

# Carbamic acid, dimethyl-, ethyl ester

<b>Other names:</b>	Ethyl dimethylcarbamate Dimethylethoxyformamide Ethyl N,N-dimethylcarbamate C2H5OC(O)N(CH3)2 (CH3)2NCOOC2H5
<b>Inchi:</b>	InChI=1S/C5H11NO2/c1-4-8-5(7)6(2)3/h4H2,1-3H3
<b>InchiKey:</b>	SUDHEDJJFGYYPL-UHFFFAOYSA-N
<b>Formula:</b>	C5H11NO2
<b>SMILES:</b>	CCOC(=O)N(C)C
<b>Mol. weight [g/mol]:</b>	117.15
<b>CAS:</b>	687-48-9

## Physical Properties

Property code	Value	Unit	Source
affp	896.60	kJ/mol	NIST Webbook
basg	865.60	kJ/mol	NIST Webbook
gf	-131.92	kJ/mol	Joback Method
hf	-323.80	kJ/mol	Joback Method
hfus	14.51	kJ/mol	Joback Method
hvap	37.92	kJ/mol	Joback Method
log10ws	-0.33		Crippen Method
logp	0.705		Crippen Method
mvol	98.730	ml/mol	McGowan Method
pc	3655.35	kPa	Joback Method
tb	402.53	K	Joback Method
tc	579.52	K	Joback Method
tf	250.74	K	Joback Method
vc	0.357	m <sup>3</sup> /kmol	Joback Method

## Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	190.41	J/mol×K	402.53	Joback Method
cpg	200.15	J/mol×K	432.03	Joback Method

cpg	209.53	J/mol×K	461.53	Joback Method
cpg	218.57	J/mol×K	491.03	Joback Method
cpg	227.26	J/mol×K	520.53	Joback Method
cpg	235.61	J/mol×K	550.03	Joback Method
cpg	243.61	J/mol×K	579.52	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=C687489&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=C687489&amp;Units=SI</a>
<b>Crippen Method:</b>	<a href="http://pubs.acs.org/doi/abs/10.1021/ci990307l">http://pubs.acs.org/doi/abs/10.1021/ci990307l</a>

## Legend

<b>affp:</b>	Proton affinity
<b>basg:</b>	Gas basicity
<b>cpg:</b>	Ideal gas heat capacity
<b>gf:</b>	Standard Gibbs free energy of formation
<b>hf:</b>	Enthalpy of formation at standard conditions
<b>hfus:</b>	Enthalpy of fusion 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
<b>tf:</b>	Normal melting (fusion) point
<b>vc:</b>	Critical Volume

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