

# Cyclopentane, 1,2-dimethyl-3-methylene

Inchi:	InChI=1S/C8H14/c1-6-4-5-7(2)8(6)3/h7-8H,1,4-5H2,2-3H3
InchiKey:	SITNPFCTAOJHRR-UHFFFAOYSA-N
Formula:	C8H14
SMILES:	C=C1CCC(C)C1C
Mol. weight [g/mol]:	110.20

## Physical Properties

Property code	Value	Unit	Source
gf	98.40	kJ/mol	Joback Method
hf	-84.07	kJ/mol	Joback Method
hfus	10.32	kJ/mol	Joback Method
hvap	33.51	kJ/mol	Joback Method
log10ws	-2.44		Crippen Method
logp	2.609		Crippen Method
mcvol	108.420	ml/mol	McGowan Method
pc	3018.96	kPa	Joback Method
rinpol	804.00		NIST Webbook
tb	392.21	K	Joback Method
tc	586.16	K	Joback Method
tf	200.26	K	Joback Method
vc	0.407	m3/kmol	Joback Method

## Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	201.97	J/molxK	392.21	Joback Method
cpg	217.01	J/molxK	424.53	Joback Method
cpg	231.41	J/molxK	456.86	Joback Method
cpg	245.18	J/molxK	489.18	Joback Method
cpg	258.34	J/molxK	521.51	Joback Method
cpg	270.90	J/molxK	553.83	Joback Method
cpg	282.87	J/molxK	586.16	Joback Method
dvisc	0.0011251	Paxs	200.26	Joback Method
dvisc	0.0007578	Paxs	232.25	Joback Method

dvisc	0.0005617	Paxs	264.24	Joback Method
dvisc	0.0004442	Paxs	296.24	Joback Method
dvisc	0.0003677	Paxs	328.23	Joback Method
dvisc	0.0003147	Paxs	360.22	Joback Method
dvisc	0.0002763	Paxs	392.21	Joback Method

## Sources

<b>McGowan Method:</b>	<a href="http://link.springer.com/article/10.1007/BF02311772">http://link.springer.com/article/10.1007/BF02311772</a>
<b>NIST Webbook:</b>	<a href="http://webbook.nist.gov/cgi/cbook.cgi?ID=R77301&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=R77301&amp;Units=SI</a>
<b>Crippen Method:</b>	<a href="http://pubs.acs.org/doi/abs/10.1021/ci9903071">http://pubs.acs.org/doi/abs/10.1021/ci9903071</a>
<b>Crippen Method:</b>	<a href="https://www.chemeo.com/doc/models/crippen_log10ws">https://www.chemeo.com/doc/models/crippen_log10ws</a>
<b>Joback Method:</b>	<a href="https://en.wikipedia.org/wiki/Joback_method">https://en.wikipedia.org/wiki/Joback_method</a>

## Legend

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

Latest version available from:

<https://www.chemeo.com/cid/64-038-9/Cyclopentane-1-2-dimethyl-3-methylene.pdf>

Generated by Cheméo on 2025-12-05 08:24:57.070796662 +0000 UTC m=+4671294.600837326.

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