

Butanoic acid, 2-chloro, 4-pentenyl ester

Inchi:	InChI=1S/C9H15ClO2/c1-3-5-6-7-12-9(11)8(10)4-2/h3,8H,1,4-7H2,2H3
InchiKey:	OADPPHWDKBRNQP-UHFFFAOYSA-N
Formula:	C9H15ClO2
SMILES:	C=CCCCOC(=O)C(Cl)CC
Mol. weight [g/mol]:	190.67

Physical Properties

Property code	Value	Unit	Source
gf	-135.55	kJ/mol	Joback Method
hf	-369.48	kJ/mol	Joback Method
hfus	21.25	kJ/mol	Joback Method
hvap	48.11	kJ/mol	Joback Method
log10ws	-2.57		Crippen Method
logp	2.513		Crippen Method
mcvol	153.050	ml/mol	McGowan Method
pc	2443.48	kPa	Joback Method
rinpol	1211.00		NIST Webbook
rinpol	1204.00		NIST Webbook
rinpol	1218.00		NIST Webbook
rinpol	1189.00		NIST Webbook
rinpol	1204.00		NIST Webbook
ripol	1671.00		NIST Webbook
ripol	1692.00		NIST Webbook
ripol	1671.00		NIST Webbook
tb	515.28	K	Joback Method
tc	701.54	K	Joback Method
tf	276.51	K	Joback Method
vc	0.588	m3/kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	333.91	J/molxK	515.28	Joback Method
cpg	346.42	J/molxK	546.32	Joback Method

cpg	358.38	J/mol×K	577.37	Joback Method
cpg	369.80	J/mol×K	608.41	Joback Method
cpg	380.68	J/mol×K	639.46	Joback Method
cpg	391.04	J/mol×K	670.50	Joback Method
cpg	400.89	J/mol×K	701.54	Joback Method
dvisc	0.0038090	Paxs	276.51	Joback Method
dvisc	0.0017815	Paxs	316.31	Joback Method
dvisc	0.0009875	Paxs	356.10	Joback Method
dvisc	0.0006163	Paxs	395.89	Joback Method
dvisc	0.0004192	Paxs	435.69	Joback Method
dvisc	0.0003042	Paxs	475.48	Joback Method
dvisc	0.0002319	Paxs	515.28	Joback Method

Sources

McGowan Method:	http://link.springer.com/article/10.1007/BF02311772
NIST Webbook:	http://webbook.nist.gov/cgi/cbook.cgi?ID=R28661&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
ripol:	Polar retention indices
tb:	Normal Boiling Point Temperature
tc:	Critical Temperature
tf:	Normal melting (fusion) point

vc: Critical Volume

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