

# 2,2',3,4',5,5',6-Heptachlorodiphenyl ether

**Inchi:** InChI=1S/C12H3Cl7O/c13-4-1-6(15)9(3-5(4)14)20-12-10(18)7(16)2-8(17)11(12)19/h1-3H  
**InchiKey:** BHJCMMSGUFRZLKK-UHFFFAOYSA-N  
**Formula:** C12H3Cl7O  
**SMILES:** Clc1cc(Cl)c(Oc2c(Cl)c(Cl)cc(Cl)c2Cl)cc1Cl  
**Mol. weight [g/mol]:** 411.33

## Physical Properties

Property code	Value	Unit	Source
gf	19.06	kJ/mol	Joback Method
hf	-140.64	kJ/mol	Joback Method
hfus	42.76	kJ/mol	Joback Method
hvap	84.60	kJ/mol	Joback Method
log10ws	-9.05		Aqueous Solubility Prediction Method
logp	8.053		Crippen Method
mcvol	223.970	ml/mol	McGowan Method
pc	2291.52	kPa	Joback Method
tb	846.61	K	Joback Method
tc	1114.35	K	Joback Method
tf	597.15	K	Joback Method
vc	0.853	m <sup>3</sup> /kmol	Joback Method

## Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	436.95	J/molxK	846.61	Joback Method
cpg	443.65	J/molxK	891.23	Joback Method
cpg	449.51	J/molxK	935.86	Joback Method
cpg	454.54	J/molxK	980.48	Joback Method
cpg	458.75	J/molxK	1025.10	Joback Method
cpg	462.15	J/molxK	1069.73	Joback Method
cpg	464.76	J/molxK	1114.35	Joback Method
dvisc	0.0003259	Paxs	597.15	Joback Method
dvisc	0.0002490	Paxs	638.73	Joback Method

dvisc	0.0001966	Paxs	680.30	Joback Method
dvisc	0.0001595	Paxs	721.88	Joback Method
dvisc	0.0001324	Paxs	763.46	Joback Method
dvisc	0.0001120	Paxs	805.03	Joback Method
dvisc	0.0000964	Paxs	846.61	Joback Method

## Sources

**Crippen Method:**

<http://pubs.acs.org/doi/abs/10.1021/ci9903071>

**Joback Method:**

[https://en.wikipedia.org/wiki/Joback\\_method](https://en.wikipedia.org/wiki/Joback_method)

**Aqueous Solubility Prediction Method:**

<http://onschallenge.wikispaces.com/file/view/AqueousDataset002.xlsx/351826032/AqueousDa>

**McGowan Method:**

<http://link.springer.com/article/10.1007/BF02311772>

## 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>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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