

Phthalic acid, octyl pent-4-enyl ester

Inchi:	InChI=1S/C21H30O4/c1-3-5-7-8-9-13-17-25-21(23)19-15-11-10-14-18(19)20(22)24-16-1
InchiKey:	IZGXXPISSXNTOK-UHFFFAOYSA-N
Formula:	C21H30O4
SMILES:	C=CCCCOC(=O)c1cccc1C(=O)OCCCCCCCC
Mol. weight [g/mol]:	346.46

Physical Properties

Property code	Value	Unit	Source
gf	-151.28	kJ/mol	Joback Method
hf	-615.88	kJ/mol	Joback Method
hfus	48.09	kJ/mol	Joback Method
hvap	82.92	kJ/mol	Joback Method
log10ws	-6.41		Crippen Method
logp	5.327		Crippen Method
mvol	293.570	ml/mol	McGowan Method
pc	1290.21	kPa	Joback Method
rinpol	2450.00		NIST Webbook
tb	860.80	K	Joback Method
tc	1063.09	K	Joback Method
tf	507.93	K	Joback Method
vc	1.133	m ³ /kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	909.38	J/molxK	860.80	Joback Method
cpg	925.25	J/molxK	894.51	Joback Method
cpg	939.98	J/molxK	928.23	Joback Method
cpg	953.61	J/molxK	961.94	Joback Method
cpg	966.16	J/molxK	995.66	Joback Method
cpg	977.66	J/molxK	1029.37	Joback Method
cpg	988.15	J/molxK	1063.09	Joback Method
dvisc	0.0005479	Paxs	507.93	Joback Method
dvisc	0.0003004	Paxs	566.74	Joback Method

dvisc	0.0001844	Paxs	625.55	Joback Method
dvisc	0.0001231	Paxs	684.37	Joback Method
dvisc	0.0000876	Paxs	743.18	Joback Method
dvisc	0.0000656	Paxs	801.99	Joback Method
dvisc	0.0000510	Paxs	860.80	Joback Method

Sources

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
NIST Webbook:	http://webbook.nist.gov/cgi/cbook.cgi?ID=U360467&Units=SI

Legend

cpg:	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:	Log10 of Water solubility in mol/l
logp:	Octanol/Water partition coefficient
mcvol:	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/61-661-0/Phthalic-acid-octyl-pent-4-enyl-ester.pdf>

Generated by Cheméo on 2024-04-29 06:41:02.757068632 +0000 UTC m=+16662111.677645996.

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