

N,N-Dimethyl-N'-isopropyl-formamidine

Inchi:	InChI=1S/C6H14N2/c1-6(2)7-5-8(3)4/h5-6H,1-4H3
InchiKey:	NCHYGMRFKDUUAI-UHFFFAOYSA-N
Formula:	C6H14N2
SMILES:	CC(C)N=CN(C)C
Mol. weight [g/mol]:	114.19

Physical Properties

Property code	Value	Unit	Source
hf	-22.70	kJ/mol	Joback Method
hvap	33.92	kJ/mol	Joback Method
log10ws	-0.68		Crippen Method
logp	0.985		Crippen Method
mcvol	111.060	ml/mol	McGowan Method
pc	2790.61	kPa	Joback Method
rinpol	806.00		NIST Webbook
rinpol	806.00		NIST Webbook
tb	425.36	K	Joback Method
tc	615.39	K	Joback Method

Sources

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
NIST Webbook:	http://webbook.nist.gov/cgi/cbook.cgi?ID=R153625&Units=SI
Crippen Method:	http://pubs.acs.org/doi/abs/10.1021/ci990307l

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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