

Data Path : Z:\voasrv\HPCHEM1\MSVOA\_Y\Data\VY013123\  
 Data File : VY012431.D  
 Acq On : 31 Jan 2023 19:28  
 Operator : KP/MD  
 Sample : VSTDCCC050  
 Misc : 5.00g/5.0mL/MSVOA\_Y/SOIL  
 ALS Vial : 23 Sample Multiplier: 1

Instrument :  
 MSVOA\_Y  
 LabSampleId :  
 VSTDCCC050

Quant Time: Feb 01 01:26:59 2023  
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_Y\methods\82Y011623S.M  
 Quant Title : SW846 8260  
 QLast Update : Tue Jan 17 04:31:47 2023  
 Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.50min  
 Max. RRF Dev : 25% Max. Rel. Area : 150%

|       | Compound                    | AvgRF | CCRF  | %Dev   | Area% | Dev(min) |
|-------|-----------------------------|-------|-------|--------|-------|----------|
| 1 I   | Pentafluorobenzene          | 1.000 | 1.000 | 0.0    | 100   | 0.00     |
| 2 T   | Dichlorodifluoromethane     | 0.437 | 0.341 | 22.0   | 92    | 0.00     |
| 3 P   | Chloromethane               | 0.403 | 0.367 | 8.9    | 100   | 0.00     |
| 4 C   | Vinyl Chloride              | 0.471 | 0.440 | 6.6#   | 101   | 0.00     |
| 5 T   | Bromomethane                | 0.364 | 0.343 | 5.8    | 103   | 0.00     |
| 6 T   | Chloroethane                | 0.307 | 0.303 | 1.3    | 106   | 0.00     |
| 7 T   | Trichlorofluoromethane      | 0.919 | 0.862 | 6.2    | 102   | 0.00     |
| 8 T   | Diethyl Ether               | 0.273 | 0.287 | -5.1   | 111   | 0.00     |
| 9 T   | 1,1,2-Trichlorotrifluoroeth | 0.447 | 0.452 | -1.1   | 111   | 0.00     |
| 10 T  | Methyl Iodide               | 0.680 | 0.626 | 7.9    | 92    | 0.00     |
| 11 T  | Tert butyl alcohol          | 0.057 | 0.052 | 8.8    | 112   | 0.00     |
| 12 CM | 1,1-Dichloroethene          | 0.459 | 0.457 | 0.4#   | 104   | 0.00     |
| 13 T  | Acrolein                    | 0.023 | 0.018 | 21.7   | 114   | 0.00     |
| 14 T  | Allyl chloride              | 0.568 | 0.618 | -8.8   | 114   | 0.00     |
| 15 T  | Acrylonitrile               | 0.119 | 0.136 | -14.3  | 116   | 0.00     |
| 16 T  | Acetone                     | 0.139 | 0.133 | 4.3    | 86    | 0.00     |
| 17 T  | Carbon Disulfide            | 1.420 | 1.281 | 9.8    | 95    | 0.00     |
| 18 T  | Methyl Acetate              | 0.312 | 0.358 | -14.7  | 119   | 0.00     |
| 19 T  | Methyl tert-butyl Ether     | 1.280 | 1.368 | -6.9   | 107   | 0.00     |
| 20 T  | Methylene Chloride          | 0.762 | 0.583 | 23.5   | 102   | 0.00     |
