

Data Path : Z:\voasrv\HPCHEM1\MSVOA\_N\Data\VN091725\  
 Data File : VN087883.D  
 Acq On : 17 Sep 2025 16:21  
 Operator : JC\MD  
 Sample : VSTDCCC020  
 Misc : 5.0mL/MSVOA\_N/WATER  
 ALS Vial : 18 Sample Multiplier: 1

Instrument :  
 MSVOA\_N  
 LabSampleId :  
 VSTDCCC020

Quant Time: Sep 18 03:14:20 2025  
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_N\methods\624N090425W.M  
 Quant Title : METHOD 624 VOLATILE ORGANIC ANALYSIS  
 QLast Update : Fri Sep 05 03:29:53 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   | 116   | 0.00     |
| 2 M  | Dichlorodifluoromethane     | 20.000  | 16.729  | 16.4  | 104   | 0.00     |
| 3 M  | Chloromethane               | 20.000  | 15.450  | 22.8  | 98    | 0.00     |
| 4 M  | Vinyl Chloride              | 20.000  | 14.544  | 27.3# | 94    | 0.00     |
| 5 M  | Bromomethane                | 20.000  | 14.522  | 27.4# | 99    | 0.00     |
| 6 M  | Chloroethane                | 20.000  | 14.544  | 27.3# | 94    | 0.00     |
| 7 M  | Trichlorofluoromethane      | 20.000  | 17.053  | 14.7  | 111   | 0.00     |
| 8 T  | Diethyl Ether               | 20.000  | 15.901  | 20.5  | 108   | 0.02     |
| 9    | 1,1,2-Trichlorotrifluoroeth | 20.000  | 17.376  | 13.1  | 112   | 0.00     |
| 10 M | 1,1-Dichloroethene          | 20.000  | 16.540  | 17.3  | 111   | 0.00     |
| 11   | Methyl Iodide               | 20.000  | 13.660  | 31.7# | 109   | 0.00     |
| 12   | Methyl Acetate              | 20.000  | 21.448  | -7.2  | 137   | 0.00     |
| 13 M | Acrolein                    | 100.000 | 114.246 | -14.2 | 153   | 0.00     |
| 14 M | Acrylonitrile               | 100.000 | 97.945  | 2.1   | 130   | -0.01    |
| 15 M | Acetone                     | 100.000 | 87.219  | 12.8  | 119   | 0.00     |
| 16 M | Carbon Disulfide            | 20.000  | 15.750  | 21.3  | 105   | 0.00     |
| 17   | Allyl chloride              | 20.000  | 17.860  | 10.7  | 121   | 0.00     |
| 18 M | Methylene Chloride          | 20.000  | 20.538  | -2.7  | 139   | 0.00     |
| 19 M | trans-1,2-Dichloroethene    | 20.000  | 18.224  | 8.9   | 119   | 0.00     |
| 20 T | Diisopropyl ether           | 20.000  | 18.007  | 10.0  | 120   | 0.00     |
| 21 M | 1,1-Dichloroethane          | 20.000  | 18.429  | 7.9   | 122   | 0.00     |
