

Data Path : Z:\voasrv\HPCHEM1\MSVOA\_D\Data\VD071921\  
 Data File : VD069739.D  
 Acq On : 19 Jul 2021 08:57  
 Operator : VA/SY  
 Sample : VSTDCCC050  
 Misc : 5.00G/5.00ml/MSVOA\_D/SOIL  
 ALS Vial : 2 Sample Multiplier: 1

Instrument :  
 MSVOA\_D  
 LabSampleId :  
 VSTDCCC050

Quant Time: Jul 20 02:08:02 2021  
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_D\Method\82D071321S.M  
 Quant Title : SW846 8260  
 QLast Update : Wed Jul 14 06:00:31 2021  
 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   | 92    | 0.00     |
| 2 T   | Dichlorodifluoromethane     | 0.498 | 0.546 | -9.6  | 89    | 0.00     |
| 3 P   | Chloromethane               | 0.705 | 0.725 | -2.8  | 88    | 0.00     |
| 4 C   | Vinyl Chloride              | 0.761 | 0.795 | -4.5# | 89    | 0.00     |
| 5 T   | Bromomethane                | 0.500 | 0.489 | 2.2   | 93    | 0.00     |
| 6 T   | Chloroethane                | 0.487 | 0.506 | -3.9  | 91    | 0.00     |
| 7 T   | Trichlorofluoromethane      | 0.959 | 1.002 | -4.5  | 92    | 0.00     |
| 8 T   | Diethyl Ether               | 0.280 | 0.277 | 1.1   | 87    | 0.00     |
| 9 T   | 1,1,2-Trichlorotrifluoroeth | 0.595 | 0.637 | -7.1  | 95    | 0.00     |
| 10 T  | Methyl Iodide               | 0.554 | 0.577 | -4.2  | 82    | 0.00     |
| 11 T  | Tert butyl alcohol          | 0.083 | 0.026 | 68.7# | 53    | 0.01     |
| 12 CM | 1,1-Dichloroethene          | 0.545 | 0.579 | -6.2# | 89    | 0.00     |
| 13 T  | Acrolein                    | 0.043 | 0.050 | -16.3 | 115   | 0.00     |
| 14 T  | Allyl chloride              | 0.843 | 0.861 | -2.1  | 87    | 0.00     |
| 15 T  | Acrylonitrile               | 0.127 | 0.130 | -2.4  | 90    | 0.00     |
| 16 T  | Acetone                     | 0.114 | 0.135 | -18.4 | 113   | 0.00     |
| 17 T  | Carbon Disulfide            | 1.857 | 2.041 | -9.9  | 89    | 0.00     |
| 18 T  | Methyl Acetate              | 0.352 | 0.290 | 17.6  | 93    | 0.00     |
| 19 T  | Methyl tert-butyl Ether     | 1.154 | 1.243 | -7.7  | 89    | 0.00     |
| 20 T  | Methylene Chloride          | 1.096 | 0.721 | 34.2# | 80    | 0.00     |
| 21 T  | trans-1,2-Dichloroethene    | 0.651 | 0.676 | -3.8  | 91    | 0.00     |
| 22 T  | Diisopropyl ether           | 1.746 | 1.930 | -10.5 | 89    | 0.00     |
| 23 T  | Vinyl Acetate               | 0.828 | 0.960 | -15.9 | 88    | 0.00     |
| 24 P  | 1,1-Dichloroethane          | 1.219 | 1.239 | -1.6  | 90    | 0.00     |
| 25 T  | 2-Butanone                  | 0.148 | 0.152 | -2.7  | 92    | 0.00     |
| 26 T  | 2,2-Dichloropropane         | 0.972 | 1.029 | -5.9  | 96    | 0.00     |
| 27 T  | cis-1,2-Dichloroethene      | 0.702 | 0.708 | -0.9  | 88    | 0.00     |
| 28 T  | Bromochloromethane          | 0.470 | 0.510 | -8.5  | 96    | 0.00     |
| 29 T  | Tetrahydrofuran             | 0.091 | 0.098 | -7.7  | 90    | 0.00     |
