

# Cedrenol acetate

## Other names:

<b>Inchi:</b>	[3R-(3«alpha»,3a«beta»,7«beta»,8a«alpha»)]-2,3,4,7,8,8a-hexahydro-3,8,8-trimethyl-1H-inden-1-yl acetate
<b>InchiKey:</b>	UTZBRSREMTWJBB-OOOXPIDESA-N
<b>Formula:</b>	C17H26O2
<b>SMILES:</b>	C=C1C(OC(C)=O)CC23CC1C(C)(C)C2CCC3C
<b>Mol. weight [g/mol]:</b>	262.39
<b>CAS:</b>	1405-92-1

## Physical Properties

Property code	Value	Unit	Source
gf	35.36	kJ/mol	Joback Method
hf	-379.23	kJ/mol	Joback Method
hfus	22.24	kJ/mol	Joback Method
hvap	59.60	kJ/mol	Joback Method
log10ws	-4.24		Crippen Method
logp	3.957		Crippen Method
mcvol	220.950	ml/mol	McGowan Method
pc	1804.63	kPa	Joback Method
rinpol	1743.00		NIST Webbook
rinpol	1743.00		NIST Webbook
tb	679.04	K	Joback Method
tc	899.33	K	Joback Method
tf	449.05	K	Joback Method
vc	0.844	m3/kmol	Joback Method

## Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	669.85	J/molxK	679.04	Joback Method
cpg	691.60	J/molxK	715.75	Joback Method
cpg	712.47	J/molxK	752.47	Joback Method
cpg	732.71	J/molxK	789.18	Joback Method
cpg	752.57	J/molxK	825.90	Joback Method
cpg	772.32	J/molxK	862.61	Joback Method

## Sources

<b>NIST Webbook:</b>	<a href="http://webbook.nist.gov/cgi/cbook.cgi?ID=C1405921&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=C1405921&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>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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