

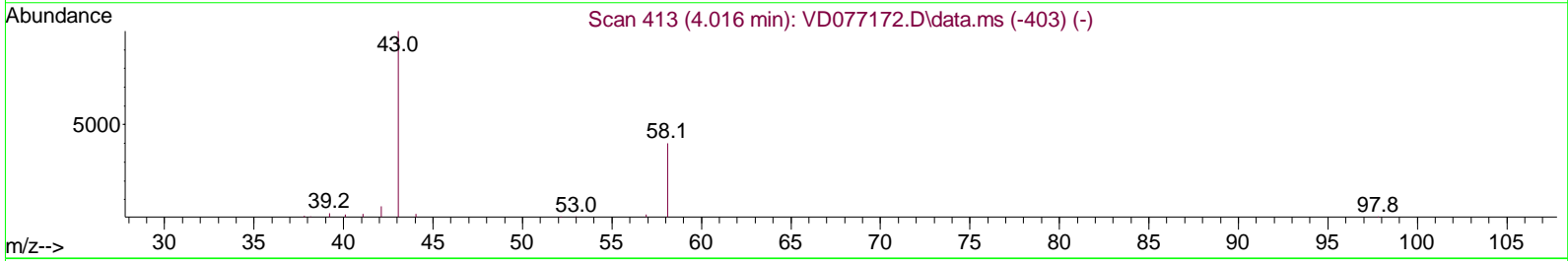
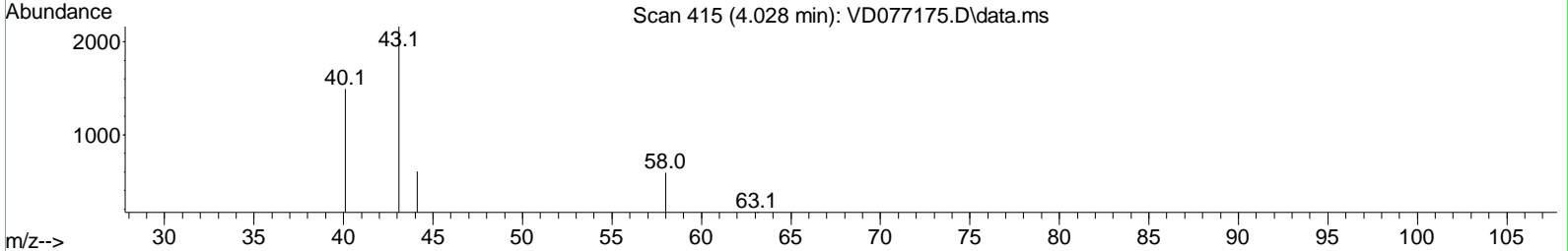
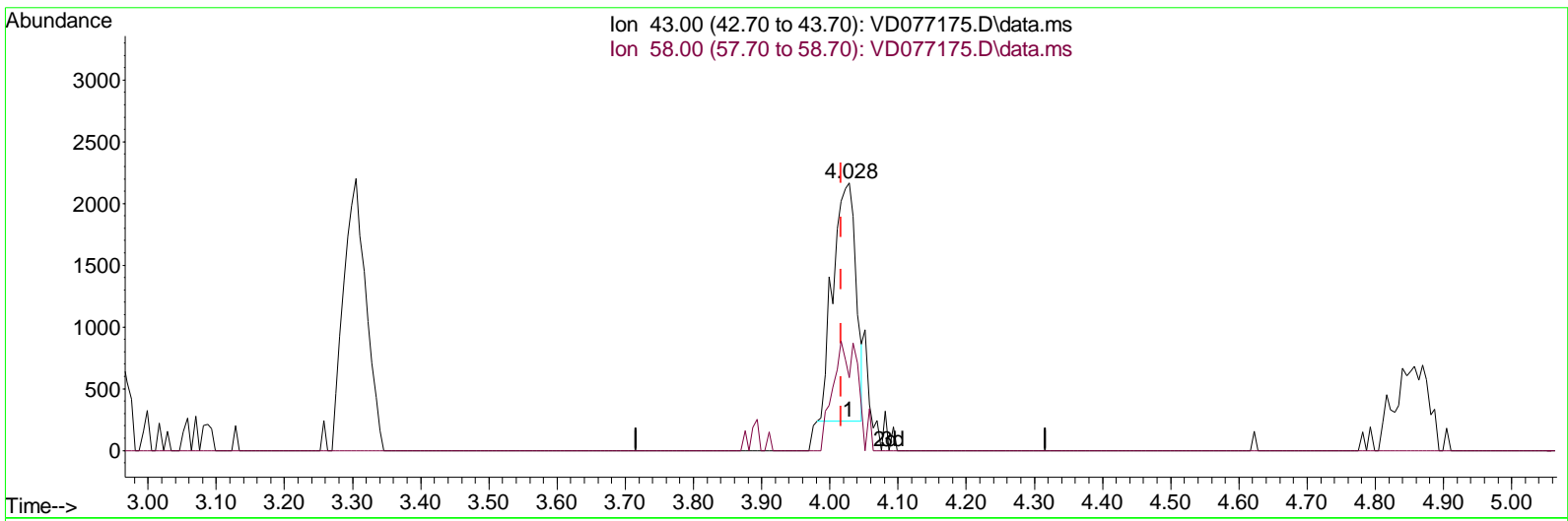
Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD092123\
 Data File : VD077175.D
 Acq On : 21 Sep 2023 12:43
 Operator : JC/SY
 Sample : 04527-02
 Misc : 5.70G/10.0ml/MSVOA_D/SOIL
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 MSVOA_D
ClientSampleId :
 EZYG9