| 30 C  | Chloroform                  | 1.212 | 1.185 | 2.2#  | 87    | 0.00     |
| 31 T  | Cyclohexane                 | 1.037 | 1.111 | -7.1  | 90    | 0.00     |
| 32 T  | 1,1,1-Trichloroethane       | 1.028 | 1.054 | -2.5  | 92    | 0.00     |
| 33 S  | 1,2-Dichloroethane-d4       | 0.624 | 0.588 | 5.8   | 87    | 0.00     |
| 34 I  | 1,4-Difluorobenzene         | 1.000 | 1.000 | 0.0   | 91    | 0.00     |
| 35 S  | Dibromofluoromethane        | 0.335 | 0.338 | -0.9  | 89    | 0.00     |
| 36 T  | 1,1-Dichloropropene         | 0.516 | 0.551 | -6.8  | 88    | 0.00     |
| 37 T  | Ethyl Acetate               | 0.195 | 0.207 | -6.2  | 90    | 0.00     |
| 38 T  | Carbon Tetrachloride        | 0.472 | 0.508 | -7.6  | 90    | 0.00     |
| 39 T  | Methylcyclohexane           | 0.559 | 0.635 | -13.6 | 86    | 0.00     |
| 40 TM | Benzene                     | 1.488 | 1.611 | -8.3  | 90    | 0.00     |
| 41 T  | Methacrylonitrile           | 0.101 | 0.095 | 5.9   | 81    | 0.00     |
| 42 TM | 1,2-Dichloroethane          | 0.404 | 0.422 | -4.5  | 90    | 0.00     |
| 43 T  | Isopropyl Acetate           | 0.357 | 0.377 | -5.6  | 90    | 0.00     |
| 44 TM | Trichloroethene             | 0.361 | 0.368 | -1.9  | 87    | 0.00     |
| 45 C  | 1,2-Dichloropropane         | 0.391 | 0.399 | -2.0# | 87    | 0.00     |
| 46 T  | Dibromomethane              | 0.191 | 0.193 | -1.0  | 86    | 0.00     |
| 47 T  | Bromodichloromethane        | 0.506 | 0.535 | -5.7  | 92    | 0.00     |
| 48 T  | Methyl methacrylate         | 0.169 | 0.177 | -4.7  | 83    | 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  
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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 | 0.002 | 0.0    | 83    | 0.00     |
| 50 S  | Toluene-d8                  | 1.303 | 1.343 | -3.1   | 88    | 0.00     |
| 51 T  | 4-Methyl-2-Pentanone        | 0.185 | 0.207 | -11.9  | 93    | 0.00     |
| 52 CM | Toluene                     | 0.895 | 1.017 | -13.6# | 91    | 0.00     |
| 53 T  | t-1,3-Dichloropropene       | 0.425 | 0.481 | -13.2  | 96    | 0.00     |
| 54 T  | cis-1,3-Dichloropropene     | 0.522 | 0.562 | -7.7   | 95    | 0.00     |
| 55 T  | 1,1,2-Trichloroethane       | 0.267 | 0.271 | -1.5   | 88    | 0.00     |
| 56 T  | Ethyl methacrylate          | 0.305 | 0.352 | -15.4  | 91    | 0.00     |
| 57 T  | 1,3-Dichloropropane         | 0.471 | 0.507 | -7.6   | 91    | 0.00     |
| 58 T  | 2-Chloroethyl Vinyl ether   | 0.136 | 0.176 | -29.4# | 109   | 0.00     |
| 59 T  | 2-Hexanone                  | 0.123 | 0.151 | -22.8  | 100   | 0.00     |
| 60 T  | Dibromochloromethane        | 0.311 | 0.336 | -8.0   | 94    | 0.00     |
| 61 T  | 1,2-Dibromoethane           | 0.248 | 0.266 | -7.3   | 93    | 0.00     |
| 62 S  | 4-Bromofluorobenzene        | 0.435 | 0.453 | -4.1   | 89    | 0.00     |
