

D-(-)-Tagatose, pentakis(trimethylsilyl) ether, methyloxime (anti)

Inchi: InChI=1S/C22H55NO6Si5/c1-24-23-19(17-25-30(2,3)4)21(28-33(11,12)13)22(29-34(14,15)16)35-36-37-38-39-40
InchiKey: ACDMQGXCXUQPRR-UHFFFAOYSA-N
Formula: C22H55NO6Si5
SMILES: CON=C(CO[Si](C)(C)C)C(O[Si](C)(C)C)C(O[Si](C)(C)C)C(CO[Si](C)(C)C)O[Si](C)(C)C
Mol. weight [g/mol]: 570.10

Physical Properties

Property code	Value	Unit	Source
log10ws	5.70		Crippen Method
logp	6.352		Crippen Method
rinpol	1835.20		NIST Webbook

Sources

Crippen Method: https://www.chemeo.com/doc/models/crippen_log10ws
NIST Webbook: <http://webbook.nist.gov/cgi/cbook.cgi?ID=U380178&Units=SI>
Crippen Method: <http://pubs.acs.org/doi/abs/10.1021/ci9903071>

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/70-078-8/D-Tagatose-pentakis-trimethylsilyl-ether-methyloxime-anti.pdf>

Generated by Cheméo on 2024-04-20 10:49:02.037871875 +0000 UTC m=+15899390.958449187.

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