

# 1-Methyl-cis-2-(trans-3-pentenyl)-cyclopropane

<b>Inchi:</b>	InChI=1S/C9H16/c1-3-4-5-6-9-7-8(9)2/h3-4,8-9H,5-7H2,1-2H3/b4-3+/t8-,9+/m1/s1
<b>InchiKey:</b>	BRJAZMJACNNPEV-BKIAHZASSA-N
<b>Formula:</b>	C9H16
<b>SMILES:</b>	CC=CCCC1CC1C
<b>Mol. weight [g/mol]:</b>	124.22

## Physical Properties

Property code	Value	Unit	Source
gf	158.16	kJ/mol	Joback Method
hf	-59.41	kJ/mol	Joback Method
hfus	18.47	kJ/mol	Joback Method
hvap	35.19	kJ/mol	Joback Method
log10ws	-2.85		Crippen Method
logp	2.999		Crippen Method
mcvol	122.510	ml/mol	McGowan Method
pc	2676.31	kPa	Joback Method
rinpol	900.00		NIST Webbook
rinpol	897.50		NIST Webbook
rinpol	900.00		NIST Webbook
tb	411.55	K	Joback Method
tc	596.57	K	Joback Method
tf	199.81	K	Joback Method
vc	0.475	m3/kmol	Joback Method

## Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	243.05	J/molxK	411.55	Joback Method
cpg	313.96	J/molxK	565.74	Joback Method
cpg	301.26	J/molxK	534.90	Joback Method
cpg	287.86	J/molxK	504.06	Joback Method
cpg	273.72	J/molxK	473.22	Joback Method
cpg	258.79	J/molxK	442.39	Joback Method
cpg	325.99	J/molxK	596.57	Joback Method

dvisc	0.0003049	Paxs	411.55	Joback Method
dvisc	0.0003318	Paxs	376.26	Joback Method
dvisc	0.0003674	Paxs	340.97	Joback Method
dvisc	0.0004165	Paxs	305.68	Joback Method
dvisc	0.0004879	Paxs	270.39	Joback Method
dvisc	0.0005994	Paxs	235.10	Joback Method
dvisc	0.0007918	Paxs	199.81	Joback Method

## Sources

<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=R137396&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=R137396&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>
<b>Crippen Method:</b>	<a href="https://www.cheméo.com/doc/models/crippen_log10ws">https://www.cheméo.com/doc/models/crippen_log10ws</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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