

# Purine, 6-amino-2-hydroxy, TMS

**Inchi:** InChI=1S/C11H21N5OSi2/c1-18(2,3)16-10-8-9(13-7-12-8)14-11(15-10)17-19(4,5)6/h7H,  
**InchiKey:** ZFXSUEONDULERV-UHFFFAOYSA-N  
**Formula:** C11H21N5OSi2  
**SMILES:** C[Si](C)(C)Nc1nc(O[Si](C)(C)C)nc2nc[nH]c12  
**Mol. weight [g/mol]:** 295.49

## Physical Properties

Property code	Value	Unit	Source
log10ws	0.30		Crippen Method
logp	2.331		Crippen Method
rinpol	2068.00		NIST Webbook
rinpol	2068.00		NIST Webbook

## Sources

**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=R386548&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/119-301-5/Purine-6-amino-2-hydroxy-TMS.pdf>

Generated by Cheméo on 2024-04-29 04:44:08.596718548 +0000 UTC m=+16655097.517295869.

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