

Ethanol, 2-[ethyl(3-methylphenyl)amino]-

Other names:	Ethanol, 2-(N-ethyl-m-toluidino)- 2-(N-Ethyl-m-toluidino)ethanol 2-(N-Ethyl-meta-toluidino)-ethanol N-Hydroxyethyl-N-ethyl-m-toluidine 3-Methyl-N-ethyl-N-«beta»-hydroxyethylaniline N-Ethyl-N-(2-hydroxyethyl)-m-toluidine N-(«beta»-Hydroxyethyl)-N-ethyl-m-toluidine N-Ethyl-N-hydroxyethyl-meta-toluidine Emery 5714 NSC 89746 2-(N-Ethyl-N-toluidino)ethanol
Inchi:	InChI=1S/C11H17NO/c1-3-12(7-8-13)11-6-4-5-10(2)9-11/h4-6,9,13H,3,7-8H2,1-2H3
InchiKey:	KRNUKKZDGDABWF-UHFFFAOYSA-N
Formula:	C11H17NO
SMILES:	CCN(CCO)c1cccc(C)c1
Mol. weight [g/mol]:	179.26
CAS:	91-88-3

Physical Properties

Property code	Value	Unit	Source
gf	118.48	kJ/mol	Joback Method
hf	-130.01	kJ/mol	Joback Method
hfus	25.01	kJ/mol	Joback Method
hvap	61.74	kJ/mol	Joback Method
log10ws	-1.96		Crippen Method
logp	1.814		Crippen Method
mcvol	157.940	ml/mol	McGowan Method
pc	2865.80	kPa	Joback Method
rinpol	1546.20		NIST Webbook
tb	587.36	K	Joback Method
tc	776.23	K	Joback Method
tf	345.96	K	Joback Method
vc	0.581	m3/kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	394.12	J/molxK	587.36	Joback Method
cpg	407.62	J/molxK	618.84	Joback Method
cpg	420.38	J/molxK	650.32	Joback Method
cpg	432.45	J/molxK	681.79	Joback Method
cpg	443.83	J/molxK	713.27	Joback Method
cpg	454.58	J/molxK	744.75	Joback Method
cpg	464.72	J/molxK	776.23	Joback Method

Pressure Dependent Properties

Property code	Value	Unit	Pressure [kPa]	Source
tbrp	387.70	K	0.10	NIST Webbook

Sources

NIST Webbook:	http://webbook.nist.gov/cgi/cbook.cgi?ID=C91883&Units=SI
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

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
rinpol:	Non-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

Latest version available from:

<https://www.cheméo.com/cid/61-655-7/Ethanol-2-ethyl-3-methylphenyl-amino.pdf>

Generated by Cheméo on 2024-04-28 19:41:04.534630288 +0000 UTC m=+16622513.455207610.

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