

Data Path : Z:\VOASRV\HPCHEM1\MSVOA L\DATA\VL120418\  
 Data File : VL032882.D  
 Acq On : 4 Dec 2018 18:44  
 Operator : SY/AP  
 Sample : J6175-03  
 Misc : 400ml/MSVOA L  
 ALS Vial : 1 Sample Multiplier: 1

Instrument :  
 MSVOA\_L  
 ClientSampleId :  
 MMSV-01

Manual Integrations  
 APPROVED

MMDadoda  
 12/10/2018 1:00:12 PM

Quant Time: Dec 05 06:33:03 2018  
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA L\METHODS\VL111418AIR.M  
 Quant Title : AIR ANALYSIS BY METHOD TO-15 Instrument: MSVOA\_L Wed Nov 14 07:00:23 2018  
 QLast Update : Wed Nov 14 07:00:23 2018  
 Response via : Initial Calibration

| Internal Standards      | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|-------------------------|-------|------|----------|-------|-------|----------|
| 1) Bromochloromethane   | 5.77  | 49   | 1425823  | 10.00 | ppbv  | 0.00     |
| 33) 1,4-Difluorobenzene | 7.30  | 114  | 3523547  | 10.00 | ppbv  | 0.00     |
| 55) Chlorobenzene-d5    | 12.25 | 117  | 3523731  | 10.00 | ppbv  | 0.00     |

|                             |        |       |          |          |      |        |
|-----------------------------|--------|-------|----------|----------|------|--------|
| System Monitoring Compounds |        |       |          |          |      |        |
| 68) 1-Bromo-4-Fluorobenzene | 14.58  | 95    | 2552475  | 9.68     | ppbv | 0.00   |
| Spiked Amount               | 10.000 | Range | 65 - 135 | Recovery | =    | 96.80% |

| Target Compounds           | R.T.  | QIon | Response | Conc   | Units  | Ovalue |
|----------------------------|-------|------|----------|--------|--------|--------|
| 2) Dichlorodifluoromethane | 3.08  | 85   | 157343   | 0.768  | ppbv   | 100    |
| 3) Chlorodifluoromethane   | 3.02  | 51   | 41101    | 0.364  | ppbv # | 89     |
| 9) Propene                 | 3.04  | 41   | 532206   | 11.765 | ppbv   | 94     |
| 10) Heptane                | 8.26  | 43   | 36762m   | 0.213  | ppbv   |        |
| 11) Trichlorofluoromethane | 4.01  | 101  | 76275    | 0.395  | ppbv   | 90     |
| 13) Ethanol                | 3.66  | 45   | 206561   | 13.941 | ppbv   | 99     |
| 15) Acetone                | 3.90  | 43   | 2012174  | 15.453 | ppbv # | 79     |
| 20) Methylene Chloride     | 4.42  | 84   | 314553   | 4.666  | ppbv   | 95     |
| 26) Hexane                 | 5.78  | 57   | 271046   | 2.006  | ppbv # | 11     |
| 27) Carbon Disulfide       | 4.62  | 76   | 87264    | 0.509  | ppbv # | 81     |
| 29) Chloroform             | 5.86  | 83   | 68207    | 0.283  | ppbv   | 92     |
| 34) 2-Butanone             | 5.33  | 43   | 10003252 | 54.736 | ppbv   | 95     |
| 36) Benzene                | 7.01  | 78   | 193451   | 0.689  | ppbv   | 99     |
| 44) 2,2,4-Trimethylpentane | 8.00  | 57   | 65103m   | 0.139  | ppbv   |        |
| 51) 2-Hexanone             | 10.27 | 43   | 1159937  | 6.308  | ppbv   | 99     |
| 52) Tetrachloroethene      | 11.37 | 164  | 6316m    | 0.065  | ppbv   |        |
| 53) Toluene                | 9.95  | 91   | 333520   | 1.139  | ppbv   | 99     |
| 58) Ethyl Benzene          | 12.87 | 91   | 83352    | 0.197  | ppbv   | 98     |
| 59) m/p-Xylene             | 13.12 | 91   | 232397m  | 0.616  | ppbv   |        |
| 60) o-Xylene               | 13.86 | 91   | 114684m  | 0.302  | ppbv   |        |
| 61) Styrene                | 13.69 | 104  | 58788    | 0.238  | ppbv   | 99     |
| 72) 4-Ethyltoluene         | 16.01 | 105  | 44836m   | 0.108  | ppbv   |        |
| 73) 1,3,5-Trimethylbenzene | 16.16 | 105  | 50089m   | 0.123  | ppbv   |        |
| 74) 1,2,4-Trimethylbenzene | 16.93 | 105  | 132729m  | 0.305  | ppbv   |        |

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

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