| 22 M | cis-1,2-Dichloroethene      | 20.000  | 20.214  | -1.1  | 133   | 0.00     |
| 23 M | tert-Butyl Alcohol          | 100.000 | 99.358  | 0.6   | 128   | 0.00     |
| 24 M | Methyl tert-Butyl Ether     | 20.000  | 20.097  | -0.5  | 137   | 0.00     |
| 25 M | Chloroform                  | 20.000  | 19.665  | 1.7   | 128   | 0.00     |
| 26   | Cyclohexane                 | 20.000  | 17.027  | 14.9  | 115   | 0.00     |
| 27 s | 1,2-Dichloroethane-d4       | 30.000  | 28.485  | 5.1   | 113   | 0.00     |
| 28 I | 1,4-Difluorobenzene         | 30.000  | 30.000  | 0.0   | 117   | 0.00     |
| 29   | 1,1-Dichloropropene         | 20.000  | 19.600  | 2.0   | 124   | 0.00     |
| 30 M | 2-Butanone                  | 100.000 | 102.336 | -2.3  | 132   | 0.00     |
| 31   | 2,2-Dichloropropane         | 20.000  | 20.732  | -3.7  | 143   | 0.01     |
| 32 M | 1,1,1-Trichloroethane       | 20.000  | 20.954  | -4.8  | 130   | 0.00     |
| 33 M | Carbon Tetrachloride        | 20.000  | 20.620  | -3.1  | 133   | 0.00     |
| 34 M | Benzene                     | 20.000  | 20.222  | -1.1  | 129   | 0.00     |
| 35   | Methacrylonitrile           | 20.000  | 19.099  | 4.5   | 114   | 0.00     |
| 36 M | 1,2-Dichloroethane          | 20.000  | 20.322  | -1.6  | 127   | 0.00     |
| 37 M | Trichloroethene             | 20.000  | 20.777  | -3.9  | 132   | 0.00     |
| 38   | Methylcyclohexane           | 20.000  | 17.826  | 10.9  | 117   | 0.00     |
| 39 M | 1,2-Dichloropropane         | 20.000  | 20.138  | -0.7  | 126   | 0.00     |
| 40   | Dibromomethane              | 20.000  | 21.899  | -9.5  | 139   | 0.00     |
| 41 M | Bromodichloromethane        | 20.000  | 21.050  | -5.3  | 132   | 0.00     |
| 42 M | Vinyl Acetate               | 100.000 | 92.264  | 7.7   | 120   | 0.00     |
| 43   | Ethyl Acetate               | 20.000  | 19.579  | 2.1   | 121   | 0.00     |
| 44   | Isopropyl Acetate           | 20.000  | 19.018  | 4.9   | 120   | 0.00     |
| 45 T | 1,4-Dioxane                 | 400.000 | 427.798 | -6.9  | 127   | 0.00     |
| 46   | Methyl methacrylate         | 20.000  | 19.253  | 3.7   | 126   | 0.00     |
| 47   | n-amyl Acetate              | 20.000  | 14.796  | 26.0# | 121   | -0.11    |
| 48 M | t-1,3-Dichloropropene       | 20.000  | 19.836  | 0.8   | 132   | 0.00     |

