

Data Path : Z:\voasrv\HPCHEM1\MSVOA\_N\Data\VN020924\  
 Data File : VN080984.D  
 Acq On : 09 Feb 2024 12:27  
 Operator : JC\MD  
 Sample : VSTDICV020  
 Misc : 5.0mL/MSVOA\_N/WATER  
 ALS Vial : 8 Sample Multiplier: 1

Instrument :  
 MSVOA\_N  
 ClientSampleId :  
 ICSVN020924

Quant Time: Feb 10 01:55:59 2024  
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_N\methods\624N020924W.M  
 Quant Title : METHOD 624 VOLATILE ORGANIC ANALYSIS  
 QLast Update : Sat Feb 10 01:52:21 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  | 90    | 0.00     |
| 2 M  | Dichlorodifluoromethane     | 20.000  | 20.721  | -3.6 | 102   | 0.00     |
| 3 M  | Chloromethane               | 20.000  | 21.047  | -5.2 | 102   | 0.00     |
| 4 M  | Vinyl Chloride              | 20.000  | 21.402  | -7.0 | 103   | 0.00     |
| 5 M  | Bromomethane                | 20.000  | 20.656  | -3.3 | 97    | 0.04     |
| 6 M  | Chloroethane                | 20.000  | 21.582  | -7.9 | 102   | 0.02     |
| 7 M  | Trichlorofluoromethane      | 20.000  | 20.686  | -3.4 | 95    | 0.01     |
| 8 T  | Diethyl Ether               | 20.000  | 19.569  | 2.2  | 92    | 0.00     |
| 9    | 1,1,2-Trichlorotrifluoroeth | 20.000  | 20.507  | -2.5 | 94    | 0.01     |
| 10 M | 1,1-Dichloroethene          | 20.000  | 20.945  | -4.7 | 96    | 0.00     |
| 11   | Methyl Iodide               | 20.000  | 18.244  | 8.8  | 89    | 0.00     |
| 12   | Methyl Acetate              | 20.000  | 19.926  | 0.4  | 96    | 0.00     |
| 13 M | Acrolein                    | 100.000 | 83.617  | 16.4 | 75    | 0.00     |
| 14 M | Acrylonitrile               | 100.000 | 101.614 | -1.6 | 96    | 0.00     |
| 15 M | Acetone                     | 100.000 | 102.676 | -2.7 | 97    | 0.00     |
| 16 M | Carbon Disulfide            | 20.000  | 19.417  | 2.9  | 89    | 0.01     |
| 17   | Allyl chloride              | 20.000  | 20.956  | -4.8 | 97    | 0.00     |
| 18 M | Methylene Chloride          | 20.000  | 19.544  | 2.3  | 89    | 0.00     |
| 19 M | trans-1,2-Dichloroethene    | 20.000  | 19.846  | 0.8  | 90    | 0.01     |
| 20 T | Diisopropyl ether           | 20.000  | 20.867  | -4.3 | 99    | 0.00     |
| 21 M | 1,1-Dichloroethane          | 20.000  | 20.533  | -2.7 | 95    | 0.00     |
