

# Morpholine, 4-(2-benzothiazolythio)-

**Other names:**

Benzothiazole, 2-(morpholinothio)-  
2-(Morpholinothio)benzothiazole  
Amax  
Benzothiazole, 2-(4-morpholinylthio)-  
Benzothiazolyl-2-sulfenmorpholide  
Morpholinylmercaptobenzothiazole  
N-(Oxydiethylene)-2-benzothiazolesulfenamide  
N-(Oxydiethylene)-2-benzothiazolesulfenamide  
N-(Oxydiethylene)benzothiazole-2-sulfenamide  
N-(Oxydiethylene)benzothiazylsulfenamide  
N,N-(Oxydiethylene)-2-benzothiazolesulfenamide  
N,N-(Oxydiethylene)-2-benzothiazylsulfenamide  
N,N-(Oxydiethylene)benzothiazole-2-sulfenamide  
NOBS Special  
Santocure MOR  
Sulfenamide M  
Sulfenax MOR  
Vulcafor BSM  
Vulkacit MOZ  
2-(4-Morpholinothio)benzothiazole  
2-(4-Morpholinylmercapto)benzothiazole  
2-Benzothiazolesulfenamide, N-morpholinyl-  
2-Benzothiazolesulfenemorpholide  
2-Benzothiazolyl N-morpholino sulfide  
2-Benzothiazolylsulfenyl morpholine  
4-(2-Benzothiazolythio)morpholine  
USAF CY-7  
2-(4-Morpholinylthio)benzothiazole  
N-Oxydiethyl-2-benzthiazolsulfenamid  
Sulfenax mob  
2-(Morpholinthio)-benzothiazole  
Accelerator NC  
Delac MOR  
OBTS  
MBS  
Meramide M  
Accel NS  
NSC 70078

**Inchi:**

InChI=1S/C11H12N2OS2/c1-2-4-10-9(3-1)12-11(15-10)16-13-5-7-14-8-6-13/h1-4H,5-8H

**InchiKey:**

MHKLKWCYGIBEQF-UHFFFAOYSA-N

**Formula:** C11H12N2OS2  
**SMILES:** c1ccc2sc(SN3CCOCC3)nc2c1  
**Mol. weight [g/mol]:** 252.36  
**CAS:** 102-77-2

## Physical Properties

Property code	Value	Unit	Source
log10ws	-3.34		Crippen Method
logp	2.636		Crippen Method
mcvol	174.600	ml/mol	McGowan Method
tf	359.45 ± 0.35	K	NIST Webbook
tf	360.05 ± 0.20	K	NIST Webbook

## Sources

**NIST Webbook:** <http://webbook.nist.gov/cgi/cbook.cgi?ID=C102772&Units=SI>  
**Crippen Method:** <http://pubs.acs.org/doi/abs/10.1021/ci990307l>  
**Crippen Method:** [https://www.chemeo.com/doc/models/crippen\\_log10ws](https://www.chemeo.com/doc/models/crippen_log10ws)  
**McGowan Method:** <http://link.springer.com/article/10.1007/BF02311772>

## Legend

**log10ws:** Log10 of Water solubility in mol/l  
**logp:** Octanol/Water partition coefficient  
**mcvol:** McGowan's characteristic volume  
**tf:** Normal melting (fusion) point

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