

Benzaldehyde, 4-hydroxy-3-methoxy, oxime, bis-TMS

Inchi: InChI=1S/C14H25NO3Si2/c1-16-14-10-12(11-15-18-20(5,6)7)8-9-13(14)17-19(2,3)4/h8-13
InchiKey: IIVLWJPBHXQTJK-RVDMUPIBSA-N
Formula: C14H25NO3Si2
SMILES: COc1cc(C=NO[Si](C)(C)C)ccc1O[Si](C)(C)C
Mol. weight [g/mol]: 311.52

Physical Properties

Property code	Value	Unit	Source
log10ws	0.37		Crippen Method
logp	4.094		Crippen Method
rinpol	1771.00		NIST Webbook

Sources

NIST Webbook: <http://webbook.nist.gov/cgi/cbook.cgi?ID=R100111&Units=SI>
Crippen Method: <http://pubs.acs.org/doi/abs/10.1021/ci990307l>
Crippen Method: https://www.chemeo.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.chemeo.com/cid/64-314-2/Benzaldehyde-4-hydroxy-3-methoxy-oxime-bis-TMS.pdf>

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