

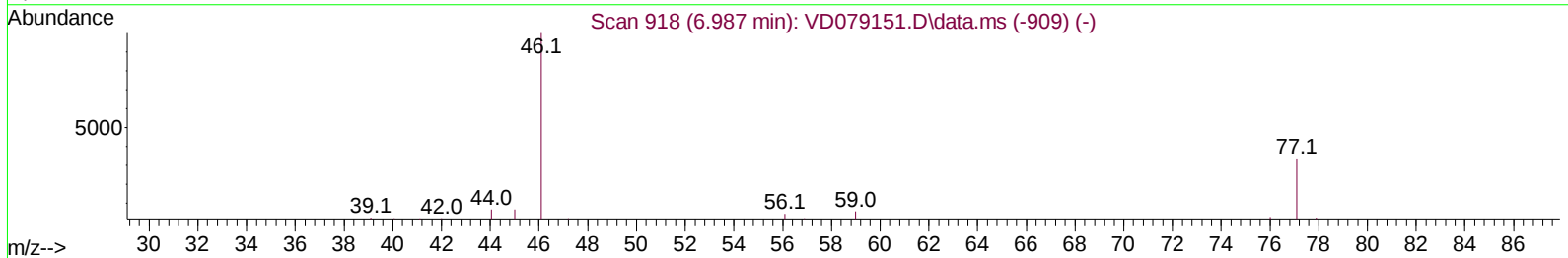
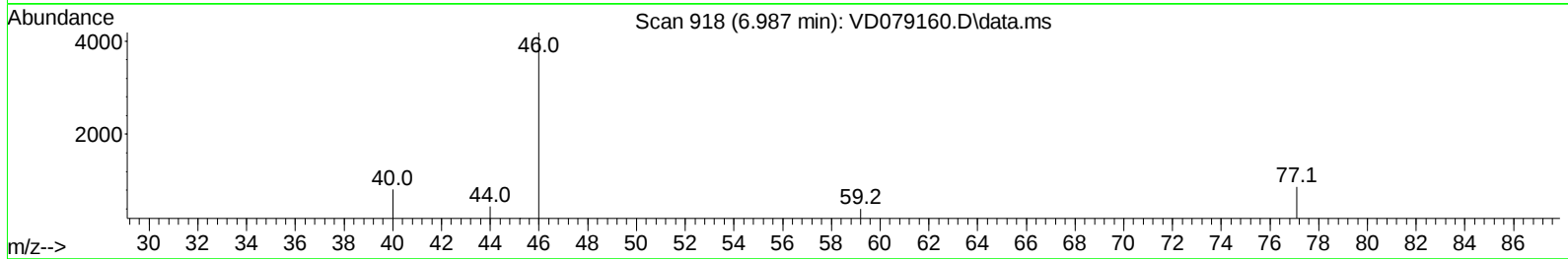
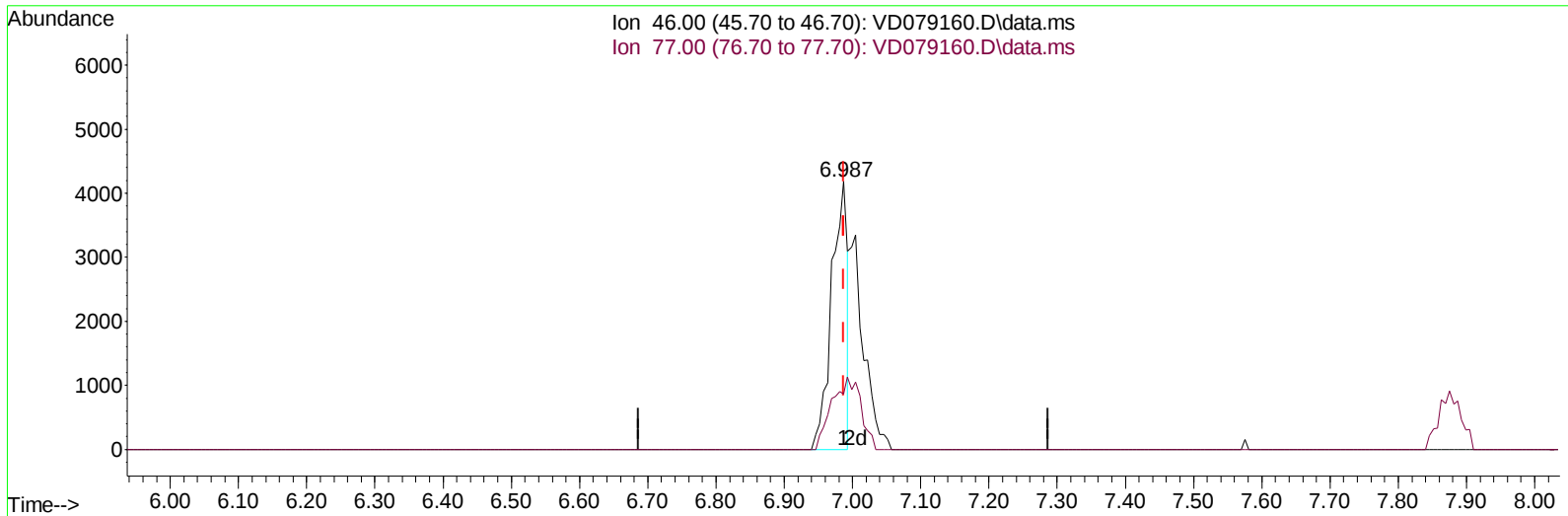
Data Path : Z:\voasrv\HPCHEM1\MSVOA\_D\Data\VD061024\  
Data File : VD079160.D  
Acq On : 10 Jun 2024 15:53  
Operator : RP/MD  
Sample : P2775-03MSD  
Misc : 5.48G/10ml/MSVOA\_D/SOIL/A  
ALS Vial : 11 Sample Multiplier: 1

