

Cyclohexyl iso-butyrate

Other names:	Cyclohexyl iso-butyrate
Inchi:	InChI=1S/C10H18O2/c1-8(2)10(11)12-9-6-4-3-5-7-9/h8-9H,3-7H2,1-2H3
InchiKey:	IYYXBWAISHVFEX-UHFFFAOYSA-N
Formula:	C10H18O2
SMILES:	CC(C)C(=O)OC1CCCCC1
Mol. weight [g/mol]:	170.25

Physical Properties

Property code	Value	Unit	Source
hf	-533.91	kJ/mol	Cheméo Relay (1.0)
hfus	12.75	kJ/mol	Joback Method
hvap	55.95	kJ/mol	Cheméo Relay (1.0)
log10ws	-2.62		Cheméo Relay (1.0)
logp	2.518		Crippen Method
mcvol	148.340	ml/mol	McGowan Method
tb	470.42	K	Cheméo Relay (1.0)
tf	221.55	K	Cheméo Relay (1.0)

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	355.54	J/mol×K	523.60	Joback Method
cpg	373.86	J/mol×K	558.58	Joback Method
cpg	391.23	J/mol×K	593.57	Joback Method
cpg	407.67	J/mol×K	628.55	Joback Method
cpg	423.20	J/mol×K	663.54	Joback Method
cpg	437.82	J/mol×K	698.52	Joback Method
cpg	451.55	J/mol×K	733.51	Joback Method
dvisc	0.0060058	Pa×s	267.00	Joback Method
dvisc	0.0023895	Pa×s	309.77	Joback Method
dvisc	0.0011889	Pa×s	352.53	Joback Method
dvisc	0.0006880	Pa×s	395.30	Joback Method
dvisc	0.0004430	Pa×s	438.07	Joback Method
dvisc	0.0003085	Pa×s	480.83	Joback Method

Sources

Crippen Method:	http://pubs.acs.org/doi/abs/10.1021/ci990307I
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=C1129471&Units=SI
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
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
tb:	Normal Boiling Point Temperature
tf:	Normal melting (fusion) point

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