

# 1,3-Cyclohexanedimethanol, 2-hydroxy-5-methyl-

Inchi:	InChI=1S/C9H18O3/c1-6-2-7(4-10)9(12)8(3-6)5-11/h6-12H,2-5H2,1H3
InchiKey:	VGNOCVYYJJSUEN-UHFFFAOYSA-N
Formula:	C9H18O3
SMILES:	CC1CC(CO)C(O)C(CO)C1
Mol. weight [g/mol]:	174.24
CAS:	90677-23-9

## Physical Properties

Property code	Value	Unit	Source
gf	-384.24	kJ/mol	Joback Method
hf	-692.48	kJ/mol	Joback Method
hfus	26.38	kJ/mol	Joback Method
hvap	85.17	kJ/mol	Joback Method
log10ws	-0.66		Crippen Method
logp	-0.006		Crippen Method
mvol	144.420	ml/mol	McGowan Method
pc	3460.21	kPa	Joback Method
tb	687.40	K	Joback Method
tc	862.11	K	Joback Method
tf	368.31	K	Joback Method
vc	0.526	m3/kmol	Joback Method

## Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	445.95	J/molxK	687.40	Joback Method
cpg	458.04	J/molxK	716.52	Joback Method
cpg	469.52	J/molxK	745.64	Joback Method
cpg	480.39	J/molxK	774.76	Joback Method
cpg	490.66	J/molxK	803.87	Joback Method
cpg	500.34	J/molxK	832.99	Joback Method
cpg	509.42	J/molxK	862.11	Joback Method
dvisc	0.0116633	Paxs	368.31	Joback Method
dvisc	0.0016561	Paxs	421.49	Joback Method

dvisc	0.0003642	Paxs	474.67	Joback Method
dvisc	0.0001087	Paxs	527.86	Joback Method
dvisc	0.0000405	Paxs	581.04	Joback Method
dvisc	0.0000178	Paxs	634.22	Joback Method
dvisc	0.0000089	Paxs	687.40	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=C90677239&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=C90677239&amp;Units=SI</a>

## Legend

<b>cpg:</b>	Ideal gas heat capacity
<b>dvisc:</b>	Dynamic viscosity
<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>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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