

Data Path : Z:\voasrv\HPCHEM1\MSVOA\_L\Data\VL101722\  
 Data File : VL039722.D  
 Acq On : 17 Oct 2022 16:48  
 Operator : SY/AP  
 Sample : N5169-03DUP  
 Misc : 400mL/MSVOA\_L  
 ALS Vial : 1 Sample Multiplier: 1

Instrument :  
 MSVOA\_L  
 ClientSampleId :  
 IA1DUP

Manual Integrations  
 APPROVED

Reviewed By :Semsettin Yesilyurt 10/18/2022  
 Supervised By :Mahesh Dadoda 10/19/2022

Quant Time: Oct 18 05:27:32 2022

Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_L\methods\VL091922AIR.M

Quant Title : AIR ANALYSIS BY METHOD T0-15 Instrument: MSVOA\_LFri Aug 2

QLast Update : Tue Oct 18 05:23:57 2022

Response via : Initial Calibration

| Compound                    | R.T.   | QIon  | Response | Conc     | Units  | Dev(Min) |
|-----------------------------|--------|-------|----------|----------|--------|----------|
| -----                       |        |       |          |          |        |          |
| Internal Standards          |        |       |          |          |        |          |
| 1) Bromochloromethane       | 5.230  | 49    | 1017320  | 10.000   | ppbv   | 0.00     |
| 33) 1,4-Difluorobenzene     | 6.687  | 114   | 2621089  | 10.000   | ppbv   | 0.00     |
| 55) Chlorobenzene-d5        | 11.497 | 117   | 2213483  | 10.000   | ppbv   | 0.00     |
|                             |        |       |          |          |        |          |
| System Monitoring Compounds |        |       |          |          |        |          |
| 68) 1-Bromo-4-Fluorobenzene | 13.809 | 95    | 1672812  | 10.124   | ppbv   | 0.00     |
| Spiked Amount               | 10.000 | Range | 65 - 135 | Recovery | =      | 101.200% |
|                             |        |       |          |          |        |          |
| Target Compounds            |        |       |          |          |        |          |
|                             |        |       |          |          | Qvalue |          |
| 2) Dichlorodifluoromethane  | 2.768  | 85    | 20158m   | 0.171    | ppbv   |          |
| 3) Chlorodifluoromethane    | 2.716  | 51    | 44678m   | 0.370    | ppbv   |          |
| 4) Chloromethane            | 2.858  | 50    | 34464m   | 0.473    | ppbv   |          |
| 9) Propene                  | 2.742  | 41    | 140221m  | 2.543    | ppbv   |          |
| 10) Heptane                 | 7.619  | 43    | 81458m   | 0.633    | ppbv   |          |
| 11) Trichlorofluoromethane  | 3.603  | 101   | 29795    | 0.204    | ppbv # | 80       |
| 13) Ethanol                 | 3.276  | 45    | 230000m  | 54.747   | ppbv   |          |
| 15) Acetone                 | 3.507  | 43    | 674438m  | 5.958    | ppbv   |          |
| 19) Isopropyl Alcohol       | 3.623  | 45    | 210006   | 3.639    | ppbv # | 1        |
| 20) Methylene Chloride      | 3.976  | 84    | 140093   | 2.677    | ppbv   | 98       |
| 26) Hexane                  | 5.250  | 57    | 221316   | 2.113    | ppbv # | 54       |
| 29) Chloroform              | 5.314  | 83    | 24544m   | 0.162    | ppbv   |          |
| 30) Cyclohexane             | 6.642  | 84    | 15137m   | 0.187    | ppbv   |          |
| 34) 2-Butanone              | 4.835  | 43    | 70664m   | 0.549    | ppbv   |          |
| 35) Carbon Tetrachloride    | 6.536  | 117   | 8445m    | 0.061    | ppbv   |          |
| 36) Benzene                 | 6.404  | 78    | 150465   | 0.751    | ppbv   | 97       |
| 44) 2,2,4-Trimethylpentane  | 7.365  | 57    | 113967m  | 0.330    | ppbv   |          |
| 52) Tetrachloroethene       | 10.648 | 164   | 2054074  | 26.868   | ppbv   | 96       |
| 53) Toluene                 | 9.256  | 91    | 840887   | 3.773    | ppbv   | 100      |
| 58) Ethyl Benzene           | 12.121 | 91    | 288571   | 1.131    | ppbv   | 98       |
| 59) m/p-Xylene              | 12.375 | 91    | 1063700  | 4.683    | ppbv   | 96       |
| 60) o-Xylene                | 13.092 | 91    | 272523   | 1.256    | ppbv   | 98       |
| 72) 4-Ethyltoluene          | 15.220 | 105   | 33704m   | 0.136    | ppbv   |          |
| 73) 1,3,5-Trimethylbenzene  | 15.384 | 105   | 25379m   | 0.112    | ppbv   |          |
| 74) 1,2,4-Trimethylbenzene  | 16.134 | 105   | 92285m   | 0.363    | ppbv   |          |
| -----                       |        |       |          |          |        |          |

(#) = qualifier out of range (m) = manual integration (+) = signals summed

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