

Data Path : Z:\voasrv\HPCHEM1\MSVOA\_X\Data\VX080624\  
 Data File : VX042875.D  
 Acq On : 06 Aug 2024 16:43  
 Operator : JC/MD  
 Sample : VSTDCCC020  
 Misc : 5.0mL/MSVOA\_X/WATER  
 ALS Vial : 18 Sample Multiplier: 1

Instrument :  
 MSVOA\_X  
 LabSampleId :  
 VSTDCCC020

Quant Time: Aug 07 05:01:34 2024  
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_X\Method\624X080624W.M  
 Quant Title : METHOD 624 VOLATILE ORGANIC ANALYSIS  
 QLast Update : Wed Aug 07 04:53:59 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  | Bromochloromethane          | 1.000 | 1.000 | 0.0     | 87    | 0.00     |
| 2 M  | Dichlorodifluoromethane     | 1.764 | 1.476 | 16.3    | 74    | 0.00     |
| 3 M  | Chloromethane               | 1.619 | 1.392 | 14.0    | 84    | 0.00     |
| 4 M  | Vinyl Chloride              | 1.729 | 1.441 | 16.7    | 78    | 0.00     |
| 5 M  | Bromomethane                | 1.054 | 0.806 | 23.5    | 70    | 0.00     |
| 6 M  | Chloroethane                | 1.028 | 0.871 | 15.3    | 76    | -0.02    |
| 7 M  | Trichlorofluoromethane      | 2.757 | 2.147 | 22.1    | 64    | -0.01    |
| 8 T  | Diethyl Ether               | 1.083 | 0.949 | 12.4    | 79    | 0.00     |
| 9    | 1,1,2-Trichlorotrifluoroeth | 1.879 | 1.590 | 15.4    | 76    | -0.01    |
| 10 M | 1,1-Dichloroethene          | 1.795 | 1.430 | 20.3    | 73    | -0.01    |
| 11   | Methyl Iodide               | 1.652 | 1.399 | 15.3    | 88    | -0.01    |
| 12   | Methyl Acetate              | 3.924 | 4.145 | -5.6    | 84    | 0.00     |
| 13 M | Acrolein                    | 0.113 | 0.093 | 17.7    | 104   | 0.00     |
| 14 M | Acrylonitrile               | 1.193 | 1.100 | 7.8     | 83    | 0.00     |
| 15 M | Acetone                     | 0.387 | 0.278 | 28.2#   | 66    | 0.01     |
| 16 M | Carbon Disulfide            | 3.736 | 2.759 | 26.2#   | 68    | -0.01    |
| 17   | Allyl chloride              | 2.905 | 2.470 | 15.0    | 71    | 0.00     |
| 18 M | Methylene Chloride          | 2.082 | 1.815 | 12.8    | 81    | -0.01    |
| 19 M | trans-1,2-Dichloroethene    | 1.828 | 1.526 | 16.5    | 74    | 0.00     |
| 20 T | Diisopropyl ether           | 6.346 | 5.643 | 11.1    | 78    | 0.00     |
| 21 M | 1,1-Dichloroethane          | 3.609 | 3.203 | 11.2    | 78    | 0.00     |
