

Data Path : Z:\voasrv\HPCHEM1\MSVOA\_D\Data\VD101424\  
 Data File : VD079970.D  
 Acq On : 14 Oct 2024 10:16  
 Operator : RP/MD  
 Sample : VSTDCCC050  
 Misc : 5.00G/5ml/MSVOA\_D/SOIL  
 ALS Vial : 2 Sample Multiplier: 1

Instrument :  
 MSVOA\_D  
 LabSampleId :  
 VSTDCCC050

Quant Time: Oct 15 03:10:03 2024  
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_D\Method\82D101124S.M  
 Quant Title : SW846 8260  
 QLast Update : Mon Oct 14 03:20:54 2024  
 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   | 95    | 0.00     |
| 2 T   | Dichlorodifluoromethane     | 0.491 | 0.469 | 4.5   | 95    | 0.00     |
| 3 P   | Chloromethane               | 0.505 | 0.477 | 5.5   | 95    | 0.00     |
| 4 C   | Vinyl Chloride              | 0.628 | 0.607 | 3.3#  | 93    | 0.00     |
| 5 T   | Bromomethane                | 0.540 | 0.453 | 16.1  | 85    | 0.00     |
| 6 T   | Chloroethane                | 0.428 | 0.398 | 7.0   | 88    | 0.00     |
| 7 T   | Trichlorofluoromethane      | 0.983 | 0.997 | -1.4  | 99    | 0.00     |
| 8 T   | Diethyl Ether               | 0.270 | 0.307 | -13.7 | 104   | 0.00     |
| 9 T   | 1,1,2-Trichlorotrifluoroeth | 0.604 | 0.611 | -1.2  | 102   | 0.00     |
| 10 T  | Methyl Iodide               | 0.592 | 0.601 | -1.5  | 87    | 0.00     |
| 11 T  | Tert butyl alcohol          | 0.035 | 0.041 | -17.1 | 112   | 0.00     |
| 12 CM | 1,1-Dichloroethene          | 0.537 | 0.530 | 1.3#  | 94    | 0.00     |
| 13 T  | Acrolein                    | 0.037 | 0.030 | 18.9  | 88    | 0.00     |
| 14 T  | Allyl chloride              | 0.798 | 0.852 | -6.8  | 99    | 0.00     |
| 15 T  | Acrylonitrile               | 0.129 | 0.148 | -14.7 | 106   | 0.00     |
| 16 T  | Acetone                     | 0.147 | 0.171 | -16.3 | 105   | 0.00     |
| 17 T  | Carbon Disulfide            | 1.497 | 1.373 | 8.3   | 82    | 0.00     |
| 18 T  | Methyl Acetate              | 0.316 | 0.357 | -13.0 | 105   | 0.00     |
| 19 T  | Methyl tert-butyl Ether     | 1.241 | 1.461 | -17.7 | 107   | 0.00     |
| 20 T  | Methylene Chloride          | 0.724 | 0.683 | 5.7   | 97    | 0.00     |
| 21 T  | trans-1,2-Dichloroethene    | 0.577 | 0.614 | -6.4  | 99    | 0.00     |
