

Data Path : Z:\voasrv\HPCHEM1\MSVOA\_Y\Data\VY062023\  
 Data File : VY014301.D  
 Acq On : 20 Jun 2023 22:31  
 Operator : KP/MD  
 Sample : VSTDCCC050  
 Misc : 5.00g/5.0mL/MSVOA\_Y/SOIL  
 ALS Vial : 31 Sample Multiplier: 1

Instrument :  
 MSVOA\_Y  
 LabSampleId :  
 VSTDCCC050

Quant Time: Jun 21 02:10:53 2023  
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_Y\methods\82Y060923S.M  
 Quant Title : SW846 8260  
 QLast Update : Fri Jun 09 18:07:13 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    | 89    | 0.00     |
| 2 T   | Dichlorodifluoromethane     | 0.557 | 0.509 | 8.6    | 84    | 0.00     |
| 3 P   | Chloromethane               | 0.707 | 0.581 | 17.8   | 78    | 0.00     |
| 4 C   | Vinyl Chloride              | 0.687 | 0.613 | 10.8#  | 82    | 0.00     |
| 5 T   | Bromomethane                | 0.567 | 0.461 | 18.7   | 83    | -0.02    |
| 6 T   | Chloroethane                | 0.467 | 0.407 | 12.8   | 84    | 0.00     |
| 7 T   | Trichlorofluoromethane      | 1.116 | 1.145 | -2.6   | 94    | 0.00     |
| 8 T   | Diethyl Ether               | 0.373 | 0.379 | -1.6   | 96    | 0.00     |
| 9 T   | 1,1,2-Trichlorotrifluoroeth | 0.563 | 0.599 | -6.4   | 94    | 0.00     |
| 10 T  | Methyl Iodide               | 0.744 | 0.697 | 6.3    | 82    | 0.00     |
| 11 T  | Tert butyl alcohol          | 0.092 | 0.109 | -18.5  | 132   | 0.00     |
| 12 CM | 1,1-Dichloroethene          | 0.571 | 0.563 | 1.4#   | 91    | 0.00     |
| 13 T  | Acrolein                    | 0.044 | 0.030 | 31.8#  | 83    | 0.00     |
| 14 T  | Allyl chloride              | 1.035 | 1.001 | 3.3    | 88    | 0.00     |
| 15 T  | Acrylonitrile               | 0.192 | 0.238 | -24.0  | 116   | 0.00     |
| 16 T  | Acetone                     | 0.210 | 0.206 | 1.9    | 103   | 0.00     |
| 17 T  | Carbon Disulfide            | 1.850 | 1.715 | 7.3    | 86    | 0.00     |
| 18 T  | Methyl Acetate              | 0.815 | 1.164 | -42.8# | 141   | 0.00     |
| 19 T  | Methyl tert-butyl Ether     | 1.684 | 1.830 | -8.7   | 96    | 0.00     |
| 20 T  | Methylene Chloride          | 0.775 | 0.874 | -12.8  | 116   | 0.00     |
| 21 T  | trans-1,2-Dichloroethene    | 0.637 | 0.651 | -2.2   | 91    | 0.00     |
| 22 T  | Diisopropyl ether           | 1.981 | 2.127 | -7.4   | 96    | 0.00     |
| 23 T  | Vinyl Acetate               | 1.169 | 1.199 | -2.6   | 92    | 0.00     |
| 24 P  | 1,1-Dichloroethane          | 1.142 | 1.190 | -4.2   | 96    | 0.00     |
| 25 T  | 2-Butanone                  | 0.258 | 0.332 | -28.7# | 128   | 0.00     |
| 26 T  | 2,2-Dichloropropane         | 1.073 | 1.017 | 5.2    | 88    | 0.00     |
| 27 T  | cis-1,2-Dichloroethene      | 0.712 | 0.753 | -5.8   | 98    | 0.00     |
| 28 T  | Bromochloromethane          | 0.465 | 0.456 | 1.9    | 95    | 0.00     |
| 29 T  | Tetrahydrofuran             | 0.164 | 0.221 | -34.8# | 131   | 0.00     |
| 30 C  | Chloroform                  | 1.212 | 1.302 | -7.4#  | 99    | 0.00     |
| 31 T  | Cyclohexane                 | 0.933 | 0.976 | -4.6   | 97    | 0.00     |
| 32 T  | 1,1,1-Trichloroethane       | 1.093 | 1.174 | -7.4   | 99    | 0.00     |
| 33 S  | 1,2-Dichloroethane-d4       | 0.696 | 0.685 | 1.6    | 100   | 0.00     |
| 34 I  | 1,4-Difluorobenzene         | 1.000 | 1.000 | 0.0    | 87    | 0.00     |
| 35 S  | Dibromofluoromethane        | 0.355 | 0.387 | -9.0   | 99    | 0.00     |
| 36 T  | 1,1-Dichloropropene         | 0.553 | 0.619 | -11.9  | 98    | 0.00     |
| 37 T  | Ethyl Acetate               | 0.381 | 0.452 | -18.6  | 107   | 0.00     |
| 38 T  | Carbon Tetrachloride        | 0.638 | 0.730 | -14.4  | 99    | 0.00     |
| 39 T  | Methylcyclohexane           | 0.613 | 0.704 | -14.8  | 97    | 0.00     |
| 40 TM | Benzene                     | 1.644 | 1.811 | -10.2  | 98    | 0.00     |
| 41 T  | Methacrylonitrile           | 0.195 | 0.244 | -25.1# | 127   | 0.00     |
| 42 TM | 1,2-Dichloroethane          | 0.581 | 0.647 | -11.4  | 101   | 0.00     |
| 43 T  | Isopropyl Acetate           | 0.669 | 0.788 | -17.8  | 106   | 0.00     |
| 44 TM | Trichloroethene             | 0.439 | 0.491 | -11.8  | 98    | 0.00     |
| 45 C  | 1,2-Dichloropropane         | 0.418 | 0.461 | -10.3# | 98    | 0.00     |
| 46 T  | Dibromomethane              | 0.257 | 0.291 | -13.2  | 103   | 0.00     |
| 47 T  | Bromodichloromethane        | 0.602 | 0.672 | -11.6  | 99    | 0.00     |
| 48 T  | Methyl methacrylate         | 0.312 | 0.398 | -27.6# | 115   | 0.00     |

