

Data Path : Z:\voasrv\HPCHEM1\MSVOA X\Data\WX051019\  
 Data File : VX009482.D  
 Acq On : 10 May 2019 14:57  
 Operator : JC/SP  
 Sample : VSTDCCC020  
 Misc : 5.0mL/MSVOA X/WATER  
 ALS Vial : 13 Sample Multiplier: 1

Instrument :  
 MSVOA\_X  
 LabSampleId :  
 VSTDCCC020

Quant Time: May 11 03:26:50 2019  
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA X\METHOD\624X050119W.M  
 Quant Title : METHOD 624 VOLATILE ORGANIC ANALYSIS  
 QLast Update : Thu May 02 06:46:57 2019  
 Response via : Initial Calibration

Min. RRF : 0.050 Min. Rel. Area : 20% Max. R.T. Dev 0.50min  
 Max. RRF Dev : 30% Max. Rel. Area : 150%

|      | Compound                    | Amount  | Calc.   | %Dev  | Area% | Dev(min) |
|------|-----------------------------|---------|---------|-------|-------|----------|
| 1 I  | Bromochloromethane          | 30.000  | 30.000  | 0.0   | 89    | 0.00     |
| 2 M  | Dichlorodifluoromethane     | 20.000  | 20.149  | -0.7  | 91    | 0.00     |
| 3 M  | Chloromethane               | 20.000  | 22.292  | -11.5 | 101   | 0.00     |
| 4 M  | Vinyl Chloride              | 20.000  | 21.395  | -7.0  | 96    | 0.00     |
| 5 M  | Bromomethane                | 20.000  | 19.274  | 3.6   | 97    | 0.00     |
| 6 M  | Chloroethane                | 20.000  | 22.296  | -11.5 | 98    | 0.00     |
| 7 M  | Trichlorofluoromethane      | 20.000  | 22.073  | -10.4 | 101   | 0.00     |
| 8 T  | Diethyl Ether               | 20.000  | 21.864  | -9.3  | 100   | 0.00     |
| 9    | 1,1,2-Trichlorotrifluoroeth | 20.000  | 22.096  | -10.5 | 100   | 0.00     |
| 10 M | 1,1-Dichloroethene          | 20.000  | 22.599  | -13.0 | 102   | 0.00     |
| 11   | Methyl Iodide               | 20.000  | 20.248  | -1.2  | 98    | 0.00     |
| 12   | Methyl Acetate              | 20.000  | 19.253  | 3.7   | 89    | 0.00     |
| 13 M | Acrolein                    | 100.000 | 99.031  | 1.0   | 97    | 0.00     |
| 14 M | Acrylonitrile               | 100.000 | 103.955 | -4.0  | 95    | 0.00     |
| 15 M | Acetone                     | 100.000 | 106.804 | -6.8  | 94    | 0.00     |
| 16 M | Carbon Disulfide            | 20.000  | 23.085  | -15.4 | 105   | 0.00     |
| 17   | Allyl chloride              | 20.000  | 20.969  | -4.8  | 96    | 0.00     |
| 18 M | Methylene Chloride          | 20.000  | 21.985  | -9.9  | 101   | 0.00     |
| 19 M | trans-1,2-Dichloroethene    | 20.000  | 22.356  | -11.8 | 102   | 0.00     |
| 20 T | Diisopropyl ether           | 20.000  | 21.617  | -8.1  | 98    | 0.00     |
| 21 M | 1,1-Dichloroethane          | 20.000  | 21.304  | -6.5  | 98    | 0.00     |
| 22 M | cis-1,2-Dichloroethene      | 20.000  | 21.889  | -9.4  | 100   | 0.00     |
| 23 M | tert-Butyl Alcohol          | 100.000 | 93.132  | 6.9   | 88    | 0.00     |
| 24 M | Methyl tert-Butyl Ether     | 20.000  | 21.521  | -7.6  | 100   | 0.00     |
| 25 M | Chloroform                  | 20.000  | 22.313  | -11.6 | 101   | 0.00     |
| 26   | Cyclohexane                 | 20.000  | 22.664  | -13.3 | 102   | 0.00     |
| 27 s | 1,2-Dichloroethane-d4       | 30.000  | 29.647  | 1.2   | 89    | 0.00     |
| 28 I | 1,4-Difluorobenzene         | 30.000  | 30.000  | 0.0   | 89    | 0.00     |
| 29   | 1,1-Dichloropropene         | 20.000  | 22.984  | -14.9 | 103   | 0.00     |
| 30 M | 2-Butanone                  | 100.000 | 105.049 | -5.0  | 94    | 0.00     |
| 31   | 2,2-Dichloropropane         | 20.000  | 21.366  | -6.8  | 97    | 0.00     |
| 32 M | 1,1,1-Trichloroethane       | 20.000  | 22.072  | -10.4 | 102   | 0.00     |
| 33 M | Carbon Tetrachloride        | 20.000  | 22.085  | -10.4 | 101   | 0.00     |
| 34 M | Benzene                     | 20.000  | 22.396  | -12.0 | 100   | 0.00     |
| 35   | Methacrylonitrile           | 20.000  | 21.697  | -8.5  | 97    | 0.00     |
| 36 M | 1,2-Dichloroethane          | 20.000  | 22.456  | -12.3 | 101   | 0.00     |
| 37 M | Trichloroethene             | 20.000  | 21.781  | -8.9  | 100   | 0.00     |
| 38   | Methylcyclohexane           | 20.000  | 23.371  | -16.9 | 106   | 0.00     |
| 39 M | 1,2-Dichloropropane         | 20.000  | 21.872  | -9.4  | 99    | 0.00     |
| 40   | Dibromomethane              | 20.000  | 22.535  | -12.7 | 101   | 0.00     |
| 41 M | Bromodichloromethane        | 20.000  | 21.640  | -8.2  | 99    | 0.00     |
| 42 M | Vinyl Acetate               | 100.000 | 112.306 | -12.3 | 100   | 0.00     |
| 43   | Ethyl Acetate               | 20.000  | 22.041  | -10.2 | 96    | 0.00     |
| 44   | Isopropyl Acetate           | 20.000  | 21.334  | -6.7  | 98    | 0.00     |
| 45 T | 1,4-Dioxane                 | 400.000 | 381.407 | 4.6   | 87    | 0.00     |

