

pentyl-d2 acetate

Inchi:	InChI=1S/C7H14O2/c1-3-4-5-6-9-7(2)8/h3-6H2,1-2H3/i2D2
InchiKey:	PGMYKACGEOXYJE-CBTSVUPCSA-N
Formula:	C7H12D2O2
SMILES:	CCCCCOC(C)=O
Mol. weight [g/mol]:	132.20

Physical Properties

Property code	Value	Unit	Source
gf	-225.86	kJ/mol	Joback Method
hf	-432.61	kJ/mol	Joback Method
hfus	16.67	kJ/mol	Joback Method
hvap	40.33	kJ/mol	Joback Method
log10ws	-1.62		Crippen Method
logp	1.740		Crippen Method
mvol	116.930	ml/mol	McGowan Method
pc	2940.89	kPa	Joback Method
ripol	1173.00		NIST Webbook
ripol	1173.00		NIST Webbook
tb	435.85	K	Joback Method
tc	612.24	K	Joback Method
tf	240.81	K	Joback Method
vc	0.452	m ³ /kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	238.65	J/molxK	435.85	Joback Method
cpg	290.48	J/molxK	582.84	Joback Method
cpg	280.85	J/molxK	553.44	Joback Method
cpg	270.86	J/molxK	524.04	Joback Method
cpg	260.50	J/molxK	494.65	Joback Method
cpg	249.76	J/molxK	465.25	Joback Method
cpg	299.73	J/molxK	612.24	Joback Method
dvisc	0.0002656	Paxs	435.85	Joback Method

dvisc	0.0003401	Paxs	403.34	Joback Method
dvisc	0.0004549	Paxs	370.84	Joback Method
dvisc	0.0006434	Paxs	338.33	Joback Method
dvisc	0.0009795	Paxs	305.82	Joback Method
dvisc	0.0016480	Paxs	273.32	Joback Method
dvisc	0.0031909	Paxs	240.81	Joback Method

Sources

NIST Webbook:	http://webbook.nist.gov/cgi/cbook.cgi?ID=R329129&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
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/71-608-8/pentyl-d2-acetate.pdf>

Generated by Cheméo on 2024-04-19 17:05:57.320588688 +0000 UTC m=+15835606.241166004.

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