

Data Path : Z:\voasrv\HPCHEM1\MSVOA\_D\Data\VD020823\  
 Data File : VD075263.D  
 Acq On : 08 Feb 2023 10:16  
 Operator : KP/SY  
 Sample : VSTDCCC050  
 Misc : 5.00G/5.00ml/MSVOA\_D/SOIL  
 ALS Vial : 2 Sample Multiplier: 1

Instrument :  
 MSVOA\_D  
 LabSampleId :  
 VSTDCCC050

Quant Time: Feb 09 05:39:51 2023  
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_D\Method\82D020723S.M  
 Quant Title : SW846 8260  
 QLast Update : Wed Feb 08 04:26:58 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                    | Amount  | Calc.   | %Dev   | Area% | Dev(min) |
|-------|-----------------------------|---------|---------|--------|-------|----------|
| 1 I   | Pentafluorobenzene          | 50.000  | 50.000  | 0.0    | 97    | 0.00     |
| 2 T   | Dichlorodifluoromethane     | 50.000  | 50.219  | -0.4   | 96    | 0.00     |
| 3 P   | Chloromethane               | 50.000  | 47.085  | 5.8    | 91    | 0.00     |
| 4 C   | Vinyl Chloride              | 50.000  | 44.099  | 11.8#  | 89    | 0.00     |
| 5 T   | Bromomethane                | 50.000  | 46.142  | 7.7    | 104   | 0.00     |
| 6 T   | Chloroethane                | 50.000  | 45.463  | 9.1    | 95    | 0.00     |
| 7 T   | Trichlorofluoromethane      | 50.000  | 47.759  | 4.5    | 97    | 0.00     |
| 8 T   | Diethyl Ether               | 50.000  | 49.108  | 1.8    | 96    | 0.00     |
| 9 T   | 1,1,2-Trichlorotrifluoroeth | 50.000  | 49.814  | 0.4    | 98    | 0.00     |
| 10 T  | Methyl Iodide               | 50.000  | 48.916  | 2.2    | 95    | 0.00     |
| 11 T  | Tert butyl alcohol          | 250.000 | 234.949 | 6.0    | 104   | 0.00     |
| 12 CM | 1,1-Dichloroethene          | 50.000  | 49.094  | 1.8#   | 98    | 0.00     |
| 13 T  | Acrolein                    | 250.000 | 202.406 | 19.0   | 92    | 0.00     |
| 14 T  | Allyl chloride              | 50.000  | 50.678  | -1.4   | 103   | 0.00     |
| 15 T  | Acrylonitrile               | 250.000 | 249.854 | 0.1    | 97    | -0.01    |
| 16 T  | Acetone                     | 250.000 | 265.087 | -6.0   | 95    | 0.00     |
| 17 T  | Carbon Disulfide            | 50.000  | 47.275  | 5.5    | 97    | 0.00     |
| 18 T  | Methyl Acetate              | 50.000  | 48.108  | 3.8    | 96    | 0.00     |
| 19 T  | Methyl tert-butyl Ether     | 50.000  | 50.727  | -1.5   | 99    | 0.00     |
| 20 T  | Methylene Chloride          | 50.000  | 50.818  | -1.6   | 95    | -0.01    |
