

# (.+/-)-2-Phenylbutyric acid, benzyldimethylsilyl ester

**Inchi:** InChI=1S/C19H24O2Si/c1-4-18(17-13-9-6-10-14-17)19(20)21-22(2,3)15-16-11-7-5-8-12-  
**InchiKey:** GYCDBLOZFHPMC-UHFFFAOYSA-N  
**Formula:** C19H24O2Si  
**SMILES:** CCC(C(=O)O[Si](C)(C)Cc1ccccc1)c1ccccc1  
**Mol. weight [g/mol]:** 312.48

## Physical Properties

Property code	Value	Unit	Source
log10ws	-2.95		Crippen Method
logp	4.710		Crippen Method
rinpol	2025.00		NIST Webbook
rinpol	2025.00		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=U375607&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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