

Data Path : Z:\voasrv\HPCHEM1\MSVOA\_N\Data\VN101723\  
 Data File : VN079577.D  
 Acq On : 17 Oct 2023 12:49  
 Operator : JC\MD  
 Sample : VSTDICV020  
 Misc : 5.0mL/MSVOA\_N/WATER  
 ALS Vial : 9 Sample Multiplier: 1

Instrument :  
 MSVOA\_N  
 ClientSampleId :  
 ICSVN101723

Quant Time: Oct 18 02:21:38 2023  
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_N\methods\624N101723W.M  
 Quant Title : METHOD 624 VOLATILE ORGANIC ANALYSIS  
 QLast Update : Wed Oct 18 02:19:15 2023  
 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  | 104   | -0.01    |
| 2 M  | Dichlorodifluoromethane     | 20.000  | 19.806  | 1.0  | 99    | 0.00     |
| 3 M  | Chloromethane               | 20.000  | 19.489  | 2.6  | 98    | 0.00     |
| 4 M  | Vinyl Chloride              | 20.000  | 20.007  | -0.0 | 102   | 0.00     |
| 5 M  | Bromomethane                | 20.000  | 19.272  | 3.6  | 99    | 0.00     |
| 6 M  | Chloroethane                | 20.000  | 19.806  | 1.0  | 101   | 0.00     |
| 7 M  | Trichlorofluoromethane      | 20.000  | 19.773  | 1.1  | 100   | 0.00     |
| 8 T  | Diethyl Ether               | 20.000  | 19.799  | 1.0  | 100   | 0.00     |
| 9    | 1,1,2-Trichlorotrifluoroeth | 20.000  | 20.095  | -0.5 | 103   | -0.01    |
| 10 M | 1,1-Dichloroethene          | 20.000  | 19.629  | 1.9  | 99    | 0.00     |
| 11   | Methyl Iodide               | 20.000  | 19.443  | 2.8  | 101   | 0.00     |
| 12   | Methyl Acetate              | 20.000  | 19.544  | 2.3  | 100   | 0.00     |
| 13 M | Acrolein                    | 100.000 | 93.968  | 6.0  | 96    | 0.00     |
| 14 M | Acrylonitrile               | 100.000 | 99.310  | 0.7  | 101   | 0.00     |
| 15 M | Acetone                     | 100.000 | 100.442 | -0.4 | 100   | 0.00     |
| 16 M | Carbon Disulfide            | 20.000  | 19.460  | 2.7  | 100   | 0.00     |
| 17   | Allyl chloride              | 20.000  | 20.215  | -1.1 | 98    | 0.00     |
| 18 M | Methylene Chloride          | 20.000  | 19.982  | 0.1  | 101   | 0.00     |
| 19 M | trans-1,2-Dichloroethene    | 20.000  | 19.966  | 0.2  | 103   | 0.00     |
| 20 T | Diisopropyl ether           | 20.000  | 20.426  | -2.1 | 100   | 0.00     |
| 21 M | 1,1-Dichloroethane          | 20.000  | 19.945  | 0.3  | 100   | 0.00     |
