

# Pentanamide, N-(3-methylphenyl)-

<b>Inchi:</b>	InChI=1S/C12H17NO/c1-3-4-8-12(14)13-11-7-5-6-10(2)9-11/h5-7,9H,3-4,8H2,1-2H3,(H,1)
<b>InchiKey:</b>	RTVXANSLXIIIKG-UHFFFAOYSA-N
<b>Formula:</b>	C12H17NO
<b>SMILES:</b>	CCCCC(O)=Nc1cccc(C)c1
<b>Mol. weight [g/mol]:</b>	191.27

## Physical Properties

Property code	Value	Unit	Source
hf	-145.75	kJ/mol	Joback Method
hvap	65.32	kJ/mol	Joback Method
log10ws	-3.59		Crippen Method
logp	3.773		Crippen Method
mcvol	167.730	ml/mol	McGowan Method
pc	2351.92	kPa	Joback Method
rinpol	1714.00		NIST Webbook
tb	674.36	K	Joback Method
tc	880.11	K	Joback Method

## 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>Joback Method:</b>	<a href="https://en.wikipedia.org/wiki/Joback_method">https://en.wikipedia.org/wiki/Joback_method</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=U306891&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=U306891&amp;Units=SI</a>

## Legend

<b>hf:</b>	Enthalpy of formation at standard conditions
<b>hvap:</b>	Enthalpy of vaporization at standard conditions
<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>pc:</b>	Critical Pressure
<b>rinpol:</b>	Non-polar retention indices
<b>tb:</b>	Normal Boiling Point Temperature
<b>tc:</b>	Critical Temperature

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