2,3-Dimethylphenol, tert-butyldimethylsilyl ether

Other names: 2,3-Dimethylphenol, tbdms derivative
Inchi: InChI=1S/C14H24OSi/c1-11-9-8-10-13(12(11)2)15-16(6,7)14(3,4)5/h8-10H,1-7H3
InchiKey: HNFBPICGQRTYTN-UHFFFAOYSA-N
Formula: C14H24OSi
SMILES: Cc1cccc(O[Si](C)(C)C(C)(C)C)c1C
Mol. weight [g/mol]: 236.43

Physical Properties

Property code	Value	Unit	Source
log10ws	-2.69		Crippen Method
logp	4.687		Crippen Method
rinpol	1501.20		NIST Webbook
rinpol	1501.20		NIST Webbook

Sources

Crippen Method: https://www.chemeo.com/doc/models/crippen_log10ws
NIST Webbook: http://webbook.nist.gov/cgi/cbook.cgi?ID=U333490&Units=SI
Crippen Method: http://pubs.acs.org/doi/abs/10.1021/ci990307l

Legend

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