

4-Butylbenzoic acid, 3-chloroprop-2-enyl ester

Inchi:

InchiKey:

Formula:

SMILES:

Mol. weight [g/mol]:

InChI=1S/C14H17ClO2/c1-2-3-5-12-6-8-13(9-7-12)14(16)17-11-4-10-15/h4,6-10H,2-3,5,13H

FDTZLBPFUIPWST-ONNFQVAWSA-N

C14H17ClO2

CCCCc1ccc(C(=O)OCC=CCl)cc1

252.74

Physical Properties

Property code	Value	Unit	Source
gf	4.15	kJ/mol	Joback Method
hf	-250.55	kJ/mol	Joback Method
hfus	32.85	kJ/mol	Joback Method
hvap	63.20	kJ/mol	Joback Method
log10ws	-4.69		Crippen Method
logp	3.938		Crippen Method
mcvol	199.740	ml/mol	McGowan Method
pc	2094.58	kPa	Joback Method
rinpol	1930.40		NIST Webbook
tb	669.26	K	Joback Method
tc	881.69	K	Joback Method
tf	383.48	K	Joback Method
vc	0.764	m3/kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	500.61	J/mol×K	669.26	Joback Method
cpg	515.20	J/mol×K	704.67	Joback Method
cpg	528.88	J/mol×K	740.07	Joback Method
cpg	541.68	J/mol×K	775.48	Joback Method
cpg	553.66	J/mol×K	810.88	Joback Method
cpg	564.85	J/mol×K	846.29	Joback Method
cpg	575.29	J/mol×K	881.69	Joback Method
dvisc	0.0012851	Pa×s	383.48	Joback Method
dvisc	0.0006964	Pa×s	431.11	Joback Method

dvisc	0.0004263	Pa×s	478.74	Joback Method
dvisc	0.0002852	Pa×s	526.37	Joback Method
dvisc	0.0002040	Pa×s	574.00	Joback Method
dvisc	0.0001535	Pa×s	621.63	Joback Method
dvisc	0.0001204	Pa×s	669.26	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=U292529&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
tb:	Normal Boiling Point Temperature
tc:	Critical Temperature
tf:	Normal melting (fusion) point
vc:	Critical Volume

Latest version available from:

<https://www.chemeo.com/cid/46-750-8/4-Butylbenzoic-acid-3-chloroprop-2-enyl-ester.pdf>

Generated by Cheméo on 2025-12-05 22:18:37.364227728 +0000 UTC m=+4721314.894268395.

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