

# Isophthalic acid, 2-chloro-5-methylphenyl isoheptyl ester

<b>Inchi:</b>	InChI=1S/C21H23ClO4/c1-14(2)6-5-11-25-20(23)16-7-4-8-17(13-16)21(24)26-19-12-15(3)
<b>InchiKey:</b>	ANXSWQNHBUZXEO-UHFFFAOYSA-N
<b>Formula:</b>	C21H23ClO4
<b>SMILES:</b>	<chem>Cc1ccc(Cl)c(OC(=O)c2cccc(C(=O)OCCCC(C)C)c2)c1</chem>
<b>Mol. weight [g/mol]:</b>	374.86

## Physical Properties

Property code	Value	Unit	Source
gf	-160.34	kJ/mol	Joback Method
hf	-548.74	kJ/mol	Joback Method
hfus	43.31	kJ/mol	Joback Method
hvap	91.19	kJ/mol	Joback Method
log10ws	-6.81		Crippen Method
logp	5.461		Crippen Method
mvol	286.350	ml/mol	McGowan Method
pc	1533.06	kPa	Joback Method
rinpol	2799.00		NIST Webbook
tb	937.75	K	Joback Method
tc	1168.01	K	Joback Method
tf	576.07	K	Joback Method
vc	1.087	m <sup>3</sup> /kmol	Joback Method

## Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	860.43	J/molxK	937.75	Joback Method
cpg	910.82	J/molxK	1129.64	Joback Method
cpg	903.35	J/molxK	1091.26	Joback Method
cpg	894.60	J/molxK	1052.88	Joback Method
cpg	884.55	J/molxK	1014.50	Joback Method
cpg	873.18	J/molxK	976.13	Joback Method
cpg	917.06	J/molxK	1168.01	Joback Method
dvisc	0.0000400	Paxs	937.75	Joback Method
dvisc	0.0000504	Paxs	877.47	Joback Method

dvisc	0.0000656	Paxs	817.19	Joback Method
dvisc	0.0000891	Paxs	756.91	Joback Method
dvisc	0.0001277	Paxs	696.63	Joback Method
dvisc	0.0001959	Paxs	636.35	Joback Method
dvisc	0.0003287	Paxs	576.07	Joback Method

## Sources

<b>Crippen Method:</b>	<a href="http://pubs.acs.org/doi/abs/10.1021/ci9903071">http://pubs.acs.org/doi/abs/10.1021/ci9903071</a>
<b>Crippen Method:</b>	<a href="https://www.chemeo.com/doc/models/crippen_log10ws">https://www.chemeo.com/doc/models/crippen_log10ws</a>
<b>Joback Method:</b>	<a href="https://en.wikipedia.org/wiki/Joback_method">https://en.wikipedia.org/wiki/Joback_method</a>
<b>McGowan Method:</b>	<a href="http://link.springer.com/article/10.1007/BF02311772">http://link.springer.com/article/10.1007/BF02311772</a>
<b>NIST Webbook:</b>	<a href="http://webbook.nist.gov/cgi/cbook.cgi?ID=U356567&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=U356567&amp;Units=SI</a>

## Legend

<b>cpg:</b>	Ideal gas heat capacity
<b>dvisc:</b>	Dynamic viscosity
<b>gf:</b>	Standard Gibbs free energy of formation
<b>hf:</b>	Enthalpy of formation at standard conditions
<b>hfus:</b>	Enthalpy of fusion at standard conditions
<b>hvap:</b>	Enthalpy of vaporization at standard conditions
<b>log10ws:</b>	Log10 of Water solubility in mol/l
<b>logp:</b>	Octanol/Water partition coefficient
<b>mcvol:</b>	McGowan's characteristic volume
<b>pc:</b>	Critical Pressure
<b>rinpol:</b>	Non-polar retention indices
<b>tb:</b>	Normal Boiling Point Temperature
<b>tc:</b>	Critical Temperature
<b>tf:</b>	Normal melting (fusion) point
<b>vc:</b>	Critical Volume

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