	378(12)	166, 527
2-Methyl-2-pentene	C_6H_{12}	340.4(0.5)	509.3(0.5)	3.26(0.01)	348(7)	170
4-Methyl-1-pentene	C_6H_{12}	327(2)	493.1(0.5)	3.18(0.07)	348(77)	170, 497
4-Methyl- <i>cis</i> -2-pentene	C_6H_{12}	329.5(0.1)	496.3(0.7)	3.24(0.01)	350(7)	170
Methyl pentyl ether	$C_6H_{14}O$	372(3)	546.5(0.2)	3.04(0.1)	395(14)	173, 174
2-Methyl-1,3-propanediol	$C_4H_{10}O_2$	494(4)	708(2)	5.4(0.4)	300(12)	55
Methyl propanoate	$C_4H_8O_2$	351.8(0.2)	530.57(0.1)	4.0(0.2)	280(98)	19, 412, 413, 415
2-Methylpropanoic acid	$C_4H_8O_2$	427.5(0.2)	605(2)	3.7(0.3)	296(10)	146
2-Methyl-1-propanol	$C_4H_{10}O$	380.99(0.07)	548(2)	4.30(0.04)	274(17)	25, 91, 148, 153, 155
2-Methyl-2-propanol	$C_4H_{10}O$	355.4(0.1)	506.2(0.1)	3.98(0.07)	283(4)	153
Methyl propyl ether	$C_4H_{10}O$	312(1)	476.2(0.2)	3.80(0.01)	281(7)	25
2-Methylpyridine	C_6H_7N	402.5(0.2)	622(1)	4.62(0.04)	306(119)	75, 528
3-Methylpyridine	C_6H_7N	417.2(0.1)	644.8(0.6)	4.63(0.03)	302(30)	90, 528
4-Methylpyridine	C_6H_7N	418.4(0.1)	645.8(0.5)	4.68(0.04)	316(67)	75, 90
<i>N</i> -Methyl-2-pyrrolidinone	C_5H_9NO	477.3(0.3)	721.7(0.4)	4.5(0.4)	330(12)	144, 242
2-Methylquinoline	$C_{10}H_9N$	520.5(0.4)	778(2)	3.91(0.02)	447(49)	530
8-Methylquinoline	$C_{10}H_9N$	520.5(0.7)	787(2)	4.22(0.02)	426(124)	530
Methyl salicylate	$C_8H_8O_3$	495.8(0.5)	709(30)	4.4(0.7)	436(17)	302
2-Methyltetrahydrofuran	$C_5H_{10}O$	353(1)	537(2)	3.74(0.06)	292(4)	290
(Methylthio)benzene	C_6H_8S	467.4(0.2)	706(4)	4.1(0.1)	374(13)	99
Methyl trifluoromethyl ether	$C_2H_3F_3O$	247.9(0.7)	377.92(0.06)	3.64(0.03)	219(2)	447, 525
Methyltris(trimethylsiloxy)silane	$C_{10}H_{30}O_3Si_4$	464.3(0.2)	597.4(0.2)	1.23(0.02)	1089(74)	34
4-Morpholinecarboxaldehyde	$C_5H_9NO_2$	511(1)	779(4)	5.0(0.4)	326(14)	54
Naphthalene	$C_{10}H_8$	491.1(0.1)	748.3(0.4)	4.06(0.04)	408(21)	79, 82, 86, 241, 251, 355, 451, 531
Neopentane	C_5H_{12}	282.65(0.06)	433.71(0.01)	3.20(0.01)	311.6(0.7)	532
Nitromethane	CH_3NO_2	374.3(0.1)	588(3)	6.0(0.2)	175(2)	533, 534
Nonadecane	$C_{19}H_{40}$	603(3)	756(5)	1.16(0.07)	1216(43)	358, 359
1-Nonadecene	$C_{19}H_{38}$	604(17)	755(8)	1.2(0.2)	1196(36)	365
1,1,1,2,2,3,3,4,4-Nonafluorohexan-5-one	$C_6H_3F_9O$	360.47	498.97(0.08)	2.20(0.02)	504(2)	308
Nonanal	$C_9H_{18}O$	468(3)	658(2)	2.7(0.1)	546(10)	242
Nonane	C_9H_{20}	424.0(0.2)	594.2(0.5)	2.29(0.05)	547(23)	34, 36, 78, 86, 131, 243–245, 247, 249, 251, 454, 456, 535, 536
1,9-Nonanediamine	$C_9H_{22}N_2$	531.8	726(7)	2.6(0.3)	600(23)	140
Nonanedioic acid	$C_9H_{16}O_4$	630.2	844(13)	2.7(0.2)	586(17)	253
Nonanoic acid	$C_9H_{18}O_2$	529(1)	712(3)	2.3(0.8)	592(16)	146
1-Nonanol	$C_9H_{20}O$	486.8(0.4)	670.6(0.5)	2.54(0.07)	555(75)	149, 152, 254, 255
2-Nonanol	$C_9H_{20}O$	466.7	649(1)	2.53(0.1)	575(11)	149, 255
3-Nonanol	$C_9H_{20}O$	468	648.0(0.3)	2.5(0.3)	577(12)	255
4-Nonanol	$C_9H_{20}O$	465.7	645.1(0.3)	2.5(0.3)	577(12)	255
2-Nonanone	$C_9H_{18}O$	467(1)	652.1(0.7)	2.5(0.1)	560(8)	49, 466
3-Nonanone	$C_9H_{18}O$	460(4)	648(4)	2.4(0.6)	560(7)	466
4-Nonanone	$C_9H_{18}O$	461(4)	643.7(0.3)	2.4(0.3)	560(7)	466
5-Nonanone	$C_9H_{18}O$	461.5(0.3)	641.4(0.3)	2.35(0.02)	560(7)	466
1-Nonene	C_9H_{18}	420.0(0.6)	594(1)	2.38(0.01)	529(2)	257
Octacosane	$C_{28}H_{58}$	705(6)	824(8)	0.8(0.1)	1916(65)	358
Octadecane	$C_{18}H_{38}$	589(2)	748(1)	1.3(0.1)	1167(41)	244
1-Octadecanol	$C_{18}H_{38}O$	624(2)	790(8)	1.28(0.1)	1157(18)	361
1-Octadecene	$C_{18}H_{36}$	588.7(1)	748(8)	1.3(0.1)	1119(34)	365

Name	Mol. Form.	T_b/K	T_c/K	P_c/MPa	$V_c/cm^3 mol^{-1}$	Ref.
