

Cyclopropanecarboxylic acid, 2-methylpentyl ester

Inchi:	InChI=1S/C10H18O2/c1-3-4-8(2)7-12-10(11)9-5-6-9/h8-9H,3-7H2,1-2H3
InchiKey:	YQWUBVGTPXCEQE-UHFFFAOYSA-N
Formula:	C10H18O2
SMILES:	CCCC(C)COC(=O)C1CC1
Mol. weight [g/mol]:	170.25

Physical Properties

Property code	Value	Unit	Source
gf	-142.29	kJ/mol	Joback Method
hf	-427.01	kJ/mol	Joback Method
hfus	19.05	kJ/mol	Joback Method
hvap	46.53	kJ/mol	Joback Method
log10ws	-2.28		Crippen Method
logp	2.376		Crippen Method
mcvol	148.340	ml/mol	McGowan Method
pc	2515.07	kPa	Joback Method
rinpol	1219.00		NIST Webbook
tb	510.79	K	Joback Method
tc	699.47	K	Joback Method
tf	277.56	K	Joback Method
vc	0.571	m3/kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	354.31	J/molxK	510.79	Joback Method
cpg	369.77	J/molxK	542.24	Joback Method
cpg	384.48	J/molxK	573.68	Joback Method
cpg	398.46	J/molxK	605.13	Joback Method
cpg	411.74	J/molxK	636.58	Joback Method
cpg	424.34	J/molxK	668.02	Joback Method
cpg	436.30	J/molxK	699.47	Joback Method
dvisc	0.0027998	Paxs	277.56	Joback Method
dvisc	0.0016887	Paxs	316.43	Joback Method

dvisc	0.0011377	Paxs	355.30	Joback Method
dvisc	0.0008286	Paxs	394.17	Joback Method
dvisc	0.0006388	Paxs	433.05	Joback Method
dvisc	0.0005141	Paxs	471.92	Joback Method
dvisc	0.0004276	Paxs	510.79	Joback Method

Sources

Joback Method:	https://en.wikipedia.org/wiki/Joback_method
McGowan Method:	http://link.springer.com/article/10.1007/BF02311772
NIST Webbook:	http://webbook.nist.gov/cgi/cbook.cgi?ID=U354658&Units=SI
Crippen Method:	http://pubs.acs.org/doi/abs/10.1021/ci9903071
Crippen Method:	https://www.cheméo.com/doc/models/crippen_log10ws

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.cheméo.com/cid/45-998-5/Cyclopropanecarboxylic-acid-2-methylpentyl-ester.pdf>

Generated by Cheméo on 2024-04-24 02:07:15.516449171 +0000 UTC m=+16213684.437026499.

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