

# Succinic acid, 3,5-dichlorophenyl ethyl ester

<b>Inchi:</b>	InChI=1S/C12H12Cl2O4/c1-2-17-11(15)3-4-12(16)18-10-6-8(13)5-9(14)7-10/h5-7H,2-4H
<b>InchiKey:</b>	JAUNYACIFVMSCA-UHFFFAOYSA-N
<b>Formula:</b>	C12H12Cl2O4
<b>SMILES:</b>	CCOC(=O)CCC(=O)Oc1cc(Cl)cc(Cl)c1
<b>Mol. weight [g/mol]:</b>	291.13

## Physical Properties

Property code	Value	Unit	Source
gf	-348.39	kJ/mol	Joback Method
hf	-598.50	kJ/mol	Joback Method
hfus	34.07	kJ/mol	Joback Method
hvap	72.99	kJ/mol	Joback Method
log10ws	-3.69		Crippen Method
logp	3.242		Crippen Method
mcvol	195.540	ml/mol	McGowan Method
pc	2377.22	kPa	Joback Method
rinpola	1962.00		NIST Webbook
rinpola	1962.00		NIST Webbook
tb	738.04	K	Joback Method
tc	957.03	K	Joback Method
tf	480.62	K	Joback Method
vc	0.746	m <sup>3</sup> /kmol	Joback Method

## Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	483.25	J/molxK	738.04	Joback Method
cpg	494.49	J/molxK	774.54	Joback Method
cpg	504.90	J/molxK	811.04	Joback Method
cpg	514.46	J/molxK	847.53	Joback Method
cpg	523.20	J/molxK	884.03	Joback Method
cpg	531.09	J/molxK	920.53	Joback Method
cpg	538.15	J/molxK	957.03	Joback Method
dvisc	0.0007350	Paxs	480.62	Joback Method

dvisc	0.0004840	Paxs	523.52	Joback Method
dvisc	0.0003395	Paxs	566.43	Joback Method
dvisc	0.0002503	Paxs	609.33	Joback Method
dvisc	0.0001922	Paxs	652.23	Joback Method
dvisc	0.0001524	Paxs	695.14	Joback Method
dvisc	0.0001242	Paxs	738.04	Joback Method

## Sources

<b>Crippen Method:</b>	<a href="https://www.chemeo.com/doc/models/crippen_log10ws">https://www.chemeo.com/doc/models/crippen_log10ws</a>
<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=U349718&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=U349718&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>

## 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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