2-Butanone, O-methyloxime

Inchi: InChI=1S/C5H11NO/c1-4-5(2)6-7-3/h4H2,1-3H3
InchiKey: CYKDEJOBIIVIQD-UHFFFAOYSA-N
Formula: C5H11NO
SMILES: CCC(C)=NOC
Mol. weight [g/mol]: 101.15

Physical Properties

Property code	Value	Unit	Source
hf	-206.32	kJ/mol	Joback Method
hvap	32.53	kJ/mol	Joback Method
log10ws	-1.16		Crippen Method
logp	1.419		Crippen Method
mcvol	92.860	ml/mol	McGowan Method
pc	3035.62	kPa	Joback Method
rinpol	687.00		NIST Webbook
rinpol	687.00		NIST Webbook
tb	412.78	K	Joback Method
tc	605.34	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=R510964&Units=SI
Crippen Method: http://pubs.acs.org/doi/abs/10.1021/ci990307l
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