

2-Pyridinealdoxime

Other names:	2-Formylpyridine ketoxime
	2-Formylpyridine oxime
	2-Hydroxyiminomethyl pyridine
	2-Pyridinaldoxime
	2-Pyridinecarbaldehyde oxime
	2-Pyridinecarboxaldehyde, oxime
	2-Pyridinecarboxaldoxime
	2-Pyridylaldoxime
	NSC 66484
	P-2-A
	Picolinaldehyde, oxime
	Picolinaldoxime
	Picolinealdehyde, oxime
	Pyridine-2-aldoxime
	Pyridine-2-carboxaldoxime
	pyridine-2-carbaldehyde oxime
	«alpha»-Picolinealdoxime
	Â«alphaÂ»-Picolinealdoxime
Inchi:	InChI=1S/C6H6N2O/c9-8-5-6-3-1-2-4-7-6/h1-5,9H
InchiKey:	MTFJSAGADRTKCI-UHFFFAOYSA-N
Formula:	C6H6N2O
SMILES:	ON=Cc1ccccn1
Mol. weight [g/mol]:	122.12
CAS:	873-69-8

Physical Properties

Property code	Value	Unit	Source
hfus	19.98	kJ/mol	Low-temperature heat capacity and standard molar enthalpy of formation of crystalline 2-pyridinealdoxime (C6H6N2O)
log10ws	-0.54		Crippen Method
logp	0.890		Crippen Method
mcvol	93.170	ml/mol	McGowan Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
hfust	19.98	kJ/mol	388.00	NIST Webbook

Sources

Crippen Method:	http://pubs.acs.org/doi/abs/10.1021/ci990307l
Crippen Method:	https://www.chemeo.com/doc/models/crippen_log10ws
Low-temperature heat capacity and standard molar enthalpy of formation	https://www.doi.org/10.1016/j.jct.2006.09.016
McGowan's method	http://link.springer.com/article/10.1007/BF02311772
Pyridinealdoxime (C ₅ H ₆ N ₂ O):	http://webbook.nist.gov/cgi/cbook.cgi?ID=C873698&Units=SI
NIST Webbook:	

Legend

hfus:	Enthalpy of fusion at standard conditions
hfust:	Enthalpy of fusion at a given temperature
log10ws:	Log10 of Water solubility in mol/l
logp:	Octanol/Water partition coefficient
mcvol:	McGowan's characteristic volume

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