

Data Path : Z:\voasrv\HPCHEM1\MSVOA\_D\Data\VD081222\  
 Data File : VD074063.D  
 Acq On : 12 Aug 2022 15:20  
 Operator : VA/SY  
 Sample : N4107-09  
 Misc : 5.06G/5.00ml/MSVOA\_D/SOIL  
 ALS Vial : 8 Sample Multiplier: 1

Instrument :  
 MSVOA\_D  
 ClientSampleId :  
 SOIL-STOCKPILE

Quant Time: Aug 12 23:47:43 2022  
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_D\Method\82D081022S.M  
 Quant Title : SW846 8260  
 QLast Update : Wed Aug 10 14:55:58 2022  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon  | Response | Conc     | Units  | Dev(Min) |
|-------------------------------|--------|-------|----------|----------|--------|----------|
| -----                         |        |       |          |          |        |          |
| Internal Standards            |        |       |          |          |        |          |
| 1) Pentafluorobenzene         | 7.967  | 168   | 1500     | 50.000   | ug/l   | 0.00     |
| 34) 1,4-Difluorobenzene       | 8.849  | 114   | 2414     | 50.000   | ug/l   | 0.00     |
| 63) Chlorobenzene-d5          | 11.632 | 117   | 2322     | 50.000   | ug/l   | 0.00     |
| 72) 1,4-Dichlorobenzene-d4    | 13.561 | 152   | 995      | 50.000   | ug/l   | 0.00     |
| System Monitoring Compounds   |        |       |          |          |        |          |
| 33) 1,2-Dichloroethane-d4     | 8.320  | 65    | 662      | 56.998   | ug/l   | 0.00     |
| Spiked Amount                 | 50.000 | Range | 50 - 163 | Recovery | =      | 114.000% |
| 35) Dibromofluoromethane      | 7.914  | 113   | 425      | 26.507   | ug/l   | 0.01     |
| Spiked Amount                 | 50.000 | Range | 54 - 147 | Recovery | =      | 53.020%# |
| 50) Toluene-d8                | 10.332 | 98    | 1600     | 36.631   | ug/l   | 0.00     |
| Spiked Amount                 | 50.000 | Range | 49 - 140 | Recovery | =      | 73.260%  |
| 62) 4-Bromofluorobenzene      | 12.614 | 95    | 862      | 43.718   | ug/l   | 0.00     |
| Spiked Amount                 | 50.000 | Range | 25 - 144 | Recovery | =      | 87.440%  |
| Target Compounds              |        |       |          |          |        |          |
|                               |        |       |          |          | Qvalue |          |
| 3) Chloromethane              | 2.203  | 50    | 149      | 2.883    | ug/l # | 40       |
| 5) Bromomethane               | 2.809  | 94    | 139      | 6.320    | ug/l # | 4        |
| 6) Chloroethane               | 2.738  | 64    | 63       | 2.758    | ug/l # | 43       |
| 16) Acetone                   | 4.150  | 43    | 56       | 23.332   | ug/l # | 46       |
| 18) Methyl Acetate            | 4.832  | 43    | 53       | 5.252    | ug/l # | 49       |
| 29) Tetrahydrofuran           | 7.708  | 42    | 67       | 28.995   | ug/l # | 31       |
| 41) Methacrylonitrile         | 7.591  | 41    | 67       | 14.119   | ug/l # | 18       |
| 48) Methyl methacrylate       | 9.561  | 41    | 55       | 6.945    | ug/l # | 6        |
| 67) Ethyl Benzene             | 11.838 | 91    | 90       | 1.043    | ug/l # | 44       |
| 76) 1,2,3-Trichloropropane    | 12.608 | 75    | 373      | 41.059   | ug/l # | 100      |
| 81) trans-1,4-Dichloro-2-b... | 12.608 | 75    | 373      | 109.586  | ug/l # | 13       |
| 86) p-Isopropyltoluene        | 13.502 | 119   | 115      | 1.787    | ug/l # | 50       |
| 95) Naphthalene               | 15.384 | 128   | 64       | 2.088    | ug/l # | 69       |
| -----                         |        |       |          |          |        |          |

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

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