

Acetic acid, 2-(2-methylpropylidene)hydrazide

Inchi: InChI=1S/C6H12N2O/c1-5(2)4-7-8-6(3)9/h4-5H,1-3H3,(H,8,9)/b7-4+
InchiKey: JQPAOGJTPACAPO-QPJJXVBHSA-N
Formula: C6H12N2O
SMILES: CC(=O)NN=CC(C)C
Mol. weight [g/mol]: 128.17
CAS: 166327-19-1

Physical Properties

Property code	Value	Unit	Source
hf	-149.34	kJ/mol	Joback Method
hvap	45.06	kJ/mol	Joback Method
log10ws	-1.21		Crippen Method
logp	0.764		Crippen Method
mcvol	112.630	ml/mol	McGowan Method
pc	3035.62	kPa	Joback Method
tb	516.96	K	Joback Method
tc	722.27	K	Joback Method

Sources

McGowan Method: <http://link.springer.com/article/10.1007/BF02311772>
NIST Webbook: <http://webbook.nist.gov/cgi/cbook.cgi?ID=C166327191&Units=SI>
Crippen Method: <http://pubs.acs.org/doi/abs/10.1021/ci990307I>
Crippen Method: https://www.chemeo.com/doc/models/crippen_log10ws
Joback Method: https://en.wikipedia.org/wiki/Joback_method

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
tb:	Normal Boiling Point Temperature
tc:	Critical Temperature

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