| 21 T  | trans-1,2-Dichloroethene    | 50.000  | 49.542  | 0.9    | 97    | -0.01    |
| 22 T  | Diisopropyl ether           | 50.000  | 51.973  | -3.9   | 99    | 0.00     |
| 23 T  | Vinyl Acetate               | 250.000 | 369.787 | -47.9# | 156   | 0.00     |
| 24 P  | 1,1-Dichloroethane          | 50.000  | 51.069  | -2.1   | 101   | -0.01    |
| 25 T  | 2-Butanone                  | 250.000 | 248.869 | 0.5    | 95    | 0.00     |
| 26 T  | 2,2-Dichloropropane         | 50.000  | 51.320  | -2.6   | 100   | 0.00     |
| 27 T  | cis-1,2-Dichloroethene      | 50.000  | 51.066  | -2.1   | 101   | 0.00     |
| 28 T  | Bromochloromethane          | 50.000  | 51.495  | -3.0   | 101   | 0.00     |
| 29 T  | Tetrahydrofuran             | 250.000 | 254.528 | -1.8   | 97    | 0.00     |
| 30 C  | Chloroform                  | 50.000  | 51.756  | -3.5#  | 102   | 0.00     |
| 31 T  | Cyclohexane                 | 50.000  | 48.613  | 2.8    | 100   | 0.00     |
| 32 T  | 1,1,1-Trichloroethane       | 50.000  | 51.545  | -3.1   | 100   | 0.00     |
| 33 S  | 1,2-Dichloroethane-d4       | 50.000  | 49.129  | 1.7    | 103   | 0.00     |
| 34 I  | 1,4-Difluorobenzene         | 50.000  | 50.000  | 0.0    | 99    | 0.00     |
| 35 S  | Dibromofluoromethane        | 50.000  | 48.121  | 3.8    | 98    | 0.00     |
| 36 T  | 1,1-Dichloropropene         | 50.000  | 50.519  | -1.0   | 99    | 0.00     |
| 37 T  | Ethyl Acetate               | 50.000  | 49.128  | 1.7    | 105   | 0.00     |
| 38 T  | Carbon Tetrachloride        | 50.000  | 53.114  | -6.2   | 103   | 0.00     |
| 39 T  | Methylcyclohexane           | 50.000  | 50.272  | -0.5   | 95    | 0.00     |
| 40 TM | Benzene                     | 50.000  | 50.140  | -0.3   | 98    | 0.00     |
| 41 T  | Methacrylonitrile           | 50.000  | 41.120  | 17.8   | 89    | -0.01    |
| 42 TM | 1,2-Dichloroethane          | 50.000  | 50.018  | -0.0   | 99    | 0.00     |
| 43 T  | Isopropyl Acetate           | 50.000  | 48.484  | 3.0    | 97    | 0.00     |
| 44 TM | Trichloroethene             | 50.000  | 49.169  | 1.7    | 97    | 0.00     |
| 45 C  | 1,2-Dichloropropane         | 50.000  | 50.242  | -0.5#  | 100   | 0.00     |
| 46 T  | Dibromomethane              | 50.000  | 49.215  | 1.6    | 99    | 0.00     |
| 47 T  | Bromodichloromethane        | 50.000  | 50.397  | -0.8   | 98    | 0.00     |
| 48 T  | Methyl methacrylate         | 50.000  | 50.705  | -1.4   | 96    | 0.00     |

