

Data Path : Z:\voasrv\HPCHEM1\MSVOA\_D\Data\VD052322\  
 Data File : VD073315.D  
 Acq On : 24 May 2022 07:41  
 Operator : VA/SY  
 Sample : VSTDCCC050  
 Misc : 5.00G/5.00ml/MSVOA\_D/SOIL  
 ALS Vial : 47 Sample Multiplier: 1

Instrument :  
 MSVOA\_D  
 LabSampleId :  
 VSTDCCC050

Quant Time: May 24 08:19:00 2022  
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_D\Method\82D051922S.M  
 Quant Title : SW846 8260  
 QLast Update : Thu May 19 15:13:42 2022  
 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    | 87    | 0.00     |
| 2 T   | Dichlorodifluoromethane     | 0.500 | 0.409 | 18.2   | 81    | 0.00     |
| 3 P   | Chloromethane               | 0.565 | 0.478 | 15.4   | 84    | 0.00     |
| 4 C   | Vinyl Chloride              | 0.517 | 0.462 | 10.6#  | 83    | 0.00     |
| 5 T   | Bromomethane                | 0.374 | 0.340 | 9.1    | 88    | 0.00     |
| 6 T   | Chloroethane                | 0.318 | 0.309 | 2.8    | 90    | 0.00     |
| 7 T   | Trichlorofluoromethane      | 0.913 | 0.859 | 5.9    | 87    | 0.00     |
| 8 T   | Diethyl Ether               | 0.251 | 0.282 | -12.4  | 99    | 0.00     |
| 9 T   | 1,1,2-Trichlorotrifluoroeth | 0.545 | 0.503 | 7.7    | 87    | -0.01    |
| 10 T  | Methyl Iodide               | 0.588 | 0.593 | -0.9   | 88    | 0.00     |
| 11 T  | Tert butyl alcohol          | 0.086 | 0.067 | 22.1   | 130   | 0.00     |
| 12 CM | 1,1-Dichloroethene          | 0.502 | 0.488 | 2.8#   | 87    | 0.00     |
| 13 T  | Acrolein                    | 0.039 | 0.032 | 17.9   | 85    | 0.00     |
| 14 T  | Allyl chloride              | 0.811 | 0.821 | -1.2   | 91    | 0.00     |
| 15 T  | Acrylonitrile               | 0.115 | 0.148 | -28.7# | 115   | 0.00     |
| 16 T  | Acetone                     | 0.096 | 0.103 | -7.3   | 107   | 0.00     |
| 17 T  | Carbon Disulfide            | 1.666 | 1.397 | 16.1   | 78    | 0.00     |
| 18 T  | Methyl Acetate              | 0.266 | 0.331 | -24.4  | 116   | 0.00     |
| 19 T  | Methyl tert-butyl Ether     | 1.145 | 1.369 | -19.6  | 105   | 0.00     |
| 20 T  | Methylene Chloride          | 0.716 | 0.779 | -8.8   | 115   | 0.00     |
| 21 T  | trans-1,2-Dichloroethene    | 0.608 | 0.626 | -3.0   | 92    | 0.00     |
| 22 T  | Diisopropyl ether           | 1.700 | 1.943 | -14.3  | 99    | 0.00     |
| 23 T  | Vinyl Acetate               | 0.907 | 1.091 | -20.3  | 104   | 0.00     |
| 24 P  | 1,1-Dichloroethane          | 1.091 | 1.157 | -6.0   | 97    | 0.00     |
| 25 T  | 2-Butanone                  | 0.140 | 0.172 | -22.9  | 113   | 0.00     |
| 26 T  | 2,2-Dichloropropane         | 0.931 | 0.782 | 16.0   | 77    | 0.00     |
| 27 T  | cis-1,2-Dichloroethene      | 0.672 | 0.717 | -6.7   | 95    | 0.00     |
| 28 T  | Bromochloromethane          | 0.376 | 0.399 | -6.1   | 94    | 0.00     |
| 29 T  | Tetrahydrofuran             | 0.091 | 0.113 | -24.2  | 111   | 0.00     |
| 30 C  | Chloroform                  | 1.139 | 1.226 | -7.6#  | 98    | 0.00     |
| 31 T  | Cyclohexane                 | 0.966 | 0.833 | 13.8   | 80    | 0.00     |
| 32 T  | 1,1,1-Trichloroethane       | 0.997 | 1.026 | -2.9   | 94    | 0.00     |
| 33 S  | 1,2-Dichloroethane-d4       | 0.553 | 0.606 | -9.6   | 100   | 0.00     |
| 34 I  | 1,4-Difluorobenzene         | 1.000 | 1.000 | 0.0    | 92    | 0.00     |
| 35 S  | Dibromofluoromethane        | 0.330 | 0.353 | -7.0   | 98    | 0.00     |
| 36 T  | 1,1-Dichloropropene         | 0.496 | 0.481 | 3.0    | 91    | 0.00     |
| 37 T  | Ethyl Acetate               | 0.196 | 0.221 | -12.8  | 108   | 0.00     |
| 38 T  | Carbon Tetrachloride        | 0.503 | 0.497 | 1.2    | 91    | 0.00     |
| 39 T  | Methylcyclohexane           | 0.553 | 0.502 | 9.2    | 82    | 0.00     |
| 40 TM | Benzene                     | 1.430 | 1.473 | -3.0   | 97    | 0.00     |
| 41 T  | Methacrylonitrile           | 0.114 | 0.139 | -21.9  | 131   | 0.00     |
| 42 TM | 1,2-Dichloroethane          | 0.398 | 0.426 | -7.0   | 101   | 0.00     |
| 43 T  | Isopropyl Acetate           | 0.362 | 0.433 | -19.6  | 110   | 0.00     |
| 44 TM | Trichloroethene             | 0.381 | 0.383 | -0.5   | 94    | 0.00     |
| 45 C  | 1,2-Dichloropropane         | 0.360 | 0.384 | -6.7#  | 100   | 0.00     |
| 46 T  | Dibromomethane              | 0.198 | 0.214 | -8.1   | 102   | 0.00     |
| 47 T  | Bromodichloromethane        | 0.506 | 0.545 | -7.7   | 101   | 0.00     |
| 48 T  | Methyl methacrylate         | 0.182 | 0.203 | -11.5  | 102   | 0.00     |

