

N,N-Diethyl-N'-propylacetamidine

Other names:	N,N-Diethyl-N'-propylethanimidamide (C ₂ H ₅) ₂ N-C(CH ₃)=N(n-C ₃ H ₇) Ethanimidamide, N,N-diethyl-N'-propyl-
Inchi:	InChI=1S/C ₉ H ₂₀ N ₂ /c1-5-8-10-9(4)11(6-2)7-3/h5-8H ₂ ,1-4H ₃
InchiKey:	AGZSVOCXEYHQGE-UHFFFAOYSA-N
Formula:	C ₉ H ₂₀ N ₂
SMILES:	CCCN=C(C)N(CC)CC
Mol. weight [g/mol]:	156.27
CAS:	151328-42-6

Physical Properties

Property code	Value	Unit	Source
affp	1037.90	kJ/mol	NIST Webbook
basg	1005.50	kJ/mol	NIST Webbook
hf	-89.13	kJ/mol	Joback Method
hvap	41.06	kJ/mol	Joback Method
log10ws	-1.82		Crippen Method
logp	2.157		Crippen Method
mcvol	153.330	ml/mol	McGowan Method
pc	2096.50	kPa	Joback Method
tb	494.32	K	Joback Method
tc	678.12	K	Joback Method

Sources

Crippen Method:	http://pubs.acs.org/doi/abs/10.1021/ci990307I
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=C151328426&Units=SI

Legend

affp:	Proton affinity
basg:	Gas basicity
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
tb:	Normal Boiling Point Temperature
tc:	Critical Temperature

Latest version available from:

<https://www.chemeo.com/cid/63-992-1/N-N-Diethyl-N-propylacetamidine.pdf>

Generated by Cheméo on 2025-12-26 02:15:12.916906136 +0000 UTC m=+6463510.446946797.

Cheméo (<https://www.chemeo.com>) is the biggest free database of chemical and physical data for the process industry.