

Dimethyl(E)-1-propenyl orthoformate

Inchi:	InChI=1S/C6H12O3/c1-4-5-9-6(7-2)8-3/h4-6H,1-3H3/b5-4+
InchiKey:	DDINHWIZOMBDDD-SNAWJCMRSA-N
Formula:	C6H12O3
SMILES:	CC=COC(OC)OC
Mol. weight [g/mol]:	132.16
CAS:	66178-19-6

Physical Properties

Property code	Value	Unit	Source
gf	-237.58	kJ/mol	Joback Method
hf	-451.89	kJ/mol	Joback Method
hfus	11.54	kJ/mol	Joback Method
hvap	35.75	kJ/mol	Joback Method
log10ws	-1.05		Crippen Method
logp	1.113		Crippen Method
mcvol	108.710	ml/mol	McGowan Method
pc	3138.51	kPa	Joback Method
tb	407.66	K	Joback Method
tc	585.73	K	Joback Method
tf	203.99	K	Joback Method
vc	0.400	m3/kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	211.99	J/mol×K	407.66	Joback Method
cpg	221.83	J/mol×K	437.34	Joback Method
cpg	231.44	J/mol×K	467.02	Joback Method
cpg	240.81	J/mol×K	496.70	Joback Method
cpg	249.93	J/mol×K	526.38	Joback Method
cpg	258.79	J/mol×K	556.06	Joback Method
cpg	267.38	J/mol×K	585.73	Joback Method
dvisc	0.0031068	Pa×s	203.99	Joback Method
dvisc	0.0012910	Pa×s	237.94	Joback Method

dvisc	0.0006679	Pa×s	271.88	Joback Method
dvisc	0.0004000	Pa×s	305.82	Joback Method
dvisc	0.0002654	Pa×s	339.77	Joback Method
dvisc	0.0001897	Pa×s	373.71	Joback Method
dvisc	0.0001434	Pa×s	407.66	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=C66178196&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
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/12-942-2/Dimethyl-E-1-propenyl-orthoformate.pdf>

Generated by Cheméo on 2025-12-05 19:00:55.463062335 +0000 UTC m=+4709452.993102988.

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