

# Cycloheptadecanol

<b>Other names:</b>	Civetol, dihydro-
<b>Inchi:</b>	InChI=1S/C17H34O/c18-17-15-13-11-9-7-5-3-1-2-4-6-8-10-12-14-16-17/h17-18H,1-16H2
<b>InchiKey:</b>	ZRLAJSMFDYQOTR-UHFFFAOYSA-N
<b>Formula:</b>	C17H34O
<b>SMILES:</b>	OC1CCCCCCCCCCCCCCC1
<b>Mol. weight [g/mol]:</b>	254.45
<b>CAS:</b>	4429-77-0

## Physical Properties

Property code	Value	Unit	Source
gf	-153.21	kJ/mol	Joback Method
hf	-559.88	kJ/mol	Joback Method
hfus	12.61	kJ/mol	Joback Method
hvap	72.44	kJ/mol	Joback Method
log10ws	-6.21		Crippen Method
logp	5.603		Crippen Method
mcvol	245.400	ml/mol	McGowan Method
pc	1927.05	kPa	Joback Method
tb	747.06	K	Joback Method
tc	981.66	K	Joback Method
tf	310.83	K	Joback Method
vc	0.852	m <sup>3</sup> /kmol	Joback Method

## Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	788.33	J/mol×K	747.06	Joback Method
cpg	815.34	J/mol×K	786.16	Joback Method
cpg	839.96	J/mol×K	825.26	Joback Method
cpg	862.15	J/mol×K	864.36	Joback Method
cpg	881.87	J/mol×K	903.46	Joback Method
cpg	899.07	J/mol×K	942.56	Joback Method
cpg	913.72	J/mol×K	981.66	Joback Method
dvisc	0.0738836	Paxs	310.83	Joback Method

dvisc	0.0017879	Paxs	383.53	Joback Method
dvisc	0.0001417	Paxs	456.24	Joback Method
dvisc	0.0000225	Paxs	528.94	Joback Method
dvisc	0.0000056	Paxs	601.65	Joback Method
dvisc	0.0000019	Paxs	674.36	Joback Method
dvisc	0.0000008	Paxs	747.06	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=C4429770&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=C4429770&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

Latest version available from:

<https://www.chemeo.com/cid/79-909-6/Cycloheptadecanol.pdf>

Generated by Cheméo on 2025-12-05 19:15:34.946469959 +0000 UTC m=+4710332.476510612.

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