

5-Methyluridine, tris(trimethylsilyl) deriv.

Inchi: InChI=1S/C19H38N2O6Si3/c1-13-11-21(19(23)20-17(13)22)18-16(27-30(8,9)10)15(26-2
InchiKey: NPQPNHPRBSPWIM-UHFFFAOYSA-N
Formula: C19H38N2O6Si3
SMILES: Cc1cn(C2OC(CO[Si](C)(C)C)C(O[Si](C)(C)C)C2O[Si](C)(C)C)c(=O)[nH]c1=O
Mol. weight [g/mol]: 474.77

Physical Properties

Property code	Value	Unit	Source
log10ws	3.51		Crippen Method
logp	2.552		Crippen Method
rinpol	2478.00		NIST Webbook
rinpol	2478.00		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=U375674&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/38-199-0/5-Methyluridine-tris-trimethylsilyl-deriv.pdf>

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