

Data Path : Z:\voasrv\HPCHEM1\MSVOA\_N\Data\VN042822\  
 Data File : VN072256.D  
 Acq On : 28 Apr 2022 22:37  
 Operator : JC\MD  
 Sample : N2617-21  
 Misc : 5.0mL/MSVOA\_N/WATER  
 ALS Vial : 26 Sample Multiplier: 1

Instrument :  
 MSVOA\_N  
 ClientSampleId :  
 TB

Quant Time: Apr 29 06:38:55 2022  
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_N\methods\82N042622W.M  
 Quant Title : SW846 8260  
 QLast Update : Tue Apr 26 18:50:11 2022  
 Response via : Initial Calibration

| Compound                    | R.T.   | QIon  | Response | Conc     | Units  | Dev(Min) |
|-----------------------------|--------|-------|----------|----------|--------|----------|
| -----                       |        |       |          |          |        |          |
| Internal Standards          |        |       |          |          |        |          |
| 1) Pentafluorobenzene       | 8.086  | 168   | 276328   | 50.000   | ug/l   | # 0.00   |
| 34) 1,4-Difluorobenzene     | 8.963  | 114   | 511739   | 50.000   | ug/l   | 0.00     |
| 63) Chlorobenzene-d5        | 11.739 | 117   | 566259   | 50.000   | ug/l   | 0.00     |
| 72) 1,4-Dichlorobenzene-d4  | 13.668 | 152   | 207866   | 50.000   | ug/l   | 0.00     |
| System Monitoring Compounds |        |       |          |          |        |          |
| 33) 1,2-Dichloroethane-d4   | 8.433  | 65    | 211218   | 49.789   | ug/l   | 0.00     |
| Spiked Amount               | 50.000 | Range | 61 - 141 | Recovery | =      | 99.580%  |
| 35) Dibromofluoromethane    | 8.022  | 113   | 153612   | 46.898   | ug/l   | 0.00     |
| Spiked Amount               | 50.000 | Range | 69 - 133 | Recovery | =      | 93.800%  |
| 50) Toluene-d8              | 10.439 | 98    | 664744   | 51.649   | ug/l   | 0.00     |
| Spiked Amount               | 50.000 | Range | 65 - 126 | Recovery | =      | 103.300% |
| 62) 4-Bromofluorobenzene    | 12.727 | 95    | 263702   | 61.729   | ug/l   | 0.00     |
| Spiked Amount               | 50.000 | Range | 58 - 135 | Recovery | =      | 123.460% |
| Target Compounds            |        |       |          |          |        |          |
|                             |        |       |          |          | Qvalue |          |
| 16) Acetone                 | 4.304  | 43    | 3828     | 2.093    | ug/l   | # 73     |
| 52) Toluene                 | 10.504 | 92    | 2426     | 0.274    | ug/l   | 93       |
| 67) Ethyl Benzene           | 11.839 | 91    | 60960    | 3.055    | ug/l   | 92       |
| 68) m/p-Xylenes             | 11.945 | 106   | 38616    | 5.212    | ug/l   | 94       |
| 69) o-Xylene                | 12.280 | 106   | 2180     | 0.303    | ug/l   | 79       |
| 78) n-propylbenzene         | 12.915 | 91    | 12289    | 0.608    | ug/l   | 96       |
| 80) 1,3,5-Trimethylbenzene  | 13.051 | 105   | 7048     | 0.461    | ug/l   | 97       |
| 84) 1,2,4-Trimethylbenzene  | 13.362 | 105   | 40212    | 2.718    | ug/l   | 96       |
| 95) Naphthalene             | 15.504 | 128   | 112315   | 13.834   | ug/l   | 98       |
| -----                       |        |       |          |          |        |          |

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

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