

Data Path : Z:\voasrv\HPCHEM1\MSVOA\_X\Data\VX102323\  
 Data File : VX038444.D  
 Acq On : 23 Oct 2023 19:43  
 Operator : JC/MD  
 Sample : 04980-03  
 Misc : 5.0mL/MSVOA\_X/WATER  
 ALS Vial : 27 Sample Multiplier: 1

Instrument :  
 MSVOA\_X  
 ClientSampleId :  
 MAIN-BAILER-SUMP

Quant Time: Oct 24 00:07:39 2023  
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_X\Method\82X100523W.M  
 Quant Title : SW846 8260  
 QLast Update : Thu Oct 05 15:27:13 2023  
 Response via : Initial Calibration

| Compound                    | R.T.   | QIon  | Response | Conc     | Units  | Dev(Min) |
|-----------------------------|--------|-------|----------|----------|--------|----------|
| -----                       |        |       |          |          |        |          |
| Internal Standards          |        |       |          |          |        |          |
| 1) Pentafluorobenzene       | 5.550  | 168   | 127654   | 50.000   | ug/l   | 0.00     |
| 34) 1,4-Difluorobenzene     | 6.763  | 114   | 233132   | 50.000   | ug/l   | 0.00     |
| 63) Chlorobenzene-d5        | 10.055 | 117   | 235895   | 50.000   | ug/l   | 0.00     |
| 72) 1,4-Dichlorobenzene-d4  | 12.024 | 152   | 125691   | 50.000   | ug/l   | 0.00     |
| System Monitoring Compounds |        |       |          |          |        |          |
| 33) 1,2-Dichloroethane-d4   | 5.958  | 65    | 90661    | 44.155   | ug/l   | 0.00     |
| Spiked Amount               | 50.000 | Range | 78 - 117 | Recovery | =      | 88.320%  |
| 35) Dibromofluoromethane    | 5.385  | 113   | 75827    | 45.139   | ug/l   | 0.00     |
| Spiked Amount               | 50.000 | Range | 75 - 124 | Recovery | =      | 90.280%  |
| 50) Toluene-d8              | 8.647  | 98    | 285421   | 44.464   | ug/l   | 0.00     |
| Spiked Amount               | 50.000 | Range | 92 - 112 | Recovery | =      | 88.920%# |
| 62) 4-Bromofluorobenzene    | 11.079 | 95    | 109924   | 45.170   | ug/l   | 0.00     |
| Spiked Amount               | 50.000 | Range | 83 - 123 | Recovery | =      | 90.340%  |
| Target Compounds            |        |       |          |          |        |          |
|                             |        |       |          |          | Qvalue |          |
| 16) Acetone                 | 2.373  | 43    | 2191     | 2.557    | ug/l   | 94       |
| 27) cis-1,2-Dichloroethene  | 4.495  | 96    | 12975    | 7.534    | ug/l   | 81       |
| 44) Trichloroethene         | 7.135  | 130   | 2569     | 1.516    | ug/l   | 90       |
| 64) Tetrachloroethene       | 9.275  | 164   | 16689    | 10.928   | ug/l   | 91       |
| 93) 1,2,4-Trichlorobenzene  | 13.585 | 180   | 16670    | 6.255    | ug/l   | 97       |
| 96) 1,2,3-Trichlorobenzene  | 13.963 | 180   | 3677     | 1.425    | ug/l   | 95       |
| -----                       |        |       |          |          |        |          |

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

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