(-)-2-Aminobutan-1-ol, ferroceneboronate derivative

Inchi: InChI=1S/C9H13BNO.C5H5.Fe/c1-2-9-7-12-10(11-9)8-5-3-4-6-8;1-2-4-5-3-1;/h3-6,9,11H

InchiKey: HUYCPQXEFXHFRS-UHFFFAOYSA-N

Formula: C14H18BFeNO

SMILES: CCC1COB(C23C4C5C6C2[Fe]56432789C3C2C7C8C39)N1

Mol. weight [g/mol]: 282.95

Physical Properties

Property code	Value	Unit	Source
rinpol	1975.00		NIST Webbook
rinpol	1975.00		NIST Webbook

Sources

NIST Webbook: http://webbook.nist.gov/cgi/cbook.cgi?ID=R250890&Units=SI

Legend

rinpol: Non-polar retention indices