# «beta»-Hydroxyphenethylamine, N-isoBOC, O-TBDMS

**Inchi:** InChI=1S/C19H33NO3Si/c1-15(2)14-22-18(21)20-13-17(16-11-9-8-10-12-16)23-24(6,7)1

**InchiKey:** SUDNYVLMMNDTSR-UHFFFAOYSA-N

**Formula:** C19H33NO3Si

**SMILES:** CC(C)COC(=O)NCC(O[Si](C)(C)C(C)(C)C)c1ccccc1

**Mol. weight [g/mol]:** 351.56

## Physical Properties

| Property code | Value | Unit | Source |
|---|---|---|---|
| log10ws | -3.26 | | Crippen Method |
| logp | 5.132 | | Crippen Method |
| rinpol | 2095.00 | | NIST Webbook |

## Sources

**Crippen Method:** http://pubs.acs.org/doi/abs/10.1021/ci990307l

**Crippen Method:** https://www.chemeo.com/doc/models/crippen_log10ws

**NIST Webbook:** http://webbook.nist.gov/cgi/cbook.cgi?ID=R392320&Units=SI

## Legend

**log10ws:** Log10 of Water solubility in mol/l

**logp:** Octanol/Water partition coefficient

**rinpol:** Non-polar retention indices