

Data Path : Z:\voasrv\HPCHEM1\MSVOA\_X\Data\WX020222\  
 Data File : VX026751.D  
 Acq On : 02 Feb 2022 20:48  
 Operator : JC/MD  
 Sample : VSTDCCC020  
 Misc : 5.0mL/MSVOA\_X/WATER  
 ALS Vial : 28 Sample Multiplier: 1

Instrument :  
 MSVOA\_X  
 LabSampleId :  
 VSTDCCC020

Quant Time: Feb 03 03:03:36 2022  
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_X\Method\624X011022W.M  
 Quant Title : METHOD 624 VOLATILE ORGANIC ANALYSIS  
 QLast Update : Mon Jan 10 12:50:55 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                    | Amount  | Calc.   | %Dev   | Area% | Dev(min) |
|------|-----------------------------|---------|---------|--------|-------|----------|
| 1 I  | Bromochloromethane          | 30.000  | 30.000  | 0.0    | 71    | 0.00     |
| 2 M  | Dichlorodifluoromethane     | 20.000  | 20.412  | -2.1   | 61    | 0.00     |
| 3 M  | Chloromethane               | 20.000  | 20.470  | -2.3   | 63    | 0.00     |
| 4 M  | Vinyl Chloride              | 20.000  | 20.008  | -0.0   | 63    | 0.00     |
| 5 M  | Bromomethane                | 20.000  | 22.415  | -12.1  | 68    | 0.00     |
| 6 M  | Chloroethane                | 20.000  | 21.488  | -7.4   | 65    | 0.00     |
| 7 M  | Trichlorofluoromethane      | 20.000  | 19.323  | 3.4    | 60    | 0.00     |
| 8 T  | Diethyl Ether               | 20.000  | 19.246  | 3.8    | 61    | 0.00     |
| 9    | 1,1,2-Trichlorotrifluoroeth | 20.000  | 20.620  | -3.1   | 64    | 0.00     |
| 10 M | 1,1-Dichloroethene          | 20.000  | 21.119  | -5.6   | 67    | 0.00     |
| 11   | Methyl Iodide               | 20.000  | 25.561  | -27.8# | 88    | 0.00     |
| 12   | Methyl Acetate              | 20.000  | 22.711  | -13.6  | 73    | 0.00     |
| 13 M | Acrolein                    | 100.000 | 112.775 | -12.8  | 90    | 0.00     |
| 14 M | Acrylonitrile               | 100.000 | 115.353 | -15.4  | 73    | 0.00     |
| 15 M | Acetone                     | 100.000 | 97.122  | 2.9    | 57    | 0.00     |
| 16 M | Carbon Disulfide            | 20.000  | 19.039  | 4.8    | 61    | 0.00     |
| 17   | Allyl chloride              | 20.000  | 20.693  | -3.5   | 68    | 0.00     |
| 18 M | Methylene Chloride          | 20.000  | 21.749  | -8.7   | 69    | 0.00     |
| 19 M | trans-1,2-Dichloroethene    | 20.000  | 21.307  | -6.5   | 68    | 0.00     |
| 20 T | Diisopropyl ether           | 20.000  | 21.289  | -6.4   | 66    | 0.00     |
