

Data Path : Z:\VOASRV\HPCHEM1\MSVOA L\DATA\VL080421\  
 Data File : VL037336.D  
 Acq On : 5 Aug 2021 9:57  
 Operator : SY/AP  
 Sample : M3239-02  
 Misc : 400mL/MSVOA L  
 ALS Vial : 1 Sample Multiplier: 1

Instrument :  
 MSVOA\_L  
 ClientSampleId :  
 IA1

Manual Integrations  
 APPROVED

MMDadoda  
 8/5/2021 1:41:30 PM

Quant Time: Aug 05 10:32:31 2021  
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA L\METHODS\VL080421AIR.M  
 Quant Title : AIR ANALYSIS BY METHOD TO-15 Instrument: MSVOA\_LWed Aug 04 22:53:01 2021  
 QLast Update : Wed Aug 04 22:53:01 2021  
 Response via : Initial Calibration

| Internal Standards      | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|-------------------------|-------|------|----------|-------|-------|----------|
| 1) Bromochloromethane   | 5.66  | 49   | 1021565  | 10.00 | ppbv  | 0.00     |
| 33) 1,4-Difluorobenzene | 7.16  | 114  | 3378453  | 10.00 | ppbv  | 0.00     |
| 55) Chlorobenzene-d5    | 12.07 | 117  | 2848603  | 10.00 | ppbv  | 0.00     |

System Monitoring Compounds

|                             |        |       |          |          |      |         |
|-----------------------------|--------|-------|----------|----------|------|---------|
| 68) 1-Bromo-4-Fluorobenzene | 14.40  | 95    | 1982158  | 10.19    | ppbv | 0.00    |
| Spiked Amount               | 10.000 | Range | 65 - 135 | Recovery | =    | 101.90% |

| Target Compounds           | R.T.  | QIon | Response | Conc   | Units  | Ovalue |
|----------------------------|-------|------|----------|--------|--------|--------|
| 2) Dichlorodifluoromethane | 3.04  | 85   | 46513    | 0.224  | ppbv   | 100    |
| 3) Chlorodifluoromethane   | 2.99  | 51   | 197286   | 0.950  | ppbv   | 94     |
| 4) Chloromethane           | 3.13  | 50   | 35398    | 0.476  | ppbv   | 89     |
| 9) Propene                 | 3.02  | 41   | 45931    | 0.612  | ppbv # | 80     |
| 10) Heptane                | 8.11  | 43   | 31676m   | 0.181  | ppbv   |        |
| 11) Trichlorofluoromethane | 3.94  | 101  | 46411m   | 0.243  | ppbv   |        |
| 13) Ethanol                | 3.61  | 45   | 194230m  | 17.170 | ppbv   |        |
| 15) Acetone                | 3.84  | 43   | 890019   | 8.021  | ppbv   | 97     |
| 17) tert-Butyl alcohol     | 4.30  | 59   | 15677m   | 0.141  | ppbv   |        |
| 19) Isopropyl Alcohol      | 3.97  | 45   | 150834   | 2.456  | ppbv # | 1      |
| 20) Methylene Chloride     | 4.34  | 84   | 969219   | 10.550 | ppbv   | 98     |
| 26) Hexane                 | 5.68  | 57   | 883186   | 5.470  | ppbv # | 48     |
| 29) Chloroform             | 5.74  | 83   | 98083    | 0.480  | ppbv   | 98     |
| 30) Cyclohexane            | 7.11  | 84   | 14189m   | 0.114  | ppbv   |        |
| 34) 2-Butanone             | 5.24  | 43   | 91645m   | 0.782  | ppbv   |        |
| 35) Carbon Tetrachloride   | 7.02  | 117  | 9660m    | 0.048  | ppbv   |        |
| 36) Benzene                | 6.88  | 78   | 125947   | 0.414  | ppbv   | 96     |
| 44) 2,2,4-Trimethylpentane | 7.86  | 57   | 93499    | 0.186  | ppbv # | 81     |
| 50) 4-Methyl-2-Pentanone   | 8.75  | 43   | 18547m   | 0.103  | ppbv   |        |
| 51) 2-Hexanone             | 10.16 | 43   | 32378m   | 0.397  | ppbv   |        |
| 52) Tetrachloroethene      | 11.20 | 164  | 7177m    | 0.055  | ppbv   |        |
| 53) Toluene                | 9.79  | 91   | 551457   | 1.536  | ppbv   | 100    |
| 58) Ethyl Benzene          | 12.70 | 91   | 53778m   | 0.129  | ppbv   |        |
| 59) m/p-Xylene             | 12.95 | 91   | 113365m  | 0.322  | ppbv   |        |
| 60) o-Xylene               | 13.68 | 91   | 37841m   | 0.118  | ppbv   |        |
| 61) Styrene                | 13.51 | 104  | 248032   | 1.413  | ppbv   | 94     |
| 69) p-Isopropyltoluene     | 17.64 | 119  | 58080m   | 0.119  | ppbv   |        |
| 74) 1,2,4-Trimethylbenzene | 16.75 | 105  | 60985    | 0.164  | ppbv # | 78     |

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

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