

Isophthalic acid, 3-methylbut-2-yl nonyl ester

Inchi:	InChI=1S/C22H34O4/c1-5-6-7-8-9-10-11-15-25-21(23)19-13-12-14-20(16-19)22(24)26-1
InchiKey:	CXPRCGJNILBOHD-UHFFFAOYSA-N
Formula:	C22H34O4
SMILES:	CCCCCCCCCOC(=O)c1cccc(C(=O)OC(C)C(C)C)c1
Mol. weight [g/mol]:	362.50

Physical Properties

Property code	Value	Unit	Source
gf	-235.58	kJ/mol	Joback Method
hf	-772.51	kJ/mol	Joback Method
hfus	44.92	kJ/mol	Joback Method
hvap	85.04	kJ/mol	Joback Method
log10ws	-6.85		Crippen Method
logp	5.795		Crippen Method
mvol	311.960	ml/mol	McGowan Method
pc	1184.16	kPa	Joback Method
rinpol	2642.00		NIST Webbook
rinpol	2642.00		NIST Webbook
tb	886.12	K	Joback Method
tc	1091.62	K	Joback Method
tf	490.96	K	Joback Method
vc	1.196	m ³ /kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	998.25	J/molxK	886.12	Joback Method
cpg	1069.24	J/molxK	1057.37	Joback Method
cpg	1057.48	J/molxK	1023.12	Joback Method
cpg	1044.54	J/molxK	988.87	Joback Method
cpg	1030.37	J/molxK	954.62	Joback Method
cpg	1014.95	J/molxK	920.37	Joback Method
cpg	1079.83	J/molxK	1091.62	Joback Method
dvisc	0.0000345	Paxs	886.12	Joback Method

dvisc	0.0000460	Paxs	820.26	Joback Method
dvisc	0.0000645	Paxs	754.40	Joback Method
dvisc	0.0000965	Paxs	688.54	Joback Method
dvisc	0.0001572	Paxs	622.68	Joback Method
dvisc	0.0002874	Paxs	556.82	Joback Method
dvisc	0.0006178	Paxs	490.96	Joback Method

Sources

Joback Method:	https://en.wikipedia.org/wiki/Joback_method
McGowan Method:	http://link.springer.com/article/10.1007/BF02311772
NIST Webbook:	http://webbook.nist.gov/cgi/cbook.cgi?ID=U344725&Units=SI
Crippen Method:	http://pubs.acs.org/doi/abs/10.1021/ci990307l
Crippen Method:	https://www.chemeo.com/doc/models/crippen_log10ws

Legend

cp_g:	Ideal gas heat capacity
dvisc:	Dynamic viscosity
gf:	Standard Gibbs free energy of formation
hf:	Enthalpy of formation at standard conditions
hfus:	Enthalpy of fusion at standard conditions
hvap:	Enthalpy of vaporization at standard conditions
log₁₀ws:	Log ₁₀ of Water solubility in mol/l
logp:	Octanol/Water partition coefficient
m_cvol:	McGowan's characteristic volume
pc:	Critical Pressure
rinpol:	Non-polar retention indices
tb:	Normal Boiling Point Temperature
tc:	Critical Temperature
tf:	Normal melting (fusion) point
vc:	Critical Volume

Latest version available from:

<https://www.chemeo.com/cid/96-073-5/Isophthalic-acid-3-methylbut-2-yl-nonyl-ester.pdf>

Generated by Cheméo on 2024-04-19 21:01:20.906214459 +0000 UTC m=+15849729.826791770.

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