

# 2-Fluorophenyl isocyanate

<b>Other names:</b>	o-Fluorophenyl isocyanate Isocyanic acid o-fluorophenyl ester Benzene, 1-fluoro-2-isocyanato-
<b>Inchi:</b>	InChI=1S/C7H4FNO/c8-6-3-1-2-4-7(6)9-5-10/h1-4H
<b>InchiKey:</b>	VZNCSZQPNIEEMN-UHFFFAOYSA-N
<b>Formula:</b>	C7H4FNO
<b>SMILES:</b>	O=C=Nc1ccccc1F
<b>Mol. weight [g/mol]:</b>	137.11
<b>CAS:</b>	16744-98-2

## Physical Properties

Property code	Value	Unit	Source
hf	-164.27	kJ/mol	Joback Method
hvap	42.83	kJ/mol	Joback Method
log10ws	-6.29		Crippen Method
logp	1.793		Crippen Method
mcvol	94.750	ml/mol	McGowan Method
pc	4189.32	kPa	Joback Method
tb	457.16	K	Joback Method
tc	670.64	K	Joback Method

## Sources

<b>Crippen Method:</b>	<a href="https://www.chemeo.com/doc/models/crippen_log10ws">https://www.chemeo.com/doc/models/crippen_log10ws</a>
<b>Joback Method:</b>	<a href="https://en.wikipedia.org/wiki/Joback_method">https://en.wikipedia.org/wiki/Joback_method</a>
<b>McGowan Method:</b>	<a href="http://link.springer.com/article/10.1007/BF02311772">http://link.springer.com/article/10.1007/BF02311772</a>
<b>NIST Webbook:</b>	<a href="http://webbook.nist.gov/cgi/cbook.cgi?ID=C16744982&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=C16744982&amp;Units=SI</a>
<b>Crippen Method:</b>	<a href="http://pubs.acs.org/doi/abs/10.1021/ci9903071">http://pubs.acs.org/doi/abs/10.1021/ci9903071</a>

## Legend

<b>hf:</b>	Enthalpy of formation at standard conditions
<b>h<sub>vap</sub>:</b>	Enthalpy of vaporization at standard conditions
<b>log<sub>10</sub>ws:</b>	Log <sub>10</sub> of Water solubility in mol/l
<b>log<sub>p</sub>:</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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