| 21 T  | trans-1,2-Dichloroethene    | 0.512 | 0.520 | -1.6   | 105   | 0.00     |
| 22 T  | Diisopropyl ether           | 1.132 | 1.282 | -13.3  | 117   | 0.00     |
| 23 T  | Vinyl Acetate               | 0.705 | 0.783 | -11.1  | 111   | 0.00     |
| 24 P  | 1,1-Dichloroethane          | 0.785 | 0.848 | -8.0   | 112   | 0.00     |
| 25 T  | 2-Butanone                  | 0.154 | 0.163 | -5.8   | 104   | 0.00     |
| 26 T  | 2,2-Dichloropropane         | 0.852 | 0.850 | 0.2    | 105   | 0.00     |
| 27 T  | cis-1,2-Dichloroethene      | 0.563 | 0.601 | -6.7   | 110   | 0.00     |
| 28 T  | Bromochloromethane          | 0.171 | 0.168 | 1.8    | 105   | 0.00     |
| 29 T  | Tetrahydrofuran             | 0.084 | 0.097 | -15.5  | 114   | 0.00     |
| 30 C  | Chloroform                  | 0.927 | 0.976 | -5.3#  | 109   | 0.00     |
| 31 T  | Cyclohexane                 | 0.729 | 0.700 | 4.0    | 111   | 0.00     |
| 32 T  | 1,1,1-Trichloroethane       | 0.918 | 0.918 | 0.0    | 104   | 0.00     |
| 33 S  | 1,2-Dichloroethane-d4       | 0.465 | 0.006 | 98.7#  | 2#    | 0.00     |
| 34 I  | 1,4-Difluorobenzene         | 1.000 | 1.000 | 0.0    | 101   | 0.00     |
| 35 S  | Dibromofluoromethane        | 0.265 | 0.003 | 98.9#  | 1#    | 0.00     |
| 36 T  | 1,1-Dichloropropene         | 0.450 | 0.449 | 0.2    | 106   | 0.00     |
| 37 T  | Ethyl Acetate               | 0.198 | 0.214 | -8.1   | 110   | 0.00     |
| 38 T  | Carbon Tetrachloride        | 0.581 | 0.564 | 2.9    | 103   | 0.00     |
| 39 T  | Methylcyclohexane           | 0.550 | 0.561 | -2.0   | 110   | 0.00     |
| 40 TM | Benzene                     | 1.299 | 1.343 | -3.4   | 109   | 0.00     |
| 41 T  | Methacrylonitrile           | 0.117 | 0.139 | -18.8  | 116   | 0.00     |
| 42 TM | 1,2-Dichloroethane          | 0.413 | 0.419 | -1.5   | 103   | 0.00     |
| 43 T  | Isopropyl Acetate           | 0.369 | 0.414 | -12.2  | 113   | 0.00     |
| 44 TM | Trichloroethene             | 0.388 | 0.398 | -2.6   | 107   | 0.00     |
| 45 C  | 1,2-Dichloropropane         | 0.271 | 0.301 | -11.1# | 116   | 0.00     |
| 46 T  | Dibromomethane              | 0.181 | 0.195 | -7.7   | 110   | 0.00     |
| 47 T  | Bromodichloromethane        | 0.457 | 0.488 | -6.8   | 110   | 0.00     |
| 48 T  | Methyl methacrylate         | 0.170 | 0.189 | -11.2  | 108   | 0.00     |

