N,N-Dimethyl-N'-nonyl-p-methylbenzamidine

Inchi: InChI=1S/C19H32N2/c1-5-6-7-8-9-10-11-16-20-19(21(3)4)18-14-12-17(2)13-15-18/h12-1
InchiKey: UTOSXRQQIVCCGQ-UHFFFAOYSA-N
Formula: C19H32N2
SMILES: CCCCCCCCCN=C(c1ccc(C)cc1)N(C)C
Mol. weight [g/mol]: 288.47

Physical Properties

Property code	Value	Unit	Source
hf	-70.47	kJ/mol	Joback Method
hvap	66.26	kJ/mol	Joback Method
log10ws	-5.27		Crippen Method
logp	5.054		Crippen Method
mcvol	270.470	ml/mol	McGowan Method
pc	1250.38	kPa	Joback Method
rinpol	2063.00		NIST Webbook
rinpol	2063.00		NIST Webbook
tb	754.78	K	Joback Method
tc	953.73	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=R159321&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