

Data Path : Z:\voasrv\HPCHEM1\MSVOA\_X\Data\VX110421\  
 Data File : VX025062.D  
 Acq On : 03 Nov 2021 20:32  
 Operator : JC/MD  
 Sample : VSTDICV020  
 Misc : 5.0mL/MSVOA\_X/WATER  
 ALS Vial : 9 Sample Multiplier: 1

Instrument :  
 MSVOA\_X  
 ClientSampleId :  
 ICSVX110421

Quant Time: Nov 03 23:55:46 2021  
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_X\Method\624X110421W.M  
 Quant Title : METHOD 624 VOLATILE ORGANIC ANALYSIS  
 QLast Update : Wed Nov 03 23:52:56 2021  
 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   | 102   | -0.01    |
| 2 M  | Dichlorodifluoromethane     | 20.000  | 21.962  | -9.8  | 106   | 0.00     |
| 3 M  | Chloromethane               | 20.000  | 21.137  | -5.7  | 96    | 0.00     |
| 4 M  | Vinyl Chloride              | 20.000  | 21.232  | -6.2  | 104   | 0.00     |
| 5 M  | Bromomethane                | 20.000  | 20.487  | -2.4  | 92    | 0.00     |
| 6 M  | Chloroethane                | 20.000  | 22.497  | -12.5 | 100   | 0.00     |
| 7 M  | Trichlorofluoromethane      | 20.000  | 21.918  | -9.6  | 106   | 0.00     |
| 8 T  | Diethyl Ether               | 20.000  | 21.613  | -8.1  | 105   | 0.00     |
| 9    | 1,1,2-Trichlorotrifluoroeth | 20.000  | 20.986  | -4.9  | 107   | 0.00     |
| 10 M | 1,1-Dichloroethene          | 20.000  | 20.252  | -1.3  | 99    | 0.00     |
| 11   | Methyl Iodide               | 20.000  | 20.850  | -4.3  | 128   | 0.00     |
| 12   | Methyl Acetate              | 20.000  | 20.673  | -3.4  | 100   | 0.00     |
| 13 M | Acrolein                    | 100.000 | 100.951 | -1.0  | 112   | 0.00     |
| 14 M | Acrylonitrile               | 100.000 | 105.735 | -5.7  | 100   | 0.00     |
| 15 M | Acetone                     | 100.000 | 99.989  | 0.0   | 93    | 0.00     |
| 16 M | Carbon Disulfide            | 20.000  | 21.043  | -5.2  | 105   | 0.00     |
| 17   | Allyl chloride              | 20.000  | 21.019  | -5.1  | 104   | 0.00     |
| 18 M | Methylene Chloride          | 20.000  | 20.507  | -2.5  | 97    | 0.00     |
| 19 M | trans-1,2-Dichloroethene    | 20.000  | 21.633  | -8.2  | 106   | 0.00     |
| 20 T | Diisopropyl ether           | 20.000  | 21.613  | -8.1  | 103   | 0.00     |
| 21 M | 1,1-Dichloroethane          | 20.000  | 21.599  | -8.0  | 103   | 0.00     |