| 22 M | cis-1,2-Dichloroethene      | 2.294 | 1.976 | 13.9    | 79    | 0.00     |
| 23 M | tert-Butyl Alcohol          | 0.550 | 0.436 | 20.7    | 73    | 0.02     |
| 24 M | Methyl tert-Butyl Ether     | 6.498 | 5.762 | 11.3    | 81    | 0.00     |
| 25 M | Chloroform                  | 3.905 | 3.379 | 13.5    | 78    | 0.00     |
| 26   | Cyclohexane                 | 2.847 | 2.276 | 20.1    | 75    | 0.00     |
| 27 s | 1,2-Dichloroethane-d4       | 2.770 | 2.545 | 8.1     | 81    | 0.00     |
| 28 I | 1,4-Difluorobenzene         | 1.000 | 1.000 | 0.0     | 79    | 0.00     |
| 29   | 1,1-Dichloropropene         | 0.387 | 0.342 | 11.6    | 72    | 0.00     |
| 30 M | 2-Butanone                  | 0.285 | 0.272 | 4.6     | 77    | 0.01     |
| 31   | 2,2-Dichloropropane         | 0.452 | 0.348 | 23.0    | 62    | 0.00     |
| 32 M | 1,1,1-Trichloroethane       | 0.535 | 0.505 | 5.6     | 77    | 0.00     |
| 33 M | Carbon Tetrachloride        | 0.458 | 0.420 | 8.3     | 74    | 0.00     |
| 34 M | Benzene                     | 1.238 | 1.199 | 3.2     | 79    | 0.00     |
| 35   | Methacrylonitrile           | 0.269 | 0.267 | 0.7     | 79    | 0.01     |
| 36 M | 1,2-Dichloroethane          | 0.457 | 0.440 | 3.7     | 77    | 0.00     |
| 37 M | Trichloroethene             | 0.313 | 0.278 | 11.2    | 72    | 0.00     |
| 38   | Methylcyclohexane           | 0.479 | 0.395 | 17.5    | 69    | 0.00     |
| 39 M | 1,2-Dichloropropane         | 0.316 | 0.300 | 5.1     | 77    | 0.00     |
| 40   | Dibromomethane              | 0.238 | 0.222 | 6.7     | 75    | 0.00     |
| 41 M | Bromodichloromethane        | 0.472 | 0.432 | 8.5     | 74    | 0.00     |
| 42 M | Vinyl Acetate               | 0.833 | 0.778 | 6.6     | 77    | 0.00     |
| 43   | Ethyl Acetate               | 0.478 | 1.875 | -292.3# | 321#  | -0.15    |
| 44   | Isopropyl Acetate           | 0.769 | 0.742 | 3.5     | 77    | 0.00     |
| 45 T | 1,4-Dioxane                 | 0.008 | 0.008 | 0.0     | 76    | 0.00     |
| 46   | Methyl methacrylate         | 0.382 | 0.353 | 7.6     | 74    | 0.00     |
| 47   | n-amyl Acetate              | 0.658 | 0.590 | 10.3    | 76    | 0.00     |
| 48 M | t-1,3-Dichloropropene       | 0.453 | 0.380 | 16.1    | 67    | 0.00     |

