

# Acetylglycine, bis-TMS

<b>Other names:</b>	Acetic acid, acetylamino, bis-TMS N-ACETYLGLYCINE TMS
<b>Inchi:</b>	InChI=1S/C10H23NO3Si2/c1-9(13-15(2,3)4)11-8-10(12)14-16(5,6)7/h8H2,1-7H3/b11-9+
<b>InchiKey:</b>	MEERPCCZBNRYNP-PKQBQFBNSA-N
<b>Formula:</b>	C10H23NO3Si2
<b>SMILES:</b>	CC(=NCC(=O)O[Si](C)(C)C)O[Si](C)(C)C
<b>Mol. weight [g/mol]:</b>	261.47

## Physical Properties

Property code	Value	Unit	Source
log10ws	2.28		Crippen Method
logp	2.634		Crippen Method
rinpol	1359.00		NIST Webbook
rinpol	1389.00		NIST Webbook
rinpol	1372.00		NIST Webbook

## Sources

<b>Crippen Method:</b>	<a href="http://pubs.acs.org/doi/abs/10.1021/ci9903071">http://pubs.acs.org/doi/abs/10.1021/ci9903071</a>
<b>Crippen Method:</b>	<a href="https://www.chemeo.com/doc/models/crippen_log10ws">https://www.chemeo.com/doc/models/crippen_log10ws</a>
<b>NIST Webbook:</b>	<a href="http://webbook.nist.gov/cgi/cbook.cgi?ID=R117189&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=R117189&amp;Units=SI</a>

## Legend

<b>log10ws:</b>	Log10 of Water solubility in mol/l
<b>logp:</b>	Octanol/Water partition coefficient
<b>rinpol:</b>	Non-polar retention indices

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