## **Carbamic acid, ethyldithio-, compound with ethylamine** Inchi: InChi=1S/C3H7NS2.C2H7N/c1-2-4-3(5)6;1-2-3/h2H2,1H3,(H2,4,5,6);2-3H2

InchiKey: Formula: SMILES: Mol. weight [g/mol]: InChI=1S/C3H7NS2.C2H7N/c1-2-4-3(5)6;1-2-3/h2H2,1H3,(H2,4,5,6);2-3H2,1H3 XSBNNDUUIRIBKE-UHFFFAOYSA-N C5H14N2S2 CCNC(=S)[S-].CC[NH3+] 166.31

## Sources

NIST Webbook:

http://webbook.nist.gov/cgi/cbook.cgi?ID=B6009211&Units=SI

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