Propanoic acid, 2-bromo-, ethyl ester

Other names:	Propionic acid, 2-bromo-, ethyl ester
	«alpha»-Bromopropionic acid ethyl ester
	Ethyl «alpha»-bromopropanoate
	Ethyl «alpha»-bromopropionate
	Ethyl 2-bromopropanoate
	Ethyl 2-bromopropionate
	Ethyl dl-2-bromopropionate
	2-Bromopropanoic acid ethyl ester
	(.+/-.)-2-Bromopropanoic acid ethyl ester
	NSC 6753
Inchi:	InChI=1S/C5H9BrO2/c1-3-8-5(7)4(2)6/h4H,3H2,1-2H3
InchiKey:	ARFLASKVLJTEJD-UHFFFAOYSA-N
Formula:	C5H9BrO2
SMILES:	CCOC(=O)C(C)Br
Mol. weight [g/mol]:	181.03
CAS:	535-11-5

Physical Properties

Property code	Value	Unit	Source
gf	-230.82	kJ/mol	Joback Method
hf	-370.28	kJ/mol	Joback Method
hfus	13.25	kJ/mol	Joback Method
hvap	41.93	kJ/mol	Joback Method
log10ws	-1.32		Crippen Method
logp	1.333		Crippen Method
mcvol	106.250	ml/mol	McGowan Method
pc	4031.24	kPa	Joback Method
rinpol	909.00		NIST Webbook
rinpol	907.00		NIST Webbook
rinpol	906.00		NIST Webbook
rinpol	931.00		NIST Webbook
ripol	1327.00		NIST Webbook
ripol	1340.00		NIST Webbook
tb	433.00	K	NIST Webbook
tb	431.20	K	NIST Webbook
tc	657.09	K	Joback Method
tf	263.07	K	Joback Method

vc	0.396	m3/kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	196.09	J/mol×K	455.81	Joback Method
cpg	236.22	J/mol×K	623.55	Joback Method
cpg	228.93	J/mol×K	590.00	Joback Method
cpg	221.27	J/mol×K	556.45	Joback Method
cpg	213.25	J/mol×K	522.90	Joback Method
cpg	204.86	J/mol×K	489.36	Joback Method
cpg	243.15	J/mol×K	657.09	Joback Method
dvisc	0.0003362	Pa×s	455.81	Joback Method
dvisc	0.0004295	Pa×s	423.69	Joback Method
dvisc	0.0005712	Pa×s	391.56	Joback Method
dvisc	0.0007993	Pa×s	359.44	Joback Method
dvisc	0.0011948	Pa×s	327.32	Joback Method
dvisc	0.0019491	Pa×s	295.19	Joback Method
dvisc	0.0035835	Pa×s	263.07	Joback Method

Sources

Crippen Method: http://pubs.acs.org/doi/abs/10.1021/ci990307l
Crippen Method: https://www.chemeo.com/doc/models/crippen_log10ws
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=C535115&Units=SI

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
ripol:	Polar retention indices
tb:	Normal Boiling Point Temperature
tc:	Critical Temperature
tf:	Normal melting (fusion) point
vc:	Critical Volume