## **VOA Compound List Grouped by Surrogates**

<u>Surrogate Definition:</u> Surrogates are compounds similar in chemical composition to the analytes of interest and spiked into environmental samples prior to preparation and analysis. They are used to evaluate extraction efficiency and matrix interference on a sample-specific basis. In order for this to work, surrogates must behave in the same manner as the corresponding analytes. Unfortunately, in practice, this equivalency is typically difficult to demonstrate and is often more assumed than empirically derived.

The matrix spike recovery is the best indicator of how a specific analyte performs in your samples. Unfortunately matrix spikes aren't performed on all sample. Surrogate recoveries will give you an overall indication of accuracy but are not analyte specific. When determining how a surrogate recovery may affect specific analytes one should consider the following:

- 1. Check associated blank and BS/BSD data. These surrogates should be within control limits. If any of these surrogates are outside of control and the same surrogates are also outside of control limits in the associated samples then the problem is procedural and not the sample matrix. The batch should be reanalyzed.
- 2. All surrogates within control limits: Even though all surrogates worked within control limits, one cannot assume all analytes worked within control limits. Some analytes are notoriously poor performers (PP). Review BS/BSD and MS/MSD data. Not all analyte types are representated by surrogates.
- 3. One or more surrogates outside of control limits: For RCP and MCP work, when a surrogate is outside of control limits reanalysis is required, unless there is obvious interference from visual inspection of the chromatogram, to ensure the problem was not due to an instrument error. If reanalysis produces the same results then matrix interference is confirmed and results are reported with surrogates outside of control limits. Some things to consider:
- 1. Does the matrix or interference only affect the surrogate? (Rare)
- 2. Does the matrix or interference only affect analytes in the part of the chromatogram with the low surrogate? (See Project Narrative for details)
- 3. Is only a certain type of analyte affected by matrix or interference? Not every type of compound is represented with a surrogate. (Example: Ketones (K) not represented)

There are 4 VOA surrogates:

These surrogates best represent the halogenated hydrocarbons (H). **Dibromofluoromethane**(SURR) is a halomethane. **1,2-Dichloroethane-d4**(SURR) is a chlorinated hydrocarbon.

These surrogates best represents the aromatics (A)

Toluene-d8 (SURR) is an aromatic hydrocarbon.

Bromofluorobenzene(Surr) is a halogenated benzene.

The following tables are the breakdown of the VOA compounds according to their corresponding surrogates. They are grouped by type (structure) of organic compound. Compounds that are not similar (in structure) to a surrogate are listed on their own table. These compounds may not behave like the surrogates. Some compounds are qualified with "PP" to indicate they are poor performers. The poor performers include those compounds that do not purge efficiently, are reactive, show low response, or poor chromatographic behavior. Due to their unique chemical characteristics, compounds with the PP qualifier may not perform like the surrogates.

## **VOA Compound List Grouped by Surrogates**

**Retention Time Analyte Type** Bromofluorobenzene (SURR) 15.79 Surrogate Toluene-d8 (SURR) 13.29 Surrogate 1,2 Dichlorobenzene 17.43 1,2,3-Trichlorobenzene 20.19 1,2,4-Trichlorobenzene 19.54 Α 1,2,4-Trimethylbenzene 16.78 1,3 Dichlorobenzene 17 Α 16.4 1,3,5-Trimethylbenzene 1,4 Dichlorobenzene 17.06 Α 2-Chlorotoluene 16.26 Α 4-Chlorotoluene 16.33 Α 4-Isopropyltoluene 17.06 **Benze**ne 11.09 Bromobenzene 16 14.71 Chlorobenzene Α Ethylbenzene 14.87 Α Isopropylbenzene 15.74 Naphthalene 19.9 Α n-Butylbenzene 17.46 Α n-Propylbenzene 16.15 Α sec-Butylbenzene 16.9 Α tert-Butylbenzene 16.69 Α Toluene 13.36 Xylene O 15.42 Α Xylene P,M 15.04 Α

