

Diglycolic acid, heptyl 2-methoxyphenyl ester

Inchi:	InChI=1S/C18H26O6/c1-3-4-5-6-9-12-23-17(19)13-22-14-18(20)24-16-11-8-7-10-15(16)2
InchiKey:	BXSPRKNXUCUIJP-UHFFFAOYSA-N
Formula:	C18H26O6
SMILES:	CCCCCCCOC(=O)COCC(=O)Oc1ccccc1OC
Mol. weight [g/mol]:	338.40

Physical Properties

Property code	Value	Unit	Source
gf	-474.38	kJ/mol	Joback Method
hf	-943.83	kJ/mol	Joback Method
hfus	43.98	kJ/mol	Joback Method
hvap	81.73	kJ/mol	Joback Method
log10ws	-3.62		Crippen Method
logp	3.131		Crippen Method
mvol	267.340	ml/mol	McGowan Method
pc	1511.67	kPa	Joback Method
rinpol	3049.00		NIST Webbook
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tb	840.32	K	Joback Method
tc	1041.15	K	Joback Method
tf	520.34	K	Joback Method
vc	1.020	m ³ /kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	818.63	J/molxK	840.32	Joback Method
cpg	833.41	J/molxK	873.79	Joback Method
cpg	847.00	J/molxK	907.26	Joback Method
cpg	859.41	J/molxK	940.74	Joback Method
cpg	870.62	J/molxK	974.21	Joback Method
cpg	880.63	J/molxK	1007.68	Joback Method
cpg	889.43	J/molxK	1041.15	Joback Method
dvisc	0.0003593	Paxs	520.34	Joback Method

dvisc	0.0002114	Paxs	573.67	Joback Method
dvisc	0.0001361	Paxs	627.00	Joback Method
dvisc	0.0000939	Paxs	680.33	Joback Method
dvisc	0.0000684	Paxs	733.66	Joback Method
dvisc	0.0000520	Paxs	786.99	Joback Method
dvisc	0.0000409	Paxs	840.32	Joback Method

Sources

McGowan Method:	http://link.springer.com/article/10.1007/BF02311772
NIST Webbook:	http://webbook.nist.gov/cgi/cbook.cgi?ID=U382246&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
h_{vap}:	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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