

Data Path : Z:\VOASRV\HPCHEM1\MSVOA\_L\DATA\VL100721\  
 Data File : VL037820.D  
 Acq On : 8 Oct 2021 6:54  
 Operator : SY/AP  
 Sample : M4079-12  
 Misc : 400mL/MSVOA\_L  
 ALS Vial : 1 Sample Multiplier: 1

Instrument :  
 MSVOA\_L  
 ClientSampleId :  
 C0P12

Manual Integrations  
 APPROVED

apatel  
 10/14/2021 1:19:12 PM

Quant Time: Oct 08 11:36:47 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   | 1260302  | 10.00 | ppbv  | -0.01    |
| 33) 1,4-Difluorobenzene | 7.16  | 114  | 3648270  | 10.00 | ppbv  | -0.01    |
| 55) Chlorobenzene-d5    | 12.07 | 117  | 3026261  | 10.00 | ppbv  | -0.02    |

System Monitoring Compounds

|                             |        |       |          |          |      |         |
|-----------------------------|--------|-------|----------|----------|------|---------|
| 68) 1-Bromo-4-Fluorobenzene | 14.40  | 95    | 2175343  | 10.28    | ppbv | -0.02   |
| Spiked Amount               | 10.000 | Range | 65 - 135 | Recovery | =    | 102.80% |

| Target Compounds           | R.T.  | QIon | Response | Conc  | Units | Qvalue |
|----------------------------|-------|------|----------|-------|-------|--------|
| 2) Dichlorodifluoromethane | 3.04  | 85   | 27288    | 0.136 | ppbv  | 98     |
| 3) Chlorodifluoromethane   | 2.98  | 51   | 23025    | 0.095 | ppbv  | 96     |
| 4) Chloromethane           | 3.13  | 50   | 9194m    | 0.104 | ppbv  |        |
| 9) Propene                 | 3.01  | 41   | 15424m   | 0.173 | ppbv  |        |
| 11) Trichlorofluoromethane | 3.94  | 101  | 36074m   | 0.201 | ppbv  |        |
| 13) Ethanol                | 3.60  | 45   | 69597m   | 9.551 | ppbv  |        |
| 15) Acetone                | 3.84  | 43   | 707298   | 6.252 | ppbv  | 96     |
| 17) tert-Butyl alcohol     | 4.30  | 59   | 31110m   | 0.267 | ppbv  |        |
| 19) Isopropyl Alcohol      | 3.97  | 45   | 38096m   | 0.541 | ppbv  |        |
| 20) Methylene Chloride     | 4.34  | 84   | 103186   | 1.805 | ppbv  | 91     |
| 26) Hexane                 | 5.68  | 57   | 138848m  | 0.888 | ppbv  |        |
| 27) Carbon Disulfide       | 4.55  | 76   | 22285m   | 0.161 | ppbv  |        |
| 29) Chloroform             | 5.74  | 83   | 24993m   | 0.112 | ppbv  |        |
| 34) 2-Butanone             | 5.25  | 43   | 72778m   | 0.494 | ppbv  |        |
| 52) Tetrachloroethene      | 11.20 | 164  | 150060   | 1.238 | ppbv  | 92     |
| 53) Toluene                | 9.79  | 91   | 94876m   | 0.275 | ppbv  |        |
| 59) m/p-Xylene             | 12.95 | 91   | 37333m   | 0.107 | ppbv  |        |

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

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