

# Cyclopentane, butylidene

Inchi:	InChI=1S/C9H16/c1-2-3-6-9-7-4-5-8-9/h6H,2-5,7-8H2,1H3
InchiKey:	SFOPXUUMZRCYNJ-UHFFFAOYSA-N
Formula:	C9H16
SMILES:	CCCC=C1CCCC1
Mol. weight [g/mol]:	124.22

## Physical Properties

Property code	Value	Unit	Source
gf	114.62	kJ/mol	Joback Method
hf	-72.24	kJ/mol	Joback Method
hfus	12.25	kJ/mol	Joback Method
hvap	36.98	kJ/mol	Joback Method
log10ws	-3.34		Crippen Method
logp	3.287		Crippen Method
mvol	122.510	ml/mol	McGowan Method
pc	2953.69	kPa	Joback Method
rinpol	946.00		NIST Webbook
rinpol	949.00		NIST Webbook
tb	431.91	K	Joback Method
tc	630.76	K	Joback Method
tf	216.69	K	Joback Method
vc	0.465	m <sup>3</sup> /kmol	Joback Method

## Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	243.31	J/molxK	431.91	Joback Method
cpg	259.39	J/molxK	465.05	Joback Method
cpg	274.62	J/molxK	498.19	Joback Method
cpg	289.04	J/molxK	531.33	Joback Method
cpg	302.68	J/molxK	564.48	Joback Method
cpg	315.57	J/molxK	597.62	Joback Method
cpg	327.76	J/molxK	630.76	Joback Method
dvisc	0.0042711	Paxs	216.69	Joback Method

dvisc	0.0019383	Paxs	252.56	Joback Method
dvisc	0.0010706	Paxs	288.43	Joback Method
dvisc	0.0006743	Paxs	324.30	Joback Method
dvisc	0.0004657	Paxs	360.17	Joback Method
dvisc	0.0003439	Paxs	396.04	Joback Method
dvisc	0.0002671	Paxs	431.91	Joback Method

## Sources

<b>NIST Webbook:</b>	<a href="http://webbook.nist.gov/cgi/cbook.cgi?ID=R10753&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=R10753&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>
<b>McGowan Method:</b>	<a href="http://link.springer.com/article/10.1007/BF02311772">http://link.springer.com/article/10.1007/BF02311772</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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