

# Pyridine, 4-ethenyl-

<b>Other names:</b>	Pyridine, 4-vinyl- 4-Vinylpyridine 4-Ethenylpyridine «gamma»-Vinylpyridine
<b>Inchi:</b>	InChI=1S/C7H7N/c1-2-7-3-5-8-6-4-7/h2-6H,1H2
<b>InchiKey:</b>	KFDVPJUYSDEJTH-UHFFFAOYSA-N
<b>Formula:</b>	C7H7N
<b>SMILES:</b>	C=Cc1ccncc1
<b>Mol. weight [g/mol]:</b>	105.14
<b>CAS:</b>	100-43-6

## Physical Properties

Property code	Value	Unit	Source
affp	944.10	kJ/mol	NIST Webbook
basg	912.30	kJ/mol	NIST Webbook
ie	9.25	eV	NIST Webbook
ie	8.90	eV	NIST Webbook
log10ws	-2.06		Crippen Method
logp	1.725		Crippen Method
mcvol	91.410	ml/mol	McGowan Method
rinpol	1026.00		NIST Webbook
rinpol	1026.00		NIST Webbook

## Pressure Dependent Properties

Property code	Value	Unit	Pressure [kPa]	Source
tbrp	336.70	K	2.00	NIST Webbook

## Sources

**Crippen Method:** [https://www.chemeo.com/doc/models/crippen\\_log10ws](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=C100436&Units=SI>  
**Crippen Method:** <http://pubs.acs.org/doi/abs/10.1021/ci990307l>

## Legend

**affp:** Proton affinity  
**basg:** Gas basicity  
**ie:** Ionization energy  
**log10ws:** Log10 of Water solubility in mol/l  
**logp:** Octanol/Water partition coefficient  
**mcvol:** McGowan's characteristic volume  
**rinpol:** Non-polar retention indices  
**tbrp:** Boiling point at reduced pressure

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