

Data Path : Z:\voasrv\HPCHEM1\MSVOA\_X\Data\VX062222\  
 Data File : VX029685.D  
 Acq On : 22 Jun 2022 09:22  
 Operator : JC/MD  
 Sample : VSTDCCC050  
 Misc : 5.0mL/MSVOA\_X/WATER  
 ALS Vial : 2 Sample Multiplier: 1

Instrument :  
 MSVOA\_X  
 LabSampleId :  
 VSTDCCC050

Quant Time: Jun 23 01:07:20 2022  
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_X\Method\82X061822W.M  
 Quant Title : SW846 8260  
 QLast Update : Mon Jun 20 01:31:01 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   | 98    | 0.00     |
| 2 T   | Dichlorodifluoromethane     | 0.606 | 0.594 | 2.0   | 90    | 0.00     |
| 3 P   | Chloromethane               | 0.652 | 0.529 | 18.9  | 82    | 0.00     |
| 4 C   | Vinyl Chloride              | 0.601 | 0.541 | 10.0# | 86    | 0.00     |
| 5 T   | Bromomethane                | 0.306 | 0.258 | 15.7  | 90    | 0.00     |
| 6 T   | Chloroethane                | 0.346 | 0.315 | 9.0   | 85    | 0.00     |
| 7 T   | Trichlorofluoromethane      | 0.837 | 0.773 | 7.6   | 88    | 0.00     |
| 8 T   | Diethyl Ether               | 0.330 | 0.308 | 6.7   | 90    | 0.00     |
| 9 T   | 1,1,2-Trichlorotrifluoroeth | 0.525 | 0.515 | 1.9   | 96    | 0.00     |
| 10 T  | Methyl Iodide               | 0.515 | 0.379 | 26.4# | 63    | 0.00     |
| 11 T  | Tert butyl alcohol          | 0.165 | 0.117 | 29.1# | 83    | 0.00     |
| 12 CM | 1,1-Dichloroethene          | 0.507 | 0.506 | 0.2#  | 97    | 0.00     |
| 13 T  | Acrolein                    | 0.036 | 0.029 | 19.4  | 82    | 0.00     |
| 14 T  | Allyl chloride              | 0.818 | 0.786 | 3.9   | 92    | 0.00     |
| 15 T  | Acrylonitrile               | 0.310 | 0.297 | 4.2   | 92    | 0.00     |
| 16 T  | Acetone                     | 0.264 | 0.219 | 17.0  | 86    | 0.00     |
| 17 T  | Carbon Disulfide            | 1.313 | 1.337 | -1.8  | 98    | 0.00     |
| 18 T  | Methyl Acetate              | 0.698 | 0.649 | 7.0   | 92    | 0.00     |
| 19 T  | Methyl tert-butyl Ether     | 1.759 | 1.716 | 2.4   | 94    | 0.00     |
| 20 T  | Methylene Chloride          | 0.588 | 0.585 | 0.5   | 100   | 0.00     |
| 21 T  | trans-1,2-Dichloroethene    | 0.554 | 0.560 | -1.1  | 97    | 0.00     |
| 22 T  | Diisopropyl ether           | 1.618 | 1.530 | 5.4   | 91    | 0.00     |
| 23 T  | Vinyl Acetate               | 1.351 | 1.315 | 2.7   | 91    | 0.00     |
| 24 P  | 1,1-Dichloroethane          | 0.973 | 0.932 | 4.2   | 93    | 0.00     |
| 25 T  | 2-Butanone                  | 0.417 | 0.373 | 10.6  | 87    | 0.00     |
| 26 T  | 2,2-Dichloropropane         | 0.757 | 0.729 | 3.7   | 91    | 0.00     |
| 27 T  | cis-1,2-Dichloroethene      | 0.647 | 0.637 | 1.5   | 95    | 0.00     |
| 28 T  | Bromochloromethane          | 0.372 | 0.363 | 2.4   | 96    | 0.00     |
| 29 T  | Tetrahydrofuran             | 0.278 | 0.249 | 10.4  | 86    | 0.00     |
| 30 C  | Chloroform                  | 1.030 | 0.981 | 4.8#  | 92    | 0.00     |
| 31 T  | Cyclohexane                 | 0.889 | 0.845 | 4.9   | 92    | 0.00     |
| 32 T  | 1,1,1-Trichloroethane       | 0.908 | 0.859 | 5.4   | 90    | 0.00     |
| 33 S  | 1,2-Dichloroethane-d4       | 0.662 | 0.587 | 11.3  | 87    | 0.00     |
| 34 I  | 1,4-Difluorobenzene         | 1.000 | 1.000 | 0.0   | 96    | 0.00     |
| 35 S  | Dibromofluoromethane        | 0.326 | 0.319 | 2.1   | 91    | 0.00     |
| 36 T  | 1,1-Dichloropropene         | 0.441 | 0.436 | 1.1   | 92    | 0.00     |
| 37 T  | Ethyl Acetate               | 0.452 | 0.438 | 3.1   | 87    | 0.00     |
| 38 T  | Carbon Tetrachloride        | 0.434 | 0.440 | -1.4  | 92    | 0.00     |
| 39 T  | Methylcyclohexane           | 0.550 | 0.571 | -3.8  | 95    | 0.00     |
| 40 TM | Benzene                     | 1.307 | 1.319 | -0.9  | 94    | 0.00     |
| 41 T  | Methacrylonitrile           | 0.238 | 0.235 | 1.3   | 87    | 0.00     |
| 42 TM | 1,2-Dichloroethane          | 0.449 | 0.432 | 3.8   | 92    | 0.00     |
| 43 T  | Isopropyl Acetate           | 0.703 | 0.684 | 2.7   | 88    | 0.00     |
| 44 TM | Trichloroethene             | 0.448 | 0.396 | 11.6  | 101   | 0.00     |
| 45 C  | 1,2-Dichloropropane         | 0.321 | 0.327 | -1.9# | 96    | 0.00     |
| 46 T  | Dibromomethane              | 0.230 | 0.242 | -5.2  | 96    | 0.00     |
| 47 T  | Bromodichloromethane        | 0.444 | 0.457 | -2.9  | 94    | 0.00     |
| 48 T  | Methyl methacrylate         | 0.345 | 0.331 | 4.1   | 87    | 0.00     |