Manual Integrations APPROVED

Reviewed By : John Carlone 09/22/2023
 Supervised By : Mahesh Dadoda 09/22/2023

Quant Time: Sep 22 01:30:45 2023
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_D\Method\SFAMDLMO90123SMA.M
 Quant Title : SFAM01.0
 QLast Update : Fri Sep 22 01:29:35 2023
 Response via : Initial Calibration



TIC: VD077175.D\data.ms

(13) Acetone (T)

4.028min (+ 0.012) 3.68 ug/L

response	4515	
Ion	Exp%	Act%
43.00	100.00	100.00
58.00	35.90	15.24
0.00	0.00	0.00
0.00	0.00	0.00

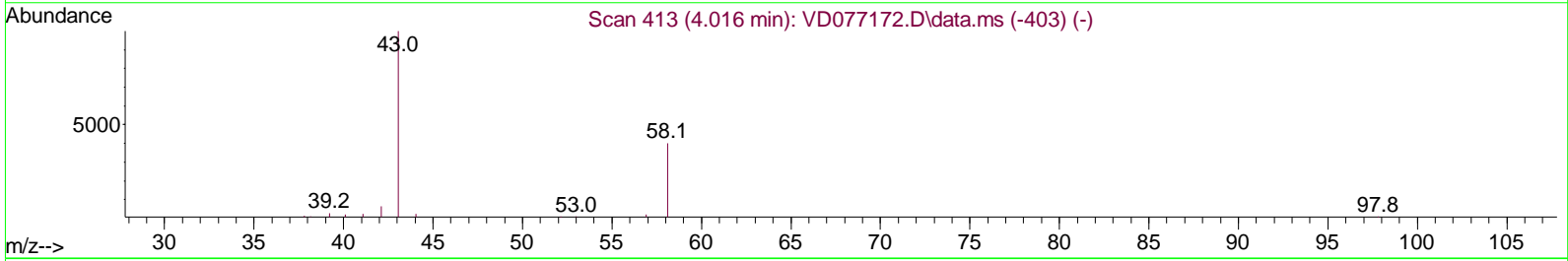
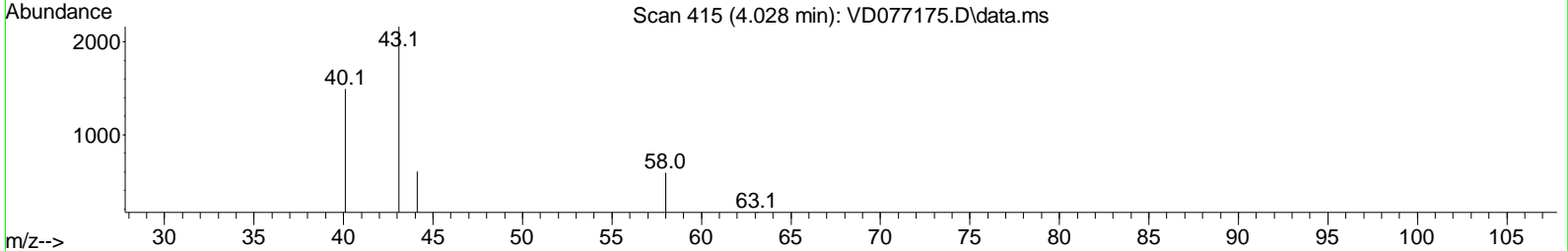
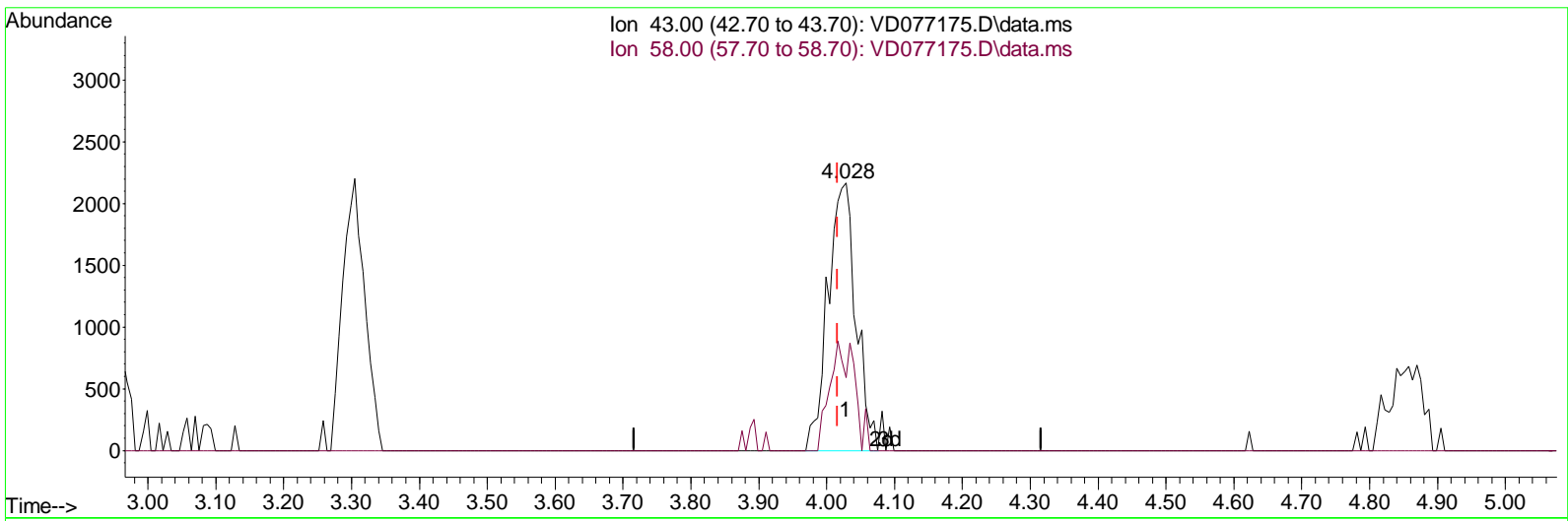
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TIC: VD077175.D\data.ms

