2,6-Dideoxy-ribo-hexonic acid, tetrakis-TMS

Inchi: InChI=1S/C18H44O5Si4/c1-15(20-24(2,3)4)18(23-27(11,12)13)16(21-25(5,6)7)14-17(19)
InchiKey: UXZPSPDODIHYMZ-UHFFFAOYSA-N
Formula: C18H44O5Si4
SMILES: CC(O[Si](C)(C)C)C(O[Si](C)(C)C)C(CC(=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	1726.00		NIST Webbook
rinpol	1726.00		NIST Webbook

Sources

NIST Webbook: http://webbook.nist.gov/cgi/cbook.cgi?ID=R100804&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