

2,2,4-Trimethylpentane-1,3-diol, ferroceneboronate derivative

Inchi: InChI=1S/C13H20BO2.C5H5.Fe/c1-10(2)12-13(3,4)9-15-14(16-12)11-7-5-6-8-11;1-2-4-5
InchiKey: SCTVLBUMJXDELG-UHFFFAOYSA-N
Formula: C18H25BFeO2
SMILES: CC(C)C1OB(C23C4C5C6C2[Fe]56432789C3C2C7C8C39)OCC1(C)C
Mol. weight [g/mol]: 340.05

Physical Properties

Property code	Value	Unit	Source
rinpol	2150.00		NIST Webbook

Sources

NIST Webbook: <http://webbook.nist.gov/cgi/cbook.cgi?ID=R250909&Units=SI>

Legend

rinpol: Non-polar retention indices

Latest version available from:

<https://www.cheméo.com/cid/12-959-4/2-2-4-Trimethylpentane-1-3-diol-ferroceneboronate-derivative.pdf>

Generated by Cheméo on 2024-05-02 02:29:55.17042655 +0000 UTC m=+16906244.091003862.

Cheméo (<https://www.cheméo.com>) is the biggest free database of chemical and physical data for the process industry.