

Cyclopentane, 1-methyl-2-methylene-

Other names:	1-Methylene-2-methylcyclopentane 2-Methyl methylenecyclopentane
Inchi:	InChI=1S/C7H12/c1-6-4-3-5-7(6)2/h7H,1,3-5H2,2H3
InchiKey:	FADJUGSUYL CNJQ-UHFFFAOYSA-N
Formula:	C7H12
SMILES:	C=C1CCCC1C
Mol. weight [g/mol]:	96.17
CAS:	41158-41-2

Physical Properties

Property code	Value	Unit	Source
gf	97.69	kJ/mol	Joback Method
hf	-43.09	kJ/mol	Joback Method
hfus	6.66	kJ/mol	Joback Method
hvap	31.59	kJ/mol	Joback Method
log10ws	-2.26		Crippen Method
logp	2.363		Crippen Method
mcvol	94.330	ml/mol	McGowan Method
pc	3497.14	kPa	Joback Method
rinpol	732.00		NIST Webbook
tb	374.00	K	Joback Method
tc	569.84	K	Joback Method
tf	193.23	K	Joback Method
vc	0.352	m3/kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	163.26	J/mol×K	374.00	Joback Method
cpg	176.80	J/mol×K	406.64	Joback Method
cpg	189.72	J/mol×K	439.28	Joback Method
cpg	202.03	J/mol×K	471.92	Joback Method
cpg	213.76	J/mol×K	504.56	Joback Method
cpg	224.91	J/mol×K	537.20	Joback Method

cpg	235.52	J/mol×K	569.84	Joback Method
dvisc	0.0017344	Pa×s	193.23	Joback Method
dvisc	0.0010433	Pa×s	223.36	Joback Method
dvisc	0.0007082	Pa×s	253.49	Joback Method
dvisc	0.0005220	Pa×s	283.62	Joback Method
dvisc	0.0004080	Pa×s	313.74	Joback Method
dvisc	0.0003329	Pa×s	343.87	Joback Method
dvisc	0.0002807	Pa×s	374.00	Joback Method

Sources

Crippen Method:	http://pubs.acs.org/doi/abs/10.1021/ci990307l
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=C41158412&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
rinpol:	Non-polar retention indices
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/41-519-0/Cyclopentane-1-methyl-2-methylene.pdf>

Generated by Cheméo on 2025-12-05 07:22:03.468167899 +0000 UTC m=+4667520.998208553.

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