

Data Path : Z:\voasrv\HPCHEM1\MSVOA\_N\Data\VN102121\  
 Data File : VN069093.D  
 Acq On : 21 Oct 2021 17:18  
 Operator : JC/MD  
 Sample : M4284-05 200X  
 Misc : 5.00mL/MSVOA\_N/WATER  
 ALS Vial : 17 Sample Multiplier: 1

Instrument :  
 MSVOA\_N  
 ClientSampleId :  
 30851-WATER

Quant Time: Oct 22 01:34:11 2021  
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_N\methods\82N101221W.M  
 Quant Title : SW846 8260  
 QLast Update : Wed Oct 13 05:44:40 2021  
 Response via : Initial Calibration

| Compound                    | R.T.   | QIon  | Response | Conc     | Units | Dev(Min)  |
|-----------------------------|--------|-------|----------|----------|-------|-----------|
| -----                       |        |       |          |          |       |           |
| Internal Standards          |        |       |          |          |       |           |
| 1) Pentafluorobenzene       | 8.091  | 168   | 322353   | 50.000   | ug/l  | 0.00      |
| 34) 1,4-Difluorobenzene     | 8.971  | 114   | 569287   | 50.000   | ug/l  | 0.00      |
| 63) Chlorobenzene-d5        | 11.746 | 117   | 529600   | 50.000   | ug/l  | 0.00      |
| 72) 1,4-Dichlorobenzene-d4  | 13.678 | 152   | 203398   | 50.000   | ug/l  | 0.00      |
| System Monitoring Compounds |        |       |          |          |       |           |
| 33) 1,2-Dichloroethane-d4   | 8.440  | 65    | 215056   | 50.600   | ug/l  | 0.00      |
| Spiked Amount               | 50.000 | Range | 61 - 141 | Recovery | =     | 101.200%  |
| 35) Dibromofluoromethane    | 8.024  | 113   | 167275   | 50.503   | ug/l  | 0.00      |
| Spiked Amount               | 50.000 | Range | 69 - 133 | Recovery | =     | 101.000%  |
| 50) Toluene-d8              | 10.443 | 98    | 707877   | 53.415   | ug/l  | 0.00      |
| Spiked Amount               | 50.000 | Range | 65 - 126 | Recovery | =     | 106.820%  |
| 62) 4-Bromofluorobenzene    | 12.736 | 95    | 247742   | 49.548   | ug/l  | 0.00      |
| Spiked Amount               | 50.000 | Range | 58 - 135 | Recovery | =     | 99.100%   |
| Target Compounds            |        |       |          |          |       |           |
| 16) Acetone                 | 4.304  | 43    | 5283     | 3.331    | ug/l  | Qvalue 99 |
| -----                       |        |       |          |          |       |           |

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

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