Instrument :  
MSVOA\_D  
ClientSampleId :  
A4BQ3MSD

## Manual IntegrationsAPPROVED

Reviewed By :Mahesh Dadoda 06/13/2024  
Supervised By :Semsettin Yesilyurt 06/13/2024

Quant Time: Jun 11 02:49:14 2024  
Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_D\Method\SFAMDLM052924SMA.M  
Quant Title : SFAM01.0  
QLast Update : Tue Jun 11 02:46:15 2024  
Response via : Initial Calibration



TIC: VD079160.D\data.ms

## (21) 2-Butanone-d5 (S)

6.987min (+ 0.000) 19.26 ug/L

response 6840

| Ion   | Exp%   | Act%   |
|-------|--------|--------|
| 46.00 | 100.00 | 100.00 |
| 77.00 | 28.20  | 48.06# |
| 0.00  | 0.00   | 0.00   |
| 0.00  | 0.00   | 0.00   |

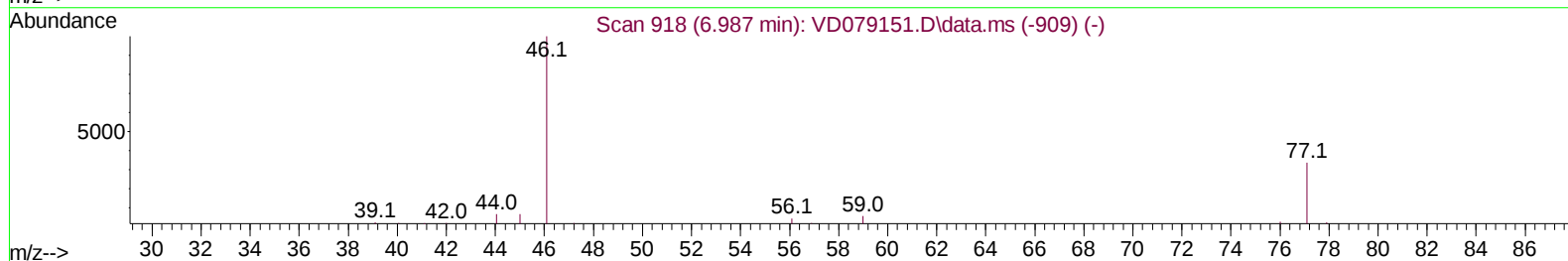
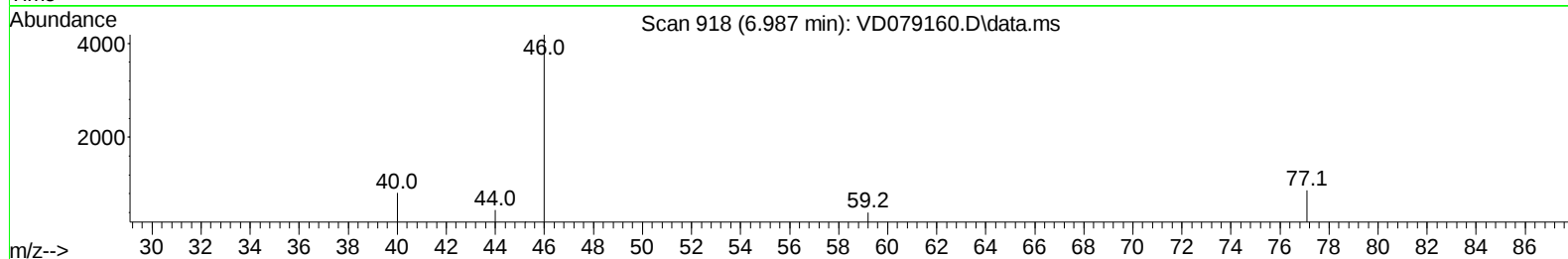
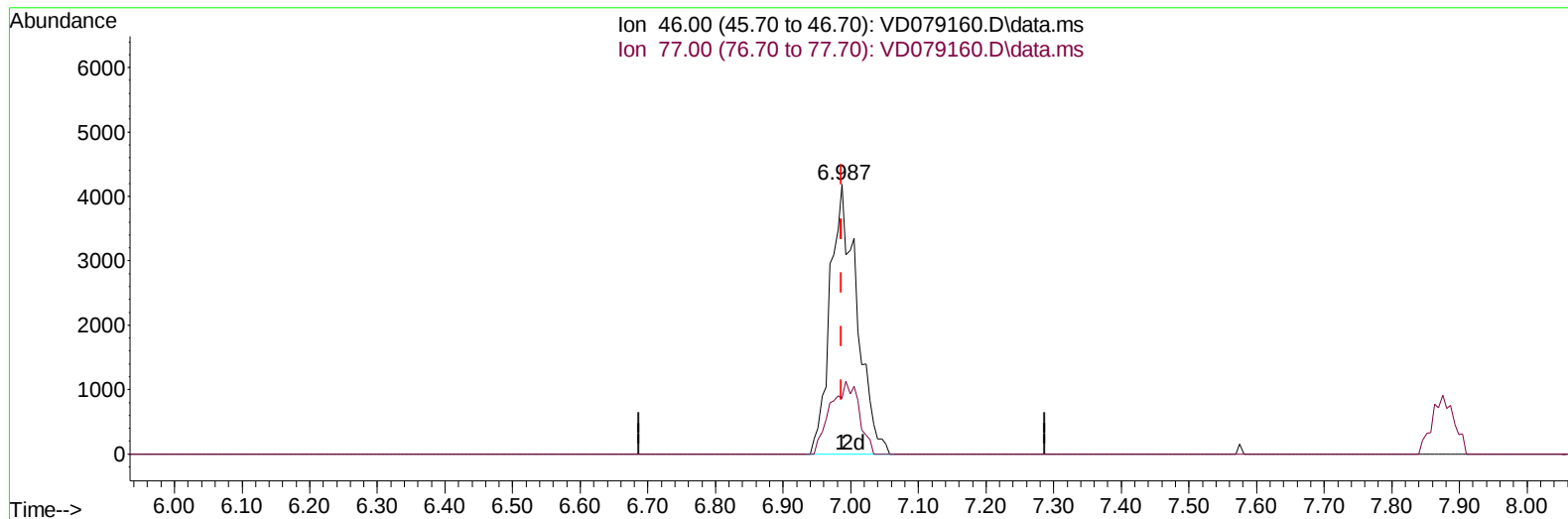
Data Path : Z:\voasrv\HPCHEM1\MSVOA\_D\Data\VD061024\  
Data File : VD079160.D  
Acq On : 10 Jun 2024 15:53  
Operator : RP/MD  
Sample : P2775-03MSD  
Misc : 5.48G/10ml/MSVOA\_D/SOIL/A  
ALS Vial : 11 Sample Multiplier: 1

