

ethyl 2-methyltetradecanoate

Inchi:	InChI=1S/C17H34O2/c1-4-6-7-8-9-10-11-12-13-14-15-16(3)17(18)19-5-2/h16H,4-15H2,1
InchiKey:	UOEPHBOLYJPFHY-UHFFFAOYSA-N
Formula:	C17H34O2
SMILES:	CCCCCCCCCCCC(C)C(=O)OCC
Mol. weight [g/mol]:	270.45

Physical Properties

Property code	Value	Unit	Source
gf	-144.10	kJ/mol	Joback Method
hf	-644.29	kJ/mol	Joback Method
hfus	39.05	kJ/mol	Joback Method
hvap	62.20	kJ/mol	Joback Method
log10ws	-5.56		Crippen Method
logp	5.497		Crippen Method
mvol	257.830	ml/mol	McGowan Method
pc	1279.16	kPa	Joback Method
ripol	2119.00		NIST Webbook
ripol	2119.00		NIST Webbook
tb	664.21	K	Joback Method
tc	833.82	K	Joback Method
tf	338.51	K	Joback Method
vc	1.006	m ³ /kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	732.67	J/molxK	664.21	Joback Method
cpg	817.65	J/molxK	805.55	Joback Method
cpg	802.22	J/molxK	777.28	Joback Method
cpg	786.03	J/molxK	749.02	Joback Method
cpg	769.05	J/molxK	720.75	Joback Method
cpg	751.27	J/molxK	692.48	Joback Method
cpg	832.34	J/molxK	833.82	Joback Method
dvisc	0.0001003	Paxs	664.21	Joback Method

dvisc	0.0001368	Paxs	609.93	Joback Method
dvisc	0.0001983	Paxs	555.64	Joback Method
dvisc	0.0003116	Paxs	501.36	Joback Method
dvisc	0.0005461	Paxs	447.08	Joback Method
dvisc	0.0011177	Paxs	392.79	Joback Method
dvisc	0.0028787	Paxs	338.51	Joback Method

Sources

NIST Webbook:	http://webbook.nist.gov/cgi/cbook.cgi?ID=R340824&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
McGowan Method:	http://link.springer.com/article/10.1007/BF02311772

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
logp:	Octanol/Water partition coefficient
m_{cvol}:	McGowan's characteristic volume
pc:	Critical Pressure
ripol:	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/82-879-6/ethyl-2-methyltetradecanoate.pdf>

Generated by Cheméo on 2024-04-19 00:19:55.730326421 +0000 UTC m=+15775244.650903736.

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