

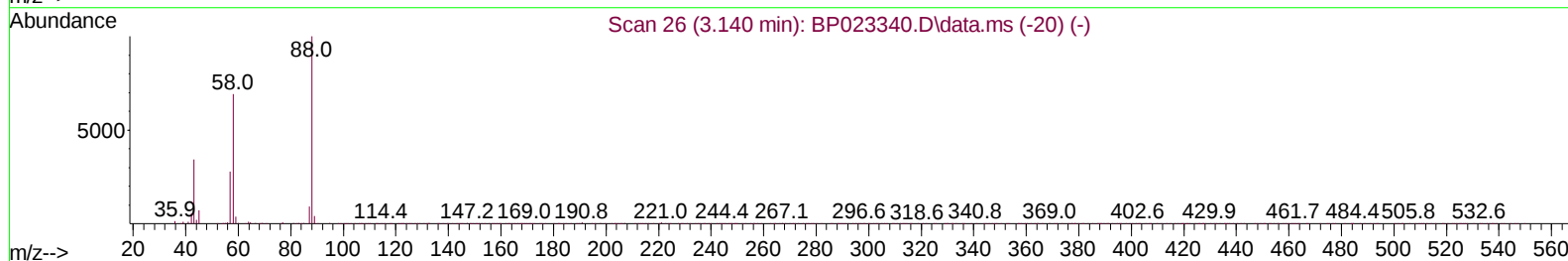
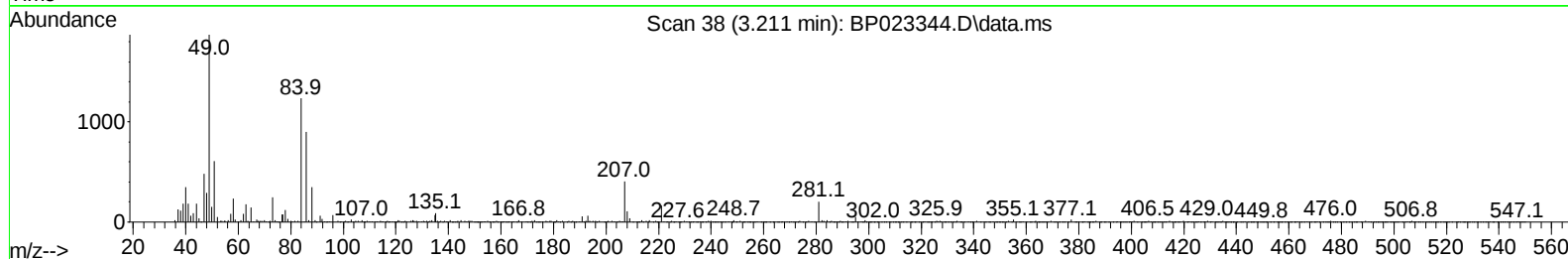
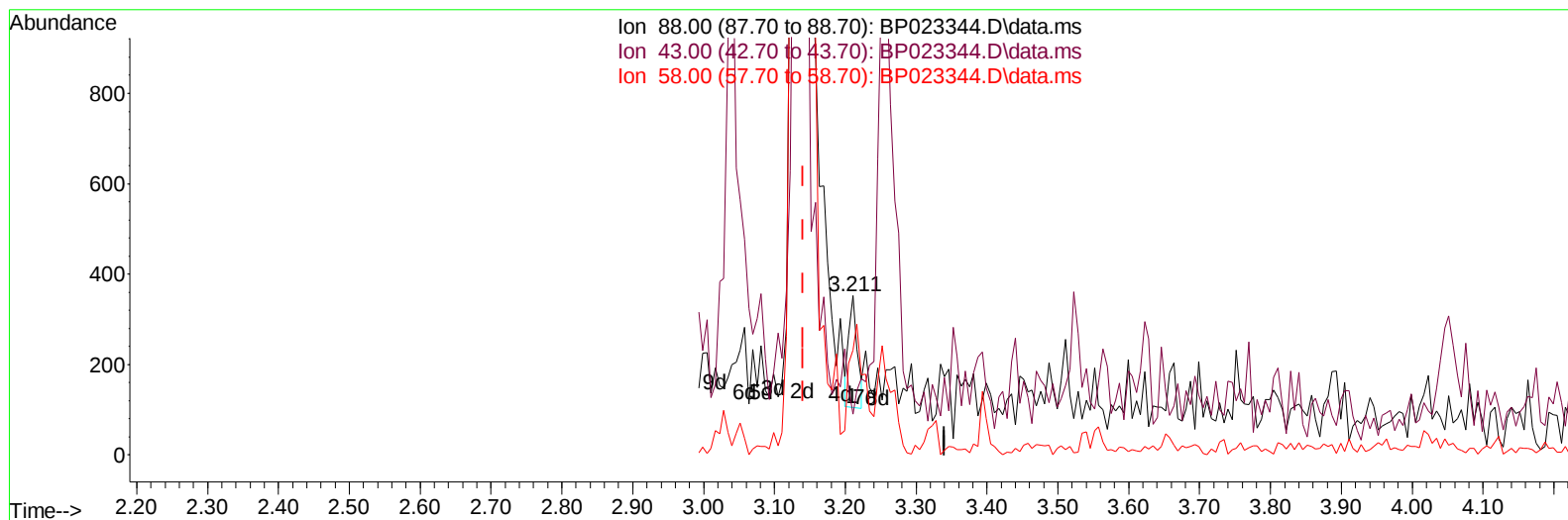
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121024\
Data File : BP023344.D
Acq On : 10 Dec 2024 13:11
Operator : RC/JU
Sample : P5107-21MS
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BHJ99MS

Manual IntegrationsAPPROVED

Quant Time: Dec 11 00:59:40 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112624.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Dec 11 00:30:30 2024
Response via : Initial Calibration

Reviewed By :Yogesh Patel 12/12/2024
Supervised By :mohammad ahmed 12/13/2024



TIC: BP023344.D\data.ms

(2) 1,4-Dioxane

3.211min (+ 0.070) 0.11 ng/uL

response 211

Ion	Exp%	Act%
88.00	100.00	100.00
43.00	28.60	25.57
58.00	66.40	65.91
0.00	0.00	0.00

