

# m-Methoxybenzoic acid, hexyl ester

<b>Inchi:</b>	InChI=1S/C14H20O3/c1-3-4-5-6-10-17-14(15)12-8-7-9-13(11-12)16-2/h7-9,11H,3-6,10H
<b>InchiKey:</b>	ROTBPRMBAFUZES-UHFFFAOYSA-N
<b>Formula:</b>	C14H20O3
<b>SMILES:</b>	CCCCCCOC(=O)c1cccc(OC)c1
<b>Mol. weight [g/mol]:</b>	236.31

## Physical Properties

Property code	Value	Unit	Source
gf	-169.14	kJ/mol	Joback Method
hf	-484.25	kJ/mol	Joback Method
hfus	29.64	kJ/mol	Joback Method
hvap	61.26	kJ/mol	Joback Method
log10ws	-3.93		Crippen Method
logp	3.432		Crippen Method
mcvol	197.670	ml/mol	McGowan Method
pc	2034.55	kPa	Joback Method
rinpol	1811.80		NIST Webbook
rinpol	1811.80		NIST Webbook
tb	650.09	K	Joback Method
tc	849.38	K	Joback Method
tf	380.87	K	Joback Method
vc	0.753	m <sup>3</sup> /kmol	Joback Method

## Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	521.01	J/molxK	650.09	Joback Method
cpg	536.82	J/molxK	683.31	Joback Method
cpg	551.77	J/molxK	716.52	Joback Method
cpg	565.89	J/molxK	749.74	Joback Method
cpg	579.17	J/molxK	782.95	Joback Method
cpg	591.63	J/molxK	816.17	Joback Method
cpg	603.27	J/molxK	849.38	Joback Method
dvisc	0.0011604	Paxs	380.87	Joback Method

dvisc	0.0006488	Paxs	425.74	Joback Method
dvisc	0.0004053	Paxs	470.61	Joback Method
dvisc	0.0002748	Paxs	515.48	Joback Method
dvisc	0.0001982	Paxs	560.35	Joback Method
dvisc	0.0001501	Paxs	605.22	Joback Method
dvisc	0.0001181	Paxs	650.09	Joback Method

## Sources

<b>Crippen Method:</b>	<a href="http://pubs.acs.org/doi/abs/10.1021/ci9903071">http://pubs.acs.org/doi/abs/10.1021/ci9903071</a>
<b>Crippen Method:</b>	<a href="https://www.chemeo.com/doc/models/crippen_log10ws">https://www.chemeo.com/doc/models/crippen_log10ws</a>
<b>Joback Method:</b>	<a href="https://en.wikipedia.org/wiki/Joback_method">https://en.wikipedia.org/wiki/Joback_method</a>
<b>McGowan Method:</b>	<a href="http://link.springer.com/article/10.1007/BF02311772">http://link.springer.com/article/10.1007/BF02311772</a>
<b>NIST Webbook:</b>	<a href="http://webbook.nist.gov/cgi/cbook.cgi?ID=U292367&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=U292367&amp;Units=SI</a>

## Legend

<b>cpg:</b>	Ideal gas heat capacity
<b>dvisc:</b>	Dynamic viscosity
<b>gf:</b>	Standard Gibbs free energy of formation
<b>hf:</b>	Enthalpy of formation at standard conditions
<b>hfus:</b>	Enthalpy of fusion at standard conditions
<b>hvap:</b>	Enthalpy of vaporization at standard conditions
<b>log10ws:</b>	Log10 of Water solubility in mol/l
<b>logp:</b>	Octanol/Water partition coefficient
<b>mcvol:</b>	McGowan's characteristic volume
<b>pc:</b>	Critical Pressure
<b>rinpol:</b>	Non-polar retention indices
<b>tb:</b>	Normal Boiling Point Temperature
<b>tc:</b>	Critical Temperature
<b>tf:</b>	Normal melting (fusion) point
<b>vc:</b>	Critical Volume

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