

Data Path : Z:\voasrv\HPCHEM1\MSVOA\_N\Data\VN050624\  
 Data File : VN082000.D  
 Acq On : 06 May 2024 09:35  
 Operator : JC\MD  
 Sample : VSTDCCC020  
 Misc : 5.0mL/MSVOA\_N/WATER  
 ALS Vial : 1 Sample Multiplier: 1

Instrument :  
 MSVOA\_N  
 LabSampleId :  
 VSTDCCC020

Quant Time: May 07 00:24:20 2024  
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_N\methods\624N042624W.M  
 Quant Title : METHOD 624 VOLATILE ORGANIC ANALYSIS  
 QLast Update : Sat Apr 27 04:37:27 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                    | Amount  | Calc.   | %Dev  | Area% | Dev(min) |
|------|-----------------------------|---------|---------|-------|-------|----------|
| 1 I  | Bromochloromethane          | 30.000  | 30.000  | 0.0   | 84    | 0.00     |
| 2 M  | Dichlorodifluoromethane     | 20.000  | 18.509  | 7.5   | 79    | 0.00     |
| 3 M  | Chloromethane               | 20.000  | 18.702  | 6.5   | 81    | 0.00     |
| 4 M  | Vinyl Chloride              | 20.000  | 19.147  | 4.3   | 83    | 0.00     |
| 5 M  | Bromomethane                | 20.000  | 18.349  | 8.3   | 77    | 0.00     |
| 6 M  | Chloroethane                | 20.000  | 20.921  | -4.6  | 87    | 0.00     |
| 7 M  | Trichlorofluoromethane      | 20.000  | 19.269  | 3.7   | 81    | 0.00     |
| 8 T  | Diethyl Ether               | 20.000  | 18.831  | 5.8   | 78    | 0.00     |
| 9    | 1,1,2-Trichlorotrifluoroeth | 20.000  | 18.230  | 8.8   | 76    | 0.02     |
| 10 M | 1,1-Dichloroethene          | 20.000  | 19.220  | 3.9   | 83    | 0.00     |
| 11   | Methyl Iodide               | 20.000  | 17.590  | 12.1  | 85    | 0.00     |
| 12   | Methyl Acetate              | 20.000  | 21.424  | -7.1  | 92    | 0.00     |
| 13 M | Acrolein                    | 100.000 | 83.954  | 16.0  | 74    | 0.00     |
| 14 M | Acrylonitrile               | 100.000 | 103.461 | -3.5  | 87    | 0.00     |
| 15 M | Acetone                     | 100.000 | 111.215 | -11.2 | 91    | 0.00     |
| 16 M | Carbon Disulfide            | 20.000  | 18.272  | 8.6   | 80    | 0.00     |
| 17   | Allyl chloride              | 20.000  | 19.922  | 0.4   | 81    | 0.00     |
| 18 M | Methylene Chloride          | 20.000  | 19.836  | 0.8   | 87    | -0.01    |
| 19 M | trans-1,2-Dichloroethene    | 20.000  | 19.145  | 4.3   | 85    | -0.01    |
| 20 T | Diisopropyl ether           | 20.000  | 19.376  | 3.1   | 82    | 0.00     |
