

# 1,3-Difluoro-5-dodecyldimethylsilyloxybenzene

**Inchi:** InChI=1S/C20H34F2OSi/c1-4-5-6-7-8-9-10-11-12-13-14-24(2,3)23-20-16-18(21)15-19(22)  
**InchiKey:** MIINNZRLPNSYGU-UHFFFAOYSA-N  
**Formula:** C20H34F2OSi  
**SMILES:** CCCCCCCCCC[Si](C)(C)Oc1cc(F)cc(F)c1  
**Mol. weight [g/mol]:** 356.57

## Physical Properties

Property code	Value	Unit	Source
log10ws	-5.75		Crippen Method
logp	7.470		Crippen Method
rinpol	2062.00		NIST Webbook
rinpol	2062.00		NIST Webbook

## Sources

**Crippen Method:** <http://pubs.acs.org/doi/abs/10.1021/ci9903071>  
**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=U299085&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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