| 22 M | cis-1,2-Dichloroethene      | 20.000  | 20.147  | -0.7 | 102   | 0.00     |
| 23 M | tert-Butyl Alcohol          | 100.000 | 89.962  | 10.0 | 92    | 0.00     |
| 24 M | Methyl tert-Butyl Ether     | 20.000  | 19.900  | 0.5  | 100   | 0.00     |
| 25 M | Chloroform                  | 20.000  | 20.283  | -1.4 | 102   | 0.00     |
| 26   | Cyclohexane                 | 20.000  | 20.512  | -2.6 | 102   | 0.00     |
| 27 s | 1,2-Dichloroethane-d4       | 30.000  | 30.275  | -0.9 | 101   | 0.00     |
| 28 I | 1,4-Difluorobenzene         | 30.000  | 30.000  | 0.0  | 102   | 0.00     |
| 29   | 1,1-Dichloropropene         | 20.000  | 19.489  | 2.6  | 102   | 0.00     |
| 30 M | 2-Butanone                  | 100.000 | 93.182  | 6.8  | 96    | 0.00     |
| 31   | 2,2-Dichloropropane         | 20.000  | 18.969  | 5.2  | 100   | 0.00     |
| 32 M | 1,1,1-Trichloroethane       | 20.000  | 19.071  | 4.6  | 100   | 0.00     |
| 33 M | Carbon Tetrachloride        | 20.000  | 19.111  | 4.4  | 99    | 0.00     |
| 34 M | Benzene                     | 20.000  | 19.041  | 4.8  | 100   | 0.00     |
| 35   | Methacrylonitrile           | 20.000  | 18.696  | 6.5  | 96    | 0.00     |
| 36 M | 1,2-Dichloroethane          | 20.000  | 19.027  | 4.9  | 99    | 0.00     |
| 37 M | Trichloroethene             | 20.000  | 19.206  | 4.0  | 104   | 0.00     |
| 38   | Methylcyclohexane           | 20.000  | 18.977  | 5.1  | 100   | 0.00     |
| 39 M | 1,2-Dichloropropane         | 20.000  | 19.067  | 4.7  | 100   | 0.00     |
| 40   | Dibromomethane              | 20.000  | 19.575  | 2.1  | 104   | 0.00     |
| 41 M | Bromodichloromethane        | 20.000  | 19.177  | 4.1  | 101   | 0.00     |
| 42 M | Vinyl Acetate               | 100.000 | 96.769  | 3.2  | 100   | 0.00     |
| 43   | Ethyl Acetate               | 20.000  | 20.373  | -1.9 | 108   | 0.00     |
| 44   | Isopropyl Acetate           | 20.000  | 18.222  | 8.9  | 99    | 0.00     |
| 45 T | 1,4-Dioxane                 | 400.000 | 371.469 | 7.1  | 94    | 0.00     |
| 46   | Methyl methacrylate         | 20.000  | 19.265  | 3.7  | 99    | 0.00     |
| 47   | n-amyl Acetate              | 20.000  | 18.713  | 6.4  | 95    | 0.00     |
| 48 M | t-1,3-Dichloropropene       | 20.000  | 19.068  | 4.7  | 99    | 0.00     |

