Cyclopentane, 1,2-dimethyl-

Other names: 1,2-Dimethylcyclopentane
Inchi: InChI=1S/C7H14/c1-6-4-3-5-7(6)2/h6-7H,3-5H2,1-2H3
InchiKey: RIRARCHMRDHZAR-UHFFFAOYSA-N
Formula: C7H14
SMILES: CC1CCCC1C
Mol. weight [g/mol]: 98.19
CAS: 2452-99-5

Physical Properties

Property code	Value	Unit	Source
gf	36.90	kJ/mol	Joback Method
hf	-147.67	kJ/mol	Joback Method
hfus	8.89	kJ/mol	Joback Method
hvap	31.12	kJ/mol	Joback Method
log10ws	-2.16		Crippen Method
logp	2.442		Crippen Method
mcvol	98.630	ml/mol	McGowan Method
pc	3291.59	kPa	Joback Method
rinpol	682.00		NIST Webbook
rinpol	689.00		NIST Webbook
rinpol	689.00		NIST Webbook
rinpol	682.00		NIST Webbook
rinpol	685.00		NIST Webbook
rinpol	677.00		NIST Webbook
rinpol	686.00		NIST Webbook
rinpol	686.00		NIST Webbook
ripol	726.00		NIST Webbook
ripol	726.00		NIST Webbook
sl	269.90	J/mol×K	NIST Webbook
tb	370.00 ± 7.00	K	NIST Webbook
tc	563.52	K	Joback Method
tf	175.31	K	Joback Method
vc	0.367	m3/kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	256.86	J/mol×K	563.52	Joback Method
cpg	244.64	J/mol×K	531.30	Joback Method
cpg	173.95	J/mol×K	370.17	Joback Method
cpg	189.42	J/mol×K	402.40	Joback Method
cpg	204.20	J/mol×K	434.62	Joback Method
cpg	218.32	J/mol×K	466.85	Joback Method
cpg	231.80	J/mol×K	499.07	Joback Method
cpl	187.40	J/mol×K	294.20	NIST Webbook
dvisc	0.0002650	Pa×s	370.17	Joback Method
dvisc	0.0003124	Pa×s	337.69	Joback Method
dvisc	0.0017720	Pa×s	175.31	Joback Method
dvisc	0.0010080	Pa×s	207.79	Joback Method
dvisc	0.0006678	Pa×s	240.26	Joback Method
dvisc	0.0004881	Pa×s	272.74	Joback Method
dvisc	0.0003813	Pa×s	305.22	Joback Method
hfust	6.41	kJ/mol	154.10	NIST Webbook
sfust	41.60	J/mol×K	154.10	NIST Webbook

Correlations

Information	Value
Property code	pvap
Equation	ln(Pvp) = A + B/(T + C)
Coeff. A	1.34204e+01
Coeff. B	-2.86939e+03
Coeff. C	-4.40100e+01
Temperature range (K), min.	262.50
Temperature range (K), max.	397.87

Sources

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=C2452995&Units=SI
The Yaws Handbook of Vapor Pressure:	https://www.sciencedirect.com/book/9780128029992/the-yaws-handbook-of-vapor-pressure
Crippen Method:	http://pubs.acs.org/doi/abs/10.1021/ci990307l

Legend

cpg:	Ideal gas heat capacity
cpl:	Liquid phase 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
hfust:	Enthalpy of fusion at a given temperature
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
pvap:	Vapor pressure
rinpol:	Non-polar retention indices
ripol:	Polar retention indices
sfust:	Entropy of fusion at a given temperature
sl:	Liquid phase molar entropy at standard conditions
tb:	Normal Boiling Point Temperature
tc:	Critical Temperature
tf:	Normal melting (fusion) point
vc:	Critical Volume