

1-Propanone, 2-chloro-1-(4-ethylphenyl)-2-methyl-

Other names:	2-Chloro-1-(4-ethylphenyl)-2-methyl-1-propanone
Inchi:	InChI=1S/C12H15ClO/c1-4-9-5-7-10(8-6-9)11(14)12(2,3)13/h5-8H,4H2,1-3H3
InchiKey:	JRCYBVHLGNUGDA-UHFFFAOYSA-N
Formula:	C12H15ClO
SMILES:	CCc1ccc(C(=O)C(C)(C)Cl)cc1
Mol. weight [g/mol]:	210.70
CAS:	55012-69-6

Physical Properties

Property code	Value	Unit	Source
gf	14.93	kJ/mol	Joback Method
hf	-203.02	kJ/mol	Joback Method
hfus	18.87	kJ/mol	Joback Method
hvap	55.08	kJ/mol	Joback Method
log10ws	-4.04		Crippen Method
logp	3.449		Crippen Method
mcvol	169.990	ml/mol	McGowan Method
pc	2477.65	kPa	Joback Method
tb	593.69	K	Joback Method
tc	820.05	K	Joback Method
tf	346.21	K	Joback Method
vc	0.643	m ³ /kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	399.66	J/molxK	593.69	Joback Method
cpg	414.77	J/molxK	631.42	Joback Method
cpg	428.81	J/molxK	669.14	Joback Method
cpg	441.82	J/molxK	706.87	Joback Method
cpg	453.88	J/molxK	744.60	Joback Method
cpg	465.06	J/molxK	782.33	Joback Method
cpg	475.42	J/molxK	820.05	Joback Method
dvisc	0.0023840	Paxs	346.21	Joback Method

dvisc	0.0012574	Paxs	387.46	Joback Method
dvisc	0.0007500	Paxs	428.70	Joback Method
dvisc	0.0004899	Paxs	469.95	Joback Method
dvisc	0.0003427	Paxs	511.20	Joback Method
dvisc	0.0002529	Paxs	552.44	Joback Method
dvisc	0.0001947	Paxs	593.69	Joback Method

Sources

NIST Webbook:	http://webbook.nist.gov/cgi/cbook.cgi?ID=C55012696&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

cp_g:	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
h_{vap}:	Enthalpy of vaporization at standard conditions
log₁₀ws:	Log ₁₀ of Water solubility in mol/l
logp:	Octanol/Water partition coefficient
mcvol:	McGowan's characteristic volume
pc:	Critical Pressure
tb:	Normal Boiling Point Temperature
tc:	Critical Temperature
tf:	Normal melting (fusion) point
vc:	Critical Volume

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