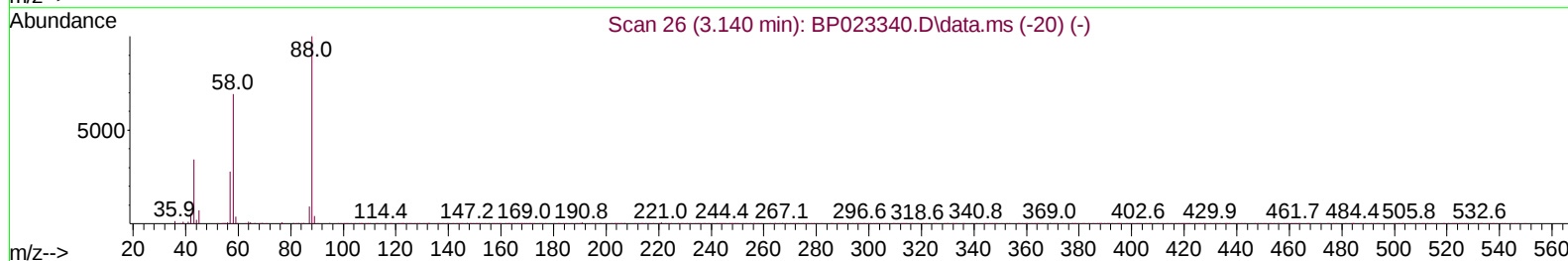
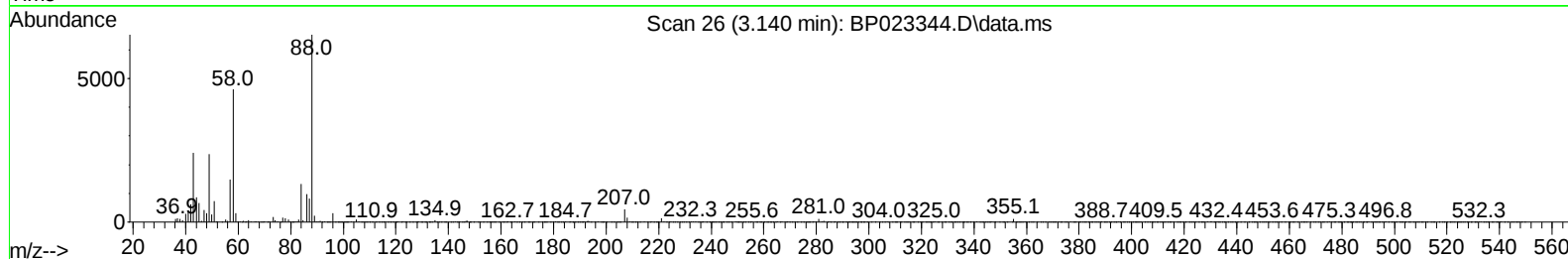
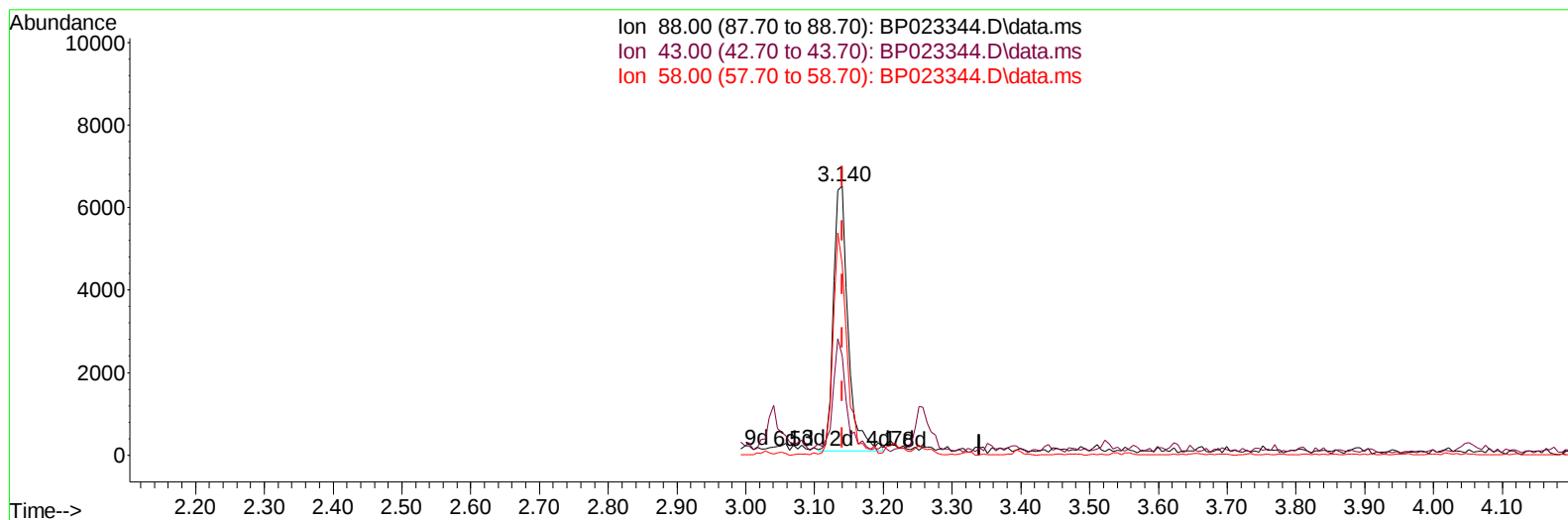
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121024\
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Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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Manual IntegrationsAPPROVED

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QLast Update : Wed Dec 11 00:30:30 2024
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TIC: BP023344.D\data.ms

(2) 1,4-Dioxane

3.140min (-0.000) 4.80 ng/uL m

response 9426

Ion	Exp%	Act%
88.00	100.00	100.00
43.00	28.60	36.96#
58.00	66.40	70.77
0.00	0.00	0.00

