

Data Path : Z:\VOASRV\HPCHEM1\MSVOA L\DATA\VL051322\  
 Data File : VL039053.D  
 Acq On : 14 May 2022 1:54  
 Operator : SY/AP  
 Sample : N2828-06  
 Misc : 400mL/MSVOA L  
 ALS Vial : 1 Sample Multiplier: 1

Instrument :  
 MSVOA L  
 ClientSampleId :  
 OA1

Manual Integrations  
 APPROVED

Reviewed By :Semsettin Yesilyurt 05/16/2022  
 Supervised By :Mahesh Dadoda 05/17/2022

Quant Time: May 14 03:58:59 2022  
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA L\METHODS\VL050522A  
 Quant Title : AIR ANALYSIS BY METHOD TO-15 Instrument:  
 QLast Update : Thu May 05 10:12:15 2022  
 Response via : Initial Calibration

| Internal Standards      | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|-------------------------|-------|------|----------|-------|-------|----------|
| 1) Bromochloromethane   | 5.65  | 49   | 880885   | 10.00 | ppbv  | -0.01    |
| 33) 1,4-Difluorobenzene | 7.16  | 114  | 2097451  | 10.00 | ppbv  | -0.02    |
| 55) Chlorobenzene-d5    | 12.06 | 117  | 1694469  | 10.00 | ppbv  | -0.01    |

## System Monitoring Compounds

|                             |        |       |          |          |      |         |
|-----------------------------|--------|-------|----------|----------|------|---------|
| 68) 1-Bromo-4-Fluorobenzene | 14.39  | 95    | 1355608  | 10.10    | ppbv | 0.00    |
| Spiked Amount               | 10.000 | Range | 65 - 135 | Recovery | =    | 101.00% |

## Target Compounds

| Target Compounds           | R.T.  | QIon | Response | Conc    | Units  | Ovalue |
|----------------------------|-------|------|----------|---------|--------|--------|
| 2) Dichlorodifluoromethane | 3.05  | 85   | 44646    | 0.323   | ppbv   | 91     |
| 3) Chlorodifluoromethane   | 2.99  | 51   | 50029    | 0.458   | ppbv   | 96     |
| 4) Chloromethane           | 3.13  | 50   | 28743    | 0.525   | ppbv   | 94     |
| 9) Propene                 | 3.01  | 41   | 165403   | 3.465   | ppbv   | 87     |
| 10) Heptane                | 8.10  | 43   | 82133    | 0.701   | ppbv   | 90     |
| 11) Trichlorofluoromethane | 3.94  | 101  | 29071    | 0.245   | ppbv   | 97     |
| 13) Ethanol                | 3.59  | 45   | 467421   | 138.368 | ppbv   | 94     |
| 15) Acetone                | 3.85  | 43   | 343838m  | 4.138   | ppbv   |        |
| 19) Isopropyl Alcohol      | 3.96  | 45   | 68301    | 1.729   | ppbv # | 84     |
| 20) Methylene Chloride     | 4.34  | 84   | 702041   | 22.417  | ppbv   | 96     |
| 26) Hexane                 | 5.67  | 57   | 926065   | 10.480  | ppbv # | 39     |
| 30) Cyclohexane            | 7.12  | 84   | 43364    | 0.586   | ppbv   | 94     |
| 34) 2-Butanone             | 5.23  | 43   | 35374m   | 0.331   | ppbv   |        |
| 35) Carbon Tetrachloride   | 7.01  | 117  | 9127m    | 0.066   | ppbv   |        |
| 36) Benzene                | 6.87  | 78   | 313079   | 1.744   | ppbv   | 98     |
| 41) Tetrahydrofuran        | 6.04  | 42   | 11324m   | 0.157   | ppbv   |        |
| 44) 2,2,4-Trimethylpentane | 7.85  | 57   | 384757   | 1.222   | ppbv # | 89     |
| 52) Tetrachloroethene      | 11.20 | 164  | 50312    | 0.784   | ppbv   | 96     |
| 53) Toluene                | 9.79  | 91   | 748792   | 3.722   | ppbv   | 100    |
| 58) Ethyl Benzene          | 12.69 | 91   | 93743    | 0.384   | ppbv   | 95     |
| 59) m/p-Xylene             | 12.94 | 91   | 323412m  | 1.554   | ppbv   |        |
| 60) o-Xylene               | 13.67 | 91   | 120592   | 0.615   | ppbv   | 95     |
| 70) n-Butylbenzene         | 18.54 | 91   | 34533m   | 0.112   | ppbv   |        |
| 72) 4-Ethyltoluene         | 15.81 | 105  | 49802    | 0.223   | ppbv # | 90     |
| 73) 1,3,5-Trimethylbenzene | 15.97 | 105  | 52351m   | 0.260   | ppbv   |        |
| 74) 1,2,4-Trimethylbenzene | 16.74 | 105  | 210490   | 0.956   | ppbv # | 66     |
| 79) Naphthalene            | 22.17 | 128  | 24521m   | 0.137   | ppbv   |        |

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

