

# (trans-1,2-Methylene)-trans-4-hexenyl-cyclopropa

<b>Inchi:</b>	InChI=1S/C10H16/c1-2-3-4-9-7-10(9)8-5-6-8/h2-3,8-10H,4-7H2,1H3/b3-2+/t9-,10-/m1/s1
<b>InchiKey:</b>	ZFAGMTQIRSTENT-HLKOZVMDSA-N
<b>Formula:</b>	C10H16
<b>SMILES:</b>	CC=CCC1CC1C1CC1
<b>Mol. weight [g/mol]:</b>	136.23

## Physical Properties

Property code	Value	Unit	Source
gf	227.33	kJ/mol	Joback Method
hf	-7.25	kJ/mol	Joback Method
hfus	19.20	kJ/mol	Joback Method
hvap	37.33	kJ/mol	Joback Method
log10ws	-2.93		Crippen Method
logp	2.999		Crippen Method
mcvol	125.740	ml/mol	McGowan Method
pc	2773.00	kPa	Joback Method
rinpol	989.50		NIST Webbook
rinpol	987.50		NIST Webbook
rinpol	989.50		NIST Webbook
tb	441.17	K	Joback Method
tc	640.04	K	Joback Method
tf	229.02	K	Joback Method
vc	0.488	m <sup>3</sup> /kmol	Joback Method

## Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	271.61	J/molxK	441.17	Joback Method
cpg	289.79	J/molxK	474.31	Joback Method
cpg	306.82	J/molxK	507.46	Joback Method
cpg	322.76	J/molxK	540.60	Joback Method
cpg	337.68	J/molxK	573.75	Joback Method
cpg	351.67	J/molxK	606.89	Joback Method
cpg	364.78	J/molxK	640.04	Joback Method

dvisc	0.0005242	Paxs	229.02	Joback Method
dvisc	0.0005401	Paxs	264.38	Joback Method
dvisc	0.0005525	Paxs	299.74	Joback Method
dvisc	0.0005625	Paxs	335.09	Joback Method
dvisc	0.0005708	Paxs	370.45	Joback Method
dvisc	0.0005777	Paxs	405.81	Joback Method
dvisc	0.0005835	Paxs	441.17	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=R138022&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=R138022&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>mvol:</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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