

# Sebacic acid, decyl 3-oxobut-2-yl ester

<b>Inchi:</b>	InChI=1S/C24H44O5/c1-4-5-6-7-8-11-14-17-20-28-23(26)18-15-12-9-10-13-16-19-24(27)
<b>InchiKey:</b>	DOHVYBJGMQNNIU-UHFFFAOYSA-N
<b>Formula:</b>	C24H44O5
<b>SMILES:</b>	CCCCCCCCCOC(=O)CCCCCCCC(=O)OC(C)C(C)=O
<b>Mol. weight [g/mol]:</b>	412.60

## Physical Properties

Property code	Value	Unit	Source
gf	-448.00	kJ/mol	Joback Method
hf	-1146.15	kJ/mol	Joback Method
hfus	61.57	kJ/mol	Joback Method
hvap	93.69	kJ/mol	Joback Method
log10ws	-6.99		Crippen Method
logp	6.312		Crippen Method
mvol	365.470	ml/mol	McGowan Method
pc	886.30	kPa	Joback Method
rinpol	2854.00		NIST Webbook
tb	954.53	K	Joback Method
tc	1171.53	K	Joback Method
tf	539.49	K	Joback Method
vc	1.427	m <sup>3</sup> /kmol	Joback Method

## Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	1239.17	J/molxK	954.53	Joback Method
cpg	1257.62	J/molxK	990.70	Joback Method
cpg	1274.50	J/molxK	1026.86	Joback Method
cpg	1289.85	J/molxK	1063.03	Joback Method
cpg	1303.71	J/molxK	1099.20	Joback Method
cpg	1316.12	J/molxK	1135.36	Joback Method
cpg	1327.12	J/molxK	1171.53	Joback Method
dvisc	0.0004535	Paxs	539.49	Joback Method
dvisc	0.0002128	Paxs	608.66	Joback Method

dvisc	0.0001166	Paxs	677.84	Joback Method
dvisc	0.0000714	Paxs	747.01	Joback Method
dvisc	0.0000475	Paxs	816.18	Joback Method
dvisc	0.0000337	Paxs	885.36	Joback Method
dvisc	0.0000251	Paxs	954.53	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=U355782&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=U355782&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>rinpolar:</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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