

3-(TRIFLUOROMETHYL)BENZOPHENONE

Other names:	Methanone, phenyl[3-(trifluoromethyl)phenyl]- m-Trifluoromethylbenzophenone
Inchi:	InChI=1S/C14H9F3O/c15-14(16,17)12-8-4-7-11(9-12)13(18)10-5-2-1-3-6-10/h1-9H
InchiKey:	IOXDAYKKVHAKSX-UHFFFAOYSA-N
Formula:	C14H9F3O
SMILES:	O=C(c1ccccc1)c1cccc(C(F)(F)F)c1
Mol. weight [g/mol]:	250.22

Physical Properties

Property code	Value	Unit	Source
ea	0.96 ± 0.09	eV	NIST Webbook
hf	-622.88	kJ/mol	Cheméo Relay (1.0)
hfus	23.13	kJ/mol	Joback Method
hvap	84.84	kJ/mol	Cheméo Relay (1.0)
ie	9.68	eV	Cheméo Relay (1.0)
log10ws	-4.23		Cheméo Relay (1.0)
logp	3.936		Crippen Method
mcvol	167.480	ml/mol	McGowan Method
pc	2581.96	kPa	Joback Method
tb	572.95	K	Cheméo Relay (1.0)
tc	803.25	K	Cheméo Relay (1.0)
tf	333.10	K	Cheméo Relay (1.0)

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	417.80	J/mol×K	626.51	Joback Method
cpg	431.75	J/mol×K	664.36	Joback Method
cpg	444.51	J/mol×K	702.20	Joback Method
cpg	456.17	J/mol×K	740.05	Joback Method
cpg	466.81	J/mol×K	777.89	Joback Method
cpg	476.52	J/mol×K	815.74	Joback Method
cpg	485.37	J/mol×K	853.58	Joback Method

Pressure Dependent Properties

Property code	Value	Unit	Pressure [kPa]	Source
tbrp	371.50 ± 0.50	K	0.05	NIST Webbook

Sources

Cheméo Relay (1.0, beta):

NIST Webbook: <http://webbook.nist.gov/cgi/cbook.cgi?ID=C728814&Units=SI>

Joback Method: https://en.wikipedia.org/wiki/Joback_method

McGowan Method: <http://link.springer.com/article/10.1007/BF02311772>

Crippen Method: <http://pubs.acs.org/doi/abs/10.1021/ci990307l>

Crippen Method: https://www.chemeo.com/doc/models/crippen_log10ws

Cheméo Relay (1.0):

Legend

cpg:	Ideal gas heat capacity
ea:	Electron affinity
hf:	Enthalpy of formation at standard conditions
hfus:	Enthalpy of fusion at standard conditions
hvap:	Enthalpy of vaporization at standard conditions
ie:	Ionization energy
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

Latest version available from:

<https://www.chemeo.com/cid/50-236-4/3-TRIFLUOROMETHYL-BENZOPHENONE.pdf>

Generated by Cheméo on 2026-07-21 15:19:15.944768193 +0000 UTC m=+158251.786653571.

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