Zink diethyldithiocarbamate

Inchi: InChI=1S/C6H13NS2.C5H11NS2.Zn/c1-3-7(4-2)5-6(8)9;1-3-6(4-2)5(7)8;/h3-5H2,1-2H3,(

InchiKey: HFANNQLVJJJDGL-UHFFFAOYSA-L

Formula: C11H22N2S4Zn

SMILES: CCN(CC)CC(=S)S[Zn]SC(=S)N(CC)CC

Mol. weight [g/mol]: 375.95

Physical Properties

Property code	Value	Unit	Source
rinpol	2728.00		NIST Webbook
rinpol	2728.00		NIST Webbook

Sources

NIST Webbook: http://webbook.nist.gov/cgi/cbook.cgi?ID=R104475&Units=SI

Legend

rinpol: Non-polar retention indices