Dimethylmalonic acid, 4-chlorophenyl pentadecyl ester

Inchi: InChI=1S/C26H41ClO4/c1-4-5-6-7-8-9-10-11-12-13-14-15-16-21-30-24(28)26(2,3)25(29)
InchiKey: UWCXKJKFJPSPST-UHFFFAOYSA-N
Formula: C26H41ClO4
SMILES: CCCCCCCCCCCCCCCCOC(=O)C(C)(C)C(=O)Oc1ccc(Cl)cc1
Mol. weight [g/mol]: 453.05

Physical Properties

Property code	Value	Unit	Source
gf	-206.11	kJ/mol	Joback Method
hf	-869.00	kJ/mol	Joback Method
hfus	59.10	kJ/mol	Joback Method
hvap	97.81	kJ/mol	Joback Method
log10ws	-8.63		Crippen Method
logp	7.906		Crippen Method
mcvol	380.560	ml/mol	McGowan Method
pc	905.06	kPa	Joback Method
rinpol	3000.00		NIST Webbook
rinpol	3000.00		NIST Webbook
tb	1012.72	K	Joback Method
tc	1240.09	K	Joback Method
tf	598.38	K	Joback Method
vc	1.470	m3/kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	1269.75	J/mol×K	1012.72	Joback Method
cpg	1286.19	J/mol×K	1050.62	Joback Method
cpg	1301.21	J/mol×K	1088.51	Joback Method
cpg	1314.90	J/mol×K	1126.41	Joback Method
cpg	1327.34	J/mol×K	1164.30	Joback Method
cpg	1338.60	J/mol×K	1202.20	Joback Method
cpg	1348.78	J/mol×K	1240.09	Joback Method
dvisc	0.0002050	Pa×s	598.38	Joback Method

dvisc	0.0001046	Pa×s	667.44	Joback Method
dvisc	0.0000606	Pa×s	736.49	Joback Method
dvisc	0.0000385	Pa×s	805.55	Joback Method
dvisc	0.0000263	Pa×s	874.61	Joback Method
dvisc	0.0000190	Pa×s	943.66	Joback Method
dvisc	0.0000144	Pa×s	1012.72	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=U361983&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