

Benzaldehyde, 3-hydroxy, oxime, bis-TMS

Inchi: InChI=1S/C13H23NO2Si2/c1-17(2,3)15-13-9-7-8-12(10-13)11-14-16-18(4,5)6/h7-11H,1-6H
InchiKey: CQRVGZXMJNCNTB-UHFFFAOYSA-N
Formula: C13H23NO2Si2
SMILES: C[Si](C)(C)ON=Cc1cccc(O[Si](C)(C)C)c1
Mol. weight [g/mol]: 281.50

Physical Properties

Property code	Value	Unit	Source
log10ws	0.49		Crippen Method
logp	4.086		Crippen Method
rinpol	1588.00		NIST Webbook
rinpol	1588.00		NIST Webbook

Sources

Crippen Method: <http://pubs.acs.org/doi/abs/10.1021/ci990307l>
Crippen Method: https://www.chemeo.com/doc/models/crippen_log10ws
NIST Webbook: <http://webbook.nist.gov/cgi/cbook.cgi?ID=R100077&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/53-670-9/Benzaldehyde-3-hydroxy-oxime-bis-TMS.pdf>

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