

# 2-Naphthalenol, 1-[(2,4-dimethylphenyl)azo]-

## Other names:

AF Red No. 5  
Brasilazina Oil Scarlet 6G  
Calco Oil Scarlet BL  
Ceres Orange RR  
Cerisol Scarlet G  
C.I. 12140  
C.I. Solvent Orange 7  
Ext D and C Red No. 14  
Fast Oil Orange II  
Fat Scarlet 2G  
FD And C red No. 32  
Grasan Orange 3R  
Japan Red 5  
Japan Red 505  
Japan Red No.5  
Lacquer Orange VR  
Oil Orange KB  
Oil Orange N Extra  
Oil Orange R  
Oil Orange 2R  
Oil Orange X  
Oil Orange XO  
Oil Red XO  
Oil Scarlet  
Oil Scarlet 371  
Oil Scarlet BL  
Oil Scarlet 6G  
Oil Scarlet YS  
Red No. 5  
Resin Scarlet 2R  
Somalia Orange A 2R  
Somalia Orange 2R  
Sudan II  
Sudan orange  
Sudan Orange RPA  
Sudan Orange RRA  
Sudan Red  
A.F. Red No. 5  
Aizen Food Red No. 5  
Brilliant oil scarlet B

Calco oil scarlet ZBL  
 Ceres oranges RR  
 Cerotinscharlach G  
 Ext. D and C Red No. 14  
 Extract D & C Red No. 14  
 Fat Red (Yellowish)  
 Fettorange B  
 Motirot G  
 Oil scarlet APYO  
 Oil scarlet L  
 Oil scarlet Y  
 Oil Red Gro  
 Oil Red RO  
 Orange Insoluble RR  
 Orange Oil KB  
 Ponceau insoluble olg  
 Pyronalrot R  
 Red B  
 Resoform Orange R  
 Rot B  
 Rot GG fettloeslich  
 Solvent Orange 7  
 Soudan II  
 Sudan ax  
 Sudan scarlet 6G  
 Sudan X  
 Waxakol vermilion L  
 1-(o-Xylylazo)-2-naphthol  
 1-(2,4-Xylylazo)-2-naphthol  
 1-((2,4-Dimethylphenyl)azo)-2-naphthalenol  
 1-Xylylazo-2-naphthol  
 2-Naphthol, 1-(2,4-xylylazo)-  
 Ext. D and C. Red. No. 14  
 Oranz rozpoustedlova 7  
 Sudan 2  
 1-(2,4-Dimethylphenylazo)-2-naphthol  
 1-[(2,4-Dimethylphenyl)diazenyl]-2-naphthol  
 Red 505  
 2-Naphthalenol, 1-[2-(2,4-dimethylphenyl)diazenyl]-  
**Inchi:** InChI=1S/C18H16N2O/c1-12-7-9-16(13(2)11-12)19-20-18-15-6-4-3-5-14(15)8-10-17(18)  
**InchiKey:** JBTHDAVBDDKSRW-UHFFFAOYSA-N  
**Formula:** C18H16N2O  
**SMILES:** Cc1ccc(N=Nc2c(O)ccc3ccccc23)c(C)c1

Mol. weight [g/mol]: 276.33  
CAS: 3118-97-6

## Physical Properties

Property code	Value	Unit	Source
chs	-9395.00 ± 20.00	kJ/mol	NIST Webbook
chs	-9576.80	kJ/mol	NIST Webbook
hf	84.78	kJ/mol	Joback Method
hfs	25.00 ± 20.00	kJ/mol	NIST Webbook
hfs	207.00	kJ/mol	NIST Webbook
hvap	83.52	kJ/mol	Joback Method
log10ws	-5.82		Crippen Method
logp	5.578		Crippen Method
mcvol	219.030	ml/mol	McGowan Method
pc	2071.76	kPa	Joback Method
tb	928.34	K	Joback Method
tc	1198.45	K	Joback Method

## Sources

**NIST Webbook:** <http://webbook.nist.gov/cgi/cbook.cgi?ID=C3118976&Units=SI>  
**Crippen Method:** <http://pubs.acs.org/doi/abs/10.1021/ci9903071>  
**Crippen Method:** [https://www.chemeo.com/doc/models/crippen\\_log10ws](https://www.chemeo.com/doc/models/crippen_log10ws)  
**Joback Method:** [https://en.wikipedia.org/wiki/Joback\\_method](https://en.wikipedia.org/wiki/Joback_method)  
**McGowan Method:** <http://link.springer.com/article/10.1007/BF02311772>

## Legend

**chs:** Standard solid enthalpy of combustion  
**hf:** Enthalpy of formation at standard conditions  
**hfs:** Solid phase enthalpy of formation 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  
**tb:** Normal Boiling Point Temperature  
**tc:** Critical Temperature

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