

VERA: Similarity and Grouping:



PREMIER

PRIORITISATION AND RISK EVALUATION
OF MEDICINES IN THE ENVIRONMENT

Molecule Input:

TotGroups: 2

['benzene', 'Ar_Cl_Br']

LogP: 5.314

MaxPartialCharge: 0.078

MinPartialCharge: -0.083

NumHAcceptors: 0

NumHDonors: 0

NumRotatableBonds: 1

NumAliphaticRings: 0

NumAromaticCarbocycles: 2

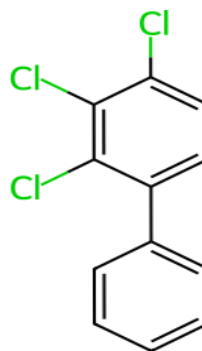
NumAromaticHeterocycles: 0

NumAromaticRings: 2

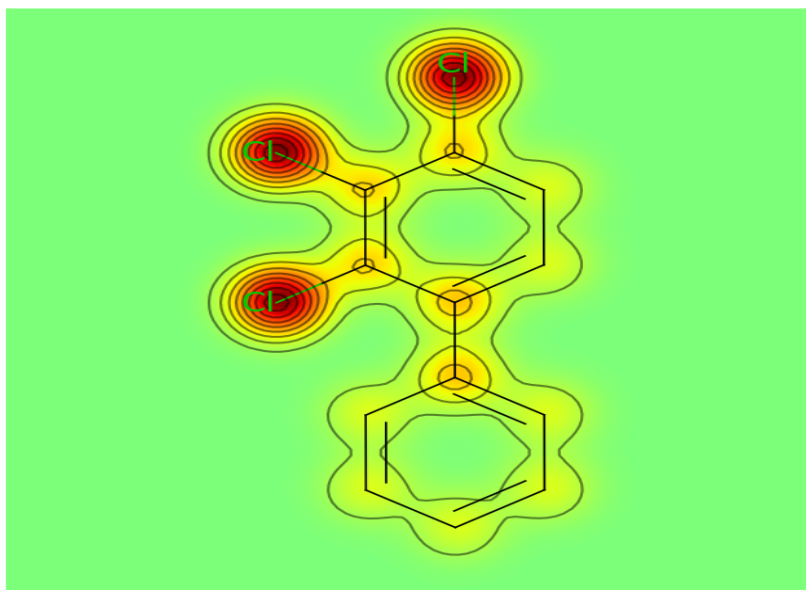
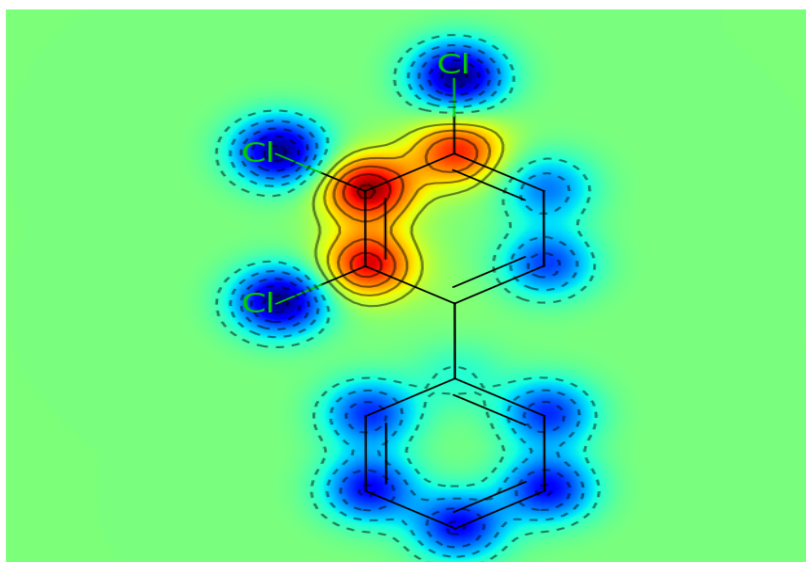
ExactMolWt: 255.961

GasteigerCharges:

CrippenContribs:



Prediction from Linear model: 0.048 -log(mg/l)



Molecules with the highest group similarity

1) SMILES: Clc1ccc(Cl)c(-c2ccccc2Cl)c1

ExpValue: 1.472 -log(mg/l)

SimilarityVega: 0.975

SimilarityGrouping: 0.875

Grp_Similarity_Mean: 0.925

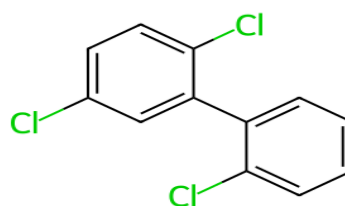
Groups in common: 3

TotGroups 3

GROUPS:

