

5-Tetradecanol

Other names:	tetradecan-5-ol
Inchi:	InChI=1S/C14H30O/c1-3-5-7-8-9-10-11-13-14(15)12-6-4-2/h14-15H,3-13H2,1-2H3
InchiKey:	UCEOYIRETCABER-UHFFFAOYSA-N
Formula:	C14H30O
SMILES:	CCCCCCCCCCC(O)CCCC
Mol. weight [g/mol]:	214.39
CAS:	21078-83-1

Physical Properties

Property code	Value	Unit	Source
gf	-72.26	kJ/mol	Joback Method
hf	-489.80	kJ/mol	Joback Method
hfus	32.58	kJ/mol	Joback Method
hvap	63.05	kJ/mol	Joback Method
log10ws	-5.06		Crippen Method
logp	4.678		Crippen Method
mcvol	213.990	ml/mol	McGowan Method
pc	1649.77	kPa	Joback Method
rinpol	1672.00		NIST Webbook
rinpol	1672.00		NIST Webbook
ripol	1974.00		NIST Webbook
tb	611.46	K	Joback Method
tc	771.56	K	Joback Method
tf	293.36	K	Joback Method
vc	0.833	m3/kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	583.00	J/mol×K	611.46	Joback Method
cpg	599.08	J/mol×K	638.14	Joback Method
cpg	614.51	J/mol×K	664.83	Joback Method
cpg	629.31	J/mol×K	691.51	Joback Method
cpg	643.50	J/mol×K	718.19	Joback Method

cpg	657.10	J/mol×K	744.87	Joback Method
cpg	670.12	J/mol×K	771.56	Joback Method
dvisc	0.0226239	Pa×s	293.36	Joback Method
dvisc	0.0038664	Pa×s	346.38	Joback Method
dvisc	0.0010562	Pa×s	399.39	Joback Method
dvisc	0.0003911	Pa×s	452.41	Joback Method
dvisc	0.0001784	Pa×s	505.43	Joback Method
dvisc	0.0000944	Pa×s	558.44	Joback Method
dvisc	0.0000558	Pa×s	611.46	Joback Method

Correlations

Information	Value
Property code	pvap
Equation	$\ln(P_{vp}) = A + B/(T + C)$
Coeff. A	1.62855e+01
Coeff. B	-5.38620e+03
Coeff. C	-9.60380e+01
Temperature range (K), min.	432.72
Temperature range (K), max.	586.85

Sources

The Yaws Handbook of Vapor Pressure:
Crippen Method:

<https://www.sciencedirect.com/book/9780128029992/the-yaws-handbook-of-vapor-pressure>
<http://pubs.acs.org/doi/abs/10.1021/ci990307l>

Crippen Method:

https://www.chemeo.com/doc/models/crippen_log10ws

Joback Method:

https://en.wikipedia.org/wiki/Joback_method

McGowan Method:

<http://link.springer.com/article/10.1007/BF02311772>

NIST Webbook:

<http://webbook.nist.gov/cgi/cbook.cgi?ID=C21078831&Units=SI>

Legend

cpg: Ideal gas heat capacity
dvisc: Dynamic viscosity
gf: Standard Gibbs free energy of formation
hf: Enthalpy of formation at standard conditions

hfus:	Enthalpy of fusion at standard conditions
hvap:	Enthalpy of vaporization at standard conditions
log10ws:	Log10 of Water solubility in mol/l
logp:	Octanol/Water partition coefficient
mcvol:	McGowan's characteristic volume
pc:	Critical Pressure
pvap:	Vapor pressure
rinpol:	Non-polar retention indices
ripol:	Polar retention indices
tb:	Normal Boiling Point Temperature
tc:	Critical Temperature
tf:	Normal melting (fusion) point
vc:	Critical Volume

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