| 21 M | 1,1-Dichloroethane          | 20.000  | 20.664  | -3.3   | 66    | 0.00     |
| 22 M | cis-1,2-Dichloroethene      | 20.000  | 21.323  | -6.6   | 69    | 0.00     |
| 23 M | tert-Butyl Alcohol          | 100.000 | 111.092 | -11.1  | 77    | -0.01    |
| 24 M | Methyl tert-Butyl Ether     | 20.000  | 20.677  | -3.4   | 65    | 0.00     |
| 25 M | Chloroform                  | 20.000  | 21.210  | -6.1   | 67    | 0.00     |
| 26   | Cyclohexane                 | 20.000  | 20.811  | -4.1   | 65    | 0.00     |
| 27 s | 1,2-Dichloroethane-d4       | 30.000  | 27.214  | 9.3    | 63    | 0.00     |
| 28 I | 1,4-Difluorobenzene         | 30.000  | 30.000  | 0.0    | 70    | 0.00     |
| 29   | 1,1-Dichloropropene         | 20.000  | 20.176  | -0.9   | 65    | 0.00     |
| 30 M | 2-Butanone                  | 100.000 | 102.546 | -2.5   | 64    | 0.00     |
| 31   | 2,2-Dichloropropane         | 20.000  | 18.185  | 9.1    | 60    | 0.00     |
| 32 M | 1,1,1-Trichloroethane       | 20.000  | 20.011  | -0.1   | 64    | 0.00     |
| 33 M | Carbon Tetrachloride        | 20.000  | 19.710  | 1.4    | 63    | 0.00     |
| 34 M | Benzene                     | 20.000  | 21.553  | -7.8   | 69    | 0.00     |
| 35   | Methacrylonitrile           | 20.000  | 21.098  | -5.5   | 67    | 0.00     |
| 36 M | 1,2-Dichloroethane          | 20.000  | 19.661  | 1.7    | 63    | 0.00     |
| 37 M | Trichloroethene             | 20.000  | 22.081  | -10.4  | 71    | 0.00     |
| 38   | Methylcyclohexane           | 20.000  | 20.385  | -1.9   | 64    | 0.00     |
| 39 M | 1,2-Dichloropropane         | 20.000  | 21.845  | -9.2   | 69    | 0.00     |
| 40   | Dibromomethane              | 20.000  | 21.689  | -8.4   | 70    | 0.00     |
| 41 M | Bromodichloromethane        | 20.000  | 20.400  | -2.0   | 65    | 0.00     |
| 42 M | Vinyl Acetate               | 100.000 | 106.406 | -6.4   | 68    | 0.00     |
| 43   | Ethyl Acetate               | 20.000  | 21.929  | -9.6   | 71    | 0.00     |
| 44   | Isopropyl Acetate           | 20.000  | 21.078  | -5.4   | 68    | 0.00     |
| 45 T | 1,4-Dioxane                 | 400.000 | 473.949 | -18.5  | 76    | 0.00     |
| 46   | Methyl methacrylate         | 20.000  | 21.359  | -6.8   | 68    | 0.00     |
| 47   | n-amyl Acetate              | 20.000  | 22.106  | -10.5  | 71    | 0.00     |
| 48 M | t-1,3-Dichloropropene       | 20.000  | 19.970  | 0.2    | 65    | 0.00     |