Ar_Cl_Br Stat: Tox

Alerts: []



Toxicity

2) SMILES: Clc1cccc(Cl)c1Cl

ExpValue: -0.338 -log(mg/l)

SimilarityVega: 0.781

SimilarityGrouping: 0.875

Grp_Similarity_Mean: 0.828

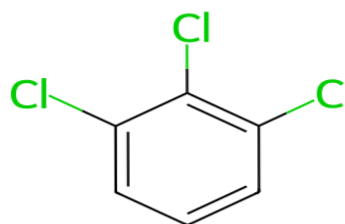
Groups in common: 3

TotGroups 3

GROUPS:

Ar_Cl_Br Stat: Tox

Alerts: []



Toxicity

3) SMILES: Clc1ccc(Cl)c(Cl)c1

ExpValue: -0.412 -log(mg/l)

SimilarityVega: 0.78

SimilarityGrouping: 0.875

Grp_Similarity_Mean: 0.828

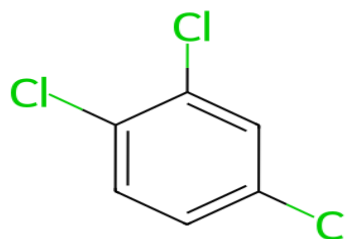
Groups in common: 3

TotGroups 3

GROUPS:

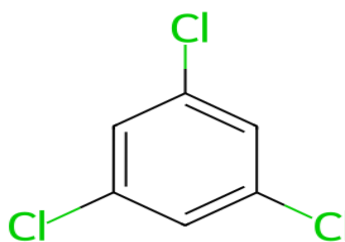
Ar_Cl_Br Stat: Tox

Alerts: []

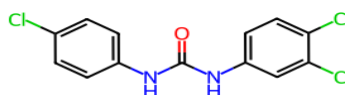


Toxicity

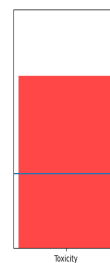
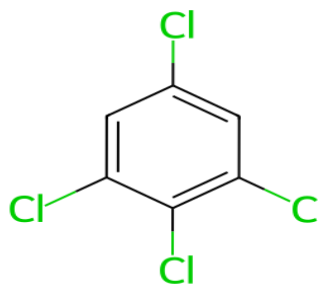
4) SMILES: Clc1cc(Cl)cc(Cl)c1
 ExpValue: -0.512 -log(mg/l)
 SimilarityVega: 0.773
 SimilarityGrouping: 0.875
 Grp_Similarity_Mean: 0.824
 Groups in common: 3
 TotGroups 3
 GROUPS:
 Ar_Cl_Br Stat: Tox
 Alerts: []



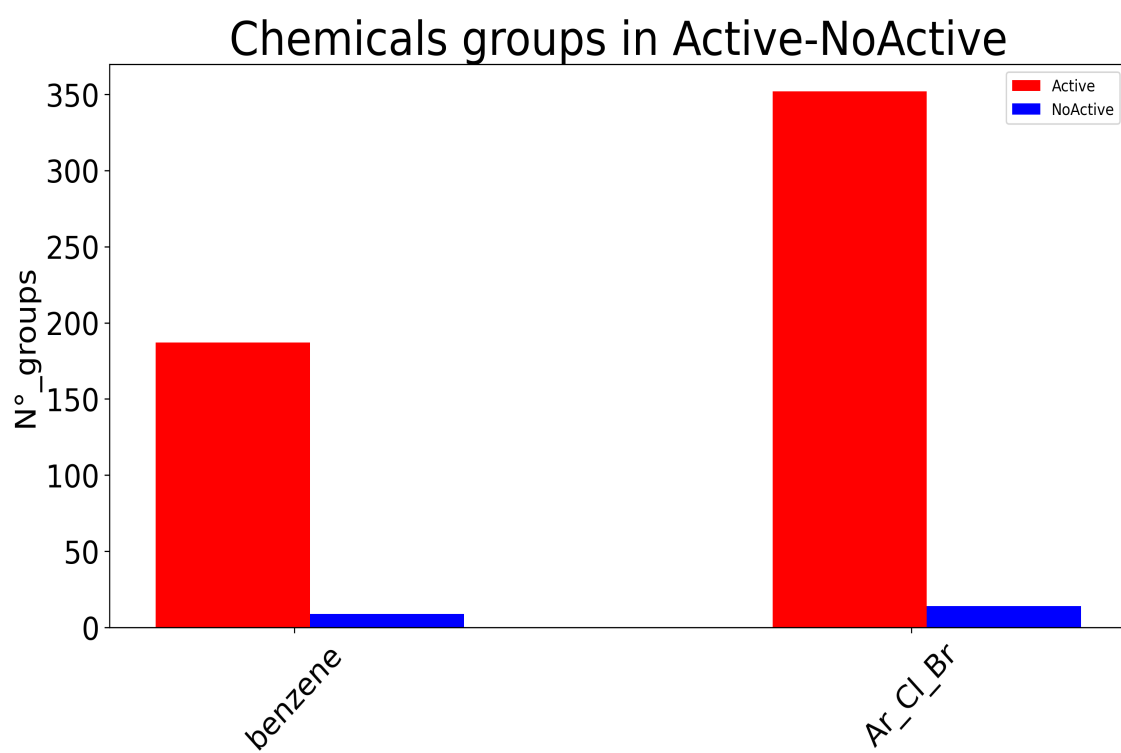
5) SMILES: O=C(Nc1ccc(Cl)cc1)Nc1ccc(Cl)c(Cl)c1
 ExpValue: 1.043 -log(mg/l)
 SimilarityVega: 0.76
 SimilarityGrouping: 0.844
 Grp_Similarity_Mean: 0.802
 Groups in common: 3
 TotGroups 4
 GROUPS:
 urea Stat: Tox
 Ar_Cl_Br Stat: Tox
 Alerts: []