1,1,1,2,2,3,3,4-Octafluorobutane	$C_4H_2F_8$	300.62(0.02)	432.0(0.1)	2.80(0.02)	360(22)	537
1,2,2,3,3,4,4,5-Octafluoro-1-pentanol	$C_5H_4F_8O$	413.1(0.3)	571(1)	2.9(0.1)	440(24)	362
Octafluorotetrahydrofuran	C_4F_8O	272.3(0.5)	399.6(0.7)	2.68(0.09)	350(30)	450
Octamethylcyclotetrasiloxane	$C_8H_{24}O_4Si_4$	448.5(0.9)	585.8(0.9)	1.33(0.01)	1006(67)	114, 538
Octamethyltrisiloxane	$C_8H_{24}O_2Si_3$	425.6(0.8)	564.1(0.2)	1.42(0.01)	882(16)	539
Octanal	$C_8H_{16}O$	447(3)	639.3(0.3)	3.0(0.3)	489(7)	163, 242
Octane	C_8H_{18}	398.8(0.1)	568.7(0.1)	2.48(0.01)	490(22)	11, 19, 33, 34, 36, 78, 83, 86, 91, 96, 131, 133, 134, 243–245, 247, 249, 318, 319, 324, 346, 465, 469, 535, 540, 541
1,8-Octanediamine	$C_8H_{20}N_2$	498.7	712(7)	2.8(0.3)	547(20)	140
Octanedioic acid	$C_8H_{14}O_4$	618.6	843(13)	3.0(0.2)	520(16)	253
Octanenitrile	$C_8H_{15}N$	475(3)	674.4(0.4)	2.85(0.03)	494(10)	35
Octanoic acid	$C_8H_{16}O_2$	513(1)	694(1)	2.9(0.3)	522(19)	145, 146, 360, 464
1-Octanol	$C_8H_{18}O$	467.8(0.8)	651(2)	2.80(0.07)	490(47)	25, 149, 152, 254, 255
2-Octanol	$C_8H_{18}O$	452.5	629.5(0.9)	2.75(0.04)	519(10)	149, 254, 255
3-Octanol	$C_8H_{18}O$	457(6)	628.4(0.3)	2.8(0.4)	515(10)	255
4-Octanol	$C_8H_{18}O$	449.5	625.1(0.3)	2.8(0.3)	516(10)	255
2-Octanone	$C_8H_{16}O$	446(3)	632.7(0.2)	2.7(0.5)	497(6)	466
3-Octanone	$C_8H_{16}O$	439(4)	627.7(0.2)	2.7(0.3)	497(6)	466
4-Octanone	$C_8H_{16}O$	439(3)	623.8(0.2)	2.7(0.3)	497(6)	466
1-Octene	C_8H_{16}	398.0(0.5)	566.58(0.05)	2.68(0.02)	464(2)	86, 257
<i>trans</i> -2-Octene	C_8H_{16}	395.5(0.5)	569.8(0.4)	2.58(0.09)	471(9)	170
<i>trans</i> -4-Octene	C_8H_{16}	394.4(0.2)	566(1)	2.55(0.06)	472(9)	170
Octylamine	$C_8H_{19}N$	451.8(0.2)	641(1)	2.82(0.03)	494(41)	542
Octylbenzene	$C_{14}H_{22}$	536(2)	725(7)	2.0(0.2)	746(17)	258
Octyl orthosilicate	$C_{32}H_{68}O_4Si$		812(16)			57
Oxazole	C_3H_3NO	342.6(0.2)	551(4)	6.8(0.2)	185(23)	54
Oxirane	C_2H_4O	283.5(0.1)	469(1)	7.2(0.2)	138(4)	543, 544
Paraldehyde	$C_6H_{12}O_3$	397(2)	563(10)	4(3)	410(15)	12
Pentacene	$C_{22}H_{14}$		1115(47)		806(23)	545
1 <i>H</i> -Pentadecafluoroheptane	C_7HF_{15}	368(2)	495.8(0.7)	1.7(0.5)	644(38)	546
Pentadecane	$C_{15}H_{32}$	543.8(0.4)	707(2)	1.54(0.09)	938(36)	243, 244, 245, 247, 249, 363, 547
Pentadecanoic acid	$C_{15}H_{30}O_2$	612(4)	777(8)	1.6(0.2)	1002(22)	360
1-Pentadecanol	$C_{15}H_{32}O$	591(2)	757(8)	1.6(0.2)	961(16)	361
1-Pentadecene	$C_{15}H_{30}$	541.5(0.4)	705(7)	1.56(0.05)	933(30)	365
Pentafluorobenzene	C_6HF_5	358(3)	530.93(0.03)	3.53(0.01)	322(22)	472, 548
3,3,4,4,4-Pentafluoro-2-butanone	$C_4H_3F_5O$	314.36(0.04)	453(1)	2.90(0.06)	333(4)	110
Pentafluoroethane	C_2HF_5	224.65	339.2(0.2)	3.63(0.01)	210(3)	180, 309, 312, 315, 447, 448, 549–554
