

# Succinic acid, di(2,2,2-trichloroethyl) ester

**Inchi:** InChI=1S/C8H8Cl6O4/c9-7(10,11)3-17-5(15)1-2-6(16)18-4-8(12,13)14/h1-4H2  
**InchiKey:** KUKFKXHVFUOECA-UHFFFAOYSA-N  
**Formula:** C8H8Cl6O4  
**SMILES:** O=C(CCC(=O)OCC(Cl)(Cl)Cl)OCC(Cl)(Cl)Cl  
**Mol. weight [g/mol]:** 380.87

## Physical Properties

Property code	Value	Unit	Source
gf	-517.26	kJ/mol	Joback Method
hf	-809.99	kJ/mol	Joback Method
hfus	32.40	kJ/mol	Joback Method
hvap	75.43	kJ/mol	Joback Method
log10ws	-4.02		Crippen Method
logp	3.593		Crippen Method
mvol	211.900	ml/mol	McGowan Method
pc	2315.84	kPa	Joback Method
rinpol	1997.00		NIST Webbook
rinpol	1997.00		NIST Webbook
tb	753.14	K	Joback Method
tc	980.64	K	Joback Method
tf	508.60	K	Joback Method
vc	0.803	m <sup>3</sup> /kmol	Joback Method

## Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	463.13	J/molxK	753.14	Joback Method
cpg	470.53	J/molxK	791.06	Joback Method
cpg	477.20	J/molxK	828.97	Joback Method
cpg	483.20	J/molxK	866.89	Joback Method
cpg	488.58	J/molxK	904.81	Joback Method
cpg	493.38	J/molxK	942.72	Joback Method
cpg	497.66	J/molxK	980.64	Joback Method
dvisc	0.0006817	Paxs	508.60	Joback Method

dvisc	0.0004207	Paxs	549.36	Joback Method
dvisc	0.0002775	Paxs	590.11	Joback Method
dvisc	0.0001932	Paxs	630.87	Joback Method
dvisc	0.0001405	Paxs	671.63	Joback Method
dvisc	0.0001060	Paxs	712.38	Joback Method
dvisc	0.0000824	Paxs	753.14	Joback Method

## Sources

<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>
<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=U349181&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=U349181&amp;Units=SI</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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