Acetone, O-propyloxime

Inchi: InChI=1S/C6H13NO/c1-4-5-8-7-6(2)3/h4-5H2,1-3H3
InchiKey: HOFOCPOTGQTZSI-UHFFFAOYSA-N
Formula: C6H13NO
SMILES: CCCON=C(C)C
Mol. weight [g/mol]: 115.17

Physical Properties

Property code	Value	Unit	Source
hf	-226.96	kJ/mol	Joback Method
hvap	34.75	kJ/mol	Joback Method
log10ws	-1.58		Crippen Method
logp	1.809		Crippen Method
mcvol	106.950	ml/mol	McGowan Method
pc	2738.28	kPa	Joback Method
rinpol	763.00		NIST Webbook
rinpol	763.00		NIST Webbook
tb	435.66	K	Joback Method
tc	626.47	K	Joback Method

Sources

NIST Webbook: http://webbook.nist.gov/cgi/cbook.cgi?ID=R511303&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
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