

4'-Butoxyacetophenone

Other names:	4-n-Butoxyacetophenone p-Butoxyacetophenone Ethanone, 1-(4-butoxyphenyl)- Acetophenone, 4'-butoxy- 1-(4-butoxyphenyl)ethan-1-one
Inchi:	InChI=1S/C12H16O2/c1-3-4-9-14-12-7-5-11(6-8-12)10(2)13/h5-8H,3-4,9H2,1-2H3
InchiKey:	YYLCJWIQUFHMY-UHFFFAOYSA-N
Formula:	C12H16O2
SMILES:	CCCCOc1ccc(C(C)=O)cc1
Mol. weight [g/mol]:	192.25
CAS:	5736-89-0

Physical Properties

Property code	Value	Unit	Source
gf	-80.98	kJ/mol	Joback Method
hf	-310.75	kJ/mol	Joback Method
hfus	23.27	kJ/mol	Joback Method
hvap	54.40	kJ/mol	Joback Method
log10ws	-3.51		Crippen Method
logp	3.068		Crippen Method
mcvol	163.620	ml/mol	McGowan Method
pc	2480.12	kPa	Joback Method
tb	581.91	K	Joback Method
tc	788.90	K	Joback Method
tf	336.10	K	Joback Method
vc	0.624	m3/kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	393.77	J/molxK	581.91	Joback Method
cpg	408.69	J/molxK	616.41	Joback Method
cpg	422.81	J/molxK	650.91	Joback Method
cpg	436.14	J/molxK	685.41	Joback Method

cpg	448.71	J/molxK	719.90	Joback Method
cpg	460.52	J/molxK	754.40	Joback Method
cpg	471.60	J/molxK	788.90	Joback Method
dvisc	0.0017122	Paxs	336.10	Joback Method
dvisc	0.0009597	Paxs	377.07	Joback Method
dvisc	0.0006026	Paxs	418.04	Joback Method
dvisc	0.0004111	Paxs	459.00	Joback Method
dvisc	0.0002986	Paxs	499.97	Joback Method
dvisc	0.0002277	Paxs	540.94	Joback Method
dvisc	0.0001803	Paxs	581.91	Joback Method

Pressure Dependent Properties

Property code	Value	Unit	Pressure [kPa]	Source
tbrp	445.50 ± 0.50	K	2.40	NIST Webbook

Sources

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
NIST Webbook:	http://webbook.nist.gov/cgi/cbook.cgi?ID=C5736890&Units=SI

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
tb:	Normal Boiling Point Temperature
tbrp:	Boiling point at reduced pressure
tc:	Critical Temperature
tf:	Normal melting (fusion) point
vc:	Critical Volume

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