

3-Deoxy-threo-pentonic acid, tetrakis-TMS

Inchi: InChI=1S/C17H42O5Si4/c1-23(2,3)19-14-15(20-24(4,5)6)13-16(21-25(7,8)9)17(18)22-26
InchiKey: OQHUZALRVKVUOV-HOTGVXAUSA-N
Formula: C17H42O5Si4
SMILES: C[Si](C)(C)OCC(CC(O[Si](C)(C)C)C(=O)O[Si](C)(C)C)O[Si](C)(C)C
Mol. weight [g/mol]: 438.85

Physical Properties

Property code	Value	Unit	Source
log10ws	4.51		Crippen Method
logp	5.046		Crippen Method
rinpol	1698.00		NIST Webbook
rinpol	1698.00		NIST Webbook

Sources

NIST Webbook: <http://webbook.nist.gov/cgi/cbook.cgi?ID=R101387&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/33-913-1/3-Deoxy-threo-pentonic-acid-tetrakis-TMS.pdf>

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