

Benzamide, 3-methyl-N-octyl-

Inchi:	InChI=1S/C16H25NO/c1-3-4-5-6-7-8-12-17-16(18)15-11-9-10-14(2)13-15/h9-11,13H,3-8
InchiKey:	PDNXPELOTIFSKU-UHFFFAOYSA-N
Formula:	C16H25NO
SMILES:	CCCCCCCCN=C(O)c1cccc(C)c1
Mol. weight [g/mol]:	247.38

Physical Properties

Property code	Value	Unit	Source
hf	-228.31	kJ/mol	Joback Method
hvap	74.22	kJ/mol	Joback Method
log10ws	-4.77		Crippen Method
logp	4.660		Crippen Method
mcvol	224.090	ml/mol	McGowan Method
pc	1671.43	kPa	Joback Method
rinpol	2187.00		NIST Webbook
rinpol	2187.00		NIST Webbook
tb	765.88	K	Joback Method
tc	963.96	K	Joback Method

Sources

Crippen Method:	http://pubs.acs.org/doi/abs/10.1021/ci9903071
Crippen Method:	https://www.chemeo.com/doc/models/crippen_log10ws
Joback Method:	https://en.wikipedia.org/wiki/Joback_method
McGowan Method:	http://link.springer.com/article/10.1007/BF02311772
NIST Webbook:	http://webbook.nist.gov/cgi/cbook.cgi?ID=U407418&Units=SI

Legend

hf:	Enthalpy of formation at standard conditions
hvap:	Enthalpy of vaporization at standard conditions

log10ws:	Log10 of Water solubility in mol/l
logp:	Octanol/Water partition coefficient
mcvol:	McGowan's characteristic volume
pc:	Critical Pressure
rinpol:	Non-polar retention indices
tb:	Normal Boiling Point Temperature
tc:	Critical Temperature

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