6) SMILES: Clc1cc(Cl)c(Cl)c(Cl)c1
 ExpValue: 0.098 -log(mg/l)
 SimilarityVega: 0.758
 SimilarityGrouping: 0.844
 Grp_Similarity_Mean: 0.801
 Groups in common: 3
 TotGroups 4
 GROUPS:
 Ar_Cl_Br Stat: Tox
 Alerts: []



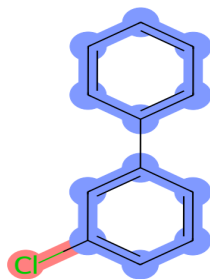
Molecule Input: Bar Chart for groups in Dataset find in target



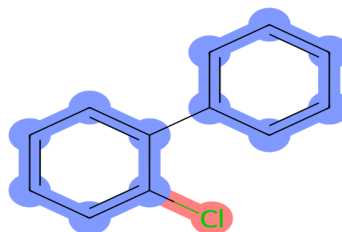
Reasoning

Orthogonal Research based on copresence of group with high prevalence Ar_Cl_Br and GRP benzene

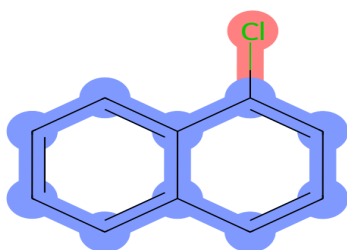
Smile: Clc1cccc(-c2ccccc2)c1
 ExpValue: -0.894 -log(mg/l)
 Similarity: 0.864



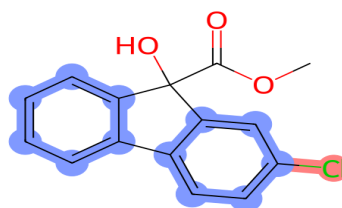
Smile: Clc1ccccc1-c1ccccc1
 ExpValue: -0.604 -log(mg/l)
 Similarity: 0.864



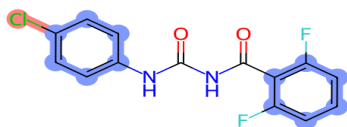
Smile: Clc1cccc2ccccc12
 ExpValue: -0.23 -log(mg/l)
 Similarity: 0.761



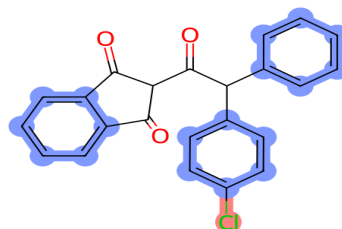
Smile: COC(=O)C1(O)c2ccccc2-c2ccc(Cl)cc21
 ExpValue: -0.158 -log(mg/l)
 Similarity: 0.749



Smile: O=C(NC(=O)c1c(F)cccc1F)Nc1ccc(Cl)cc1
 ExpValue: -2.631 -log(mg/l)
 Similarity: 0.722

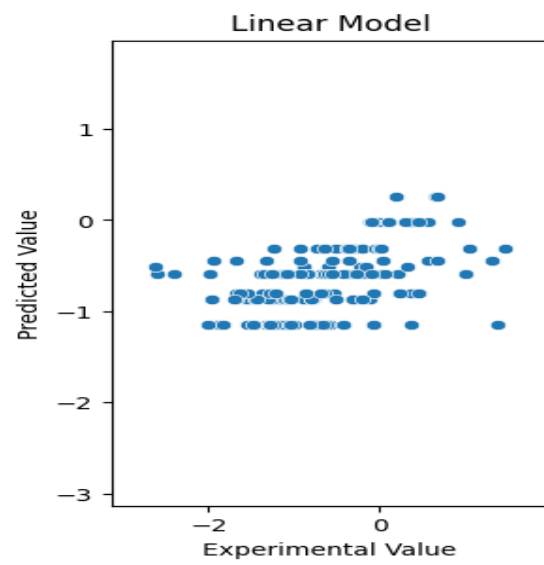


Smile: O=C1c2ccccc2C(=O)C1C(=O)C(c1ccccc1)c1ccc(Cl)cc1
 ExpValue: 0.328 -log(mg/l)
 Similarity: 0.678