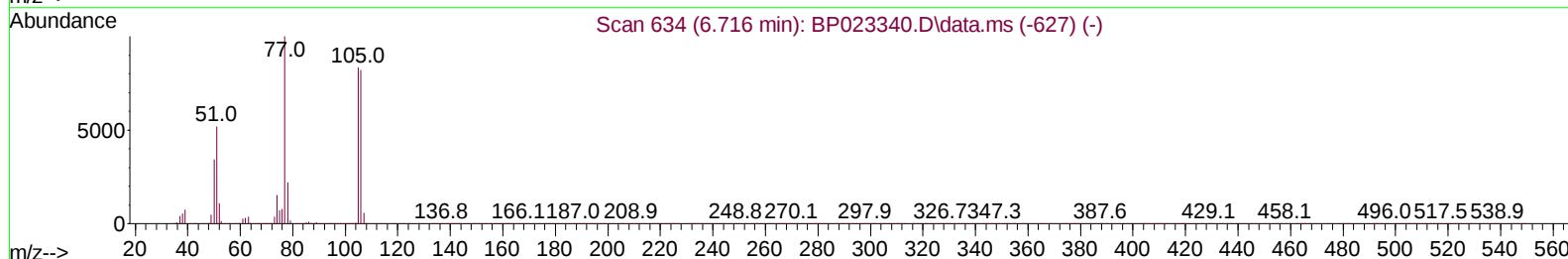
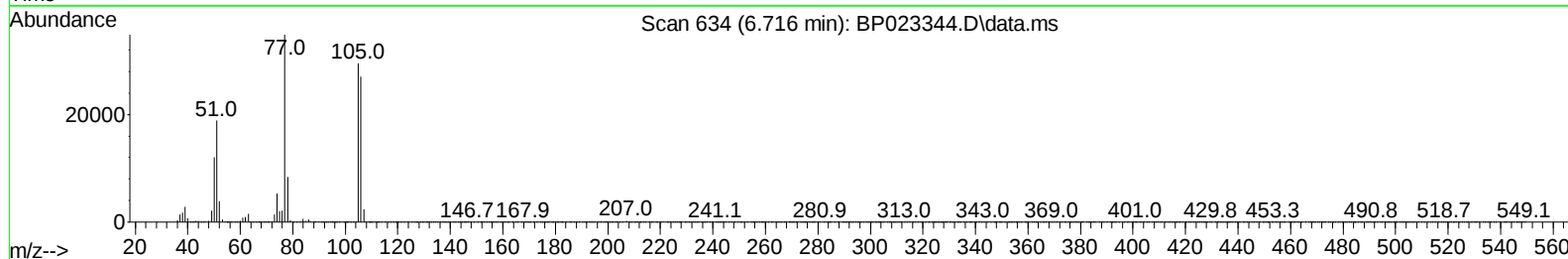
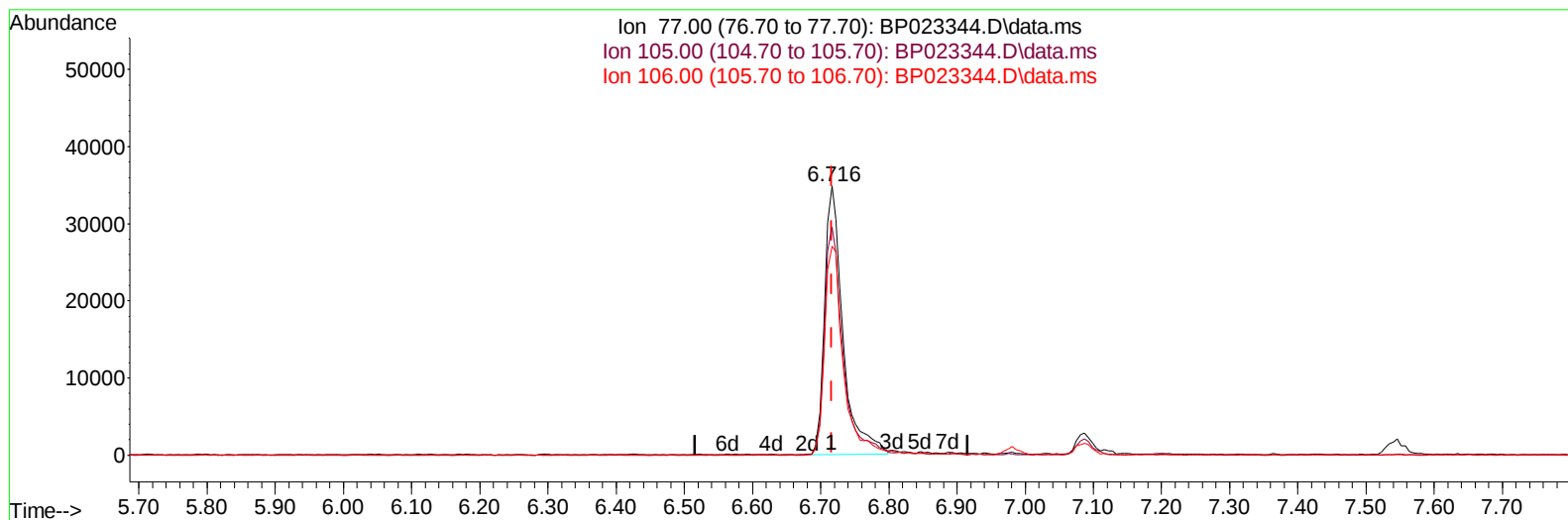
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121024\
 Data File : BP023344.D
 Acq On : 10 Dec 2024 13:11
 Operator : RC/JU
 Sample : P5107-21MS
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BHJ99MS

Manual IntegrationsAPPROVED

Quant Time: Dec 11 01:03:12 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112624.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Dec 11 00:30:30 2024
 Response via : Initial Calibration

Reviewed By :Yogesh Patel 12/12/2024
 Supervised By :mohammad ahmed 12/13/2024



TIC: BP023344.D\data.ms

(6) Benzaldehyde

6.716min (-0.000) 22.21 ng/u1

response 65107

Ion	Exp%	Act%
77.00	100.00	100.00
105.00	88.30	84.81
106.00	84.20	77.63
0.00	0.00	0.00