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 Sample : VSTDCCC050  
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 ALS Vial : 2 Sample Multiplier: 1

Instrument :  
 MSVOA\_X  
 LabSampleId :  
 VSTDCCC050

Quant Time: Jun 23 01:07:20 2022  
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_X\Method\82X061822W.M  
 Quant Title : SW846 8260  
 QLast Update : Mon Jun 20 01:31:01 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) |
|-------|-----------------------------|-------|-------|-------|-------|----------|
| 49 T  | 1,4-Dioxane                 | 0.009 | 0.007 | 22.2  | 79    | 0.00     |
| 50 S  | Toluene-d8                  | 1.190 | 1.180 | 0.8   | 90    | 0.00     |
| 51 T  | 4-Methyl-2-Pentanone        | 0.453 | 0.429 | 5.3   | 87    | 0.00     |
| 52 CM | Toluene                     | 0.837 | 0.866 | -3.5# | 94    | 0.00     |
| 53 T  | t-1,3-Dichloropropene       | 0.449 | 0.492 | -9.6  | 94    | 0.00     |
| 54 T  | cis-1,3-Dichloropropene     | 0.502 | 0.549 | -9.4  | 95    | 0.00     |
| 55 T  | 1,1,2-Trichloroethane       | 0.338 | 0.354 | -4.7  | 96    | 0.00     |
| 56 T  | Ethyl methacrylate          | 0.511 | 0.547 | -7.0  | 94    | 0.00     |
| 57 T  | 1,3-Dichloropropane         | 0.564 | 0.577 | -2.3  | 95    | 0.00     |
| 58 T  | 2-Chloroethyl Vinyl ether   | 0.240 | 0.222 | 7.5   | 83    | 0.00     |
| 59 T  | 2-Hexanone                  | 0.343 | 0.326 | 5.0   | 86    | 0.00     |
| 60 T  | Dibromochloromethane        | 0.329 | 0.365 | -10.9 | 98    | 0.00     |
| 61 T  | 1,2-Dibromoethane           | 0.351 | 0.373 | -6.3  | 96    | 0.00     |
| 62 S  | 4-Bromofluorobenzene        | 0.449 | 0.464 | -3.3  | 93    | 0.00     |
| 63 I  | Chlorobenzene-d5            | 1.000 | 1.000 | 0.0   | 95    | 0.00     |
| 64 T  | Tetrachloroethene           | 0.362 | 0.436 | -20.4 | 113   | 0.00     |
| 65 PM | Chlorobenzene               | 0.981 | 1.032 | -5.2  | 97    | 0.00     |
| 66 T  | 1,1,1,2-Tetrachloroethane   | 0.352 | 0.371 | -5.4  | 94    | 0.00     |
| 67 C  | Ethyl Benzene               | 1.765 | 1.815 | -2.8# | 94    | 0.00     |
| 68 T  | m/p-Xylenes                 | 0.688 | 0.717 | -4.2  | 96    | 0.00     |
| 69 T  | o-Xylene                    | 0.680 | 0.700 | -2.9  | 95    | 0.00     |
| 70 T  | Styrene                     | 1.113 | 1.190 | -6.9  | 95    | 0.00     |
| 71 P  | Bromoform                   | 0.250 | 0.285 | -14.0 | 98    | 0.00     |