(13) Acetone (T)

4.028min (+ 0.012) 5.07 ug/L m

response	6225	
Ion	Exp%	Act%
43.00	100.00	100.00
58.00	35.90	11.05
0.00	0.00	0.00
0.00	0.00	0.00

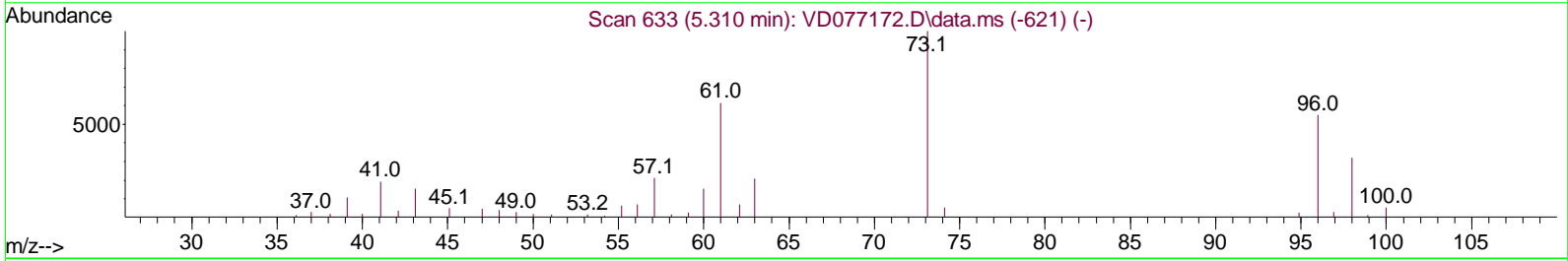
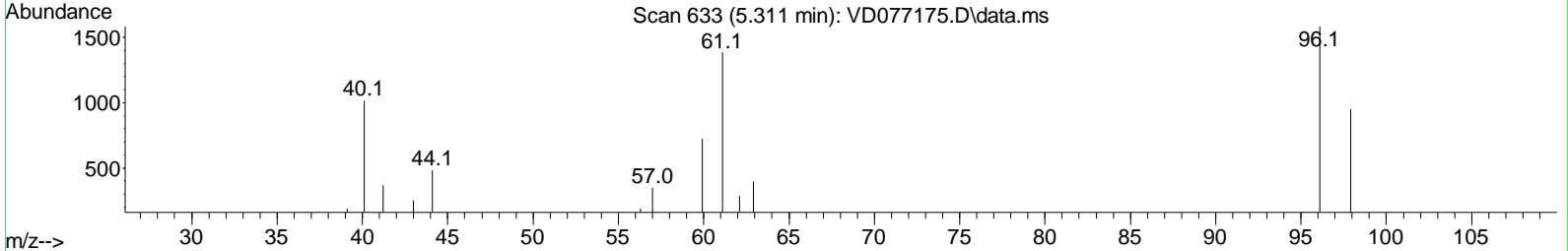
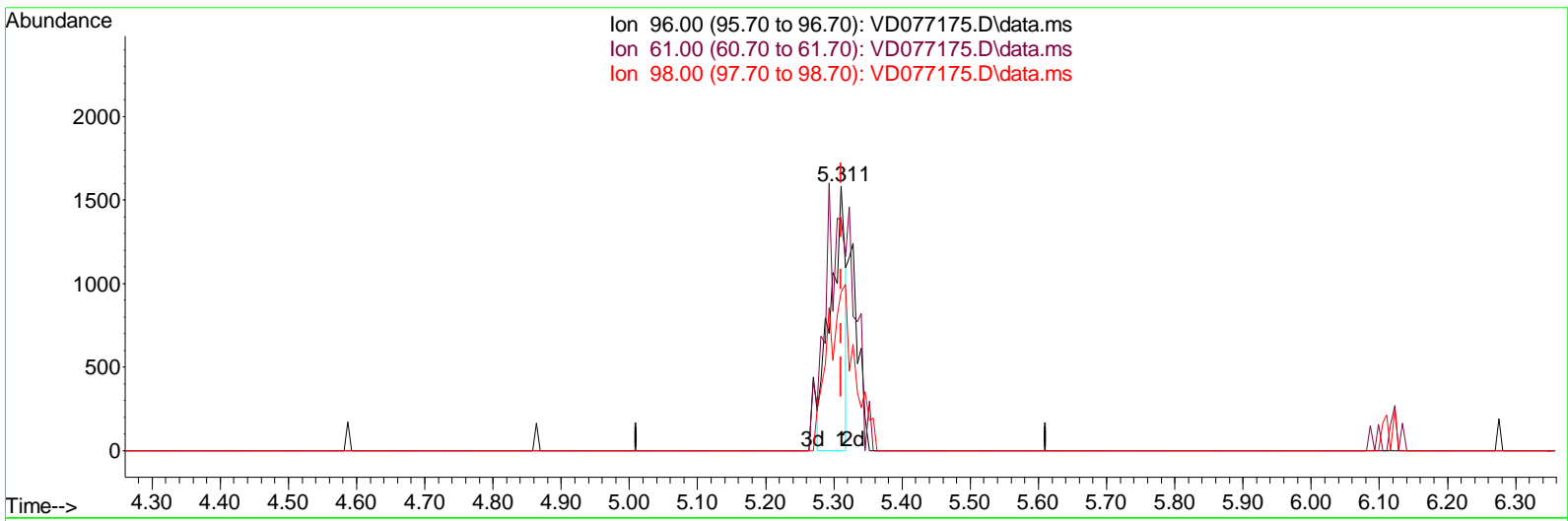
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 Operator : JC/SY
 Sample : 04527-02
 Misc : 5.70G/10.0ml/MSVOA_D/SOIL
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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TIC: VD077175.D\data.ms