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 ALS Vial : 31 Sample Multiplier: 1

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 MSVOA\_Y  
 LabSampleId :  
 VSTDCCC050

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 Quant Title : SW846 8260  
 QLast Update : Fri Jun 09 18:07:13 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) |
|-------|-----------------------------|-------|-------|--------|-------|----------|
| 49 T  | 1,4-Dioxane                 | 0.003 | 0.005 | -66.7# | 127   | -0.01    |
| 50 S  | Toluene-d8                  | 1.326 | 1.335 | -0.7   | 97    | 0.00     |
| 51 T  | 4-Methyl-2-Pentanone        | 0.363 | 0.494 | -36.1# | 123   | 0.00     |
| 52 CM | Toluene                     | 1.019 | 1.153 | -13.2# | 99    | 0.00     |
| 53 T  | t-1,3-Dichloropropene       | 0.634 | 0.705 | -11.2  | 98    | 0.00     |
| 54 T  | cis-1,3-Dichloropropene     | 0.691 | 0.759 | -9.8   | 98    | 0.00     |
| 55 T  | 1,1,2-Trichloroethane       | 0.341 | 0.406 | -19.1  | 106   | 0.00     |
| 56 T  | Ethyl methacrylate          | 0.470 | 0.572 | -21.7  | 104   | 0.00     |
| 57 T  | 1,3-Dichloropropane         | 0.583 | 0.677 | -16.1  | 102   | 0.00     |
| 58 T  | 2-Chloroethyl Vinyl ether   | 0.177 | 0.283 | -59.9# | 122   | 0.00     |
| 59 T  | 2-Hexanone                  | 0.262 | 0.360 | -37.4# | 126   | 0.00     |
| 60 T  | Dibromochloromethane        | 0.436 | 0.514 | -17.9  | 105   | 0.00     |
| 61 T  | 1,2-Dibromoethane           | 0.336 | 0.395 | -17.6  | 106   | 0.00     |
| 62 S  | 4-Bromofluorobenzene        | 0.471 | 0.562 | -19.3  | 112   | 0.00     |
| 63 I  | Chlorobenzene-d5            | 1.000 | 1.000 | 0.0    | 94    | 0.00     |
| 64 T  | Tetrachloroethene           | 0.426 | 0.529 | -24.2  | 117   | 0.00     |
| 65 PM | Chlorobenzene               | 1.201 | 1.272 | -5.9   | 100   | 0.00     |
| 66 T  | 1,1,1,2-Tetrachloroethane   | 0.473 | 0.508 | -7.4   | 100   | 0.00     |
| 67 C  | Ethyl Benzene               | 2.103 | 2.315 | -10.1# | 101   | 0.00     |
| 68 T  | m/p-Xylenes                 | 0.809 | 0.891 | -10.1  | 101   | 0.00     |
| 69 T  | o-Xylene                    | 0.765 | 0.854 | -11.6  | 102   | 0.00     |
| 70 T  | Styrene                     | 1.305 | 1.475 | -13.0  | 103   | 0.00     |
| 71 P  | Bromoform                   | 0.324 | 0.382 | -17.9  | 112   | 0.00     |
| 72 I  | 1,4-Dichlorobenzene-d4      | 1.000 | 1.000 | 0.0    | 103   | 0.00     |
