

# Dihydrochalcone, 2',6',4-trihydroxy-4'-methoxy, tris-TMS

Other names:	Dihydrochalcone, 2',4,6'-trihydroxy-4'-methoxy, TMS
Inchi:	InChI=1S/C25H40O5Si3/c1-27-21-17-24(29-32(5,6)7)22(25(18-21)30-33(8,9)10)15-16-23
InchiKey:	SBHSCBLSVDDBS-UHFFFAOYSA-N
Formula:	C25H40O5Si3
SMILES:	COc1cc(O[Si](C)(C)C)c(CCC(=O)c2ccc(O[Si](C)(C)C)cc2)c(O[Si](C)(C)C)c1
Mol. weight [g/mol]:	504.84

## Physical Properties

Property code	Value	Unit	Source
log10ws	-1.31		Crippen Method
logp	7.152		Crippen Method
rinpol	2700.00		NIST Webbook
rinpol	2705.00		NIST Webbook
rinpol	2712.00		NIST Webbook
rinpol	2712.00		NIST Webbook

## Sources

Crippen Method:	<a href="http://pubs.acs.org/doi/abs/10.1021/ci9903071">http://pubs.acs.org/doi/abs/10.1021/ci9903071</a>
Crippen Method:	<a href="https://www.chemeo.com/doc/models/crippen_log10ws">https://www.chemeo.com/doc/models/crippen_log10ws</a>
NIST Webbook:	<a href="http://webbook.nist.gov/cgi/cbook.cgi?ID=R46324&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=R46324&amp;Units=SI</a>

## 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/46-639-2/Dihydrochalcone-2-6-4-trihydroxy-4-methoxy-tris-TMS.pdf>

Generated by Cheméo on 2024-04-17 14:15:16.778192793 +0000 UTC m=+15652565.698770109.

Cheméo (<https://www.chemeo.com>) is the biggest free database of chemical and physical data for the process industry.