

Ethyl 2,6-dimethylbenzoate

Inchi:	InChI=1S/C11H14O2/c1-4-13-11(12)10-8(2)6-5-7-9(10)3/h5-7H,4H2,1-3H3
InchiKey:	NBBSIMLSNFENN-UHFFFAOYSA-N
Formula:	C11H14O2
SMILES:	CCOC(=O)c1c(C)cccc1C
Mol. weight [g/mol]:	178.23
CAS:	36596-67-5

Physical Properties

Property code	Value	Unit	Source
gf	-99.03	kJ/mol	Joback Method
hf	-301.58	kJ/mol	Joback Method
hfus	20.30	kJ/mol	Joback Method
hvap	52.84	kJ/mol	Joback Method
log10ws	-3.08		Crippen Method
logp	2.480		Crippen Method
mcvol	149.530	ml/mol	McGowan Method
pc	2693.00	kPa	Joback Method
tb	564.01	K	Joback Method
tc	775.37	K	Joback Method
tf	337.35	K	Joback Method
vc	0.568	m ³ /kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	345.82	J/mol×K	564.01	Joback Method
cpg	359.69	J/mol×K	599.24	Joback Method
cpg	372.86	J/mol×K	634.46	Joback Method
cpg	385.33	J/mol×K	669.69	Joback Method
cpg	397.11	J/mol×K	704.92	Joback Method
cpg	408.21	J/mol×K	740.15	Joback Method
cpg	418.63	J/mol×K	775.37	Joback Method
dvisc	0.0013555	Paxs	337.35	Joback Method
dvisc	0.0008295	Paxs	375.13	Joback Method

dvisc	0.0005553	Paxs	412.90	Joback Method
dvisc	0.0003976	Paxs	450.68	Joback Method
dvisc	0.0002998	Paxs	488.46	Joback Method
dvisc	0.0002354	Paxs	526.23	Joback Method
dvisc	0.0001909	Paxs	564.01	Joback Method

Sources

NIST Webbook:	http://webbook.nist.gov/cgi/cbook.cgi?ID=C36596675&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

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
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/34-044-5/Ethyl-2-6-dimethylbenzoate.pdf>

Generated by Cheméo on 2024-04-26 07:44:34.987791003 +0000 UTC m=+16406723.908368318.

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