Zink dipentyldithiocarbamate

Inchi: InChI=1S/C12H25NS2.C11H23NS2.Zn/c1-3-5-7-9-13(11-12(14)15)10-8-6-4-2;1-3-5-7-9-

InchiKey: ATDAFKHZUZYDHX-UHFFFAOYSA-L

Formula: C23H46N2S4Zn

SMILES: CCCCCN(CCCCC)C(=S)[S-].CCCCCN(CCCCC)CC(=S)[S-].[Zn]

Mol. weight [g/mol]: 544.26

Physical Properties

Property code	Value	Unit	Source
rinpol	3706.00		NIST Webbook

Sources

NIST Webbook: http://webbook.nist.gov/cgi/cbook.cgi?ID=R104519&Units=SI

Legend

rinpol: Non-polar retention indices