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 ALS Vial : 2 Sample Multiplier: 1

Instrument :  
 MSVOA\_D  
 LabSampleId :  
 VSTDCCC050

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 Quant Title : SW846 8260  
 QLast Update : Wed Feb 08 04:26:58 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                    | Amount   | Calc.   | %Dev  | Area% | Dev(min) |
|-------|-----------------------------|----------|---------|-------|-------|----------|
| 49 T  | 1,4-Dioxane                 | 1000.000 | 957.044 | 4.3   | 91    | 0.00     |
| 50 S  | Toluene-d8                  | 50.000   | 48.507  | 3.0   | 99    | 0.00     |
| 51 T  | 4-Methyl-2-Pentanone        | 250.000  | 250.805 | -0.3  | 96    | 0.00     |
| 52 CM | Toluene                     | 50.000   | 51.124  | -2.2# | 98    | 0.00     |
| 53 T  | t-1,3-Dichloropropene       | 50.000   | 54.089  | -8.2  | 103   | 0.00     |
| 54 T  | cis-1,3-Dichloropropene     | 50.000   | 54.110  | -8.2  | 103   | 0.00     |
| 55 T  | 1,1,2-Trichloroethane       | 50.000   | 51.422  | -2.8  | 99    | 0.00     |
| 56 T  | Ethyl methacrylate          | 50.000   | 52.978  | -6.0  | 98    | 0.00     |
| 57 T  | 1,3-Dichloropropane         | 50.000   | 51.223  | -2.4  | 99    | 0.00     |
| 58 T  | 2-Chloroethyl Vinyl ether   | 250.000  | 284.610 | -13.8 | 125   | 0.00     |
| 59 T  | 2-Hexanone                  | 250.000  | 257.602 | -3.0  | 96    | 0.00     |
| 60 T  | Dibromochloromethane        | 50.000   | 51.088  | -2.2  | 97    | 0.00     |
| 61 T  | 1,2-Dibromoethane           | 50.000   | 52.113  | -4.2  | 102   | 0.00     |
| 62 S  | 4-Bromofluorobenzene        | 50.000   | 49.478  | 1.0   | 100   | 0.00     |
| 63 I  | Chlorobenzene-d5            | 50.000   | 50.000  | 0.0   | 96    | 0.00     |
| 64 T  | Tetrachloroethene           | 50.000   | 50.571  | -1.1  | 98    | 0.00     |
| 65 PM | Chlorobenzene               | 50.000   | 50.494  | -1.0  | 97    | 0.00     |
| 66 T  | 1,1,1,2-Tetrachloroethane   | 50.000   | 51.667  | -3.3  | 96    | 0.00     |
| 67 C  | Ethyl Benzene               | 50.000   | 51.887  | -3.8# | 97    | 0.00     |
| 68 T  | m/p-Xylenes                 | 100.000  | 106.700 | -6.7  | 98    | 0.00     |
| 69 T  | o-Xylene                    | 50.000   | 53.797  | -7.6  | 100   | 0.00     |
| 70 T  | Styrene                     | 50.000   | 53.501  | -7.0  | 100   | 0.00     |
| 71 P  | Bromoform                   | 50.000   | 52.389  | -4.8  | 96    | 0.00     |
| 72 I  | 1,4-Dichlorobenzene-d4      | 50.000   | 50.000  | 0.0   | 99    | 0.00     |
| 73 T  | Isopropylbenzene            | 50.000   | 51.123  | -2.2  | 98    | 0.00     |
| 74 T  | N-ethyl acetate             | 50.000   | 51.037  | -2.1  | 99    | 0.00     |
| 75 P  | 1,1,2,2-Tetrachloroethane   | 50.000   | 49.332  | 1.3   | 98    | 0.00     |
| 76 T  | 1,2,3-Trichloropropane      | 50.000   | 47.308  | 5.4   | 102   | 0.00     |
| 77 T  | Bromobenzene                | 50.000   | 49.352  | 1.3   | 96    | 0.00     |
| 78 T  | n-propylbenzene             | 50.000   | 51.517  | -3.0  | 97    | 0.00     |
| 79 T  | 2-Chlorotoluene             | 50.000   | 50.752  | -1.5  | 98    | 0.00     |
| 80 T  | 1,3,5-Trimethylbenzene      | 50.000   | 51.216  | -2.4  | 97    | 0.00     |
| 81 T  | trans-1,4-Dichloro-2-butene | 50.000   | 59.630  | -19.3 | 124   | 0.00     |
| 82 T  | 4-Chlorotoluene             | 50.000   | 50.594  | -1.2  | 98    | 0.00     |
| 83 T  | tert-Butylbenzene           | 50.000   | 51.205  | -2.4  | 96    | 0.00     |
| 84 T  | 1,2,4-Trimethylbenzene      | 50.000   | 51.709  | -3.4  | 98    | 0.00     |
| 85 T  | sec-Butylbenzene            | 50.000   | 51.935  | -3.9  | 99    | 0.00     |
| 86 T  | p-Isopropyltoluene          | 50.000   | 52.664  | -5.3  | 98    | 0.00     |
| 87 T  | 1,3-Dichlorobenzene         | 50.000   | 51.110  | -2.2  | 99    | 0.00     |
| 88 T  | 1,4-Dichlorobenzene         | 50.000   | 50.659  | -1.3  | 100   | 0.00     |
| 89 T  | n-Butylbenzene              | 50.000   | 52.075  | -4.2  | 99    | 0.00     |
| 90 T  | Hexachloroethane            | 50.000   | 51.008  | -2.0  | 99    | 0.00     |
| 91 T  | 1,2-Dichlorobenzene         | 50.000   | 51.072  | -2.1  | 99    | 0.00     |
| 92 T  | 1,2-Dibromo-3-Chloropropane | 50.000   | 54.049  | -8.1  | 104   | 0.00     |
| 93 T  | 1,2,4-Trichlorobenzene      | 50.000   | 50.448  | -0.9  | 98    | 0.00     |
| 94 T  | Hexachlorobutadiene         | 50.000   | 50.419  | -0.8  | 94    | 0.00     |
| 95 T  | Naphthalene                 | 50.000   | 49.897  | 0.2   | 96    | 0.00     |

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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               | Amount | Calc.  | %Dev | Area% | Dev(min) |
|------|------------------------|--------|--------|------|-------|----------|
| 96 T | 1,2,3-Trichlorobenzene | 50.000 | 50.709 | -1.4 | 98    | 0.00     |

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