

Data Path : Z:\voasrv\HPCHEM1\MSVOA\_N\Data\VN120525\  
 Data File : VN088469.D  
 Acq On : 05 Dec 2025 14:55  
 Operator : JC\MD  
 Sample : VSTDCCC020  
 Misc : 5.0mL/MSVOA\_N/WATER  
 ALS Vial : 9 Sample Multiplier: 1

Instrument :  
 MSVOA\_N  
 LabSampleId :  
 VSTDCCC020

Quant Time: Dec 05 23:39:21 2025  
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_N\methods\624N111325W.M  
 Quant Title : METHOD 624 VOLATILE ORGANIC ANALYSIS  
 QLast Update : Fri Nov 14 00:49:21 2025  
 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   | 99    | 0.00     |
| 2 M  | Dichlorodifluoromethane     | 20.000  | 15.561  | 22.2  | 69    | 0.00     |
| 3 M  | Chloromethane               | 20.000  | 17.490  | 12.6  | 82    | 0.00     |
| 4 M  | Vinyl Chloride              | 20.000  | 16.195  | 19.0  | 77    | 0.00     |
| 5 M  | Bromomethane                | 20.000  | 13.239  | 33.8# | 66    | 0.00     |
| 6 M  | Chloroethane                | 20.000  | 18.089  | 9.6   | 85    | 0.00     |
| 7 M  | Trichlorofluoromethane      | 20.000  | 16.546  | 17.3  | 77    | 0.00     |
| 8 T  | Diethyl Ether               | 20.000  | 19.495  | 2.5   | 97    | 0.00     |
| 9    | 1,1,2-Trichlorotrifluoroeth | 20.000  | 16.443  | 17.8  | 77    | -0.01    |
| 10 M | 1,1-Dichloroethene          | 20.000  | 16.613  | 16.9  | 78    | 0.00     |
| 11   | Methyl Iodide               | 20.000  | 12.667  | 36.7# | 63    | 0.00     |
| 12   | Methyl Acetate              | 20.000  | 20.998  | -5.0  | 102   | 0.00     |
| 13 M | Acrolein                    | 100.000 | 103.391 | -3.4  | 114   | 0.00     |
| 14 M | Acrylonitrile               | 100.000 | 96.007  | 4.0   | 94    | 0.00     |
| 15 M | Acetone                     | 100.000 | 95.814  | 4.2   | 93    | 0.00     |
| 16 M | Carbon Disulfide            | 20.000  | 16.356  | 18.2  | 76    | 0.00     |
| 17   | Allyl chloride              | 20.000  | 16.688  | 16.6  | 80    | 0.00     |
| 18 M | Methylene Chloride          | 20.000  | 19.210  | 3.9   | 93    | 0.00     |
| 19 M | trans-1,2-Dichloroethene    | 20.000  | 16.623  | 16.9  | 78    | 0.00     |
| 20 T | Diisopropyl ether           | 20.000  | 20.335  | -1.7  | 99    | 0.00     |
| 21 M | 1,1-Dichloroethane          | 20.000  | 18.816  | 5.9   | 90    | 0.00     |
