

# 5-ethyl-4-methyl-2-pentyl-thiazole

<b>Inchi:</b>	InChI=1S/C11H19NS/c1-4-6-7-8-11-12-9(3)10(5-2)13-11/h4-8H2,1-3H3
<b>InchiKey:</b>	YHXYVNQEOGPEKA-UHFFFAOYSA-N
<b>Formula:</b>	C11H19NS
<b>SMILES:</b>	CCCCC1nc(C)c(CC)s1
<b>Mol. weight [g/mol]:</b>	197.34

## Physical Properties

Property code	Value	Unit	Source
log10ws	-4.33		Crippen Method
logp	3.747		Crippen Method
mcvol	172.720	ml/mol	McGowan Method
rinsol	1454.00		NIST Webbook

## Sources

<b>Crippen Method:</b>	<a href="http://pubs.acs.org/doi/abs/10.1021/ci9903071">http://pubs.acs.org/doi/abs/10.1021/ci9903071</a>
<b>Crippen Method:</b>	<a href="https://www.chemeo.com/doc/models/crippen_log10ws">https://www.chemeo.com/doc/models/crippen_log10ws</a>
<b>McGowan Method:</b>	<a href="http://link.springer.com/article/10.1007/BF02311772">http://link.springer.com/article/10.1007/BF02311772</a>
<b>NIST Webbook:</b>	<a href="http://webbook.nist.gov/cgi/cbook.cgi?ID=R498369&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=R498369&amp;Units=SI</a>

## Legend

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

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