

Data Path : Z:\voasrv\HPCHEM1\MSVOA\_X\Data\VX022521\  
 Data File : VX020918.D  
 Acq On : 25 Feb 2021 18:04  
 Operator : JC/MD  
 Sample : VSTDCCC020  
 Misc : 5.0mL/MSVOA\_X/WATER  
 ALS Vial : 17 Sample Multiplier: 1

Instrument :  
 MSVOA\_X  
 LabSampleId :  
 VSTDCCC020

Quant Time: Feb 25 18:27:50 2021  
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_X\Method\624X022521W.M  
 Quant Title : METHOD 624 VOLATILE ORGANIC ANALYSIS  
 QLast Update : Thu Feb 25 13:26:35 2021  
 Response via : Initial Calibration

Min. RRF : 0.050 Min. Rel. Area : 20% Max. R.T. Dev 0.50min  
 Max. RRF Dev : 30% Max. Rel. Area : 150%

|      | Compound                    | AvgRF | CCRF   | %Dev  | Area% | Dev(min) |
|------|-----------------------------|-------|--------|-------|-------|----------|
| 1 I  | Bromochloromethane          | 1.000 | 1.000  | 0.0   | 99    | 0.00     |
| 2 M  | Dichlorodifluoromethane     | 1.689 | 1.688  | 0.1   | 91    | 0.00     |
| 3 M  | Chloromethane               | 1.805 | 1.817  | -0.7  | 93    | 0.00     |
| 4 M  | Vinyl Chloride              | 2.137 | 2.137  | 0.0   | 91    | 0.00     |
| 5 M  | Bromomethane                | 0.805 | 0.727  | 9.7   | 91    | 0.00     |
| 6 M  | Chloroethane                | 1.354 | 1.361  | -0.5  | 93    | -0.01    |
| 7 M  | Trichlorofluoromethane      | 2.760 | 2.892  | -4.8  | 95    | -0.01    |
| 8 T  | Diethyl Ether               | 1.245 | 1.326  | -6.5  | 97    | 0.00     |
| 9    | 1,1,2-Trichlorotrifluoroeth | 1.721 | 1.788  | -3.9  | 93    | 0.00     |
| 10 M | 1,1-Dichloroethene          | 1.774 | 1.828  | -3.0  | 93    | 0.00     |
| 11   | Methyl Iodide               | 2.198 | 2.307  | -5.0  | 106   | 0.00     |
| 12   | Methyl Acetate              | 1.961 | 2.099  | -7.0  | 99    | 0.00     |
| 13 M | Acrolein                    | 0.279 | 0.226  | 19.0  | 83    | 0.00     |
| 14 M | Acrylonitrile               | 0.978 | 1.033  | -5.6  | 98    | 0.00     |
| 15 M | Acetone                     | 0.294 | 0.289  | 1.7   | 85    | 0.00     |
| 16 M | Carbon Disulfide            | 5.007 | 5.081  | -1.5  | 91    | 0.00     |
| 17   | Allyl chloride              | 2.797 | 2.852  | -2.0  | 94    | 0.00     |
| 18 M | Methylene Chloride          | 2.004 | 2.163  | -7.9  | 98    | 0.00     |
| 19 M | trans-1,2-Dichloroethene    | 1.967 | 2.056  | -4.5  | 95    | 0.00     |