| 22 M | cis-1,2-Dichloroethene      | 20.000  | 18.504  | 7.5   | 90    | 0.00     |
| 23 M | tert-Butyl Alcohol          | 100.000 | 80.577  | 19.4  | 82    | -0.02    |
| 24 M | Methyl tert-Butyl Ether     | 20.000  | 19.581  | 2.1   | 97    | 0.00     |
| 25 M | Chloroform                  | 20.000  | 19.615  | 1.9   | 96    | 0.00     |
| 26   | Cyclohexane                 | 20.000  | 13.588  | 32.1# | 62    | -0.01    |
| 27 s | 1,2-Dichloroethane-d4       | 30.000  | 32.064  | -6.9  | 106   | 0.00     |
| 28 I | 1,4-Difluorobenzene         | 30.000  | 30.000  | 0.0   | 85    | 0.00     |
| 29   | 1,1-Dichloropropene         | 20.000  | 19.072  | 4.6   | 80    | 0.00     |
| 30 M | 2-Butanone                  | 100.000 | 105.023 | -5.0  | 93    | -0.01    |
| 31   | 2,2-Dichloropropane         | 20.000  | 19.780  | 1.1   | 84    | 0.00     |
| 32 M | 1,1,1-Trichloroethane       | 20.000  | 20.073  | -0.4  | 85    | 0.00     |
| 33 M | Carbon Tetrachloride        | 20.000  | 18.817  | 5.9   | 80    | 0.00     |
| 34 M | Benzene                     | 20.000  | 20.705  | -3.5  | 88    | 0.00     |
| 35   | Methacrylonitrile           | 20.000  | 21.354  | -6.8  | 95    | 0.00     |
| 36 M | 1,2-Dichloroethane          | 20.000  | 23.358  | -16.8 | 101   | 0.00     |
| 37 M | Trichloroethene             | 20.000  | 19.642  | 1.8   | 84    | 0.00     |
| 38   | Methylcyclohexane           | 20.000  | 14.015  | 29.9# | 58    | 0.00     |
| 39 M | 1,2-Dichloropropane         | 20.000  | 22.065  | -10.3 | 96    | 0.00     |
| 40   | Dibromomethane              | 20.000  | 22.670  | -13.4 | 97    | 0.00     |
| 41 M | Bromodichloromethane        | 20.000  | 23.106  | -15.5 | 103   | 0.00     |
| 42 M | Vinyl Acetate               | 100.000 | 118.589 | -18.6 | 104   | 0.00     |
| 43   | Ethyl Acetate               | 20.000  | 21.606  | -8.0  | 93    | 0.00     |
| 44   | Isopropyl Acetate           | 20.000  | 21.928  | -9.6  | 96    | 0.00     |
| 45 T | 1,4-Dioxane                 | 400.000 | 398.484 | 0.4   | 82    | 0.00     |
| 46   | Methyl methacrylate         | 20.000  | 19.999  | 0.0   | 90    | 0.00     |
| 47   | n-amyl Acetate              | 20.000  | 22.325  | -11.6 | 106   | -0.12    |
| 48 M | t-1,3-Dichloropropene       | 20.000  | 20.091  | -0.5  | 92    | 0.00     |

