

Isophthalic acid, di(2-methylphenyl) ester

Inchi:	InChI=1S/C22H18O4/c1-15-8-3-5-12-19(15)25-21(23)17-10-7-11-18(14-17)22(24)26-20-
InchiKey:	LIBNNSNDAMJTFG-UHFFFAOYSA-N
Formula:	C22H18O4
SMILES:	<chem>Cc1ccccc1OC(=O)c1cccc(C(=O)Oc2ccccc2C)c1</chem>
Mol. weight [g/mol]:	346.38

Physical Properties

Property code	Value	Unit	Source
gf	-25.14	kJ/mol	Joback Method
hf	-311.83	kJ/mol	Joback Method
hfus	39.27	kJ/mol	Joback Method
hvap	91.69	kJ/mol	Joback Method
log10ws	-6.60		Crippen Method
logp	4.742		Crippen Method
mvol	264.440	ml/mol	McGowan Method
pc	1906.90	kPa	Joback Method
rinpol	2990.00		NIST Webbook
rinpol	2990.00		NIST Webbook
tb	950.32	K	Joback Method
tc	1201.35	K	Joback Method
tf	598.84	K	Joback Method
vc	0.992	m ³ /kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	790.76	J/molxK	950.32	Joback Method
cpg	835.38	J/molxK	1159.51	Joback Method
cpg	829.28	J/molxK	1117.67	Joback Method
cpg	821.82	J/molxK	1075.83	Joback Method
cpg	812.95	J/molxK	1034.00	Joback Method
cpg	802.62	J/molxK	992.16	Joback Method
cpg	840.17	J/molxK	1201.35	Joback Method
dvisc	0.0000458	Paxs	950.32	Joback Method

dvisc	0.0000566	Paxs	891.74	Joback Method
dvisc	0.0000721	Paxs	833.16	Joback Method
dvisc	0.0000952	Paxs	774.58	Joback Method
dvisc	0.0001316	Paxs	716.00	Joback Method
dvisc	0.0001927	Paxs	657.42	Joback Method
dvisc	0.0003041	Paxs	598.84	Joback Method

Sources

McGowan Method:	http://link.springer.com/article/10.1007/BF02311772
NIST Webbook:	http://webbook.nist.gov/cgi/cbook.cgi?ID=U344353&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

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
log_p:	Octanol/Water partition coefficient
m_{cvol}:	McGowan's characteristic volume
pc:	Critical Pressure
rin_{pol}:	Non-polar retention indices
tb:	Normal Boiling Point Temperature
tc:	Critical Temperature
tf:	Normal melting (fusion) point
vc:	Critical Volume

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