

Data Path : Z:\voasrv\HPCHEM1\MSVOA\_D\Data\VD090525\  
 Data File : VD080520.D  
 Acq On : 05 Sep 2025 12:21  
 Operator : RP/MD  
 Sample : VSTDCCC050  
 Misc : 5.00G/5ml/MSVOA\_D/SOIL  
 ALS Vial : 8 Sample Multiplier: 1

Instrument :  
 MSVOA\_D  
 LabSampleId :  
 VSTDCCC050

Quant Time: Sep 06 03:01:28 2025  
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_D\Method\82D082725S.M  
 Quant Title : SW846 8260  
 QLast Update : Thu Aug 28 03:02:37 2025  
 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   | 79    | 0.00     |
| 2 T   | Dichlorodifluoromethane     | 0.426 | 0.372 | 12.7  | 81    | 0.00     |
| 3 P   | Chloromethane               | 0.672 | 0.635 | 5.5   | 82    | 0.00     |
| 4 C   | Vinyl Chloride              | 0.694 | 0.708 | -2.0# | 86    | 0.00     |
| 5 T   | Bromomethane                | 0.390 | 0.369 | 5.4   | 84    | 0.00     |
| 6 T   | Chloroethane                | 0.453 | 0.482 | -6.4  | 88    | 0.00     |
| 7 T   | Trichlorofluoromethane      | 0.808 | 0.814 | -0.7  | 85    | 0.00     |
| 8 T   | Diethyl Ether               | 0.320 | 0.294 | 8.1   | 77    | 0.00     |
| 9 T   | 1,1,2-Trichlorotrifluoroeth | 0.524 | 0.519 | 1.0   | 83    | 0.00     |
| 10 T  | Methyl Iodide               | 0.605 | 0.578 | 4.5   | 74    | 0.00     |
| 11 T  | Tert butyl alcohol          | 0.041 | 0.039 | 4.9   | 82    | 0.00     |
| 12 CM | 1,1-Dichloroethene          | 0.546 | 0.529 | 3.1#  | 78    | 0.00     |
| 13 T  | Acrolein                    | 0.043 | 0.023 | 46.5# | 44#   | 0.00     |
| 14 T  | Allyl chloride              | 0.948 | 0.866 | 8.6   | 74    | 0.00     |
| 15 T  | Acrylonitrile               | 0.143 | 0.136 | 4.9   | 78    | 0.00     |
| 16 T  | Acetone                     | 0.132 | 0.129 | 2.3   | 78    | 0.00     |
| 17 T  | Carbon Disulfide            | 1.846 | 1.695 | 8.2   | 76    | 0.00     |
| 18 T  | Methyl Acetate              | 0.372 | 0.372 | 0.0   | 87    | -0.01    |
| 19 T  | Methyl tert-butyl Ether     | 1.433 | 1.390 | 3.0   | 79    | 0.00     |
| 20 T  | Methylene Chloride          | 0.725 | 0.641 | 11.6  | 77    | 0.00     |
| 21 T  | trans-1,2-Dichloroethene    | 0.625 | 0.594 | 5.0   | 78    | 0.00     |
