

Data Path : Z:\VOASRV\HPCHEM1\MSVOA\_L\DATA\VL092721\  
 Data File : VL037683.D  
 Acq On : 27 Sep 2021 16:43  
 Operator : SY/AP  
 Sample : M3896-02  
 Misc : 400mL/MSVOA\_L  
 ALS Vial : 7 Sample Multiplier: 1

Instrument :  
 MSVOA\_L  
 ClientSampleId :  
 MID

Manual Integrations  
 APPROVED

MMDadoda  
 10/1/2021 11:05:04 AM

Quant Time: Sep 28 20:12:53 2021  
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA\_L\METHODS\VL091621AIR.M  
 Quant Title : AIR ANALYSIS BY METHOD T0-15 Instrument: MSVOA\_L Thu Sep 1  
 QLast Update : Thu Sep 16 13:42:46 2021  
 Response via : Initial Calibration

| Internal Standards      | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|-------------------------|-------|------|----------|-------|-------|----------|
| 1) Bromochloromethane   | 5.67  | 49   | 1522362  | 10.00 | ppbv  | 0.02     |
| 33) 1,4-Difluorobenzene | 7.19  | 114  | 4367710  | 10.00 | ppbv  | 0.03     |
| 55) Chlorobenzene-d5    | 12.10 | 117  | 3853007  | 10.00 | ppbv  | 0.03     |

| System Monitoring Compounds |        |       |          |          |      |        |
|-----------------------------|--------|-------|----------|----------|------|--------|
| 68) 1-Bromo-4-Fluorobenzene | 14.42  | 95    | 2681330  | 9.63     | ppbv | 0.03   |
| Spiked Amount               | 10.000 | Range | 65 - 135 | Recovery | =    | 96.30% |

| Target Compounds           | R.T.  | QIon | Response | Conc   | Units  | Qvalue |
|----------------------------|-------|------|----------|--------|--------|--------|
| 2) Dichlorodifluoromethane | 3.06  | 85   | 38757    | 0.141  | ppbv   | 95     |
| 4) Chloromethane           | 3.16  | 50   | 31838m   | 0.300  | ppbv   |        |
| 9) Propene                 | 3.03  | 41   | 42742m   | 0.422  | ppbv   |        |
| 10) Heptane                | 8.13  | 43   | 75851m   | 0.321  | ppbv   |        |
| 11) Trichlorofluoromethane | 3.96  | 101  | 55786    | 0.252  | ppbv   | 93     |
| 13) Ethanol                | 3.62  | 45   | 159292   | 16.078 | ppbv   | 92     |
| 15) Acetone                | 3.85  | 43   | 3737119  | 23.627 | ppbv   | 99     |
| 18) 1,1-Dichloroethene     | 4.29  | 96   | 8530m    | 0.112  | ppbv   |        |
| 19) Isopropyl Alcohol      | 3.98  | 45   | 828170   | 9.623  | ppbv # | 92     |
| 20) Methylene Chloride     | 4.35  | 84   | 371744   | 4.923  | ppbv   | 95     |
| 24) 1,1-Dichloroethane     | 5.01  | 63   | 60063m   | 0.383  | ppbv   |        |
| 26) Hexane                 | 5.69  | 57   | 349440   | 1.920  | ppbv # | 31     |
| 29) Chloroform             | 5.76  | 83   | 28466m   | 0.107  | ppbv   |        |
| 32) 1,1,1-Trichloroethane  | 6.52  | 97   | 867260   | 3.506  | ppbv   | 98     |
| 34) 2-Butanone             | 5.26  | 43   | 479551   | 2.702  | ppbv   | 92     |
| 35) Carbon Tetrachloride   | 7.03  | 117  | 8320m    | 0.034  | ppbv   |        |
| 41) Tetrahydrofuran        | 6.07  | 42   | 93178m   | 1.117  | ppbv   |        |
| 51) 2-Hexanone             | 10.15 | 43   | 167597m  | 1.686  | ppbv   |        |
| 53) Toluene                | 9.82  | 91   | 361515m  | 0.893  | ppbv   |        |
| 58) Ethyl Benzene          | 12.72 | 91   | 182154   | 0.356  | ppbv   | 99     |
| 59) m/p-Xylene             | 12.98 | 91   | 445187m  | 0.999  | ppbv   |        |
| 60) o-Xylene               | 13.70 | 91   | 246337m  | 0.581  | ppbv   |        |
| 61) Styrene                | 13.54 | 104  | 96130m   | 0.460  | ppbv   |        |
| 74) 1,2,4-Trimethylbenzene | 16.76 | 105  | 108721m  | 0.227  | ppbv   |        |

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

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