

Isophthalic acid, 2-chloroethyl ethyl ester

Inchi:	InChI=1S/C12H13ClO4/c1-2-16-11(14)9-4-3-5-10(8-9)12(15)17-7-6-13/h3-5,8H,2,6-7H2,
InchiKey:	JOCXGHTXBALZLR-UHFFFAOYSA-N
Formula:	C12H13ClO4
SMILES:	CCOC(=O)c1cccc(C(=O)OCCCl)c1
Mol. weight [g/mol]:	256.68

Physical Properties

Property code	Value	Unit	Source
gf	-326.83	kJ/mol	Joback Method
hf	-571.29	kJ/mol	Joback Method
hfus	30.26	kJ/mol	Joback Method
hvap	67.94	kJ/mol	Joback Method
log10ws	-2.95		Crippen Method
logp	2.259		Crippen Method
mcvol	183.300	ml/mol	McGowan Method
pc	2502.50	kPa	Joback Method
rinpol	1942.00		NIST Webbook
rinpol	1942.00		NIST Webbook
tb	695.63	K	Joback Method
tc	910.73	K	Joback Method
tf	438.18	K	Joback Method
vc	0.697	m3/kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	461.40	J/molxK	695.63	Joback Method
cpg	473.77	J/molxK	731.48	Joback Method
cpg	485.31	J/molxK	767.33	Joback Method
cpg	496.02	J/molxK	803.18	Joback Method
cpg	505.89	J/molxK	839.03	Joback Method
cpg	514.95	J/molxK	874.88	Joback Method
cpg	523.19	J/molxK	910.73	Joback Method
dvisc	0.0009861	Paxs	438.18	Joback Method

dvisc	0.0006136	Paxs	481.09	Joback Method
dvisc	0.0004127	Paxs	524.00	Joback Method
dvisc	0.0002947	Paxs	566.90	Joback Method
dvisc	0.0002207	Paxs	609.81	Joback Method
dvisc	0.0001717	Paxs	652.72	Joback Method
dvisc	0.0001377	Paxs	695.63	Joback Method

Sources

Joback Method:	https://en.wikipedia.org/wiki/Joback_method
McGowan Method:	http://link.springer.com/article/10.1007/BF02311772
NIST Webbook:	http://webbook.nist.gov/cgi/cbook.cgi?ID=U345868&Units=SI
Crippen Method:	http://pubs.acs.org/doi/abs/10.1021/ci990307l
Crippen Method:	https://www.chemeo.com/doc/models/crippen_log10ws

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/94-200-5/Isophthalic-acid-2-chloroethyl-ethyl-ester.pdf>

Generated by Cheméo on 2024-04-24 22:24:52.672670732 +0000 UTC m=+16286741.593248045.

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