

# Cyclopropanemethanol, «alpha»-methyl-«alpha»-propyl-

<b>Other names:</b>	2-Pentanol, 2-cyclopropyl- 2-Cyclopropyl-2-pentanol
<b>Inchi:</b>	InChI=1S/C8H16O/c1-3-6-8(2,9)7-4-5-7/h7,9H,3-6H2,1-2H3
<b>InchiKey:</b>	WEXFBWATVRHNFT-UHFFFAOYSA-N
<b>Formula:</b>	C8H16O
<b>SMILES:</b>	CCCC(C)(O)C1CC1
<b>Mol. weight [g/mol]:</b>	128.21
<b>CAS:</b>	24230-08-8

## Physical Properties

Property code	Value	Unit	Source
gf	-56.75	kJ/mol	Joback Method
hf	-296.63	kJ/mol	Joback Method
hfus	11.28	kJ/mol	Joback Method
hvap	48.70	kJ/mol	Joback Method
log10ws	-2.20		Crippen Method
logp	1.948		Crippen Method
mvol	118.590	ml/mol	McGowan Method
pc	3284.05	kPa	Joback Method
tb	436.70 ± 0.50	K	NIST Webbook
tc	659.77	K	Joback Method
tf	261.10	K	Joback Method
vc	0.449	m <sup>3</sup> /kmol	Joback Method

## Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	278.31	J/mol×K	478.13	Joback Method
cpg	338.31	J/mol×K	629.50	Joback Method
cpg	327.71	J/mol×K	599.22	Joback Method
cpg	316.45	J/mol×K	568.95	Joback Method
cpg	304.50	J/mol×K	538.68	Joback Method
cpg	291.80	J/mol×K	508.40	Joback Method
cpg	348.30	J/mol×K	659.77	Joback Method

dvisc	0.0003243	Paxs	478.13	Joback Method
dvisc	0.0004982	Paxs	441.96	Joback Method
dvisc	0.0008262	Paxs	405.79	Joback Method
dvisc	0.0015128	Paxs	369.62	Joback Method
dvisc	0.0031582	Paxs	333.44	Joback Method
dvisc	0.0078871	Paxs	297.27	Joback Method
dvisc	0.0253814	Paxs	261.10	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=C24230088&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=C24230088&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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