

# 5-Amino-2-methoxyphenol, trimethylsilyl ether

<b>Other names:</b>	5-Amino-2-methoxyphenol, tms derivative
<b>Inchi:</b>	InChI=1S/C10H17NO2Si/c1-12-9-6-5-8(11)7-10(9)13-14(2,3)4/h5-7H,11H2,1-4H3
<b>InchiKey:</b>	DLAVHTHUUJYFMFF-UHFFFAOYSA-N
<b>Formula:</b>	C10H17NO2Si
<b>SMILES:</b>	COc1ccc(N)cc1O[Si](C)(C)C
<b>Mol. weight [g/mol]:</b>	211.33

## Physical Properties

Property code	Value	Unit	Source
log10ws	-0.23		Crippen Method
logp	2.491		Crippen Method
rinpol	1509.60		NIST Webbook
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## Sources

<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=U352888&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=U352888&amp;Units=SI</a>
<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>

## 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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<https://www.chemeo.com/cid/121-808-0/5-Amino-2-methoxyphenol-trimethylsilyl-ether.pdf>

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