| Analyte                               | Retention Time | Туре      |
|---------------------------------------|----------------|-----------|
| 1,2-Dichloroethane-d4(SURR)           | 10.44          | Surrogate |
| Dibromofluoromethane(SURR)            | 9.93           | Surrogate |
| 1,1,1,2-Tetrachloroethane             | 14.64          | H         |
| 1,1,1-Trichloroethane                 | 10.64          | H         |
| 1,1,2,2-Tetrachloroethane             | 15.41          | H         |
| 1,1,2-Trichloro-1,2,2-trifluoroethane | 7.86           | H         |
| 1,1,2-Trichloroethane                 | 13.16          | H         |
| 1,1-Dichloroethane                    | 8.91           | Н         |
| 1,2,3-Trichloropropane                | 15.54          | H         |
| 1,2-Dibromo-3-Chloropropane           | 17.92          | H PP      |
| 1,2-Dibromoethane                     | 13.94          | H         |
| 1,2-Dichloroethane                    | 10.53          | H         |
| 1,2-Dichloropropane                   | 11.74          | H         |
| 1,3-Dichloropropane                   | 13.4           | H         |
| 1-Chlorobutane                        | 10.61          | H         |
| 2,2-Dichloropropane                   | 9.89           | H         |
| Bromochloromethane                    | 9.49           | H         |
| Bromodichloromethane                  | 11.83          | H         |
| Bromomethane                          | 5.91           | Н         |
| Chloroethane                          | 6.12           | Н         |
| Chloromethane                         | 4.99           | H         |
| Dibromochloromethane                  | 13.7           | Н         |
| Dibromomethane                        | 11.7           | Н         |
| Dichlorodifluoromethane               | 4.69           | H         |
| Hexachloroethane                      | 18.01          | Н         |
| lodomethane                           | 7.67           | Н         |
| Pentachloroethane                     | 16.52          | Н         |
| Trichlorofluoromethane                | 6.92           | H         |

A = Aromatic Hydrocarbon Compounds

H = Halogenated Hydrocarbon Compounds

PP = Poor Performers

## **VOA Compound List Grouped by Surrogates**

Analyte Retention Time Type

| Allalyte                                 | Retention Time | rype    |
|--|----------------|---------|
| No Representative Surrogate              | ED             | - 10 mm |
| 2-Butanone                               | 9.38           | K PP    |
| 2-Hexanone                               | 13.52          | KPP     |
| 4-Methyl-2-Pentanone                     | 12.61          | K PP    |
| Acetone                                  | 7.05           | K PP    |
| 1,1-Dichloroethene                       | 7.62           |         |
| 1,1-Dichloropropene                      | 10.84          |         |
| 1,4-Dioxane                              | 11.93          | PP      |
| 1-Chlorohexane                           | 14.61          |         |
| 2-Chloroethyl vinyl ether                | 12.24          | PP      |
| 2-Nitropropane                           | 12.61          | PP      |
| Acrolein                                 | 6.89           | PP      |
| Acrylonitrile                            | 7.66           | PP      |
| Allyl Chloride                           | 7.91           |         |
| Bromoform                                | 15.19          |         |
| Carbon Disulfide                         | 8.11           | PP      |
| Carbon Tetrachloride                     | 11.06          |         |
| Chloroform                               | 9.8            |         |
| Chloroprene                              | 9.33           |         |
| cis-1,2 Dichloroethene                   | 9.56           |         |
| cis-1,3-Dichloropropene                  | 12.52          |         |
| cis-1,4-Dichloro-2-butene                | 15.25          |         |
| Cyclohexane                              | 10.97          |         |
| Diethyl ether                            | 7.22           |         |
| Di-isopropyl ether                       | 9.39           |         |
| Ethyl Methacrylate                       | 13.37          | PP      |
| Ethyl tertiary-butyl ether               | 9.84           |         |
| Hexachlorobutadiene                      | 19.93          |         |
| Methacrylonitrile                        | 9.75           | PP      |
| Methyl Acetate                           | 7.84           |         |
| Methyl Acrylate                          | 9.82           |         |
| Methyl Cyclohexane                       | 12.36          |         |
| Methyl Methacrylate                      | 11.94          |         |
| Methyl tert-Butyl Ether                  | 8.73           |         |
| Methylene Chloride                       | 7.79           |         |
| Styrene                                  | 15.35          |         |
| Tertiary-amyl methyl ether               | 11.23          |         |
|  | 7.62           | - 1 V   |
| Tertiary-butyl Alcohol Tetrachloroethene | 14.09          |         |
|  |                | PP      |
| Tetrahydrofuran                          | 10.17          | Pr -    |
| trans-1,2-Dichloroethene                 | 8.61           |         |
| trans-1,3-Dichloropropene                | 12.98          | DD      |
| Trans-1,4-Dichloro-2-Butene              | 15.54          | PP      |
| Trichloroethene                          | 11.78          |         |
| Vinyl Acetate                            | 9.02           | +       |
| Vinyl Chloride                           | 5.27           |         |

K = Ketone Compounds

PP = Poor Performers