

2,5-di-O-acetyl-1,4-Anhydro-3-O-methyl-D-ribitol

Inchi:	InChI=1S/C10H16O6/c1-6(11)14-4-8-10(13-3)9(5-15-8)16-7(2)12/h8-10H,4-5H2,1-3H3/t8
InchiKey:	SULWLWFNHXEWRP-KXUCPTDWSA-N
Formula:	C10H16O6
SMILES:	COC1C(COC(C)=O)OCC1OC(C)=O
Mol. weight [g/mol]:	232.23

Physical Properties

Property code	Value	Unit	Source
gf	-604.51	kJ/mol	Joback Method
hf	-983.75	kJ/mol	Joback Method
hfus	32.47	kJ/mol	Joback Method
hvap	62.72	kJ/mol	Joback Method
log10ws	-0.14		Crippen Method
logp	-0.105		Crippen Method
mcvol	167.520	ml/mol	McGowan Method
pc	2505.01	kPa	Joback Method
rinpol	1480.59		NIST Webbook
rinpol	1480.59		NIST Webbook
tb	636.09	K	Joback Method
tc	837.04	K	Joback Method
tf	398.00	K	Joback Method
vc	0.622	m3/kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	464.89	J/mol×K	636.09	Joback Method
cpg	480.18	J/mol×K	669.58	Joback Method
cpg	494.68	J/mol×K	703.07	Joback Method
cpg	508.36	J/mol×K	736.57	Joback Method
cpg	521.20	J/mol×K	770.06	Joback Method
cpg	533.18	J/mol×K	803.55	Joback Method
cpg	544.28	J/mol×K	837.04	Joback Method
dvisc	0.0014137	Pa×s	398.00	Joback Method

dvisc	0.0009483	Pa×s	437.68	Joback Method
dvisc	0.0006797	Pa×s	477.36	Joback Method
dvisc	0.0005128	Pa×s	517.05	Joback Method
dvisc	0.0004027	Pa×s	556.73	Joback Method
dvisc	0.0003266	Pa×s	596.41	Joback Method
dvisc	0.0002719	Pa×s	636.09	Joback Method

Sources

McGowan Method:	http://link.springer.com/article/10.1007/BF02311772
NIST Webbook:	http://webbook.nist.gov/cgi/cbook.cgi?ID=R180637&Units=SI
Crippen Method:	http://pubs.acs.org/doi/abs/10.1021/ci990307l
Crippen Method:	https://www.cheméo.com/doc/models/crippen_log10ws
Joback Method:	https://en.wikipedia.org/wiki/Joback_method

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.cheméo.com/cid/86-061-9/2-5-di-O-acetyl-1-4-Anhydro-3-O-methyl-D-ribitol.pdf>

Generated by Cheméo on 2025-12-05 18:08:10.000454469 +0000 UTC m=+4706287.530495134.

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