3-Methoxyacetophenone

Other names:	1-(3-methoxyphenyl)ethanone
	1-Acetyl-3-methoxybenzene
	3'-methoxyacetophenone
	3-Acetylanisole
	3-Methoxy-1-acetylbenzene
	Acetophenone, 3'-methoxy-
	Acetophenone, m-methoxy-
	Ethanone, 1-(3-methoxyphenyl)-
	Methyl 3-methoxyphenyl ketone
	NSC 65593
	m-Acetanisole
	m-Acetylanisole
	m-Metoxyacetophenone
	m-methoxyacetophenone
	meta-Methoxyacetophenone
	methyl (3-methoxyphenyl) ketone
Inchi:	InChI=1S/C9H10O2/c1-7(10)8-4-3-5-9(6-8)11-2/h3-6H,1-2H3
InchiKey:	BAYUSCHCCGXLAY-UHFFFAOYSA-N
Formula:	C9H10O2
SMILES:	COc1cccc(C(C)=O)c1
Mol. weight [g/mol]:	150.17
CAS:	586-37-8

Physical Properties

Property code	Value	Unit	Source
affp	871.20	kJ/mol	NIST Webbook
basg	839.30	kJ/mol	NIST Webbook
gf	-106.24	kJ/mol	Joback Method
hf	-248.83	kJ/mol	Joback Method
hfus	15.50	kJ/mol	Joback Method
hvap	47.72	kJ/mol	Joback Method
ie	8.53 ± 0.05	eV	NIST Webbook
log10ws	-2.25		Crippen Method
logp	1.898		Crippen Method
mcvol	121.350	ml/mol	McGowan Method
pc	3380.21	kPa	Joback Method
rinpol	1321.00		NIST Webbook

rinpol	1279.00		NIST Webbook
rinpol	1295.00		NIST Webbook
rinpol	1295.00		NIST Webbook
rinpol	1321.00		NIST Webbook
rinpol	1301.00		NIST Webbook
rinpol	1282.00		NIST Webbook
rinpol	1277.00		NIST Webbook
rinpol	1279.00		NIST Webbook
rinpol	1298.00		NIST Webbook
ripol	2148.00		NIST Webbook
ripol	2148.00		NIST Webbook
tb	513.20	K	NIST Webbook
tc	730.64	K	Joback Method
tf	302.29	K	Joback Method
vc	0.456	m3/kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	256.24	J/mol×K	513.27	Joback Method
cpg	268.48	J/mol×K	549.50	Joback Method
cpg	280.07	J/mol×K	585.73	Joback Method
cpg	291.03	J/mol×K	621.95	Joback Method
cpg	301.35	J/mol×K	658.18	Joback Method
cpg	311.06	J/mol×K	694.41	Joback Method
cpg	320.16	J/mol×K	730.64	Joback Method
dvisc	0.0017319	Pa×s	302.29	Joback Method
dvisc	0.0010297	Pa×s	337.45	Joback Method
dvisc	0.0006754	Pa×s	372.62	Joback Method
dvisc	0.0004764	Pa×s	407.78	Joback Method
dvisc	0.0003552	Pa×s	442.94	Joback Method
dvisc	0.0002765	Pa×s	478.11	Joback Method
dvisc	0.0002227	Pa×s	513.27	Joback Method
hvapt	67.80	kJ/mol	298.15	Standard molar enthalpy of formation of methoxyacetophenone isomers

Pressure Dependent Properties

Property code	Value	Unit	Pressure [kPa]	Source
tbrp	398.70	K	1.60	NIST Webbook

Sources

Crippen Method:	http://pubs.acs.org/doi/abs/10.1021/ci990307l
Crippen Method:	https://www.chemeo.com/doc/models/crippen_log10ws
Standard molar enthalpy of formation of methoxyacetophenone isomers:	https://www.doi.org/10.1016/j.jct.2014.03.027
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=C586378&Units=SI

Legend

affp:	Proton affinity
basg:	Gas basicity
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
hvapt:	Enthalpy of vaporization at a given temperature
ie:	Ionization energy
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
ripol:	Polar retention indices
tb:	Normal Boiling Point Temperature
tbrp:	Boiling point at reduced pressure
tc:	Critical Temperature
tf:	Normal melting (fusion) point

vc: Critical Volume