

Acetyl-o-phenetidine

Inchi:	InChI=1S/C8H11NO.C3H6O/c1-2-10-8-6-4-3-5-7(8)9;1-3(2)4/h3-6H,2,9H2,1H3;1-2H3
InchiKey:	WRNMZLSLAABCMS-UHFFFAOYSA-N
Formula:	C11H17NO2
SMILES:	CC(C)=O.CCOc1ccccc1N
Mol. weight [g/mol]:	195.26

Sources

NIST Webbook: <http://webbook.nist.gov/cgi/cbook.cgi?ID=B6001492&Units=SI>

Latest version available from:

<https://www.cheméo.com/cid/11-314-0/Acetyl-o-phenetidine.pdf>

Generated by Cheméo on 2024-04-25 22:17:20.20160072 +0000 UTC m=+16372689.122178042.

Cheméo (<https://www.cheméo.com>) is the biggest free database of chemical and physical data for the process industry.