

# Hydroxy-3-endo-bornanone-2

<b>Inchi:</b>	InChI=1S/C10H16O2/c1-9(2)6-4-5-10(9,3)8(12)7(6)11/h6-7,11H,4-5H2,1-3H3
<b>InchiKey:</b>	AXMKZEOEXSKFJI-UHFFFAOYSA-N
<b>Formula:</b>	C10H16O2
<b>SMILES:</b>	CC12CCC(C(O)C1=O)C2(C)C
<b>Mol. weight [g/mol]:</b>	168.23
<b>CAS:</b>	21488-68-6

## Physical Properties

Property code	Value	Unit	Source
gf	-143.09	kJ/mol	Joback Method
hf	-410.42	kJ/mol	Joback Method
hfus	8.97	kJ/mol	Joback Method
hvap	55.86	kJ/mol	Joback Method
log10ws	-1.73		Crippen Method
logp	1.373		Crippen Method
mcvol	137.480	ml/mol	McGowan Method
pc	3295.37	kPa	Joback Method
tb	597.09	K	Joback Method
tc	811.66	K	Joback Method
tf	403.18	K	Joback Method
vc	0.521	m3/kmol	Joback Method

## Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	384.29	J/molxK	597.09	Joback Method
cpg	399.17	J/molxK	632.85	Joback Method
cpg	413.28	J/molxK	668.61	Joback Method
cpg	426.81	J/molxK	704.38	Joback Method
cpg	439.96	J/molxK	740.14	Joback Method
cpg	452.93	J/molxK	775.90	Joback Method
cpg	465.91	J/molxK	811.66	Joback Method

# Sources

<b>Crippen Method:</b>	<a href="http://pubs.acs.org/doi/abs/10.1021/ci990307I">http://pubs.acs.org/doi/abs/10.1021/ci990307I</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=C21488686&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=C21488686&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>mccvol:</b>	McGowan's characteristic volume
<b>pc:</b>	Critical Pressure
<b>tb:</b>	Normal Boiling Point Temperature
<b>tc:</b>	Critical Temperature
<b>tf:</b>	Normal melting (fusion) point
<b>vc:</b>	Critical Volume

Latest version available from:

<https://www.chemeo.com/cid/13-436-3/Hydroxy-3-endo-bornanone-2.pdf>

Generated by Cheméo on 2024-05-02 04:04:52.989481592 +0000 UTC m=+16911941.910058903.

Cheméo (<https://www.chemeo.com>) is the biggest free database of chemical and physical data for the process industry.