

3,4-DIHYDROXYPHENYLETHYLAMINE, TRIS(TRIMETHYLSILYL)- DERIV.

Other names: 3,4-Dihydroxyphenylethylamine, tris(trimethylsilyl)- deriv.

Dopamine, N,O,O-tris-TMS

Dopamine, TMS

Dopamine, 3tms derivative

Inchi: InChI=1S/C17H35NO2Si3/c1-21(2,3)18-13-12-15-10-11-16(19-22(4,5)6)17(14-15)20-23(

InchiKey: IDQRGLQSFWNZMQ-UHFFFAOYSA-N

Formula: C17H35NO2Si3

SMILES: C[Si](C)(C)NCCc1ccc(O[Si](C)(C)C)c(O[Si](C)(C)C)c1

Mol. weight [g/mol]: 369.73

Physical Properties

Property code	Value	Unit	Source
logp	5.081		Crippen Method
rinpol	2104.00		NIST Webbook
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Sources

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

Cheméo Relay (1.0):

Cheméo Relay (1.0, beta):

NIST Webbook: <http://webbook.nist.gov/cgi/cbook.cgi?ID=C68595540&Units=SI>

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

Legend

logp: Octanol/Water partition coefficient

rinpol: Non-polar retention indices

Latest version available from:

<https://www.chemeo.com/cid/122-833-1/3-4-DIHYDROXYPHENYLETHYLAMINE-TRIS-TRIMETHYLSILYL-DERIV.p>

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