

N''-(4-nitro-phenyl)-N,N,N',N'-tetramethyl-guanidine

Inchi:	InChI=1S/C11H16N4O2/c1-13(2)11(14(3)4)12-9-5-7-10(8-6-9)15(16)17/h5-8H,1-4H3
InchiKey:	JHJHOUGVVOFMET-UHFFFAOYSA-N
Formula:	C11H16N4O2
SMILES:	CN(C)C(=Nc1ccc([N+](=O)[O-])cc1)N(C)C
Mol. weight [g/mol]:	236.27

Physical Properties

Property code	Value	Unit	Source
hf	151.42	kJ/mol	Joback Method
hvap	67.09	kJ/mol	Joback Method
log10ws	-2.08		Crippen Method
logp	1.706		Crippen Method
mcvol	185.150	ml/mol	McGowan Method
pc	2379.54	kPa	Joback Method
rinpol	2146.00		NIST Webbook
rinpol	2146.00		NIST Webbook
tb	736.02	K	Joback Method
tc	975.76	K	Joback Method

Sources

NIST Webbook:	http://webbook.nist.gov/cgi/cbook.cgi?ID=R153096&Units=SI
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

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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