1,1,1,2,2-Pentafluoropentan-3-one	$C_5H_3F_5O$	335.24	475.5(0.1)	2.64(0.01)	356(1)	308
1,1,1,2,2-Pentafluoropropane	$C_3H_3F_5$	255.1(0.3)	380.1(0.4)	3.14(0.02)	273(3)	555
1,1,1,3,3-Pentafluoropropane	$C_3H_3F_5$	288.5	427.20(0.07)	3.66(0.02)	262(14)	203
1,1,2,2,3-Pentafluoropropane	$C_3H_3F_5$	298.2	447.57(0.06)	3.96(0.02)	258(13)	203
1,1,1,2,2-Pentafluoro-3-(1,1,2,2-tetrafluoroethoxy)propane	$C_5H_3F_9O$	343.4	473.0(0.1)	2.24(0.01)	457(2)	113
Pentafluoro(trifluoromethoxy)ethane	C_3F_8O	249.5(0.3)	356.8(0.1)	2.4(0.5)	319(5)	442
Pentafluoro(trifluoromethyl)sulfur	CF_8S	252.4(0.2)	381.2(0.1)	3.4(0.1)	284(4)	442
Pentanal	$C_5H_{10}O$	376(2)	567(3)	3.1(0.3)	313(11)	104, 124
Pentane	C_5H_{12}	309.21(0.07)	469.7(0.1)	3.37(0.01)	310(1)	19, 34, 36, 52, 64, 77, 84–86, 127, 131, 133, 134, 201, 202, 243, 244, 254, 286, 324, 359, 394, 456, 490, 556–562
Pentanedioic acid	$C_5H_8O_4$	546(10)	840(13)	4.3(0.5)	343(13)	253
Pentanenitrile	C_5H_9N	413(1)	610.3(0.2)	3.58(0.05)	320(8)	35
Pentanoic acid	$C_5H_{10}O_2$	459.2(0.3)	639(2)	3.6(0.1)	347(15)	11, 41, 145, 146, 464
1-Pentanol	$C_5H_{12}O$	410.8(0.4)	587.9(0.4)	3.9(0.3)	331(9)	25, 34, 149, 329, 456, 465, 563

Name	Mol. Form.	T_b/K	T_c/K	P_c/MPa	$V_c/cm^3 mol^{-1}$	Ref.
2-Pentanol	C ₅ H ₁₂ O	392.2(0.5)	560.4(0.2)	4.2(0.4)	340(3)	149, 254, 465
3-Pentanol	C ₅ H ₁₂ O	396(2)	559.6(0.3)	4.9(0.9)	325(2)	465
2-Pentanone	C ₅ H ₁₀ O	375.3(0.1)	561.0(0.2)	3.70(0.06)	324(4)	25, 172
3-Pentanone	C ₅ H ₁₀ O	375.0(0.1)	561.4(0.2)	3.73(0.07)	319(11)	25
1-Pentene	C ₅ H ₁₀	310.0(0.2)	464.74(0.04)	3.55(0.02)	301.0(0.1)	86, 470, 559
<i>cis</i> -2-Pentene	C ₅ H ₁₀	303.1(0.3)	474.9(0.4)	3.69(0.02)	301(7)	564
Pentyl acetate	C ₇ H ₁₄ O ₂	422.5(0.3)	600(2)	2.79(0.03)	466(123)	141, 162, 166
Pentylbenzene	C ₁₁ H ₁₆	476(3)	675(7)	2.6(0.3)	559(12)	258
Pentyl benzoate	C ₁₂ H ₁₆ O ₂	533(3)	736(14)	2.2(0.2)	661(14)	169
Pentyl formate	C ₆ H ₁₂ O ₂	399(3)	576(4)	3.5(0.8)	453(35)	415
Perfluoroacetone	C ₃ F ₆ O	245.7(0.4)	357.2(0.1)	2.85(0.01)	329(7)	197, 470
Perfluorobutane	C ₄ F ₁₀	271.0(0.8)	386.3(0.2)	2.33(0.02)	380(18)	565, 566
Perfluorocyclobutane	C ₄ F ₈	267.3	388.4(0.1)	2.78(0.01)	316(8)	265, 470, 567, 568
Perfluorocyclohexane	C ₆ F ₁₂	325.95 s	457.1(0.5)	2.24(0.03)	424(32)	569
Perfluorocyclohexene	C ₆ F ₁₀	324.8(0.1)	461.7(0.7)	2.6(0.1)	434(28)	546
Perfluorodecane	C ₁₀ F ₂₂	408(3)	542.4(0.4)	1.45(0.03)	892(7)	240, 570
Perfluorodimethoxymethane	C ₃ F ₈ O ₂	263.1(0.7)	372.4(0.2)	2.34(0.01)	370(30)	116, 571
Perfluoro-2,3-dimethylbutane	C ₆ F ₁₄	332.9(0.3)	463.0(0.1)	1.95(0.03)	523(18)	240, 572
Perfluoroethyl ethyl ether	C ₄ H ₅ F ₅ O	301(3)	431.23(0.08)	2.53(0.01)	366(3)	176
Perfluoroethyl 2,2,2-trifluoroethyl ether	C ₄ H ₂ F ₈ O	301.04	421.68(0.08)	2.33(0.01)	409(3)	176
Perfluoroheptane	C ₇ F ₁₆	355.6(0.2)	477(3)	1.63(0.01)	603(14)	232, 566, 570, 573–576
