

Benzophenone, 2,4-dihydroxy-4'-methyl

Other names:	2,4-dihydroxy-4'-methylbenzophenone
Inchi:	InChI=1S/C14H12O3/c1-9-2-4-10(5-3-9)14(17)12-7-6-11(15)8-13(12)16/h2-8,15-16H,1H
InchiKey:	FGCATPMSNCVLQK-UHFFFAOYSA-N
Formula:	C14H12O3
SMILES:	<chem>Cc1ccc(C(=O)c2ccc(O)cc2O)cc1</chem>
Mol. weight [g/mol]:	228.24
CAS:	40444-43-7

Physical Properties

Property code	Value	Unit	Source
gf	-155.97	kJ/mol	Joback Method
hf	-337.90	kJ/mol	Joback Method
hfus	32.87	kJ/mol	Joback Method
hvap	84.75	kJ/mol	Joback Method
log10ws	-3.00		Crippen Method
logp	2.637		Crippen Method
mcvol	173.910	ml/mol	McGowan Method
pc	4140.93	kPa	Joback Method
tb	793.17	K	Joback Method
tc	1053.98	K	Joback Method
tf	586.27	K	Joback Method
vc	0.541	m3/kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	481.65	J/molxK	793.17	Joback Method
cpg	493.46	J/molxK	836.64	Joback Method
cpg	504.80	J/molxK	880.11	Joback Method
cpg	515.87	J/molxK	923.58	Joback Method
cpg	526.90	J/molxK	967.05	Joback Method
cpg	538.11	J/molxK	1010.52	Joback Method
cpg	549.70	J/molxK	1053.98	Joback Method
dvisc	0.0000163	Paxs	586.27	Joback Method

dvisc	0.0000085	Paxs	620.75	Joback Method
dvisc	0.0000047	Paxs	655.24	Joback Method
dvisc	0.0000028	Paxs	689.72	Joback Method
dvisc	0.0000017	Paxs	724.20	Joback Method
dvisc	0.0000011	Paxs	758.69	Joback Method
dvisc	0.0000008	Paxs	793.17	Joback Method

Sources

Crippen Method:	http://pubs.acs.org/doi/abs/10.1021/ci9903071
Crippen Method:	https://www.cheméo.com/doc/models/crippen_log10ws
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=C40444437&Units=SI

Legend

cp_g:	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
h_{vap}:	Enthalpy of vaporization at standard conditions
log₁₀ws:	Log ₁₀ of Water solubility in mol/l
log_p:	Octanol/Water partition coefficient
mc_{vol}:	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.cheméo.com/cid/39-025-1/Benzophenone-2-4-dihydroxy-4-methyl.pdf>

Generated by Cheméo on 2024-04-24 09:44:39.537029841 +0000 UTC m=+16241128.457607169.

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