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 Sample : VSTDCCC020  
 Misc : 5.0mL/MSVOA\_N/WATER  
 ALS Vial : 18 Sample Multiplier: 1

Instrument :  
 MSVOA\_N  
 LabSampleId :  
 VSTDCCC020

Quant Time: Sep 18 03:14:20 2025  
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_N\methods\624N090425W.M  
 Quant Title : METHOD 624 VOLATILE ORGANIC ANALYSIS  
 QLast Update : Fri Sep 05 03:29:53 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  | 19.979  | 0.1   | 134   | 0.00     |
| 50 M | 1,1,2-Trichloroethane       | 20.000  | 22.327  | -11.6 | 143   | 0.00     |
| 51   | Ethyl methacrylate          | 20.000  | 19.940  | 0.3   | 137   | 0.00     |
| 52   | 1,3-Dichloropropane         | 20.000  | 21.468  | -7.3  | 135   | 0.00     |
| 53 M | Dibromochloromethane        | 20.000  | 22.739  | -13.7 | 156   | 0.00     |
| 54 M | 1,2-Dibromoethane           | 20.000  | 22.136  | -10.7 | 143   | 0.00     |
| 55 M | 2-Chloroethyl vinyl ether   | 100.000 | 92.170  | 7.8   | 116   | 0.00     |
| 56 M | Bromoform                   | 20.000  | 23.489  | -17.4 | 161   | 0.00     |
| 57 I | Chlorobenzene-d5            | 30.000  | 30.000  | 0.0   | 121   | 0.00     |
| 58 M | 4-Methyl-2-Pentanone        | 100.000 | 101.406 | -1.4  | 133   | 0.00     |
| 59 M | 2-Hexanone                  | 100.000 | 96.401  | 3.6   | 129   | 0.00     |
| 60 S | 4-Bromofluorobenzene        | 30.000  | 30.278  | -0.9  | 120   | 0.00     |
| 61 M | Tetrachloroethene           | 20.000  | 20.313  | -1.6  | 131   | 0.00     |
| 62 M | Toluene                     | 20.000  | 19.619  | 1.9   | 128   | 0.00     |
| 63 S | Toluene-d8                  | 30.000  | 28.681  | 4.4   | 114   | 0.00     |
| 64 M | Chlorobenzene               | 20.000  | 20.237  | -1.2  | 135   | 0.00     |
| 65   | 1,1,1,2-Tetrachloroethane   | 20.000  | 21.515  | -7.6  | 144   | 0.00     |
| 66 M | Ethyl Benzene               | 20.000  | 19.184  | 4.1   | 130   | 0.00     |
| 67 M | m/p-Xylenes                 | 40.000  | 40.618  | -1.5  | 139   | 0.00     |
| 68 M | o-Xylene                    | 20.000  | 20.678  | -3.4  | 138   | 0.00     |
| 69 M | Styrene                     | 20.000  | 20.180  | -0.9  | 139   | 0.00     |
| 70   | Isopropylbenzene            | 20.000  | 20.470  | -2.3  | 142   | 0.00     |
| 71 M | 1,1,2,2-Tetrachloroethane   | 20.000  | 21.921  | -9.6  | 140   | 0.00     |
| 72   | 1,2,3-Trichloropropane      | 20.000  | 20.958  | -4.8  | 125   | 0.00     |
| 73   | Bromobenzene                | 20.000  | 22.538  | -12.7 | 151   | 0.00     |
| 74   | n-propylbenzene             | 20.000  | 19.794  | 1.0   | 133   | 0.00     |
| 75   | 2-Chlorotoluene             | 20.000  | 20.533  | -2.7  | 135   | 0.00     |
| 76   | 1,3,5-Trimethylbenzene      | 20.000  | 20.675  | -3.4  | 139   | 0.00     |
| 77   | t-1,4-Dichloro-2-butene     | 20.000  | 20.038  | -0.2  | 141   | 0.00     |
| 78   | 4-Chlorotoluene             | 20.000  | 20.570  | -2.9  | 135   | 0.00     |
| 79   | tert-butylbenzene           | 20.000  | 20.916  | -4.6  | 141   | 0.00     |
| 80   | 1,2,4-Trimethylbenzene      | 20.000  | 20.601  | -3.0  | 137   | 0.00     |
| 81   | sec-Butylbenzene            | 20.000  | 20.558  | -2.8  | 137   | 0.00     |
| 82   | p-Isopropyltoluene          | 20.000  | 20.914  | -4.6  | 142   | 0.00     |
| 83 M | 1,3-Dichlorobenzene         | 20.000  | 21.657  | -8.3  | 140   | 0.00     |
| 84 M | 1,4-Dichlorobenzene         | 20.000  | 22.673  | -13.4 | 149   | 0.00     |
| 85   | n-Butylbenzene              | 20.000  | 20.422  | -2.1  | 141   | 0.00     |
| 86 T | Hexachloroethane            | 20.000  | 21.030  | -5.2  | 138   | 0.00     |
| 87 M | 1,2-Dichlorobenzene         | 20.000  | 22.821  | -14.1 | 149   | 0.00     |
| 88   | 1,2-Dibromo-3-Chloropropane | 20.000  | 22.956  | -14.8 | 142   | 0.00     |
| 89   | 1,2,4-Trichlorobenzene      | 20.000  | 22.001  | -10.0 | 150   | 0.00     |
| 90   | Hexachlorobutadiene         | 20.000  | 23.438  | -17.2 | 147   | 0.00     |
| 91 M | Naphthalene                 | 20.000  | 20.825  | -4.1  | 145   | 0.00     |
| 92   | 1,2,3-Trichlorobenzene      | 20.000  | 22.001  | -10.0 | 150   | 0.00     |

(#) = Out of Range

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