| 20 T | Diisopropyl ether           | 5.671 | 5.951  | -4.9  | 96    | 0.00     |
| 21 M | 1,1-Dichloroethane          | 3.363 | 3.533  | -5.1  | 96    | 0.00     |
| 22 M | cis-1,2-Dichloroethene      | 2.249 | 2.383  | -6.0  | 96    | 0.00     |
| 23 M | tert-Butyl Alcohol          | 0.444 | 0.461  | -3.8  | 97    | 0.01     |
| 24 M | Methyl tert-Butyl Ether     | 6.219 | 6.671  | -7.3  | 97    | 0.00     |
| 25 M | Chloroform                  | 3.444 | 3.667  | -6.5  | 97    | 0.00     |
| 26   | Cyclohexane                 | 3.045 | 3.111  | -2.2  | 92    | 0.00     |
| 27 s | 1,2-Dichloroethane-d4       | 2.113 | 2.067  | 2.2   | 96    | 0.00     |
| 28 I | 1,4-Difluorobenzene         | 1.000 | 1.000  | 0.0   | 94    | 0.00     |
| 29   | 1,1-Dichloropropene         | 0.470 | 0.502  | -6.8  | 93    | 0.00     |
| 30 M | 2-Butanone                  | 0.232 | 0.244  | -5.2  | 92    | 0.00     |
| 31   | 2,2-Dichloropropane         | 0.542 | 0.579  | -6.8  | 93    | 0.00     |
| 32 M | 1,1,1-Trichloroethane       | 0.528 | 0.572  | -8.3  | 95    | 0.00     |
| 33 M | Carbon Tetrachloride        | 0.458 | 0.492  | -7.4  | 94    | 0.00     |
| 34 M | Benzene                     | 1.413 | 1.540  | -9.0  | 95    | 0.00     |
| 35   | Methacrylonitrile           | 0.250 | 0.265  | -6.0  | 95    | 0.00     |
| 36 M | 1,2-Dichloroethane          | 0.475 | 0.530  | -11.6 | 98    | 0.00     |
| 37 M | Trichloroethene             | 0.382 | 0.417  | -9.2  | 94    | 0.00     |
| 38   | Methylcyclohexane           | 0.598 | 0.641  | -7.2  | 92    | 0.00     |
| 39 M | 1,2-Dichloropropane         | 0.349 | 0.383  | -9.7  | 98    | 0.00     |
| 40   | Dibromomethane              | 0.240 | 0.261  | -8.8  | 96    | 0.00     |
| 41 M | Bromodichloromethane        | 0.502 | 0.546  | -8.8  | 95    | 0.00     |
| 42 M | Vinyl Acetate               | 0.866 | 0.945  | -9.1  | 97    | 0.00     |
| 43   | Ethyl Acetate               | 0.442 | 0.488  | -10.4 | 100   | 0.00     |
| 44   | Isopropyl Acetate           | 0.749 | 0.813  | -8.5  | 97    | 0.00     |
| 45 T | 1,4-Dioxane                 | 0.009 | 0.010# | -11.1 | 100   | 0.00     |
| 46   | Methyl methacrylate         | 0.372 | 0.402  | -8.1  | 97    | 0.00     |
| 47   | n-amyl Acetate              | 0.616 | 0.655  | -6.3  | 95    | 0.00     |
| 48 M | t-1,3-Dichloropropene       | 0.560 | 0.600  | -7.1  | 96    | 0.00     |