| 22 T  | Diisopropyl ether           | 1.950 | 1.818 | 6.8   | 75    | 0.00     |
| 23 T  | Vinyl Acetate               | 1.110 | 1.052 | 5.2   | 76    | 0.00     |
| 24 P  | 1,1-Dichloroethane          | 1.129 | 1.088 | 3.6   | 79    | 0.00     |
| 25 T  | 2-Butanone                  | 0.187 | 0.178 | 4.8   | 78    | 0.00     |
| 26 T  | 2,2-Dichloropropane         | 0.894 | 0.882 | 1.3   | 80    | 0.00     |
| 27 T  | cis-1,2-Dichloroethene      | 0.746 | 0.707 | 5.2   | 76    | 0.00     |
| 28 T  | Bromochloromethane          | 0.460 | 0.446 | 3.0   | 78    | 0.00     |
| 29 T  | Tetrahydrofuran             | 0.118 | 0.114 | 3.4   | 79    | 0.00     |
| 30 C  | Chloroform                  | 1.079 | 1.078 | 0.1#  | 80    | 0.00     |
| 31 T  | Cyclohexane                 | 1.057 | 0.967 | 8.5   | 78    | 0.00     |
| 32 T  | 1,1,1-Trichloroethane       | 0.899 | 0.920 | -2.3  | 83    | 0.00     |
| 33 S  | 1,2-Dichloroethane-d4       | 0.564 | 0.485 | 14.0  | 76    | 0.00     |
| 34 I  | 1,4-Difluorobenzene         | 1.000 | 1.000 | 0.0   | 76    | 0.00     |
| 35 S  | Dibromofluoromethane        | 0.331 | 0.293 | 11.5  | 76    | 0.00     |
| 36 T  | 1,1-Dichloropropene         | 0.438 | 0.453 | -3.4  | 80    | 0.00     |
| 37 T  | Ethyl Acetate               | 0.211 | 0.219 | -3.8  | 78    | 0.00     |
| 38 T  | Carbon Tetrachloride        | 0.417 | 0.456 | -9.4  | 85    | 0.00     |
| 39 T  | Methylcyclohexane           | 0.583 | 0.603 | -3.4  | 80    | 0.00     |
| 40 TM | Benzene                     | 1.394 | 1.412 | -1.3  | 77    | 0.00     |
| 41 T  | Methacrylonitrile           | 0.125 | 0.143 | -14.4 | 86    | 0.00     |
| 42 TM | 1,2-Dichloroethane          | 0.338 | 0.355 | -5.0  | 81    | 0.00     |
| 43 T  | Isopropyl Acetate           | 0.429 | 0.436 | -1.6  | 79    | 0.00     |
| 44 TM | Trichloroethene             | 0.338 | 0.351 | -3.8  | 80    | 0.00     |
| 45 C  | 1,2-Dichloropropane         | 0.346 | 0.350 | -1.2# | 78    | 0.00     |
| 46 T  | Dibromomethane              | 0.185 | 0.197 | -6.5  | 81    | 0.00     |
| 47 T  | Bromodichloromethane        | 0.452 | 0.481 | -6.4  | 82    | 0.00     |
| 48 T  | Methyl methacrylate         | 0.213 | 0.216 | -1.4  | 77    | 0.00     |

