

2-Methylcyclopropanemethanol

Other names:	Cyclopropanemethanol, 2-methyl-
Inchi:	InChI=1S/C5H10O/c1-4-2-5(4)3-6/h4-6H,2-3H2,1H3
InchiKey:	SHEINYPABNPRPM-UHFFFAOYSA-N
Formula:	C5H10O
SMILES:	CC1CC1CO
Mol. weight [g/mol]:	86.13
CAS:	6077-72-1

Physical Properties

Property code	Value	Unit	Source
gf	-92.56	kJ/mol	Joback Method
hf	-246.30	kJ/mol	Joback Method
hfus	12.00	kJ/mol	Joback Method
hvap	43.01	kJ/mol	Joback Method
log10ws	-0.59		Crippen Method
logp	0.635		Crippen Method
mcvol	76.320	ml/mol	McGowan Method
pc	4391.59	kPa	Joback Method
tb	406.20	K	NIST Webbook
tc	582.49	K	Joback Method
tf	220.63	K	Joback Method
vc	0.290	m ³ /kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	154.30	J/mol×K	408.05	Joback Method
cpg	198.08	J/mol×K	553.42	Joback Method
cpg	190.17	J/mol×K	524.34	Joback Method
cpg	181.86	J/mol×K	495.27	Joback Method
cpg	173.12	J/mol×K	466.20	Joback Method
cpg	163.94	J/mol×K	437.12	Joback Method
cpg	205.60	J/mol×K	582.49	Joback Method
dvisc	0.0004267	Paxs	408.05	Joback Method

dvisc	0.0006072	Paxs	376.81	Joback Method
dvisc	0.0009212	Paxs	345.58	Joback Method
dvisc	0.0015181	Paxs	314.34	Joback Method
dvisc	0.0027935	Paxs	283.10	Joback Method
dvisc	0.0059796	Paxs	251.87	Joback Method
dvisc	0.0158779	Paxs	220.63	Joback Method

Sources

Crippen Method:	http://pubs.acs.org/doi/abs/10.1021/ci9903071
Crippen Method:	https://www.chemeo.com/doc/models/crippen_log10ws
Joback Method:	https://en.wikipedia.org/wiki/Joback_method
McGowan Method:	http://link.springer.com/article/10.1007/BF02311772
NIST Webbook:	http://webbook.nist.gov/cgi/cbook.cgi?ID=C6077721&Units=SI

Legend

cpg:	Ideal gas heat capacity
dvisc:	Dynamic viscosity
gf:	Standard Gibbs free energy of formation
hf:	Enthalpy of formation at standard conditions
hfus:	Enthalpy of fusion at standard conditions
hvap:	Enthalpy of vaporization at standard conditions
log10ws:	Log10 of Water solubility in mol/l
logp:	Octanol/Water partition coefficient
mcvol:	McGowan's characteristic volume
pc:	Critical Pressure
tb:	Normal Boiling Point Temperature
tc:	Critical Temperature
tf:	Normal melting (fusion) point
vc:	Critical Volume

Latest version available from:

<https://www.chemeo.com/cid/12-842-3/2-Methylcyclopropanemethanol.pdf>

Generated by Cheméo on 2024-04-19 19:08:21.227830974 +0000 UTC m=+15842950.148408289.

Cheméo (<https://www.chemeo.com>) is the biggest free database of chemical and physical data for the process industry.