

# 7«alpha»-hydroxymanool

<b>Inchi:</b>	InChI=1S/C20H34O2/c1-7-19(5,22)11-10-15-12-17(21)18(3,4)16-9-8-14(2)13-20(15,16)6
<b>InchiKey:</b>	QBGCJSNIPHJJC-V-AZDWTMOSSA-N
<b>Formula:</b>	C20H34O2
<b>SMILES:</b>	<chem>C=CC(C)(O)CCC1CC(O)C(C)(C)C2CCC(=C)CC12C</chem>
<b>Mol. weight [g/mol]:</b>	306.48

## Physical Properties

Property code	Value	Unit	Source
gf	26.63	kJ/mol	Joback Method
hf	-469.25	kJ/mol	Joback Method
hfus	24.37	kJ/mol	Joback Method
hvap	88.95	kJ/mol	Joback Method
log10ws	-5.48		Crippen Method
logp	4.473		Crippen Method
mvol	274.080	ml/mol	McGowan Method
pc	1570.96	kPa	Joback Method
rinpol	2235.00		NIST Webbook
rinpol	2235.00		NIST Webbook
tb	851.00	K	Joback Method
tc	1054.41	K	Joback Method
tf	508.02	K	Joback Method
vc	1.022	m <sup>3</sup> /kmol	Joback Method

## Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	936.83	J/mol×K	851.00	Joback Method
cpg	957.92	J/mol×K	884.90	Joback Method
cpg	979.01	J/mol×K	918.80	Joback Method
cpg	1000.30	J/mol×K	952.71	Joback Method
cpg	1022.02	J/mol×K	986.61	Joback Method
cpg	1044.37	J/mol×K	1020.51	Joback Method
cpg	1067.56	J/mol×K	1054.41	Joback Method

# Sources

<b>NIST Webbook:</b>	<a href="http://webbook.nist.gov/cgi/cbook.cgi?ID=R333485&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=R333485&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>hvp:</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>rinp:</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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