Perfluoro-1-heptene	C ₇ F ₁₄	354(4)	478.2(0.7)	1.7(0.3)	555(34)	546
1 <i>H</i> -Perfluorohexane	C ₆ HF ₁₃	345(4)	471.8(0.7)	2.0(0.3)	504(32)	546
Perfluorohexane	C ₆ F ₁₄	330.3(0.2)	451(3)	1.88(0.02)	552(77)	198, 470, 546, 570, 572, 573, 577
Perfluoro-1-hexene	C ₆ F ₁₂	330.2	454.3(0.7)	1.9(0.7)	462(31)	546
Perfluoroisobutane	C ₄ F ₁₀	273	395.4(0.7)		396(22)	578
Perfluoroisopentane	C ₅ F ₁₂	303.26(0.03)	423(19)	2.12(0.02)	465(8)	240
Perfluoroisopropyl methyl ether	C ₄ H ₃ F ₇ O	302(1)	433.30(0.08)	2.55(0.01)	369(3)	176
Perfluoromethylcyclohexane	C ₇ F ₁₄	349.4(0.2)	486.5(1)	2.02(0.01)	561(4)	325, 326, 327, 566, 569
Perfluoromethylcyclopentane	C ₆ F ₁₂	321.58(0.01)	451.43(0.04)	2.17(0.01)	419(37)	579
Perfluoro-2-methylpentane	C ₆ F ₁₄	330.8(0.3)	454.6(0.2)	1.87(0.02)	585(10)	572, 580
Perfluoro-3-methylpentane	C ₆ F ₁₄	331(9)	450(1)	1.69(0.01)	511(36)	572
Perfluoronaphthalene	C ₁₀ F ₈	473(8)	673(1)	2.9(0.6)	464(29)	546
Perfluorononane	C ₉ F ₂₀	390(3)	524.0(0.1)	1.56(0.04)	846(48)	570
Perfluorooctane	C ₈ F ₁₈	378(2)	502.3(0.1)	1.66(0.02)	738(17)	34, 240, 570, 573
Perfluorooxetane	C ₃ F ₆ O	244.5(0.6)	361.8(0.5)	3.10(0.02)	274(24)	116, 571
1 <i>H</i> -Perfluoropentane	C ₅ HF ₁₁	319(2)	443.9(0.7)	2.2(0.3)	413(28)	546
Perfluoropentane	C ₅ F ₁₂	302.3(0.2)	421.8(0.1)	2.04(0.02)	463(7)	570, 573
Perfluoropropane	C ₃ F ₈	236.3(0.3)	345.03(0.08)	2.67(0.01)	301(12)	187, 305, 470, 581
Perfluoropropyl methyl ether	C ₄ H ₃ F ₇ O	307(1)	437.7(0.1)	2.48(0.01)	382(3)	176
Perfluorotoluene	C ₇ F ₈	377.8(0.3)	534.4(0.2)	2.70(0.02)	423(26)	25
Perfluorotributylamine	C ₁₂ F ₂₇ N	451(2)	566(4)	1.24(0.09)	1196(66)	582
Perfluorovaleric acid	C ₅ HF ₉ O ₂	415.3(0.3)	545.5(0.2)	2.10(0.05)	496(29)	583
Phenol	C ₆ H ₆ O	455.0(0.1)	694.3(0.1)	5.5(0.2)	283(10)	218, 584
Phenyl acetate	C ₈ H ₈ O ₂	468(1)	686(2)	3.60(0.06)	407(12)	55
4-Phenyl-1-butanol	C ₁₀ H ₁₄ O	537(3)	746(14)	3.1(0.2)	493(19)	584
Phenyl isocyanate	C ₇ H ₅ NO	439.4(0.4)	657(10)	3.6(0.1)	342(15)	54
3-Phenyl-1-propanol	C ₉ H ₁₂ O	514(4)	732(14)	3.4(0.5)	450(14)	584
(1 <i>S</i>)-(-)- α -Pinene	C ₁₀ H ₁₆	429.1(0.4)	644(2)	3.4(0.2)	472(10)	496
Piperazine	C ₄ H ₁₀ N ₂	421.78(0.05)	660(5)	5.4(0.4)	283(16)	331, 489
Piperidine	C ₅ H ₁₁ N	379.34(0.09)	594.14(0.02)	4.7(0.1)	294(10)	39, 218
Propanal	C ₃ H ₆ O	321.1(0.2)	503.7(0.8)	5.04(0.03)	218(9)	104, 124, 163, 242
Propane	C ₃ H ₈	231.04(0.09)	369.9(0.1)	4.25(0.01)	199(6)	29, 125, 127, 131–134, 252, 322, 323, 336, 371, 387, 470, 486, 509, 540, 585–596
1,3-Propanediamine	C ₃ H ₁₀ N ₂	412.3(0.7)	632(7)	5.7(0.6)	257(13)	140
1,2-Propanediol	C ₃ H ₈ O ₂	460.4(0.2)	676(1)	5.9(0.2)	237(10)	32
1,3-Propanediol	C ₃ H ₈ O ₂	487.8(0.3)	718(2)	6.7(0.2)	255(12)	55

Name	Mol. Form.	T_b/K	T_c/K	P_c/MPa	$V_c/cm^3 mol^{-1}$	Ref.
