

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
rropol	2728.00		NIST Webbook
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Sources

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

Legend

rropol: Non-polar retention indices

Latest version available from:

<https://www.cheméo.com/cid/124-811-3/Zink-diethyldithiocarbamate.pdf>

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