

Pentanoic acid, 4-hydroxy-2-propyl, TMS, # 2

Inchi: InChI=1S/C14H32O3Si2/c1-9-10-13(14(15)17-19(6,7)8)11-12(2)16-18(3,4)5/h12-13H,9-14H
InchiKey: YWWMPRJWXQPBOQ-UHFFFAOYSA-N
Formula: C14H32O3Si2
SMILES: CCCC(CC(C)O[Si](C)(C)C)C(=O)O[Si](C)(C)C
Mol. weight [g/mol]: 304.57

Physical Properties

Property code	Value	Unit	Source
log10ws	0.40		Crippen Method
logp	4.411		Crippen Method
rinpol	1405.00		NIST Webbook

Sources

Crippen Method: <http://pubs.acs.org/doi/abs/10.1021/ci990307l>
Crippen Method: https://www.cheméo.com/doc/models/crippen_log10ws
NIST Webbook: <http://webbook.nist.gov/cgi/cbook.cgi?ID=R168008&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.cheméo.com/cid/62-719-5/Pentanoic-acid-4-hydroxy-2-propyl-TMS-2.pdf>

Generated by Cheméo on 2024-05-11 10:06:10.179130772 +0000 UTC m=+17711219.099708093.

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