

4-oxodecanal

Inchi:	InChI=1S/C10H18O2/c1-2-3-4-5-7-10(12)8-6-9-11/h9H,2-8H2,1H3
InchiKey:	UKCSOARNKKNOCN-UHFFFAOYSA-N
Formula:	C10H18O2
SMILES:	CCCCCCC(=O)CCC=O
Mol. weight [g/mol]:	170.25

Physical Properties

Property code	Value	Unit	Source
gf	-195.12	kJ/mol	Joback Method
hf	-447.89	kJ/mol	Joback Method
hfus	25.54	kJ/mol	Joback Method
hvap	51.32	kJ/mol	Joback Method
log10ws	-2.57		Crippen Method
logp	2.505		Crippen Method
mcvol	154.900	ml/mol	McGowan Method
pc	2400.57	kPa	Joback Method
rinpol	1333.00		NIST Webbook
rinpol	1303.00		NIST Webbook
rinpol	1304.00		NIST Webbook
tb	530.73	K	Joback Method
tc	708.49	K	Joback Method
tf	294.39	K	Joback Method
vc	0.619	m3/kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	365.55	J/molxK	530.73	Joback Method
cpg	426.10	J/molxK	678.86	Joback Method
cpg	415.09	J/molxK	649.24	Joback Method
cpg	403.55	J/molxK	619.61	Joback Method
cpg	391.46	J/molxK	589.98	Joback Method
cpg	378.79	J/molxK	560.36	Joback Method
cpg	436.58	J/molxK	708.49	Joback Method

dvisc	0.0003179	Paxs	530.73	Joback Method
dvisc	0.0004104	Paxs	491.34	Joback Method
dvisc	0.0005539	Paxs	451.95	Joback Method
dvisc	0.0007918	Paxs	412.56	Joback Method
dvisc	0.0012203	Paxs	373.17	Joback Method
dvisc	0.0020830	Paxs	333.78	Joback Method
dvisc	0.0041027	Paxs	294.39	Joback Method

Sources

Crippen Method:	http://pubs.acs.org/doi/abs/10.1021/ci990307l
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=R224405&Units=SI

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
mvol:	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.chemeo.com/cid/50-905-1/4-oxodecanal.pdf>

Generated by Cheméo on 2024-04-27 02:50:26.635192782 +0000 UTC m=+16475475.555770093.

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