

# Cyclohexane, 1-ethenyl-1-methyl-2-(1-methylethenyl)-4-(1-methyl-2-propenyl)-

<b>Other names:</b>	o-Menth-8-ene, 4-isopropylidene-1-vinyl- Elixene 1-Methyl-4-(1-methylethylidene)-2-(1-methylvinyl)-1-vinylcyclohexane
<b>Inchi:</b>	InChI=1S/C15H24/c1-7-15(6)9-8-13(11(2)3)10-14(15)12(4)5/h7,14H,1,4,8-10H2,2-3,5-6H
<b>InchiKey:</b>	BQSLMQNYHVFRDT-UHFFFAOYSA-N
<b>Formula:</b>	C15H24
<b>SMILES:</b>	C=CC1(C)CCC(=C(C)C)CC1C(=C)C
<b>Mol. weight [g/mol]:</b>	204.35
<b>CAS:</b>	3242-08-8

## Physical Properties

Property code	Value	Unit	Source
gf	290.71	kJ/mol	Joback Method
hf	3.60	kJ/mol	Joback Method
hfus	16.36	kJ/mol	Joback Method
hvap	47.56	kJ/mol	Joback Method
log10ws	-5.07		Crippen Method
logp	4.891		Crippen Method
mcvol	198.450	ml/mol	McGowan Method
pc	1861.11	kPa	Joback Method
rinpol	1488.00		NIST Webbook
rinpol	1492.00		NIST Webbook
rinpol	1449.00		NIST Webbook
rinpol	1449.00		NIST Webbook
rinpol	1445.00		NIST Webbook
rinpol	1511.00		NIST Webbook
rinpol	1492.00		NIST Webbook
ripol	1514.00		NIST Webbook
tb	557.48	K	Joback Method
tc	771.21	K	Joback Method
tf	264.77	K	Joback Method
vc	0.752	m <sup>3</sup> /kmol	Joback Method

# Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	490.96	J/mol×K	557.48	Joback Method
cpg	512.54	J/mol×K	593.10	Joback Method
cpg	532.84	J/mol×K	628.72	Joback Method
cpg	551.99	J/mol×K	664.35	Joback Method
cpg	570.13	J/mol×K	699.97	Joback Method
cpg	587.38	J/mol×K	735.59	Joback Method
cpg	603.88	J/mol×K	771.21	Joback Method

## Sources

<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>
<b>NIST Webbook:</b>	<a href="http://webbook.nist.gov/cgi/cbook.cgi?ID=C3242088&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=C3242088&amp;Units=SI</a>

## Legend

<b>cpg:</b>	Ideal gas heat capacity
<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>ripol:</b>	Non-polar retention indices
<b>ripol:</b>	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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