

Aziprotryne

Other names:

1,3,5-Triazin-2-amine, 4-azido-N-(1-methylethyl)-6-(methylthio)-s-Triazine, 2-azido-4-(isopropylamino)-6-(methylthio)-Azeprotryne
2-Azido-4-isopropylamino-6-methylthio-s-triazine
2-Azido-4-isopropylamino-6-methylthio-1,3,5-triazine
4-Azido-N-(1-methylethyl)-6-(methylthio)-1,3,5-triazin-2-amine
Aziprotryn
Azirpotryne
Brasoran
Brassoron
C 7019
CIBA C 7019
Isopropylamino-4-azido-6-methylthio-1,3,5-triazin
Mesoranil
Mezaronil
Mezuron

Inchi:

InChI=1S/C7H11N7S/c1-4(2)9-5-10-6(13-14-8)12-7(11-5)15-3/h4H,1-3H3,(H,9,10,11,12)

InchiKey:

AFIIBUOYKYSPKB-UHFFFAOYSA-N

Formula:

C7H11N7S

SMILES:

CSc1nc(N=[N+]=[N-])nc(NC(C)C)n1

Mol. weight [g/mol]:

225.27

CAS:

4658-28-0

Physical Properties

Property code	Value	Unit	Source
log10ws	-8.13		Crippen Method
logp	2.356		Crippen Method
mcvol	163.340	ml/mol	McGowan Method
rinpol	1766.00		NIST Webbook
rinpol	1779.00		NIST Webbook
rinpol	1796.00		NIST Webbook
rinpol	1766.00		NIST Webbook

Sources

Crippen Method:	http://pubs.acs.org/doi/abs/10.1021/ci990307I
Crippen Method:	https://www.chemeo.com/doc/models/crippen_log10ws
McGowan Method:	http://link.springer.com/article/10.1007/BF02311772
NIST Webbook:	http://webbook.nist.gov/cgi/cbook.cgi?ID=C4658280&Units=SI

Legend

log10ws:	Log10 of Water solubility in mol/l
logp:	Octanol/Water partition coefficient
mcvol:	McGowan's characteristic volume
rinpol:	Non-polar retention indices

Latest version available from:

<https://www.chemeo.com/cid/44-612-3/Aziprotryne.pdf>

Generated by Cheméo on 2024-04-24 21:54:57.804006626 +0000 UTC m=+16284946.724583941.

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