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 Sample : VSTDCCC050  
 Misc : 5.00G/5.00ml/MSVOA\_D/SOIL  
 ALS Vial : 47 Sample Multiplier: 1

Instrument :  
 MSVOA\_D  
 LabSampleId :  
 VSTDCCC050

Quant Time: May 24 08:19:00 2022  
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_D\Method\82D051922S.M  
 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 | 0.003 | -50.0# | 116   | 0.00     |
| 50 S  | Toluene-d8                  | 1.216 | 1.294 | -6.4   | 96    | 0.00     |
| 51 T  | 4-Methyl-2-Pentanone        | 0.188 | 0.237 | -26.1# | 115   | 0.00     |
| 52 CM | Toluene                     | 0.904 | 0.958 | -6.0#  | 96    | 0.00     |
| 53 T  | t-1,3-Dichloropropene       | 0.457 | 0.494 | -8.1   | 102   | 0.00     |
| 54 T  | cis-1,3-Dichloropropene     | 0.547 | 0.573 | -4.8   | 97    | 0.00     |
| 55 T  | 1,1,2-Trichloroethane       | 0.267 | 0.304 | -13.9  | 107   | 0.00     |
| 56 T  | Ethyl methacrylate          | 0.297 | 0.375 | -26.3# | 109   | 0.00     |
| 57 T  | 1,3-Dichloropropane         | 0.449 | 0.509 | -13.4  | 104   | 0.00     |
| 58 T  | 2-Chloroethyl Vinyl ether   | 0.122 | 0.157 | -28.7# | 110   | 0.00     |
| 59 T  | 2-Hexanone                  | 0.124 | 0.160 | -29.0# | 116   | 0.00     |
| 60 T  | Dibromochloromethane        | 0.334 | 0.381 | -14.1  | 107   | 0.00     |
| 61 T  | 1,2-Dibromoethane           | 0.256 | 0.291 | -13.7  | 108   | 0.00     |
| 62 S  | 4-Bromofluorobenzene        | 0.427 | 0.477 | -11.7  | 102   | 0.00     |
| 63 I  | Chlorobenzene-d5            | 1.000 | 1.000 | 0.0    | 94    | 0.00     |
| 64 T  | Tetrachloroethene           | 0.327 | 0.324 | 0.9    | 96    | 0.00     |
| 65 PM | Chlorobenzene               | 1.026 | 1.069 | -4.2   | 100   | 0.00     |
| 66 T  | 1,1,1,2-Tetrachloroethane   | 0.380 | 0.415 | -9.2   | 104   | 0.00     |
| 67 C  | Ethyl Benzene               | 1.801 | 1.900 | -5.5#  | 97    | 0.00     |
| 68 T  | m/p-Xylenes                 | 0.700 | 0.744 | -6.3   | 96    | 0.00     |
| 69 T  | o-Xylene                    | 0.643 | 0.698 | -8.6   | 99    | 0.00     |
| 70 T  | Styrene                     | 1.106 | 1.236 | -11.8  | 100   | 0.00     |
| 71 P  | Bromoform                   | 0.196 | 0.227 | -15.8  | 111   | 0.00     |
| 72 I  | 1,4-Dichlorobenzene-d4      | 1.000 | 1.000 | 0.0    | 95    | 0.00     |
