

2,4,5-Trifluorobenzyl alcohol, bromomethyl dimethylsilyl ether

Inchi: InChI=1S/C10H12BrF3OSi/c1-16(2,6-11)15-5-7-3-9(13)10(14)4-8(7)12/h3-4H,5-6H2,1-2H
InchiKey: DSCGVIHXSAFDMU-UHFFFAOYSA-N
Formula: C₁₀H₁₂BrF₃OSi
SMILES: C[Si](C)(CBr)OCc1cc(F)c(F)cc1F
Mol. weight [g/mol]: 313.19

Physical Properties

Property code	Value	Unit	Source
log10ws	-2.26		Crippen Method
logp	3.760		Crippen Method
rinpol	1455.00		NIST Webbook

Sources

Crippen Method: https://www.chemeo.com/doc/models/crippen_log10ws
NIST Webbook: <http://webbook.nist.gov/cgi/cbook.cgi?ID=U376066&Units=SI>
Crippen Method: <http://pubs.acs.org/doi/abs/10.1021/ci9903071>

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/67-294-2/2-4-5-Trifluorobenzyl-alcohol-bromomethyl-dimethylsilyl-ether.pdf>

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