

# Diethyl 3-methylcyclopropane-1,2-dicarboxylate

Inchi:	InChI=1S/C10H16O4/c1-4-13-9(11)7-6(3)8(7)10(12)14-5-2/h6-8H,4-5H2,1-3H3
InchiKey:	IRISXSVBOADFLC-UHFFFAOYSA-N
Formula:	C10H16O4
SMILES:	CCOC(=O)C1C(C)C1C(=O)OCC
Mol. weight [g/mol]:	200.23
CAS:	4104-67-0

## Physical Properties

Property code	Value	Unit	Source
gf	-389.19	kJ/mol	Joback Method
hf	-707.21	kJ/mol	Joback Method
hfus	27.51	kJ/mol	Joback Method
hvap	55.46	kJ/mol	Joback Method
log10ws	-0.90		Crippen Method
logp	0.995		Crippen Method
mcvol	155.780	ml/mol	McGowan Method
pc	2445.89	kPa	Joback Method
tb	578.18	K	Joback Method
tc	769.71	K	Joback Method
tf	356.24	K	Joback Method
vc	0.599	m3/kmol	Joback Method

## Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	404.49	J/molxK	578.18	Joback Method
cpg	418.98	J/molxK	610.10	Joback Method
cpg	432.80	J/molxK	642.02	Joback Method
cpg	445.95	J/molxK	673.94	Joback Method
cpg	458.45	J/molxK	705.87	Joback Method
cpg	470.29	J/molxK	737.79	Joback Method
cpg	481.48	J/molxK	769.71	Joback Method
dvisc	0.0016078	Paxs	356.24	Joback Method
dvisc	0.0012420	Paxs	393.23	Joback Method

dvisc	0.0010031	Paxs	430.22	Joback Method
dvisc	0.0008379	Paxs	467.21	Joback Method
dvisc	0.0007187	Paxs	504.20	Joback Method
dvisc	0.0006295	Paxs	541.19	Joback Method
dvisc	0.0005608	Paxs	578.18	Joback Method

## Sources

<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=C4104670&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=C4104670&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>
<b>Crippen Method:</b>	<a href="https://www.chemeo.com/doc/models/crippen_log10ws">https://www.chemeo.com/doc/models/crippen_log10ws</a>

## Legend

<b>cpg:</b>	Ideal gas heat capacity
<b>dvisc:</b>	Dynamic viscosity
<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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