

Cyclopropanecarboxylic acid, 2,4,5-trichlorophenyl ester

Inchi:	InChI=1S/C10H7Cl3O2/c11-6-3-8(13)9(4-7(6)12)15-10(14)5-1-2-5/h3-5H,1-2H2
InchiKey:	JTXKTAJEYVQEMR-UHFFFAOYSA-N
Formula:	C10H7Cl3O2
SMILES:	O=C(Oc1cc(Cl)c(Cl)cc1Cl)C1CC1
Mol. weight [g/mol]:	265.52

Physical Properties

Property code	Value	Unit	Source
gf	-92.12	kJ/mol	Joback Method
hf	-266.83	kJ/mol	Joback Method
hfus	28.04	kJ/mol	Joback Method
hvap	64.34	kJ/mol	Joback Method
log10ws	-4.33		Crippen Method
logp	3.962		Crippen Method
mcvol	161.300	ml/mol	McGowan Method
pc	2995.87	kPa	Joback Method
rinsol	1820.00		NIST Webbook
tb	665.14	K	Joback Method
tc	908.07	K	Joback Method
tf	446.30	K	Joback Method
vc	0.616	m ³ /kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	354.10	J/molxK	665.14	Joback Method
cpg	364.59	J/molxK	705.63	Joback Method
cpg	374.28	J/molxK	746.12	Joback Method
cpg	383.20	J/molxK	786.61	Joback Method
cpg	391.42	J/molxK	827.10	Joback Method
cpg	398.97	J/molxK	867.58	Joback Method
cpg	405.92	J/molxK	908.07	Joback Method
dvisc	0.0014407	Paxs	446.30	Joback Method
dvisc	0.0011180	Paxs	482.77	Joback Method

dvisc	0.0008990	Paxs	519.25	Joback Method
dvisc	0.0007439	Paxs	555.72	Joback Method
dvisc	0.0006301	Paxs	592.19	Joback Method
dvisc	0.0005441	Paxs	628.67	Joback Method
dvisc	0.0004775	Paxs	665.14	Joback Method

Sources

McGowan Method:	http://link.springer.com/article/10.1007/BF02311772
NIST Webbook:	http://webbook.nist.gov/cgi/cbook.cgi?ID=U354664&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

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/20-098-1/Cyclopropanecarboxylic-acid-2-4-5-trichlorophenyl-ester.pdf>

Generated by Cheméo on 2024-04-20 15:41:40.566126061 +0000 UTC m=+15916949.486703377.

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