

Benzophenone, 2,4-dihydroxy-4'-phenyl-

Inchi:	InChI=1S/C19H14O3/c20-16-10-11-17(18(21)12-16)19(22)15-8-6-14(7-9-15)13-4-2-1-3-5
InchiKey:	UTLCXVVKPSSBSU-UHFFFAOYSA-N
Formula:	C19H14O3
SMILES:	O=C(c1ccc(-c2ccccc2)cc1)c1ccc(O)cc1O
Mol. weight [g/mol]:	290.31
CAS:	36130-57-1

Physical Properties

Property code	Value	Unit	Source
gf	-1.46	kJ/mol	Joback Method
hf	-204.57	kJ/mol	Joback Method
hfus	39.87	kJ/mol	Joback Method
hvap	98.15	kJ/mol	Joback Method
log10ws	-5.13		Crippen Method
logp	3.996		Crippen Method
mvol	220.600	ml/mol	McGowan Method
pc	3448.03	kPa	Joback Method
tb	934.25	K	Joback Method
tc	1210.58	K	Joback Method
tf	669.04	K	Joback Method
vc	0.714	m3/kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	654.84	J/molxK	934.25	Joback Method
cpg	668.82	J/molxK	980.31	Joback Method
cpg	682.85	J/molxK	1026.36	Joback Method
cpg	697.23	J/molxK	1072.42	Joback Method
cpg	712.26	J/molxK	1118.47	Joback Method
cpg	728.25	J/molxK	1164.53	Joback Method
cpg	745.51	J/molxK	1210.58	Joback Method
dvisc	0.0000042	Paxs	669.04	Joback Method
dvisc	0.0000021	Paxs	713.24	Joback Method

dvisc	0.0000012	Paxs	757.44	Joback Method
dvisc	0.0000007	Paxs	801.64	Joback Method
dvisc	0.0000004	Paxs	845.85	Joback Method
dvisc	0.0000003	Paxs	890.05	Joback Method
dvisc	0.0000002	Paxs	934.25	Joback Method

Sources

Crippen Method:	http://pubs.acs.org/doi/abs/10.1021/ci9903071
Crippen Method:	https://www.chemeo.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=C36130571&Units=SI

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
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/78-331-8/Benzophenone-2-4-dihydroxy-4-phenyl.pdf>

Generated by Cheméo on 2024-04-25 17:25:29.441677906 +0000 UTC m=+16355178.362255227.

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