

Butanal, dimethylhydrazone

Other names:	Butyraldehyde, dimethylhydrazone
Inchi:	InChI=1S/C6H14N2/c1-4-5-6-7-8(2)3/h6H,4-5H2,1-3H3
InchiKey:	SEEKWDSXURZRIV-UHFFFAOYSA-N
Formula:	C6H14N2
SMILES:	CCCC=NN(C)C
Mol. weight [g/mol]:	114.19
CAS:	10424-98-3

Physical Properties

Property code	Value	Unit	Source
hf	-17.42	kJ/mol	Joback Method
hvap	34.31	kJ/mol	Joback Method
log10ws	-1.06		Crippen Method
logp	1.334		Crippen Method
mcvol	111.060	ml/mol	McGowan Method
pc	2767.17	kPa	Joback Method
rinpol	860.00		NIST Webbook
rinpol	860.00		NIST Webbook
rinpol	860.00		NIST Webbook
rinpol	878.90		NIST Webbook
tb	425.80	K	Joback Method
tc	611.49	K	Joback Method

Sources

NIST Webbook:	http://webbook.nist.gov/cgi/cbook.cgi?ID=C10424983&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
McGowan Method:	http://link.springer.com/article/10.1007/BF02311772

Legend

hf:	Enthalpy of formation at standard conditions
h_{vap}:	Enthalpy of vaporization at standard conditions
log₁₀ws:	Log10 of Water solubility in mol/l
log_p:	Octanol/Water partition coefficient
mcvol:	McGowan's characteristic volume
pc:	Critical Pressure
r_{inpol}:	Non-polar retention indices
tb:	Normal Boiling Point Temperature
tc:	Critical Temperature

Latest version available from:

<https://www.cheméo.com/cid/64-789-6/Butanal-dimethylhydrazone.pdf>

Generated by Cheméo on 2025-12-06 01:49:54.390892857 +0000 UTC m=+4733991.920933521.

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