

Data Path : Z:\voasrv\HPCHEM1\MSVOA\_N\Data\VN072225\  
 Data File : VN087397.D  
 Acq On : 22 Jul 2025 18:11  
 Operator : JC\MD  
 Sample : VSTDCCC020  
 Misc : 5.0mL/MSVOA\_N/WATER  
 ALS Vial : 10 Sample Multiplier: 1

Instrument :  
 MSVOA\_N  
 LabSampleId :  
 VSTDCCC020

Quant Time: Jul 23 05:58:50 2025  
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_N\methods\624N062525W.M  
 Quant Title : METHOD 624 VOLATILE ORGANIC ANALYSIS  
 QLast Update : Wed Jun 25 10:49:56 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                    | AvgRF | CCRF  | %Dev  | Area% | Dev(min) |
|------|-----------------------------|-------|-------|-------|-------|----------|
| 1 I  | Bromochloromethane          | 1.000 | 1.000 | 0.0   | 77    | -0.02    |
| 2 M  | Dichlorodifluoromethane     | 1.806 | 1.113 | 38.4# | 52    | -0.01    |
| 3 M  | Chloromethane               | 1.844 | 1.262 | 31.6# | 62    | -0.01    |
| 4 M  | Vinyl Chloride              | 2.005 | 1.413 | 29.5# | 60    | -0.01    |
| 5 M  | Bromomethane                | 1.064 | 0.663 | 37.7# | 53    | -0.01    |
| 6 M  | Chloroethane                | 1.075 | 1.220 | -13.5 | 97    | -0.02    |
| 7 M  | Trichlorofluoromethane      | 2.459 | 2.797 | -13.7 | 93    | -0.02    |
| 8 T  | Diethyl Ether               | 1.208 | 1.130 | 6.5   | 81    | -0.02    |
| 9    | 1,1,2-Trichlorotrifluoroeth | 1.823 | 1.479 | 18.9  | 68    | -0.04    |
| 10 M | 1,1-Dichloroethene          | 1.817 | 1.411 | 22.3  | 65    | -0.04    |
| 11   | Methyl Iodide               | 1.628 | 0.779 | 52.1# | 46#   | -0.02    |
| 12   | Methyl Acetate              | 2.851 | 2.459 | 13.7  | 72    | -0.03    |
| 13 M | Acrolein                    | 0.438 | 0.393 | 10.3  | 84    | -0.03    |
| 14 M | Acrylonitrile               | 1.330 | 0.983 | 26.1# | 62    | -0.02    |
| 15 M | Acetone                     | 0.388 | 0.343 | 11.6  | 72    | -0.03    |
| 16 M | Carbon Disulfide            | 5.313 | 3.201 | 39.8# | 51    | -0.03    |
| 17   | Allyl chloride              | 2.997 | 2.005 | 33.1# | 56    | -0.04    |
| 18 M | Methylene Chloride          | 2.105 | 1.633 | 22.4  | 65    | -0.04    |
| 19 M | trans-1,2-Dichloroethene    | 1.987 | 1.445 | 27.3# | 61    | -0.03    |
| 20 T | Diisopropyl ether           | 6.359 | 5.072 | 20.2  | 64    | -0.03    |
