

4-Hydroxy-2-methylbenzaldehyde

Other names:	Benzaldehyde, 4-hydroxy-2-methyl-
Inchi:	InChI=1S/C8H8O2/c1-6-4-8(10)3-2-7(6)5-9/h2-5,10H,1H3
InchiKey:	JDWWIEFMFPWBST-UHFFFAOYSA-N
Formula:	C8H8O2
SMILES:	<chem>Cc1cc(O)ccc1C=O</chem>
Mol. weight [g/mol]:	136.15
CAS:	41438-18-0

Physical Properties

Property code	Value	Unit	Source
gf	-134.88	kJ/mol	Joback Method
hf	-246.28	kJ/mol	Joback Method
hfus	18.20	kJ/mol	Joback Method
hvap	56.07	kJ/mol	Joback Method
log10ws	-1.74		Crippen Method
logp	1.513		Crippen Method
mcvol	107.260	ml/mol	McGowan Method
pc	4775.98	kPa	Joback Method
tb	543.38	K	Joback Method
tc	774.17	K	Joback Method
tf	372.58	K	Joback Method
vc	0.358	m3/kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	238.56	J/molxK	543.38	Joback Method
cpg	281.20	J/molxK	735.71	Joback Method
cpg	273.85	J/molxK	697.24	Joback Method
cpg	265.98	J/molxK	658.78	Joback Method
cpg	257.53	J/molxK	620.31	Joback Method
cpg	248.41	J/molxK	581.85	Joback Method
cpg	288.11	J/molxK	774.17	Joback Method
dvisc	0.0000748	Paxs	543.38	Joback Method

dvisc	0.0001096	Paxs	514.91	Joback Method
dvisc	0.0001678	Paxs	486.45	Joback Method
dvisc	0.0002710	Paxs	457.98	Joback Method
dvisc	0.0004664	Paxs	429.51	Joback Method
dvisc	0.0008670	Paxs	401.05	Joback Method
dvisc	0.0017717	Paxs	372.58	Joback Method

Sources

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=C41438180&Units=SI
Crippen Method:	http://pubs.acs.org/doi/abs/10.1021/ci990307l
Crippen Method:	https://www.chemeo.com/doc/models/crippen_log10ws

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/81-321-5/4-Hydroxy-2-methylbenzaldehyde.pdf>

Generated by Cheméo on 2024-04-25 16:29:08.996673672 +0000 UTC m=+16351797.917250988.

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