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 Sample : VSTDCCC050  
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 ALS Vial : 8 Sample Multiplier: 1

Instrument :  
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 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 | 0.003 | -50.0# | 84    | 0.00     |
| 50 S  | Toluene-d8                  | 1.306 | 1.160 | 11.2   | 75    | 0.00     |
| 51 T  | 4-Methyl-2-Pentanone        | 0.214 | 0.222 | -3.7   | 80    | 0.00     |
| 52 CM | Toluene                     | 0.870 | 0.899 | -3.3#  | 79    | 0.00     |
| 53 T  | t-1,3-Dichloropropene       | 0.460 | 0.470 | -2.2   | 78    | 0.00     |
| 54 T  | cis-1,3-Dichloropropene     | 0.550 | 0.560 | -1.8   | 78    | 0.00     |
| 55 T  | 1,1,2-Trichloroethane       | 0.248 | 0.262 | -5.6   | 81    | 0.00     |
| 56 T  | Ethyl methacrylate          | 0.346 | 0.355 | -2.6   | 78    | 0.00     |
| 57 T  | 1,3-Dichloropropane         | 0.436 | 0.453 | -3.9   | 80    | 0.00     |
| 58 T  | 2-Chloroethyl Vinyl ether   | 0.134 | 0.140 | -4.5   | 78    | 0.00     |
| 59 T  | 2-Hexanone                  | 0.152 | 0.158 | -3.9   | 79    | 0.00     |
| 60 T  | Dibromochloromethane        | 0.313 | 0.331 | -5.8   | 81    | 0.00     |
| 61 T  | 1,2-Dibromoethane           | 0.236 | 0.254 | -7.6   | 81    | 0.00     |
| 62 S  | 4-Bromofluorobenzene        | 0.425 | 0.363 | 14.6   | 76    | 0.00     |
| 63 I  | Chlorobenzene-d5            | 1.000 | 1.000 | 0.0    | 76    | 0.00     |
| 64 T  | Tetrachloroethene           | 0.309 | 0.335 | -8.4   | 83    | 0.00     |
| 65 PM | Chlorobenzene               | 1.065 | 1.127 | -5.8   | 80    | 0.00     |
| 66 T  | 1,1,1,2-Tetrachloroethane   | 0.340 | 0.363 | -6.8   | 81    | 0.00     |
| 67 C  | Ethyl Benzene               | 1.844 | 1.949 | -5.7#  | 80    | 0.00     |
| 68 T  | m/p-Xylenes                 | 0.710 | 0.751 | -5.8   | 80    | 0.00     |
| 69 T  | o-Xylene                    | 0.678 | 0.713 | -5.2   | 79    | 0.00     |
| 70 T  | Styrene                     | 1.153 | 1.220 | -5.8   | 79    | 0.00     |
| 71 P  | Bromoform                   | 0.193 | 0.215 | -11.4  | 85    | 0.00     |
| 72 I  | 1,4-Dichlorobenzene-d4      | 1.000 | 1.000 | 0.0    | 79    | 0.00     |
| 73 T  | Isopropylbenzene            | 3.663 | 3.780 | -3.2   | 81    | 0.00     |
| 74 T  | N-amyl acetate              | 0.972 | 0.981 | -0.9   | 79    | 0.00     |
| 75 P  | 1,1,2,2-Tetrachloroethane   | 0.681 | 0.714 | -4.8   | 83    | 0.00     |
| 76 T  | 1,2,3-Trichloropropane      | 0.408 | 0.425 | -4.2   | 83    | 0.00     |
| 77 T  | Bromobenzene                | 0.848 | 0.875 | -3.2   | 81    | 0.00     |
| 78 T  | n-propylbenzene             | 4.448 | 4.585 | -3.1   | 81    | 0.00     |
| 79 T  | 2-Chlorotoluene             | 2.535 | 2.546 | -0.4   | 79    | 0.00     |
| 80 T  | 1,3,5-Trimethylbenzene      | 2.977 | 3.078 | -3.4   | 82    | 0.00     |
| 81 T  | trans-1,4-Dichloro-2-butene | 0.269 | 0.268 | 0.4    | 79    | 0.00     |
| 82 T  | 4-Chlorotoluene             | 2.607 | 2.671 | -2.5   | 81    | 0.00     |
| 83 T  | tert-Butylbenzene           | 2.583 | 2.721 | -5.3   | 83    | 0.00     |
| 84 T  | 1,2,4-Trimethylbenzene      | 2.997 | 3.091 | -3.1   | 81    | 0.00     |
| 85 T  | sec-Butylbenzene            | 3.855 | 4.005 | -3.9   | 82    | 0.00     |
| 86 T  | p-Isopropyltoluene          | 3.224 | 3.380 | -4.8   | 83    | 0.00     |
| 87 T  | 1,3-Dichlorobenzene         | 1.634 | 1.731 | -5.9   | 83    | 0.00     |
| 88 T  | 1,4-Dichlorobenzene         | 1.641 | 1.693 | -3.2   | 82    | 0.00     |
| 89 T  | n-Butylbenzene              | 3.072 | 3.210 | -4.5   | 82    | 0.00     |
| 90 T  | Hexachloroethane            | 0.624 | 0.647 | -3.7   | 84    | 0.00     |
| 91 T  | 1,2-Dichlorobenzene         | 1.454 | 1.510 | -3.9   | 82    | 0.00     |
| 92 T  | 1,2-Dibromo-3-Chloropropane | 0.102 | 0.106 | -3.9   | 86    | 0.00     |
| 93 T  | 1,2,4-Trichlorobenzene      | 0.911 | 0.958 | -5.2   | 84    | 0.00     |
| 94 T  | Hexachlorobutadiene         | 0.414 | 0.452 | -9.2   | 89    | 0.00     |
| 95 T  | Naphthalene                 | 1.938 | 2.011 | -3.8   | 83    | 0.00     |

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Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_D\Method\82D082725S.M  
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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 | 0.780 | 0.832 | -6.7 | 85    | 0.00     |

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