Isophthalic acid, 2-ethylphenyl isobutyl ester

Inchi: InChI=1S/C20H22O4/c1-4-15-8-5-6-11-18(15)24-20(22)17-10-7-9-16(12-17)19(21)23-13
InchiKey: POKKUATZZIRARQ-UHFFFAOYSA-N
Formula: C20H22O4
SMILES: CCc1ccccc1OC(=O)c1cccc(C(=O)OCC(C)C)c1
Mol. weight [g/mol]: 326.39

Physical Properties

Property code	Value	Unit	Source
gf	-147.20	kJ/mol	Joback Method
hf	-500.89	kJ/mol	Joback Method
hfus	36.91	kJ/mol	Joback Method
hvap	83.91	kJ/mol	Joback Method
log10ws	-5.62		Crippen Method
logp	4.281		Crippen Method
mcvol	260.020	ml/mol	McGowan Method
pc	1727.46	kPa	Joback Method
rinpol	2445.00		NIST Webbook
rinpol	2445.00		NIST Webbook
tb	872.46	K	Joback Method
tc	1099.91	K	Joback Method
tf	522.36	K	Joback Method
vc	0.982	m3/kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	780.90	J/mol×K	872.46	Joback Method
cpg	795.15	J/mol×K	910.37	Joback Method
cpg	808.06	J/mol×K	948.28	Joback Method
cpg	819.68	J/mol×K	986.18	Joback Method
cpg	830.02	J/mol×K	1024.09	Joback Method
cpg	839.13	J/mol×K	1062.00	Joback Method
cpg	847.04	J/mol×K	1099.91	Joback Method
dvisc	0.0004963	Pa×s	522.36	Joback Method

dvisc	0.0002824	Pa×s	580.71	Joback Method
dvisc	0.0001781	Pa×s	639.06	Joback Method
dvisc	0.0001213	Pa×s	697.41	Joback Method
dvisc	0.0000877	Pa×s	755.76	Joback Method
dvisc	0.0000664	Pa×s	814.11	Joback Method
dvisc	0.0000522	Pa×s	872.46	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=U356761&Units=SI
Crippen Method:	http://pubs.acs.org/doi/abs/10.1021/ci990307l

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