

Carbonic acid, (1R)-(-)-menthyl heptyl ester

Inchi: InChI=1S/C18H34O3/c1-5-6-7-8-9-12-20-18(19)21-17-13-16(14(2)3)11-10-15(17)4/h14-17
InchiKey: NAEWNDFLNZKAME-UHFFFAOYSA-N
Formula: C18H34O3
SMILES: CCCCCCOC(=O)OC1CC(C(C)C)CCC1C
Mol. weight [g/mol]: 298.46

Physical Properties

Property code	Value	Unit	Source
gf	-231.65	kJ/mol	Joback Method
hf	-783.51	kJ/mol	Joback Method
hfus	36.80	kJ/mol	Joback Method
hvap	66.65	kJ/mol	Joback Method
log10ws	-5.57		Crippen Method
logp	5.571		Crippen Method
mvol	266.930	ml/mol	McGowan Method
pc	1303.29	kPa	Joback Method
rinpol	1960.00		NIST Webbook
rinpol	1960.00		NIST Webbook
tb	719.72	K	Joback Method
tc	908.95	K	Joback Method
tf	370.91	K	Joback Method
vc	1.010	m ³ /kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	823.23	J/molxK	719.72	Joback Method
cpg	844.75	J/molxK	751.26	Joback Method
cpg	865.07	J/molxK	782.80	Joback Method
cpg	884.20	J/molxK	814.34	Joback Method
cpg	902.13	J/molxK	845.88	Joback Method
cpg	918.89	J/molxK	877.42	Joback Method
cpg	934.47	J/molxK	908.95	Joback Method
dvisc	0.0018011	Paxs	370.91	Joback Method

dvisc	0.0008175	Paxs	429.04	Joback Method
dvisc	0.0004480	Paxs	487.18	Joback Method
dvisc	0.0002791	Paxs	545.32	Joback Method
dvisc	0.0001905	Paxs	603.45	Joback Method
dvisc	0.0001390	Paxs	661.59	Joback Method
dvisc	0.0001068	Paxs	719.72	Joback Method

Sources

McGowan Method:	http://link.springer.com/article/10.1007/BF02311772
NIST Webbook:	http://webbook.nist.gov/cgi/cbook.cgi?ID=U392436&Units=SI
Crippen Method:	http://pubs.acs.org/doi/abs/10.1021/ci9903071
Crippen Method:	https://www.cheméo.com/doc/models/crippen_log10ws
Joback Method:	https://en.wikipedia.org/wiki/Joback_method

Legend

cp_g:	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
h_{vap}:	Enthalpy of vaporization at standard conditions
log₁₀ws:	Log ₁₀ of Water solubility in mol/l
log_p:	Octanol/Water partition coefficient
m_{cvol}:	McGowan's characteristic volume
pc:	Critical Pressure
rin_{pol}:	Non-polar retention indices
tb:	Normal Boiling Point Temperature
tc:	Critical Temperature
tf:	Normal melting (fusion) point
vc:	Critical Volume

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