

2'-Hydroxy-4'-methoxyacetophenone, tert-butyldimethylsilyl ether

Other names: 2'-Hydroxy-4'-methoxyacetophenone, tbdms derivative
Inchi: InChI=1S/C15H24O3Si/c1-11(16)13-9-8-12(17-5)10-14(13)18-19(6,7)15(2,3)4/h8-10H,1-
InchiKey: HYDIABVZVIXOKN-UHFFFAOYSA-N
Formula: C15H24O3Si
SMILES: COc1ccc(C(C)=O)c(O[Si](C)(C)C(C)(C)C)c1
Mol. weight [g/mol]: 280.43

Physical Properties

Property code	Value	Unit	Source
log10ws	-2.52		Crippen Method
logp	4.282		Crippen Method
rinpol	1871.20		NIST Webbook

Sources

Crippen Method: https://www.chemeo.com/doc/models/crippen_log10ws
NIST Webbook: <http://webbook.nist.gov/cgi/cbook.cgi?ID=U352931&Units=SI>
Crippen Method: <http://pubs.acs.org/doi/abs/10.1021/ci990307l>

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/31-016-9/2-Hydroxy-4-methoxyacetophenone-tert-butyldimethylsilyl-ether.pdf>

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