| 21 M | 1,1-Dichloroethane          | 20.000  | 20.101  | -0.5  | 84    | 0.00     |
| 22 M | cis-1,2-Dichloroethene      | 20.000  | 20.145  | -0.7  | 87    | 0.00     |
| 23 M | tert-Butyl Alcohol          | 100.000 | 96.919  | 3.1   | 78    | 0.02     |
| 24 M | Methyl tert-Butyl Ether     | 20.000  | 19.129  | 4.4   | 79    | 0.00     |
| 25 M | Chloroform                  | 20.000  | 19.917  | 0.4   | 84    | 0.00     |
| 26   | Cyclohexane                 | 20.000  | 19.559  | 2.2   | 85    | 0.00     |
| 27 s | 1,2-Dichloroethane-d4       | 30.000  | 31.290  | -4.3  | 89    | 0.00     |
| 28 I | 1,4-Difluorobenzene         | 30.000  | 30.000  | 0.0   | 87    | 0.00     |
| 29   | 1,1-Dichloropropene         | 20.000  | 18.012  | 9.9   | 79    | 0.00     |
| 30 M | 2-Butanone                  | 100.000 | 99.053  | 0.9   | 85    | 0.00     |
| 31   | 2,2-Dichloropropane         | 20.000  | 22.705  | -13.5 | 102   | 0.00     |
| 32 M | 1,1,1-Trichloroethane       | 20.000  | 19.302  | 3.5   | 84    | 0.00     |
| 33 M | Carbon Tetrachloride        | 20.000  | 18.396  | 8.0   | 80    | 0.00     |
| 34 M | Benzene                     | 20.000  | 19.400  | 3.0   | 84    | 0.00     |
| 35   | Methacrylonitrile           | 20.000  | 19.811  | 0.9   | 88    | 0.00     |
| 36 M | 1,2-Dichloroethane          | 20.000  | 18.479  | 7.6   | 82    | 0.00     |
| 37 M | Trichloroethene             | 20.000  | 17.948  | 10.3  | 79    | 0.00     |
| 38   | Methylcyclohexane           | 20.000  | 18.640  | 6.8   | 81    | 0.00     |
| 39 M | 1,2-Dichloropropane         | 20.000  | 19.686  | 1.6   | 89    | 0.00     |
| 40   | Dibromomethane              | 20.000  | 19.527  | 2.4   | 87    | 0.00     |
| 41 M | Bromodichloromethane        | 20.000  | 18.874  | 5.6   | 85    | 0.00     |
| 42 M | Vinyl Acetate               | 100.000 | 94.393  | 5.6   | 84    | 0.00     |
| 43   | Ethyl Acetate               | 20.000  | 19.750  | 1.3   | 88    | 0.00     |
| 44   | Isopropyl Acetate           | 20.000  | 19.179  | 4.1   | 84    | 0.00     |
| 45 T | 1,4-Dioxane                 | 400.000 | 385.683 | 3.6   | 84    | 0.00     |
| 46   | Methyl methacrylate         | 20.000  | 19.563  | 2.2   | 86    | 0.00     |
| 47   | n-amyl Acetate              | 20.000  | 18.192  | 9.0   | 81    | 0.00     |
| 48 M | t-1,3-Dichloropropene       | 20.000  | 18.614  | 6.9   | 84    | 0.00     |