(17) trans-1,2-Dichloroethene (T)

5.311min (+ 0.000) 0.58 ug/L

response	2348	
Ion	Exp%	Act%
96.00	100.00	100.00
61.00	121.60	87.43
98.00	62.40	59.95
0.00	0.00	0.00

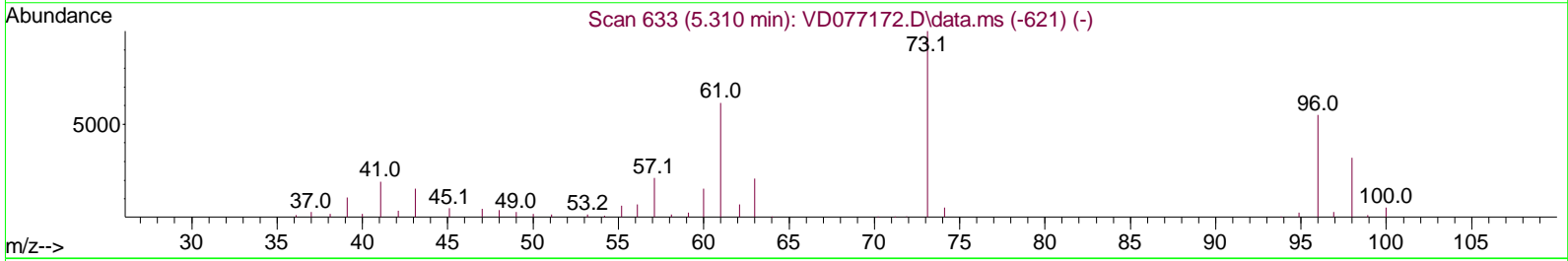
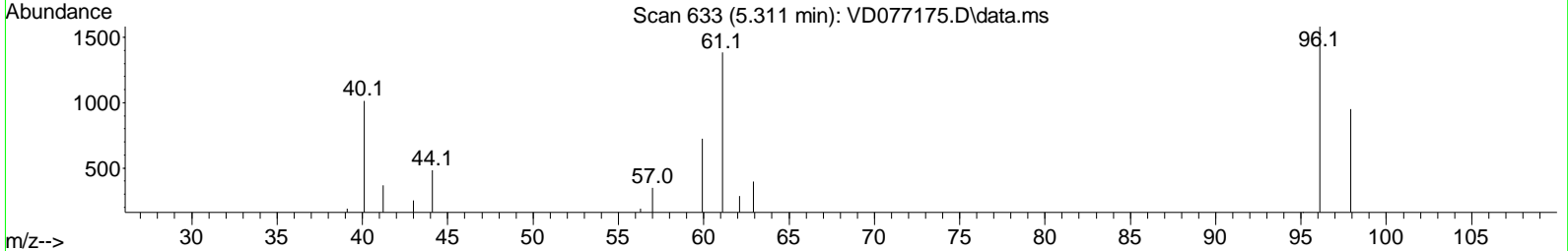
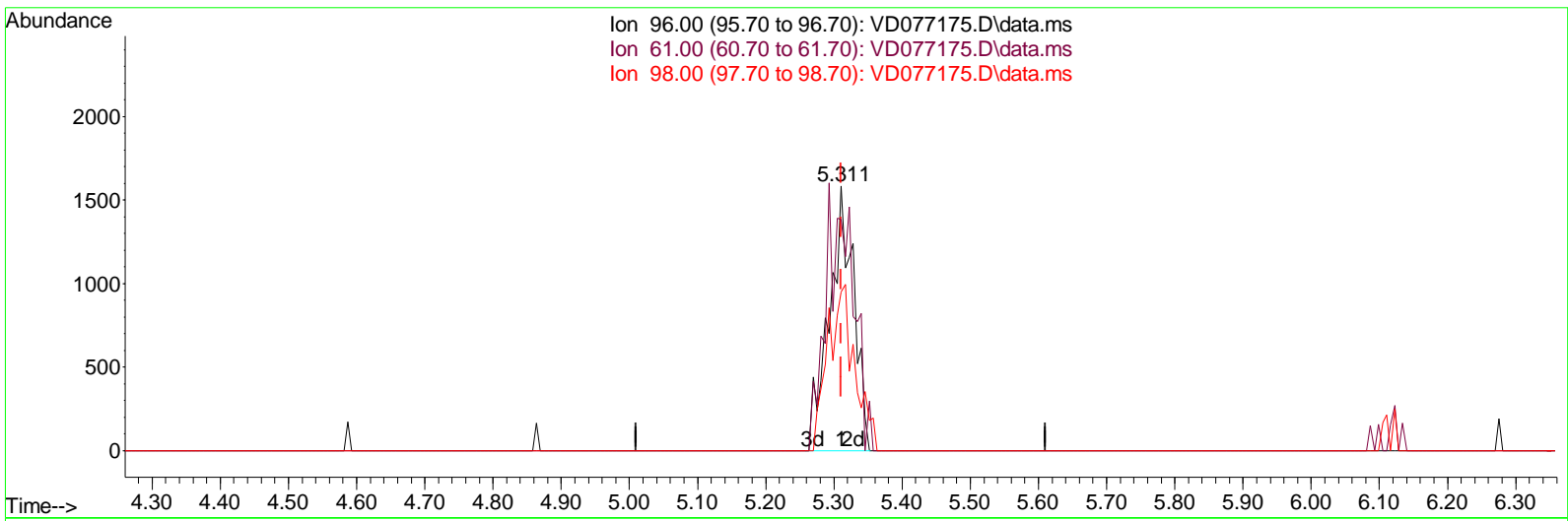
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TIC: VD077175.D\data.ms

(17) trans-1,2-Dichloroethene (T)

5.311min (+ 0.000) 0.97 ug/L m

response	3902	
Ion	Exp%	Act%
96.00	100.00	100.00
61.00	121.60	87.43
98.00	62.40	59.95
0.00	0.00	0.00

