

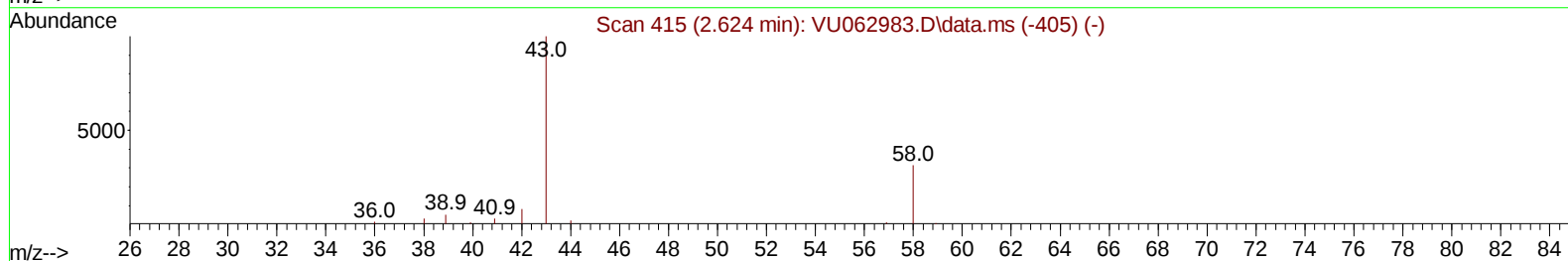
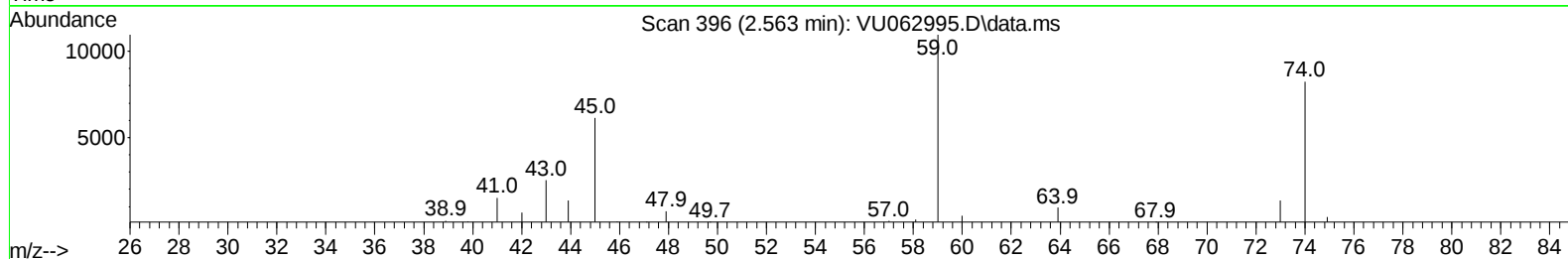
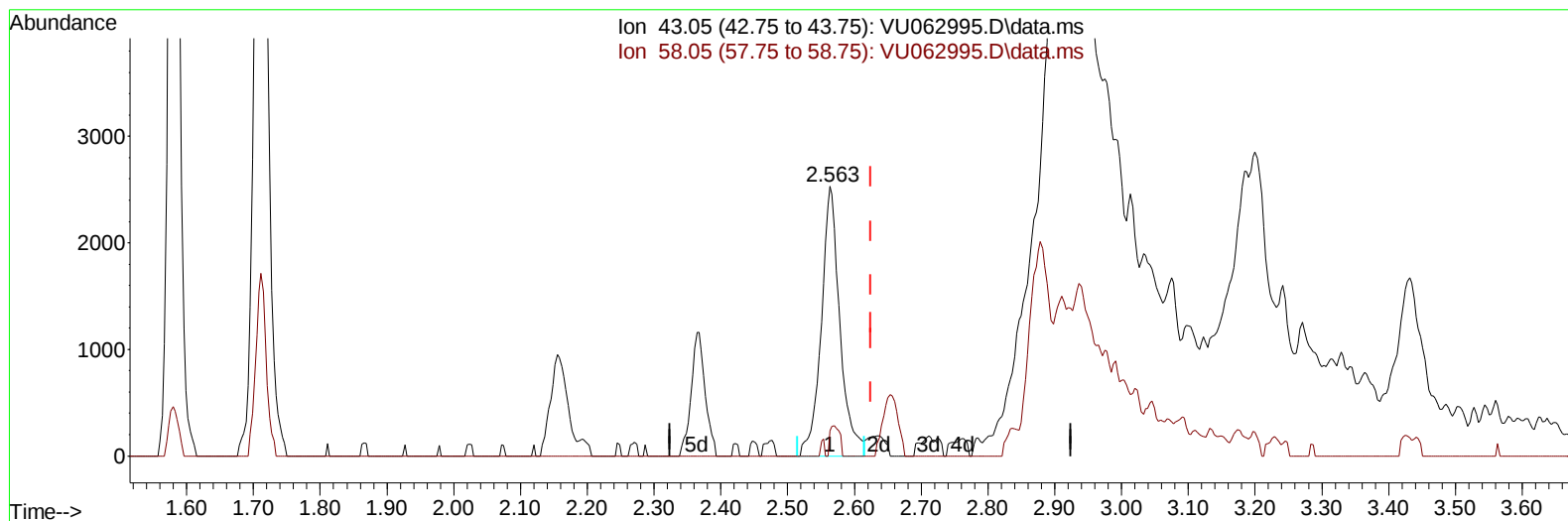
Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU012425\
Data File : VU062995.D
Acq On : 24 Jan 2025 16:03
Operator : MD/SY
Sample : Q1174-07
Misc : 25mL/MSVOA_U/WATER
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
E2937

