

Phthalic acid, nonyl 2,3,6-trifluorobenzyl ester

Inchi:	InChI=1S/C24H27F3O4/c1-2-3-4-5-6-7-10-15-30-23(28)17-11-8-9-12-18(17)24(29)31-16
InchiKey:	KGYQQLRVQSUSGR-UHFFFAOYSA-N
Formula:	C24H27F3O4
SMILES:	CCCCCCCCCOC(=O)c1ccccc1C(=O)OCc1c(F)ccc(F)c1F
Mol. weight [g/mol]:	436.46

Physical Properties

Property code	Value	Unit	Source
gf	-714.77	kJ/mol	Joback Method
hf	-1189.44	kJ/mol	Joback Method
hfus	59.26	kJ/mol	Joback Method
hvap	92.08	kJ/mol	Joback Method
log10ws	-8.41		Crippen Method
logp	6.368		Crippen Method
mvol	321.690	ml/mol	McGowan Method
pc	1139.03	kPa	Joback Method
rinpol	2817.00		NIST Webbook
rinpol	2817.00		NIST Webbook
tb	972.19	K	Joback Method
tc	1190.84	K	Joback Method
tf	609.25	K	Joback Method
vc	1.266	m ³ /kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	1034.30	J/molxK	972.19	Joback Method
cpg	1047.57	J/molxK	1008.63	Joback Method
cpg	1059.46	J/molxK	1045.07	Joback Method
cpg	1070.01	J/molxK	1081.52	Joback Method
cpg	1079.26	J/molxK	1117.96	Joback Method
cpg	1087.22	J/molxK	1154.40	Joback Method
cpg	1093.95	J/molxK	1190.84	Joback Method

Sources

Crippen Method:	https://www.chemeo.com/doc/models/crippen_log10ws
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=U377792&Units=SI
Crippen Method:	http://pubs.acs.org/doi/abs/10.1021/ci990307I

Legend

cpg:	Ideal gas heat capacity
gf:	Standard Gibbs free energy of formation
hf:	Enthalpy of formation at standard conditions
hfus:	Enthalpy of fusion at standard conditions
hvp:	Enthalpy of vaporization at standard conditions
log10ws:	Log10 of Water solubility in mol/l
logp:	Octanol/Water partition coefficient
mcvol:	McGowan's characteristic volume
pc:	Critical Pressure
rinp:	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/124-733-0/Phthalic-acid-nonyl-2-3-6-trifluorobenzyl-ester.pdf>

Generated by Cheméo on 2024-05-03 15:45:33.395216839 +0000 UTC m=+17040382.315794156.

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