

# Cytidine, etho-oxime-TMS derivative

**Inchi:** InChI=1S/C22H48N4O5Si4/c1-14-27-24-22-23-17(25-32(2,3)4)15-16-26(22)20-18(29-33)  
**InchiKey:** RNTKEVOGXFBPRB-TZWRJOPASA-N  
**Formula:** C22H48N4O5Si4  
**SMILES:** CCON=c1nc(N[Si](C)(C)C)ccn1C1OC(O[Si](C)(C)C)C(O[Si](C)(C)C)C1O[Si](C)(C)C  
**Mol. weight [g/mol]:** 560.98

## Physical Properties

Property code	Value	Unit	Source
log10ws	3.16		Crippen Method
logp	5.129		Crippen Method
rinpol	2545.00		NIST Webbook
rinpol	2545.00		NIST Webbook

## Sources

**Crippen Method:** <http://pubs.acs.org/doi/abs/10.1021/ci9903071>  
**Crippen Method:** [https://www.chemeo.com/doc/models/crippen\\_log10ws](https://www.chemeo.com/doc/models/crippen_log10ws)  
**NIST Webbook:** <http://webbook.nist.gov/cgi/cbook.cgi?ID=R245744&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/45-569-1/Cytidine-etho-oxime-TMS-derivative.pdf>

Generated by Cheméo on 2024-04-19 18:15:16.76942608 +0000 UTC m=+15839765.690003392.

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