| 73 T  | Isopropylbenzene            | 3.564 | 3.804 | -6.7   | 97    | 0.00     |
| 74 T  | N-amyl acetate              | 0.778 | 0.917 | -17.9  | 110   | 0.00     |
| 75 P  | 1,1,2,2-Tetrachloroethane   | 0.657 | 0.766 | -16.6  | 113   | 0.00     |
| 76 T  | 1,2,3-Trichloropropane      | 0.491 | 0.544 | -10.8  | 107   | 0.00     |
| 77 T  | Bromobenzene                | 0.839 | 0.900 | -7.3   | 103   | 0.00     |
| 78 T  | n-propylbenzene             | 4.475 | 4.680 | -4.6   | 95    | 0.00     |
| 79 T  | 2-Chlorotoluene             | 2.584 | 2.737 | -5.9   | 99    | 0.00     |
| 80 T  | 1,3,5-Trimethylbenzene      | 3.019 | 3.236 | -7.2   | 98    | 0.00     |
| 81 T  | trans-1,4-Dichloro-2-butene | 0.204 | 0.216 | -5.9   | 98    | 0.00     |
| 82 T  | 4-Chlorotoluene             | 2.708 | 2.908 | -7.4   | 100   | 0.00     |
| 83 T  | tert-Butylbenzene           | 2.545 | 2.751 | -8.1   | 99    | 0.00     |
| 84 T  | 1,2,4-Trimethylbenzene      | 2.983 | 3.253 | -9.1   | 101   | 0.00     |
| 85 T  | sec-Butylbenzene            | 3.886 | 4.097 | -5.4   | 97    | 0.00     |
| 86 T  | p-Isopropyltoluene          | 3.167 | 3.382 | -6.8   | 98    | 0.00     |
| 87 T  | 1,3-Dichlorobenzene         | 1.691 | 1.776 | -5.0   | 101   | 0.00     |
| 88 T  | 1,4-Dichlorobenzene         | 1.707 | 1.773 | -3.9   | 102   | 0.00     |
| 89 T  | n-Butylbenzene              | 3.018 | 3.124 | -3.5   | 94    | 0.00     |
| 90 T  | Hexachloroethane            | 0.655 | 0.685 | -4.6   | 101   | 0.00     |
| 91 T  | 1,2-Dichlorobenzene         | 1.463 | 1.597 | -9.2   | 104   | 0.00     |
| 92 T  | 1,2-Dibromo-3-Chloropropane | 0.101 | 0.115 | -13.9  | 113   | 0.00     |
| 93 T  | 1,2,4-Trichlorobenzene      | 0.831 | 0.910 | -9.5   | 101   | 0.00     |
| 94 T  | Hexachlorobutadiene         | 0.468 | 0.467 | 0.2    | 99    | 0.00     |
| 95 T  | Naphthalene                 | 1.466 | 1.835 | -25.2# | 114   | 0.00     |

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MSVOA\_D  
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Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_D\Method\82D051922S.M  
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) |
|------|------------------------|-------|-------|-------|-------|----------|
| 96 T | 1,2,3-Trichlorobenzene | 0.737 | 0.820 | -11.3 | 103   | 0.00     |

(#) = Out of Range

SPCC's out = 0 CCC's out = 6