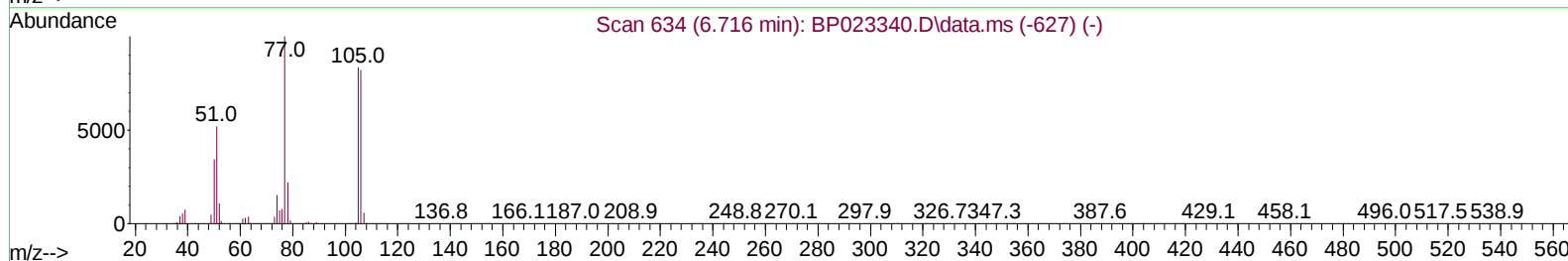
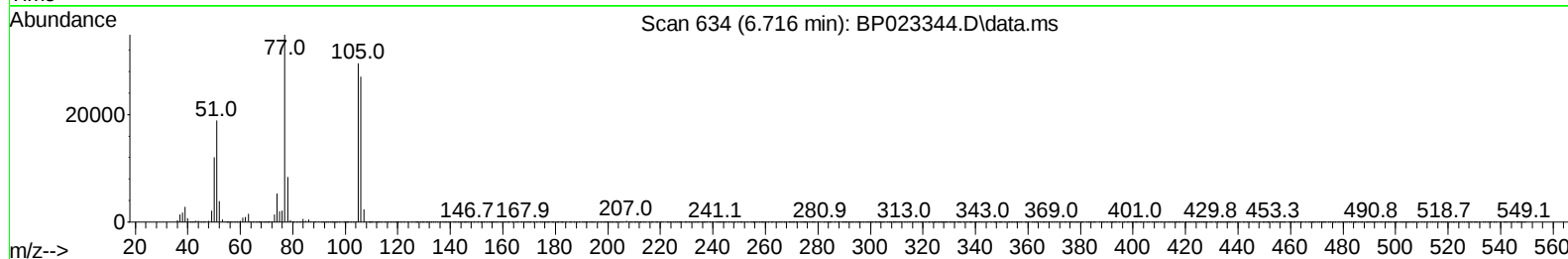
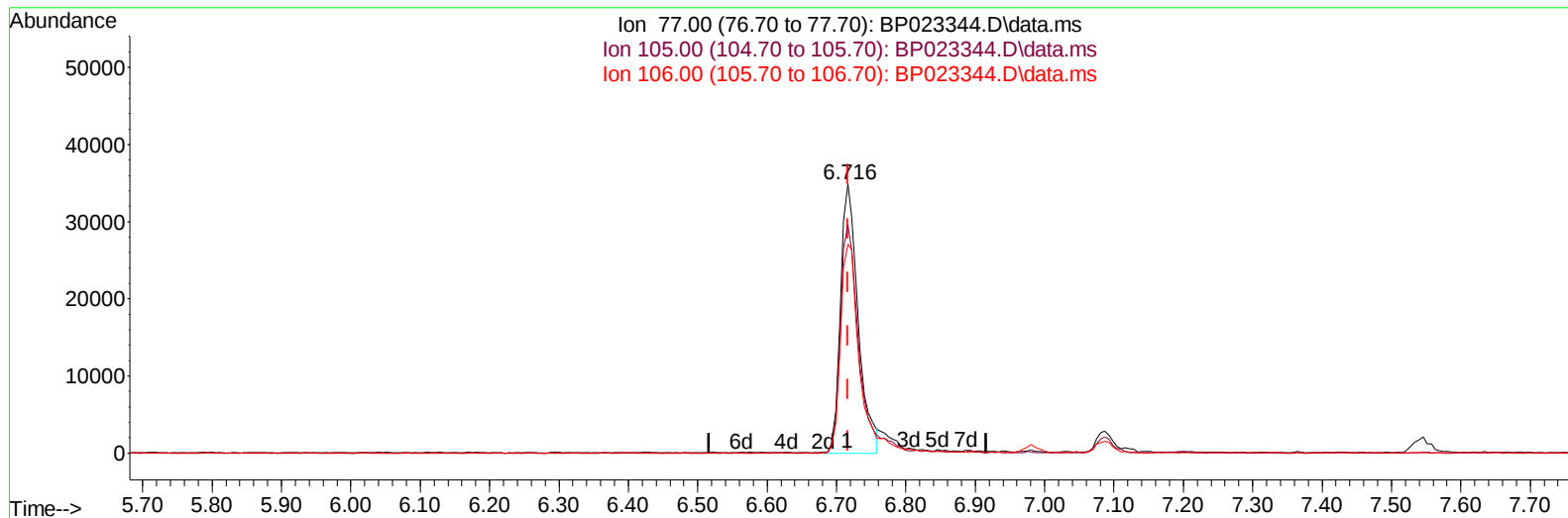
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121024\
Data File : BP023344.D
Acq On : 10 Dec 2024 13:11
Operator : RC/JU
Sample : P5107-21MS
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BHJ99MS

Manual IntegrationsAPPROVED

Quant Time: Dec 11 01:03:12 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112624.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Dec 11 00:30:30 2024
Response via : Initial Calibration

Reviewed By :Yogesh Patel 12/12/2024
Supervised By :mohammad ahmed 12/13/2024



TIC: BP023344.D\data.ms

(6) Benzaldehyde

6.716min (-0.000) 20.93 ng/u1 m

response 61369

Ion	Exp%	Act%
77.00	100.00	100.00
105.00	88.30	84.81
106.00	84.20	77.63
0.00	0.00	0.00

