

«beta»-2-Desoxyribose, TMS

Inchi: InChI=1S/C14H34O4Si3/c1-19(2,3)15-11-13-12(17-20(4,5)6)10-14(16-13)18-21(7,8)9/h1
InchiKey: BTOQHLYLAKILOAR-HZSPNIEDSA-N
Formula: C14H34O4Si3
SMILES: C[Si](C)(C)OCC1OC(O[Si](C)(C)C)CC1O[Si](C)(C)C
Mol. weight [g/mol]: 350.67

Physical Properties

Property code	Value	Unit	Source
logp	4.024		Crippen Method
rinpol	1502.00		NIST Webbook
rinpol	1517.00		NIST Webbook
rinpol	1502.00		NIST Webbook
rinpol	1517.00		NIST Webbook

Sources

NIST Webbook: <http://webbook.nist.gov/cgi/cbook.cgi?ID=R154371&Units=SI>
Crippen Method: <http://pubs.acs.org/doi/abs/10.1021/ci990307l>
Crippen Method: https://www.cheméo.com/doc/models/crippen_log10ws
Cheméo Relay (1.0):
Cheméo Relay (1.0, beta):

Legend

logp: Octanol/Water partition coefficient
rinpol: Non-polar retention indices

Latest version available from:

<https://www.cheméo.com/cid/113-872-8/beta-2-Desoxyribose-TMS.pdf>

Generated by Cheméo on 2026-07-24 18:41:20.189271759 +0000 UTC m=+429576.031157181.

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