2-(M-phenoxyphenoxy)-3-(2-trimethylsilylethyl) pyrazine

Inchi: InChI=1S/C21H24N2O2Si/c1-26(2,3)15-12-20-21(23-14-13-22-20)25-19-11-7-10-18(16-1

InchiKey: XZZSMZQDSFLGGA-UHFFFAOYSA-N

Formula: C21H24N2O2Si

SMILES: C[Si](C)(C)CCc1nccnc1Oc1cccc(Oc2ccccc2)c1

Mol. weight [g/mol]: 364.51

Physical Properties

Property code	Value	Unit	Source
log10ws	-3.80		Crippen Method
logp	5.942		Crippen Method

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=B6002916&Units=SI

Legend

log10ws: Log10 of Water solubility in mol/l

logp: Octanol/Water partition coefficient