

2,3-Dichlorophenol, tert-butyldimethylsilyl ether

Other names:	2,3-dichlorophenol, TBDMS 2,3-Dichlorophenol, tbdms derivative
Inchi:	InChI=1S/C12H18Cl2OSi/c1-12(2,3)16(4,5)15-10-8-6-7-9(13)11(10)14/h6-8H,1-5H3
InchiKey:	PIBTZLDTFZSVPX-UHFFFAOYSA-N
Formula:	C12H18Cl2OSi
SMILES:	CC(C)(C)[Si](C)(C)Oc1cccc(Cl)c1Cl
Mol. weight [g/mol]:	277.26

Physical Properties

Property code	Value	Unit	Source
log10ws	-3.11		Crippen Method
logp	5.377		Crippen Method
rinpol	1663.30		NIST Webbook

Sources

NIST Webbook:	http://webbook.nist.gov/cgi/cbook.cgi?ID=U333420&Units=SI
Crippen Method:	http://pubs.acs.org/doi/abs/10.1021/ci9903071
Crippen Method:	https://www.chemeo.com/doc/models/crippen_log10ws

Legend

log10ws:	Log10 of Water solubility in mol/l
logp:	Octanol/Water partition coefficient
rinpol:	Non-polar retention indices

Latest version available from:

<https://www.chemeo.com/cid/60-636-9/2-3-Dichlorophenol-tert-butyldimethylsilyl-ether.pdf>

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