| 63 I  | Chlorobenzene-d5            | 1.000 | 1.000 | 0.0    | 90    | 0.00     |
| 64 T  | Tetrachloroethene           | 0.305 | 0.308 | -1.0   | 87    | 0.00     |
| 65 PM | Chlorobenzene               | 0.999 | 1.072 | -7.3   | 91    | 0.00     |
| 66 T  | 1,1,1,2-Tetrachloroethane   | 0.357 | 0.386 | -8.1   | 93    | 0.00     |
| 67 C  | Ethyl Benzene               | 1.769 | 2.000 | -13.1# | 91    | 0.00     |
| 68 T  | m/p-Xylenes                 | 0.669 | 0.782 | -16.9  | 93    | 0.00     |
| 69 T  | o-Xylene                    | 0.609 | 0.706 | -15.9  | 92    | 0.00     |
| 70 T  | Styrene                     | 1.063 | 1.250 | -17.6  | 93    | 0.00     |
| 71 P  | Bromoform                   | 0.173 | 0.192 | -11.0  | 96    | 0.00     |
| 72 I  | 1,4-Dichlorobenzene-d4      | 1.000 | 1.000 | 0.0    | 95    | 0.00     |
| 73 T  | Isopropylbenzene            | 3.560 | 3.934 | -10.5  | 93    | 0.00     |
| 74 T  | N-amyl acetate              | 0.774 | 0.804 | -3.9   | 91    | 0.00     |
| 75 P  | 1,1,2,2-Tetrachloroethane   | 0.709 | 0.731 | -3.1   | 96    | 0.00     |
| 76 T  | 1,2,3-Trichloropropane      | 0.486 | 0.499 | -2.7   | 93    | 0.00     |
| 77 T  | Bromobenzene                | 0.790 | 0.825 | -4.4   | 92    | 0.00     |
| 78 T  | n-propylbenzene             | 4.514 | 5.114 | -13.3  | 94    | 0.00     |
| 79 T  | 2-Chlorotoluene             | 2.654 | 2.851 | -7.4   | 92    | 0.00     |
| 80 T  | 1,3,5-Trimethylbenzene      | 3.020 | 3.383 | -12.0  | 94    | 0.00     |
| 81 T  | trans-1,4-Dichloro-2-butene | 0.181 | 0.204 | -12.7  | 106   | 0.00     |
| 82 T  | 4-Chlorotoluene             | 2.785 | 2.993 | -7.5   | 94    | 0.00     |
| 83 T  | tert-Butylbenzene           | 2.462 | 2.737 | -11.2  | 95    | 0.00     |
| 84 T  | 1,2,4-Trimethylbenzene      | 2.996 | 3.347 | -11.7  | 94    | 0.00     |
| 85 T  | sec-Butylbenzene            | 3.869 | 4.369 | -12.9  | 94    | 0.00     |
| 86 T  | p-Isopropyltoluene          | 3.151 | 3.575 | -13.5  | 95    | 0.00     |
| 87 T  | 1,3-Dichlorobenzene         | 1.640 | 1.732 | -5.6   | 94    | 0.00     |
| 88 T  | 1,4-Dichlorobenzene         | 1.638 | 1.707 | -4.2   | 94    | 0.00     |
| 89 T  | n-Butylbenzene              | 3.122 | 3.524 | -12.9  | 95    | 0.00     |
| 90 T  | Hexachloroethane            | 0.641 | 0.689 | -7.5   | 95    | 0.00     |
| 91 T  | 1,2-Dichlorobenzene         | 1.414 | 1.502 | -6.2   | 95    | 0.00     |
| 92 T  | 1,2-Dibromo-3-Chloropropane | 0.106 | 0.107 | -0.9   | 94    | 0.00     |
| 93 T  | 1,2,4-Trichlorobenzene      | 0.786 | 0.853 | -8.5   | 95    | 0.00     |
| 94 T  | Hexachlorobutadiene         | 0.474 | 0.504 | -6.3   | 95    | 0.00     |
| 95 T  | Naphthalene                 | 1.426 | 1.539 | -7.9   | 94    | 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 | 0.675 | 0.747 | -10.7 | 98    | 0.00     |

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