Manual IntegrationsAPPROVED

Quant Time: Jan 24 19:07:05 2025
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMUTR010225WMA.M
Quant Title : TRACE VOA SFAM1.0
QLast Update : Fri Jan 24 19:02:27 2025
Response via : Initial Calibration

Reviewed By :John Carlone 01/27/2025
Supervised By :Mahesh Dadoda 01/27/2025



TIC: VU062995.D\data.ms

(13) Acetone (T)

2.563min (-0.061) 4.79 ug/L

response 4920

Ion	Exp%	Act%
43.05	100.00	100.00
58.05	32.20	5.83
0.00	0.00	0.00
0.00	0.00	0.00

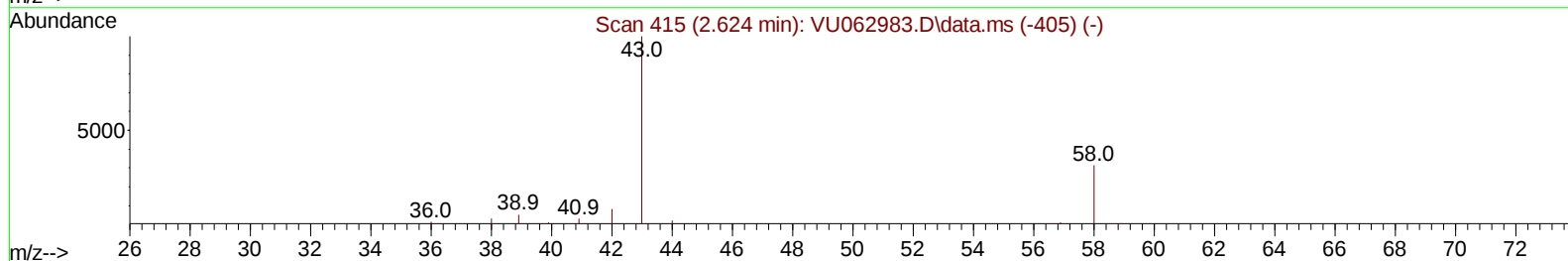
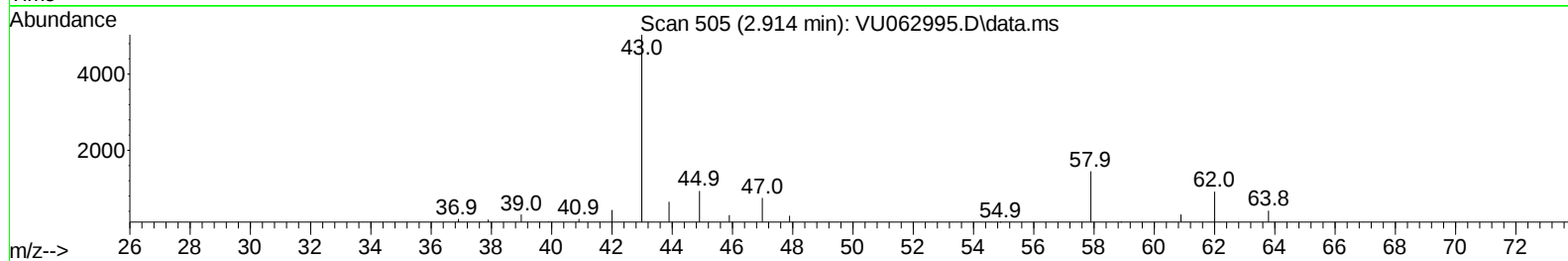
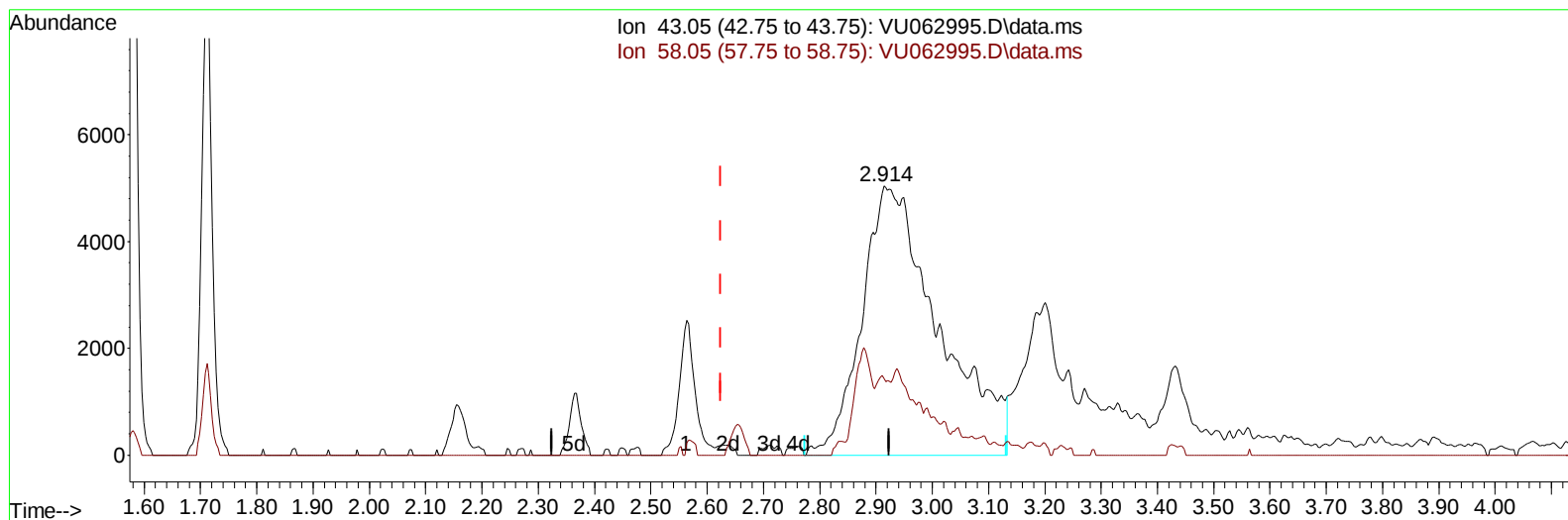
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(13) Acetone (T)

