

# Cyclopropanecarboxylic acid, 4-methoxyphenyl ester

Inchi:	InChI=1S/C11H12O3/c1-13-9-4-6-10(7-5-9)14-11(12)8-2-3-8/h4-8H,2-3H2,1H3
InchiKey:	JUNVKRVSFQQOPN-UHFFFAOYSA-N
Formula:	C11H12O3
SMILES:	COc1ccc(OC(=O)C2CC2)cc1
Mol. weight [g/mol]:	192.21

## Physical Properties

Property code	Value	Unit	Source
gf	-133.65	kJ/mol	Joback Method
hf	-349.53	kJ/mol	Joback Method
hfus	20.01	kJ/mol	Joback Method
hvap	54.50	kJ/mol	Joback Method
log10ws	-2.40		Crippen Method
logp	2.011		Crippen Method
mcvol	144.540	ml/mol	McGowan Method
pc	3089.85	kPa	Joback Method
rinpol	1536.00		NIST Webbook
tb	588.19	K	Joback Method
tc	811.44	K	Joback Method
tf	365.00	K	Joback Method
vc	0.542	m3/kmol	Joback Method

## Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	356.08	J/molxK	588.19	Joback Method
cpg	370.58	J/molxK	625.40	Joback Method
cpg	384.16	J/molxK	662.61	Joback Method
cpg	396.86	J/molxK	699.81	Joback Method
cpg	408.72	J/molxK	737.02	Joback Method
cpg	419.76	J/molxK	774.23	Joback Method
cpg	430.01	J/molxK	811.44	Joback Method
dvisc	0.0014390	Paxs	365.00	Joback Method
dvisc	0.0010344	Paxs	402.20	Joback Method

dvisc	0.0007863	Paxs	439.40	Joback Method
dvisc	0.0006238	Paxs	476.60	Joback Method
dvisc	0.0005118	Paxs	513.79	Joback Method
dvisc	0.0004313	Paxs	550.99	Joback Method
dvisc	0.0003714	Paxs	588.19	Joback Method

## Sources

<b>NIST Webbook:</b>	<a href="http://webbook.nist.gov/cgi/cbook.cgi?ID=U307524&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=U307524&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>
<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>

## 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>rinpol:</b>	Non-polar retention indices
<b>tb:</b>	Normal Boiling Point Temperature
<b>tc:</b>	Critical Temperature
<b>tf:</b>	Normal melting (fusion) point
<b>vc:</b>	Critical Volume

Latest version available from:

<https://www.chemeo.com/cid/12-201-4/Cyclopropanecarboxylic-acid-4-methoxyphenyl-ester.pdf>

Generated by Cheméo on 2024-04-19 14:37:54.464515222 +0000 UTC m=+15826723.385092534.

Cheméo (<https://www.chemeo.com>) is the biggest free database of chemical and physical data for the process industry.