

# Isophthalic acid, 2-methylbutyl octyl ester

<b>Inchi:</b>	InChI=1S/C21H32O4/c1-4-6-7-8-9-10-14-24-20(22)18-12-11-13-19(15-18)21(23)25-16-1
<b>InchiKey:</b>	NEBHRYPBZSVUFW-UHFFFAOYSA-N
<b>Formula:</b>	C21H32O4
<b>SMILES:</b>	CCCCCCCCOC(=O)c1cccc(C(=O)OCC(C)CC)c1
<b>Mol. weight [g/mol]:</b>	348.48

## Physical Properties

Property code	Value	Unit	Source
gf	-241.56	kJ/mol	Joback Method
hf	-746.59	kJ/mol	Joback Method
hfus	45.85	kJ/mol	Joback Method
hvap	83.20	kJ/mol	Joback Method
log10ws	-6.32		Crippen Method
logp	5.407		Crippen Method
mvol	297.870	ml/mol	McGowan Method
pc	1259.27	kPa	Joback Method
rinpol	2602.00		NIST Webbook
rinpol	2602.00		NIST Webbook
tb	863.68	K	Joback Method
tc	1066.32	K	Joback Method
tf	494.69	K	Joback Method
vc	1.145	m <sup>3</sup> /kmol	Joback Method

## Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	937.65	J/molxK	863.68	Joback Method
cpg	1008.08	J/molxK	1032.54	Joback Method
cpg	996.30	J/molxK	998.77	Joback Method
cpg	983.39	J/molxK	965.00	Joback Method
cpg	969.33	J/molxK	931.23	Joback Method
cpg	954.09	J/molxK	897.45	Joback Method
cpg	1018.76	J/molxK	1066.32	Joback Method
dvisc	0.0000436	Paxs	863.68	Joback Method

dvisc	0.0000571	Paxs	802.18	Joback Method
dvisc	0.0000783	Paxs	740.68	Joback Method
dvisc	0.0001136	Paxs	679.18	Joback Method
dvisc	0.0001775	Paxs	617.69	Joback Method
dvisc	0.0003060	Paxs	556.19	Joback Method
dvisc	0.0006043	Paxs	494.69	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=U343990&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=U343990&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

Latest version available from:

<https://www.chemeo.com/cid/87-228-3/Isophthalic-acid-2-methylbutyl-octyl-ester.pdf>

Generated by Cheméo on 2024-04-27 05:53:01.260784205 +0000 UTC m=+16486430.181361526.

Cheméo (<https://www.chemeo.com>) is the biggest free database of chemical and physical data for the process industry.