

Data Path : Z:\voasrv\HPCHEM1\MSVOA\_X\Data\VX093022\  
 Data File : VX031822.D  
 Acq On : 30 Sep 2022 19:26  
 Operator : JC/MD  
 Sample : VSTDCCC020  
 Misc : 5.0mL/MSVOA\_X/WATER  
 ALS Vial : 19 Sample Multiplier: 1

Instrument :  
 MSVOA\_X  
 LabSampleId :  
 VSTDCCC020

Quant Time: Oct 03 01:08:03 2022  
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_X\Method\624X093022W.M  
 Quant Title : METHOD 624 VOLATILE ORGANIC ANALYSIS  
 QLast Update : Mon Oct 03 00:56:55 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  | Bromochloromethane          | 1.000 | 1.000 | 0.0  | 97    | 0.00     |
| 2 M  | Dichlorodifluoromethane     | 1.251 | 1.189 | 5.0  | 90    | 0.00     |
| 3 M  | Chloromethane               | 1.311 | 1.390 | -6.0 | 94    | 0.00     |
| 4 M  | Vinyl Chloride              | 1.758 | 1.729 | 1.6  | 88    | 0.00     |
| 5 M  | Bromomethane                | 1.609 | 1.570 | 2.4  | 95    | 0.00     |
| 6 M  | Chloroethane                | 1.550 | 1.410 | 9.0  | 86    | 0.00     |
| 7 M  | Trichlorofluoromethane      | 3.758 | 3.748 | 0.3  | 89    | 0.00     |
| 8 T  | Diethyl Ether               | 1.273 | 1.267 | 0.5  | 87    | 0.00     |
| 9    | 1,1,2-Trichlorotrifluoroeth | 1.783 | 1.774 | 0.5  | 91    | 0.00     |
| 10 M | 1,1-Dichloroethene          | 1.647 | 1.598 | 3.0  | 90    | 0.00     |
| 11   | Methyl Iodide               | 1.803 | 1.546 | 14.3 | 95    | 0.00     |
| 12   | Methyl Acetate              | 1.882 | 1.937 | -2.9 | 93    | 0.00     |
| 13 M | Acrolein                    | 0.167 | 0.151 | 9.6  | 93    | 0.00     |
| 14 M | Acrylonitrile               | 0.799 | 0.808 | -1.1 | 91    | 0.00     |
| 15 M | Acetone                     | 0.216 | 0.220 | -1.9 | 88    | 0.00     |
| 16 M | Carbon Disulfide            | 3.613 | 3.227 | 10.7 | 88    | 0.00     |
| 17   | Allyl chloride              | 1.994 | 1.902 | 4.6  | 87    | 0.00     |
| 18 M | Methylene Chloride          | 1.885 | 1.882 | 0.2  | 91    | 0.00     |
| 19 M | trans-1,2-Dichloroethene    | 1.907 | 1.765 | 7.4  | 86    | 0.00     |
| 20 T | Diisopropyl ether           | 4.582 | 4.564 | 0.4  | 89    | 0.00     |
| 21 M | 1,1-Dichloroethane          | 2.989 | 2.980 | 0.3  | 92    | 0.00     |