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 ALS Vial : 23 Sample Multiplier: 1

Instrument :  
 MSVOA\_Y  
 LabSampleId :  
 VSTDCCC050

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 Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.50min  
 Max. RRF Dev : 25% Max. Rel. Area : 150%

|       | Compound                    | AvgRF | CCRF  | %Dev   | Area% | Dev(min) |
|-------|-----------------------------|-------|-------|--------|-------|----------|
| 49 T  | 1,4-Dioxane                 | 0.002 | 0.003 | -50.0# | 107   | 0.00     |
| 50 S  | Toluene-d8                  | 1.044 | 0.019 | 98.2#  | 2#    | 0.00     |
| 51 T  | 4-Methyl-2-Pentanone        | 0.197 | 0.222 | -12.7  | 113   | 0.00     |
| 52 CM | Toluene                     | 0.865 | 0.890 | -2.9#  | 107   | 0.00     |
| 53 T  | t-1,3-Dichloropropene       | 0.478 | 0.506 | -5.9   | 107   | 0.00     |
| 54 T  | cis-1,3-Dichloropropene     | 0.514 | 0.545 | -6.0   | 108   | 0.00     |
| 55 T  | 1,1,2-Trichloroethane       | 0.258 | 0.284 | -10.1  | 112   | 0.00     |
| 56 T  | Ethyl methacrylate          | 0.338 | 0.381 | -12.7  | 111   | 0.00     |
| 57 T  | 1,3-Dichloropropane         | 0.421 | 0.463 | -10.0  | 112   | 0.00     |
| 58 T  | 2-Chloroethyl Vinyl ether   | 0.112 | 0.113 | -0.9   | 105   | 0.00     |
| 59 T  | 2-Hexanone                  | 0.150 | 0.161 | -7.3   | 103   | 0.00     |
| 60 T  | Dibromochloromethane        | 0.342 | 0.371 | -8.5   | 109   | 0.00     |
| 61 T  | 1,2-Dibromoethane           | 0.249 | 0.270 | -8.4   | 110   | 0.00     |
| 62 S  | 4-Bromofluorobenzene        | 0.373 | 0.014 | 96.2#  | 4#    | 0.00     |
| 63 I  | Chlorobenzene-d5            | 1.000 | 1.000 | 0.0    | 103   | 0.00     |
| 64 T  | Tetrachloroethene           | 0.456 | 0.496 | -8.8   | 113   | 0.00     |
| 65 PM | Chlorobenzene               | 1.023 | 1.052 | -2.8   | 109   | 0.00     |
| 66 T  | 1,1,1,2-Tetrachloroethane   | 0.386 | 0.406 | -5.2   | 109   | 0.00     |
| 67 C  | Ethyl Benzene               | 1.828 | 1.878 | -2.7#  | 108   | 0.00     |
| 68 T  | m/p-Xylenes                 | 0.723 | 0.738 | -2.1   | 107   | 0.00     |
| 69 T  | o-Xylene                    | 0.690 | 0.710 | -2.9   | 107   | 0.00     |
| 70 T  | Styrene                     | 1.154 | 1.212 | -5.0   | 108   | 0.00     |
| 71 P  | Bromoform                   | 0.247 | 0.265 | -7.3   | 109   | 0.00     |
| 72 I  | 1,4-Dichlorobenzene-d4      | 1.000 | 1.000 | 0.0    | 105   | 0.00     |
| 73 T  | Isopropylbenzene            | 3.605 | 3.625 | -0.6   | 109   | 0.00     |
| 74 T  | N-amyl acetate              | 0.707 | 0.743 | -5.1   | 109   | 0.00     |
| 75 P  | 1,1,2,2-Tetrachloroethane   | 0.568 | 0.598 | -5.3   | 112   | 0.00     |
| 76 T  | 1,2,3-Trichloropropane      | 0.460 | 0.506 | -10.0  | 105   | 0.00     |
| 77 T  | Bromobenzene                | 0.868 | 0.870 | -0.2   | 107   | 0.00     |
| 78 T  | n-propylbenzene             | 4.219 | 4.218 | 0.0    | 110   | 0.00     |
| 79 T  | 2-Chlorotoluene             | 2.385 | 2.394 | -0.4   | 109   | 0.00     |
| 80 T  | 1,3,5-Trimethylbenzene      | 3.134 | 3.123 | 0.4    | 109   | 0.00     |
| 81 T  | trans-1,4-Dichloro-2-butene | 0.223 | 0.232 | -4.0   | 108   | 0.00     |
| 82 T  | 4-Chlorotoluene             | 2.506 | 2.491 | 0.6    | 109   | 0.00     |
| 83 T  | tert-Butylbenzene           | 2.732 | 2.717 | 0.5    | 111   | 0.00     |
| 84 T  | 1,2,4-Trimethylbenzene      | 3.091 | 3.069 | 0.7    | 107   | 0.00     |
| 85 T  | sec-Butylbenzene            | 3.889 | 3.897 | -0.2   | 110   | 0.00     |
| 86 T  | p-Isopropyltoluene          | 3.398 | 3.379 | 0.6    | 109   | 0.00     |
| 87 T  | 1,3-Dichlorobenzene         | 1.725 | 1.702 | 1.3    | 107   | 0.00     |
| 88 T  | 1,4-Dichlorobenzene         | 1.721 | 1.674 | 2.7    | 107   | 0.00     |
| 89 T  | n-Butylbenzene              | 2.945 | 2.945 | 0.0    | 110   | 0.00     |
| 90 T  | Hexachloroethane            | 0.613 | 0.608 | 0.8    | 110   | 0.00     |
| 91 T  | 1,2-Dichlorobenzene         | 1.540 | 1.541 | -0.1   | 109   | 0.00     |
| 92 T  | 1,2-Dibromo-3-Chloropropane | 0.117 | 0.121 | -3.4   | 109   | 0.00     |
| 93 T  | 1,2,4-Trichlorobenzene      | 0.987 | 0.963 | 2.4    | 108   | 0.00     |
| 94 T  | Hexachlorobutadiene         | 0.594 | 0.575 | 3.2    | 111   | 0.00     |
| 95 T  | Naphthalene                 | 1.947 | 2.009 | -3.2   | 109   | 0.00     |

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Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.50min  
Max. RRF Dev : 25% Max. Rel. Area : 150%

|      | Compound               | AvgRF | CCRF  | %Dev | Area% | Dev(min) |
|------|------------------------|-------|-------|------|-------|----------|
| 96 T | 1,2,3-Trichlorobenzene | 0.849 | 0.826 | 2.7  | 109   | 0.00     |

(#) = Out of Range                      SPCC's out = 0    CCC's out = 6