

Carbonic acid, ethyl 4-methoxyphenyl ester

Inchi:	InChI=1S/C10H12O4/c1-3-13-10(11)14-9-6-4-8(12-2)5-7-9/h4-7H,3H2,1-2H3
InchiKey:	GBSOFIMKLOZOHN-UHFFFAOYSA-N
Formula:	C10H12O4
SMILES:	CCOC(=O)Oc1ccc(OC)cc1
Mol. weight [g/mol]:	196.20

Physical Properties

Property code	Value	Unit	Source
gf	-307.82	kJ/mol	Joback Method
hf	-533.91	kJ/mol	Joback Method
hfus	20.47	kJ/mol	Joback Method
hvap	54.77	kJ/mol	Joback Method
log10ws	-2.39		Crippen Method
logp	2.230		Crippen Method
mcvol	147.180	ml/mol	McGowan Method
pc	2921.84	kPa	Joback Method
rinsol	1505.00		NIST Webbook
tb	580.99	K	Joback Method
tc	790.21	K	Joback Method
tf	358.02	K	Joback Method
vc	0.547	m ³ /kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	347.86	J/molxK	580.99	Joback Method
cpg	360.73	J/molxK	615.86	Joback Method
cpg	372.99	J/molxK	650.73	Joback Method
cpg	384.61	J/molxK	685.60	Joback Method
cpg	395.59	J/molxK	720.47	Joback Method
cpg	405.91	J/molxK	755.34	Joback Method
cpg	415.54	J/molxK	790.21	Joback Method
dvisc	0.0010140	Paxs	358.02	Joback Method
dvisc	0.0006243	Paxs	395.18	Joback Method

dvisc	0.0004178	Paxs	432.34	Joback Method
dvisc	0.0002979	Paxs	469.50	Joback Method
dvisc	0.0002233	Paxs	506.67	Joback Method
dvisc	0.0001740	Paxs	543.83	Joback Method
dvisc	0.0001401	Paxs	580.99	Joback Method

Sources

NIST Webbook:	http://webbook.nist.gov/cgi/cbook.cgi?ID=U357919&Units=SI
Crippen Method:	http://pubs.acs.org/doi/abs/10.1021/ci9903071
Crippen Method:	https://www.chemeo.com/doc/models/crippen_log10ws
Joback Method:	https://en.wikipedia.org/wiki/Joback_method
McGowan Method:	http://link.springer.com/article/10.1007/BF02311772

Legend

cpg:	Ideal gas heat capacity
dvisc:	Dynamic viscosity
gf:	Standard Gibbs free energy of formation
hf:	Enthalpy of formation at standard conditions
hfus:	Enthalpy of fusion at standard conditions
hvap:	Enthalpy of vaporization at standard conditions
log10ws:	Log10 of Water solubility in mol/l
logp:	Octanol/Water partition coefficient
mcvol:	McGowan's characteristic volume
pc:	Critical Pressure
rinpol:	Non-polar retention indices
tb:	Normal Boiling Point Temperature
tc:	Critical Temperature
tf:	Normal melting (fusion) point
vc:	Critical Volume

Latest version available from:

<https://www.chemeo.com/cid/29-251-1/Carbonic-acid-ethyl-4-methoxyphenyl-ester.pdf>

Generated by Cheméo on 2024-04-26 21:30:18.127792081 +0000 UTC m=+16456267.048369397.

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