

Pentanoyl chloride

Other names:	UN 2502 Valeroyl chloride Valeryl chloride n-Valeroyl chloride
Inchi:	InChI=1S/C5H9ClO/c1-2-3-4-5(6)7/h2-4H2,1H3
InchiKey:	XGISHOFUAFNYQF-UHFFFAOYSA-N
Formula:	C5H9ClO
SMILES:	CCCCC(=O)Cl
Mol. weight [g/mol]:	120.58
CAS:	638-29-9

Physical Properties

Property code	Value	Unit	Source
gf	-149.63	kJ/mol	Joback Method
hf	-274.85	kJ/mol	Joback Method
hfus	14.50	kJ/mol	Joback Method
hvap	37.86	kJ/mol	Joback Method
log10ws	-1.84		Crippen Method
logp	1.942		Crippen Method
mcvol	95.120	ml/mol	McGowan Method
pc	3577.07	kPa	Joback Method
rinsol	809.40		NIST Webbook
tb	399.20	K	NIST Webbook
tb	400.40 ± 0.60	K	NIST Webbook
tc	591.48	K	Joback Method
tf	225.96	K	Joback Method
vc	0.370	m ³ /kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	166.17	J/mol×K	405.10	Joback Method
cpg	211.83	J/mol×K	591.48	Joback Method
cpg	205.06	J/mol×K	560.42	Joback Method

cpg	197.97	J/molxK	529.35	Joback Method
cpg	190.54	J/molxK	498.29	Joback Method
cpg	182.77	J/molxK	467.23	Joback Method
cpg	174.65	J/molxK	436.16	Joback Method
cpl	187.90	J/molxK	298.00	NIST Webbook
dvisc	0.0003596	Paxs	405.10	Joback Method
dvisc	0.0004551	Paxs	375.24	Joback Method
dvisc	0.0005998	Paxs	345.39	Joback Method
dvisc	0.0008330	Paxs	315.53	Joback Method
dvisc	0.0012390	Paxs	285.67	Joback Method
dvisc	0.0020219	Paxs	255.82	Joback Method
dvisc	0.0037554	Paxs	225.96	Joback Method

Correlations

Information	Value
Property code	pvap
Equation	$\ln(P_{vp}) = A + B/(T + C)$
Coeff. A	1.66775e+01
Coeff. B	-4.93981e+03
Coeff. C	-1.25600e+01
Temperature range (K), min.	307.00
Temperature range (K), max.	447.17

Sources

McGowan Method:	http://link.springer.com/article/10.1007/BF02311772
NIST Webbook:	http://webbook.nist.gov/cgi/cbook.cgi?ID=C638299&Units=SI
The Yaws Handbook of Vapor Pressure:	https://www.sciencedirect.com/book/9780128029992/the-yaws-handbook-of-vapor-pressure
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

Legend

cpg: Ideal gas heat capacity

cpl:	Liquid phase 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
pvap:	Vapor 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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