

Propanoic acid, 2,3-dibromo, 3-methylbutyl ester

Inchi:	InChI=1S/C8H14Br2O2/c1-6(2)3-4-12-8(11)7(10)5-9/h6-7H,3-5H2,1-2H3
InchiKey:	JMGSCPGYIYOURH-UHFFFAOYSA-N
Formula:	C8H14Br2O2
SMILES:	CC(C)CCOC(=O)C(Br)CBr
Mol. weight [g/mol]:	302.00

Physical Properties

Property code	Value	Unit	Source
gf	-193.68	kJ/mol	Joback Method
hf	-411.15	kJ/mol	Joback Method
hfus	22.79	kJ/mol	Joback Method
hvap	54.65	kJ/mol	Joback Method
log10ws	-2.77		Crippen Method
logp	2.734		Crippen Method
mcvol	166.020	ml/mol	McGowan Method
pc	3135.00	kPa	Joback Method
rinpola	1383.00		NIST Webbook
tb	590.17	K	Joback Method
tc	801.33	K	Joback Method
tf	341.68	K	Joback Method
vc	0.620	m ³ /kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	358.39	J/molxK	590.17	Joback Method
cpg	370.14	J/molxK	625.36	Joback Method
cpg	381.23	J/molxK	660.56	Joback Method
cpg	391.68	J/molxK	695.75	Joback Method
cpg	401.52	J/molxK	730.94	Joback Method
cpg	410.76	J/molxK	766.13	Joback Method
cpg	419.43	J/molxK	801.33	Joback Method
dvisc	0.0027803	Paxs	341.68	Joback Method
dvisc	0.0014410	Paxs	383.10	Joback Method

dvisc	0.0008491	Paxs	424.51	Joback Method
dvisc	0.0005496	Paxs	465.92	Joback Method
dvisc	0.0003819	Paxs	507.34	Joback Method
dvisc	0.0002804	Paxs	548.75	Joback Method
dvisc	0.0002150	Paxs	590.17	Joback Method

Sources

McGowan Method:	http://link.springer.com/article/10.1007/BF02311772
NIST Webbook:	http://webbook.nist.gov/cgi/cbook.cgi?ID=R30193&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

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

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