

# Benzenesulfonic acid, hydrazide

Other names:

Benzenesulfohydrazide  
Benzenesulfonic hydrazide  
Benzenesulfonohydrazide  
Benzenesulfonyl hydrazide  
Benzenesulfonylhydrazine  
Benzensulfonylhydrazine  
Celogen BSH  
ChKhZ 9  
Genitron BSH  
Phenylsulfohydrazide  
Phenylsulfonyl hydrazide  
Phenylsulfonylhydrazine  
Porofo BSH  
Porofo ChKhZ 9  
Porofo-BSH-Pulver  
BSH  
Porofo  
Benzene sulphonohydrazide  
Hydrazide BSG  
Nitropore OBSH  
NSC 643

Inchi:

InchiKey:

Formula:

SMILES:

Mol. weight [g/mol]:

CAS:

InChI=1S/C6H8N2O2S/c7-8-11(9,10)6-4-2-1-3-5-6/h1-5,8H,7H2

VJRITMATACIYAF-UHFFFAOYSA-N

C6H8N2O2S

NNS(=O)(=O)c1ccccc1

172.21

80-17-1

## Physical Properties

| Property code | Value    | Unit   | Source         |
|---------------|----------|--------|----------------|
| chs           | -3607.00 | kJ/mol | NIST Webbook   |
| gf            | -200.65  | kJ/mol | Joback Method  |
| hf            | -296.73  | kJ/mol | Joback Method  |
| hfus          | 27.01    | kJ/mol | Joback Method  |
| hvap          | 66.94    | kJ/mol | Joback Method  |
| log10ws       | -1.35    |        | Crippen Method |

|       |         |         |                |
|-------|---------|---------|----------------|
| logp  | -0.161  |         | Crippen Method |
| mcvol | 119.690 | ml/mol  | McGowan Method |
| pc    | 6524.67 | kPa     | Joback Method  |
| tb    | 533.84  | K       | Joback Method  |
| tc    | 756.16  | K       | Joback Method  |
| tf    | 358.28  | K       | Joback Method  |
| vc    | 0.454   | m3/kmol | Joback Method  |

## Temperature Dependent Properties

| Property code | Value  | Unit    | Temperature [K] | Source        |
|---------------|--------|---------|-----------------|---------------|
| cpg           | 261.94 | J/mol×K | 533.84          | Joback Method |
| cpg           | 273.39 | J/mol×K | 570.89          | Joback Method |
| cpg           | 284.06 | J/mol×K | 607.95          | Joback Method |
| cpg           | 293.96 | J/mol×K | 645.00          | Joback Method |
| cpg           | 303.10 | J/mol×K | 682.06          | Joback Method |
| cpg           | 311.50 | J/mol×K | 719.11          | Joback Method |
| cpg           | 319.17 | J/mol×K | 756.16          | Joback Method |

## Sources

|                 |                                                                                                                                         |
|-----------------|-----------------------------------------------------------------------------------------------------------------------------------------|
| Crippen Method: | <a href="http://pubs.acs.org/doi/abs/10.1021/ci990307l">http://pubs.acs.org/doi/abs/10.1021/ci990307l</a>                               |
| Crippen Method: | <a href="https://www.chemeo.com/doc/models/crippen_log10ws">https://www.chemeo.com/doc/models/crippen_log10ws</a>                       |
| Joback Method:  | <a href="https://en.wikipedia.org/wiki/Joback_method">https://en.wikipedia.org/wiki/Joback_method</a>                                   |
| McGowan Method: | <a href="http://link.springer.com/article/10.1007/BF02311772">http://link.springer.com/article/10.1007/BF02311772</a>                   |
| NIST Webbook:   | <a href="http://webbook.nist.gov/cgi/cbook.cgi?ID=C80171&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=C80171&amp;Units=SI</a> |

## Legend

|       |                                                 |
|-------|-------------------------------------------------|
| chs:  | Standard solid enthalpy of combustion           |
| 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 |

|                 |                                     |
|-----------------|-------------------------------------|
| <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                     |

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