

# Isophthalic acid, nonyl trans-hex-3-enyl ester

<b>Inchi:</b>	InChI=1S/C23H34O4/c1-3-5-7-9-10-11-13-18-27-23(25)21-16-14-15-20(19-21)22(24)26-
<b>InchiKey:</b>	UOINNQINQRTPH-SOFGYWHQSA-N
<b>Formula:</b>	C23H34O4
<b>SMILES:</b>	CCC=CCCOC(=O)c1cccc(C(=O)OCCCCCCCCC)c1
<b>Mol. weight [g/mol]:</b>	374.51

## Physical Properties

Property code	Value	Unit	Source
gf	-142.06	kJ/mol	Joback Method
hf	-665.37	kJ/mol	Joback Method
hfus	54.75	kJ/mol	Joback Method
hvap	88.00	kJ/mol	Joback Method
log10ws	-7.25		Crippen Method
logp	6.107		Crippen Method
mcvol	321.750	ml/mol	McGowan Method
pc	1136.73	kPa	Joback Method
rinpol	2740.00		NIST Webbook
tb	914.04	K	Joback Method
tc	1122.42	K	Joback Method
tf	527.15	K	Joback Method
vc	1.244	m3/kmol	Joback Method

## Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	1030.46	J/molxK	914.04	Joback Method
cpg	1046.83	J/molxK	948.77	Joback Method
cpg	1062.01	J/molxK	983.50	Joback Method
cpg	1076.03	J/molxK	1018.23	Joback Method
cpg	1088.95	J/molxK	1052.96	Joback Method
cpg	1100.82	J/molxK	1087.69	Joback Method
cpg	1111.67	J/molxK	1122.42	Joback Method
dvisc	0.0003985	Paxs	527.15	Joback Method
dvisc	0.0002069	Paxs	591.63	Joback Method

dvisc	0.0001222	Paxs	656.11	Joback Method
dvisc	0.0000793	Paxs	720.59	Joback Method
dvisc	0.0000552	Paxs	785.08	Joback Method
dvisc	0.0000406	Paxs	849.56	Joback Method
dvisc	0.0000312	Paxs	914.04	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=U356692&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=U356692&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.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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