

# 2-(P-trimethylsilyloxy)-3-methyl pyrazine

<b>Inchi:</b>	InChI=1S/C14H18N2OSi/c1-11-14(16-10-9-15-11)17-12-5-7-13(8-6-12)18(2,3)4/h5-10H,
<b>InchiKey:</b>	OWTOIHCPFZCFMS-UHFFFAOYSA-N
<b>Formula:</b>	C14H18N2OSi
<b>SMILES:</b>	Cc1nccnc1Oc1ccc([Si](C)(C)C)cc1
<b>Mol. weight [g/mol]:</b>	258.39
<b>CAS:</b>	116295-61-5

## Physical Properties

Property code	Value	Unit	Source
log10ws	-5.78		Crippen Method
logp	3.123		Crippen Method

## 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.cheméo.com/doc/models/crippen_log10ws">https://www.cheméo.com/doc/models/crippen_log10ws</a>
<b>NIST Webbook:</b>	<a href="http://webbook.nist.gov/cgi/cbook.cgi?ID=C116295615&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=C116295615&amp;Units=SI</a>

## Legend

<b>log10ws:</b>	Log10 of Water solubility in mol/l
<b>logp:</b>	Octanol/Water partition coefficient

Latest version available from:

<https://www.cheméo.com/cid/120-904-4/2-P-trimethylsilyloxy-3-methyl-pyrazine.pdf>

Generated by Cheméo on 2024-04-30 22:12:48.30563092 +0000 UTC m=+16804417.226208236.

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