2-Hydroxy-5-methoxybenzaldehyde, acetate

Inchi: InChI=1S/C10H10O4/c1-7(12)14-10-4-3-9(13-2)5-8(10)6-11/h3-6H,1-2H3
InchiKey: AJNKRQSECUQOPV-UHFFFAOYSA-N
Formula: C10H10O4
SMILES: COc1ccc(OC(C)=O)c(C=O)c1
Mol. weight [g/mol]: 194.18

Physical Properties

Property code	Value	Unit	Source
gf	-311.97	kJ/mol	Joback Method
hf	-498.74	kJ/mol	Joback Method
hfus	21.18	kJ/mol	Joback Method
hvap	59.74	kJ/mol	Joback Method
log10ws	-2.15		Crippen Method
logp	1.433		Crippen Method
mcvol	142.880	ml/mol	McGowan Method
pc	3177.55	kPa	Joback Method
rinpol	1558.80		NIST Webbook
rinpol	1558.80		NIST Webbook
tb	612.21	K	Joback Method
tc	827.41	K	Joback Method
tf	390.31	K	Joback Method
vc	0.546	m3/kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	338.79	J/mol×K	612.21	Joback Method
cpg	350.19	J/mol×K	648.08	Joback Method
cpg	360.97	J/mol×K	683.94	Joback Method
cpg	371.12	J/mol×K	719.81	Joback Method
cpg	380.62	J/mol×K	755.68	Joback Method
cpg	389.47	J/mol×K	791.54	Joback Method
cpg	397.65	J/mol×K	827.41	Joback Method
dvisc	0.0011188	Pa×s	390.31	Joback Method

dvisc	0.0007397	Pa×s	427.29	Joback Method
dvisc	0.0005224	Pa×s	464.28	Joback Method
dvisc	0.0003883	Pa×s	501.26	Joback Method
dvisc	0.0003007	Pa×s	538.24	Joback Method
dvisc	0.0002406	Pa×s	575.23	Joback Method
dvisc	0.0001978	Pa×s	612.21	Joback Method

Sources

Crippen Method:	http://pubs.acs.org/doi/abs/10.1021/ci990307l
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=U352898&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
rinpol:	Non-polar retention indices
tb:	Normal Boiling Point Temperature
tc:	Critical Temperature
tf:	Normal melting (fusion) point
vc:	Critical Volume