

Phthalic acid, hexyl 2-methoxybenzyl ester

Inchi: InChI=1S/C22H26O5/c1-3-4-5-10-15-26-21(23)18-12-7-8-13-19(18)22(24)27-16-17-11-6
InchiKey: ZAOSZPSCZFQDGP-UHFFFAOYSA-N
Formula: C22H26O5
SMILES: CCCCCOC(=O)c1ccccc1C(=O)OCc1ccccc1OC
Mol. weight [g/mol]: 370.44

Physical Properties

Property code	Value	Unit	Source
gf	-232.92	kJ/mol	Joback Method
hf	-669.11	kJ/mol	Joback Method
hfus	46.80	kJ/mol	Joback Method
hvap	91.16	kJ/mol	Joback Method
log10ws	-6.28		Crippen Method
logp	4.789		Crippen Method
mcvol	294.070	ml/mol	McGowan Method
pc	1453.46	kPa	Joback Method
rinpol	3167.00		NIST Webbook
rinpol	3167.00		NIST Webbook
tb	941.08	K	Joback Method
tc	1164.13	K	Joback Method
tf	582.13	K	Joback Method
vc	1.117	m ³ /kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	924.20	J/molxK	941.08	Joback Method
cpg	976.76	J/molxK	1126.96	Joback Method
cpg	969.08	J/molxK	1089.78	Joback Method
cpg	960.00	J/molxK	1052.61	Joback Method
cpg	949.51	J/molxK	1015.43	Joback Method
cpg	937.58	J/molxK	978.26	Joback Method
cpg	983.08	J/molxK	1164.13	Joback Method
dvisc	0.0000323	Paxs	941.08	Joback Method

dvisc	0.0000407	Paxs	881.25	Joback Method
dvisc	0.0000531	Paxs	821.43	Joback Method
dvisc	0.0000721	Paxs	761.61	Joback Method
dvisc	0.0001032	Paxs	701.78	Joback Method
dvisc	0.0001580	Paxs	641.95	Joback Method
dvisc	0.0002639	Paxs	582.13	Joback Method

Sources

McGowan Method:	http://link.springer.com/article/10.1007/BF02311772
NIST Webbook:	http://webbook.nist.gov/cgi/cbook.cgi?ID=U382495&Units=SI
Crippen Method:	http://pubs.acs.org/doi/abs/10.1021/ci990307l
Crippen Method:	https://www.cheméo.com/doc/models/crippen_log10ws
Joback Method:	https://en.wikipedia.org/wiki/Joback_method

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

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