Propanenitrile	C_3H_5N	370.4(0.4)	561.3(0.2)	4.26(0.07)	211(7)	35
1-Propanethiol	C_3H_8S	340.8(0.1)	536.6(0.6)	4.7(0.1)	286(11)	143, 144
Propanoic acid	$C_3H_6O_2$	414.6(0.2)	603(3)	4.5(0.7)	232(12)	39, 145, 146, 493, 597, 598
1-Propanol	C_3H_8O	370.19(0.09)	536.8(0.2)	5.1(0.1)	220(22)	19, 34, 80, 148–150, 152, 153, 599, 600
2-Propanol	C_3H_8O	355.36(0.09)	508.3(0.2)	4.7(0.1)	226(1)	25, 91, 148, 149, 153, 324, 440, 520, 601, 602
Propene	C_3H_6	225.5(0.1)	364.9(0.5)	4.59(0.02)	184(11)	67, 123, 172, 181, 427, 603–610
2-Propoxyethanol	$C_5H_{12}O_2$	425(3)	614.7(0.7)	3.65(0.09)	364(13)	11, 41
1-Propoxy-2-propanol	$C_6H_{14}O_2$	423.3(0.7)	605(1)	3.1(0.1)	417(16)	32
Propyl acetate	$C_5H_{10}O_2$	374.1(0.2)	549.69(0.08)	3.37(0.07)	346(24)	19, 411, 413, 415
Propylamine	C_3H_9N	320.36(0.08)	499.2(0.4)	4.77(0.06)	230(12)	170, 193
Propylbenzene	C_9H_{12}	432.3(0.5)	638.3(0.1)	3.20(0.02)	441(6)	79, 86, 88
Propyl benzoate	$C_{10}H_{12}O_2$	504(2)	710(14)	2.6(0.3)	530(10)	169
Propyl butanoate	$C_7H_{14}O_2$	417(2)	593(1)	2.72(0.06)	463(10)	162, 166
Propylcyclohexane	C_9H_{18}	429.8(0.3)	630.8(0.9)	2.87(0.04)	489(13)	170, 171
Propylene carbonate	$C_4H_6O_3$	514.8(0.7)	763(2)	4.1(0.2)	256.5(0.1)	55
1,2-Propylene glycol 1- <i>tert</i> -butyl ether	$C_7H_{16}O_2$	425.2	601(4)	2.7(0.1)	468(15)	99
1,2-Propylene glycol monomethyl ether acetate	$C_6H_{12}O_3$	419.1(0.4)	598(1)	3.1(0.2)	432(16)	41, 144
Propyl formate	$C_4H_8O_2$	353.8(0.2)	538.1(0.1)	4.07(0.02)	281(31)	19, 412, 413, 415
Propyl isobutanoate	$C_7H_{14}O_2$	407(4)	582(11)	3(1)	463(9)	162, 167
Propyl 3-methylbutanoate	$C_8H_{16}O_2$	428(3)	609(6)	2.5(0.8)	523(9)	147
Propyl propanoate	$C_6H_{12}O_2$	395.3(0.1)	569(3)	3.1(0.1)	403(6)	162, 166
Propyne	C_3H_4	250	402(2)	5.63(0.06)	160(10)	605, 611
Pyrazine	$C_4H_4N_2$	389.4(0.1)	627(1)	6.49(0.03)	225(16)	612
Pyridine	C_5H_5N	388.3(0.1)	619(2)	5.63(0.07)	248(12)	16, 17, 60, 90, 246, 290, 520, 613
Pyrrole	C_4H_5N	402.89(0.04)	639.7(0.2)	8.0(0.2)	222(15)	241
Pyrrolidine	C_4H_9N	359.8(0.1)	568.6(0.2)	5.69(0.08)	259(3)	241, 290
Quinoline	C_9H_7N	510.2(0.5)	782(3)	4.75(0.1)	382(17)	218
Resorcinol	$C_6H_6O_2$	553(2)	836(10)	6.3(0.3)	292(10)	489
Stearic acid	$C_{18}H_{36}O_2$	644(3)	803(8)	1.3(0.2)	1251(27)	360
Styrene	C_8H_8	418.4(0.6)	635(2)	3.9(0.2)	357(15)	481
Succinic acid	$C_4H_6O_4$	507(3)	851(20)		308(21)	253
<i>m</i> -Terphenyl	$C_{18}H_{14}$	648(1)	883(7)	2.2(0.2)	747(37)	108, 358
<i>o</i> -Terphenyl	$C_{18}H_{14}$	610(5)	857(6)	2.9(0.1)	737(37)	108
<i>p</i> -Terphenyl	$C_{18}H_{14}$	649	913(22)	2.5(0.5)	713(37)	108, 614
Tetrabutyl silicate	$C_{16}H_{36}O_4Si$		682(14)			57
Tetrachloromethane	CCl_4	349.8(0.2)	556.5(0.3)	4.57(0.07)	276(9)	19, 27, 469, 615–619
Tetracosane	$C_{24}H_{50}$	664(5)	800(5)	0.9(0.1)	1585(55)	358, 359
Tetradecamethylcycloheptasiloxane	$C_{14}H_{42}O_7Si_7$	548.4(0.2)	683.2(0.2)	0.99(0.02)	1634(110)	34
Tetradecane	$C_{14}H_{30}$	526.6(0.4)	693(1)	1.56(0.08)	870(49)	34, 244
Tetradecanedioic acid	$C_{14}H_{26}O_4$	639(10)		1.9(0.2)		253
Tetradecanoic acid	$C_{14}H_{28}O_2$	599(1)	763(8)	1.6(0.2)	921(20)	360
1-Tetradecanol	$C_{14}H_{30}O$	569.0(0.4)	743(7)	1.70(0.04)	887(15)	361
2-Tetradecanone	$C_{14}H_{28}O$	562(6)	728(9)	1.6(0.5)	896(26)	256
3-Tetradecanone	$C_{14}H_{28}O$	552(7)	727(6)	1.6(0.5)	896(26)	256
4-Tetradecanone	$C_{14}H_{28}O$	552(7)	725(6)	1.6(0.5)	900(27)	256
7-Tetradecanone	$C_{14}H_{28}O$	552(8)	723(8)	1.6(0.6)	904(27)	256
1-Tetradecene	$C_{14}H_{28}$	524.2(0.4)	691(7)	1.58(0.07)	851(24)	365
Tetradecyl orthosilicate	$C_{40}H_{84}O_4Si$		849(16)			57
Tetraethoxysilane	$C_8H_{20}O_4Si$	441(1)	587(12)	2.0(0.4)	701.3(0.2)	57
Tetraethylene glycol	$C_8H_{18}O_5$	588(7)	800(30)	2.8(0.7)	608(31)	283
Tetraethylsilane	$C_8H_{20}Si$	426.5(0.7)	606(2)	2.297(0.01)	596.4(0.2)	620
1,2,3,4-Tetrafluorobenzene	$C_6H_2F_4$	367.4(0.8)	550.8(0.2)	3.791(0.01)	312(22)	25
1,2,3,5-Tetrafluorobenzene	$C_6H_2F_4$	357.4(0.8)	535.2(0.2)	3.75(0.01)	311(22)	25
1,2,4,5-Tetrafluorobenzene	$C_6H_2F_4$	363.3(0.3)	543.3(0.2)	3.80(0.01)	309(22)	25
1,1,2,2-Tetrafluoro-2-(2,2-difluoromethoxy)ethane	$C_4H_4F_6O$	352.13	501.08(0.08)	3.09(0.02)	356(1)	113

Name	Mol. Form.	T_b/K	T_c/K	P_c/MPa	$V_c/cm^3 mol^{-1}$	Ref.