| 22 M | cis-1,2-Dichloroethene      | 20.000  | 20.467  | -2.3 | 93    | 0.00     |
| 23 M | tert-Butyl Alcohol          | 100.000 | 108.853 | -8.9 | 100   | 0.00     |
| 24 M | Methyl tert-Butyl Ether     | 20.000  | 20.233  | -1.2 | 94    | 0.00     |
| 25 M | Chloroform                  | 20.000  | 20.944  | -4.7 | 95    | 0.00     |
| 26   | Cyclohexane                 | 20.000  | 21.443  | -7.2 | 97    | 0.00     |
| 27 s | 1,2-Dichloroethane-d4       | 30.000  | 30.716  | -2.4 | 94    | 0.00     |
| 28 I | 1,4-Difluorobenzene         | 30.000  | 30.000  | 0.0  | 89    | 0.00     |
| 29   | 1,1-Dichloropropene         | 20.000  | 20.670  | -3.4 | 94    | 0.00     |
| 30 M | 2-Butanone                  | 100.000 | 101.785 | -1.8 | 95    | 0.00     |
| 31   | 2,2-Dichloropropane         | 20.000  | 21.568  | -7.8 | 99    | 0.00     |
| 32 M | 1,1,1-Trichloroethane       | 20.000  | 20.789  | -3.9 | 96    | 0.00     |
| 33 M | Carbon Tetrachloride        | 20.000  | 20.823  | -4.1 | 95    | 0.00     |
| 34 M | Benzene                     | 20.000  | 20.756  | -3.8 | 95    | 0.00     |
| 35   | Methacrylonitrile           | 20.000  | 20.742  | -3.7 | 100   | 0.00     |
| 36 M | 1,2-Dichloroethane          | 20.000  | 20.933  | -4.7 | 95    | 0.00     |
| 37 M | Trichloroethene             | 20.000  | 21.097  | -5.5 | 94    | 0.00     |
| 38   | Methylcyclohexane           | 20.000  | 20.854  | -4.3 | 96    | 0.00     |
| 39 M | 1,2-Dichloropropane         | 20.000  | 20.606  | -3.0 | 94    | 0.00     |
| 40   | Dibromomethane              | 20.000  | 21.008  | -5.0 | 96    | 0.00     |
| 41 M | Bromodichloromethane        | 20.000  | 20.293  | -1.5 | 95    | 0.00     |
| 42 M | Vinyl Acetate               | 100.000 | 101.941 | -1.9 | 97    | 0.00     |
| 43   | Ethyl Acetate               | 20.000  | 20.557  | -2.8 | 94    | 0.00     |
| 44   | Isopropyl Acetate           | 20.000  | 20.537  | -2.7 | 96    | 0.00     |
| 45 T | 1,4-Dioxane                 | 400.000 | 435.018 | -8.8 | 100   | 0.00     |
| 46   | Methyl methacrylate         | 20.000  | 19.986  | 0.1  | 95    | 0.00     |
| 47   | n-amyl Acetate              | 20.000  | 19.672  | 1.6  | 92    | 0.00     |
| 48 M | t-1,3-Dichloropropene       | 20.000  | 20.610  | -3.0 | 97    | 0.00     |

