# [(2-Ethyl-5-methylfuran-3,4-diyl)bis(oxy)]bis(trime

**Inchi:** InChI=1S/C13H26O3Si2/c1-9-11-13(16-18(6,7)8)12(10(2)14-11)15-17(3,4)5/h9H2,1-8H3
**InchiKey:** KKARWYSMFMFIGP-UHFFFAOYSA-N
**Formula:** C13H26O3Si2
**SMILES:** CCc1oc(C)c(O[Si](C)(C)C)c1O[Si](C)(C)C
**Mol. weight [g/mol]:** 286.51

## Physical Properties

| Property code | Value | Unit | Source |
|---|---|---|---|
| log10ws | -4.50 | | Crippen Method |
| logp | 4.578 | | Crippen Method |
| rinpol | 1349.90 | | NIST Webbook |

## Sources

**Crippen Method:** https://www.chemeo.com/doc/models/crippen_log10ws
**NIST Webbook:** http://webbook.nist.gov/cgi/cbook.cgi?ID=U333831&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