

Data Path : Z:\voasrv\HPCHEM1\MSVOA\_U\Data\VU061021\  
 Data File : VU044168.D  
 Acq On : 10 Jun 2021 20:05  
 Operator : SY/MD  
 Sample : M2658-14  
 Misc : 5.0mL/MSVOA\_U/WATER  
 ALS Vial : 28 Sample Multiplier: 1

Instrument :  
 MSVOA\_U  
 ClientSampleId :  
 G-WALL-1

Quant Time: Jun 11 01:22:26 2021  
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_U\Method\82U060921W.M  
 Quant Title : SW846 8260  
 QLast Update : Thu Jun 10 05:11:37 2021  
 Response via : Initial Calibration

| Compound                    | R.T. QIon |     | Response | Conc       | Units  | Dev(Min) |
|-----------------------------|-----------|-----|----------|------------|--------|----------|
| -----                       |           |     |          |            |        |          |
| Internal Standards          |           |     |          |            |        |          |
| 1) Pentafluorobenzene       | 5.382     | 168 | 145914   | 50.000     | ug/l   | 0.00     |
| 34) 1,4-Difluorobenzene     | 6.256     | 114 | 287284   | 50.000     | ug/l   | 0.00     |
| 63) Chlorobenzene-d5        | 9.423     | 117 | 303805   | 50.000     | ug/l   | 0.00     |
| 72) 1,4-Dichlorobenzene-d4  | 11.819    | 152 | 144622   | 50.000     | ug/l   | 0.00     |
| System Monitoring Compounds |           |     |          |            |        |          |
| 33) 1,2-Dichloroethane-d4   | 5.710     | 65  | 124448   | 53.632     | ug/l   | 0.00     |
| Spiked Amount               | 50.000    |     | Recovery | = 107.260% |        |          |
| 35) Dibromofluoromethane    | 5.298     | 113 | 100491   | 50.285     | ug/l   | 0.00     |
| Spiked Amount               | 50.000    |     | Recovery | = 100.580% |        |          |
| 50) Toluene-d8              | 7.906     | 98  | 391652   | 51.030     | ug/l   | 0.00     |
| Spiked Amount               | 50.000    |     | Recovery | = 102.060% |        |          |
| 62) 4-Bromofluorobenzene    | 10.639    | 95  | 144193   | 46.881     | ug/l   | 0.00     |
| Spiked Amount               | 50.000    |     | Recovery | = 93.760%  |        |          |
| Target Compounds            |           |     |          |            |        |          |
|                             |           |     |          |            |        | Qvalue   |
| 16) Acetone                 | 2.633     | 43  | 27647    | 21.115     | ug/l   | 98       |
| 20) Methylene Chloride      | 3.054     | 84  | 13674    | 5.472      | ug/l   | 97       |
| 29) Tetrahydrofuran         | 5.070     | 42  | 1093     | 1.038      | ug/l # | 90       |
| 43) Isopropyl Acetate       | 5.922     | 43  | 5868     | 1.081      | ug/l   | 93       |
| 51) 4-Methyl-2-Pentanone    | 7.800     | 43  | 8933     | 2.273      | ug/l   | 93       |
| -----                       |           |     |          |            |        |          |

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

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