Acetone, 2-hydroxyethylimine

Inchi: InChI=1S/C5H11NO/c1-5(2)6-3-4-7/h7H,3-4H2,1-2H3
InchiKey: PIOHLQLULGTQAC-UHFFFAOYSA-N
Formula: C5H11NO
SMILES: CC(C)=NCCO
Mol. weight [g/mol]: 101.15

Physical Properties

Property code	Value	Unit	Source
hf	-226.33	kJ/mol	Joback Method
hvap	46.80	kJ/mol	Joback Method
log10ws	-0.34		Crippen Method
logp	0.460		Crippen Method
mcvol	92.860	ml/mol	McGowan Method
pc	3435.91	kPa	Joback Method
rinpol	773.00		NIST Webbook
rinpol	773.00		NIST Webbook
tb	482.54	K	Joback Method
tc	666.81	K	Joback Method

Sources

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
NIST Webbook: http://webbook.nist.gov/cgi/cbook.cgi?ID=R511274&Units=SI
Crippen Method: http://pubs.acs.org/doi/abs/10.1021/ci990307l

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