

Cyclobutanecarboxylic acid, 2-chlorophenyl ester

Inchi:	InChI=1S/C11H11ClO2/c12-9-6-1-2-7-10(9)14-11(13)8-4-3-5-8/h1-2,6-8H,3-5H2
InchiKey:	QQMAENAUZRPEMZ-UHFFFAOYSA-N
Formula:	C11H11ClO2
SMILES:	O=C(Oc1ccccc1Cl)C1CCC1
Mol. weight [g/mol]:	210.66

Physical Properties

Property code	Value	Unit	Source
gf	-52.68	kJ/mol	Joback Method
hf	-239.21	kJ/mol	Joback Method
hfus	20.92	kJ/mol	Joback Method
hvap	56.64	kJ/mol	Joback Method
log10ws	-3.38		Crippen Method
logp	3.046		Crippen Method
mvol	150.910	ml/mol	McGowan Method
pc	3107.10	kPa	Joback Method
rinpol	1527.00		NIST Webbook
rinpol	1527.00		NIST Webbook
tb	607.47	K	Joback Method
tc	845.52	K	Joback Method
tf	369.17	K	Joback Method
vc	0.566	m ³ /kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	358.89	J/molxK	607.47	Joback Method
cpg	373.56	J/molxK	647.14	Joback Method
cpg	387.15	J/molxK	686.82	Joback Method
cpg	399.72	J/molxK	726.49	Joback Method
cpg	411.32	J/molxK	766.17	Joback Method
cpg	421.99	J/molxK	805.84	Joback Method
cpg	431.79	J/molxK	845.52	Joback Method
dvisc	0.0018581	Paxs	369.17	Joback Method

dvisc	0.0012308	Paxs	408.89	Joback Method
dvisc	0.0008770	Paxs	448.60	Joback Method
dvisc	0.0006603	Paxs	488.32	Joback Method
dvisc	0.0005188	Paxs	528.04	Joback Method
dvisc	0.0004216	Paxs	567.75	Joback Method
dvisc	0.0003521	Paxs	607.47	Joback Method

Sources

McGowan Method:	http://link.springer.com/article/10.1007/BF02311772
NIST Webbook:	http://webbook.nist.gov/cgi/cbook.cgi?ID=U299073&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

Legend

cp_g:	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
h_{vap}:	Enthalpy of vaporization at standard conditions
log₁₀ws:	Log ₁₀ of Water solubility in mol/l
log_p:	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

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