

2-Butenoyl chloride

Other names:	Crotonyl chloride Crotonic acid chloride (2E)-2-butenoyl chloride crotonoyl chloride
Inchi:	InChI=1S/C4H5ClO/c1-2-3-4(5)6/h2-3H,1H3
InchiKey:	RJUIDDKTATZJFE-UHFFFAOYSA-N
Formula:	C4H5ClO
SMILES:	CC=CC(=O)Cl
Mol. weight [g/mol]:	104.53
CAS:	10487-71-5

Physical Properties

Property code	Value	Unit	Source
gf	-77.83	kJ/mol	Joback Method
hf	-136.99	kJ/mol	Joback Method
hfus	12.11	kJ/mol	Joback Method
hvap	35.59	kJ/mol	Joback Method
log10ws	-1.28		Crippen Method
logp	1.328		Crippen Method
mcvol	76.730	ml/mol	McGowan Method
pc	4305.56	kPa	Joback Method
rinpola	796.00		NIST Webbook
tb	394.00 ± 4.00	K	NIST Webbook
tb	394.00 ± 4.00	K	NIST Webbook
tc	584.10	K	Joback Method
tf	209.61	K	Joback Method
vc	0.294	m ³ /kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	116.29	J/mol×K	386.38	Joback Method
cpg	122.83	J/mol×K	419.33	Joback Method
cpg	128.99	J/mol×K	452.29	Joback Method

cpg	134.79	J/molxK	485.24	Joback Method
cpg	140.25	J/molxK	518.20	Joback Method
cpg	145.38	J/molxK	551.15	Joback Method
cpg	150.21	J/molxK	584.10	Joback Method
dvisc	0.0031414	Paxs	209.61	Joback Method
dvisc	0.0016640	Paxs	239.07	Joback Method
dvisc	0.0010133	Paxs	268.53	Joback Method
dvisc	0.0006807	Paxs	298.00	Joback Method
dvisc	0.0004911	Paxs	327.46	Joback Method
dvisc	0.0003740	Paxs	356.92	Joback Method
dvisc	0.0002969	Paxs	386.38	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=C10487715&Units=SI
Crippen Method:	http://pubs.acs.org/doi/abs/10.1021/ci9903071

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

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