

# Benzenecarboximidic acid, N-methyl-, methyl ester

**Inchi:** InChI=1S/C9H11NO/c1-10-9(11-2)8-6-4-3-5-7-8/h3-7H,1-2H3/b10-9-  
**InchiKey:** ODCHIEXXFGKZNZ-KTKRTIGZSA-N  
**Formula:** C9H11NO  
**SMILES:** CN=C(OC)c1ccccc1  
**Mol. weight [g/mol]:** 149.19  
**CAS:** 1775-61-7

## Physical Properties

Property code	Value	Unit	Source
hf	-52.35	kJ/mol	Joback Method
hvap	43.71	kJ/mol	Joback Method
log10ws	-1.54		Crippen Method
logp	1.709		Crippen Method
mcvol	125.460	ml/mol	McGowan Method
pc	2865.80	kPa	Joback Method
tb	530.98	K	Joback Method
tc	760.95	K	Joback Method

## Sources

**Joback Method:** [https://en.wikipedia.org/wiki/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=C1775617&Units=SI>  
**Crippen Method:** <http://pubs.acs.org/doi/abs/10.1021/ci990307I>  
**Crippen Method:** [https://www.chemeo.com/doc/models/crippen\\_log10ws](https://www.chemeo.com/doc/models/crippen_log10ws)

## Legend

**hf:** Enthalpy of formation at standard conditions  
**hvap:** Enthalpy of vaporization at standard conditions  
**log10ws:** Log10 of Water solubility in mol/l

<b>logp:</b>	Octanol/Water partition coefficient
<b>mcvol:</b>	McGowan's characteristic volume
<b>pc:</b>	Critical Pressure
<b>tb:</b>	Normal Boiling Point Temperature
<b>tc:</b>	Critical Temperature

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