| 22 T  | Diisopropyl ether           | 1.731 | 2.016 | -16.5 | 105   | 0.00     |
| 23 T  | Vinyl Acetate               | 0.972 | 1.180 | -21.4 | 107   | 0.00     |
| 24 P  | 1,1-Dichloroethane          | 1.140 | 1.184 | -3.9  | 101   | 0.00     |
| 25 T  | 2-Butanone                  | 0.180 | 0.212 | -17.8 | 107   | 0.00     |
| 26 T  | 2,2-Dichloropropane         | 0.986 | 1.034 | -4.9  | 105   | 0.00     |
| 27 T  | cis-1,2-Dichloroethene      | 0.717 | 0.764 | -6.6  | 99    | 0.00     |
| 28 T  | Bromochloromethane          | 0.487 | 0.488 | -0.2  | 95    | 0.00     |
| 29 T  | Tetrahydrofuran             | 0.096 | 0.119 | -24.0 | 110   | 0.00     |
| 30 C  | Chloroform                  | 1.208 | 1.258 | -4.1# | 100   | 0.00     |
| 31 T  | Cyclohexane                 | 0.864 | 0.838 | 3.0   | 94    | 0.00     |
| 32 T  | 1,1,1-Trichloroethane       | 1.000 | 1.034 | -3.4  | 100   | 0.00     |
| 33 S  | 1,2-Dichloroethane-d4       | 0.569 | 0.578 | -1.6  | 93    | 0.00     |
| 34 I  | 1,4-Difluorobenzene         | 1.000 | 1.000 | 0.0   | 97    | 0.00     |
| 35 S  | Dibromofluoromethane        | 0.361 | 0.354 | 1.9   | 92    | 0.00     |
| 36 T  | 1,1-Dichloropropene         | 0.484 | 0.494 | -2.1  | 98    | 0.00     |
| 37 T  | Ethyl Acetate               | 0.234 | 0.256 | -9.4  | 104   | 0.00     |
| 38 T  | Carbon Tetrachloride        | 0.537 | 0.541 | -0.7  | 98    | 0.00     |
| 39 T  | Methylcyclohexane           | 0.543 | 0.565 | -4.1  | 96    | 0.00     |
| 40 TM | Benzene                     | 1.530 | 1.591 | -4.0  | 99    | 0.00     |
| 41 T  | Methacrylonitrile           | 0.123 | 0.146 | -18.7 | 114   | 0.00     |
| 42 TM | 1,2-Dichloroethane          | 0.421 | 0.453 | -7.6  | 104   | 0.00     |
| 43 T  | Isopropyl Acetate           | 0.431 | 0.488 | -13.2 | 108   | 0.00     |
| 44 TM | Trichloroethene             | 0.378 | 0.383 | -1.3  | 98    | 0.00     |
| 45 C  | 1,2-Dichloropropane         | 0.388 | 0.424 | -9.3# | 104   | 0.00     |
| 46 T  | Dibromomethane              | 0.214 | 0.236 | -10.3 | 104   | 0.00     |
| 47 T  | Bromodichloromethane        | 0.568 | 0.612 | -7.7  | 103   | 0.00     |
| 48 T  | Methyl methacrylate         | 0.194 | 0.234 | -20.6 | 107   | 0.00     |