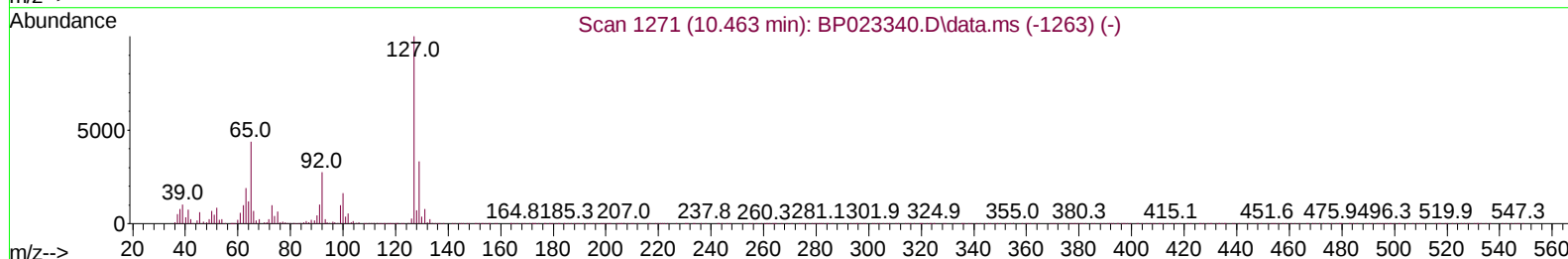
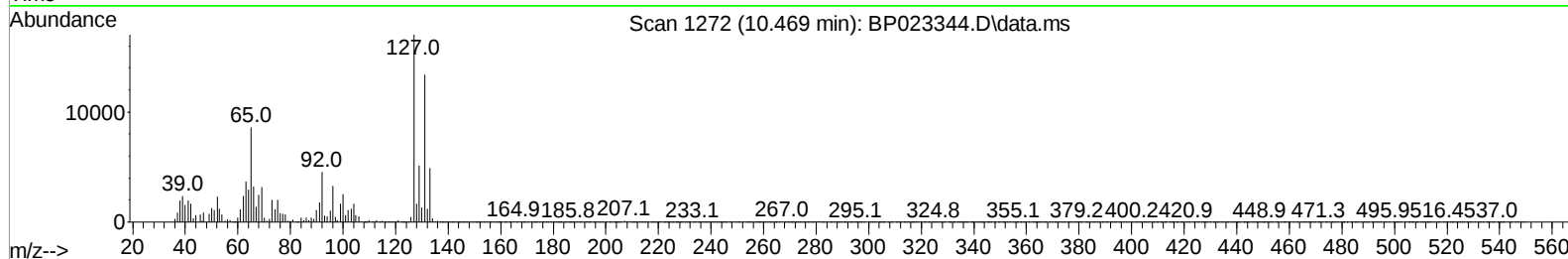
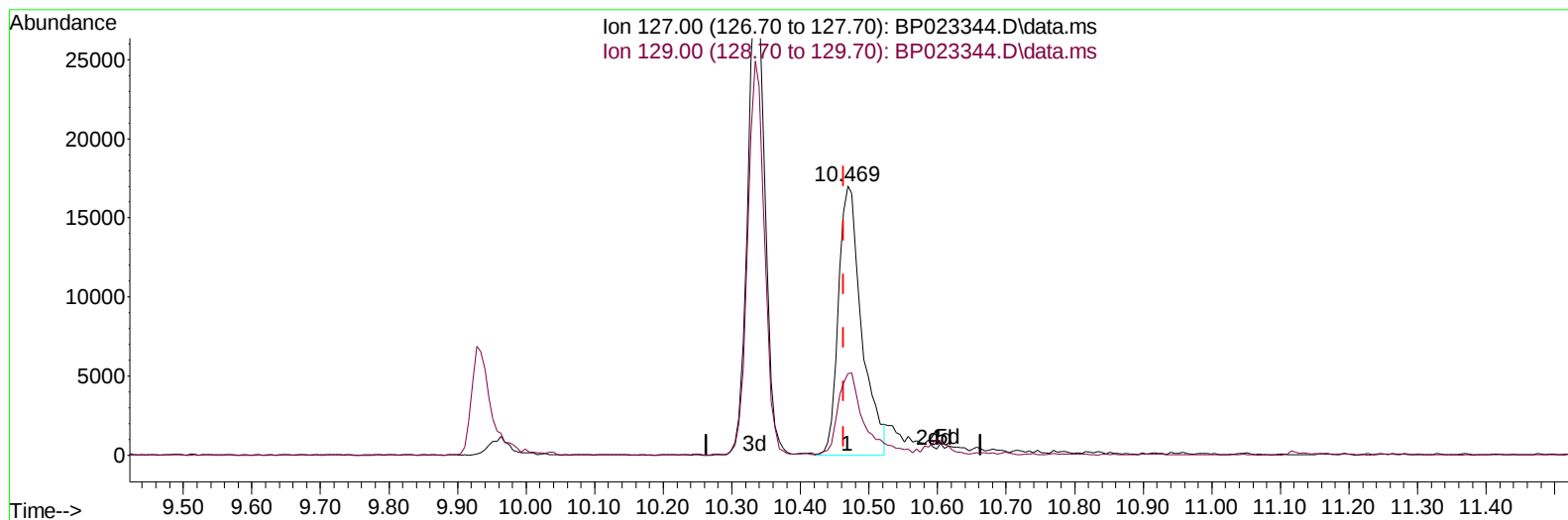
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121024\
Data File : BP023344.D
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Operator : RC/JU
Sample : P5107-21MS
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BHJ99MS

Manual IntegrationsAPPROVED

Quant Time: Dec 11 00:59:40 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112624.MA.M
Quant Title : SVOA CALIBRATION
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Reviewed By :Yogesh Patel 12/12/2024
Supervised By :mohammad ahmed 12/13/2024



