

# Fumaric acid, di(naphth-1-yl) ester

<b>Inchi:</b>	InChI=1S/C24H16O4/c25-23(27-21-13-5-9-17-7-1-3-11-19(17)21)15-16-24(26)28-22-14-
<b>InchiKey:</b>	RFNWEPPOTCGGNE-FOCLMDBBSA-N
<b>Formula:</b>	C24H16O4
<b>SMILES:</b>	O=C(C=CC(=O)Oc1cccc2ccccc12)Oc1cccc2ccccc12
<b>Mol. weight [g/mol]:</b>	368.38

## Physical Properties

Property code	Value	Unit	Source
gf	182.44	kJ/mol	Joback Method
hf	-78.81	kJ/mol	Joback Method
hfus	45.03	kJ/mol	Joback Method
hvap	96.44	kJ/mol	Joback Method
log10ws	-7.21		Crippen Method
logp	5.060		Crippen Method
mcvol	273.160	ml/mol	McGowan Method
pc	1964.82	kPa	Joback Method
rinpol	3272.00		NIST Webbook
rinpol	3272.00		NIST Webbook
tb	1006.54	K	Joback Method
tc	1265.54	K	Joback Method
tf	642.76	K	Joback Method
vc	1.036	m <sup>3</sup> /kmol	Joback Method

## Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	819.18	J/molxK	1006.54	Joback Method
cpg	831.34	J/molxK	1049.71	Joback Method
cpg	842.82	J/molxK	1092.87	Joback Method
cpg	853.79	J/molxK	1136.04	Joback Method
cpg	864.43	J/molxK	1179.21	Joback Method
cpg	874.90	J/molxK	1222.37	Joback Method
cpg	885.39	J/molxK	1265.54	Joback Method
dvisc	0.0005651	Paxs	642.76	Joback Method

dvisc	0.0003960	Paxs	703.39	Joback Method
dvisc	0.0002936	Paxs	764.02	Joback Method
dvisc	0.0002274	Paxs	824.65	Joback Method
dvisc	0.0001825	Paxs	885.28	Joback Method
dvisc	0.0001506	Paxs	945.91	Joback Method
dvisc	0.0001272	Paxs	1006.54	Joback Method

## Sources

<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=U405821&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=U405821&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.cheméo.com/doc/models/crippen_log10ws">https://www.cheméo.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>

## 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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