

Hydroquinone, bis(tert-butyldimethylsilyl) ether

Other names:	Hydroquinone, 2tbdms derivative
Inchi:	InChI=1S/C18H34O2Si2/c1-17(2,3)21(7,8)19-15-11-13-16(14-12-15)20-22(9,10)18(4,5)6
InchiKey:	FMHSADBDISJXNQ-UHFFFAOYSA-N
Formula:	C18H34O2Si2
SMILES:	CC(C)(C)[Si](C)(C)Oc1ccc(O[Si](C)(C)C(C)(C)C)cc1
Mol. weight [g/mol]:	338.63

Physical Properties

Property code	Value	Unit	Source
log10ws	-2.00		Crippen Method
logp	6.455		Crippen Method
rinpol	1864.70		NIST Webbook
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Sources

Crippen Method:	https://www.chemeo.com/doc/models/crippen_log10ws
NIST Webbook:	http://webbook.nist.gov/cgi/cbook.cgi?ID=U352775&Units=SI
Crippen Method:	http://pubs.acs.org/doi/abs/10.1021/ci990307l

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/115-078-8/Hydroquinone-bis-tert-butyldimethylsilyl-ether.pdf>

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