

# Aniline, 2-nitro-4-trifluoromethyl-

<b>Other names:</b>	p-Toluidine, «alpha», «alpha», «alpha»-trifluoro-2-nitro-4-Amino-3-nitrobenzotrifluoride 4-Trifluoromethyl 2-nitroaniline «alpha», «alpha», «alpha»-Trifluoro-2-nitro-p-toluidine Benzenamine, 2-nitro-4-(trifluoromethyl)- 2-Nitro-4-(trifluoromethyl)aniline
<b>Inchi:</b>	InChI=1S/C7H5F3N2O2/c8-7(9,10)4-1-2-5(11)6(3-4)12(13)14/h1-3H,11H2
<b>InchiKey:</b>	ATXBGHLILIABGX-UHFFFAOYSA-N
<b>Formula:</b>	C7H5F3N2O2
<b>SMILES:</b>	<chem>Nc1ccc(C(F)(F)F)cc1[N+](=O)[O-]</chem>
<b>Mol. weight [g/mol]:</b>	206.12
<b>CAS:</b>	400-98-6

## Physical Properties

Property code	Value	Unit	Source
gf	-378.38	kJ/mol	Joback Method
hf	-548.27	kJ/mol	Joback Method
hfus	25.53	kJ/mol	Joback Method
hvap	58.26	kJ/mol	Joback Method
log10ws	-2.86		Crippen Method
logp	2.196		Crippen Method
mcvol	118.440	ml/mol	McGowan Method
pc	3736.23	kPa	Joback Method
tb	615.15	K	Joback Method
tc	851.35	K	Joback Method
tf	451.17	K	Joback Method
vc	0.473	m3/kmol	Joback Method

## Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	298.10	J/molxK	615.15	Joback Method
cpg	307.46	J/molxK	654.52	Joback Method
cpg	315.99	J/molxK	693.88	Joback Method

cpg	323.75	J/mol×K	733.25	Joback Method
cpg	330.80	J/mol×K	772.62	Joback Method
cpg	337.20	J/mol×K	811.98	Joback Method
cpg	343.01	J/mol×K	851.35	Joback Method

## Sources

<b>Crippen Method:</b>	<a href="http://pubs.acs.org/doi/abs/10.1021/ci9903071">http://pubs.acs.org/doi/abs/10.1021/ci9903071</a>
<b>Crippen Method:</b>	<a href="https://www.chemeo.com/doc/models/crippen_log10ws">https://www.chemeo.com/doc/models/crippen_log10ws</a>
<b>Joback Method:</b>	<a href="https://en.wikipedia.org/wiki/Joback_method">https://en.wikipedia.org/wiki/Joback_method</a>
<b>McGowan Method:</b>	<a href="http://link.springer.com/article/10.1007/BF02311772">http://link.springer.com/article/10.1007/BF02311772</a>
<b>NIST Webbook:</b>	<a href="http://webbook.nist.gov/cgi/cbook.cgi?ID=C400986&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=C400986&amp;Units=SI</a>

## Legend

<b>cpg:</b>	Ideal gas heat capacity
<b>gf:</b>	Standard Gibbs free energy of formation
<b>hf:</b>	Enthalpy of formation at standard conditions
<b>hfus:</b>	Enthalpy of fusion at standard conditions
<b>hvap:</b>	Enthalpy of vaporization at standard conditions
<b>log10ws:</b>	Log10 of Water solubility in mol/l
<b>logp:</b>	Octanol/Water partition coefficient
<b>mcvol:</b>	McGowan's characteristic volume
<b>pc:</b>	Critical Pressure
<b>tb:</b>	Normal Boiling Point Temperature
<b>tc:</b>	Critical Temperature
<b>tf:</b>	Normal melting (fusion) point
<b>vc:</b>	Critical Volume

Latest version available from:

<https://www.chemeo.com/cid/42-630-5/Aniline-2-nitro-4-trifluoromethyl.pdf>

Generated by Cheméo on 2024-04-23 15:56:04.729292974 +0000 UTC m=+16177013.649870288.

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