

Cyclopropanecarboxylic acid, 3-(2,2-dichlorovinyl)-2,2-dimethyl-, (3-phenoxyphenyl)methyl ester, (1R-cis)-

Other names:

(+)-cis-Permethrin
(1R)-cis-Permethrin

Permethrin cis

cis-Permethrin

Inchi: InChI=1S/C21H20Cl2O3/c1-21(2)17(12-18(22)23)19(21)20(24)25-13-14-7-6-10-16(11-14)

InchiKey: RLLPVAHGXHCWKJ-HKUYNNGSSA-N

Formula: C21H20Cl2O3

SMILES: CC1(C)C(C=C(Cl)Cl)C1C(=O)OCc1cccc(Oc2ccccc2)c1

Mol. weight [g/mol]: 391.29

CAS: 54774-45-7

Physical Properties

Property code	Value	Unit	Source
gf	89.86	kJ/mol	Joback Method
hf	-268.89	kJ/mol	Joback Method
hfus	43.08	kJ/mol	Joback Method
hvap	86.07	kJ/mol	Joback Method
log10ws	-6.52		Crippen Method
logp	6.113		Crippen Method
mcvol	281.860	ml/mol	McGowan Method
pc	1644.42	kPa	Joback Method
tb	913.47	K	Joback Method
tc	1160.83	K	Joback Method
tf	560.34	K	Joback Method
vc	1.069	m3/kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	837.42	J/molxK	913.47	Joback Method
cpg	855.30	J/molxK	954.70	Joback Method
cpg	872.99	J/molxK	995.92	Joback Method
cpg	890.73	J/molxK	1037.15	Joback Method
cpg	908.75	J/molxK	1078.38	Joback Method

cpg	927.28	J/mol×K	1119.60	Joback Method
cpg	946.56	J/mol×K	1160.83	Joback Method

Sources

NIST Webbook:	http://webbook.nist.gov/cgi/cbook.cgi?ID=C54774457&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
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
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/70-726-8/Cyclopropanecarboxylic-acid-3-2-2-dichlorovinyl-2-2-dimethyl-3-phenoxyphen>

Generated by Cheméo on 2024-04-23 17:13:03.594823068 +0000 UTC m=+16181632.515400379.

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