

Chloromethyl 8-chlorododecanoate

Other names:	8-Chlorododecanoic acid, chloromethyl ester
Inchi:	InChI=1S/C13H24Cl2O2/c1-2-3-8-12(15)9-6-4-5-7-10-13(16)17-11-14/h12H,2-11H2,1H3
InchiKey:	JVKMUBVORYFQKQ-UHFFFAOYSA-N
Formula:	C13H24Cl2O2
SMILES:	CCCCC(Cl)CCCCCCC(=O)OCCI
Mol. weight [g/mol]:	283.23
CAS:	80419-04-1

Physical Properties

Property code	Value	Unit	Source
gf	-201.64	kJ/mol	Joback Method
hf	-593.21	kJ/mol	Joback Method
hfus	37.08	kJ/mol	Joback Method
hvap	62.07	kJ/mol	Joback Method
log10ws	-5.04		Crippen Method
logp	4.864		Crippen Method
mcvol	225.950	ml/mol	McGowan Method
pc	1618.07	kPa	Joback Method
rinpol	1878.00		NIST Webbook
rinpol	1887.00		NIST Webbook
rinpol	1919.00		NIST Webbook
rinpol	1878.00		NIST Webbook
rinpol	1891.00		NIST Webbook
ripol	2476.00		NIST Webbook
ripol	2481.00		NIST Webbook
ripol	2453.00		NIST Webbook
tb	647.55	K	Joback Method
tc	828.32	K	Joback Method
tf	353.27	K	Joback Method
vc	0.879	m ³ /kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
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cpg	579.55	J/molxK	647.55	Joback Method
cpg	594.71	J/molxK	677.68	Joback Method
cpg	609.14	J/molxK	707.81	Joback Method
cpg	622.86	J/molxK	737.94	Joback Method
cpg	635.88	J/molxK	768.06	Joback Method
cpg	648.22	J/molxK	798.19	Joback Method
cpg	659.91	J/molxK	828.32	Joback Method
dvisc	0.0025326	Paxs	353.27	Joback Method
dvisc	0.0011518	Paxs	402.32	Joback Method
dvisc	0.0006217	Paxs	451.36	Joback Method
dvisc	0.0003786	Paxs	500.41	Joback Method
dvisc	0.0002520	Paxs	549.46	Joback Method
dvisc	0.0001792	Paxs	598.50	Joback Method
dvisc	0.0001343	Paxs	647.55	Joback Method

Sources

NIST Webbook:	http://webbook.nist.gov/cgi/cbook.cgi?ID=C80419041&Units=SI
Crippen Method:	http://pubs.acs.org/doi/abs/10.1021/ci990307I
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
rinpolar:	Non-polar retention indices
ripolar:	Polar retention indices
tb:	Normal Boiling Point Temperature

tc: Critical Temperature
tf: Normal melting (fusion) point
vc: Critical Volume

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