| 72 I  | 1,4-Dichlorobenzene-d4      | 1.000 | 1.000 | 0.0   | 98    | 0.00     |
| 73 T  | Isopropylbenzene            | 3.749 | 3.651 | 2.6   | 95    | 0.00     |
| 74 T  | N-amyl acetate              | 1.302 | 1.267 | 2.7   | 91    | 0.00     |
| 75 P  | 1,1,2,2-Tetrachloroethane   | 1.212 | 1.090 | 10.1  | 89    | 0.00     |
| 76 T  | 1,2,3-Trichloropropane      | 1.089 | 1.068 | 1.9   | 94    | 0.00     |
| 77 T  | Bromobenzene                | 0.856 | 0.871 | -1.8  | 99    | 0.00     |
| 78 T  | n-propylbenzene             | 4.222 | 4.201 | 0.5   | 95    | 0.00     |
| 79 T  | 2-Chlorotoluene             | 2.614 | 2.531 | 3.2   | 95    | 0.00     |
| 80 T  | 1,3,5-Trimethylbenzene      | 3.157 | 3.105 | 1.6   | 95    | 0.00     |
| 81 T  | trans-1,4-Dichloro-2-butene | 0.327 | 0.355 | -8.6  | 96    | 0.00     |
| 82 T  | 4-Chlorotoluene             | 2.981 | 2.911 | 2.3   | 96    | 0.00     |
| 83 T  | tert-Butylbenzene           | 3.078 | 2.981 | 3.2   | 94    | 0.00     |
| 84 T  | 1,2,4-Trimethylbenzene      | 3.149 | 3.098 | 1.6   | 95    | 0.00     |
| 85 T  | sec-Butylbenzene            | 3.815 | 3.824 | -0.2  | 95    | 0.00     |
| 86 T  | p-Isopropyltoluene          | 3.194 | 3.226 | -1.0  | 96    | 0.00     |
| 87 T  | 1,3-Dichlorobenzene         | 1.630 | 1.668 | -2.3  | 98    | 0.00     |
| 88 T  | 1,4-Dichlorobenzene         | 1.627 | 1.649 | -1.4  | 97    | 0.00     |
| 89 T  | n-Butylbenzene              | 2.696 | 2.821 | -4.6  | 98    | 0.00     |
| 90 T  | Hexachloroethane            | 0.457 | 0.497 | -8.8  | 97    | 0.00     |
| 91 T  | 1,2-Dichlorobenzene         | 1.602 | 1.627 | -1.6  | 98    | 0.00     |
| 92 T  | 1,2-Dibromo-3-Chloropropane | 0.269 | 0.249 | 7.4   | 90    | 0.00     |
| 93 T  | 1,2,4-Trichlorobenzene      | 0.977 | 1.055 | -8.0  | 101   | 0.00     |
| 94 T  | Hexachlorobutadiene         | 0.400 | 0.435 | -8.7  | 102   | 0.00     |
| 95 T  | Naphthalene                 | 3.507 | 3.780 | -7.8  | 100   | 0.00     |

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LabSampleId :  
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Quant Title : SW846 8260  
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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 | 0.978 | 1.033 | -5.6 | 100   | 0.00     |

(#) = Out of Range

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