

2-Hydroxyethyl caprate

Inchi:	InChI=1S/C12H24O3/c1-2-3-4-5-6-7-8-9-12(14)15-11-10-13/h13H,2-11H2,1H3
InchiKey:	IGPIBLIHYRGADN-UHFFFAOYSA-N
Formula:	C12H24O3
SMILES:	CCCCCCCCC(=O)OCCO
Mol. weight [g/mol]:	216.32

Physical Properties

Property code	Value	Unit	Source
gf	-320.58	kJ/mol	Joback Method
hf	-688.04	kJ/mol	Joback Method
hfus	33.71	kJ/mol	Joback Method
hvap	68.14	kJ/mol	Joback Method
log10ws	-2.97		Crippen Method
logp	2.663		Crippen Method
mcvol	193.250	ml/mol	McGowan Method
pc	2007.30	kPa	Joback Method
rinpol	1595.00		NIST Webbook
rinpol	1595.00		NIST Webbook
tb	642.43	K	Joback Method
tc	808.95	K	Joback Method
tf	357.98	K	Joback Method
vc	0.750	m3/kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	530.16	J/molxK	642.43	Joback Method
cpg	592.71	J/molxK	781.19	Joback Method
cpg	581.31	J/molxK	753.44	Joback Method
cpg	569.37	J/molxK	725.69	Joback Method
cpg	556.86	J/molxK	697.94	Joback Method
cpg	543.80	J/molxK	670.18	Joback Method
cpg	603.57	J/molxK	808.95	Joback Method
dvisc	0.0000535	Paxs	642.43	Joback Method

dvisc	0.0000829	Paxs	595.02	Joback Method
dvisc	0.0001385	Paxs	547.61	Joback Method
dvisc	0.0002551	Paxs	500.20	Joback Method
dvisc	0.0005340	Paxs	452.80	Joback Method
dvisc	0.0013282	Paxs	405.39	Joback Method
dvisc	0.0042058	Paxs	357.98	Joback Method

Sources

NIST Webbook:	http://webbook.nist.gov/cgi/cbook.cgi?ID=R540455&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
McGowan Method:	http://link.springer.com/article/10.1007/BF02311772

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
rinpol:	Non-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/22-954-8/2-Hydroxyethyl-caprate.pdf>

Generated by Cheméo on 2024-04-17 01:18:13.680366875 +0000 UTC m=+15605942.600944192.

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