

3,6-Dideoxy-ribo-hexonic acid, tetrakis-TMS

Inchi: InChI=1S/C18H44O5Si4/c1-15(20-24(2,3)4)16(21-25(5,6)7)14-17(22-26(8,9)10)18(19)23
InchiKey: XOPZJDYYVRROEQ-LRUHZDSUSA-N
Formula: C18H44O5Si4
SMILES: CC(O[Si](C)(C)C)C(CC(O[Si](C)(C)C)C(=O)O[Si](C)(C)C)O[Si](C)(C)C
Mol. weight [g/mol]: 452.88

Physical Properties

Property code	Value	Unit	Source
log10ws	3.98		Crippen Method
logp	5.435		Crippen Method
rinpol	1710.00		NIST Webbook

Sources

NIST Webbook: <http://webbook.nist.gov/cgi/cbook.cgi?ID=R101154&Units=SI>
Crippen Method: <http://pubs.acs.org/doi/abs/10.1021/ci9903071>
Crippen Method: https://www.cheméo.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.cheméo.com/cid/54-847-2/3-6-Dideoxy-ribo-hexonic-acid-tetrakis-TMS.pdf>

Generated by Cheméo on 2024-04-28 00:00:41.149914026 +0000 UTC m=+16551690.070491342.

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