| 73 T  | Isopropylbenzene            | 3.814 | 4.022 | -5.5   | 102   | 0.00     |
| 74 T  | N-amyl acetate              | 1.209 | 1.271 | -5.1   | 105   | 0.00     |
| 75 P  | 1,1,2,2-Tetrachloroethane   | 0.852 | 0.944 | -10.8  | 114   | 0.00     |
| 76 T  | 1,2,3-Trichloropropane      | 0.656 | 0.811 | -23.6  | 119   | 0.00     |
| 77 T  | Bromobenzene                | 1.003 | 1.026 | -2.3   | 103   | 0.00     |
| 78 T  | n-propylbenzene             | 4.659 | 5.006 | -7.4   | 105   | 0.00     |
| 79 T  | 2-Chlorotoluene             | 2.681 | 2.814 | -5.0   | 105   | 0.00     |
| 80 T  | 1,3,5-Trimethylbenzene      | 3.266 | 3.501 | -7.2   | 105   | 0.00     |
| 81 T  | trans-1,4-Dichloro-2-butene | 0.303 | 0.328 | -8.3   | 108   | 0.00     |
| 82 T  | 4-Chlorotoluene             | 2.841 | 2.960 | -4.2   | 104   | 0.00     |
| 83 T  | tert-Butylbenzene           | 2.775 | 3.027 | -9.1   | 107   | 0.00     |
| 84 T  | 1,2,4-Trimethylbenzene      | 3.233 | 3.522 | -8.9   | 107   | 0.00     |
| 85 T  | sec-Butylbenzene            | 4.035 | 4.422 | -9.6   | 106   | 0.00     |
| 86 T  | p-Isopropyltoluene          | 3.453 | 3.716 | -7.6   | 105   | 0.00     |
| 87 T  | 1,3-Dichlorobenzene         | 1.914 | 2.008 | -4.9   | 106   | 0.00     |
| 88 T  | 1,4-Dichlorobenzene         | 1.929 | 2.005 | -3.9   | 106   | 0.00     |
| 89 T  | n-Butylbenzene              | 3.203 | 3.505 | -9.4   | 107   | 0.00     |
| 90 T  | Hexachloroethane            | 0.688 | 0.741 | -7.7   | 109   | 0.00     |
| 91 T  | 1,2-Dichlorobenzene         | 1.719 | 1.840 | -7.0   | 109   | 0.00     |
| 92 T  | 1,2-Dibromo-3-Chloropropane | 0.162 | 0.192 | -18.5  | 132   | 0.00     |
| 93 T  | 1,2,4-Trichlorobenzene      | 1.089 | 1.135 | -4.2   | 106   | 0.00     |
| 94 T  | Hexachlorobutadiene         | 0.629 | 0.687 | -9.2   | 109   | 0.00     |
| 95 T  | Naphthalene                 | 2.210 | 2.553 | -15.5  | 119   | 0.00     |

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Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_Y\methods\82Y060923S.M  
Quant Title : SW846 8260  
QLast Update : Fri Jun 09 18:07:13 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) |
|------|------------------------|-------|-------|------|-------|----------|
| 96 T | 1,2,3-Trichlorobenzene | 0.984 | 1.052 | -6.9 | 112   | 0.00     |

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