| 22 M | cis-1,2-Dichloroethene      | 2.304 | 2.245 | 2.6  | 92    | 0.00     |
| 23 M | tert-Butyl Alcohol          | 0.298 | 0.281 | 5.7  | 82    | -0.01    |
| 24 M | Methyl tert-Butyl Ether     | 5.661 | 5.424 | 4.2  | 88    | 0.00     |
| 25 M | Chloroform                  | 3.670 | 3.658 | 0.3  | 91    | 0.00     |
| 26   | Cyclohexane                 | 2.466 | 2.370 | 3.9  | 87    | 0.00     |
| 27 s | 1,2-Dichloroethane-d4       | 2.258 | 2.353 | -4.2 | 94    | 0.00     |
| 28 I | 1,4-Difluorobenzene         | 1.000 | 1.000 | 0.0  | 96    | 0.00     |
| 29   | 1,1-Dichloropropene         | 0.465 | 0.442 | 4.9  | 88    | 0.00     |
| 30 M | 2-Butanone                  | 0.189 | 0.192 | -1.6 | 88    | 0.00     |
| 31   | 2,2-Dichloropropane         | 0.466 | 0.401 | 13.9 | 83    | 0.00     |
| 32 M | 1,1,1-Trichloroethane       | 0.616 | 0.576 | 6.5  | 89    | 0.00     |
| 33 M | Carbon Tetrachloride        | 0.564 | 0.524 | 7.1  | 89    | 0.00     |
| 34 M | Benzene                     | 1.376 | 1.343 | 2.4  | 90    | 0.00     |
| 35   | Methacrylonitrile           | 0.198 | 0.197 | 0.5  | 90    | 0.00     |
| 36 M | 1,2-Dichloroethane          | 0.485 | 0.486 | -0.2 | 90    | 0.00     |
| 37 M | Trichloroethene             | 0.449 | 0.416 | 7.3  | 91    | 0.00     |
| 38   | Methylcyclohexane           | 0.569 | 0.538 | 5.4  | 89    | 0.00     |
| 39 M | 1,2-Dichloropropane         | 0.315 | 0.318 | -1.0 | 94    | 0.00     |
| 40   | Dibromomethane              | 0.275 | 0.270 | 1.8  | 89    | 0.00     |
| 41 M | Bromodichloromethane        | 0.499 | 0.482 | 3.4  | 92    | 0.00     |
| 42 M | Vinyl Acetate               | 0.683 | 0.677 | 0.9  | 89    | 0.00     |
| 43   | Ethyl Acetate               | 0.380 | 0.401 | -5.5 | 93    | 0.00     |
| 44   | Isopropyl Acetate           | 0.605 | 0.591 | 2.3  | 89    | 0.00     |
| 45 T | 1,4-Dioxane                 | 0.009 | 0.009 | 0.0  | 84    | 0.00     |
| 46   | Methyl methacrylate         | 0.282 | 0.290 | -2.8 | 91    | 0.00     |
| 47   | n-amyl Acetate              | 0.519 | 0.496 | 4.4  | 88    | 0.00     |
| 48 M | t-1,3-Dichloropropene       | 0.492 | 0.429 | 12.8 | 85    | 0.00     |

