N,N-Dimethyl-N'-(3-bromophenyl)-propionamidine

Inchi: InChI=1S/C11H15BrN2/c1-4-11(14(2)3)13-10-7-5-6-9(12)8-10/h5-8H,4H2,1-3H3
InchiKey: IKPWJLPMELAVPF-UHFFFAOYSA-N
Formula: C11H15BrN2
SMILES: CCC(=Nc1cccc(Br)c1)N(C)C
Mol. weight [g/mol]: 255.15

Physical Properties

Property code	Value	Unit	Source
hf	120.98	kJ/mol	Joback Method
hvap	54.89	kJ/mol	Joback Method
log10ws	-3.52		Crippen Method
logp	3.451		Crippen Method
mcvol	175.250	ml/mol	McGowan Method
pc	2522.65	kPa	Joback Method
rinpol	1807.00		NIST Webbook
tb	637.90	K	Joback Method
tc	872.34	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=R161889&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