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 Misc : 5.0mL/MSVOA\_X/WATER  
 ALS Vial : 17 Sample Multiplier: 1

Instrument :  
 MSVOA\_X  
 LabSampleId :  
 VSTDCCC020

Quant Time: Feb 25 18:27:50 2021  
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_X\Method\624X022521W.M  
 Quant Title : METHOD 624 VOLATILE ORGANIC ANALYSIS  
 QLast Update : Thu Feb 25 13:26:35 2021  
 Response via : Initial Calibration

Min. RRF : 0.050 Min. Rel. Area : 20% Max. R.T. Dev 0.50min  
 Max. RRF Dev : 30% Max. Rel. Area : 150%

|      | Compound                    | AvgRF | CCRF  | %Dev  | Area% | Dev(min) |
|------|-----------------------------|-------|-------|-------|-------|----------|
| 49 T | cis-1,3-Dichloropropene     | 0.611 | 0.658 | -7.7  | 95    | 0.00     |
| 50 M | 1,1,2-Trichloroethane       | 0.353 | 0.390 | -10.5 | 96    | 0.00     |
| 51   | Ethyl methacrylate          | 0.552 | 0.596 | -8.0  | 97    | 0.00     |
| 52   | 1,3-Dichloropropane         | 0.597 | 0.661 | -10.7 | 97    | 0.00     |
| 53 M | Dibromochloromethane        | 0.374 | 0.414 | -10.7 | 99    | 0.00     |
| 54 M | 1,2-Dibromoethane           | 0.367 | 0.416 | -13.4 | 102   | 0.00     |
| 55 M | 2-Chloroethyl vinyl ether   | 0.261 | 0.268 | -2.7  | 92    | 0.00     |
| 56 M | Bromoform                   | 0.264 | 0.283 | -7.2  | 93    | 0.00     |
| 57 I | Chlorobenzene-d5            | 1.000 | 1.000 | 0.0   | 96    | 0.00     |
| 58 M | 4-Methyl-2-Pentanone        | 0.491 | 0.527 | -7.3  | 98    | 0.00     |
| 59 M | 2-Hexanone                  | 0.374 | 0.391 | -4.5  | 93    | 0.00     |
| 60 S | 4-Bromofluorobenzene        | 0.496 | 0.485 | 2.2   | 93    | 0.00     |
| 61 M | Tetrachloroethene           | 0.370 | 0.389 | -5.1  | 92    | 0.00     |
| 62 M | Toluene                     | 1.723 | 1.841 | -6.8  | 95    | 0.00     |
| 63 S | Toluene-d8                  | 1.366 | 1.326 | 2.9   | 93    | 0.00     |
| 64 M | Chlorobenzene               | 1.058 | 1.083 | -2.4  | 92    | 0.00     |
| 65   | 1,1,1,2-Tetrachloroethane   | 0.387 | 0.406 | -4.9  | 96    | 0.00     |
| 66 M | Ethyl Benzene               | 1.920 | 1.992 | -3.8  | 92    | 0.00     |
| 67 M | m/p-Xylenes                 | 0.713 | 0.740 | -3.8  | 91    | 0.00     |
| 68 M | o-Xylene                    | 0.679 | 0.735 | -8.2  | 95    | 0.00     |
| 69 M | Styrene                     | 1.185 | 1.262 | -6.5  | 93    | 0.00     |
| 70   | Isopropylbenzene            | 1.829 | 1.905 | -4.2  | 92    | 0.00     |
| 71 M | 1,1,2,2-Tetrachloroethane   | 0.573 | 0.628 | -9.6  | 97    | 0.00     |
| 72   | 1,2,3-Trichloropropane      | 0.517 | 0.533 | -3.1  | 95    | 0.00     |
| 73   | Bromobenzene                | 0.428 | 0.460 | -7.5  | 96    | 0.00     |
| 74   | n-propylbenzene             | 2.096 | 2.237 | -6.7  | 93    | 0.00     |
| 75   | 2-Chlorotoluene             | 1.251 | 1.332 | -6.5  | 92    | 0.00     |
| 76   | 1,3,5-Trimethylbenzene      | 1.508 | 1.629 | -8.0  | 94    | 0.00     |
| 77   | t-1,4-Dichloro-2-butene     | 0.216 | 0.218 | -0.9  | 94    | 0.00     |
| 78   | 4-Chlorotoluene             | 1.427 | 1.511 | -5.9  | 93    | 0.00     |
| 79   | tert-butylbenzene           | 1.444 | 1.522 | -5.4  | 94    | 0.00     |
| 80   | 1,2,4-Trimethylbenzene      | 1.531 | 1.620 | -5.8  | 92    | 0.00     |
| 81   | sec-Butylbenzene            | 1.746 | 1.879 | -7.6  | 94    | 0.00     |
| 82   | p-Isopropyltoluene          | 1.580 | 1.647 | -4.2  | 92    | 0.00     |
| 83 M | 1,3-Dichlorobenzene         | 0.785 | 0.834 | -6.2  | 92    | 0.00     |
| 84 M | 1,4-Dichlorobenzene         | 0.769 | 0.812 | -5.6  | 92    | 0.00     |
| 85   | n-Butylbenzene              | 1.427 | 1.498 | -5.0  | 93    | 0.00     |
| 86 T | Hexachloroethane            | 0.298 | 0.302 | -1.3  | 89    | 0.00     |
| 87 M | 1,2-Dichlorobenzene         | 0.741 | 0.825 | -11.3 | 95    | 0.00     |
| 88   | 1,2-Dibromo-3-Chloropropane | 0.129 | 0.138 | -7.0  | 98    | 0.00     |
| 89   | 1,2,4-Trichlorobenzene      | 0.488 | 0.538 | -10.2 | 99    | 0.00     |
| 90   | Hexachlorobutadiene         | 0.200 | 0.202 | -1.0  | 89    | 0.00     |
| 91 M | Naphthalene                 | 1.710 | 1.820 | -6.4  | 95    | 0.00     |
| 92   | 1,2,3-Trichlorobenzene      | 0.489 | 0.488 | 0.2   | 85    | 0.00     |

(#) = Out of Range

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