Local Linear Model

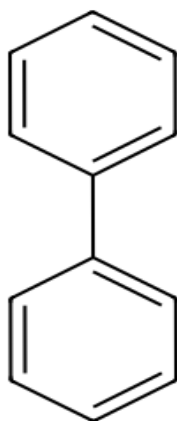
The target assessment was made by using Linear model model not by this Local model



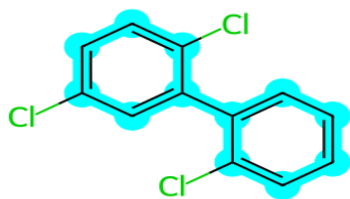
Mean square error: 0.46

Scaffold

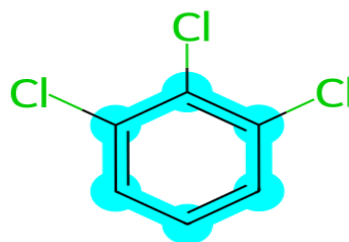
Scaffold analysis



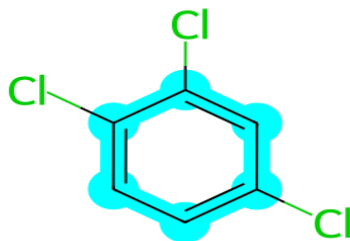
Smile: Clc1ccc(Cl)c(-c2ccccc2Cl)c1
ExpValue: 1.472 -log(mg/l)
Similarity: 0.975



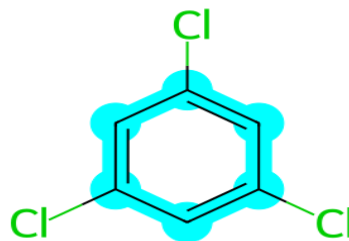
Smile: Clc1cccc(Cl)c1Cl
ExpValue: -0.338 -log(mg/l)
Similarity: 0.781



Smile: Clc1ccc(Cl)c(Cl)c1
ExpValue: -0.412 -log(mg/l)
Similarity: 0.78



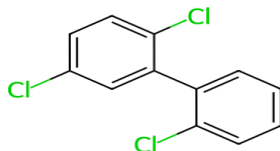
Smile: Clc1cc(Cl)cc(Cl)c1
ExpValue: -0.512 -log(mg/l)
Similarity: 0.773



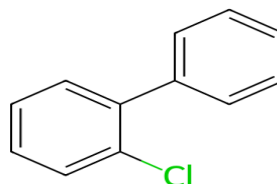
Descriptors

Calculator Similarity descriptor

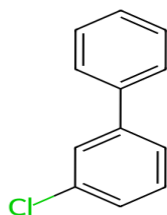
Smile: Clc1ccc(Cl)c(-c2ccccc2Cl)c1
 ExpValue: 1.472 -log(mg/l)
 DistanceCityBlock: 0.285
 LogP:5.313800000000002
 MW: 255.961333264



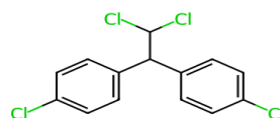
Smile: Clc1ccccc1-c1ccccc1
 ExpValue: -0.604 -log(mg/l)
 DistanceCityBlock: 2.931
 LogP:4.007000000000002
 MW: 188.039277968



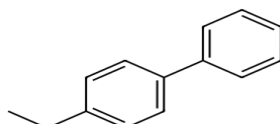
Smile: Clc1cccc(-c2ccccc2)c1
 ExpValue: -0.894 -log(mg/l)
 DistanceCityBlock: 3.002
 LogP:4.007000000000002
 MW: 188.039277968



Smile: Clc1ccc(C(c2ccc(Cl)cc2)C(Cl)Cl)cc1
 ExpValue: -0.642 -log(mg/l)
 DistanceCityBlock: 3.631
 LogP:5.929000000000002
 MW: 317.95366104



Smile: CCc1ccc(-c2ccccc2)cc1
 ExpValue: 0.223 -log(mg/l)
 DistanceCityBlock: 4.554
 LogP:3.916000000000002
 MW: 182.109550448



Smile: Clc1cccc2ccccc12
 ExpValue: -0.23 -log(mg/l)
 DistanceCityBlock: 4.638
 LogP:3.4932000000000016
 MW: 162.023627904

