

# Dodecanoic acid, 2-methyl-, methyl ester

<b>Other names:</b>	Methyl 2-methyldodecanoate
<b>Inchi:</b>	InChI=1S/C14H28O2/c1-4-5-6-7-8-9-10-11-12-13(2)14(15)16-3/h13H,4-12H2,1-3H3
<b>InchiKey:</b>	IMZMJDATCHRVSX-UHFFFAOYSA-N
<b>Formula:</b>	C14H28O2
<b>SMILES:</b>	CCCCCCCCCCC(C)C(=O)OC
<b>Mol. weight [g/mol]:</b>	228.37
<b>CAS:</b>	55554-08-0

## Physical Properties

Property code	Value	Unit	Source
gf	-169.36	kJ/mol	Joback Method
hf	-582.37	kJ/mol	Joback Method
hfus	31.28	kJ/mol	Joback Method
hvap	55.53	kJ/mol	Joback Method
log10ws	-4.30		Crippen Method
logp	4.326		Crippen Method
mcvol	215.560	ml/mol	McGowan Method
pc	1589.81	kPa	Joback Method
tb	523.00 ± 5.00	K	NIST Webbook
tc	766.26	K	Joback Method
tf	304.70	K	Joback Method
vc	0.838	m <sup>3</sup> /kmol	Joback Method

## Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	567.35	J/mol×K	595.57	Joback Method
cpg	646.34	J/mol×K	737.81	Joback Method
cpg	631.92	J/mol×K	709.36	Joback Method
cpg	616.82	J/mol×K	680.91	Joback Method
cpg	601.04	J/mol×K	652.47	Joback Method
cpg	584.55	J/mol×K	624.02	Joback Method
cpg	660.10	J/mol×K	766.26	Joback Method
dvisc	0.0001432	Paxs	595.57	Joback Method

dvisc	0.0001938	Paxs	547.09	Joback Method
dvisc	0.0002782	Paxs	498.61	Joback Method
dvisc	0.0004318	Paxs	450.13	Joback Method
dvisc	0.0007451	Paxs	401.66	Joback Method
dvisc	0.0014934	Paxs	353.18	Joback Method
dvisc	0.0037349	Paxs	304.70	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=C55554080&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=C55554080&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>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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