

Data Path : Z:\VOASRV\HPCHEM1\MSVOA L\DATA\VL111418\  
 Data File : VL032754.D  
 Acq On : 14 Nov 2018 11:45  
 Operator : SY/AP  
 Sample : J5891-01  
 Misc : 400ml/MSVOA L  
 ALS Vial : 1 Sample Multiplier: 1

Instrument :  
 MSVOA\_L  
 ClientSampleId :  
 IASC1

Manual Integrations  
 APPROVED

MMDadoda  
 11/16/2018 1:30:31 PM

Quant Time: Nov 15 05:54:06 2018  
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA L\METHODS\VL111418AIR.M  
 Quant Title : AIR ANALYSIS BY METHOD TO-15 Instrument: MSVOA\_LWed Nov 14 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   | 1328910  | 10.00 | ppbv  | 0.00     |
| 33) 1,4-Difluorobenzene | 7.30  | 114  | 3276052  | 10.00 | ppbv  | 0.00     |
| 55) Chlorobenzene-d5    | 12.24 | 117  | 3057145  | 10.00 | ppbv  | 0.00     |

|                             |        |       |          |          |      |        |
|-----------------------------|--------|-------|----------|----------|------|--------|
| System Monitoring Compounds |        |       |          |          |      |        |
| 68) 1-Bromo-4-Fluorobenzene | 14.58  | 95    | 2256604  | 9.87     | ppbv | 0.00   |
| Spiked Amount               | 10.000 | Range | 65 - 135 | Recovery | =    | 98.70% |

| Target Compounds           | R.T.  | QIon | Response | Conc    | Units  | Ovalue |
|----------------------------|-------|------|----------|---------|--------|--------|
| 2) Dichlorodifluoromethane | 3.08  | 85   | 106842   | 0.559   | ppbv   | 99     |
| 3) Chlorodifluoromethane   | 3.02  | 51   | 104974m  | 0.996   | ppbv   |        |
| 4) Chloromethane           | 3.18  | 50   | 31810    | 0.545   | ppbv   | 94     |
| 9) Propene                 | 3.05  | 41   | 107707m  | 2.555   | ppbv   |        |
| 10) Heptane                | 8.25  | 43   | 79755    | 0.495   | ppbv   | 92     |
| 11) Trichlorofluoromethane | 4.01  | 101  | 53982    | 0.300   | ppbv   | 93     |
| 13) Ethanol                | 3.66  | 45   | 1415455  | 102.497 | ppbv   | 98     |
| 15) Acetone                | 3.91  | 43   | 741819   | 6.113   | ppbv   | 96     |
| 19) Isopropyl Alcohol      | 4.04  | 45   | 229318   | 2.836   | ppbv # | 97     |
| 20) Methylene Chloride     | 4.41  | 84   | 299889   | 4.773   | ppbv   | 97     |
| 26) Hexane                 | 5.79  | 57   | 476955   | 3.788   | ppbv # | 44     |
| 30) Cyclohexane            | 7.26  | 84   | 23490m   | 0.213   | ppbv   |        |
| 34) 2-Butanone             | 5.34  | 43   | 96412    | 0.567   | ppbv   | 99     |
| 35) Carbon Tetrachloride   | 7.14  | 117  | 21579    | 0.082   | ppbv # | 80     |
| 36) Benzene                | 7.01  | 78   | 119182   | 0.457   | ppbv   | 94     |
| 41) Tetrahydrofuran        | 6.17  | 42   | 17751m   | 0.182   | ppbv   |        |
| 44) 2,2,4-Trimethylpentane | 7.99  | 57   | 379361   | 0.870   | ppbv # | 85     |
| 52) Tetrachloroethene      | 11.37 | 164  | 4834m    | 0.053   | ppbv   |        |
| 53) Toluene                | 9.95  | 91   | 3204888  | 11.776  | ppbv   | 100    |
| 58) Ethyl Benzene          | 12.87 | 91   | 206015   | 0.560   | ppbv   | 96     |
| 59) m/p-Xylene             | 13.13 | 91   | 677354   | 2.068   | ppbv   | 99     |
| 60) o-Xylene               | 13.85 | 91   | 267129   | 0.811   | ppbv   | 99     |
| 64) n-propylbenzene        | 15.73 | 120  | 26150m   | 0.223   | ppbv   |        |
| 72) 4-Ethyltoluene         | 16.00 | 105  | 123927   | 0.343   | ppbv # | 90     |
| 73) 1,3,5-Trimethylbenzene | 16.16 | 105  | 133016   | 0.377   | ppbv   | 92     |
| 74) 1,2,4-Trimethylbenzene | 16.93 | 105  | 445781   | 1.180   | ppbv # | 68     |
| 79) Naphthalene            | 22.43 | 128  | 45860m   | 0.185   | ppbv   |        |

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

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