4-Acetoxybenzaldehyde

Other names:	p-Acetoxybenzaldehyde
	Benzaldehyde, 4-(acetyloxy)-
	p-Hydroxybenzaldehyde acetate
	Benzaldehyde, p-hydroxy-, acetate
	4-formylphenyl acetate
Inchi:	InChI=1S/C9H8O3/c1-7(11)12-9-4-2-8(6-10)3-5-9/h2-6H,1H3
InchiKey:	SEVSMVUOKAMPDO-UHFFFAOYSA-N
Formula:	C9H8O3
SMILES:	CC(=O)Oc1ccc(C=O)cc1
Mol. weight [g/mol]:	164.16
CAS:	878-00-2

Physical Properties

Property code	Value	Unit	Source
gf	-205.76	kJ/mol	Joback Method
hf	-334.41	kJ/mol	Joback Method
hfus	17.79	kJ/mol	Joback Method
hvap	54.44	kJ/mol	Joback Method
log10ws	-2.03		Crippen Method
logp	1.424		Crippen Method
mcvol	122.920	ml/mol	McGowan Method
pc	3690.97	kPa	Joback Method
rinpol	1294.00		NIST Webbook
rinpol	228.40		NIST Webbook
tb	561.93	K	Joback Method
tc	782.61	K	Joback Method
tf	344.29	K	Joback Method
vc	0.472	m3/kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	271.43	J/mol×K	561.93	Joback Method
cpg	282.25	J/mol×K	598.71	Joback Method

cpg	292.43	J/mol×K	635.49	Joback Method
cpg	301.97	J/mol×K	672.27	Joback Method
cpg	310.88	J/mol×K	709.05	Joback Method
cpg	319.18	J/mol×K	745.83	Joback Method
cpg	326.87	J/mol×K	782.61	Joback Method
dvisc	0.0018131	Pa×s	344.29	Joback Method
dvisc	0.0011322	Pa×s	380.56	Joback Method
dvisc	0.0007674	Pa×s	416.84	Joback Method
dvisc	0.0005535	Pa×s	453.11	Joback Method
dvisc	0.0004191	Pa×s	489.38	Joback Method
dvisc	0.0003297	Pa×s	525.66	Joback Method
dvisc	0.0002676	Pa×s	561.93	Joback Method

Pressure Dependent Properties

Property code	Value	Unit	Pressure [kPa]	Source
tbrp	425.70	K	2.30	NIST Webbook
tbrp	444.00 ± 1.00	K	4.70	NIST Webbook

Sources

McGowan Method: http://link.springer.com/article/10.1007/BF02311772
NIST Webbook: http://webbook.nist.gov/cgi/cbook.cgi?ID=C878002&Units=SI
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

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
tbrp:	Boiling point at reduced pressure
tc:	Critical Temperature
tf:	Normal melting (fusion) point
vc:	Critical Volume