2.914min (+ 0.290) 46.29 ug/L m

response 47584

Ion	Exp%	Act%
43.05	100.00	100.00
58.05	32.20	0.60
0.00	0.00	0.00
0.00	0.00	0.00

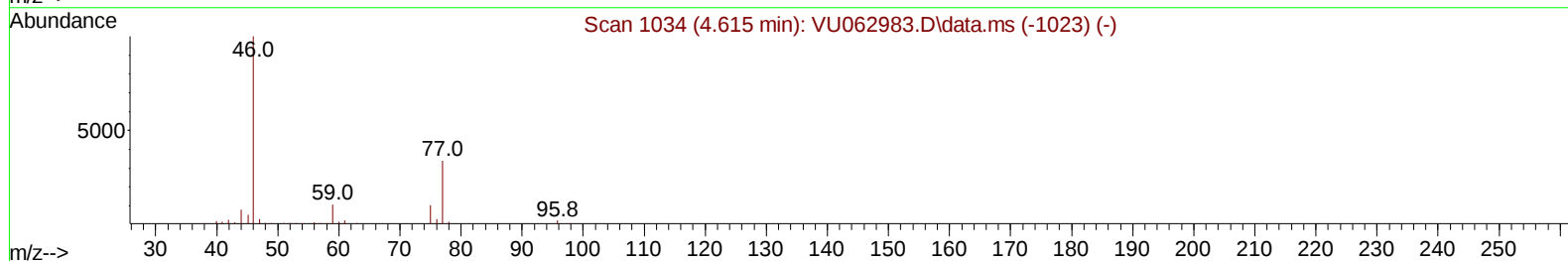
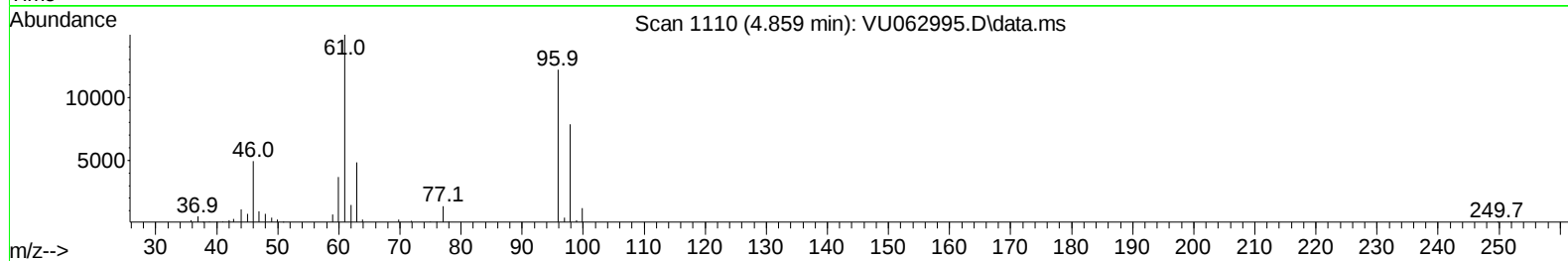
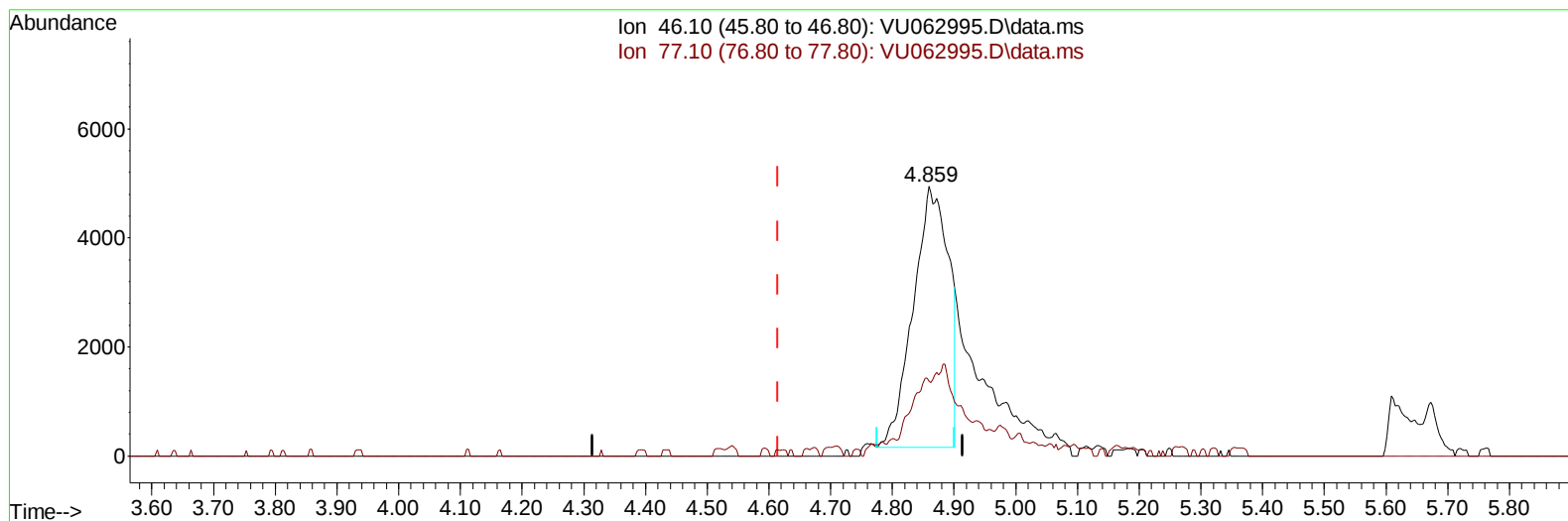
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TIC: VU062995.D\data.ms