Instrument :  
MSVOA\_D  
ClientSampleId :  
A4BQ3MSD

Manual IntegrationsAPPROVED

Quant Time: Jun 11 02:49:14 2024  
Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_D\Method\SFAMDLM052924SMA.M  
Quant Title : SFAM01.0  
QLast Update : Tue Jun 11 02:46:15 2024  
Response via : Initial Calibration

Reviewed By :Mahesh Dadoda 06/13/2024  
Supervised By :Semsettin Yesilyurt 06/13/2024



TIC: VD079160.D\data.ms

**(21) 2-Butanone-d5 (S)**

**6.987min (+ 0.000) 32.30 ug/L m**

**response 11471**

| Ion   | Exp%   | Act%   |
|-------|--------|--------|
| 46.00 | 100.00 | 100.00 |
| 77.00 | 28.20  | 28.65  |
| 0.00  | 0.00   | 0.00   |
| 0.00  | 0.00   | 0.00   |

Data Path : Z:\voasrv\HPCHEM1\MSVOA\_D\Data\VD061024\  
 Data File : VD079160.D  
 Acq On : 10 Jun 2024 15:53  
 Operator : RP/MD  
 Sample : P2775-03MSD  
 Misc : 5.48G/10mL/MSVOA\_D/SOIL/A  
 ALS Vial : 11 Sample Multiplier: 1

Instrument :  
 MSVOA\_D  
 ClientSampleId :  
 A4BQ3MSD

Manual IntegrationsAPPROVED

Quant Time: Jun 11 02:49:14 2024  
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_D\Method\SFAMDLM052924SMA.M  
 Quant Title : SFAM01.0  
 QLast Update : Tue Jun 11 02:46:15 2024  
 Response via : Initial Calibration

Reviewed By :Mahesh Dadoda 06/13/2024  
 Supervised By :Semsettin Yesilyurt 06/13/2024