| 21 M | 1,1-Dichloroethane          | 3.801 | 2.815 | 25.9# | 61    | -0.02    |
| 22 M | cis-1,2-Dichloroethene      | 2.313 | 1.756 | 24.1  | 65    | -0.02    |
| 23 M | tert-Butyl Alcohol          | 0.500 | 0.369 | 26.2# | 63    | -0.02    |
| 24 M | Methyl tert-Butyl Ether     | 6.374 | 5.122 | 19.6  | 69    | -0.02    |
| 25 M | Chloroform                  | 3.641 | 3.265 | 10.3  | 76    | -0.02    |
| 26   | Cyclohexane                 | 3.159 | 1.698 | 46.2# | 46#   | -0.02    |
| 27 s | 1,2-Dichloroethane-d4       | 2.260 | 2.435 | -7.7  | 81    | -0.02    |
| 28 I | 1,4-Difluorobenzene         | 1.000 | 1.000 | 0.0   | 62    | -0.02    |
| 29   | 1,1-Dichloropropene         | 0.468 | 0.416 | 11.1  | 65    | -0.02    |
| 30 M | 2-Butanone                  | 0.315 | 0.323 | -2.5  | 70    | -0.02    |
| 31   | 2,2-Dichloropropane         | 0.564 | 0.586 | -3.9  | 70    | -0.02    |
| 32 M | 1,1,1-Trichloroethane       | 0.546 | 0.595 | -9.0  | 74    | -0.02    |
| 33 M | Carbon Tetrachloride        | 0.460 | 0.508 | -10.4 | 79    | -0.02    |
| 34 M | Benzene                     | 1.481 | 1.367 | 7.7   | 65    | -0.02    |
| 35   | Methacrylonitrile           | 0.318 | 0.304 | 4.4   | 66    | -0.03    |
| 36 M | 1,2-Dichloroethane          | 0.479 | 0.578 | -20.7 | 81    | -0.02    |
| 37 M | Trichloroethene             | 0.342 | 0.315 | 7.9   | 67    | -0.02    |
| 38   | Methylcyclohexane           | 0.537 | 0.376 | 30.0# | 50#   | -0.02    |
| 39 M | 1,2-Dichloropropane         | 0.372 | 0.361 | 3.0   | 65    | -0.02    |
| 40   | Dibromomethane              | 0.253 | 0.274 | -8.3  | 73    | -0.02    |
| 41 M | Bromodichloromethane        | 0.510 | 0.573 | -12.4 | 74    | -0.02    |
| 42 M | Vinyl Acetate               | 1.026 | 0.909 | 11.4  | 60    | -0.02    |
| 43   | Ethyl Acetate               | 0.569 | 0.621 | -9.1  | 69    | -0.02    |
| 44   | Isopropyl Acetate           | 0.928 | 0.967 | -4.2  | 72    | -0.02    |
| 45 T | 1,4-Dioxane                 | 0.006 | 0.007 | -16.7 | 77    | -0.02    |
| 46   | Methyl methacrylate         | 0.431 | 0.433 | -0.5  | 68    | -0.02    |
| 47   | n-amyl Acetate              | 0.567 | 0.579 | -2.1  | 71    | -0.02    |
| 48 M | t-1,3-Dichloropropene       | 0.569 | 0.553 | 2.8   | 67    | -0.02    |

