

3-Deoxy-ribo-hexonic acid, pentakis-TMS

Inchi: InChI=1S/C21H52O6Si5/c1-28(2,3)23-17-20(26-31(10,11)12)18(24-29(4,5)6)16-19(25-30)21-22
InchiKey: VXOUZNJJSOAZDPY-AQNXPMDSA-N
Formula: C₂₁H₅₂O₆Si₅
SMILES: C[Si](C)(C)OCC(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]: 541.06

Physical Properties

Property code	Value	Unit	Source
log10ws	5.59		Crippen Method
logp	6.266		Crippen Method
rinpol	1920.00		NIST Webbook

Sources

NIST Webbook: <http://webbook.nist.gov/cgi/cbook.cgi?ID=R101347&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/45-924-6/3-Deoxy-ribo-hexonic-acid-pentakis-TMS.pdf>

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