

Data Path : Z:\VOASRV\HPCHEM1\MSVOA\_L\DATA\VL100721\  
 Data File : VL037818.D  
 Acq On : 8 Oct 2021 5:35  
 Operator : SY/AP  
 Sample : M4079-03  
 Misc : 400mL/MSVOA\_L  
 ALS Vial : 1 Sample Multiplier: 1

Instrument :  
 MSVOA\_L  
 ClientSampleId :  
 C0P13

Manual Integrations  
 APPROVED

apatel  
 10/14/2021 1:19:15 PM

Quant Time: Oct 08 11:32:28 2021  
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA\_L\METHODS\VL100421AIR.M  
 Quant Title : AIR ANALYSIS BY METHOD TO-15 Instrument: MSVOA\_L Mon Oct 0  
 QLast Update : Mon Oct 04 17:43:54 2021  
 Response via : Initial Calibration

| Internal Standards      | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|-------------------------|-------|------|----------|-------|-------|----------|
| 1) Bromochloromethane   | 5.65  | 49   | 1179892  | 10.00 | ppbv  | -0.01    |
| 33) 1,4-Difluorobenzene | 7.17  | 114  | 3557292  | 10.00 | ppbv  | 0.00     |
| 55) Chlorobenzene-d5    | 12.07 | 117  | 3118520  | 10.00 | ppbv  | -0.02    |

| System Monitoring Compounds |        | R.T.  | QIon     | Response | Conc | Units   | Dev(Min) |
|-----------------------------|--------|-------|----------|----------|------|---------|----------|
| 68) 1-Bromo-4-Fluorobenzene | 14.40  | 95    | 2181851  | 10.00    | ppbv | -0.02   |          |
| Spiked Amount               | 10.000 | Range | 65 - 135 | Recovery | =    | 100.00% |          |

| Target Compounds           | R.T.  | QIon | Response | Conc   | Units  | Qvalue |
|----------------------------|-------|------|----------|--------|--------|--------|
| 2) Dichlorodifluoromethane | 3.04  | 85   | 34567m   | 0.183  | ppbv   |        |
| 3) Chlorodifluoromethane   | 2.98  | 51   | 48603    | 0.214  | ppbv   | 99     |
| 4) Chloromethane           | 3.13  | 50   | 25024    | 0.301  | ppbv # | 86     |
| 9) Propene                 | 3.00  | 41   | 44951m   | 0.539  | ppbv   |        |
| 11) Trichlorofluoromethane | 3.94  | 101  | 40096    | 0.238  | ppbv # | 72     |
| 13) Ethanol                | 3.60  | 45   | 98352m   | 14.417 | ppbv   |        |
| 15) Acetone                | 3.84  | 43   | 768548   | 7.256  | ppbv # | 88     |
| 17) tert-Butyl alcohol     | 4.31  | 59   | 17466m   | 0.160  | ppbv   |        |
| 19) Isopropyl Alcohol      | 3.98  | 45   | 25709m   | 0.390  | ppbv   |        |
| 20) Methylene Chloride     | 4.34  | 84   | 209894   | 3.922  | ppbv   | 98     |
| 26) Hexane                 | 5.68  | 57   | 141702   | 0.968  | ppbv # | 45     |
| 29) Chloroform             | 5.74  | 83   | 133491   | 0.638  | ppbv   | 89     |
| 34) 2-Butanone             | 5.24  | 43   | 152874   | 1.065  | ppbv # | 89     |
| 35) Carbon Tetrachloride   | 7.02  | 117  | 8122m    | 0.040  | ppbv   |        |
| 36) Benzene                | 6.87  | 78   | 34252m   | 0.113  | ppbv   |        |
| 52) Tetrachloroethene      | 11.21 | 164  | 1367697  | 11.570 | ppbv   | 98     |
| 53) Toluene                | 9.79  | 91   | 48015m   | 0.143  | ppbv   |        |
| 58) Ethyl Benzene          | 12.70 | 91   | 60312m   | 0.144  | ppbv   |        |
| 59) m/p-Xylene             | 12.95 | 91   | 212763m  | 0.592  | ppbv   |        |
| 60) o-Xylene               | 13.67 | 91   | 63857m   | 0.185  | ppbv   |        |
| 74) 1,2,4-Trimethylbenzene | 16.75 | 105  | 57440m   | 0.154  | ppbv   |        |
| 79) Naphthalene            | 22.18 | 128  | 56202m   | 0.246  | ppbv   |        |

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

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