

2-Ethyl-1-dodecanol

Inchi:	InChI=1S/C14H30O/c1-3-5-6-7-8-9-10-11-12-14(4-2)13-15/h14-15H,3-13H2,1-2H3
InchiKey:	SBAZXXVUILTVKK-UHFFFAOYSA-N
Formula:	C14H30O
SMILES:	CCCCCCCCCCC(CC)CO
Mol. weight [g/mol]:	214.39
CAS:	19780-33-7

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	-4.70		Crippen Method
logp	4.536		Crippen Method
mcvol	213.990	ml/mol	McGowan Method
pc	1649.77	kPa	Joback Method
ripol	2090.00		NIST Webbook
ripol	2090.00		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

Sources

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=C19780337&Units=SI
Crippen Method:	http://pubs.acs.org/doi/abs/10.1021/ci990307l

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
ripol:	Polar retention indices
tb:	Normal Boiling Point Temperature
tc:	Critical Temperature
tf:	Normal melting (fusion) point
vc:	Critical Volume

Latest version available from:

<https://www.chemeo.com/cid/74-163-9/2-Ethyl-1-dodecanol.pdf>

Generated by Cheméo on 2025-12-22 22:10:47.445093809 +0000 UTC m=+6189644.975134460.

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