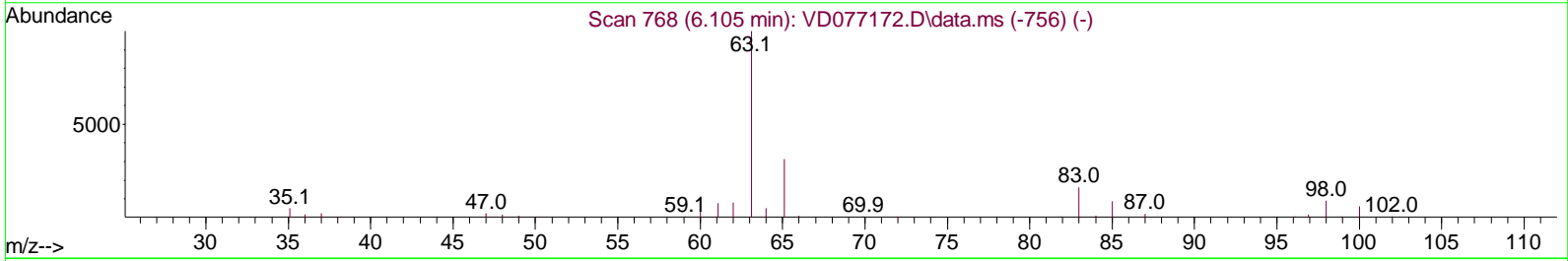
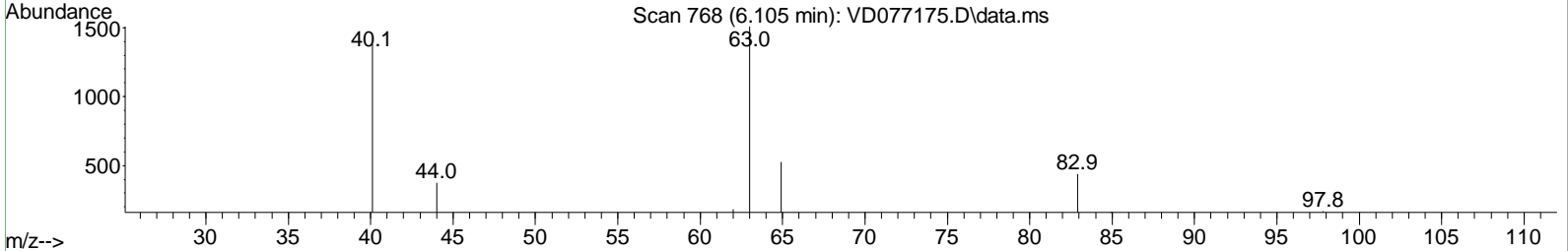
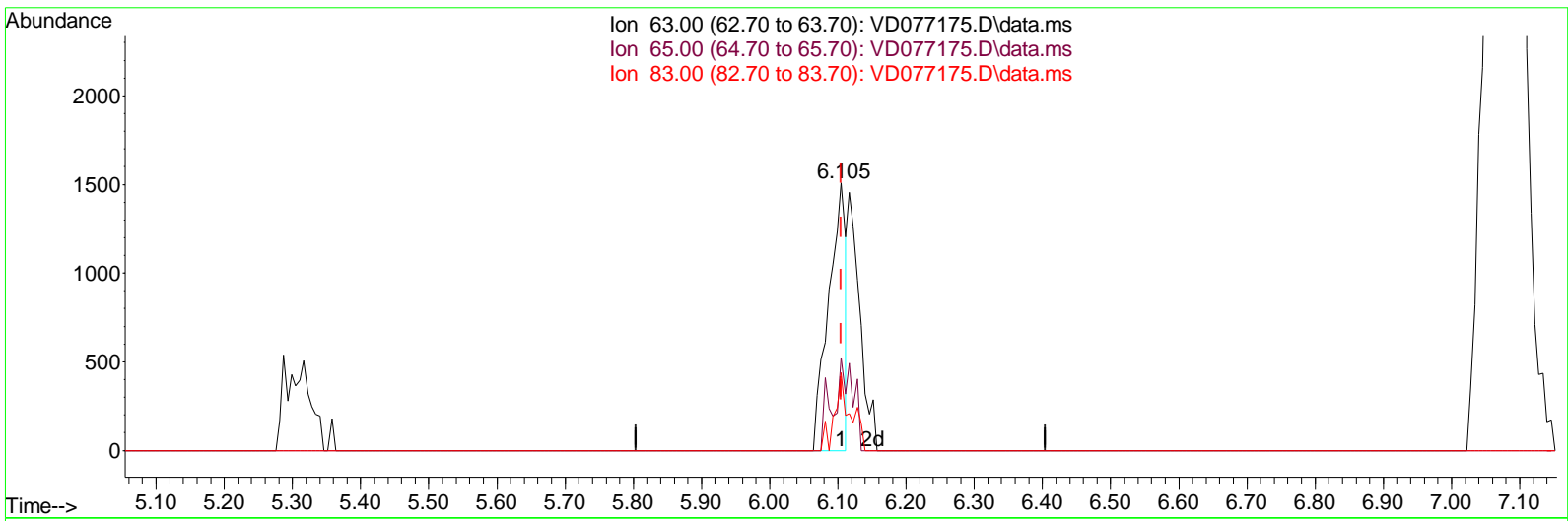
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TIC: VD077175.D\data.ms

(19) 1,1-Dichloroethane (T)

6.105min (+ 0.000) 0.38 ug/L

response	2589	
Ion	Exp%	Act%
63.00	100.00	100.00
65.00	32.20	34.81
83.00	14.20	29.11#
0.00	0.00	0.00

