3,5-Dimethoxy-4-hydroxycinnamaldehyde

Other names:	3-(4-Hydroxy-3,5-dimethoxyphenyl)-2-propenal
	2-Propenal, 3-(4-hydroxy-3,5-dimethoxyphenyl)
	Sinapic aldehyde
	Sinapinaldehyde
	2-Propenal, 3-(4-hydroxy-3,5-dimethoxyphenyl), (E)-
Inchi:	InChI=1S/C11H12O4/c1-14-9-6-8(4-3-5-12)7-10(15-2)11(9)13/h3-7,13H,1-2H3/b4-3+
InchiKey:	CDICDSOGTRCHMG-ONEGZZNKSA-N
Formula:	C11H12O4
SMILES:	COc1cc(C=CC=O)cc(OC)c1O
Mol. weight [g/mol]:	208.21
CAS:	87345-53-7

Physical Properties

Property code	Value	Unit	Source
gf	-249.03	kJ/mol	Joback Method
hf	-466.89	kJ/mol	Joback Method
hfus	28.16	kJ/mol	Joback Method
hvap	68.19	kJ/mol	Joback Method
log10ws	-1.78		Crippen Method
logp	1.621		Crippen Method
mcvol	156.970	ml/mol	McGowan Method
pc	3333.53	kPa	Joback Method
rinpol	1909.00		NIST Webbook
rinpol	1915.00		NIST Webbook
ripol	3458.00		NIST Webbook
tb	666.00	K	Joback Method
tc	888.57	K	Joback Method
tf	458.29	K	Joback Method
vc	0.542	m3/kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	398.69	J/mol×K	666.00	Joback Method

cpg	410.06	J/mol×K	703.10	Joback Method
cpg	420.78	J/mol×K	740.19	Joback Method
cpg	430.90	J/mol×K	777.29	Joback Method
cpg	440.47	J/mol×K	814.38	Joback Method
cpg	449.53	J/mol×K	851.48	Joback Method
cpg	458.15	J/mol×K	888.57	Joback Method
dvisc	0.0002831	Pa×s	458.29	Joback Method
dvisc	0.0001487	Pa×s	492.91	Joback Method
dvisc	0.0000850	Pa×s	527.53	Joback Method
dvisc	0.0000520	Pa×s	562.14	Joback Method
dvisc	0.0000337	Pa×s	596.76	Joback Method
dvisc	0.0000229	Pa×s	631.38	Joback Method
dvisc	0.0000162	Pa×s	666.00	Joback Method

Sources

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=C87345537&Units=SI
Crippen Method:	http://pubs.acs.org/doi/abs/10.1021/ci990307l

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
ripol:	Polar retention indices
tb:	Normal Boiling Point Temperature
tc:	Critical Temperature

tf:	Normal melting (fusion) point
vc:	Critical Volume

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