

# MELIBIOSE 8TMS-1

<b>Inchi:</b>	InChI=1S/C36H86O11Si8/c1-48(2,3)38-26-28-30(42-50(7,8)9)31(43-51(10,11)12)33(45-5
<b>InchiKey:</b>	SWPQHIFGBMZXPV-GQKLMCKXSA-N
<b>Formula:</b>	C36H86O11Si8
<b>SMILES:</b>	C[Si](C)(C)OCC1OC(OCC2OC(O[Si](C)(C)C)C(O[Si](C)(C)C)C(O[Si](C)(C)C)C2O[Si](C)
<b>Mol. weight [g/mol]:</b>	919.75

## Physical Properties

Property code	Value	Unit	Source
log10ws	8.84		Crippen Method
logp	9.476		Crippen Method
rmpol	2850.00		NIST Webbook
rmpol	2878.00		NIST Webbook
rmpol	2850.00		NIST Webbook

## 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=R439226&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=R439226&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>rmpol:</b>	Non-polar retention indices

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