N,N-Dimethyl-N'-phenyl-benzamidine

Inchi: InChI=1S/C15H16N2/c1-17(2)15(13-9-5-3-6-10-13)16-14-11-7-4-8-12-14/h3-12H,1-2H3
InchiKey: YRCVSLXBPYHZDA-UHFFFAOYSA-N
Formula: C15H16N2
SMILES: CN(C)C(=Nc1ccccc1)c1ccccc1
Mol. weight [g/mol]: 224.30

Physical Properties

Property code	Value	Unit	Source
hf	260.09	kJ/mol	Joback Method
hvap	58.97	kJ/mol	Joback Method
log10ws	-3.24		Crippen Method
logp	3.327		Crippen Method
mcvol	190.350	ml/mol	McGowan Method
pc	2271.90	kPa	Joback Method
rinpol	1842.00		NIST Webbook
rinpol	1842.00		NIST Webbook
tb	684.96	K	Joback Method
tc	934.54	K	Joback Method

Sources

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=R159410&Units=SI
Crippen Method: http://pubs.acs.org/doi/abs/10.1021/ci990307l
Crippen Method: https://www.chemeo.com/doc/models/crippen_log10ws

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