(20) 2-Butanone-d5 (S)

4.859min (+ 0.245) 14.64 ug/L

response 18994

Ion	Exp%	Act%
46.10	100.00	100.00
77.10	30.90	12.39#
0.00	0.00	0.00
0.00	0.00	0.00

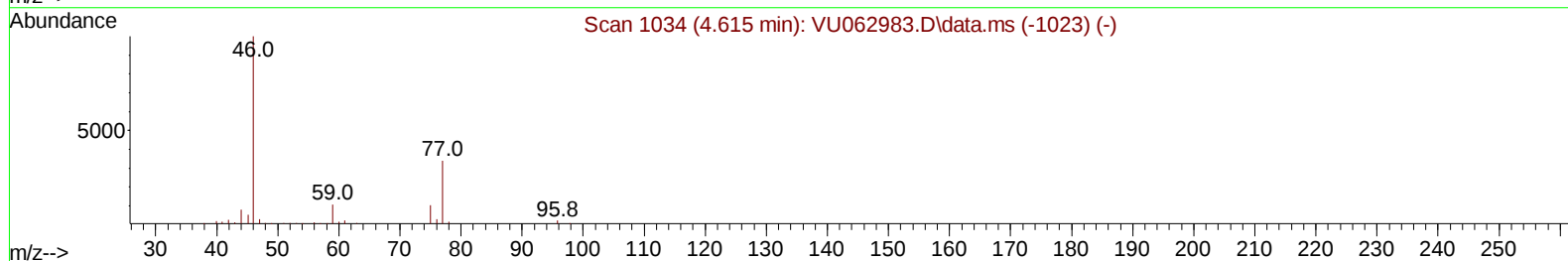
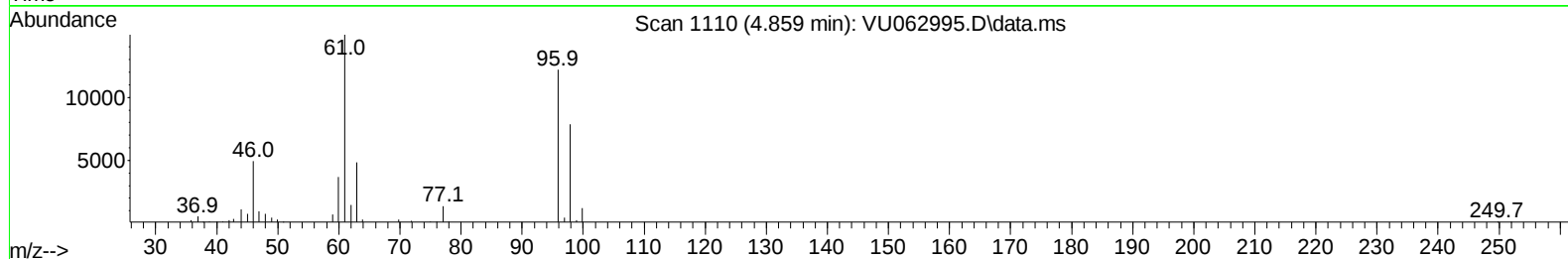
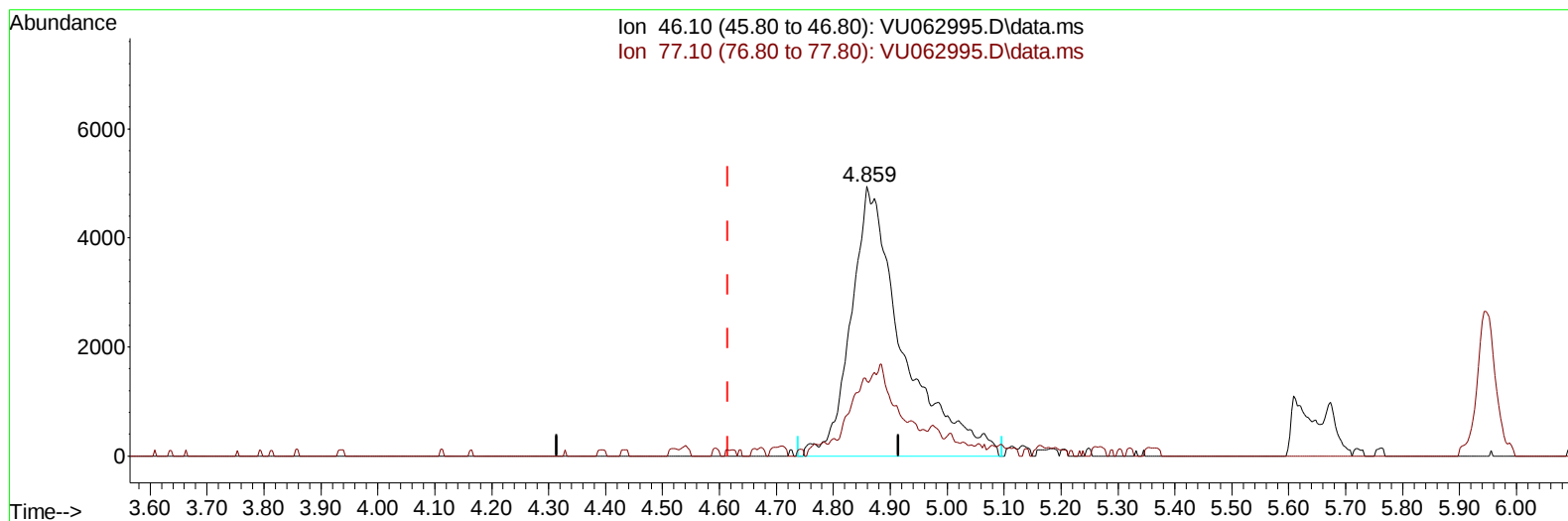
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TIC: VU062995.D\data.ms

