N,N-Dimethyl-2-phenyl-N'-(3-methylphenyl)-acetar

Inchi: InChI=1S/C17H20N2/c1-14-8-7-11-16(12-14)18-17(19(2)3)13-15-9-5-4-6-10-15/h4-12H,1
InchiKey: MPEXXWGRQYCECW-ISLYRVAYSA-N
Formula: C17H20N2
SMILES: Cc1cccc(N=C(Cc2ccccc2)N(C)C)c1
Mol. weight [g/mol]: 252.35

Physical Properties

Property code	Value	Unit	Source
hf	207.34	kJ/mol	Joback Method
hvap	64.09	kJ/mol	Joback Method
log10ws	-4.03		Crippen Method
logp	3.829		Crippen Method
mcvol	218.530	ml/mol	McGowan Method
pc	1882.17	kPa	Joback Method
rinpol	2118.00		NIST Webbook
rinpol	2118.00		NIST Webbook
tb	735.70	K	Joback Method
tc	975.66	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=R162164&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