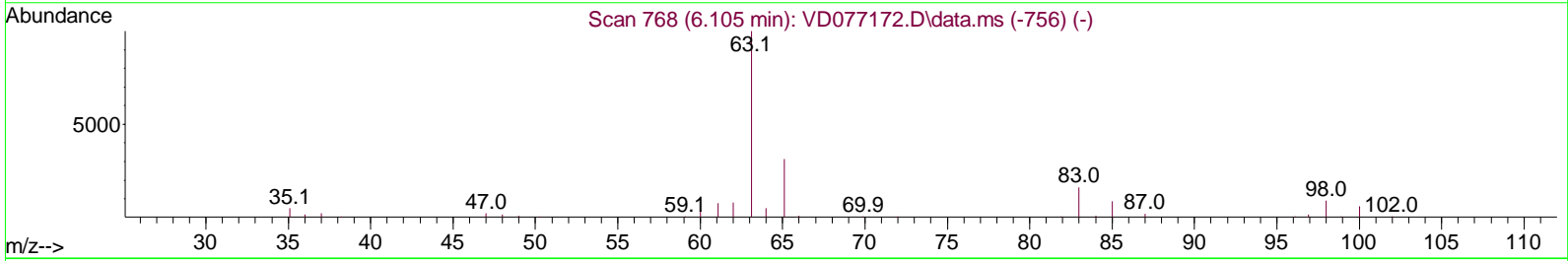
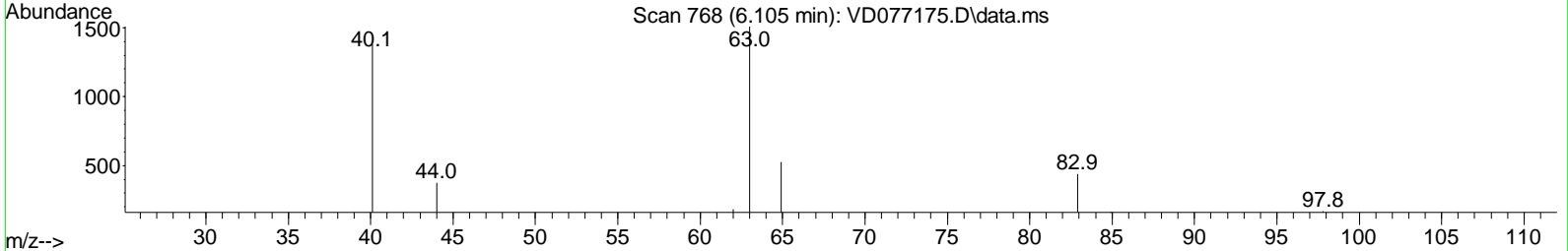
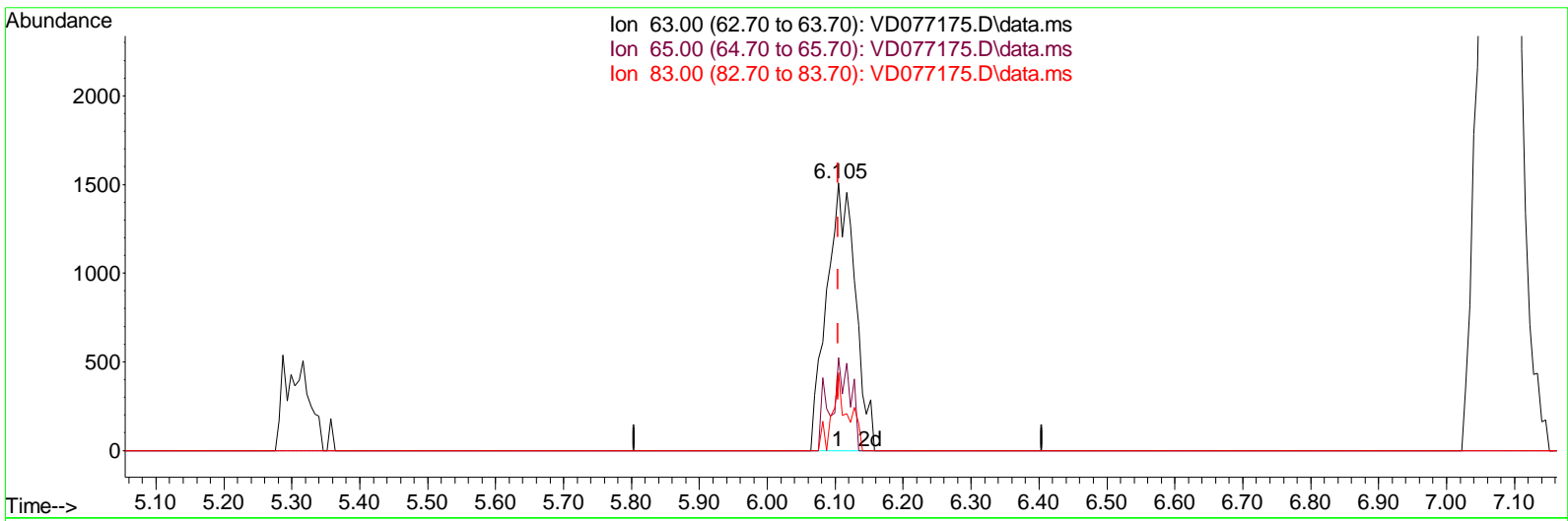
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 Operator : JC/SY
 Sample : 04527-02
 Misc : 5.70G/10.0ml/MSVOA_D/SOIL
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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TIC: VD077175.D\data.ms

