

Acetamide, N-tetrahydrofurfuryl-2-(2-thienyl)-

Inchi:	InChI=1S/C11H15NO2S/c13-11(7-10-4-2-6-15-10)12-8-9-3-1-5-14-9/h2,4,6,9H,1,3,5,7-8
InchiKey:	FBBGQQWNZLUCCK-UHFFFAOYSA-N
Formula:	C11H15NO2S
SMILES:	O=C(Cc1cccs1)NCC1CCCO1
Mol. weight [g/mol]:	225.31

Physical Properties

Property code	Value	Unit	Source
log10ws	-2.18		Crippen Method
logp	1.586		Crippen Method
mcvol	169.300	ml/mol	McGowan Method
rinpol	1872.00		NIST Webbook
rinpol	1872.00		NIST Webbook

Sources

Crippen Method:	https://www.chemeo.com/doc/models/crippen_log10ws
McGowan Method:	http://link.springer.com/article/10.1007/BF02311772
NIST Webbook:	http://webbook.nist.gov/cgi/cbook.cgi?ID=U306900&Units=SI
Crippen Method:	http://pubs.acs.org/doi/abs/10.1021/ci990307l

Legend

log10ws:	Log10 of Water solubility in mol/l
logp:	Octanol/Water partition coefficient
mcvol:	McGowan's characteristic volume
rinpol:	Non-polar retention indices

Latest version available from:

<https://www.chemeo.com/cid/124-383-9/Acetamide-N-tetrahydrofurfuryl-2-2-thienyl.pdf>

Generated by Cheméo on 2024-04-29 05:54:35.911032815 +0000 UTC m=+16659324.831610126.

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