

1-Pentamine, N-methyl-

Other names:	Pentylamine, N-methyl- Methylamylamine N-Amyl-N-methylamine N-Methylamylamine N-Methylpentanamine N-Methylpentylamine N-Methyl-N-amylamine Amylmethylamine N-Methyl-n-pentylamine
Inchi:	InChI=1S/C6H15N/c1-3-4-5-6-7-2/h7H,3-6H2,1-2H3
InchiKey:	UOIWOHLIGKIYFE-UHFFFAOYSA-N
Formula:	C6H15N
SMILES:	CCCCCNC
Mol. weight [g/mol]:	101.19
CAS:	25419-06-1

Physical Properties

Property code	Value	Unit	Source
gf	89.03	kJ/mol	Joback Method
hf	-113.70	kJ/mol	Joback Method
hfus	16.39	kJ/mol	Joback Method
hvap	35.39	kJ/mol	Joback Method
log10ws	-1.52		Crippen Method
logp	1.396		Crippen Method
mcvol	105.380	ml/mol	McGowan Method
pc	3152.62	kPa	Joback Method
rinpol	706.00		NIST Webbook
rinpol	706.00		NIST Webbook
rinpol	706.00		NIST Webbook
tb	390.15 ± 3.00	K	NIST Webbook
tc	556.73	K	Joback Method
tf	210.04	K	Joback Method
vc	0.406	m ³ /kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	199.23	J/mol×K	386.85	Joback Method
cpg	210.62	J/mol×K	415.16	Joback Method
cpg	221.60	J/mol×K	443.48	Joback Method
cpg	232.16	J/mol×K	471.79	Joback Method
cpg	242.32	J/mol×K	500.10	Joback Method
cpg	252.08	J/mol×K	528.41	Joback Method
cpg	261.47	J/mol×K	556.73	Joback Method

Sources

McGowan Method:	http://link.springer.com/article/10.1007/BF02311772
NIST Webbook:	http://webbook.nist.gov/cgi/cbook.cgi?ID=C25419061&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

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

Latest version available from:

<https://www.cheméo.com/cid/77-767-6/1-Pentanamine-N-methyl.pdf>

Generated by Cheméo on 2024-04-25 03:41:24.038174723 +0000 UTC m=+16305732.958752036.

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