

# Sebacic acid, hexyl 4-methoxybenzyl ester

<b>Inchi:</b>	InChI=1S/C24H38O5/c1-3-4-5-12-19-28-23(25)13-10-8-6-7-9-11-14-24(26)29-20-21-15-
<b>InchiKey:</b>	MTHPBINMLLLNQA-UHFFFAOYSA-N
<b>Formula:</b>	C24H38O5
<b>SMILES:</b>	CCCCCOC(=O)CCCCCCCC(=O)OCc1ccc(OC)cc1
<b>Mol. weight [g/mol]:</b>	406.56

## Physical Properties

Property code	Value	Unit	Source
gf	-318.86	kJ/mol	Joback Method
hf	-935.45	kJ/mol	Joback Method
hfus	58.33	kJ/mol	Joback Method
hvap	92.68	kJ/mol	Joback Method
log10ws	-6.89		Crippen Method
logp	5.983		Crippen Method
mcvol	346.010	ml/mol	McGowan Method
pc	1020.08	kPa	Joback Method
rinqol	3111.00		NIST Webbook
tb	955.18	K	Joback Method
tc	1169.41	K	Joback Method
tf	565.73	K	Joback Method
vc	1.337	m <sup>3</sup> /kmol	Joback Method

## Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	1149.17	J/molxK	955.18	Joback Method
cpg	1165.55	J/molxK	990.89	Joback Method
cpg	1180.40	J/molxK	1026.59	Joback Method
cpg	1193.74	J/molxK	1062.30	Joback Method
cpg	1205.60	J/molxK	1098.00	Joback Method
cpg	1216.01	J/molxK	1133.71	Joback Method
cpg	1224.99	J/molxK	1169.41	Joback Method
dvisc	0.0002663	Paxs	565.73	Joback Method
dvisc	0.0001432	Paxs	630.64	Joback Method

dvisc	0.0000864	Paxs	695.55	Joback Method
dvisc	0.0000569	Paxs	760.45	Joback Method
dvisc	0.0000400	Paxs	825.36	Joback Method
dvisc	0.0000296	Paxs	890.27	Joback Method
dvisc	0.0000228	Paxs	955.18	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=U354365&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=U354365&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.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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