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 ALS Vial : 13 Sample Multiplier: 1

Instrument :  
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 LabSampleId :  
 VSTDCCC020

Quant Time: May 11 03:26:50 2019  
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA X\METHOD\624X050119W.M  
 Quant Title : METHOD 624 VOLATILE ORGANIC ANALYSIS  
 QLast Update : Thu May 02 06:46:57 2019  
 Response via : Initial Calibration

Min. RRF : 0.050 Min. Rel. Area : 20% Max. R.T. Dev 0.50min  
 Max. RRF Dev : 30% Max. Rel. Area : 150%

|      | Compound                    | Amount  | Calc.   | %Dev  | Area% | Dev(min) |
|------|-----------------------------|---------|---------|-------|-------|----------|
| 46   | Methyl methacrylate         | 20.000  | 21.728  | -8.6  | 99    | 0.00     |
| 47   | n-amyl Acetate              | 20.000  | 21.434  | -7.2  | 100   | 0.00     |
| 48 M | t-1,3-Dichloropropene       | 20.000  | 21.304  | -6.5  | 100   | 0.00     |
| 49 T | cis-1,3-Dichloropropene     | 20.000  | 21.657  | -8.3  | 100   | 0.00     |
| 50 M | 1,1,2-Trichloroethane       | 20.000  | 21.694  | -8.5  | 97    | 0.00     |
| 51   | Ethyl methacrylate          | 20.000  | 21.459  | -7.3  | 98    | 0.00     |
| 52   | 1,3-Dichloropropane         | 20.000  | 21.960  | -9.8  | 99    | 0.00     |
| 53 M | Dibromochloromethane        | 20.000  | 21.777  | -8.9  | 101   | 0.00     |
| 54 M | 1,2-Dibromoethane           | 20.000  | 22.272  | -11.4 | 101   | 0.00     |
| 55 M | 2-Chloroethyl vinyl ether   | 100.000 | 85.591  | 14.4  | 74    | 0.00     |
| 56 M | Bromoform                   | 20.000  | 20.680  | -3.4  | 99    | 0.00     |
| 57 I | Chlorobenzene-d5            | 30.000  | 30.000  | 0.0   | 91    | 0.00     |
| 58 M | 4-Methyl-2-Pentanone        | 100.000 | 105.786 | -5.8  | 97    | 0.00     |
| 59 M | 2-Hexanone                  | 100.000 | 103.695 | -3.7  | 95    | 0.00     |
| 60 S | 4-Bromofluorobenzene        | 30.000  | 29.278  | 2.4   | 88    | 0.00     |
| 61 M | Tetrachloroethene           | 20.000  | 22.681  | -13.4 | 103   | 0.00     |
| 62 M | Toluene                     | 20.000  | 22.582  | -12.9 | 102   | 0.00     |
| 63 S | Toluene-d8                  | 30.000  | 29.644  | 1.2   | 88    | 0.00     |
| 64 M | Chlorobenzene               | 20.000  | 22.179  | -10.9 | 103   | 0.00     |
| 65   | 1,1,1,2-Tetrachloroethane   | 20.000  | 22.149  | -10.7 | 104   | 0.00     |
| 66 M | Ethyl Benzene               | 20.000  | 22.406  | -12.0 | 103   | 0.00     |
| 67 M | m/p-Xylenes                 | 40.000  | 45.087  | -12.7 | 103   | 0.00     |