1,1,1,2-Tetrafluoroethane	$C_2H_2F_4$	247.0(0.1)	374.2(0.2)	4.06(0.01)	200(2)	113, 182, 183, 188, 278–280, 305, 313, 315, 447, 448, 478, 621–631
1,1,2,2-Tetrafluoroethane	$C_2H_2F_4$	253(1)	391.75(0.08)	4.61(0.01)	192(1)	178, 632
Tetrafluoroethene	C_2F_4	197(1)	307(1)	3.94(0.05)	183(3)	633, 634
1,2,2,2-Tetrafluoroethyl difluoromethyl ether	$C_3H_2F_6O$	296(2)	428.95(0.08)	3.05(0.01)	315(2)	176
1,1,2,2-Tetrafluoroethyl 1,1,1-trifluoroethyl ether	$C_4H_3F_7O$	329.37	463.89(0.07)	2.71(0.01)	373(1)	113
Tetrafluoromethane	CF_4	145.2(0.1)	227.54(0.03)	3.73(0.03)	140(1)	506, 635
1,1,2,2-Tetrafluoro-3-methoxypropane	$C_4H_6F_4O$	347.4(0.1)	505.4(0.1)	3.28(0.01)	331(1)	113
1,2,2,3-Tetrafluoro-1-propanol	$C_3H_4F_4O$	386.4(0.4)	554(2)	3.3(0.2)	280(15)	362
1,1,2,2-Tetrafluoro-3-(1,1,2,2-tetrafluoroethoxy)propane	$C_5H_4F_8O$	366.32	510.07(0.08)	2.58(0.01)	440(2)	113
1,1,1,2-Tetrafluoro-2-(trifluoromethoxy)ethane	C_3HF_7O	264(2)	377.26(0.06)	2.62(0.01)	321(2)	525
3,4,4,4-Tetrafluoro-3-(trifluoromethyl)-2-butanone	$C_5H_3F_7O$	328.76(0.05)	468(1)	2.50(0.06)	409(5)	110
4,4,5,5-Tetrafluoro-2-(trifluoromethyl)-1,3-dioxolane	$C_4HF_7O_2$	304.6(0.1)	435(1)	2.62(0.07)	376(4)	110
Tetrahexoxysilane	$C_{24}H_{30}O_4Si$		757(16)			57
Tetrahydrofuran	C_4H_8O	339.1(0.1)	540(1)	5.29(0.06)	223(2)	23, 241, 290
1,2,3,4-Tetrahydronaphthalene	$C_{10}H_{12}$	480.3(0.3)	720(1)	3.6(0.1)	431(40)	144, 242, 636
Tetrahydropyran	$C_5H_{10}O$	361.1(0.4)	572.0(0.3)	4.8(0.2)	278(18)	75
Tetrahydrothiophene	C_4H_8S	394.2(0.2)	632.0(0.2)	5.4(0.6)	276(19)	143, 241
1,2,4,5-Tetraisopropylbenzene	$C_{18}H_{30}$	532	703(1)	1.65(0.02)	983(83)	317
Tetramethoxysilane	$C_4H_{12}O_4Si$	393.2(0.7)	558(12)	2.8(0.6)	464(22)	57
1,2,4,5-Tetramethylbenzene	$C_{10}H_{14}$	470(1)	676(2)	2.9(0.3)	489(11)	39
2,2,3,3-Tetramethylhexane	$C_{10}H_{22}$	433(2)	623.0(0.5)	2.51(0.06)	574(14)	536
2,2,5,5-Tetramethylhexane	$C_{10}H_{22}$	410(2)	581.4(0.5)	2.19(0.01)	600(14)	536
2,2,3,3-Tetramethylpentane	C_9H_{20}	413.3(0.4)	607.5(0.5)	2.74(0.03)	514(15)	536
2,2,3,4-Tetramethylpentane	C_9H_{20}	406.1(0.8)	592.6(0.5)	2.60(0.03)	517(17)	536
2,2,4,4-Tetramethylpentane	C_9H_{20}	395(1)	574.6(0.5)	2.49(0.01)	532(16)	536
2,3,3,4-Tetramethylpentane	C_9H_{20}	414.6(0.7)	607.5(0.5)	2.72(0.04)	517(17)	536
Tetramethylsilane	$C_4H_{12}Si$	299.8(0.5)	449(2)	2.82(0.01)	362(7)	637–639
Tetramethylstannane	$C_4H_{12}Sn$	350(1)	521.77(0.02)	2.98(0.01)	109(109)	640, 641
Tetranoxysilane	$C_3H_7O_4Si$		830(16)			57
Tetrapropyl silicate	$C_{12}H_{28}O_4Si$		649(12)			57
Thiacyclohexane	$C_5H_{10}S$	414.88(0.04)	684(44)	6.50(0.07)	284(11)	104
Thiobis(trifluoromethane)	C_2F_4S	251.3(0.3)	376.8(0.1)	3.2(0.5)	216(19)	442
Thiophene	C_4H_4S	357.2(0.1)	579.4(0.2)	5.7(0.2)	230(3)	241, 290
Thymol	$C_{10}H_{14}O$	506(3)	698(10)	3(2)	528(21)	302
