

# 1-Methyl-2-dimethyl-(isopropyl)-silyloxybenzene

**Inchi:** InChI=1S/C12H20OSi/c1-10(2)14(4,5)13-12-9-7-6-8-11(12)3/h6-10H,1-5H3  
**InchiKey:** BQMFAMDAADWPHG-UHFFFAOYSA-N  
**Formula:** C12H20OSi  
**SMILES:** Cc1ccccc1O[Si](C)(C)C(C)C  
**Mol. weight [g/mol]:** 208.37  
**CAS:** 62790-80-1

## Physical Properties

Property code	Value	Unit	Source
log10ws	-1.79		Crippen Method
logp	3.989		Crippen Method
rinpol	1323.20		NIST Webbook

## Sources

**NIST Webbook:** <http://webbook.nist.gov/cgi/cbook.cgi?ID=C62790801&Units=SI>  
**Crippen Method:** <http://pubs.acs.org/doi/abs/10.1021/ci9903071>  
**Crippen Method:** [https://www.cheméo.com/doc/models/crippen\\_log10ws](https://www.cheméo.com/doc/models/crippen_log10ws)

## 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.cheméo.com/cid/62-560-1/1-Methyl-2-dimethyl-isopropyl-silyloxybenzene.pdf>

Generated by Cheméo on 2024-04-24 14:35:05.601633645 +0000 UTC m=+16258554.522210957.

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