ortho-Formylphenoxyacetic acid

Other names:	2-Formylphenoxyacetic acid
	Acetic acid, (2-formylphenoxy)-
	o-Formylphenoxyacetic acid
	Acetic acid, (o-formylphenoxy)-
	2-(2-Formylphenoxy)acetic acid
	2-Carboxymethoxybenzaldehyde
Inchi:	InChI=1S/C9H8O4/c10-5-7-3-1-2-4-8(7)13-6-9(11)12/h1-5H,6H2,(H,11,12)
InchiKey:	ANWMNLAAFDCKMT-UHFFFAOYSA-N
Formula:	C9H8O4
SMILES:	O=Cc1ccccc1OCC(=O)O
Mol. weight [g/mol]:	180.16
CAS:	6280-80-4

Physical Properties

Property code	Value	Unit	Source
gf	-342.58	kJ/mol	Joback Method
hf	-486.64	kJ/mol	Joback Method
hfus	21.88	kJ/mol	Joback Method
hvap	71.12	kJ/mol	Joback Method
log10ws	-1.35		Crippen Method
logp	0.962		Crippen Method
mcvol	128.790	ml/mol	McGowan Method
pc	4140.93	kPa	Joback Method
tb	654.11	K	Joback Method
tc	859.10	K	Joback Method
tf	405.11	K	Joback Method
vc	0.491	m3/kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	314.59	J/mol×K	654.11	Joback Method
cpg	323.40	J/mol×K	688.27	Joback Method
cpg	331.64	J/mol×K	722.44	Joback Method

cpg	339.34	J/mol×K	756.60	Joback Method
cpg	346.49	J/mol×K	790.77	Joback Method
cpg	353.11	J/mol×K	824.93	Joback Method
cpg	359.21	J/mol×K	859.10	Joback Method
dvisc	0.0019234	Pa×s	405.11	Joback Method
dvisc	0.0008776	Pa×s	446.61	Joback Method
dvisc	0.0004576	Pa×s	488.11	Joback Method
dvisc	0.0002643	Pa×s	529.61	Joback Method
dvisc	0.0001653	Pa×s	571.11	Joback Method
dvisc	0.0001102	Pa×s	612.61	Joback Method
dvisc	0.0000773	Pa×s	654.11	Joback Method

Sources

McGowan Method:	http://link.springer.com/article/10.1007/BF02311772
NIST Webbook:	http://webbook.nist.gov/cgi/cbook.cgi?ID=C6280804&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
tb:	Normal Boiling Point Temperature
tc:	Critical Temperature
tf:	Normal melting (fusion) point
vc:	Critical Volume

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