

4-Methoxyphenol, tert-butyldimethylsilyl ether

Other names:	4-Methoxyphenol, tbdms derivative
Inchi:	InChI=1S/C13H22O2Si/c1-13(2,3)16(5,6)15-12-9-7-11(14-4)8-10-12/h7-10H,1-6H3
InchiKey:	HDPHSOHXTXSDJI-UHFFFAOYSA-N
Formula:	C13H22O2Si
SMILES:	COc1ccc(O[Si](C)(C)C(C)(C)C)cc1
Mol. weight [g/mol]:	238.40

Physical Properties

Property code	Value	Unit	Source
log10ws	-1.85		Crippen Method
logp	4.079		Crippen Method
rinpol	1533.00		NIST Webbook
rinpol	1508.00		NIST Webbook
rinpol	1533.00		NIST Webbook

Sources

NIST Webbook:	http://webbook.nist.gov/cgi/cbook.cgi?ID=U352896&Units=SI
Crippen Method:	http://pubs.acs.org/doi/abs/10.1021/ci990307l
Crippen Method:	https://www.chemeo.com/doc/models/crippen_log10ws

Legend

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

Latest version available from:

<https://www.chemeo.com/cid/124-959-0/4-Methoxyphenol-tert-butyldimethylsilyl-ether.pdf>

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