

Benzoic acid, 3-acetyloxy-, tert.-butyldimethylsilyl ester

Inchi: InChI=1S/C15H22O4Si/c1-11(16)18-13-9-7-8-12(10-13)14(17)19-20(5,6)15(2,3)4/h7-10H
InchiKey: BREGBMWULNVKBK-UHFFFAOYSA-N
Formula: C15H22O4Si
SMILES: CC(=O)Oc1cccc(C(=O)O[Si](C)(C)C(C)(C)C)c1
Mol. weight [g/mol]: 294.42

Physical Properties

Property code	Value	Unit	Source
log10ws	-2.17		Crippen Method
logp	3.774		Crippen Method
rinpol	1836.00		NIST Webbook
rinpol	1836.00		NIST Webbook

Sources

Crippen Method: <http://pubs.acs.org/doi/abs/10.1021/ci9903071>
Crippen Method: https://www.chemeo.com/doc/models/crippen_log10ws
NIST Webbook: <http://webbook.nist.gov/cgi/cbook.cgi?ID=U375027&Units=SI>

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/30-398-7/Benzoic-acid-3-acetyloxy-tert-butyldimethylsilyl-ester.pdf>

Generated by Cheméo on 2025-05-21 21:23:54.345663108 +0000 UTC m=+3171679.846107333.

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