

Methanal, diethylhydrazone

Inchi: InChI=1S/C5H12N2/c1-4-7(5-2)6-3/h3-5H2,1-2H3
InchiKey: ZTLZQVMNFGEQRQ-UHFFFAOYSA-N
Formula: C5H12N2
SMILES: C=NN(CC)CC
Mol. weight [g/mol]: 100.16

Physical Properties

Property code	Value	Unit	Source
hf	11.43	kJ/mol	Joback Method
hvap	31.45	kJ/mol	Joback Method
log10ws	-0.64		Crippen Method
logp	0.944		Crippen Method
mcvol	96.970	ml/mol	McGowan Method
pc	3035.62	kPa	Joback Method
rinpol	760.00		NIST Webbook
tb	395.44	K	Joback Method
tc	577.13	K	Joback Method

Sources

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=R511883&Units=SI>
Crippen Method: <http://pubs.acs.org/doi/abs/10.1021/ci990307I>
Crippen Method: https://www.chemeo.com/doc/models/crippen_log10ws

Legend

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
rinpol:	Non-polar retention indices
tb:	Normal Boiling Point Temperature
tc:	Critical Temperature

Latest version available from:

<https://www.cheméo.com/cid/58-319-4/Methanal-diethylhydrazone.pdf>

Generated by Cheméo on 2024-04-28 01:16:51.182986009 +0000 UTC m=+16556260.103563324.

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