

# N,N-Dimethyl-N'-benzyl-propionamide

<b>Inchi:</b>	InChI=1S/C12H18N2/c1-4-12(14(2)3)13-10-11-8-6-5-7-9-11/h5-9H,4,10H2,1-3H3/b13-12
<b>InchiKey:</b>	GTOOPRWBIUWQHJ-OUKQBFOZSA-N
<b>Formula:</b>	C12H18N2
<b>SMILES:</b>	CCC(=NCc1ccccc1)N(C)C
<b>Mol. weight [g/mol]:</b>	190.28

## Physical Properties

Property code	Value	Unit	Source
hf	85.48	kJ/mol	Joback Method
hvap	50.02	kJ/mol	Joback Method
log10ws	-2.67		Crippen Method
logp	2.557		Crippen Method
mcvol	171.840	ml/mol	McGowan Method
pc	2181.56	kPa	Joback Method
rinpol	1617.00		NIST Webbook
tb	589.64	K	Joback Method
tc	806.67	K	Joback Method

## Sources

<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=R162000&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=R162000&amp;Units=SI</a>
<b>Crippen Method:</b>	<a href="http://pubs.acs.org/doi/abs/10.1021/ci990307I">http://pubs.acs.org/doi/abs/10.1021/ci990307I</a>
<b>Crippen Method:</b>	<a href="https://www.chemeo.com/doc/models/crippen_log10ws">https://www.chemeo.com/doc/models/crippen_log10ws</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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