

Data Path : Z:\voasrv\HPCHEM1\MSVOA\_L\Data\VL061422\  
 Data File : VL039297.D  
 Acq On : 15 Jun 2022 13:07  
 Operator : SY/AP  
 Sample : VSTDCCC010  
 Misc : 400mL/MSVOA\_L  
 ALS Vial : 1 Sample Multiplier: 1

Instrument :  
 MSVOA\_L  
 LabSampleId :  
 VSTDCCC010

Quant Time: Jun 16 01:17:37 2022  
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA\_L\METHODS\VL053122AIR.M  
 Quant Title : AIR ANALYSIS BY METHOD TO-15 Instrument: MSVOA\_LWed Jun 01 01:22:18 2022  
 QLast Update : Wed Jun 01 01:22:18 2022  
 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                    | AvgRF | CCRF   | %Dev  | Area% | Dev(min) |
|------|-----------------------------|-------|--------|-------|-------|----------|
| 1 I  | Bromochloromethane          | 1.000 | 1.000  | 0.0   | 98    | -0.02    |
| 2 T  | Dichlorodifluoromethane     | 1.756 | 1.389  | 20.9  | 91    | -0.02    |
| 3    | Chlorodifluoromethane       | 1.272 | 1.066  | 16.2  | 92    | -0.01    |
| 4    | Chloromethane               | 0.654 | 0.512  | 21.7  | 84    | -0.01    |
| 5 T  | Vinyl Chloride              | 0.611 | 0.547  | 10.5  | 92    | -0.01    |
| 6 T  | Bromomethane                | 0.301 | 0.257  | 14.6  | 94    | -0.01    |
| 7    | Chloroethane                | 0.213 | 0.202  | 5.2   | 99    | -0.02    |
| 8 T  | Dichlorotetrafluoroethane   | 1.605 | 1.354  | 15.6  | 91    | -0.02    |
| 9 T  | Propene                     | 0.555 | 0.460  | 17.1  | 93    | -0.01    |
| 10 T | Heptane                     | 1.294 | 1.221  | 5.6   | 98    | -0.02    |
| 11 T | Trichlorofluoromethane      | 1.590 | 1.386  | 12.8  | 97    | -0.02    |
| 12 T | 1,1,2-Trichlorotrifluoroeth | 1.051 | 1.029  | 2.1   | 105   | -0.02    |
| 13   | Ethanol                     | 0.030 | 0.030# | 0.0   | 106   | -0.02    |
| 14 T | Bromoethene                 | 0.473 | 0.408  | 13.7  | 90    | 0.00     |
| 15 T | Acetone                     | 1.042 | 1.134  | -8.8  | 122   | -0.02    |
| 16 T | 1,3-Butadiene               | 0.560 | 0.539  | 3.8   | 101   | -0.01    |
| 17   | tert-Butyl alcohol          | 0.808 | 0.935  | -15.7 | 115   | -0.01    |
| 18 T | 1,1-Dichloroethene          | 0.472 | 0.454  | 3.8   | 102   | -0.01    |
| 19 T | Isopropyl Alcohol           | 0.489 | 0.502  | -2.7  | 101   | 0.00     |
| 20 T | Methylene Chloride          | 0.439 | 0.375  | 14.6  | 110   | -0.01    |
| 21 T | Allyl Chloride              | 0.703 | 0.686  | 2.4   | 102   | -0.02    |
| 22 T | trans-1,2-Dichloroethene    | 0.454 | 0.439  | 3.3   | 98    | -0.02    |
| 23 T | Vinyl Acetate               | 0.862 | 0.715  | 17.1  | 78    | -0.02    |
| 24 T | 1,1-Dichloroethane          | 1.181 | 1.009  | 14.6  | 94    | -0.01    |
| 25 T | Ethyl Acetate               | 2.348 | 2.355  | -0.3  | 107   | -0.02    |
| 26 T | Hexane                      | 1.032 | 0.917  | 11.1  | 95    | -0.01    |
| 27 T | Carbon Disulfide            | 1.069 | 1.060  | 0.8   | 97    | -0.02    |
| 28 T | Methyl tert-Butyl Ether     | 1.139 | 0.996  | 12.6  | 96    | -0.02    |
| 29 T | Chloroform                  | 1.736 | 1.428  | 17.7  | 88    | -0.02    |
| 30 T | Cyclohexane                 | 0.867 | 0.776  | 10.5  | 97    | -0.02    |
| 31 T | cis-1,2-Dichloroethene      | 1.052 | 0.898  | 14.6  | 89    | -0.01    |
| 32 T | 1,1,1-Trichloroethane       | 1.699 | 1.446  | 14.9  | 87    | -0.02    |
| 33 I | 1,4-Difluorobenzene         | 1.000 | 1.000  | 0.0   | 110   | -0.02    |
| 34 T | 2-Butanone                  | 0.543 | 0.451  | 16.9  | 92    | -0.02    |
| 35 T | Carbon Tetrachloride        | 0.730 | 0.545  | 25.3  | 85    | -0.02    |
| 36 T | Benzene                     | 0.812 | 0.679  | 16.4  | 97    | -0.02    |
| 37 T | 1,2-Dichloroethane          | 0.534 | 0.395  | 26.0  | 84    | -0.02    |
| 38 T | Trichloroethene             | 0.373 | 0.297  | 20.4  | 86    | -0.02    |
| 39 T | 1,2-Dichloropropane         | 0.323 | 0.254  | 21.4  | 92    | -0.02    |
| 40 T | 1,4-Dioxane                 | 0.106 | 0.091  | 14.2  | 97    | -0.02    |
| 41 T | Tetrahydrofuran             | 0.327 | 0.256  | 21.7  | 89    | -0.02    |
| 42 T | Bromodichloromethane        | 0.742 | 0.559  | 24.7  | 86    | -0.02    |
| 43   | Methyl Methacrylate         | 0.316 | 0.292  | 7.6   | 99    | -0.02    |
| 44 T | 2,2,4-Trimethylpentane      | 1.420 | 1.153  | 18.8  | 96    | -0.02    |
| 45 T | t-1,3-Dichloropropene       | 0.329 | 0.220  | 33.1# | 63    | -0.02    |
| 46 T | cis-1,3-Dichloropropene     | 0.435 | 0.314  | 27.8  | 73    | -0.02    |
| 47 T | 1,1,2-Trichloroethane       | 0.337 | 0.290  | 13.9  | 98    | -0.01    |
| 48 T | Dibromochloromethane        | 0.602 | 0.316  | 47.5# | 57    | -0.01    |