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 Sample : VSTDCCC020  
 Misc : 5.0mL/MSVOA\_N/WATER  
 ALS Vial : 10 Sample Multiplier: 1

Instrument :  
 MSVOA\_N  
 LabSampleId :  
 VSTDCCC020

Quant Time: Jul 23 05:58:50 2025  
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_N\methods\624N062525W.M  
 Quant Title : METHOD 624 VOLATILE ORGANIC ANALYSIS  
 QLast Update : Wed Jun 25 10:49:56 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                    | AvgRF | CCRF  | %Dev   | Area% | Dev(min) |
|------|-----------------------------|-------|-------|--------|-------|----------|
| 49 T | cis-1,3-Dichloropropene     | 0.605 | 0.591 | 2.3    | 65    | -0.02    |
| 50 M | 1,1,2-Trichloroethane       | 0.359 | 0.371 | -3.3   | 69    | -0.02    |
| 51   | Ethyl methacrylate          | 0.547 | 0.523 | 4.4    | 69    | -0.03    |
| 52   | 1,3-Dichloropropane         | 0.617 | 0.612 | 0.8    | 66    | -0.02    |
| 53 M | Dibromochloromethane        | 0.377 | 0.443 | -17.5  | 80    | -0.02    |
| 54 M | 1,2-Dibromoethane           | 0.371 | 0.387 | -4.3   | 70    | -0.02    |
| 55 M | 2-Chloroethyl vinyl ether   | 0.313 | 0.321 | -2.6   | 69    | -0.02    |
| 56 M | Bromoform                   | 0.259 | 0.291 | -12.4  | 77    | -0.02    |
| 57 I | Chlorobenzene-d5            | 1.000 | 1.000 | 0.0    | 64    | -0.02    |
| 58 M | 4-Methyl-2-Pentanone        | 0.603 | 0.684 | -13.4  | 80    | -0.02    |
| 59 M | 2-Hexanone                  | 0.389 | 0.440 | -13.1  | 87    | -0.02    |
| 60 S | 4-Bromofluorobenzene        | 0.463 | 0.482 | -4.1   | 69    | -0.02    |
| 61 M | Tetrachloroethene           | 0.297 | 0.272 | 8.4    | 65    | -0.02    |
| 62 M | Toluene                     | 1.694 | 1.559 | 8.0    | 64    | -0.02    |
| 63 S | Toluene-d8                  | 1.347 | 1.384 | -2.7   | 63    | -0.02    |
| 64 M | Chlorobenzene               | 1.081 | 0.986 | 8.8    | 64    | -0.02    |
| 65   | 1,1,1,2-Tetrachloroethane   | 0.357 | 0.391 | -9.5   | 77    | -0.02    |
| 66 M | Ethyl Benzene               | 1.855 | 1.577 | 15.0   | 61    | -0.02    |
| 67 M | m/p-Xylenes                 | 0.711 | 0.619 | 12.9   | 62    | -0.02    |
| 68 M | o-Xylene                    | 0.678 | 0.578 | 14.7   | 61    | -0.02    |
| 69 M | Styrene                     | 1.164 | 1.042 | 10.5   | 64    | -0.02    |
| 70   | Isopropylbenzene            | 1.689 | 1.500 | 11.2   | 65    | -0.02    |
| 71 M | 1,1,2,2-Tetrachloroethane   | 0.608 | 0.604 | 0.7    | 68    | -0.02    |
| 72   | 1,2,3-Trichloropropane      | 0.524 | 0.507 | 3.2    | 70    | -0.01    |
| 73   | Bromobenzene                | 0.410 | 0.409 | 0.2    | 72    | -0.02    |
| 74   | n-propylbenzene             | 2.063 | 1.948 | 5.6    | 69    | -0.02    |
| 75   | 2-Chlorotoluene             | 1.246 | 1.161 | 6.8    | 69    | -0.02    |
| 76   | 1,3,5-Trimethylbenzene      | 1.375 | 1.304 | 5.2    | 71    | -0.02    |
| 77   | t-1,4-Dichloro-2-butene     | 0.232 | 0.163 | 29.7#  | 53    | -0.01    |
| 78   | 4-Chlorotoluene             | 1.276 | 1.258 | 1.4    | 73    | -0.02    |
| 79   | tert-butylbenzene           | 1.156 | 1.094 | 5.4    | 71    | -0.02    |
| 80   | 1,2,4-Trimethylbenzene      | 1.380 | 1.335 | 3.3    | 72    | -0.02    |
| 81   | sec-Butylbenzene            | 1.680 | 1.597 | 4.9    | 69    | -0.02    |
| 82   | p-Isopropyltoluene          | 1.398 | 1.321 | 5.5    | 67    | -0.02    |
| 83 M | 1,3-Dichlorobenzene         | 0.789 | 0.802 | -1.6   | 71    | -0.02    |
| 84 M | 1,4-Dichlorobenzene         | 0.794 | 0.775 | 2.4    | 69    | -0.02    |
| 85   | n-Butylbenzene              | 1.275 | 1.234 | 3.2    | 68    | -0.02    |
| 86 T | Hexachloroethane            | 0.265 | 0.285 | -7.5   | 80    | -0.02    |
| 87 M | 1,2-Dichlorobenzene         | 0.744 | 0.775 | -4.2   | 74    | -0.02    |
| 88   | 1,2-Dibromo-3-Chloropropane | 0.137 | 0.141 | -2.9   | 74    | -0.01    |
| 89   | 1,2,4-Trichlorobenzene      | 0.421 | 0.443 | -5.2   | 75    | -0.02    |
| 90   | Hexachlorobutadiene         | 0.125 | 0.190 | -52.0# | 108   | -0.02    |
| 91 M | Naphthalene                 | 1.656 | 1.418 | 14.4   | 63    | -0.02    |
| 92   | 1,2,3-Trichlorobenzene      | 0.421 | 0.443 | -5.2   | 75    | -0.02    |

(#) = Out of Range

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