

Data Path : Z:\voasrv\HPCHEM1\MSVOA\_N\Data\VN111225\  
 Data File : VN088271.D  
 Acq On : 12 Nov 2025 13:13  
 Operator : JC\MD  
 Sample : IBLK  
 Misc : 5.0mL/MSVOA\_N/WATER  
 ALS Vial : 9 Sample Multiplier: 1

Instrument :  
 MSVOA\_N  
 ClientSampleId :  
 IBLK

Quant Time: Nov 14 00:11:36 2025  
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_N\methods\82N102325W.M  
 Quant Title : SW846 8260  
 QLast Update : Mon Nov 03 01:16:06 2025  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon  | Response | Conc     | Units       | Dev(Min) |
|-------------------------------|--------|-------|----------|----------|-------------|----------|
| -----                         |        |       |          |          |             |          |
| Internal Standards            |        |       |          |          |             |          |
| 1) Pentafluorobenzene         | 8.206  | 168   | 184602   | 50.000   | ug/l        | 0.00     |
| 34) 1,4-Difluorobenzene       | 9.082  | 114   | 429826   | 50.000   | ug/l        | 0.00     |
| 63) Chlorobenzene-d5          | 11.841 | 117   | 433884   | 50.000   | ug/l        | 0.00     |
| 72) 1,4-Dichlorobenzene-d4    | 13.770 | 152   | 223288   | 50.000   | ug/l        | # 0.00   |
| System Monitoring Compounds   |        |       |          |          |             |          |
| 33) 1,2-Dichloroethane-d4     | 8.559  | 65    | 200992   | 62.964   | ug/l        | 0.00     |
| Spiked Amount                 | 50.000 | Range | 74 - 125 | Recovery | = 125.920%# |          |
| 35) Dibromofluoromethane      | 8.147  | 113   | 150578   | 52.146   | ug/l        | 0.00     |
| Spiked Amount                 | 50.000 | Range | 75 - 124 | Recovery | = 104.300%  |          |
| 50) Toluene-d8                | 10.541 | 98    | 567375   | 50.994   | ug/l        | 0.00     |
| Spiked Amount                 | 50.000 | Range | 86 - 113 | Recovery | = 101.980%  |          |
| 62) 4-Bromofluorobenzene      | 12.829 | 95    | 226803   | 57.719   | ug/l        | 0.00     |
| Spiked Amount                 | 50.000 | Range | 77 - 121 | Recovery | = 115.440%  |          |
| Target Compounds              |        |       |          |          |             |          |
|                               |        |       |          |          | Qvalue      |          |
| 5) Bromomethane               | 2.995  | 94    | 981      | 0.614    | ug/l        | 86       |
| 11) Tert butyl alcohol        | 5.530  | 59    | 134      | 0.244    | ug/l #      | 71       |
| 16) Acetone                   | 4.430  | 43    | 2971     | 1.853    | ug/l #      | 85       |
| 17) Carbon Disulfide          | 4.712  | 76    | 1415     | 0.201    | ug/l #      | 74       |
| 18) Methyl Acetate            | 5.030  | 43    | 1782     | 0.543    | ug/l #      | 54       |
| 27) cis-1,2-Dichloroethene    | 7.477  | 96    | 686      | 0.234    | ug/l #      | 1        |
| 29) Tetrahydrofuran           | 7.830  | 42    | 365      | 0.265    | ug/l #      | 34       |
| 31) Cyclohexane               | 8.200  | 56    | 6593     | 1.475    | ug/l #      | 42       |
| 36) 1,1-Dichloropropene       | 8.212  | 75    | 20000    | 4.744    | ug/l #      | 53       |
| 38) Carbon Tetrachloride      | 8.435  | 117   | 1267     | 0.291    | ug/l #      | 23       |
| 48) Methyl methacrylate       | 9.677  | 41    | 1328     | 0.344    | ug/l #      | 63       |
| 55) 1,1,2-Trichloroethane     | 10.988 | 97    | 1543     | 0.514    | ug/l #      | 25       |
| 59) 2-Hexanone                | 11.224 | 43    | 812      | 0.247    | ug/l        | 70       |
| 61) 1,2-Dibromoethane         | 11.500 | 107   | 1411     | 0.456    | ug/l        | 81       |
| 68) m/p-Xylenes               | 12.053 | 106   | 3248     | 0.501    | ug/l #      | 14       |
| 71) Bromoform                 | 12.535 | 173   | 2430     | 1.009    | ug/l #      | 28       |
| 74) N-amyl acetate            | 12.506 | 43    | 1311     | 0.268    | ug/l #      | 38       |
| 78) n-propylbenzene           | 13.059 | 91    | 4812     | 0.229    | ug/l        | 99       |
| 79) 2-Chlorotoluene           | 13.094 | 91    | 2970     | 0.248    | ug/l #      | 42       |
| 82) 4-Chlorotoluene           | 13.270 | 91    | 63883    | 4.927    | ug/l #      | 45       |
| 84) 1,2,4-Trimethylbenzene    | 13.400 | 105   | 3566     | 0.249    | ug/l        | 75       |
| 85) sec-Butylbenzene          | 13.588 | 105   | 4888     | 0.260    | ug/l        | 86       |
| 86) p-Isopropyltoluene        | 13.682 | 119   | 31108    | 2.066    | ug/l #      | 50       |
| 87) 1,3-Dichlorobenzene       | 13.694 | 146   | 2531     | 0.341    | ug/l #      | 24       |
| 88) 1,4-Dichlorobenzene       | 13.759 | 146   | 2102     | 0.264    | ug/l #      | 1        |
| 89) n-Butylbenzene            | 14.094 | 91    | 8149     | 0.546    | ug/l #      | 40       |
| 90) Hexachloroethane          | 14.329 | 117   | 1575     | 0.540    | ug/l        | 65       |
| 92) 1,2-Dibromo-3-Chloropr... | 14.429 | 75    | 693      | 0.555    | ug/l        | 72       |
| 93) 1,2,4-Trichlorobenzene    | 15.364 | 180   | 1082     | 0.250    | ug/l #      | 59       |
| 94) Hexachlorobutadiene       | 15.476 | 225   | 605      | 0.375    | ug/l #      | 63       |
| -----                         |        |       |          |          |             |          |

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

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