(20) 2-Butanone-d5 (S)

4.859min (+ 0.245) 24.06 ug/L m

response 31215

Ion	Exp%	Act%
46.10	100.00	100.00
77.10	30.90	7.54#
0.00	0.00	0.00
0.00	0.00	0.00

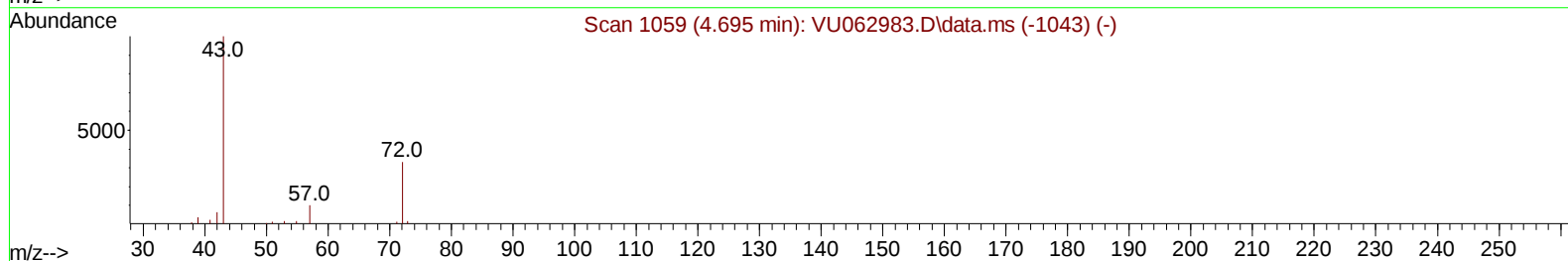
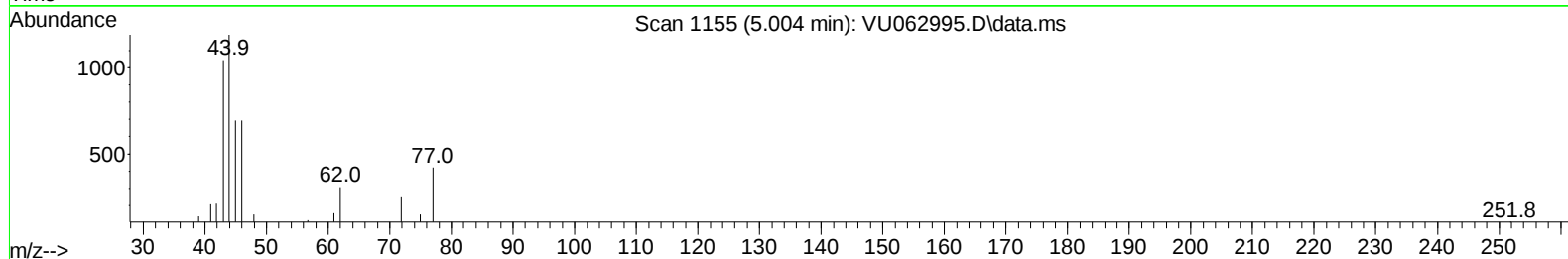
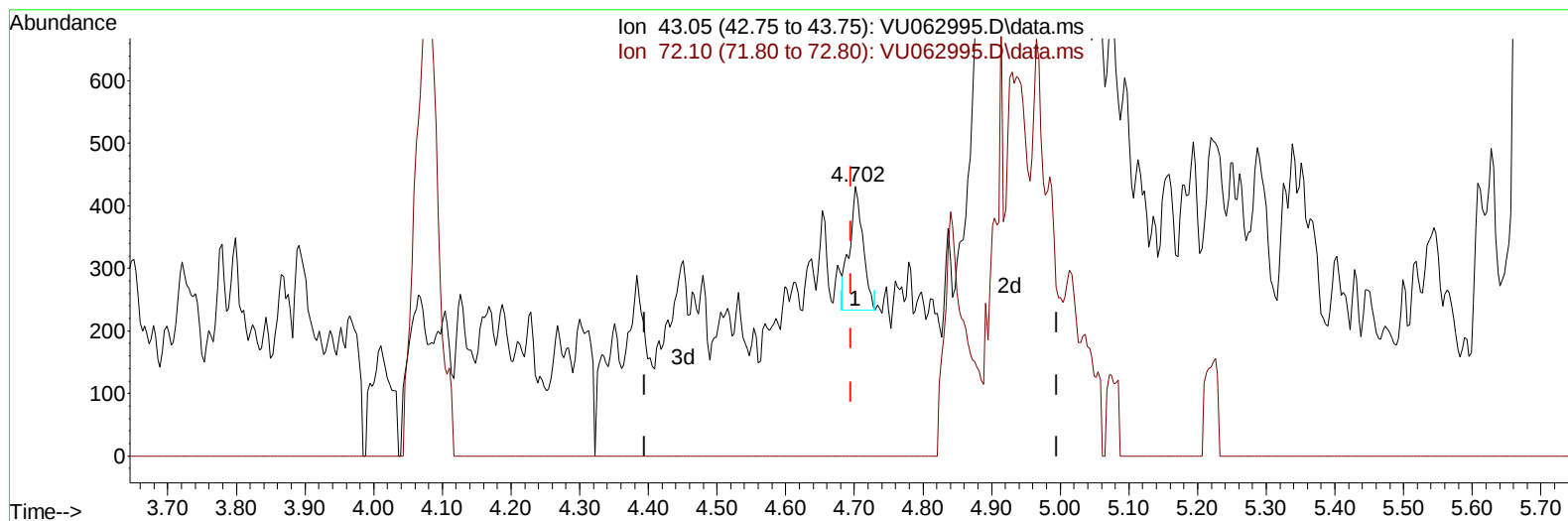
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Operator : MD/SY
Sample : Q1174-07
Misc : 25mL/MSVOA_U/WATER
ALS Vial : 15 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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Manual IntegrationsAPPROVED

