

# 2-(N-Ethyl-N-m-tolyl)aminoethanol, benzyldimethylsilyl ether

**Inchi:** InChI=1S/C20H29NOSi/c1-5-21(20-13-9-10-18(2)16-20)14-15-22-23(3,4)17-19-11-7-6-8  
**InchiKey:** WEFRWTFEDATRKY-UHFFFAOYSA-N  
**Formula:** C20H29NOSi  
**SMILES:** CCN(CCO[Si](C)(C)Cc1ccccc1)c1cccc(C)c1  
**Mol. weight [g/mol]:** 327.54

## Physical Properties

Property code	Value	Unit	Source
log10ws	-2.79		Crippen Method
logp	4.825		Crippen Method
rinpol	2286.00		NIST Webbook

## Sources

**NIST Webbook:** <http://webbook.nist.gov/cgi/cbook.cgi?ID=U375541&Units=SI>  
**Crippen Method:** <http://pubs.acs.org/doi/abs/10.1021/ci9903071>  
**Crippen Method:** [https://www.chemeo.com/doc/models/crippen\\_log10ws](https://www.chemeo.com/doc/models/crippen_log10ws)

## Legend

**log10ws:** Log10 of Water solubility in mol/l  
**logp:** Octanol/Water partition coefficient  
**rinpol:** Non-polar retention indices

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