

Data Path : Z:\voasrv\HPCHEM1\MSVOA\_U\Data\VU061021\  
 Data File : VU044165.D  
 Acq On : 10 Jun 2021 18:54  
 Operator : SY/MD  
 Sample : M2658-05  
 Misc : 5.0mL/MSVOA\_U/WATER  
 ALS Vial : 25 Sample Multiplier: 1

Instrument :  
 MSVOA\_U  
 ClientSampleId :  
 P-WALL

Quant Time: Jun 11 01:21:56 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  | 143306   | 50.000     | ug/l   | 0.00     |
| 34) 1,4-Difluorobenzene     | 6.256  | 114  | 285362   | 50.000     | ug/l   | 0.00     |
| 63) Chlorobenzene-d5        | 9.423  | 117  | 308908   | 50.000     | ug/l   | 0.00     |
| 72) 1,4-Dichlorobenzene-d4  | 11.819 | 152  | 152092   | 50.000     | ug/l   | 0.00     |
| System Monitoring Compounds |        |      |          |            |        |          |
| 33) 1,2-Dichloroethane-d4   | 5.710  | 65   | 123363   | 54.132     | ug/l   | 0.00     |
| Spiked Amount               | 50.000 |      | Recovery | = 108.260% |        |          |
| 35) Dibromofluoromethane    | 5.298  | 113  | 98572    | 49.657     | ug/l   | 0.00     |
| Spiked Amount               | 50.000 |      | Recovery | = 99.320%  |        |          |
| 50) Toluene-d8              | 7.906  | 98   | 389943   | 51.150     | ug/l   | 0.00     |
| Spiked Amount               | 50.000 |      | Recovery | = 102.300% |        |          |
| 62) 4-Bromofluorobenzene    | 10.639 | 95   | 152138   | 49.797     | ug/l   | 0.00     |
| Spiked Amount               | 50.000 |      | Recovery | = 99.600%  |        |          |
| Target Compounds            |        |      |          |            |        |          |
|                             |        |      |          |            |        | Qvalue   |
| 16) Acetone                 | 2.633  | 43   | 28473    | 22.247     | ug/l   | 98       |
| 20) Methylene Chloride      | 3.054  | 84   | 12390    | 4.953      | ug/l   | 96       |
| 29) Tetrahydrofuran         | 5.070  | 42   | 1080     | 1.044      | ug/l # | 82       |
| 43) Isopropyl Acetate       | 5.922  | 43   | 13676    | 2.535      | ug/l   | 98       |
| 51) 4-Methyl-2-Pentanone    | 7.803  | 43   | 5655     | 1.448      | ug/l   | 98       |
| -----                       |        |      |          |            |        |          |

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

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