Toluene	C_7H_8	383.75(0.07)	591.9(0.2)	4.13(0.02)	314(7)	34, 63, 68, 74, 78, 79, 84, 86, 88, 90, 96, 163, 172, 258, 481, 636, 642–646
Triaccontane	$C_{30}H_{62}$	724(7)	843(8)	0.6(0.1)	2055(69)	358
Tribromomethane	$CHBr_3$	422.3(0.5)	682(1)	5.8(0.2)	261(12)	647
Trichloroacetyl chloride	C_2Cl_4O	391.3(0.3)	604(2)	4.21(0.03)	331(54)	410
Trichloroethylsilane	$C_2H_5Cl_3Si$	371.8(0.7)	559.9(0.6)	3.34(0.04)	403(5)	266
Trichlorofluoromethane	CCl_3F	296.8(0.6)	471.1(0.2)	4.40(0.03)	248.0(0.9)	33, 34, 191, 271, 571
Trichloromethane	$CHCl_3$	334.3(0.1)	536.0(0.4)	5.5(0.2)	237(7)	18, 27, 28, 117
Trichloromethylsilane	CH_2Cl_3Si	339(2)	517.7(0.3)	3.52(0.03)	329(9)	264, 648
1,3,5-Trichloro-2,4,6-trifluorobenzene	$C_6Cl_3F_3$	472(27)	684.7(0.4)	3.3(0.1)	443(27)	25
1,1,2-Trichloro-1,2,2-trifluoroethane	$C_2Cl_3F_3$	320.8(0.2)	487.4(0.2)	3.40(0.02)	325(1)	191, 265, 271, 649
Tricosane	$C_{23}H_{48}$	654(9)	790(8)	0.9(0.1)	1527(53)	359
Tridecane	$C_{13}H_{28}$	508.5(0.4)	676(1)	1.68(0.04)	824(30)	34, 244, 249
1-Tridecanol	$C_{13}H_{28}O$	560(8)	732(7)	1.8(0.2)	828(14)	361
2-Tridecanone	$C_{13}H_{26}O$	541(1)	717(6)	1.8(0.2)	820(24)	256
3-Tridecanone	$C_{13}H_{26}O$	539(7)	716(5)	1.7(0.5)	823(24)	256
4-Tridecanone	$C_{13}H_{26}O$	539(7)	712(6)	1.7(0.5)	823(24)	256
5-Tridecanone	$C_{13}H_{26}O$	539(8)	710(8)	1.7(0.7)	826(17)	256
6-Tridecanone	$C_{13}H_{26}O$	539(7)	709(5)	1.7(0.5)	826(24)	256
7-Tridecanone	$C_{13}H_{26}O$	539(9)	708(5)	1.7(0.5)	830(24)	256

Name	Mol. Form.	T_b/K	T_c/K	P_c/MPa	$V_c/cm^3 mol^{-1}$	Ref.
1-Tridecene	$C_{13}H_{26}$	506.0(0.7)	673(7)	1.74(0.05)	770(17)	365
Tridecylbenzene	$C_{19}H_{32}$	613(4)	790(8)	1.5(0.1)	1079(43)	258
Triethylamine	$C_6H_{15}N$	361.9(0.2)	535.6(0.3)	3.1(0.3)	392(25)	117, 286
1,3,5-Triethylbenzene	$C_{12}H_{18}$	489.0(0.9)	679(2)	2.32(0.01)	624(60)	332
Triethylene glycol	$C_6H_{14}O_4$	561.8(0.2)	775(30)	3.3(0.2)	454(25)	283
Trifluoroacetone	C_2F_3N	204.3(0.8)	311.1(0.4)	3.61(0.04)	202(4)	470
1,2,3-Trifluorobenzene	$C_6H_3F_3$	368	560.3(0.4)	4.1(0.4)	296(20)	119
1,2,4-Trifluorobenzene	$C_6H_3F_3$	363	551.1(0.4)	4.1(0.6)	297(20)	119
1,3,5-Trifluorobenzene	$C_6H_3F_3$	350.1(0.5)	530.9(0.4)	3.8(0.2)	300(20)	119
1,1,1-Trifluoroethane	$C_2H_3F_3$	225.9(0.1)	345.89(0.07)	3.77(0.01)	195(15)	179, 203, 309, 478, 479, 549, 624, 650–653
2,2,2-Trifluoroethanol	$C_2H_3F_3O$	346.9(0.3)	498.57(0.05)	4.81(0.01)	211(12)	654, 655
2,2,2-Trifluoroethyl methyl ether	$C_5H_5F_3O$	304.77	448.98(0.08)	3.51(0.06)	277(3)	176
Trifluoroiodomethane	CF_3I	251.3(0.6)	396.44(0.06)	3.95(0.01)	231(3)	656–659
Trifluoromethane	CHF_3	191.1(0.1)	299.00(0.02)	4.82(0.01)	133(1)	310, 606, 660–664
Trifluoromethyl difluoromethyl ether	C_2HF_5O	238.1(0.2)	354.49(0.06)	3.36(0.02)	226(20)	203
Trifluoromethyl 1,1,2,2-tetrafluoroethyl ether	C_3HF_7O	270(1)	387.8(0.5)	2.65(0.01)	341(22)	116, 571
3,3,3-Trifluoropropene	$C_3H_3F_3$	246(4)	378.6(0.5)	3.61(0.08)	229(14)	440
