

# 4-Methylcatechol, bis(trimethylsilyl) ether

<b>Other names:</b>	[(4-Methyl-1,2-phenylene)bis(oxy)]bis(trimethylsilane) 4-Methyl-1,2-dihydroxybenzene, TMS 4-Methyl-1,2-dihydroxybenzene,bisTMS ether 1,2-Benzenediol, 4-methyl, bis-TMS 4-Methylcatechol, 2tms derivative
<b>Inchi:</b>	InChI=1S/C13H24O2Si2/c1-11-8-9-12(14-16(2,3)4)13(10-11)15-17(5,6)7/h8-10H,1-7H3
<b>InchiKey:</b>	OZIRHEHKTRSGJX-UHFFFAOYSA-N
<b>Formula:</b>	C13H24O2Si2
<b>SMILES:</b>	Cc1ccc(O[Si](C)(C)C)c(O[Si](C)(C)C)c1
<b>Mol. weight [g/mol]:</b>	268.50

## Physical Properties

Property code	Value	Unit	Source
log10ws	0.04		Crippen Method
logp	4.422		Crippen Method
rinpol	1388.00		NIST Webbook
rinpol	1382.30		NIST Webbook

## Sources

<b>NIST Webbook:</b>	<a href="http://webbook.nist.gov/cgi/cbook.cgi?ID=U332799&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=U332799&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>
<b>Crippen Method:</b>	<a href="https://www.chemeo.com/doc/models/crippen_log10ws">https://www.chemeo.com/doc/models/crippen_log10ws</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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