

DECANOIC ACID, CYCLOHEXYL ESTER

Other names:	Cyclohexyl caprate
Inchi:	InChI=1S/C16H30O2/c1-2-3-4-5-6-7-11-14-16(17)18-15-12-9-8-10-13-15/h15H,2-14H2,1
InchiKey:	FNQCJNQUVQXXJN-UHFFFAOYSA-N
Formula:	C16H30O2
SMILES:	CCCCCCCCCCC(=O)OC1CCCCC1
Mol. weight [g/mol]:	254.41

Physical Properties

Property code	Value	Unit	Source
af	0.7350		Cheméo Relay (1.0)
hf	-656.87	kJ/mol	Cheméo Relay (1.0)
hfus	31.82	kJ/mol	Joback Method
hvap	82.33	kJ/mol	Cheméo Relay (1.0)
ie	9.63	eV	Cheméo Relay (1.0)
log10ws	-5.28		Cheméo Relay (1.0)
logp	5.003		Crippen Method
mcvol	232.880	ml/mol	McGowan Method
pc	1596.17	kPa	Joback Method
tb	575.58	K	Cheméo Relay (1.0)
tc	779.56	K	Cheméo Relay (1.0)
tf	263.60	K	Cheméo Relay (1.0)
vc	0.879	m3/kmol	Cheméo Relay (1.0)

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	670.95	J/mol×K	661.32	Joback Method
cpg	778.03	J/mol×K	851.42	Joback Method
cpg	762.73	J/mol×K	819.74	Joback Method
cpg	746.44	J/mol×K	788.05	Joback Method
cpg	729.14	J/mol×K	756.37	Joback Method
cpg	710.81	J/mol×K	724.69	Joback Method
cpg	691.42	J/mol×K	693.00	Joback Method
dvisc	0.0003783	Pa×s	505.47	Joback Method

dvisc	0.0001289	Pa×s	661.32	Joback Method
dvisc	0.0001736	Pa×s	609.37	Joback Method
dvisc	0.0002471	Pa×s	557.42	Joback Method
dvisc	0.0028997	Pa×s	349.62	Joback Method
dvisc	0.0006384	Pa×s	453.52	Joback Method
dvisc	0.0012337	Pa×s	401.57	Joback Method

Sources

Crippen Method:	http://pubs.acs.org/doi/abs/10.1021/ci990307l
Crippen Method:	https://www.chemeo.com/doc/models/crippen_log10ws
Cheméo Relay (1.0):	
Cheméo Relay (1.0, beta):	
NIST Webbook:	http://webbook.nist.gov/cgi/cbook.cgi?ID=C1551402&Units=SI
Joback Method:	https://en.wikipedia.org/wiki/Joback_method
McGowan Method:	http://link.springer.com/article/10.1007/BF02311772

Legend

af:	Acentric Factor
cpg:	Ideal gas heat capacity
dvisc:	Dynamic viscosity
hf:	Enthalpy of formation at standard conditions
hfus:	Enthalpy of fusion at standard conditions
hvap:	Enthalpy of vaporization at standard conditions
ie:	Ionization energy
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

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