

# Cyclopentene,3-methoxy-

<b>Inchi:</b>	InChI=1S/C6H10O/c1-7-6-4-2-3-5-6/h2,4,6H,3,5H2,1H3
<b>InchiKey:</b>	PUNAFIBTAHETMN-UHFFFAOYSA-N
<b>Formula:</b>	C6H10O
<b>SMILES:</b>	COC1C=CCC1
<b>Mol. weight [g/mol]:</b>	98.14
<b>CAS:</b>	39819-74-4

## Physical Properties

Property code	Value	Unit	Source
gf	-38.85	kJ/mol	Joback Method
hf	-181.13	kJ/mol	Joback Method
hfus	7.64	kJ/mol	Joback Method
hvap	31.91	kJ/mol	Joback Method
ie	9.45 ± 0.05	eV	NIST Webbook
log10ws	-1.28		Crippen Method
logp	1.351		Crippen Method
mcvol	86.110	ml/mol	McGowan Method
pc	3930.78	kPa	Joback Method
tb	373.54	K	Joback Method
tc	570.82	K	Joback Method
tf	191.27	K	Joback Method
vc	0.317	m <sup>3</sup> /kmol	Joback Method

## Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	149.07	J/mol×K	373.54	Joback Method
cpg	161.23	J/mol×K	406.42	Joback Method
cpg	172.86	J/mol×K	439.30	Joback Method
cpg	183.96	J/mol×K	472.18	Joback Method
cpg	194.55	J/mol×K	505.06	Joback Method
cpg	204.65	J/mol×K	537.94	Joback Method
cpg	214.25	J/mol×K	570.82	Joback Method
dvisc	0.0019874	Paxs	191.27	Joback Method

dvisc	0.0011102	Paxs	221.65	Joback Method
dvisc	0.0007136	Paxs	252.03	Joback Method
dvisc	0.0005045	Paxs	282.40	Joback Method
dvisc	0.0003815	Paxs	312.78	Joback Method
dvisc	0.0003031	Paxs	343.16	Joback Method
dvisc	0.0002500	Paxs	373.54	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=C39819744&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=C39819744&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>ie:</b>	Ionization energy
<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>tb:</b>	Normal Boiling Point Temperature
<b>tc:</b>	Critical Temperature
<b>tf:</b>	Normal melting (fusion) point
<b>vc:</b>	Critical Volume

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