

# Pentazocine TMS derivative

<b>Other names:</b>	Pentazocine O-TMS
<b>Inchi:</b>	InChI=1S/C22H35NOSi/c1-16(2)10-12-23-13-11-22(4)17(3)21(23)14-18-8-9-19(15-20(18
<b>InchiKey:</b>	KMKNLGHEFMHPCX-UHFFFAOYSA-N
<b>Formula:</b>	C22H35NOSi
<b>SMILES:</b>	CC(C)=CCN1CCC2(C)c3cc(O[Si](C)(C)C)ccc3CC1C2C
<b>Mol. weight [g/mol]:</b>	357.60

## Physical Properties

Property code	Value	Unit	Source
log10ws	-3.61		Crippen Method
logp	5.391		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=U137224&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=U137224&amp;Units=SI</a>

## Legend

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

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