

6-Tridecanol, 3,9-diethyl-

Other names:	3,9-Diethyl-6-hydroxytridecane 3,9-Diethyl-6-tridecanol 3,9-Diethyltridecanol-6
Inchi:	InChI=1S/C17H36O/c1-5-9-10-16(8-4)12-14-17(18)13-11-15(6-2)7-3/h15-18H,5-14H2,1-
InchiKey:	NWDHHLKROQOLMQ-UHFFFAOYSA-N
Formula:	C17H36O
SMILES:	CCCCC(CC)CCC(O)CCC(CC)CC
Mol. weight [g/mol]:	256.47
CAS:	123-24-0

Physical Properties

Property code	Value	Unit	Source
gf	-51.88	kJ/mol	Joback Method
hf	-562.28	kJ/mol	Joback Method
hfus	33.30	kJ/mol	Joback Method
hvap	68.95	kJ/mol	Joback Method
log10ws	-5.83		Crippen Method
logp	5.560		Crippen Method
mcvol	256.260	ml/mol	McGowan Method
pc	1337.84	kPa	Joback Method
tb	582.15	K	NIST Webbook
tc	844.95	K	Joback Method
tf	297.17	K	Joback Method
vc	0.989	m ³ /kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	750.73	J/mol×K	679.22	Joback Method
cpg	833.35	J/mol×K	817.33	Joback Method
cpg	818.31	J/mol×K	789.71	Joback Method
cpg	802.55	J/mol×K	762.09	Joback Method
cpg	786.05	J/mol×K	734.46	Joback Method
cpg	768.79	J/mol×K	706.84	Joback Method

cpg	847.71	J/mol×K	844.95	Joback Method
dvisc	0.0000274	Paxs	679.22	Joback Method
dvisc	0.0000485	Paxs	615.55	Joback Method
dvisc	0.0000978	Paxs	551.87	Joback Method
dvisc	0.0002370	Paxs	488.19	Joback Method
dvisc	0.0007487	Paxs	424.52	Joback Method
dvisc	0.0035502	Paxs	360.84	Joback Method
dvisc	0.0327970	Paxs	297.17	Joback Method

Sources

McGowan Method:	http://link.springer.com/article/10.1007/BF02311772
NIST Webbook:	http://webbook.nist.gov/cgi/cbook.cgi?ID=C123240&Units=SI
Crippen Method:	http://pubs.acs.org/doi/abs/10.1021/ci9903071
Crippen Method:	https://www.chemeo.com/doc/models/crippen_log10ws
Joback Method:	https://en.wikipedia.org/wiki/Joback_method

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
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/28-849-9/6-Tridecanol-3-9-diethyl.pdf>

Generated by Cheméo on 2024-04-24 03:44:21.905876375 +0000 UTC m=+16219510.826453726.

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