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 Misc : 5.0mL/MSVOA\_X/WATER  
 ALS Vial : 18 Sample Multiplier: 1

Instrument :  
 MSVOA\_X  
 LabSampleId :  
 VSTDCCC020

Quant Time: Aug 07 05:01:34 2024  
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_X\Method\624X080624W.M  
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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 | cis-1,3-Dichloropropene     | 0.490 | 0.436 | 11.0  | 71    | 0.00     |
| 50 M | 1,1,2-Trichloroethane       | 0.325 | 0.324 | 0.3   | 79    | 0.00     |
| 51   | Ethyl methacrylate          | 0.532 | 0.490 | 7.9   | 76    | 0.00     |
| 52   | 1,3-Dichloropropane         | 0.533 | 0.505 | 5.3   | 75    | 0.00     |
| 53 M | Dibromochloromethane        | 0.378 | 0.333 | 11.9  | 72    | 0.00     |
| 54 M | 1,2-Dibromoethane           | 0.338 | 0.326 | 3.6   | 77    | 0.00     |
| 55 M | 2-Chloroethyl vinyl ether   | 0.246 | 0.252 | -2.4  | 83    | 0.00     |
| 56 M | Bromoform                   | 0.271 | 0.228 | 15.9  | 69    | 0.00     |
| 57 I | Chlorobenzene-d5            | 1.000 | 1.000 | 0.0   | 81    | 0.00     |
| 58 M | 4-Methyl-2-Pentanone        | 0.565 | 0.567 | -0.4  | 80    | 0.00     |
| 59 M | 2-Hexanone                  | 0.447 | 0.435 | 2.7   | 79    | 0.00     |
| 60 S | 4-Bromofluorobenzene        | 0.489 | 0.469 | 4.1   | 78    | 0.00     |
| 61 M | Tetrachloroethene           | 0.298 | 0.278 | 6.7   | 76    | 0.00     |
| 62 M | Toluene                     | 1.469 | 1.385 | 5.7   | 77    | 0.00     |
| 63 S | Toluene-d8                  | 1.355 | 1.399 | -3.2  | 82    | 0.00     |
| 64 M | Chlorobenzene               | 0.940 | 0.874 | 7.0   | 76    | 0.00     |
| 65   | 1,1,1,2-Tetrachloroethane   | 0.341 | 0.308 | 9.7   | 74    | 0.00     |
| 66 M | Ethyl Benzene               | 1.633 | 1.471 | 9.9   | 74    | 0.00     |
| 67 M | m/p-Xylenes                 | 0.617 | 0.556 | 9.9   | 74    | 0.00     |
| 68 M | o-Xylene                    | 0.609 | 0.556 | 8.7   | 74    | 0.00     |
| 69 M | Styrene                     | 1.048 | 0.943 | 10.0  | 74    | 0.00     |
| 70   | Isopropylbenzene            | 1.602 | 1.457 | 9.1   | 74    | 0.00     |
| 71 M | 1,1,2,2-Tetrachloroethane   | 0.578 | 0.551 | 4.7   | 78    | 0.00     |
| 72   | 1,2,3-Trichloropropane      | 0.472 | 0.448 | 5.1   | 80    | 0.00     |
| 73   | Bromobenzene                | 0.407 | 0.361 | 11.3  | 74    | 0.00     |
| 74   | n-propylbenzene             | 1.874 | 1.622 | 13.4  | 71    | 0.00     |
| 75   | 2-Chlorotoluene             | 1.144 | 1.030 | 10.0  | 75    | 0.00     |
| 76   | 1,3,5-Trimethylbenzene      | 1.358 | 1.229 | 9.5   | 75    | 0.00     |
| 77   | t-1,4-Dichloro-2-butene     | 0.168 | 0.123 | 26.8# | 64    | 0.00     |
| 78   | 4-Chlorotoluene             | 1.312 | 1.130 | 13.9  | 72    | 0.00     |
| 79   | tert-butylbenzene           | 1.393 | 1.245 | 10.6  | 75    | 0.00     |
| 80   | 1,2,4-Trimethylbenzene      | 1.340 | 1.201 | 10.4  | 75    | 0.00     |
| 81   | sec-Butylbenzene            | 1.708 | 1.492 | 12.6  | 72    | 0.00     |
| 82   | p-Isopropyltoluene          | 1.430 | 1.207 | 15.6  | 70    | 0.00     |
| 83 M | 1,3-Dichlorobenzene         | 0.749 | 0.644 | 14.0  | 72    | 0.00     |
| 84 M | 1,4-Dichlorobenzene         | 0.753 | 0.638 | 15.3  | 70    | 0.00     |
| 85   | n-Butylbenzene              | 1.289 | 0.984 | 23.7  | 63    | 0.00     |
| 86 T | Hexachloroethane            | 0.258 | 0.210 | 18.6  | 70    | 0.00     |
| 87 M | 1,2-Dichlorobenzene         | 0.746 | 0.656 | 12.1  | 74    | 0.00     |
| 88   | 1,2-Dibromo-3-Chloropropane | 0.131 | 0.119 | 9.2   | 79    | 0.00     |
| 89   | 1,2,4-Trichlorobenzene      | 0.486 | 0.373 | 23.3  | 65    | 0.00     |
| 90   | Hexachlorobutadiene         | 0.210 | 0.172 | 18.1  | 70    | 0.00     |
| 91 M | Naphthalene                 | 1.635 | 1.401 | 14.3  | 73    | 0.00     |
| 92   | 1,2,3-Trichlorobenzene      | 0.490 | 0.399 | 18.6  | 68    | 0.00     |

(#) = Out of Range

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