| 22 M | cis-1,2-Dichloroethene      | 20.000  | 21.255  | -6.3  | 104   | 0.00     |
| 23 M | tert-Butyl Alcohol          | 100.000 | 94.061  | 5.9   | 87    | 0.00     |
| 24 M | Methyl tert-Butyl Ether     | 20.000  | 21.237  | -6.2  | 101   | 0.00     |
| 25 M | Chloroform                  | 20.000  | 20.870  | -4.4  | 99    | 0.00     |
| 26   | Cyclohexane                 | 20.000  | 21.096  | -5.5  | 102   | 0.00     |
| 27 s | 1,2-Dichloroethane-d4       | 30.000  | 29.451  | 1.8   | 101   | 0.00     |
| 28 I | 1,4-Difluorobenzene         | 30.000  | 30.000  | 0.0   | 101   | 0.00     |
| 29   | 1,1-Dichloropropene         | 20.000  | 21.699  | -8.5  | 100   | 0.00     |
| 30 M | 2-Butanone                  | 100.000 | 106.169 | -6.2  | 97    | 0.00     |
| 31   | 2,2-Dichloropropane         | 20.000  | 21.454  | -7.3  | 99    | 0.00     |
| 32 M | 1,1,1-Trichloroethane       | 20.000  | 21.436  | -7.2  | 104   | 0.00     |
| 33 M | Carbon Tetrachloride        | 20.000  | 21.748  | -8.7  | 106   | 0.00     |
| 34 M | Benzene                     | 20.000  | 21.761  | -8.8  | 103   | 0.00     |
| 35   | Methacrylonitrile           | 20.000  | 21.265  | -6.3  | 101   | -0.08    |
| 36 M | 1,2-Dichloroethane          | 20.000  | 21.171  | -5.9  | 100   | 0.00     |
| 37 M | Trichloroethene             | 20.000  | 21.376  | -6.9  | 107   | 0.00     |
| 38   | Methylcyclohexane           | 20.000  | 22.409  | -12.0 | 107   | 0.00     |
| 39 M | 1,2-Dichloropropane         | 20.000  | 20.917  | -4.6  | 105   | 0.00     |
| 40   | Dibromomethane              | 20.000  | 20.907  | -4.5  | 102   | 0.00     |
| 41 M | Bromodichloromethane        | 20.000  | 20.429  | -2.1  | 100   | 0.00     |
| 42 M | Vinyl Acetate               | 100.000 | 105.204 | -5.2  | 98    | 0.00     |
| 43   | Ethyl Acetate               | 20.000  | 20.913  | -4.6  | 97    | 0.00     |
| 44   | Isopropyl Acetate           | 20.000  | 21.786  | -8.9  | 106   | 0.00     |
| 45 T | 1,4-Dioxane                 | 400.000 | 449.905 | -12.5 | 106   | -0.01    |
| 46   | Methyl methacrylate         | 20.000  | 20.212  | -1.1  | 96    | 0.00     |
| 47   | n-amyl Acetate              | 20.000  | 20.457  | -2.3  | 100   | 0.00     |
| 48 M | t-1,3-Dichloropropene       | 20.000  | 19.990  | 0.1   | 98    | 0.00     |

