(CH3)2N-CH=N-(1,1-dimethylpropyl)

Inchi: InChI=1S/C8H18N2/c1-6-8(2,3)9-7-10(4)5/h7H,6H2,1-5H3
InchiKey: HNKZWZOXHKVDEA-UHFFFAOYSA-N
Formula: C8H18N2
SMILES: CCC(C)(C)N=CN(C)C
Mol. weight [g/mol]: 142.24
CAS: 133835-17-3

Physical Properties

Property code	Value	Unit	Source
affp	1022.00	kJ/mol	NIST Webbook
basg	989.60	kJ/mol	NIST Webbook
hf	-67.45	kJ/mol	Joback Method
hvap	37.46	kJ/mol	Joback Method
log10ws	-1.51		Crippen Method
logp	1.765		Crippen Method
mcvol	139.240	ml/mol	McGowan Method
pc	2324.78	kPa	Joback Method
tb	468.33	K	Joback Method
tc	662.25	K	Joback Method

Sources

McGowan Method: http://link.springer.com/article/10.1007/BF02311772
NIST Webbook: http://webbook.nist.gov/cgi/cbook.cgi?ID=C133835173&Units=SI
Crippen Method: http://pubs.acs.org/doi/abs/10.1021/ci990307l
Crippen Method: https://www.chemeo.com/doc/models/crippen_log10ws
Joback Method: https://en.wikipedia.org/wiki/Joback_method

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