

2-Cyclopentene-1-carboxylic acid

Inchi:	InChI=1S/C6H8O2/c7-6(8)5-3-1-2-4-5/h1,3,5H,2,4H2,(H,7,8)
InchiKey:	MOMBAXHNIPLMSI-UHFFFAOYSA-N
Formula:	C6H8O2
SMILES:	O=C(O)C1C=CCC1
Mol. weight [g/mol]:	112.13
CAS:	2348-89-2

Physical Properties

Property code	Value	Unit	Source
gf	-199.59	kJ/mol	Joback Method
hf	-313.72	kJ/mol	Joback Method
hfus	12.14	kJ/mol	Joback Method
hvap	52.92	kJ/mol	Joback Method
log10ws	-0.94		Crippen Method
logp	1.037		Crippen Method
mcvol	87.680	ml/mol	McGowan Method
pc	4938.44	kPa	Joback Method
ripol	1642.00		NIST Webbook
tb	497.17	K	Joback Method
tc	697.05	K	Joback Method
tf	279.79	K	Joback Method
vc	0.324	m ³ /kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	190.10	J/molxK	497.17	Joback Method
cpg	234.08	J/molxK	663.73	Joback Method
cpg	226.35	J/molxK	630.42	Joback Method
cpg	218.12	J/molxK	597.11	Joback Method
cpg	209.35	J/molxK	563.80	Joback Method
cpg	200.02	J/molxK	530.48	Joback Method
cpg	241.31	J/molxK	697.05	Joback Method
dvisc	0.0002496	Paxs	497.17	Joback Method

dvisc	0.0003799	Paxs	460.94	Joback Method
dvisc	0.0006212	Paxs	424.71	Joback Method
dvisc	0.0011131	Paxs	388.48	Joback Method
dvisc	0.0022488	Paxs	352.25	Joback Method
dvisc	0.0053385	Paxs	316.02	Joback Method
dvisc	0.0158532	Paxs	279.79	Joback Method

Sources

McGowan Method:	http://link.springer.com/article/10.1007/BF02311772
NIST Webbook:	http://webbook.nist.gov/cgi/cbook.cgi?ID=C2348892&Units=SI
Crippen Method:	http://pubs.acs.org/doi/abs/10.1021/ci9903071
Crippen Method:	https://www.cheméo.com/doc/models/crippen_log10ws
Joback Method:	https://en.wikipedia.org/wiki/Joback_method

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
ripol:	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.cheméo.com/cid/81-047-0/2-Cyclopentene-1-carboxylic-acid.pdf>

Generated by Cheméo on 2024-04-18 23:19:36.10753277 +0000 UTC m=+15771625.028110091.

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