

# 1-methyl-cis-2-(1-ethenyl)-cyclopropane

<b>Inchi:</b>	InChI=1S/C6H10/c1-3-6-4-5(6)2/h3,5-6H,1,4H2,2H3/t5-,6+/m1/s1
<b>InchiKey:</b>	JVVPJIPOOZHNTM-RITPCOANSA-N
<b>Formula:</b>	C6H10
<b>SMILES:</b>	C=CC1CC1C
<b>Mol. weight [g/mol]:</b>	82.14

## Physical Properties

Property code	Value	Unit	Source
gf	140.52	kJ/mol	Joback Method
hf	10.72	kJ/mol	Joback Method
hfus	9.22	kJ/mol	Joback Method
hvap	27.88	kJ/mol	Joback Method
log10ws	-1.60		Crippen Method
logp	1.828		Crippen Method
mcvol	80.240	ml/mol	McGowan Method
pc	3650.93	kPa	Joback Method
rinpol	605.40		NIST Webbook
rinpol	609.80		NIST Webbook
tb	335.43	K	Joback Method
tc	516.94	K	Joback Method
tf	169.32	K	Joback Method
vc	0.308	m3/kmol	Joback Method

## Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	128.08	J/molxK	335.43	Joback Method
cpg	139.86	J/molxK	365.68	Joback Method
cpg	151.04	J/molxK	395.93	Joback Method
cpg	161.64	J/molxK	426.19	Joback Method
cpg	171.70	J/molxK	456.44	Joback Method
cpg	181.23	J/molxK	486.69	Joback Method
cpg	190.26	J/molxK	516.94	Joback Method
dvisc	0.0002452	Paxs	169.32	Joback Method

dvisc	0.0002358	Paxs	197.00	Joback Method
dvisc	0.0002290	Paxs	224.69	Joback Method
dvisc	0.0002239	Paxs	252.38	Joback Method
dvisc	0.0002198	Paxs	280.06	Joback Method
dvisc	0.0002165	Paxs	307.75	Joback Method
dvisc	0.0002138	Paxs	335.43	Joback Method

## Sources

<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>
<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=R137267&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=R137267&amp;Units=SI</a>
<b>Crippen Method:</b>	<a href="http://pubs.acs.org/doi/abs/10.1021/ci990307l">http://pubs.acs.org/doi/abs/10.1021/ci990307l</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

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