

D-Arabinopyranose, tetrakis(trimethylsilyl) ether (isomer 2)

Inchi: InChI=1S/C17H42O5Si4/c1-23(2,3)19-14-13-18-17(22-26(10,11)12)16(21-25(7,8)9)15(14)
InchiKey: KEOUSSOURMHEKN-UHFFFAOYSA-N
Formula: C17H42O5Si4
SMILES: C[Si](C)(C)OC1COC(O[Si](C)(C)C)C(O[Si](C)(C)C)C1O[Si](C)(C)C
Mol. weight [g/mol]: 438.85

Physical Properties

Property code	Value	Unit	Source
log10ws	4.58		Crippen Method
logp	4.854		Crippen Method
rinpol	1636.50		NIST Webbook
rinpol	1636.50		NIST Webbook

Sources

Crippen Method: <http://pubs.acs.org/doi/abs/10.1021/ci9903071>
Crippen Method: https://www.chemeo.com/doc/models/crippen_log10ws
NIST Webbook: <http://webbook.nist.gov/cgi/cbook.cgi?ID=U380136&Units=SI>

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/33-373-1/D-Arabinopyranose-tetrakis-trimethylsilyl-ether-isomer-2.pdf>

Generated by Cheméo on 2024-04-26 17:37:53.435446369 +0000 UTC m=+16442322.356023691.

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