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TIC: VU062995.D\data.ms

(21) 2-Butanone (T)

4.702min (+ 0.007) 0.19 ug/L

response 265

Ion	Exp%	Act%
43.05	100.00	100.00
72.10	34.10	312.08#
0.00	0.00	0.00
0.00	0.00	0.00

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Instrument :
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Manual IntegrationsAPPROVED

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Difluorobenzene	6.386	114	114627	5.000	ug/L	0.15
28) Chlorobenzene-d5	9.444	117	104320	5.000	ug/L	0.04
58) 1,4-Dichlorobenzene-d4	11.811	152	48846	5.000	ug/L	0.00
System Monitoring Compounds						
4) Vinyl Chloride-d3	1.708	65	21909	3.451	ug/L	0.12
Spiked Amount 5.000	Range 40	- 130	Recovery	=	69.000%	
7) Chloroethane-d5	2.068	69	14484	2.836	ug/L	0.16
Spiked Amount 5.000	Range 65	- 130	Recovery	=	56.800%#	
11) 1,1-Dichloroethene-d2	2.747	65	10174	3.094	ug/L	0.19
Spiked Amount 5.000	Range 60	- 125	Recovery	=	61.800%	
20) 2-Butanone-d5	4.859	46	31215m	24.058	ug/L	0.24
Spiked Amount 50.000	Range 40	- 130	Recovery	=	48.120%	
24) Chloroform-d	5.216	84	46188	2.957	ug/L	0.17
Spiked Amount 5.000	Range 70	- 125	Recovery	=	59.200%#	
26) 1,2-Dichloroethane-d4	5.882	65	34267	4.045	ug/L	0.19
Spiked Amount 5.000	Range 70	- 130	Recovery	=	80.800%	
32) Benzene-d6	5.901	84	114980	4.509	ug/L	0.19
Spiked Amount 5.000	Range 70	- 125	Recovery	=	90.200%	
36) 1,2-Dichloropropane-d6	6.808	67	36348	5.141	ug/L	0.13
Spiked Amount 5.000	Range 60	- 140	Recovery	=	102.800%	
41) Toluene-d8	7.962	98	107001	4.307	ug/L	0.07
Spiked Amount 5.000	Range 70	- 130	Recovery	=	86.200%	
43) trans-1,3-Dichloroprop...	8.232	79	14089	3.998	ug/L	0.06
Spiked Amount 5.000	Range 55	- 130	Recovery	=	80.000%	
46) 2-Hexanone-d5	8.679	63	46429	41.887	ug/L	0.05
Spiked Amount 50.000	Range 45	- 130	Recovery	=	83.780%	
56) 1,1,2,2-Tetrachloroeth...	10.759	84	26726	4.514	ug/L	0.02
Spiked Amount 5.000	Range 65	- 120	Recovery	=	90.200%	
66) 1,2-Dichlorobenzene-d4	12.190	152	36140	4.447	ug/L	0.00
Spiked Amount 5.000	Range 80	- 120	Recovery	=	89.000%	
Target Compounds						Qvalue
5) vinyl chloride	1.715	62	34194	4.144	ug/L	95
8) Chloroethane	2.091	64	6996	1.411	ug/L	99
13) Acetone	2.914	43	47584m	46.294	ug/L	
14) Carbon disulfide	2.978	76	35309	1.554	ug/L	97
19) 1,1-Dichloroethane	4.075	63	793047	58.182	ug/L	96
21) 2-Butanone	4.959	43	18700m	13.661	ug/L	
22) cis-1,2-Dichloroethene	4.843	96	51353	6.186	ug/L	97
33) Benzene	5.949	78	25182	0.840	ug/L	100
40) 4-Methyl-2-pentanone	7.862	43	35497	10.748	ug/L #	93
42) Toluene	8.029	91	77368	2.356	ug/L	99
52) Ethylbenzene	9.592	91	48804	1.309	ug/L	96
53) m,p-Xylene	9.711	106	56449	4.050	ug/L	98
54) o-Xylene	10.116	106	17674	1.324	ug/L	99
63) 1,2,4-Trimethylbenzene	11.467	105	31691	1.145	ug/L	98

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

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