2,2'-(TRIMETHYLENEDIOXY)BIS(1,3,2-DIOXABOR

Other names: 1,3,2-Dioxaborinane, 2,2'-[1,3-propanediylbis(oxy)]bis-
2,2'-[propane-1,3-diylbis(oxy)]bis-1,3,2-dioxaborinane
Inchi: InChI=1S/C9H18B2O6/c1-4-12-10(13-5-1)16-8-3-9-17-11-14-6-2-7-15-11/h1-9H2
InchiKey: FZNHHPNJQSIEPT-UHFFFAOYSA-N
Formula: C9H18B2O6
SMILES: C1COB(OCCCOB2OCCCO2)OC1
Mol. weight [g/mol]: 243.86

Physical Properties

Property code	Value	Unit	Source
logp	0.253		Crippen Method

Sources

Cheméo Relay (1.0):
Cheméo Relay (1.0, beta):
NIST Webbook: http://webbook.nist.gov/cgi/cbook.cgi?ID=C20905355&Units=SI
Crippen Method: http://pubs.acs.org/doi/abs/10.1021/ci990307l
Crippen Method: https://www.chemeo.com/doc/models/crippen_log10ws

Legend

logp: Octanol/Water partition coefficient