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 ALS Vial : 19 Sample Multiplier: 1

Instrument :  
 MSVOA\_X  
 LabSampleID :  
 VSTDCCC020

Quant Time: Oct 03 01:08:03 2022  
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_X\Method\624X093022W.M  
 Quant Title : METHOD 624 VOLATILE ORGANIC ANALYSIS  
 QLast Update : Mon Oct 03 00:56:55 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 | cis-1,3-Dichloropropene     | 0.540 | 0.485 | 10.2 | 86    | 0.00     |
| 50 M | 1,1,2-Trichloroethane       | 0.398 | 0.389 | 2.3  | 90    | 0.00     |
| 51   | Ethyl methacrylate          | 0.517 | 0.493 | 4.6  | 88    | 0.00     |
| 52   | 1,3-Dichloropropane         | 0.611 | 0.602 | 1.5  | 89    | 0.00     |
| 53 M | Dibromochloromethane        | 0.444 | 0.410 | 7.7  | 92    | 0.00     |
| 54 M | 1,2-Dibromoethane           | 0.441 | 0.425 | 3.6  | 90    | 0.00     |
| 55 M | 2-Chloroethyl vinyl ether   | 0.253 | 0.259 | -2.4 | 89    | 0.00     |
| 56 M | Bromoform                   | 0.325 | 0.271 | 16.6 | 87    | 0.00     |
| 57 I | Chlorobenzene-d5            | 1.000 | 1.000 | 0.0  | 94    | 0.00     |
| 58 M | 4-Methyl-2-Pentanone        | 0.403 | 0.427 | -6.0 | 90    | 0.00     |
| 59 M | 2-Hexanone                  | 0.302 | 0.319 | -5.6 | 90    | 0.00     |
| 60 S | 4-Bromofluorobenzene        | 0.555 | 0.560 | -0.9 | 92    | 0.00     |
| 61 M | Tetrachloroethene           | 0.482 | 0.442 | 8.3  | 87    | 0.00     |
| 62 M | Toluene                     | 1.718 | 1.699 | 1.1  | 90    | 0.00     |
| 63 S | Toluene-d8                  | 1.524 | 1.568 | -2.9 | 94    | 0.00     |
| 64 M | Chlorobenzene               | 1.172 | 1.131 | 3.5  | 89    | 0.00     |
| 65   | 1,1,1,2-Tetrachloroethane   | 0.444 | 0.416 | 6.3  | 90    | 0.00     |
| 66 M | Ethyl Benzene               | 1.936 | 1.911 | 1.3  | 89    | 0.00     |
| 67 M | m/p-Xylenes                 | 0.806 | 0.791 | 1.9  | 88    | 0.00     |
| 68 M | o-Xylene                    | 0.791 | 0.770 | 2.7  | 88    | 0.00     |
| 69 M | Styrene                     | 1.285 | 1.241 | 3.4  | 87    | 0.00     |
| 70   | Isopropylbenzene            | 2.057 | 2.038 | 0.9  | 89    | 0.00     |
| 71 M | 1,1,2,2-Tetrachloroethane   | 0.586 | 0.601 | -2.6 | 91    | 0.00     |
| 72   | 1,2,3-Trichloropropane      | 0.508 | 0.522 | -2.8 | 89    | 0.00     |
| 73   | Bromobenzene                | 0.552 | 0.523 | 5.3  | 88    | 0.00     |
| 74   | n-propylbenzene             | 2.261 | 2.236 | 1.1  | 88    | 0.00     |
| 75   | 2-Chlorotoluene             | 1.359 | 1.360 | -0.1 | 89    | 0.00     |
| 76   | 1,3,5-Trimethylbenzene      | 1.734 | 1.728 | 0.3  | 88    | 0.00     |
| 77   | t-1,4-Dichloro-2-butene     | 0.147 | 0.114 | 22.4 | 81    | 0.00     |
| 78   | 4-Chlorotoluene             | 1.550 | 1.555 | -0.3 | 89    | 0.00     |
| 79   | tert-butylbenzene           | 1.743 | 1.686 | 3.3  | 87    | 0.00     |
| 80   | 1,2,4-Trimethylbenzene      | 1.719 | 1.759 | -2.3 | 91    | 0.00     |
| 81   | sec-Butylbenzene            | 2.134 | 2.082 | 2.4  | 88    | 0.00     |
| 82   | p-Isopropyltoluene          | 1.876 | 1.789 | 4.6  | 86    | 0.00     |
| 83 M | 1,3-Dichlorobenzene         | 1.044 | 0.984 | 5.7  | 88    | 0.00     |
| 84 M | 1,4-Dichlorobenzene         | 1.070 | 1.021 | 4.6  | 88    | 0.00     |
| 85   | n-Butylbenzene              | 1.446 | 1.389 | 3.9  | 89    | 0.00     |
| 86 T | Hexachloroethane            | 0.294 | 0.258 | 12.2 | 85    | 0.00     |
| 87 M | 1,2-Dichlorobenzene         | 1.041 | 1.004 | 3.6  | 89    | 0.00     |
| 88   | 1,2-Dibromo-3-Chloropropane | 0.118 | 0.117 | 0.8  | 91    | 0.00     |
| 89   | 1,2,4-Trichlorobenzene      | 0.674 | 0.595 | 11.7 | 89    | 0.00     |
| 90   | Hexachlorobutadiene         | 0.271 | 0.233 | 14.0 | 88    | 0.00     |
| 91 M | Naphthalene                 | 2.156 | 2.057 | 4.6  | 87    | 0.00     |
| 92   | 1,2,3-Trichlorobenzene      | 0.667 | 0.616 | 7.6  | 92    | 0.00     |

(#) = Out of Range

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