

# Bisphenol C

**Other names:**

2,2-Bis(3-methyl-4-hydroxyphenyl)propane  
Phenol, 4,4'-(1-methylethylidene)bis[2-methyl-  
o-Cresol, 4,4'-isopropylidenedi-  
Dicresylolpropane  
Nonox DCP  
2,2-Bis(4-hydroxy-3-methylphenyl)propane  
3,3'-Dimethyldian  
4,4'-Isopropylidenebis[2-methylphenol]  
4,4'-Isopropylidenedi-o-cresol  
o-Cresol, 4,4'-(2,2-propylene)bis-  
3,3'-Dimethylbisphenol A  
NSC 408489

**Inchi:**

InChI=1S/C17H20O2/c1-11-9-13(5-7-15(11)18)17(3,4)14-6-8-16(19)12(2)10-14/h5-10,18

**InchiKey:**

YMTYZTXUZLQUSF-UHFFFAOYSA-N

**Formula:**

C17H20O2

**SMILES:**

Cc1cc(C(C)(C)c2ccc(O)c(C)c2)ccc1O

**Mol. weight [g/mol]:**

256.34

**CAS:**

79-97-0

## Physical Properties

Property code	Value	Unit	Source
gf	-8.58	kJ/mol	Joback Method
hf	-307.46	kJ/mol	Joback Method
hfus	31.24	kJ/mol	Joback Method
hvap	84.04	kJ/mol	Joback Method
log10ws	-4.13		Crippen Method
logp	4.041		Crippen Method
mcvol	214.610	ml/mol	McGowan Method
pc	2767.17	kPa	Joback Method
tb	809.69	K	Joback Method
tc	1064.09	K	Joback Method
tf	585.09	K	Joback Method
vc	0.693	m <sup>3</sup> /kmol	Joback Method

# Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	633.88	J/molxK	809.69	Joback Method
cpg	649.15	J/molxK	852.09	Joback Method
cpg	663.85	J/molxK	894.49	Joback Method
cpg	678.23	J/molxK	936.89	Joback Method
cpg	692.53	J/molxK	979.29	Joback Method
cpg	706.98	J/molxK	1021.69	Joback Method
cpg	721.84	J/molxK	1064.09	Joback Method
dvisc	0.0000113	Paxs	585.09	Joback Method
dvisc	0.0000054	Paxs	622.52	Joback Method
dvisc	0.0000028	Paxs	659.96	Joback Method
dvisc	0.0000016	Paxs	697.39	Joback Method
dvisc	0.0000009	Paxs	734.82	Joback Method
dvisc	0.0000006	Paxs	772.26	Joback Method
dvisc	0.0000004	Paxs	809.69	Joback Method

# Pressure Dependent Properties

Property code	Value	Unit	Pressure [kPa]	Source
tbrp	514.20	K	1.60	NIST Webbook

## Sources

McGowan Method:

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

NIST Webbook:

<http://webbook.nist.gov/cgi/cbook.cgi?ID=C79970&Units=SI>

Crippen Method:

<http://pubs.acs.org/doi/abs/10.1021/ci990307I>

Crippen Method:

[https://www.chemeo.com/doc/models/crippen\\_log10ws](https://www.chemeo.com/doc/models/crippen_log10ws)

Joback Method:

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

## 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>tbrp:</b>	Boiling point at reduced pressure
<b>tc:</b>	Critical Temperature
<b>tf:</b>	Normal melting (fusion) point
<b>vc:</b>	Critical Volume

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