

# dehydrocarveol

<b>Inchi:</b>	InChI=1S/C10H14O/c1-7(2)9-5-4-8(3)10(11)6-9/h4-5,10-11H,1,6H2,2-3H3
<b>InchiKey:</b>	SUSDZWDMNJCYTN-UHFFFAOYSA-N
<b>Formula:</b>	C10H14O
<b>SMILES:</b>	C=C(C)C1=CC=C(C)C(O)C1
<b>Mol. weight [g/mol]:</b>	150.22

## Physical Properties

Property code	Value	Unit	Source
gf	40.90	kJ/mol	Joback Method
hf	-139.38	kJ/mol	Joback Method
hfus	16.66	kJ/mol	Joback Method
hvap	56.28	kJ/mol	Joback Method
log10ws	-2.84		Crippen Method
logp	2.200		Crippen Method
mcvol	133.870	ml/mol	McGowan Method
pc	3114.04	kPa	Joback Method
ripol	1593.00		NIST Webbook
ripol	1591.00		NIST Webbook
ripol	1941.00		NIST Webbook
ripol	1949.00		NIST Webbook
ripol	1945.00		NIST Webbook
ripol	1943.00		NIST Webbook
tb	544.77	K	Joback Method
tc	743.50	K	Joback Method
tf	281.50	K	Joback Method
vc	0.501	m <sup>3</sup> /kmol	Joback Method

## Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	312.23	J/molxK	544.77	Joback Method
cpg	325.47	J/molxK	577.89	Joback Method
cpg	338.02	J/molxK	611.01	Joback Method
cpg	349.91	J/molxK	644.14	Joback Method

cpg	361.14	J/mol×K	677.26	Joback Method
cpg	371.76	J/mol×K	710.38	Joback Method
cpg	381.77	J/mol×K	743.50	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=R197190&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=R197190&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>rinpol:</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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