

2-Chloroethanol, benzyldimethylsilyl ether

Inchi: InChI=1S/C11H17ClOSi/c1-14(2,13-9-8-12)10-11-6-4-3-5-7-11/h3-7H,8-10H2,1-2H3
InchiKey: VNCFRCZTGNMXIC-UHFFFAOYSA-N
Formula: C11H17ClOSi
SMILES: C[Si](C)(Cc1ccccc1)OCCCl
Mol. weight [g/mol]: 228.79

Physical Properties

Property code	Value	Unit	Source
log10ws	-0.91		Crippen Method
logp	3.229		Crippen Method
rinpol	1484.00		NIST Webbook
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Sources

Crippen Method: <http://pubs.acs.org/doi/abs/10.1021/ci9903071>
Crippen Method: https://www.chemeo.com/doc/models/crippen_log10ws
NIST Webbook: <http://webbook.nist.gov/cgi/cbook.cgi?ID=U375539&Units=SI>

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/39-148-5/2-Chloroethanol-benzyldimethylsilyl-ether.pdf>

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