Propanamine, 2-methyl, N,N-dipropyl

Other names: Dipropyl isobutylamine
Inchi: InChI=1S/C10H23N/c1-5-7-11(8-6-2)9-10(3)4/h10H,5-9H2,1-4H3
InchiKey: PSZIEXSRZMVQEN-UHFFFAOYSA-N
Formula: C10H23N
SMILES: CCCN(CCC)CC(C)C
Mol. weight [g/mol]: 157.30

Physical Properties

Property code	Value	Unit	Source
gf	141.66	kJ/mol	Joback Method
hf	-187.48	kJ/mol	Joback Method
hfus	21.15	kJ/mol	Joback Method
hvap	39.51	kJ/mol	Joback Method
log10ws	-2.34		Crippen Method
logp	2.764		Crippen Method
mcvol	161.740	ml/mol	McGowan Method
pc	2119.73	kPa	Joback Method
rinpol	961.00		NIST Webbook
rinpol	961.00		NIST Webbook
tb	440.20	K	Joback Method
tc	604.82	K	Joback Method
tf	219.93	K	Joback Method
vc	0.608	m3/kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	347.56	J/mol×K	440.20	Joback Method
cpg	363.91	J/mol×K	467.64	Joback Method
cpg	379.60	J/mol×K	495.07	Joback Method
cpg	394.66	J/mol×K	522.51	Joback Method
cpg	409.10	J/mol×K	549.94	Joback Method
cpg	422.95	J/mol×K	577.38	Joback Method
cpg	436.22	J/mol×K	604.82	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=R12953&Units=SI
Crippen Method:	http://pubs.acs.org/doi/abs/10.1021/ci990307l

Legend

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
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