

Phthalic acid, 4-fluorobenzyl pentyl ester

Inchi:	InChI=1S/C20H21FO4/c1-2-3-6-13-24-19(22)17-7-4-5-8-18(17)20(23)25-14-15-9-11-16(2)
InchiKey:	WLKNEFYSOXOKAW-UHFFFAOYSA-N
Formula:	C20H21FO4
SMILES:	CCCCCOC(=O)c1ccccc1C(=O)OCc1ccc(F)cc1
Mol. weight [g/mol]:	344.38

Physical Properties

Property code	Value	Unit	Source
gf	-339.57	kJ/mol	Joback Method
hf	-691.72	kJ/mol	Joback Method
hfus	43.51	kJ/mol	Joback Method
hvap	83.48	kJ/mol	Joback Method
log10ws	-6.07		Crippen Method
logp	4.530		Crippen Method
mcvol	261.790	ml/mol	McGowan Method
pc	1653.80	kPa	Joback Method
rinpol	2444.00		NIST Webbook
rinpol	2444.00		NIST Webbook
tb	872.17	K	Joback Method
tc	1090.48	K	Joback Method
tf	537.95	K	Joback Method
vc	1.006	m ³ /kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	787.68	J/mol×K	872.17	Joback Method
cpg	801.27	J/mol×K	908.56	Joback Method
cpg	813.65	J/mol×K	944.94	Joback Method
cpg	824.84	J/mol×K	981.33	Joback Method
cpg	834.86	J/mol×K	1017.71	Joback Method
cpg	843.76	J/mol×K	1054.10	Joback Method
cpg	851.56	J/mol×K	1090.48	Joback Method

Sources

NIST Webbook:	http://webbook.nist.gov/cgi/cbook.cgi?ID=U377744&Units=SI
Crippen Method:	http://pubs.acs.org/doi/abs/10.1021/ci9903071
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

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/56-612-0/Phthalic-acid-4-fluorobenzyl-pentyl-ester.pdf>

Generated by Cheméo on 2024-04-27 18:05:30.084646745 +0000 UTC m=+16530379.005224067.

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