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

Instrument :  
 MSVOA\_N  
 ClientSampleId :  
 ICSVN101723

Quant Time: Oct 18 02:21:38 2023  
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_N\methods\624N101723W.M  
 Quant Title : METHOD 624 VOLATILE ORGANIC ANALYSIS  
 QLast Update : Wed Oct 18 02:19:15 2023  
 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  | 19.118 | 4.4  | 100   | 0.00     |
| 50 M | 1,1,2-Trichloroethane       | 20.000  | 19.385 | 3.1  | 102   | 0.00     |
| 51   | Ethyl methacrylate          | 20.000  | 19.176 | 4.1  | 98    | 0.00     |
| 52   | 1,3-Dichloropropane         | 20.000  | 18.737 | 6.3  | 99    | 0.00     |
| 53 M | Dibromochloromethane        | 20.000  | 19.213 | 3.9  | 99    | 0.00     |
| 54 M | 1,2-Dibromoethane           | 20.000  | 19.092 | 4.5  | 101   | 0.00     |
| 55 M | 2-Chloroethyl vinyl ether   | 100.000 | 86.156 | 13.8 | 94    | 0.00     |
| 56 M | Bromoform                   | 20.000  | 18.514 | 7.4  | 96    | 0.00     |
| 57 I | Chlorobenzene-d5            | 30.000  | 30.000 | 0.0  | 101   | 0.00     |
| 58 M | 4-Methyl-2-Pentanone        | 100.000 | 99.463 | 0.5  | 98    | 0.00     |
| 59 M | 2-Hexanone                  | 100.000 | 97.963 | 2.0  | 97    | 0.00     |
| 60 S | 4-Bromofluorobenzene        | 30.000  | 30.842 | -2.8 | 102   | 0.00     |
| 61 M | Tetrachloroethene           | 20.000  | 20.033 | -0.2 | 100   | 0.00     |
| 62 M | Toluene                     | 20.000  | 19.832 | 0.8  | 101   | 0.00     |
| 63 S | Toluene-d8                  | 30.000  | 30.092 | -0.3 | 100   | 0.00     |
| 64 M | Chlorobenzene               | 20.000  | 19.781 | 1.1  | 102   | 0.00     |
| 65   | 1,1,1,2-Tetrachloroethane   | 20.000  | 19.493 | 2.5  | 99    | 0.00     |
| 66 M | Ethyl Benzene               | 20.000  | 19.979 | 0.1  | 101   | 0.00     |
| 67 M | m/p-Xylenes                 | 40.000  | 40.808 | -2.0 | 102   | 0.00     |
| 68 M | o-Xylene                    | 20.000  | 20.739 | -3.7 | 105   | 0.00     |
| 69 M | Styrene                     | 20.000  | 20.740 | -3.7 | 102   | 0.00     |
| 70   | Isopropylbenzene            | 20.000  | 20.114 | -0.6 | 101   | 0.00     |
| 71 M | 1,1,2,2-Tetrachloroethane   | 20.000  | 19.336 | 3.3  | 98    | 0.00     |
| 72   | 1,2,3-Trichloropropane      | 20.000  | 18.981 | 5.1  | 97    | 0.00     |
| 73   | Bromobenzene                | 20.000  | 19.848 | 0.8  | 99    | 0.00     |
| 74   | n-propylbenzene             | 20.000  | 20.377 | -1.9 | 102   | 0.00     |
| 75   | 2-Chlorotoluene             | 20.000  | 20.460 | -2.3 | 103   | 0.00     |
| 76   | 1,3,5-Trimethylbenzene      | 20.000  | 20.667 | -3.3 | 103   | 0.00     |
| 77   | t-1,4-Dichloro-2-butene     | 20.000  | 19.675 | 1.6  | 97    | 0.00     |
| 78   | 4-Chlorotoluene             | 20.000  | 20.226 | -1.1 | 100   | 0.00     |
| 79   | tert-butylbenzene           | 20.000  | 20.099 | -0.5 | 102   | 0.00     |
| 80   | 1,2,4-Trimethylbenzene      | 20.000  | 20.465 | -2.3 | 102   | 0.00     |
| 81   | sec-Butylbenzene            | 20.000  | 20.170 | -0.9 | 102   | 0.00     |
| 82   | p-Isopropyltoluene          | 20.000  | 20.434 | -2.2 | 102   | 0.00     |
| 83 M | 1,3-Dichlorobenzene         | 20.000  | 20.734 | -3.7 | 103   | 0.00     |
| 84 M | 1,4-Dichlorobenzene         | 20.000  | 20.124 | -0.6 | 102   | 0.00     |
| 85   | n-Butylbenzene              | 20.000  | 20.007 | -0.0 | 100   | 0.00     |
| 86 T | Hexachloroethane            | 20.000  | 19.014 | 4.9  | 99    | 0.00     |
| 87 M | 1,2-Dichlorobenzene         | 20.000  | 20.072 | -0.4 | 97    | 0.00     |
| 88   | 1,2-Dibromo-3-Chloropropane | 20.000  | 18.829 | 5.9  | 90    | 0.00     |
| 89   | 1,2,4-Trichlorobenzene      | 20.000  | 17.760 | 11.2 | 87    | 0.00     |
| 90   | Hexachlorobutadiene         | 20.000  | 18.637 | 6.8  | 93    | 0.00     |
| 91 M | Naphthalene                 | 20.000  | 15.114 | 24.4 | 76    | 0.00     |
| 92   | 1,2,3-Trichlorobenzene      | 20.000  | 17.760 | 11.2 | 87    | 0.00     |

(#) = Out of Range

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