

# Terephthalic acid, dec-4-enyl heptyl ester

<b>Inchi:</b>	InChI=1S/C25H38O4/c1-3-5-7-9-10-11-13-15-21-29-25(27)23-18-16-22(17-19-23)24(26)
<b>InchiKey:</b>	CTTFTEXFDCPDMJ-ZHACJKMWSA-N
<b>Formula:</b>	C25H38O4
<b>SMILES:</b>	CCCCC=CCCCOC(=O)c1ccc(C(=O)OCCCCCCC)cc1
<b>Mol. weight [g/mol]:</b>	402.57

## Physical Properties

Property code	Value	Unit	Source
gf	-125.22	kJ/mol	Joback Method
hf	-706.65	kJ/mol	Joback Method
hfus	59.93	kJ/mol	Joback Method
hvap	92.45	kJ/mol	Joback Method
log10ws	-8.09		Crippen Method
logp	6.887		Crippen Method
mvol	349.930	ml/mol	McGowan Method
pc	1002.71	kPa	Joback Method
rinpol	3080.00		NIST Webbook
rinpol	3080.00		NIST Webbook
tb	959.80	K	Joback Method
tc	1175.26	K	Joback Method
tf	549.69	K	Joback Method
vc	1.355	m <sup>3</sup> /kmol	Joback Method

## Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	1153.14	J/molxK	959.80	Joback Method
cpg	1170.06	J/molxK	995.71	Joback Method
cpg	1185.67	J/molxK	1031.62	Joback Method
cpg	1200.03	J/molxK	1067.53	Joback Method
cpg	1213.20	J/molxK	1103.44	Joback Method
cpg	1225.23	J/molxK	1139.35	Joback Method
cpg	1236.18	J/molxK	1175.26	Joback Method
dvisc	0.0003167	Paxs	549.69	Joback Method

dvisc	0.0001611	Paxs	618.04	Joback Method
dvisc	0.0000938	Paxs	686.39	Joback Method
dvisc	0.0000602	Paxs	754.74	Joback Method
dvisc	0.0000416	Paxs	823.10	Joback Method
dvisc	0.0000304	Paxs	891.45	Joback Method
dvisc	0.0000233	Paxs	959.80	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=U356353&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=U356353&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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