| Compound                      | R.T.     | QIon  | Response | Conc   | Units    | Dev(Min) |
|-------------------------------|----------|-------|----------|--------|----------|----------|
| -----                         |          |       |          |        |          |          |
| Internal Standards            |          |       |          |        |          |          |
| 1) 1,4-Difluorobenzene        | 8.781    | 114   | 150894   | 25.000 | ug/L     | 0.00     |
| 28) Chlorobenzene-d5          | 11.581   | 117   | 122394   | 25.000 | ug/L     | 0.00     |
| 58) 1,4-Dichlorobenzene-d4    | 13.516   | 152   | 40184    | 25.000 | ug/L     | 0.00     |
| System Monitoring Compounds   |          |       |          |        |          |          |
| 4) Vinyl Chloride-d3          | 2.270    | 65    | 95088    | 20.478 | ug/L     | 0.00     |
| Spiked Amount 25.000          | Range 30 | - 150 | Recovery | =      | 81.920%  |          |
| 7) Chloroethane-d5            | 2.799    | 69    | 96610    | 21.517 | ug/L     | 0.00     |
| Spiked Amount 25.000          | Range 30 | - 150 | Recovery | =      | 86.080%  |          |
| 11) 1,1-Dichloroethene-d2     | 3.917    | 65    | 16219    | 19.676 | ug/L     | 0.00     |
| Spiked Amount 25.000          | Range 45 | - 110 | Recovery | =      | 78.720%  |          |
| 21) 2-Butanone-d5             | 6.987    | 46    | 11471m   | 32.298 | ug/L     | 0.00     |
| Spiked Amount 50.000          | Range 20 | - 135 | Recovery | =      | 64.600%  |          |
| 24) Chloroform-d              | 7.569    | 84    | 93443    | 22.148 | ug/L     | 0.00     |
| Spiked Amount 25.000          | Range 40 | - 150 | Recovery | =      | 88.600%  |          |
| 26) 1,2-Dichloroethane-d4     | 8.228    | 65    | 43351    | 22.274 | ug/L     | 0.00     |
| Spiked Amount 25.000          | Range 70 | - 130 | Recovery | =      | 89.080%  |          |
| 32) Benzene-d6                | 8.199    | 84    | 184892   | 27.250 | ug/L     | 0.00     |
| Spiked Amount 25.000          | Range 20 | - 135 | Recovery | =      | 109.000% |          |
| 36) 1,2-Dichloropropane-d6    | 9.210    | 67    | 51714    | 26.459 | ug/L     | 0.00     |
| Spiked Amount 25.000          | Range 70 | - 120 | Recovery | =      | 105.840% |          |
| 41) Toluene-d8                | 10.269   | 98    | 153001   | 24.083 | ug/L     | 0.00     |
| Spiked Amount 25.000          | Range 30 | - 130 | Recovery | =      | 96.320%  |          |
| 43) trans-1,3-Dichloroprop... | 10.522   | 79    | 19052    | 23.192 | ug/L     | 0.00     |
| Spiked Amount 25.000          | Range 30 | - 135 | Recovery | =      | 92.760%  |          |
| 47) 2-Hexanone-d5             | 10.875   | 63    | 10068    | 39.847 | ug/L     | 0.00     |
| Spiked Amount 50.000          | Range 20 | - 135 | Recovery | =      | 79.700%  |          |
| 56) 1,1,2,2-Tetrachloroeth... | 12.651   | 84    | 36189    | 20.904 | ug/L     | 0.00     |
| Spiked Amount 25.000          | Range 45 | - 120 | Recovery | =      | 83.600%  |          |
| 66) 1,2-Dichlorobenzene-d4    | 13.816   | 152   | 27663    | 19.947 | ug/L     | 0.00     |
| Spiked Amount 25.000          | Range 75 | - 120 | Recovery | =      | 79.800%  |          |
| Target Compounds              |          |       |          |        |          |          |
|                               |          |       |          |        | Qvalue   |          |
| 2) Dichlorodifluoromethane    | 1.934    | 85    | 48899    | 20.471 | ug/L     | 97       |
| 3) Chloromethane              | 2.146    | 50    | 92992    | 29.209 | ug/L     | 100      |
| 5) Vinyl chloride             | 2.281    | 62    | 127799   | 24.321 | ug/L     | 95       |
| 6) Bromomethane               | 2.693    | 94    | 86069    | 21.217 | ug/L     | 99       |
| 8) Chloroethane               | 2.834    | 64    | 93849    | 24.368 | ug/L     | 91       |
| 9) Trichlorofluoromethane     | 3.170    | 101   | 83355    | 20.823 | ug/L     | 99       |
| 10) 1,1,2-Trichloro-1,2,2-... | 3.964    | 101   | 48670    | 20.276 | ug/L #   | 86       |
| 12) 1,1-Dichloroethene        | 3.940    | 96    | 45630    | 21.835 | ug/L     | 90       |
| 13) Acetone                   | 4.022    | 43    | 12664    | 24.629 | ug/L     | 95       |
| 14) Carbon disulfide          | 4.269    | 76    | 146184   | 20.053 | ug/L     | 98       |
| 15) Methyl Acetate            | 4.564    | 43    | 11608    | 16.569 | ug/L     | 97       |
| 16) Methylene chloride        | 4.799    | 84    | 71760    | 26.690 | ug/L     | 95       |
| 17) trans-1,2-Dichloroethene  | 5.316    | 96    | 52005    | 23.028 | ug/L     | 90       |
| 18) Methyl tert-butyl Ether   | 5.316    | 73    | 91357    | 22.809 | ug/L #   | 91       |
| 19) 1,1-Dichloroethane        | 6.111    | 63    | 92703    | 25.496 | ug/L     | 98       |
| 20) cis-1,2-Dichloroethene    | 7.081    | 96    | 57025    | 24.250 | ug/L     | 96       |
| 22) 2-Butanone                | 7.087    | 43    | 16564    | 30.665 | ug/L     | 98       |
| 23) Bromochloromethane        | 7.428    | 128   | 28450    | 23.863 | ug/L     | 89       |