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

Instrument :  
 MSVOA\_N  
 LabSampleId :  
 VSTDCCC020

Quant Time: Dec 05 23:39:21 2025  
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_N\methods\624N111325W.M  
 Quant Title : METHOD 624 VOLATILE ORGANIC ANALYSIS  
 QLast Update : Fri Nov 14 00:49:21 2025  
 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  | 21.211  | -6.1  | 92    | 0.00     |
| 50 M | 1,1,2-Trichloroethane       | 20.000  | 23.102  | -15.5 | 101   | 0.00     |
| 51   | Ethyl methacrylate          | 20.000  | 20.954  | -4.8  | 97    | 0.00     |
| 52   | 1,3-Dichloropropane         | 20.000  | 22.266  | -11.3 | 97    | 0.00     |
| 53 M | Dibromochloromethane        | 20.000  | 22.041  | -10.2 | 98    | 0.00     |
| 54 M | 1,2-Dibromoethane           | 20.000  | 21.326  | -6.6  | 95    | 0.00     |
| 55 M | 2-Chloroethyl vinyl ether   | 100.000 | 97.998  | 2.0   | 89    | 0.00     |
| 56 M | Bromoform                   | 20.000  | 21.050  | -5.3  | 96    | -0.01    |
| 57 I | Chlorobenzene-d5            | 30.000  | 30.000  | 0.0   | 92    | 0.00     |
| 58 M | 4-Methyl-2-Pentanone        | 100.000 | 108.787 | -8.8  | 99    | 0.00     |
| 59 M | 2-Hexanone                  | 100.000 | 104.116 | -4.1  | 99    | 0.00     |
| 60 S | 4-Bromofluorobenzene        | 30.000  | 30.133  | -0.4  | 94    | 0.00     |
| 61 M | Tetrachloroethene           | 20.000  | 18.774  | 6.1   | 81    | 0.00     |
| 62 M | Toluene                     | 20.000  | 20.044  | -0.2  | 89    | 0.00     |
| 63 S | Toluene-d8                  | 30.000  | 30.378  | -1.3  | 91    | 0.00     |
| 64 M | Chlorobenzene               | 20.000  | 19.648  | 1.8   | 90    | 0.00     |
| 65   | 1,1,1,2-Tetrachloroethane   | 20.000  | 21.055  | -5.3  | 95    | 0.00     |
| 66 M | Ethyl Benzene               | 20.000  | 18.197  | 9.0   | 82    | 0.00     |
| 67 M | m/p-Xylenes                 | 40.000  | 37.761  | 5.6   | 85    | 0.00     |
| 68 M | o-Xylene                    | 20.000  | 18.420  | 7.9   | 85    | 0.00     |
| 69 M | Styrene                     | 20.000  | 18.783  | 6.1   | 88    | 0.00     |
| 70   | Isopropylbenzene            | 20.000  | 17.212  | 13.9  | 78    | 0.00     |
| 71 M | 1,1,2,2-Tetrachloroethane   | 20.000  | 20.533  | -2.7  | 94    | 0.00     |
| 72   | 1,2,3-Trichloropropane      | 20.000  | 20.049  | -0.2  | 100   | 0.00     |
| 73   | Bromobenzene                | 20.000  | 20.129  | -0.6  | 93    | 0.00     |
| 74   | n-propylbenzene             | 20.000  | 17.439  | 12.8  | 78    | 0.00     |
| 75   | 2-Chlorotoluene             | 20.000  | 18.025  | 9.9   | 82    | 0.00     |
| 76   | 1,3,5-Trimethylbenzene      | 20.000  | 17.463  | 12.7  | 79    | 0.00     |
| 77   | t-1,4-Dichloro-2-butene     | 20.000  | 18.166  | 9.2   | 89    | 0.00     |
| 78   | 4-Chlorotoluene             | 20.000  | 18.722  | 6.4   | 85    | 0.00     |
| 79   | tert-butylbenzene           | 20.000  | 16.143  | 19.3  | 74    | 0.00     |
| 80   | 1,2,4-Trimethylbenzene      | 20.000  | 17.778  | 11.1  | 82    | 0.00     |
| 81   | sec-Butylbenzene            | 20.000  | 15.892  | 20.5  | 71    | 0.00     |
| 82   | p-Isopropyltoluene          | 20.000  | 16.032  | 19.8  | 72    | 0.00     |
| 83 M | 1,3-Dichlorobenzene         | 20.000  | 19.295  | 3.5   | 87    | 0.00     |
| 84 M | 1,4-Dichlorobenzene         | 20.000  | 18.835  | 5.8   | 87    | 0.00     |
| 85   | n-Butylbenzene              | 20.000  | 15.123  | 24.4  | 66    | 0.00     |
| 86 T | Hexachloroethane            | 20.000  | 15.037  | 24.8  | 69    | 0.00     |
| 87 M | 1,2-Dichlorobenzene         | 20.000  | 18.806  | 6.0   | 85    | 0.00     |
| 88   | 1,2-Dibromo-3-Chloropropane | 20.000  | 18.673  | 6.6   | 90    | 0.00     |
| 89   | 1,2,4-Trichlorobenzene      | 20.000  | 17.255  | 13.7  | 81    | 0.00     |
| 90   | Hexachlorobutadiene         | 20.000  | 15.075  | 24.6  | 67    | 0.00     |
| 91 M | Naphthalene                 | 20.000  | 16.901  | 15.5  | 80    | 0.00     |
| 92   | 1,2,3-Trichlorobenzene      | 20.000  | 17.255  | 13.7  | 81    | 0.00     |

(#) = Out of Range

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