

# 2,6-Dichlorobenzyl alcohol, benzyldimethylsilyl ether

**Inchi:** InChI=1S/C16H18Cl2OSi/c1-20(2,12-13-7-4-3-5-8-13)19-11-14-15(17)9-6-10-16(14)18/h  
**InchiKey:** PSTOIZGTWGDDCK-UHFFFAOYSA-N  
**Formula:** C16H18Cl2OSi  
**SMILES:** C[Si](C)(Cc1cccc1)OCc1c(Cl)cccc1Cl  
**Mol. weight [g/mol]:** 325.31

## Physical Properties

Property code	Value	Unit	Source
log10ws	-3.82		Crippen Method
logp	5.497		Crippen Method
rinpol	2149.00		NIST Webbook
rinpol	2149.00		NIST Webbook

## Sources

**Crippen Method:** [https://www.chemeo.com/doc/models/crippen\\_log10ws](https://www.chemeo.com/doc/models/crippen_log10ws)  
**NIST Webbook:** <http://webbook.nist.gov/cgi/cbook.cgi?ID=U376094&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

Latest version available from:

<https://www.chemeo.com/cid/119-649-0/2-6-Dichlorobenzyl-alcohol-benzyldimethylsilyl-ether.pdf>

Generated by Cheméo on 2024-04-29 12:34:13.848377518 +0000 UTC m=+16683302.768954829.

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