2(1H)-Pyridone, 6-methyl-

Other names:	6-Methyl-2-pyridinol
	2(1H)-Pyridinone, 6-methyl-
	2-Methyl-6-pyridone
	6-Methyl-2-hydroxypyridine
	6-Methyl-2-pyridinone
	6-Methyl-2-pyridone
	2-Hydroxy-6-methylpyridine
	6-Methyl-2-pyridyl alcohol
	6-Methyl-2(1H)-pyridone
	2-Methyl-6-hydroxypyridine
	6-Methyl-2(1H)-pyridinone
	6-methylpyridin-2-ol
Inchi:	InChI=1S/C6H7NO/c1-5-3-2-4-6(8)7-5/h2-4H,1H3,(H,7,8)
InchiKey:	JEAVIRYCMBDJIU-UHFFFAOYSA-N
Formula:	C6H7NO
SMILES:	Cc1cccc(=O)[nH]1
Mol. weight [g/mol]:	109.13
CAS:	3279-76-3

Physical Properties

Property code	Value	Unit	Source
chs	-3149.20 ± 1.90	kJ/mol	NIST Webbook
hf	-120.30 ± 2.50	kJ/mol	NIST Webbook
hfs	-212.30 ± 2.10	kJ/mol	NIST Webbook
hsub	92.00 ± 1.30	kJ/mol	NIST Webbook
hsub	92.00 ± 1.30	kJ/mol	NIST Webbook
ie	8.19 ± 0.03	eV	NIST Webbook
log10ws	-0.59		Crippen Method
logp	0.201		Crippen Method
mcvol	87.490	ml/mol	McGowan Method

Sources

Crippen Method: https://www.chemeo.com/doc/models/crippen_log10ws

McGowan Method:	http://link.springer.com/article/10.1007/BF02311772
NIST Webbook:	http://webbook.nist.gov/cgi/cbook.cgi?ID=C3279763&Units=SI
Crippen Method:	http://pubs.acs.org/doi/abs/10.1021/ci990307l

Legend

chs:	Standard solid enthalpy of combustion
hf:	Enthalpy of formation at standard conditions
hfs:	Solid phase enthalpy of formation at standard conditions
hsub:	Enthalpy of sublimation at standard conditions
ie:	Ionization energy
log10ws:	Log10 of Water solubility in mol/l
logp:	Octanol/Water partition coefficient
mcvol:	McGowan's characteristic volume