

# Sebacic acid, 3-methylphenyl octyl ester

<b>Inchi:</b>	InChI=1S/C25H40O4/c1-3-4-5-6-11-14-20-28-24(26)18-12-9-7-8-10-13-19-25(27)29-23-
<b>InchiKey:</b>	LLASFIGGWUYVGP-UHFFFAOYSA-N
<b>Formula:</b>	C25H40O4
<b>SMILES:</b>	CCCCCCCCOC(=O)CCCCCCCC(=O)Oc1cccc(C)c1
<b>Mol. weight [g/mol]:</b>	404.58

## Physical Properties

Property code	Value	Unit	Source
gf	-205.44	kJ/mol	Joback Method
hf	-823.87	kJ/mol	Joback Method
hfus	59.73	kJ/mol	Joback Method
hvap	92.49	kJ/mol	Joback Method
log10ws	-7.82		Crippen Method
logp	6.925		Crippen Method
mcvol	354.230	ml/mol	McGowan Method
pc	970.49	kPa	Joback Method
rinsol	3069.00		NIST Webbook
tb	955.64	K	Joback Method
tc	1169.97	K	Joback Method
tf	554.77	K	Joback Method
vc	1.375	m <sup>3</sup> /kmol	Joback Method

## Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	1181.32	J/molxK	955.64	Joback Method
cpg	1253.73	J/molxK	1134.25	Joback Method
cpg	1241.97	J/molxK	1098.53	Joback Method
cpg	1228.89	J/molxK	1062.81	Joback Method
cpg	1214.46	J/molxK	1027.08	Joback Method
cpg	1198.61	J/molxK	991.36	Joback Method
cpg	1264.21	J/molxK	1169.97	Joback Method
dvisc	0.0000265	Paxs	955.64	Joback Method
dvisc	0.0000346	Paxs	888.83	Joback Method

dvisc	0.0000471	Paxs	822.02	Joback Method
dvisc	0.0000678	Paxs	755.20	Joback Method
dvisc	0.0001048	Paxs	688.39	Joback Method
dvisc	0.0001777	Paxs	621.58	Joback Method
dvisc	0.0003423	Paxs	554.77	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=U354930&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=U354930&amp;Units=SI</a>
<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.cheméo.com/doc/models/crippen_log10ws">https://www.cheméo.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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