Data Path : Z:\voasrv\HPCHEM1\MSVOA\_D\Data\VD061024\  
 Data File : VD079160.D  
 Acq On : 10 Jun 2024 15:53  
 Operator : RP/MD  
 Sample : P2775-03MSD  
 Misc : 5.48G/10mL/MSVOA\_D/SOIL/A  
 ALS Vial : 11 Sample Multiplier: 1

Instrument :  
 MSVOA\_D  
 ClientSampleId :  
 A4BQ3MSD

Manual IntegrationsAPPROVED

Quant Time: Jun 11 02:49:14 2024  
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_D\Method\SFAMDLM052924SMA.M  
 Quant Title : SFAM01.0  
 QLast Update : Tue Jun 11 02:46:15 2024  
 Response via : Initial Calibration

Reviewed By :Mahesh Dadoda 06/13/2024  
 Supervised By :Semsettin Yesilyurt 06/13/2024

| Compound                      | R.T.   | QIon | Response | Conc   | Units  | Dev(Min) |
|-------------------------------|--------|------|----------|--------|--------|----------|
| 25) Chloroform                | 7.593  | 83   | 102580   | 24.849 | ug/L   | 100      |
| 27) 1,2-Dichloroethane        | 8.328  | 62   | 56541    | 23.594 | ug/L # | 94       |
| 29) Cyclohexane               | 7.875  | 56   | 43214    | 19.569 | ug/L   | 98       |
| 30) 1,1,1-Trichloroethane     | 7.793  | 97   | 81731    | 28.205 | ug/L   | 98       |
| 31) Carbon tetrachloride      | 7.987  | 117  | 70401    | 26.144 | ug/L   | 96       |
| 33) Benzene                   | 8.252  | 78   | 220645   | 30.268 | ug/L   | 100      |
| 34) Trichloroethene           | 9.028  | 95   | 48906    | 26.048 | ug/L   | 97       |
| 35) Methylcyclohexane         | 9.269  | 83   | 43355    | 13.936 | ug/L   | 98       |
| 37) 1,2-Dichloropropane       | 9.305  | 63   | 52504    | 29.405 | ug/L   | 100      |
| 38) Bromodichloromethane      | 9.587  | 83   | 71153    | 28.665 | ug/L   | 90       |
| 39) cis-1,3-Dichloropropene   | 10.016 | 75   | 70131    | 25.160 | ug/L   | 100      |
| 40) 4-Methyl-2-pentanone      | 10.157 | 43   | 35810    | 48.043 | ug/L   | 98       |
| 42) Toluene                   | 10.334 | 91   | 216181   | 26.706 | ug/L   | 98       |
| 44) trans-1,3-Dichloropropene | 10.551 | 75   | 59395    | 25.199 | ug/L   | 96       |
| 45) 1,1,2-Trichloroethane     | 10.734 | 97   | 40397    | 25.788 | ug/L   | 96       |
| 46) Tetrachloroethene         | 10.810 | 164  | 33375    | 20.188 | ug/L   | 95       |
| 48) 2-Hexanone                | 10.922 | 43   | 23029    | 35.704 | ug/L   | 95       |
| 49) Dibromochloromethane      | 11.075 | 129  | 47003    | 25.338 | ug/L   | 92       |