| 68 M | o-Xylene                    | 20.000  | 22.180  | -10.9 | 103   | 0.00     |
| 69 M | Styrene                     | 20.000  | 21.838  | -9.2  | 101   | 0.00     |
| 70   | Isopropylbenzene            | 20.000  | 22.376  | -11.9 | 102   | 0.00     |
| 71 M | 1,1,2,2-Tetrachloroethane   | 20.000  | 21.444  | -7.2  | 100   | 0.00     |
| 72   | 1,2,3-Trichloropropane      | 20.000  | 21.655  | -8.3  | 99    | 0.00     |
| 73   | Bromobenzene                | 20.000  | 22.075  | -10.4 | 102   | 0.00     |
| 74   | n-propylbenzene             | 20.000  | 22.369  | -11.8 | 103   | 0.00     |
| 75   | 2-Chlorotoluene             | 20.000  | 22.083  | -10.4 | 102   | 0.00     |
| 76   | 1,3,5-Trimethylbenzene      | 20.000  | 22.375  | -11.9 | 103   | 0.00     |
| 77   | t-1,4-Dichloro-2-butene     | 20.000  | 20.687  | -3.4  | 103   | 0.00     |
| 78   | 4-Chlorotoluene             | 20.000  | 22.194  | -11.0 | 104   | 0.00     |
| 79   | tert-butylbenzene           | 20.000  | 22.352  | -11.8 | 103   | 0.00     |
| 80   | 1,2,4-Trimethylbenzene      | 20.000  | 22.153  | -10.8 | 103   | 0.00     |
| 81   | sec-Butylbenzene            | 20.000  | 22.493  | -12.5 | 104   | 0.00     |
| 82   | p-Isopropyltoluene          | 20.000  | 22.592  | -13.0 | 105   | 0.00     |
| 83 M | 1,3-Dichlorobenzene         | 20.000  | 22.108  | -10.5 | 105   | 0.00     |
| 84 M | 1,4-Dichlorobenzene         | 20.000  | 21.851  | -9.3  | 105   | 0.00     |
| 85   | n-Butylbenzene              | 20.000  | 22.247  | -11.2 | 107   | 0.00     |
| 86 T | Hexachloroethane            | 20.000  | 21.223  | -6.1  | 99    | 0.00     |
| 87 M | 1,2-Dichlorobenzene         | 20.000  | 22.055  | -10.3 | 103   | 0.00     |
| 88   | 1,2-Dibromo-3-Chloropropane | 20.000  | 20.810  | -4.0  | 98    | 0.00     |
| 89   | 1,2,4-Trichlorobenzene      | 20.000  | 22.158  | -10.8 | 108   | 0.00     |
| 90   | Hexachlorobutadiene         | 20.000  | 23.289  | -16.4 | 111   | 0.00     |

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QLast Update : Thu May 02 06:46:57 2019  
Response via : Initial Calibration

Min. RRF : 0.050 Min. Rel. Area : 20% Max. R.T. Dev 0.50min  
Max. RRF Dev : 30% Max. Rel. Area : 150%

|      | Compound               | Amount | Calc.  | %Dev  | Area% | Dev(min) |
|------|------------------------|--------|--------|-------|-------|----------|
| 91 M | Naphthalene            | 20.000 | 21.837 | -9.2  | 101   | 0.00     |
| 92   | 1,2,3-Trichlorobenzene | 20.000 | 22.639 | -13.2 | 106   | 0.00     |

(#) = Out of Range

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