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 Operator : JC/MD  
 Sample : VSTDCCC020  
 Misc : 5.0mL/MSVOA\_X/WATER  
 ALS Vial : 28 Sample Multiplier: 1

Instrument :  
 MSVOA\_X  
 LabSampleId :  
 VSTDCCC020

Quant Time: Feb 03 03:03:36 2022  
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_X\Method\624X011022W.M  
 Quant Title : METHOD 624 VOLATILE ORGANIC ANALYSIS  
 QLast Update : Mon Jan 10 12:50:55 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                    | Amount  | Calc.   | %Dev  | Area% | Dev(min) |
|------|-----------------------------|---------|---------|-------|-------|----------|
| 49 T | cis-1,3-Dichloropropene     | 20.000  | 20.888  | -4.4  | 68    | 0.00     |
| 50 M | 1,1,2-Trichloroethane       | 20.000  | 22.518  | -12.6 | 71    | 0.00     |
| 51   | Ethyl methacrylate          | 20.000  | 21.908  | -9.5  | 70    | 0.00     |
| 52   | 1,3-Dichloropropane         | 20.000  | 22.128  | -10.6 | 70    | 0.00     |
| 53 M | Dibromochloromethane        | 20.000  | 20.629  | -3.1  | 65    | 0.00     |
| 54 M | 1,2-Dibromoethane           | 20.000  | 21.562  | -7.8  | 69    | 0.00     |
| 55 M | 2-Chloroethyl vinyl ether   | 100.000 | 102.012 | -2.0  | 69    | 0.00     |
| 56 M | Bromoform                   | 20.000  | 20.061  | -0.3  | 66    | 0.00     |
| 57 I | Chlorobenzene-d5            | 30.000  | 30.000  | 0.0   | 72    | 0.00     |
| 58 M | 4-Methyl-2-Pentanone        | 100.000 | 113.203 | -13.2 | 71    | 0.00     |
| 59 M | 2-Hexanone                  | 100.000 | 107.510 | -7.5  | 66    | 0.00     |
| 60 S | 4-Bromofluorobenzene        | 30.000  | 28.315  | 5.6   | 68    | 0.00     |
| 61 M | Tetrachloroethene           | 20.000  | 22.882  | -14.4 | 73    | 0.00     |
| 62 M | Toluene                     | 20.000  | 21.631  | -8.2  | 70    | 0.00     |
| 63 S | Toluene-d8                  | 30.000  | 29.278  | 2.4   | 69    | 0.00     |
| 64 M | Chlorobenzene               | 20.000  | 22.052  | -10.3 | 70    | 0.00     |
| 65   | 1,1,1,2-Tetrachloroethane   | 20.000  | 21.626  | -8.1  | 69    | 0.00     |
| 66 M | Ethyl Benzene               | 20.000  | 21.026  | -5.1  | 67    | 0.00     |
| 67 M | m/p-Xylenes                 | 40.000  | 43.259  | -8.1  | 69    | 0.00     |
| 68 M | o-Xylene                    | 20.000  | 22.297  | -11.5 | 71    | 0.00     |
| 69 M | Styrene                     | 20.000  | 21.404  | -7.0  | 69    | 0.00     |
| 70   | Isopropylbenzene            | 20.000  | 21.401  | -7.0  | 68    | 0.00     |
| 71 M | 1,1,2,2-Tetrachloroethane   | 20.000  | 21.739  | -8.7  | 71    | 0.00     |
| 72   | 1,2,3-Trichloropropane      | 20.000  | 21.707  | -8.5  | 70    | 0.00     |
| 73   | Bromobenzene                | 20.000  | 22.067  | -10.3 | 71    | 0.00     |
| 74   | n-propylbenzene             | 20.000  | 20.889  | -4.4  | 67    | 0.00     |
| 75   | 2-Chlorotoluene             | 20.000  | 21.085  | -5.4  | 68    | 0.00     |
| 76   | 1,3,5-Trimethylbenzene      | 20.000  | 21.264  | -6.3  | 68    | 0.00     |
| 77   | t-1,4-Dichloro-2-butene     | 20.000  | 20.338  | -1.7  | 72    | 0.00     |
| 78   | 4-Chlorotoluene             | 20.000  | 20.577  | -2.9  | 66    | 0.00     |
| 79   | tert-butylbenzene           | 20.000  | 21.134  | -5.7  | 68    | 0.00     |
| 80   | 1,2,4-Trimethylbenzene      | 20.000  | 21.158  | -5.8  | 67    | 0.00     |
| 81   | sec-Butylbenzene            | 20.000  | 20.951  | -4.8  | 67    | 0.00     |
| 82   | p-Isopropyltoluene          | 20.000  | 20.874  | -4.4  | 67    | 0.00     |
| 83 M | 1,3-Dichlorobenzene         | 20.000  | 21.226  | -6.1  | 70    | 0.00     |
| 84 M | 1,4-Dichlorobenzene         | 20.000  | 21.627  | -8.1  | 71    | 0.00     |
| 85   | n-Butylbenzene              | 20.000  | 19.644  | 1.8   | 64    | 0.00     |
| 86 T | Hexachloroethane            | 20.000  | 19.588  | 2.1   | 65    | 0.00     |
| 87 M | 1,2-Dichlorobenzene         | 20.000  | 21.886  | -9.4  | 71    | 0.00     |
| 88   | 1,2-Dibromo-3-Chloropropane | 20.000  | 20.446  | -2.2  | 68    | 0.00     |
| 89   | 1,2,4-Trichlorobenzene      | 20.000  | 20.604  | -3.0  | 70    | 0.00     |
| 90   | Hexachlorobutadiene         | 20.000  | 20.257  | -1.3  | 66    | 0.00     |
| 91 M | Naphthalene                 | 20.000  | 21.393  | -7.0  | 71    | 0.00     |
| 92   | 1,2,3-Trichlorobenzene      | 20.000  | 21.128  | -5.6  | 70    | 0.00     |

(#) = Out of Range

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