Trimethylamine	C_3H_9N	275.9(0.2)	433.0(0.6)	4.08(0.04)	254(6)	322, 323, 665
1,2,3-Trimethylbenzene	C_9H_{12}	449.1(0.4)	664.4(0.1)	3.45(0.03)	423(11)	79
1,2,4-Trimethylbenzene	C_9H_{12}	442.5(0.3)	649.1(0.1)	3.3(0.1)	436(12)	79, 86, 96, 251
1,3,5-Trimethylbenzene	C_9H_{12}	437.8(0.3)	637.31(0.1)	3.13(0.05)	435(12)	79
3,7,7-Trimethyl-bicyclo[4.1.0]hept-3-ene	$C_{10}H_{16}$	445(2)	658(2)	2.9(0.5)	487(10)	496
2,2,3-Trimethylbutane	C_7H_{16}	353.9(0.1)	531.3(0.5)	2.96(0.03)	401(13)	346, 349
Trimethylchlorosilane	C_3H_9ClSi	330.8(0.4)	497.7(0.6)	3.20(0.03)	366(6)	266
1 α ,3 α ,5 β -1,3,5-Trimethylcyclohexane	C_9H_{18}	414(2)	602(2)	2.6(0.3)	494(14)	74
3,3,5-Trimethylheptane	$C_{10}H_{22}$	430(3)	609.5(0.5)	2.32(0.05)	583(18)	536
2,2,5-Trimethylhexane	C_9H_{20}	397(2)	570(2)	2.46(0.03)	547(18)	350
2,2,3-Trimethylpentane	C_8H_{18}	382.9(0.4)	563.5(0.4)	2.73(0.02)	442(16)	346
2,2,4-Trimethylpentane	C_8H_{18}	372.3(0.2)	543.9(0.4)	2.57(0.02)	475(20)	33, 34, 86, 159, 346, 666, 667
2,3,3-Trimethylpentane	C_8H_{18}	387.8(0.3)	573.5(0.4)	2.82(0.03)	454(14)	346
2,3,4-Trimethylpentane	C_8H_{18}	386.5(0.3)	566.4(0.4)	2.72(0.02)	462(12)	346
<i>cis</i> -Tri(methylphenyl)trisiloxane	$C_{21}H_{24}O_3Si_3$		824(8)			353
<i>trans</i> -2,4,6-Trimethyl-2,4,6-triphenylcyclotrisiloxane	$C_{21}H_{24}O_3Si_3$		839(8)			353
Undecafluorocyclohexane	C_6HF_{11}	335.2	477.7(0.7)			546
Undecane	$C_{11}H_{24}$	469.0(0.3)	638.8(0.2)	2.01(0.03)	683(20)	34, 86, 131, 243, 249, 363, 535
Undecanoic acid	$C_{11}H_{22}O_2$	553	728(7)	2.1(0.2)	741(20)	360
1-Undecanol	$C_{11}H_{24}O$	519(2)	703.0(0.6)	2.15(0.07)	707(12)	149
2-Undecanone	$C_{11}H_{22}O$	506.2(0.3)	688(2)	2.08(0.01)	692(20)	256
3-Undecanone	$C_{11}H_{22}O$	500	685(2)	2.0(0.4)	692(20)	256
4-Undecanone	$C_{11}H_{22}O$	501(3)	681(2)	2.0(0.2)	692(20)	256
5-Undecanone	$C_{11}H_{22}O$	500	679(2)	2.0(0.2)	692(20)	256
6-Undecanone	$C_{11}H_{22}O$	500.5(0.5)	678(2)	2.02(0.01)	692(20)	256
Undecylbenzene	$C_{17}H_{28}$	585(3)	763(8)	1.6(0.1)	946(35)	258
Vinyl acetate	$C_4H_6O_2$	345.8(0.3)	519.2(0.2)	4.17(0.03)	269(7)	156, 440
<i>m</i> -Xylene	C_8H_{10}	412.2(0.4)	616.9(0.3)	3.54(0.01)	377(7)	34, 68, 79, 88, 90, 106, 668
<i>o</i> -Xylene	C_8H_{10}	417.5(0.4)	630.26(0.1)	3.74(0.01)	372(40)	34, 68, 78, 79, 88, 669
<i>p</i> -Xylene	C_8H_{10}	411.4(0.5)	616.17(0.09)	3.55(0.02)	372(35)	34, 68, 74, 79, 88, 90, 669
2,3-Xylenol	$C_8H_{10}O$	490.03(0.05)	723(1)	4.1(0.3)	397(15)	218
2,4-Xylenol	$C_8H_{10}O$	484.09(0.03)	708(1)	3.5(0.3)	389(15)	218
2,5-Xylenol	$C_8H_{10}O$	484.29(0.08)	707(1)	3.9(0.1)	397(15)	25
2,6-Xylenol	$C_8H_{10}O$	474.18(0.05)	701(1)	3.8(0.1)	396(15)	218
3,4-Xylenol	$C_8H_{10}O$	500.46(0.05)	730(1)	4.9(0.5)	388(15)	218
3,5-Xylenol	$C_8H_{10}O$	494.86(0.05)	716(1)	3.8(0.2)	396(15)	218

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