TIC: BP023344.D\data.ms

(32) 4-Chloroaniline

10.469min (+ 0.006) 7.67 ng/ul

response 39879

Ion	Exp%	Act%
127.00	100.00	100.00
129.00	32.40	30.31
0.00	0.00	0.00
0.00	0.00	0.00

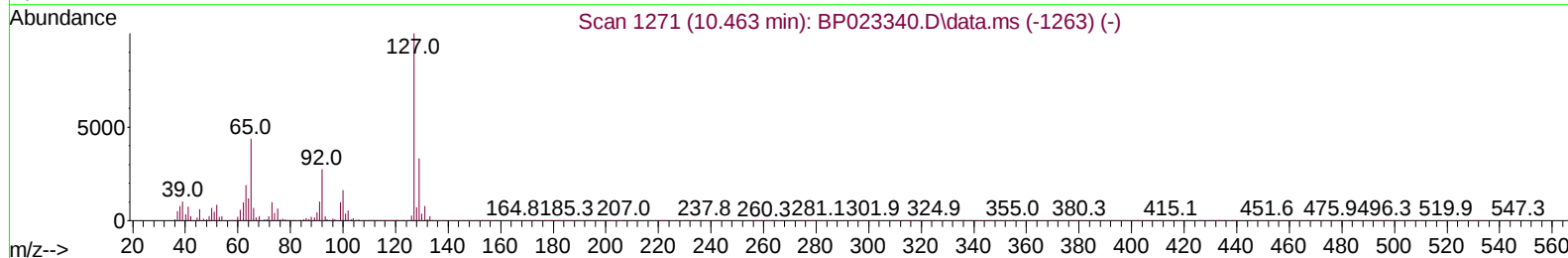
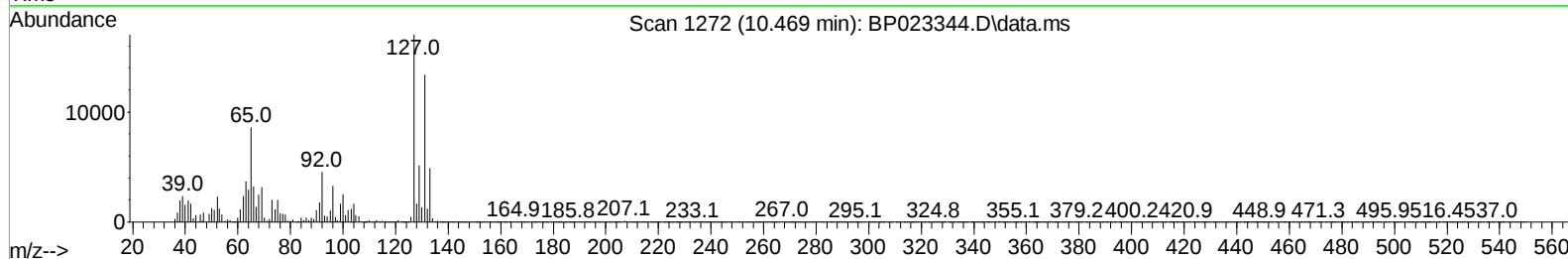
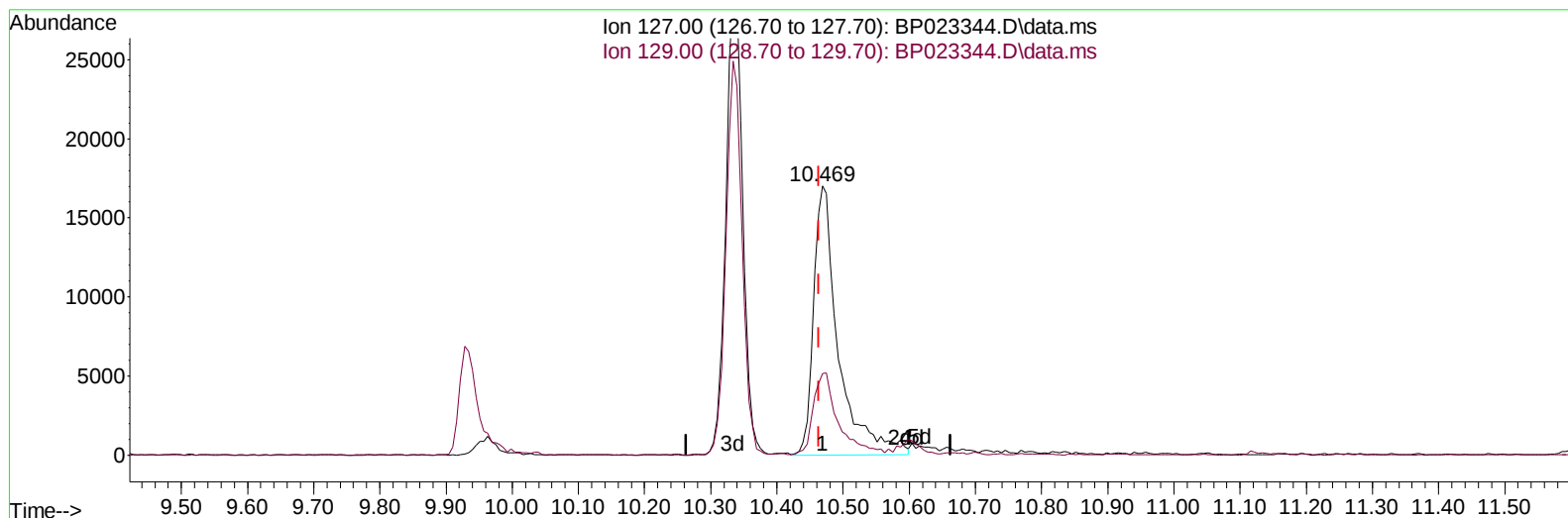
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121024\
Data File : BP023344.D
Acq On : 10 Dec 2024 13:11
Operator : RC/JU
Sample : P5107-21MS
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BHJ99MS

Manual IntegrationsAPPROVED

Quant Time: Dec 11 00:59:40 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112624.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Dec 11 00:30:30 2024
Response via : Initial Calibration

