

D-(+)-Galactose, pentakis(trimethylsilyl) ether, benzyloxime (isomer 1)

Inchi: InChI=1S/C28H59NO6Si5/c1-36(2,3)31-23-26(33-38(7,8)9)28(35-40(13,14)15)27(34-39)
InchiKey: HTPNKYHYJRZYRL-UHFFFAOYSA-N
Formula: C28H59NO6Si5
SMILES: C[Si](C)(C)OCC(O[Si](C)(C)C)C(O[Si](C)(C)C)C(O[Si](C)(C)C)C(O[Si](C)(C)C)C(C=NOCc1ccccc1)O[Si](C)(C)C
Mol. weight [g/mol]: 646.20

Physical Properties

Property code	Value	Unit	Source
log10ws	3.48		Crippen Method
logp	7.921		Crippen Method
rinpol	2416.50		NIST Webbook

Sources

NIST Webbook: <http://webbook.nist.gov/cgi/cbook.cgi?ID=U380343&Units=SI>
Crippen Method: <http://pubs.acs.org/doi/abs/10.1021/ci9903071>
Crippen Method: https://www.cheméo.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.cheméo.com/cid/70-348-8/D-Galactose-pentakis-trimethylsilyl-ether-benzyloxime-isomer-1.pdf>

Generated by Cheméo on 2024-04-23 17:09:39.543947219 +0000 UTC m=+16181428.464524533.

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