

4-Toluoylacetonitrile

Inchi:	InChI=1S/C10H9NO/c1-8-2-4-9(5-3-8)10(12)6-7-11/h2-5H,6H2,1H3
InchiKey:	AIECDYDQPCANJK-UHFFFAOYSA-N
Formula:	C10H9NO
SMILES:	<chem>Cc1ccc(C(=O)CC#N)cc1</chem>
Mol. weight [g/mol]:	159.18
CAS:	7391-28-8

Physical Properties

Property code	Value	Unit	Source
gf	140.36	kJ/mol	Joback Method
hf	27.63	kJ/mol	Joback Method
hfus	18.41	kJ/mol	Joback Method
hvap	58.02	kJ/mol	Joback Method
log10ws	-2.89		Crippen Method
logp	2.091		Crippen Method
mvol	130.950	ml/mol	McGowan Method
pc	3005.73	kPa	Joback Method
tb	615.81	K	Joback Method
tc	847.63	K	Joback Method
tf	356.32	K	Joback Method
vc	0.519	m3/kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	300.20	J/molxK	615.81	Joback Method
cpg	311.13	J/molxK	654.45	Joback Method
cpg	321.30	J/molxK	693.08	Joback Method
cpg	330.75	J/molxK	731.72	Joback Method
cpg	339.51	J/molxK	770.36	Joback Method
cpg	347.62	J/molxK	809.00	Joback Method
cpg	355.11	J/molxK	847.63	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=C7391288&Units=SI
Crippen Method:	http://pubs.acs.org/doi/abs/10.1021/ci9903071

Legend

cpg:	Ideal gas heat capacity
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
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/79-616-1/4-Toluoylacetonitrile.pdf>

Generated by Cheméo on 2024-04-27 17:49:42.884721749 +0000 UTC m=+16529431.805299068.

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