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 ALS Vial : 2 Sample Multiplier: 1

Instrument :  
 MSVOA\_D  
 LabSampleId :  
 VSTDCCC050

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 Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_D\Method\82D101124S.M  
 Quant Title : SW846 8260  
 QLast Update : Mon Oct 14 03:20:54 2024  
 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                  | 1.286 | 1.276 | 0.8    | 90    | 0.00     |
| 51 T  | 4-Methyl-2-Pentanone        | 0.221 | 0.271 | -22.6  | 111   | 0.00     |
| 52 CM | Toluene                     | 0.936 | 1.005 | -7.4#  | 98    | 0.00     |
| 53 T  | t-1,3-Dichloropropene       | 0.497 | 0.555 | -11.7  | 106   | 0.00     |
| 54 T  | cis-1,3-Dichloropropene     | 0.586 | 0.640 | -9.2   | 102   | 0.00     |
| 55 T  | 1,1,2-Trichloroethane       | 0.301 | 0.324 | -7.6   | 103   | 0.00     |
| 56 T  | Ethyl methacrylate          | 0.351 | 0.425 | -21.1  | 107   | 0.00     |
| 57 T  | 1,3-Dichloropropane         | 0.485 | 0.536 | -10.5  | 102   | 0.00     |
| 58 T  | 2-Chloroethyl Vinyl ether   | 0.130 | 0.162 | -24.6  | 104   | 0.00     |
| 59 T  | 2-Hexanone                  | 0.163 | 0.200 | -22.7  | 107   | 0.00     |
| 60 T  | Dibromochloromethane        | 0.391 | 0.415 | -6.1   | 101   | 0.00     |
| 61 T  | 1,2-Dibromoethane           | 0.277 | 0.295 | -6.5   | 101   | 0.00     |
| 62 S  | 4-Bromofluorobenzene        | 0.449 | 0.473 | -5.3   | 96    | 0.00     |
| 63 I  | Chlorobenzene-d5            | 1.000 | 1.000 | 0.0    | 98    | 0.00     |
| 64 T  | Tetrachloroethene           | 0.353 | 0.361 | -2.3   | 99    | 0.00     |
| 65 PM | Chlorobenzene               | 1.164 | 1.238 | -6.4   | 104   | 0.00     |
| 66 T  | 1,1,1,2-Tetrachloroethane   | 0.418 | 0.441 | -5.5   | 105   | 0.00     |
| 67 C  | Ethyl Benzene               | 1.968 | 2.137 | -8.6#  | 102   | 0.00     |
| 68 T  | m/p-Xylenes                 | 0.755 | 0.829 | -9.8   | 102   | 0.00     |
| 69 T  | o-Xylene                    | 0.699 | 0.775 | -10.9  | 105   | 0.00     |
| 70 T  | Styrene                     | 1.244 | 1.408 | -13.2  | 103   | 0.00     |
| 71 P  | Bromoform                   | 0.251 | 0.271 | -8.0   | 104   | 0.00     |
| 72 I  | 1,4-Dichlorobenzene-d4      | 1.000 | 1.000 | 0.0    | 99    | 0.00     |
| 73 T  | Isopropylbenzene            | 3.401 | 3.817 | -12.2  | 109   | 0.00     |
| 74 T  | N-amyl acetate              | 0.823 | 0.952 | -15.7  | 107   | 0.00     |
| 75 P  | 1,1,2,2-Tetrachloroethane   | 0.690 | 0.756 | -9.6   | 107   | 0.00     |
| 76 T  | 1,2,3-Trichloropropane      | 0.552 | 0.500 | 9.4    | 91    | 0.00     |
| 77 T  | Bromobenzene                | 0.880 | 0.928 | -5.5   | 103   | 0.00     |
| 78 T  | n-propylbenzene             | 4.229 | 4.629 | -9.5   | 105   | 0.00     |
| 79 T  | 2-Chlorotoluene             | 2.457 | 2.633 | -7.2   | 103   | 0.00     |
| 80 T  | 1,3,5-Trimethylbenzene      | 2.867 | 3.102 | -8.2   | 105   | 0.00     |
| 81 T  | trans-1,4-Dichloro-2-butene | 0.247 | 0.285 | -15.4  | 110   | 0.00     |
| 82 T  | 4-Chlorotoluene             | 2.639 | 2.837 | -7.5   | 104   | 0.00     |
| 83 T  | tert-Butylbenzene           | 2.457 | 2.772 | -12.8  | 109   | 0.00     |
| 84 T  | 1,2,4-Trimethylbenzene      | 2.879 | 3.234 | -12.3  | 106   | 0.00     |
| 85 T  | sec-Butylbenzene            | 3.776 | 4.132 | -9.4   | 106   | 0.00     |
| 86 T  | p-Isopropyltoluene          | 3.150 | 3.520 | -11.7  | 107   | 0.00     |
| 87 T  | 1,3-Dichlorobenzene         | 1.800 | 1.900 | -5.6   | 105   | 0.00     |
| 88 T  | 1,4-Dichlorobenzene         | 1.799 | 1.893 | -5.2   | 107   | 0.00     |
| 89 T  | n-Butylbenzene              | 3.047 | 3.400 | -11.6  | 107   | 0.00     |
| 90 T  | Hexachloroethane            | 0.696 | 0.704 | -1.1   | 104   | 0.00     |
| 91 T  | 1,2-Dichlorobenzene         | 1.573 | 1.667 | -6.0   | 105   | 0.00     |
| 92 T  | 1,2-Dibromo-3-Chloropropane | 0.115 | 0.119 | -3.5   | 105   | 0.00     |
| 93 T  | 1,2,4-Trichlorobenzene      | 0.986 | 1.094 | -11.0  | 109   | 0.00     |
| 94 T  | Hexachlorobutadiene         | 0.543 | 0.589 | -8.5   | 106   | 0.00     |
| 95 T  | Naphthalene                 | 1.758 | 2.097 | -19.3  | 112   | 0.00     |

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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) |
|------|------------------------|-------|-------|-------|-------|----------|
| 96 T | 1,2,3-Trichlorobenzene | 0.873 | 0.999 | -14.4 | 111   | 0.00     |

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