

Data Path : Z:\voasrv\HPCHEM1\MSVOA\_N\Data\VN103123\  
 Data File : VN079758.D  
 Acq On : 31 Oct 2023 20:55  
 Operator : JC\MD  
 Sample : 05107-19  
 Misc : 5.0mL/MSVOA\_N/WATER  
 ALS Vial : 26 Sample Multiplier: 1

Instrument :  
 MSVOA\_N  
 ClientSampleId :  
 1051

Quant Time: Nov 01 01:27:37 2023  
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_N\methods\82N103023W.M  
 Quant Title : SW846 8260  
 QLast Update : Tue Oct 31 07:23:24 2023  
 Response via : Initial Calibration

| Compound                    | R.T.   | QIon  | Response | Conc     | Units  | Dev(Min)  |
|-----------------------------|--------|-------|----------|----------|--------|-----------|
| -----                       |        |       |          |          |        |           |
| Internal Standards          |        |       |          |          |        |           |
| 1) Pentafluorobenzene       | 8.230  | 168   | 303494   | 50.000   | ug/l   | 0.00      |
| 34) 1,4-Difluorobenzene     | 9.106  | 114   | 570702   | 50.000   | ug/l   | 0.00      |
| 63) Chlorobenzene-d5        | 11.870 | 117   | 595231   | 50.000   | ug/l   | 0.00      |
| 72) 1,4-Dichlorobenzene-d4  | 13.794 | 152   | 289712   | 50.000   | ug/l   | # 0.00    |
| System Monitoring Compounds |        |       |          |          |        |           |
| 33) 1,2-Dichloroethane-d4   | 8.582  | 65    | 220333   | 45.208   | ug/l   | 0.00      |
| Spiked Amount               | 50.000 | Range | 74 - 125 | Recovery | =      | 90.420%   |
| 35) Dibromofluoromethane    | 8.171  | 113   | 205917   | 50.495   | ug/l   | 0.00      |
| Spiked Amount               | 50.000 | Range | 75 - 124 | Recovery | =      | 101.000%  |
| 50) Toluene-d8              | 10.571 | 98    | 691493   | 45.487   | ug/l   | 0.00      |
| Spiked Amount               | 50.000 | Range | 86 - 113 | Recovery | =      | 90.980%   |
| 62) 4-Bromofluorobenzene    | 12.853 | 95    | 311534   | 65.075   | ug/l   | 0.00      |
| Spiked Amount               | 50.000 | Range | 83 - 123 | Recovery | =      | 130.160%# |
| Target Compounds            |        |       |          |          |        |           |
|                             |        |       |          |          | Qvalue |           |
| 16) Acetone                 | 4.430  | 43    | 124694   | 59.801   | ug/l   | 94        |
| 25) 2-Butanone              | 7.488  | 43    | 43951    | 15.345   | ug/l   | # 80      |
| 40) Benzene                 | 8.612  | 78    | 1050816  | 58.852   | ug/l   | 100       |
| 51) 4-Methyl-2-Pentanone    | 10.447 | 43    | 10139    | 1.700    | ug/l   | 89        |
| 52) Toluene                 | 10.635 | 92    | 649806   | 59.613   | ug/l   | 99        |
| 67) Ethyl Benzene           | 11.970 | 91    | 149729   | 6.451    | ug/l   | 97        |
| 68) m/p-Xylenes             | 12.070 | 106   | 283769   | 33.056   | ug/l   | 97        |
| 69) o-Xylene                | 12.400 | 106   | 171119   | 20.730   | ug/l   | 94        |
| 73) Isopropylbenzene        | 12.694 | 105   | 12320    | 0.469    | ug/l   | 99        |
| 78) n-propylbenzene         | 13.041 | 91    | 19292    | 0.638    | ug/l   | 95        |
| 80) 1,3,5-Trimethylbenzene  | 13.176 | 105   | 67387    | 3.150    | ug/l   | 99        |
| 84) 1,2,4-Trimethylbenzene  | 13.482 | 105   | 210382   | 10.195   | ug/l   | 100       |
| 95) Naphthalene             | 15.641 | 128   | 5789147  | 658.834  | ug/l   | 99        |
| -----                       |        |       |          |          |        |           |

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

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