

# Fumaric acid, 3-heptyl octadecyl ester

<b>Inchi:</b>	InChI=1S/C29H54O4/c1-4-7-9-10-11-12-13-14-15-16-17-18-19-20-21-22-26-32-28(30)24
<b>InchiKey:</b>	HHQRPSYJJSXGLN-OCOZRVBESA-N
<b>Formula:</b>	C29H54O4
<b>SMILES:</b>	CCCCCCCCCCCCCCCCCOC(=O)C=CC(=O)OC(CC)CCCC
<b>Mol. weight [g/mol]:</b>	466.74

## Physical Properties

Property code	Value	Unit	Source
gf	-196.76	kJ/mol	Joback Method
hf	-1019.55	kJ/mol	Joback Method
hfus	73.12	kJ/mol	Joback Method
hvap	98.03	kJ/mol	Joback Method
log10ws	-9.65		Crippen Method
logp	8.859		Crippen Method
mcvol	430.050	ml/mol	McGowan Method
pc	674.65	kPa	Joback Method
rinpola	3174.00		NIST Webbook
tb	1019.22	K	Joback Method
tc	1266.06	K	Joback Method
tf	540.83	K	Joback Method
vc	1.681	m <sup>3</sup> /kmol	Joback Method

## Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	1511.74	J/molxK	1019.22	Joback Method
cpg	1606.58	J/molxK	1224.92	Joback Method
cpg	1591.09	J/molxK	1183.78	Joback Method
cpg	1573.97	J/molxK	1142.64	Joback Method
cpg	1555.11	J/molxK	1101.50	Joback Method
cpg	1534.40	J/molxK	1060.36	Joback Method
cpg	1620.54	J/molxK	1266.06	Joback Method
dvisc	0.0000107	Paxs	1019.22	Joback Method
dvisc	0.0000147	Paxs	939.49	Joback Method

dvisc	0.0000214	Paxs	859.76	Joback Method
dvisc	0.0000336	Paxs	780.02	Joback Method
dvisc	0.0000585	Paxs	700.29	Joback Method
dvisc	0.0001173	Paxs	620.56	Joback Method
dvisc	0.0002886	Paxs	540.83	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=U348695&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=U348695&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>
<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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