Reviewed By :Yogesh Patel 12/12/2024
Supervised By :mohammad ahmed 12/13/2024



TIC: BP023344.D\data.ms

(32) 4-Chloroaniline

10.469min (+ 0.006) 8.58 ng/ul m

response 44601

Ion	Exp%	Act%
127.00	100.00	100.00
129.00	32.40	30.31
0.00	0.00	0.00
0.00	0.00	0.00

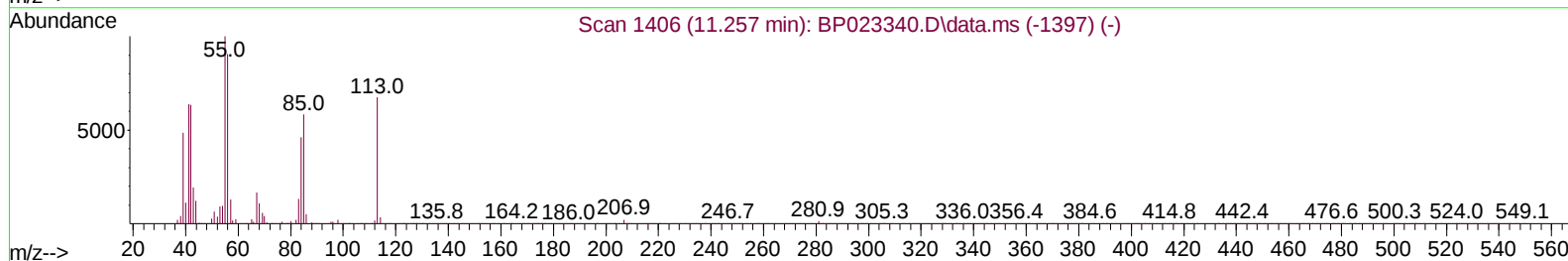
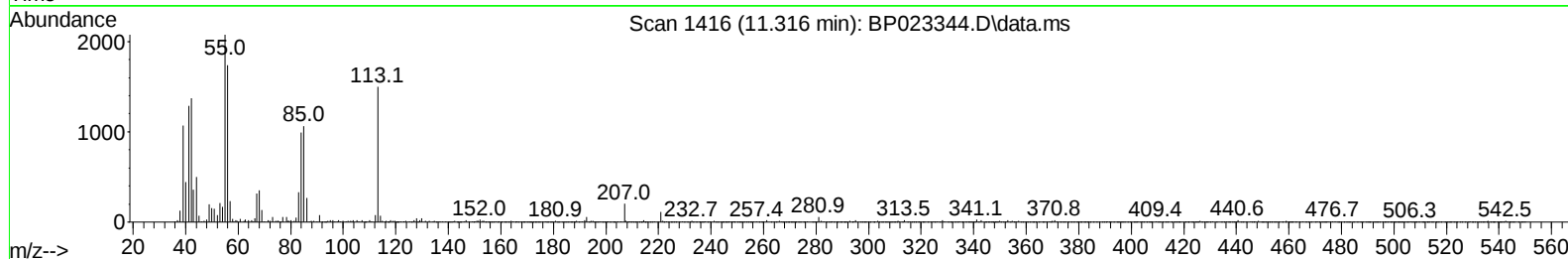
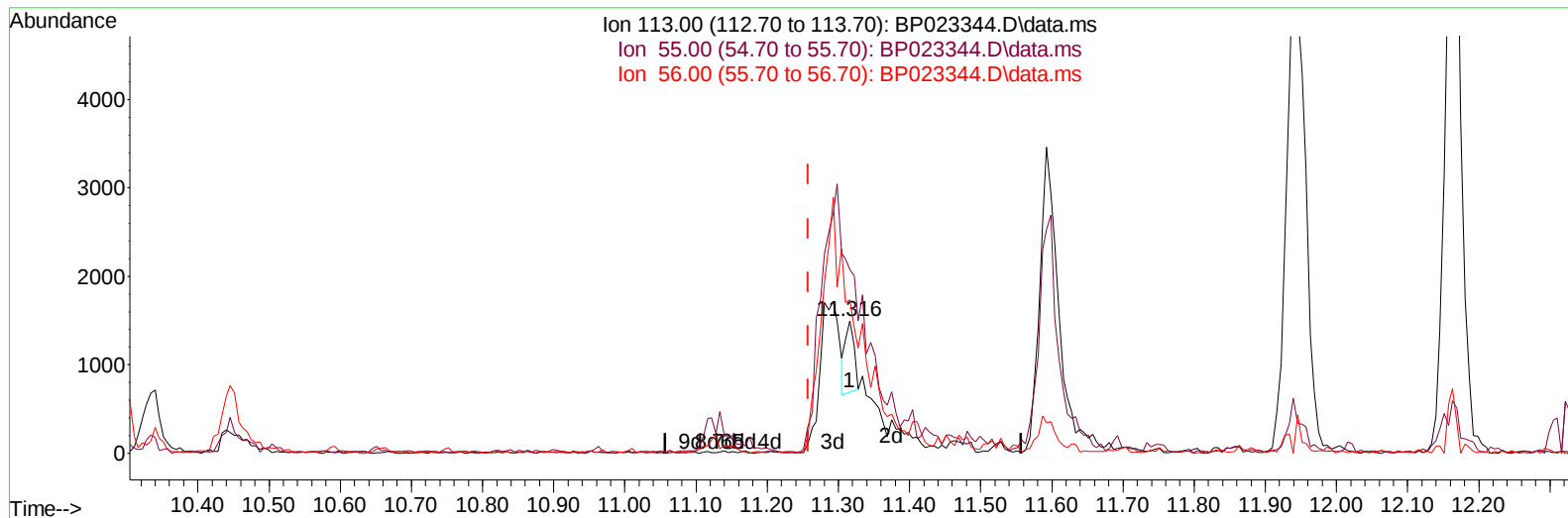
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Data File : BP023344.D
Acq On : 10 Dec 2024 13:11
Operator : RC/JU
Sample : P5107-21MS
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BHJ99MS

Manual IntegrationsAPPROVED

Quant Time: Dec 11 00:59:40 2024
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Reviewed By :Yogesh Patel 12/12/2024
Supervised By :mohammad ahmed 12/13/2024



