1,2,3,6,8-pentabromo-dibenzofuran

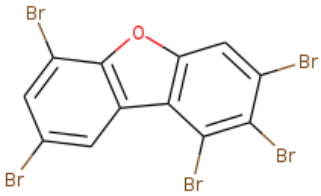
InChI: InChI=1S/C12H3Br5O/c13-4-1-5-9-8(18-12(5)7(15)2-4)3-6(14)10(16)11(9)17/h1-3H

InChI Key: OWGWMEKSWJIYNM-UHFFFAOYSA-N

Formula: C12H3Br5O

SMILES: Brc1cc2c(oc3cc(Br)c(Br)c(Br)c32)c(Br)c1

Molecular Weight: 562.67



Physical Properties

Property	Value	Unit	Source
$logP_{oct/wat}$	7.399		Crippen Method

Sources

NIST Webbook: http://webbook.nist.gov/cgi/inchi/InChI=1S/C12H3Br5O/c13-4-1-5-9-8(18-12(5)7(15)2-4)3-6(14)10(16)11(9)17/h1-3H

Crippen Method: http://pubs.acs.org/doi/abs/10.1021/ci990307l

Legend

***log*P$_{oct/wat}$**: Octanol/Water partition coefficient .