

# Benzene, 1-isocyanato-4-methoxy-

<b>Other names:</b>	4-Methoxyphenylisocyanate p-Methoxyphenyl isocyanate p-Methoxyphenyl 1socyanate p-Anisyl isocyanate Isocyanic acid, p-methoxyphenyl ester
<b>Inchi:</b>	InChI=1S/C8H7NO2/c1-11-8-4-2-7(3-5-8)9-6-10/h2-5H,1H3
<b>InchiKey:</b>	FMDGXCSMDZMDHZ-UHFFFAOYSA-N
<b>Formula:</b>	C8H7NO2
<b>SMILES:</b>	COc1ccc(N=C=O)cc1
<b>Mol. weight [g/mol]:</b>	149.15
<b>CAS:</b>	5416-93-3

## Physical Properties

Property code	Value	Unit	Source
hf	-121.02	kJ/mol	Joback Method
hvap	48.28	kJ/mol	Joback Method
log10ws	-6.07		Crippen Method
logp	1.662		Crippen Method
mcvol	112.940	ml/mol	McGowan Method
pc	3843.54	kPa	Joback Method
tb	503.19	K	Joback Method
tc	723.14	K	Joback Method

## Sources

**Crippen Method:** [https://www.chemeo.com/doc/models/crippen\\_log10ws](https://www.chemeo.com/doc/models/crippen_log10ws)

**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=C5416933&Units=SI>

**Crippen Method:** <http://pubs.acs.org/doi/abs/10.1021/ci990307l>

# Legend

<b>hf:</b>	Enthalpy of formation at standard conditions
<b>hvap:</b>	Enthalpy of vaporization at standard conditions
<b>log10ws:</b>	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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