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 ALS Vial : 8 Sample Multiplier: 1

Instrument :  
 MSVOA\_N  
 ClientSampleId :  
 ICVVN020924

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 Quant Title : METHOD 624 VOLATILE ORGANIC ANALYSIS  
 QLast Update : Sat Feb 10 01:52:21 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  | 20.169  | -0.8 | 93    | 0.00     |
| 50 M | 1,1,2-Trichloroethane       | 20.000  | 19.909  | 0.5  | 93    | 0.00     |
| 51   | Ethyl methacrylate          | 20.000  | 19.537  | 2.3  | 92    | 0.00     |
| 52   | 1,3-Dichloropropane         | 20.000  | 20.449  | -2.2 | 94    | 0.00     |
| 53 M | Dibromochloromethane        | 20.000  | 19.199  | 4.0  | 91    | 0.00     |
| 54 M | 1,2-Dibromoethane           | 20.000  | 20.139  | -0.7 | 95    | 0.00     |
| 55 M | 2-Chloroethyl vinyl ether   | 100.000 | 97.922  | 2.1  | 92    | 0.00     |
| 56 M | Bromoform                   | 20.000  | 18.099  | 9.5  | 87    | 0.00     |
| 57 I | Chlorobenzene-d5            | 30.000  | 30.000  | 0.0  | 89    | 0.00     |
| 58 M | 4-Methyl-2-Pentanone        | 100.000 | 105.309 | -5.3 | 96    | 0.00     |
| 59 M | 2-Hexanone                  | 100.000 | 107.938 | -7.9 | 100   | 0.00     |
| 60 S | 4-Bromofluorobenzene        | 30.000  | 31.070  | -3.6 | 90    | 0.00     |
| 61 M | Tetrachloroethene           | 20.000  | 21.175  | -5.9 | 93    | 0.00     |
| 62 M | Toluene                     | 20.000  | 20.945  | -4.7 | 93    | 0.00     |
| 63 S | Toluene-d8                  | 30.000  | 30.898  | -3.0 | 90    | 0.00     |
| 64 M | Chlorobenzene               | 20.000  | 20.739  | -3.7 | 92    | 0.00     |
| 65   | 1,1,1,2-Tetrachloroethane   | 20.000  | 20.229  | -1.1 | 90    | 0.00     |
| 66 M | Ethyl Benzene               | 20.000  | 21.128  | -5.6 | 93    | 0.00     |
| 67 M | m/p-Xylenes                 | 40.000  | 41.382  | -3.5 | 90    | 0.00     |
| 68 M | o-Xylene                    | 20.000  | 20.632  | -3.2 | 93    | 0.00     |
| 69 M | Styrene                     | 20.000  | 20.490  | -2.4 | 92    | 0.00     |
| 70   | Isopropylbenzene            | 20.000  | 20.663  | -3.3 | 91    | 0.00     |
| 71 M | 1,1,2,2-Tetrachloroethane   | 20.000  | 20.253  | -1.3 | 91    | 0.00     |
| 72   | 1,2,3-Trichloropropane      | 20.000  | 20.683  | -3.4 | 96    | 0.00     |
| 73   | Bromobenzene                | 20.000  | 20.519  | -2.6 | 90    | 0.00     |
| 74   | n-propylbenzene             | 20.000  | 20.687  | -3.4 | 91    | 0.00     |
| 75   | 2-Chlorotoluene             | 20.000  | 21.056  | -5.3 | 92    | 0.00     |
| 76   | 1,3,5-Trimethylbenzene      | 20.000  | 20.753  | -3.8 | 91    | 0.00     |
| 77   | t-1,4-Dichloro-2-butene     | 20.000  | 19.811  | 0.9  | 93    | 0.00     |
| 78   | 4-Chlorotoluene             | 20.000  | 20.931  | -4.7 | 91    | 0.00     |
| 79   | tert-butylbenzene           | 20.000  | 20.265  | -1.3 | 87    | 0.00     |
| 80   | 1,2,4-Trimethylbenzene      | 20.000  | 21.202  | -6.0 | 93    | 0.00     |
| 81   | sec-Butylbenzene            | 20.000  | 20.845  | -4.2 | 91    | 0.00     |
| 82   | p-Isopropyltoluene          | 20.000  | 20.328  | -1.6 | 91    | 0.00     |
| 83 M | 1,3-Dichlorobenzene         | 20.000  | 20.280  | -1.4 | 87    | 0.00     |
| 84 M | 1,4-Dichlorobenzene         | 20.000  | 20.054  | -0.3 | 87    | 0.00     |
| 85   | n-Butylbenzene              | 20.000  | 20.785  | -3.9 | 93    | 0.00     |
| 86 T | Hexachloroethane            | 20.000  | 19.245  | 3.8  | 87    | 0.00     |
| 87 M | 1,2-Dichlorobenzene         | 20.000  | 20.382  | -1.9 | 92    | 0.00     |
| 88   | 1,2-Dibromo-3-Chloropropane | 20.000  | 21.125  | -5.6 | 103   | 0.00     |
| 89   | 1,2,4-Trichlorobenzene      | 20.000  | 19.258  | 3.7  | 84    | 0.00     |
| 90   | Hexachlorobutadiene         | 20.000  | 19.067  | 4.7  | 85    | 0.00     |
| 91 M | Naphthalene                 | 20.000  | 19.142  | 4.3  | 86    | 0.00     |
| 92   | 1,2,3-Trichlorobenzene      | 20.000  | 19.258  | 3.7  | 84    | 0.00     |

(#) = Out of Range

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