

Dimethylmalonic acid, ethyl 2-ethylhexyl ester

Inchi:

InchiKey:

Formula:

SMILES:

Mol. weight [g/mol]:

InChI=1S/C15H28O4/c1-6-9-10-12(7-2)11-19-14(17)15(4,5)13(16)18-8-3/h12H,6-11H2,1

UAVPSEKGESQSEI-UHFFFAOYSA-N

C15H28O4

CCCCC(CC)COC(=O)C(C)(C)C(=O)OCC

272.38

Physical Properties

Property code	Value	Unit	Source
gf	-392.02	kJ/mol	Joback Method
hf	-856.56	kJ/mol	Joback Method
hfus	29.24	kJ/mol	Joback Method
hvap	65.61	kJ/mol	Joback Method
log10ws	-3.34		Crippen Method
logp	3.335		Crippen Method
mcvol	237.090	ml/mol	McGowan Method
pc	1551.22	kPa	Joback Method
rinpol	1591.00		NIST Webbook
rinpol	1591.00		NIST Webbook
tb	691.51	K	Joback Method
tc	876.12	K	Joback Method
tf	390.55	K	Joback Method
vc	0.906	m3/kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	680.11	J/mol×K	691.51	Joback Method
cpg	696.84	J/mol×K	722.28	Joback Method
cpg	712.70	J/mol×K	753.05	Joback Method
cpg	727.68	J/mol×K	783.82	Joback Method
cpg	741.83	J/mol×K	814.59	Joback Method
cpg	755.14	J/mol×K	845.35	Joback Method
cpg	767.65	J/mol×K	876.12	Joback Method
dvisc	0.0017437	Pa×s	390.55	Joback Method

dvisc	0.0007852	Pa×s	440.71	Joback Method
dvisc	0.0004162	Pa×s	490.87	Joback Method
dvisc	0.0002481	Pa×s	541.03	Joback Method
dvisc	0.0001615	Pa×s	591.19	Joback Method
dvisc	0.0001124	Pa×s	641.35	Joback Method
dvisc	0.0000825	Pa×s	691.51	Joback Method

Sources

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=U361630&Units=SI
Crippen Method:	http://pubs.acs.org/doi/abs/10.1021/ci990307l
Crippen Method:	https://www.chemeo.com/doc/models/crippen_log10ws

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

Latest version available from:

<https://www.chemeo.com/cid/68-813-4/Dimethylmalonic-acid-ethyl-2-ethylhexyl-ester.pdf>

Generated by Cheméo on 2025-12-23 18:09:50.774719284 +0000 UTC m=+6261588.304759939.

Cheméo (<https://www.chemeo.com>) is the biggest free database of chemical and physical data for the process industry.