| 50) 1,2-Dibromoethane         | 11.175 | 107  | 35519    | 24.672 | ug/L   | 95       |
| 51) Chlorobenzene             | 11.604 | 112  | 121021   | 22.470 | ug/L   | 99       |
| 52) Ethylbenzene              | 11.681 | 91   | 183834   | 21.772 | ug/L   | 99       |
| 53) m,p-Xylene                | 11.793 | 106  | 73163    | 22.085 | ug/L   | 94       |
| 54) o-Xylene                  | 12.122 | 106  | 70295    | 21.963 | ug/L   | 98       |
| 55) Styrene                   | 12.134 | 104  | 118874   | 20.947 | ug/L   | 96       |
| 57) 1,1,2,2-Tetrachloroethane | 12.669 | 83   | 36448    | 20.786 | ug/L   | 92       |
| 59) Bromoform                 | 12.298 | 173  | 24365    | 33.083 | ug/L   | 99       |
| 60) Isopropylbenzene          | 12.422 | 105  | 150117   | 28.253 | ug/L   | 99       |
| 61) 1,2,3-Trichloropropane    | 12.722 | 75   | 25645    | 36.148 | ug/L   | 98       |
| 62) 1,3,5-Trimethylbenzene    | 12.904 | 105  | 86938    | 23.930 | ug/L   | 99       |
| 63) 1,2,4-Trimethylbenzene    | 13.210 | 105  | 96789    | 24.909 | ug/L   | 100      |
| 64) 1,3-Dichlorobenzene       | 13.457 | 146  | 60362    | 23.648 | ug/L   | 99       |
| 65) 1,4-Dichlorobenzene       | 13.534 | 146  | 61017    | 22.560 | ug/L   | 98       |
| 67) 1,2-Dichlorobenzene       | 13.834 | 146  | 49471    | 20.965 | ug/L   | 96       |
| 68) 1,2-Dibromo-3-chloropr... | 14.445 | 75   | 4066     | 27.879 | ug/L # | 86       |
| 69) 1,3,5-Trichlorobenzene    | 14.592 | 180  | 21534    | 13.376 | ug/L   | 99       |
| 70) 1,2,4-trichlorobenzene    | 15.098 | 180  | 13977    | 10.454 | ug/L # | 83       |
| 71) Naphthalene               | 15.328 | 128  | 30409    | 12.161 | ug/L   | 99       |
| 72) 1,2,3-Trichlorobenzene    | 15.522 | 180  | 10327    | 8.937  | ug/L   | 96       |

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

Data Path : Z:\voasrv\HPCHEM1\MSVOA\_D\Data\VD061024\  
Data File : VD079160.D  
Acq On : 10 Jun 2024 15:53  
Operator : RP/MD  
Sample : P2775-03MSD  
Misc : 5.48G/10ml/MSVOA\_D/SOIL/A  
ALS Vial : 11 Sample Multiplier: 1

Instrument :  
MSVOA\_D  
ClientSampleId :  
A4BQ3MSD

## Manual IntegrationsAPPROVED

Reviewed By :Mahesh Dadoda 06/13/2024  
Supervised By :Semsettin Yesilyurt 06/13/2024

Quant Time: Jun 11 02:49:14 2024  
Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_D\Method\SFAMDLM052924SMA.M  
Quant Title : SFAM01.0  
QLast Update : Tue Jun 11 02:46:15 2024  
Response via : Initial Calibration