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 LabSampleId :  
 VSTDCCC010

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 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                  | AvgRF | CCRF  | %Dev  | Area% | Dev(min) |
|------|---------------------------|-------|-------|-------|-------|----------|
| 49 T | Bromoform                 | 0.501 | 0.419 | 16.4  | 90    | -0.02    |
| 50 T | 4-Methyl-2-Pentanone      | 0.813 | 0.723 | 11.1  | 97    | -0.02    |
| 51 T | 2-Hexanone                | 0.633 | 0.467 | 26.2  | 81    | -0.02    |
| 52 T | Tetrachloroethene         | 0.317 | 0.271 | 14.5  | 101   | -0.02    |
| 53 T | Toluene                   | 0.935 | 0.815 | 12.8  | 99    | -0.02    |
| 54 T | 1,2-Dibromoethane         | 0.453 | 0.294 | 35.1# | 72    | -0.02    |
| 55 I | Chlorobenzene-d5          | 1.000 | 1.000 | 0.0   | 109   | -0.02    |
| 56   | 1,1,1,2-Tetrachloroethane | 0.473 | 0.402 | 15.0  | 102   | -0.02    |
| 57 T | Chlorobenzene             | 0.752 | 0.659 | 12.4  | 105   | -0.02    |
| 58 T | Ethyl Benzene             | 1.377 | 1.171 | 15.0  | 101   | -0.02    |
| 59 T | m/p-Xylene                | 1.213 | 1.005 | 17.1  | 99    | -0.02    |
| 60 T | o-Xylene                  | 1.140 | 0.967 | 15.2  | 102   | -0.02    |
| 61 T | Styrene                   | 0.610 | 0.565 | 7.4   | 104   | -0.02    |
| 62   | Isopropylbenzene          | 1.694 | 1.519 | 10.3  | 109   | -0.02    |
| 63 T | 1,1,2,2-Tetrachloroethane | 0.640 | 0.615 | 3.9   | 120   | -0.02    |
| 64   | n-propylbenzene           | 0.416 | 0.401 | 3.6   | 112   | -0.02    |
| 65   | tert-Butylbenzene         | 1.549 | 1.393 | 10.1  | 110   | -0.02    |
| 66 T | Benzyl Chloride           | 0.201 | 0.089 | 55.7# | 40    | -0.02    |
| 67   | sec-Butylbenzene          | 2.114 | 1.911 | 9.6   | 109   | -0.02    |
| 68 S | 1-Bromo-4-Fluorobenzene   | 0.827 | 0.781 | 5.6   | 105   | -0.02    |
| 69   | p-Isopropyltoluene        | 1.783 | 1.664 | 6.7   | 113   | -0.02    |
| 70   | n-Butylbenzene            | 1.830 | 1.644 | 10.2  | 105   | -0.02    |
| 71   | 2-Chlorotoluene           | 1.292 | 1.124 | 13.0  | 105   | -0.02    |
| 72 T | 4-Ethyltoluene            | 1.336 | 1.216 | 9.0   | 107   | -0.02    |
| 73 T | 1,3,5-Trimethylbenzene    | 1.205 | 1.093 | 9.3   | 106   | -0.01    |
| 74 T | 1,2,4-Trimethylbenzene    | 1.345 | 1.192 | 11.4  | 106   | -0.02    |
| 75 T | 1,3-Dichlorobenzene       | 0.761 | 0.708 | 7.0   | 109   | -0.02    |
| 76 T | 1,4-Dichlorobenzene       | 0.759 | 0.683 | 10.0  | 108   | -0.02    |
| 77 T | 1,2-Dichlorobenzene       | 0.767 | 0.696 | 9.3   | 110   | -0.02    |
| 78 T | Hexachloro-1,3-Butadiene  | 0.701 | 0.560 | 20.1  | 103   | -0.02    |
| 79 T | Naphthalene               | 1.035 | 1.029 | 0.6   | 99    | -0.02    |
| 80 T | Naphthalene,2-methyl-     | 0.385 | 0.258 | 33.0# | 58    | -0.01    |
| 81 T | 1,2,4-Trichlorobenzene    | 0.656 | 0.588 | 10.4  | 101   | -0.02    |

(#) = Out of Range

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