

Data Path : Z:\VOASRV\HPCHEM1\MSVOA L\DATA\VL021820\  
 Data File : VL034629.D  
 Acq On : 18 Feb 2020 18:21  
 Operator : SY/AP  
 Sample : L1532-05  
 Misc : 400ml/MSVOA L  
 ALS Vial : 1 Sample Multiplier: 1

Instrument :  
 MSVOA\_L  
 ClientSampleId :  
 IR-PR-IA

Manual Integrations  
 APPROVED

MMDadoda  
 2/21/2020 3:25:53 PM

Quant Time: Feb 19 06:48:55 2020  
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA L\METHODS\VL013020AIR.M  
 Quant Title : AIR ANALYSIS BY METHOD TO-15 Instrument: MSVOA\_L Thu Jan 30 2020  
 QLast Update : Thu Jan 30 17:16:44 2020  
 Response via : Initial Calibration

| Internal Standards      | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|-------------------------|-------|------|----------|-------|-------|----------|
| 1) Bromochloromethane   | 5.65  | 49   | 901264   | 10.00 | ppbv  | -0.01    |
| 33) 1,4-Difluorobenzene | 7.17  | 114  | 1854668  | 10.00 | ppbv  | -0.01    |
| 55) Chlorobenzene-d5    | 12.11 | 117  | 1848130  | 10.00 | ppbv  | 0.00     |

|                             |        |       |          |          |      |         |
|-----------------------------|--------|-------|----------|----------|------|---------|
| System Monitoring Compounds |        |       |          |          |      |         |
| 68) 1-Bromo-4-Fluorobenzene | 14.45  | 95    | 1599004  | 10.24    | ppbv | 0.00    |
| Spiked Amount               | 10.000 | Range | 65 - 135 | Recovery | =    | 102.40% |

| Target Compounds           | R.T.  | QIon | Response | Conc   | Units  | Ovalue |
|----------------------------|-------|------|----------|--------|--------|--------|
| 2) Dichlorodifluoromethane | 3.01  | 85   | 47007    | 0.237  | ppbv   | 89     |
| 3) Chlorodifluoromethane   | 2.95  | 51   | 35778m   | 0.329  | ppbv   |        |
| 4) Chloromethane           | 3.10  | 50   | 29767    | 0.447  | ppbv   | 92     |
| 9) Propene                 | 2.98  | 41   | 33964    | 0.801  | ppbv # | 72     |
| 11) Trichlorofluoromethane | 3.92  | 101  | 43958    | 0.244  | ppbv   | 89     |
| 13) Ethanol                | 3.57  | 45   | 913256   | 80.409 | ppbv   | 99     |
| 15) Acetone                | 3.82  | 43   | 616348   | 4.653  | ppbv   | 100    |
| 19) Isopropyl Alcohol      | 3.94  | 45   | 55129    | 0.713  | ppbv # | 1      |
| 20) Methylene Chloride     | 4.32  | 84   | 88137    | 1.747  | ppbv   | 92     |
| 26) Hexane                 | 5.67  | 57   | 100203   | 1.116  | ppbv # | 44     |
| 29) Chloroform             | 5.74  | 83   | 16568m   | 0.104  | ppbv   |        |
| 35) Carbon Tetrachloride   | 7.02  | 117  | 12461m   | 0.070  | ppbv   |        |
| 36) Benzene                | 6.89  | 78   | 20148m   | 0.122  | ppbv   |        |
| 52) Tetrachloroethene      | 11.24 | 164  | 11703m   | 0.177  | ppbv   |        |
| 53) Toluene                | 9.82  | 91   | 497055   | 2.743  | ppbv   | 95     |

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

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