

2-Propanamine, N,N,2-trimethyl-

Other names:	t-C ₄ H ₉ N(CH ₃) ₂ N,N-Dimethyl-t-butylamine
Inchi:	InChI=1S/C ₆ H ₁₅ N/c1-6(2,3)7(4)5/h1-5H3
InchiKey:	OXQMIXBVXHWDPX-UHFFFAOYSA-N
Formula:	C ₆ H ₁₅ N
SMILES:	CN(C)C(C)(C)C
Mol. weight [g/mol]:	101.19
CAS:	918-02-5

Physical Properties

Property code	Value	Unit	Source
affp	979.60	kJ/mol	NIST Webbook
basg	948.60	kJ/mol	NIST Webbook
gf	113.26	kJ/mol	Joback Method
hf	-108.39	kJ/mol	Joback Method
hfus	6.90	kJ/mol	Joback Method
hvap	34.80	kJ/mol	NIST Webbook
ie	8.08	eV	NIST Webbook
log10ws	-1.02		Crippen Method
logp	1.347		Crippen Method
mcvol	105.380	ml/mol	McGowan Method
pc	3166.83	kPa	Joback Method
rinpola	693.00		NIST Webbook
rinpola	693.00		NIST Webbook
tb	362.80 ± 1.50	K	NIST Webbook
tb	362.15 ± 2.00	K	NIST Webbook
tb	363.80 ± 1.00	K	NIST Webbook
tb	362.65 ± 2.00	K	NIST Webbook
tb	363.65 ± 2.00	K	NIST Webbook
tc	519.39	K	Joback Method
tf	183.15 ± 5.00	K	NIST Webbook
vc	0.379	m ³ /kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	183.93	J/mol×K	345.89	Joback Method
cpg	197.42	J/mol×K	374.81	Joback Method
cpg	210.25	J/mol×K	403.72	Joback Method
cpg	222.45	J/mol×K	432.64	Joback Method
cpg	234.02	J/mol×K	461.56	Joback Method
cpg	245.01	J/mol×K	490.48	Joback Method
cpg	255.44	J/mol×K	519.39	Joback Method

Sources

McGowan Method:	http://link.springer.com/article/10.1007/BF02311772
NIST Webbook:	http://webbook.nist.gov/cgi/cbook.cgi?ID=C918025&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

Legend

affp:	Proton affinity
basg:	Gas basicity
cpg:	Ideal gas heat capacity
gf:	Standard Gibbs free energy of formation
hf:	Enthalpy of formation at standard conditions
hfus:	Enthalpy of fusion at standard conditions
hvap:	Enthalpy of vaporization at standard conditions
ie:	Ionization energy
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

tf: Normal melting (fusion) point

vc: Critical Volume

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