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

Instrument :  
 MSVOA\_N  
 LabSampleId :  
 VSTDCCC020

Quant Time: May 07 00:24:20 2024  
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_N\methods\624N042624W.M  
 Quant Title : METHOD 624 VOLATILE ORGANIC ANALYSIS  
 QLast Update : Sat Apr 27 04:37:27 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                    | Amount  | Calc.   | %Dev | Area% | Dev(min) |
|------|-----------------------------|---------|---------|------|-------|----------|
| 49 T | cis-1,3-Dichloropropene     | 20.000  | 19.341  | 3.3  | 85    | 0.00     |
| 50 M | 1,1,2-Trichloroethane       | 20.000  | 20.201  | -1.0 | 88    | 0.00     |
| 51   | Ethyl methacrylate          | 20.000  | 17.914  | 10.4 | 79    | 0.00     |
| 52   | 1,3-Dichloropropane         | 20.000  | 19.076  | 4.6  | 84    | 0.00     |
| 53 M | Dibromochloromethane        | 20.000  | 17.952  | 10.2 | 82    | 0.00     |
| 54 M | 1,2-Dibromoethane           | 20.000  | 18.804  | 6.0  | 82    | 0.00     |
| 55 M | 2-Chloroethyl vinyl ether   | 100.000 | 104.557 | -4.6 | 91    | 0.00     |
| 56 M | Bromoform                   | 20.000  | 17.742  | 11.3 | 78    | 0.00     |
| 57 I | Chlorobenzene-d5            | 30.000  | 30.000  | 0.0  | 87    | 0.00     |
| 58 M | 4-Methyl-2-Pentanone        | 100.000 | 98.871  | 1.1  | 84    | 0.00     |
| 59 M | 2-Hexanone                  | 100.000 | 100.240 | -0.2 | 86    | 0.00     |
| 60 S | 4-Bromofluorobenzene        | 30.000  | 28.862  | 3.8  | 85    | 0.00     |
| 61 M | Tetrachloroethene           | 20.000  | 17.347  | 13.3 | 73    | 0.00     |
| 62 M | Toluene                     | 20.000  | 19.227  | 3.9  | 83    | 0.00     |
| 63 S | Toluene-d8                  | 30.000  | 30.661  | -2.2 | 88    | 0.00     |
| 64 M | Chlorobenzene               | 20.000  | 18.951  | 5.2  | 81    | 0.00     |
| 65   | 1,1,1,2-Tetrachloroethane   | 20.000  | 18.758  | 6.2  | 80    | 0.00     |
| 66 M | Ethyl Benzene               | 20.000  | 19.379  | 3.1  | 81    | 0.00     |
| 67 M | m/p-Xylenes                 | 40.000  | 38.477  | 3.8  | 82    | 0.00     |
| 68 M | o-Xylene                    | 20.000  | 19.582  | 2.1  | 82    | 0.00     |
| 69 M | Styrene                     | 20.000  | 19.048  | 4.8  | 83    | 0.00     |
| 70   | Isopropylbenzene            | 20.000  | 18.952  | 5.2  | 80    | 0.00     |
| 71 M | 1,1,2,2-Tetrachloroethane   | 20.000  | 20.229  | -1.1 | 84    | 0.00     |
| 72   | 1,2,3-Trichloropropane      | 20.000  | 19.515  | 2.4  | 85    | 0.00     |
| 73   | Bromobenzene                | 20.000  | 18.390  | 8.0  | 81    | 0.00     |
| 74   | n-propylbenzene             | 20.000  | 19.771  | 1.1  | 84    | 0.00     |
| 75   | 2-Chlorotoluene             | 20.000  | 18.321  | 8.4  | 79    | 0.00     |
| 76   | 1,3,5-Trimethylbenzene      | 20.000  | 18.967  | 5.2  | 80    | 0.00     |
| 77   | t-1,4-Dichloro-2-butene     | 20.000  | 20.728  | -3.6 | 91    | 0.00     |
| 78   | 4-Chlorotoluene             | 20.000  | 19.048  | 4.8  | 83    | 0.00     |
| 79   | tert-butylbenzene           | 20.000  | 18.807  | 6.0  | 79    | 0.00     |
| 80   | 1,2,4-Trimethylbenzene      | 20.000  | 18.819  | 5.9  | 81    | 0.00     |
| 81   | sec-Butylbenzene            | 20.000  | 19.196  | 4.0  | 83    | 0.00     |
| 82   | p-Isopropyltoluene          | 20.000  | 19.922  | 0.4  | 84    | 0.00     |
| 83 M | 1,3-Dichlorobenzene         | 20.000  | 19.584  | 2.1  | 83    | 0.00     |
| 84 M | 1,4-Dichlorobenzene         | 20.000  | 19.461  | 2.7  | 84    | 0.00     |
| 85   | n-Butylbenzene              | 20.000  | 20.207  | -1.0 | 89    | 0.00     |
| 86 T | Hexachloroethane            | 20.000  | 19.598  | 2.0  | 84    | 0.00     |
| 87 M | 1,2-Dichlorobenzene         | 20.000  | 18.788  | 6.1  | 81    | 0.00     |
| 88   | 1,2-Dibromo-3-Chloropropane | 20.000  | 19.006  | 5.0  | 78    | 0.00     |
| 89   | 1,2,4-Trichlorobenzene      | 20.000  | 18.238  | 8.8  | 80    | 0.00     |
| 90   | Hexachlorobutadiene         | 20.000  | 16.162  | 19.2 | 74    | 0.00     |
| 91 M | Naphthalene                 | 20.000  | 18.230  | 8.8  | 78    | 0.00     |
| 92   | 1,2,3-Trichlorobenzene      | 20.000  | 18.238  | 8.8  | 80    | 0.00     |

(#) = Out of Range

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