(19) 1,1-Dichloroethane (T)

6.105min (+ 0.000) 0.65 ug/L m

response	4423	
Ion	Exp%	Act%
63.00	100.00	100.00
65.00	32.20	34.81
83.00	14.20	29.11#
0.00	0.00	0.00

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 Response via : Initial Calibrati on

Compound	R. T.	QI on	Response	Conc	Units	Dev(Mi n)
Internal Standards						
1) 1, 4-Di fl uorobenzene	8. 775	114	275866	25. 000	ug/L	0. 00
28) Chl orobenzene-d5	11. 581	117	272575	25. 000	ug/L	0. 00
58) 1, 4-Di chl orobenzene-d4	13. 516	152	126028	25. 000	ug/L	0. 00
System Moni toring Compounds						
4) Vi nyl Chl ori de-d3	2. 275	65	147377	21. 681	ug/L	0. 00
Spi ked Amount 25. 000	Range 30 - 150		Recovery =	86. 720%		
7) Chl oroethane-d5	2. 799	69	133039	23. 800	ug/L	0. 00
Spi ked Amount 25. 000	Range 30 - 150		Recovery =	95. 200%		
11) 1, 1-Di chl oroethene-d2	3. 911	65	36262	23. 484	ug/L	0. 00
Spi ked Amount 25. 000	Range 45 - 110		Recovery =	93. 920%		
21) 2-Butanone-d5	6. 981	46	48136	56. 211	ug/L	0. 00
Spi ked Amount 50. 000	Range 20 - 135		Recovery =	112. 420%		
24) Chl oroform-d	7. 563	84	150645	19. 957	ug/L	0. 00
Spi ked Amount 25. 000	Range 40 - 150		Recovery =	79. 840%		
26) 1, 2-Di chl oroethane-d4	8. 228	65	98669	27. 579	ug/L	0. 00
Spi ked Amount 25. 000	Range 70 - 130		Recovery =	110. 320%		
32) Benzene-d6	8. 199	84	375514	24. 943	ug/L	0. 00
Spi ked Amount 25. 000	Range 20 - 135		Recovery =	99. 760%		
36) 1, 2-Di chl oropropane-d6	9. 210	67	118415	27. 428	ug/L	0. 00
Spi ked Amount 25. 000	Range 70 - 120		Recovery =	109. 720%		
41) Tol uene-d8	10. 269	98	326619	23. 701	ug/L	0. 00
Spi ked Amount 25. 000	Range 30 - 130		Recovery =	94. 800%		
43) trans-1, 3-Di chl oroprop. . .	10. 528	79	47398	24. 095	ug/L	0. 00
Spi ked Amount 25. 000	Range 30 - 135		Recovery =	96. 360%		
47) 2-Hexanone-d5	10. 875	63	37878	54. 932	ug/L	0. 00
Spi ked Amount 50. 000	Range 20 - 135		Recovery =	109. 860%		
56) 1, 1, 2, 2-Tetrachl oroeth. . .	12. 651	84	52590	13. 664	ug/L	0. 00
Spi ked Amount 25. 000	Range 45 - 120		Recovery =	54. 640%		
66) 1, 2-Di chl orobenzene-d4	13. 816	152	103978	25. 704	ug/L	0. 00
Spi ked Amount 25. 000	Range 75 - 120		Recovery =	102. 800%		
Target Compounds						
13) Acetone	4. 028	43	6225m	5. 067	ug/L	
14) Carbon di sul fi de	4. 269	76	36959	3. 082	ug/L #	94
16) Methyl ene chl ori de	4. 805	84	38821	6. 914	ug/L	95
17) trans-1, 2-Di chl oroethene	5. 311	96	3902m	0. 967	ug/L	
19) 1, 1-Di chl oroethane	6. 105	63	4423m	0. 650	ug/L	
20) ci s-1, 2-Di chl oroethene	7. 075	96	89498	19. 381	ug/L	96
25) Chl oroform	7. 587	83	42427	5. 587	ug/L	96
34) Tri chl oroethene	9. 022	95	13282	3. 059	ug/L	89
42) Tol uene	10. 328	91	7808	0. 440	ug/L	98
46) Tetrachl oroethene	10. 810	164	108863	30. 672	ug/L	96
52) Ethyl benzene	11. 681	91	6372	0. 332	ug/L	96
53) m, p-Xyl ene	11. 787	106	7825	1. 056	ug/L	75
54) o-Xyl ene	12. 116	106	4517	0. 631	ug/L	82
60) I sopropyl benzene	12. 416	105	4670	0. 288	ug/L #	81
62) 1, 3, 5-Tri methyl benzene	12. 898	105	21317	1. 697	ug/L	100
63) 1, 2, 4-Tri methyl benzene	13. 210	105	44715	3. 633	ug/L	96

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Compound	R.T.	QI on	Response	Conc	Units	Dev(Min)
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 (#) = qualifier out of range (m) = manual integration (+) = signals summed

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