

2-Propenoic acid, 3-(4-methoxyphenyl)-, (E)-

Other names:	trans-2-Propenoic acid, 3-(4-methoxyphenyl)- (E)-p-Methoxy-cinnamic acid trans-4-methoxycinnamic acid
Inchi:	InChI=1S/C10H10O3/c1-13-9-5-2-8(3-6-9)4-7-10(11)12/h2-7H,1H3,(H,11,12)/b7-4+
InchiKey:	AFDXODALSZRGIH-QPJJXVBHSA-N
Formula:	C10H10O3
SMILES:	COc1ccc(C=CC(=O)O)cc1
Mol. weight [g/mol]:	178.18
CAS:	943-89-5

Physical Properties

Property code	Value	Unit	Source
gf	-154.42	kJ/mol	Joback Method
hf	-304.48	kJ/mol	Joback Method
hfus	22.39	kJ/mol	Joback Method
hsub	134.00 ± 1.00	kJ/mol	NIST Webbook
hvap	66.59	kJ/mol	Joback Method
log10ws	-1.93		Crippen Method
logp	1.793		Crippen Method
mcvol	137.010	ml/mol	McGowan Method
pc	3577.07	kPa	Joback Method
rinpol	1733.00		NIST Webbook
rinpol	1733.00		NIST Webbook
tb	632.49	K	Joback Method
tc	839.79	K	Joback Method
tf	369.30	K	Joback Method
vc	0.510	m3/kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	328.56	J/mol×K	632.49	Joback Method
cpg	338.85	J/mol×K	667.04	Joback Method
cpg	348.51	J/mol×K	701.59	Joback Method

cpg	357.56	J/mol×K	736.14	Joback Method
cpg	366.02	J/mol×K	770.69	Joback Method
cpg	373.93	J/mol×K	805.24	Joback Method
cpg	381.31	J/mol×K	839.79	Joback Method
dvisc	0.0023160	Pa×s	369.30	Joback Method
dvisc	0.0008973	Pa×s	413.17	Joback Method
dvisc	0.0004170	Pa×s	457.03	Joback Method
dvisc	0.0002217	Pa×s	500.89	Joback Method
dvisc	0.0001304	Pa×s	544.76	Joback Method
dvisc	0.0000831	Pa×s	588.62	Joback Method
dvisc	0.0000563	Pa×s	632.49	Joback Method

Sources

NIST Webbook:	http://webbook.nist.gov/cgi/cbook.cgi?ID=C943895&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
McGowan Method:	http://link.springer.com/article/10.1007/BF02311772

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
hsub:	Enthalpy of sublimation 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

Latest version available from:

<https://www.cheméo.com/cid/85-955-8/2-Propenoic-acid-3-4-methoxyphenyl-E.pdf>

Generated by Cheméo on 2025-12-05 13:54:22.447373683 +0000 UTC m=+4691059.977414358.

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