

# 4,5-Octanediol

<b>Other names:</b>	4,5-dihydroxyoctane
<b>Inchi:</b>	InChI=1S/C8H18O2/c1-3-5-7(9)8(10)6-4-2/h7-10H,3-6H2,1-2H3
<b>InchiKey:</b>	YOZZZCLQJVMZGY-UHFFFAOYSA-N
<b>Formula:</b>	C8H18O2
<b>SMILES:</b>	CCCC(O)C(O)CCC
<b>Mol. weight [g/mol]:</b>	146.23
<b>CAS:</b>	22607-10-9

## Physical Properties

Property code	Value	Unit	Source
gf	-262.04	kJ/mol	Joback Method
hf	-523.47	kJ/mol	Joback Method
hfus	17.61	kJ/mol	Joback Method
hvap	65.98	kJ/mol	Joback Method
log10ws	-1.92		Crippen Method
logp	1.308		Crippen Method
mcvol	135.320	ml/mol	McGowan Method
pc	3135.00	kPa	Joback Method
tb	565.92	K	Joback Method
tc	727.66	K	Joback Method
tf	271.56	K	Joback Method
vc	0.509	m <sup>3</sup> /kmol	Joback Method

## Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	345.58	J/mol×K	565.92	Joback Method
cpg	356.16	J/mol×K	592.88	Joback Method
cpg	366.31	J/mol×K	619.83	Joback Method
cpg	376.04	J/mol×K	646.79	Joback Method
cpg	385.36	J/mol×K	673.75	Joback Method
cpg	394.29	J/mol×K	700.71	Joback Method
cpg	402.84	J/mol×K	727.66	Joback Method
dvisc	0.2642980	Paxs	271.56	Joback Method

dvisc	0.0200295	Paxs	320.62	Joback Method
dvisc	0.0030104	Paxs	369.68	Joback Method
dvisc	0.0007054	Paxs	418.74	Joback Method
dvisc	0.0002241	Paxs	467.80	Joback Method
dvisc	0.0000885	Paxs	516.86	Joback Method
dvisc	0.0000411	Paxs	565.92	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=C22607109&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=C22607109&amp;Units=SI</a>

## Legend

<b>cp<sub>g</sub>:</b>	Ideal gas heat capacity
<b>dvisc:</b>	Dynamic viscosity
<b>g<sub>f</sub>:</b>	Standard Gibbs free energy of formation
<b>h<sub>f</sub>:</b>	Enthalpy of formation at standard conditions
<b>h<sub>fus</sub>:</b>	Enthalpy of fusion at standard conditions
<b>h<sub>vap</sub>:</b>	Enthalpy of vaporization at standard conditions
<b>log<sub>10</sub>w<sub>s</sub>:</b>	Log <sub>10</sub> of Water solubility in mol/l
<b>log<sub>p</sub>:</b>	Octanol/Water partition coefficient
<b>mc<sub>vol</sub>:</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/81-407-0/4-5-Octanediol.pdf>

Generated by Cheméo on 2024-04-25 09:26:17.602058211 +0000 UTC m=+16326426.522635524.

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