

# Succinic acid, 3,5-dimethylphenyl 3,4-dimethylphenyl ester

<b>Inchi:</b>	InChI=1S/C20H22O4/c1-13-9-14(2)11-18(10-13)24-20(22)8-7-19(21)23-17-6-5-15(3)16(4)
<b>InchiKey:</b>	JMQFAKAKZBTIGP-UHFFFAOYSA-N
<b>Formula:</b>	C20H22O4
<b>SMILES:</b>	<chem>Cc1cc(C)cc(OC(=O)CCC(=O)Oc2ccc(C)c(C)c2)c1</chem>
<b>Mol. weight [g/mol]:</b>	326.39

## Physical Properties

Property code	Value	Unit	Source
gf	-164.02	kJ/mol	Joback Method
hf	-518.55	kJ/mol	Joback Method
hfus	39.66	kJ/mol	Joback Method
hvap	85.63	kJ/mol	Joback Method
log10ws	-5.65		Crippen Method
logp	4.211		Crippen Method
mcvol	260.020	ml/mol	McGowan Method
pc	1674.16	kPa	Joback Method
rinsol	2659.00		NIST Webbook
tb	882.86	K	Joback Method
tc	1108.90	K	Joback Method
tf	562.40	K	Joback Method
vc	0.988	m <sup>3</sup> /kmol	Joback Method

## Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	778.05	J/molxK	882.86	Joback Method
cpg	834.60	J/molxK	1071.23	Joback Method
cpg	825.84	J/molxK	1033.56	Joback Method
cpg	815.81	J/molxK	995.88	Joback Method
cpg	804.52	J/molxK	958.21	Joback Method
cpg	791.94	J/molxK	920.53	Joback Method
cpg	842.13	J/molxK	1108.90	Joback Method
dvisc	0.0000590	Paxs	882.86	Joback Method
dvisc	0.0000724	Paxs	829.45	Joback Method

dvisc	0.0000912	Paxs	776.04	Joback Method
dvisc	0.0001190	Paxs	722.63	Joback Method
dvisc	0.0001620	Paxs	669.22	Joback Method
dvisc	0.0002326	Paxs	615.81	Joback Method
dvisc	0.0003578	Paxs	562.40	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=U360732&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=U360732&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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