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 Sample : VSTDICV020  
 Misc : 5.0mL/MSVOA\_X/WATER  
 ALS Vial : 9 Sample Multiplier: 1

Instrument :  
 MSVOA\_X  
 ClientSampleId :  
 ICSVVX110421

Quant Time: Nov 03 23:55:46 2021  
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_X\Method\624X110421W.M  
 Quant Title : METHOD 624 VOLATILE ORGANIC ANALYSIS  
 QLast Update : Wed Nov 03 23:52:56 2021  
 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  | 21.436  | -7.2  | 104   | 0.00     |
| 50 M | 1,1,2-Trichloroethane       | 20.000  | 20.927  | -4.6  | 102   | 0.00     |
| 51   | Ethyl methacrylate          | 20.000  | 21.174  | -5.9  | 104   | 0.00     |
| 52   | 1,3-Dichloropropane         | 20.000  | 20.850  | -4.3  | 99    | 0.00     |
| 53 M | Dibromochloromethane        | 20.000  | 21.232  | -6.2  | 105   | 0.00     |
| 54 M | 1,2-Dibromoethane           | 20.000  | 20.838  | -4.2  | 101   | 0.00     |
| 55 M | 2-Chloroethyl vinyl ether   | 100.000 | 103.976 | -4.0  | 105   | 0.00     |
| 56 M | Bromoform                   | 20.000  | 22.063  | -10.3 | 112   | 0.00     |
| 57 I | Chlorobenzene-d5            | 30.000  | 30.000  | 0.0   | 99    | 0.00     |
| 58 M | 4-Methyl-2-Pentanone        | 100.000 | 110.655 | -10.7 | 100   | 0.00     |
| 59 M | 2-Hexanone                  | 100.000 | 110.431 | -10.4 | 100   | 0.00     |
| 60 S | 4-Bromofluorobenzene        | 30.000  | 30.841  | -2.8  | 100   | 0.00     |
| 61 M | Tetrachloroethene           | 20.000  | 22.337  | -11.7 | 107   | 0.00     |
| 62 M | Toluene                     | 20.000  | 21.774  | -8.9  | 102   | 0.00     |
| 63 S | Toluene-d8                  | 30.000  | 30.977  | -3.3  | 101   | 0.00     |
| 64 M | Chlorobenzene               | 20.000  | 21.404  | -7.0  | 100   | 0.00     |
| 65   | 1,1,1,2-Tetrachloroethane   | 20.000  | 21.620  | -8.1  | 99    | 0.00     |
| 66 M | Ethyl Benzene               | 20.000  | 21.822  | -9.1  | 101   | 0.00     |
| 67 M | m/p-Xylenes                 | 40.000  | 44.706  | -11.8 | 106   | 0.00     |
| 68 M | o-Xylene                    | 20.000  | 22.996  | -15.0 | 107   | 0.00     |
| 69 M | Styrene                     | 20.000  | 22.274  | -11.4 | 107   | 0.00     |
| 70   | Isopropylbenzene            | 20.000  | 22.164  | -10.8 | 104   | 0.00     |
| 71 M | 1,1,2,2-Tetrachloroethane   | 20.000  | 21.608  | -8.0  | 100   | 0.00     |
| 72   | 1,2,3-Trichloropropane      | 20.000  | 21.741  | -8.7  | 100   | 0.00     |
| 73   | Bromobenzene                | 20.000  | 21.437  | -7.2  | 103   | 0.00     |
| 74   | n-propylbenzene             | 20.000  | 21.653  | -8.3  | 100   | 0.00     |
| 75   | 2-Chlorotoluene             | 20.000  | 21.754  | -8.8  | 102   | 0.00     |
| 76   | 1,3,5-Trimethylbenzene      | 20.000  | 21.924  | -9.6  | 102   | 0.00     |
| 77   | t-1,4-Dichloro-2-butene     | 20.000  | 20.721  | -3.6  | 103   | 0.00     |
| 78   | 4-Chlorotoluene             | 20.000  | 21.584  | -7.9  | 100   | 0.00     |
| 79   | tert-butylbenzene           | 20.000  | 21.774  | -8.9  | 101   | 0.00     |
| 80   | 1,2,4-Trimethylbenzene      | 20.000  | 22.145  | -10.7 | 101   | 0.00     |
| 81   | sec-Butylbenzene            | 20.000  | 22.229  | -11.1 | 103   | 0.00     |
| 82   | p-Isopropyltoluene          | 20.000  | 22.037  | -10.2 | 102   | 0.00     |
| 83 M | 1,3-Dichlorobenzene         | 20.000  | 21.983  | -9.9  | 105   | 0.00     |
| 84 M | 1,4-Dichlorobenzene         | 20.000  | 21.842  | -9.2  | 105   | 0.00     |
| 85   | n-Butylbenzene              | 20.000  | 21.441  | -7.2  | 102   | 0.00     |
| 86 T | Hexachloroethane            | 20.000  | 21.570  | -7.9  | 105   | 0.00     |
| 87 M | 1,2-Dichlorobenzene         | 20.000  | 21.691  | -8.5  | 101   | 0.00     |
| 88   | 1,2-Dibromo-3-Chloropropane | 20.000  | 20.485  | -2.4  | 100   | 0.00     |
| 89   | 1,2,4-Trichlorobenzene      | 20.000  | 21.942  | -9.7  | 102   | 0.00     |
| 90   | Hexachlorobutadiene         | 20.000  | 22.038  | -10.2 | 105   | 0.00     |
| 91 M | Naphthalene                 | 20.000  | 21.381  | -6.9  | 101   | 0.00     |
| 92   | 1,2,3-Trichlorobenzene      | 20.000  | 22.087  | -10.4 | 104   | 0.00     |

(#) = Out of Range

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