

1,2-Propanediol, ferrocenylboronate

Inchi: InChI=1S/C8H10BO2.C5H5.Fe/c1-7-6-10-9(11-7)8-4-2-3-5-8;1-2-4-5-3-1;/h2-5,7H,6H2,1
InchiKey: MAHJAEGTFDKVN-X-UHFFFAOYSA-N
Formula: C13H15BFeO2
SMILES: CC1COB(C23C4C5C6C2[Fe]56432789C3C2C7C8C39)O1
Mol. weight [g/mol]: 269.91

Physical Properties

Property code	Value	Unit	Source
rinp	1770.00		NIST Webbook

Sources

NIST Webbook: <http://webbook.nist.gov/cgi/cbook.cgi?ID=R163083&Units=SI>

Legend

rinp: Non-polar retention indices

Latest version available from:

<https://www.cheméo.com/cid/14-171-6/1-2-Propanediol-ferrocenylboronate.pdf>

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