

# 2,3,6-Trifluorobenzyl alcohol, bromomethyl dimethylsilyl ether

**Inchi:** InChI=1S/C10H12BrF3OSi/c1-16(2,6-11)15-5-7-8(12)3-4-9(13)10(7)14/h3-4H,5-6H2,1-2H  
**InchiKey:** VIZSKEQTZWMTOS-UHFFFAOYSA-N  
**Formula:** C10H12BrF3OSi  
**SMILES:** C[Si](C)(CBr)OCc1c(F)ccc(F)c1F  
**Mol. weight [g/mol]:** 313.19

## Physical Properties

Property code	Value	Unit	Source
log10ws	-2.26		Crippen Method
logp	3.760		Crippen Method
rinpol	1476.00		NIST Webbook

## Sources

**Crippen Method:** <http://pubs.acs.org/doi/abs/10.1021/ci990307l>  
**Crippen Method:** [https://www.cheméo.com/doc/models/crippen\\_log10ws](https://www.cheméo.com/doc/models/crippen_log10ws)  
**NIST Webbook:** <http://webbook.nist.gov/cgi/cbook.cgi?ID=U376063&Units=SI>

## Legend

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

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