

# Malonic acid, butyl 2-methylpentyl ester

<b>Inchi:</b>	InChI=1S/C13H24O4/c1-4-6-8-16-12(14)9-13(15)17-10-11(3)7-5-2/h11H,4-10H2,1-3H3
<b>InchiKey:</b>	GIZWADKCVVPYSE-UHFFFAOYSA-N
<b>Formula:</b>	C13H24O4
<b>SMILES:</b>	CCCCOC(=O)CC(=O)OCC(C)CCC
<b>Mol. weight [g/mol]:</b>	244.33

## Physical Properties

Property code	Value	Unit	Source
gf	-411.70	kJ/mol	Joback Method
hf	-806.53	kJ/mol	Joback Method
hfus	31.48	kJ/mol	Joback Method
hvap	62.46	kJ/mol	Joback Method
log10ws	-2.75		Crippen Method
logp	2.699		Crippen Method
mvol	208.910	ml/mol	McGowan Method
pc	1786.37	kPa	Joback Method
rinpol	1529.00		NIST Webbook
rinpol	1529.00		NIST Webbook
tb	648.98	K	Joback Method
tc	828.08	K	Joback Method
tf	365.59	K	Joback Method
vc	0.805	m <sup>3</sup> /kmol	Joback Method

## Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	568.82	J/molxK	648.98	Joback Method
cpg	584.20	J/molxK	678.83	Joback Method
cpg	598.88	J/molxK	708.68	Joback Method
cpg	612.86	J/molxK	738.53	Joback Method
cpg	626.13	J/molxK	768.38	Joback Method
cpg	638.71	J/molxK	798.23	Joback Method
cpg	650.59	J/molxK	828.08	Joback Method
dvisc	0.0019834	Paxs	365.59	Joback Method

dvisc	0.0009630	Paxs	412.82	Joback Method
dvisc	0.0005423	Paxs	460.05	Joback Method
dvisc	0.0003399	Paxs	507.29	Joback Method
dvisc	0.0002306	Paxs	554.52	Joback Method
dvisc	0.0001663	Paxs	601.75	Joback Method
dvisc	0.0001258	Paxs	648.98	Joback Method

## Sources

<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=U349300&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=U349300&amp;Units=SI</a>
<b>Crippen Method:</b>	<a href="http://pubs.acs.org/doi/abs/10.1021/ci990307l">http://pubs.acs.org/doi/abs/10.1021/ci990307l</a>
<b>Crippen Method:</b>	<a href="https://www.chemeo.com/doc/models/crippen_log10ws">https://www.chemeo.com/doc/models/crippen_log10ws</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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