TIC: BP023344.D\data.ms

(34) Caprolactam

11.316min (+ 0.059) 0.50 ng/ul

response 687

Ion	Exp%	Act%
113.00	100.00	100.00
55.00	159.40	138.61
56.00	125.90	115.97
0.00	0.00	0.00

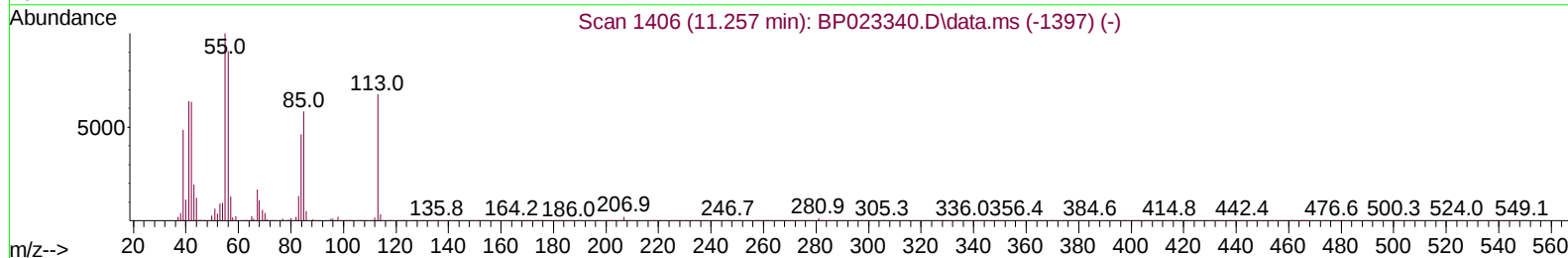
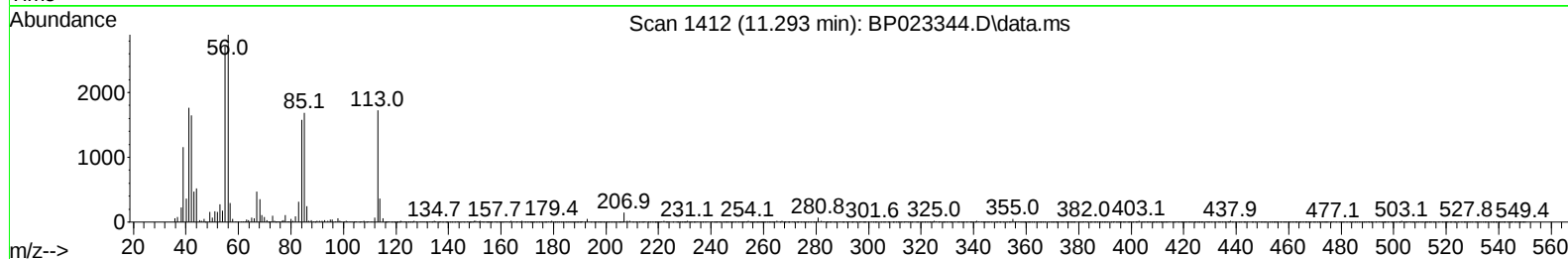
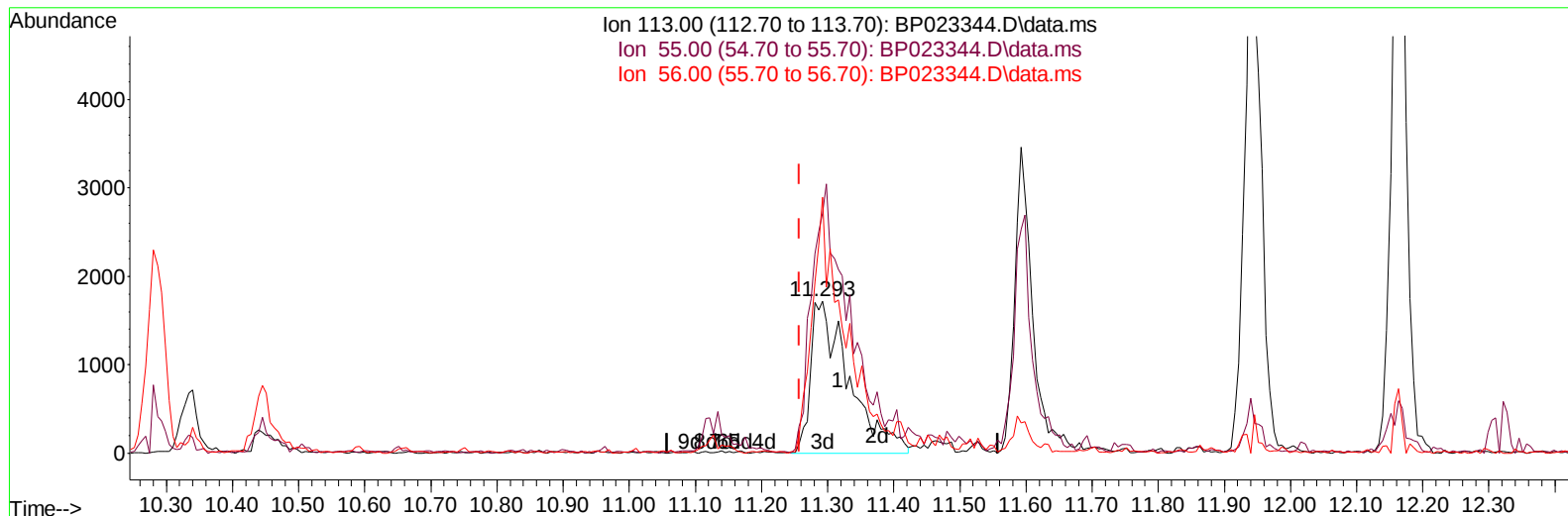
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Data File : BP023344.D
Acq On : 10 Dec 2024 13:11
Operator : RC/JU
Sample : P5107-21MS
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BHJ99MS

Manual IntegrationsAPPROVED

Quant Time: Dec 11 00:59:40 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112624.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Dec 11 00:30:30 2024
Response via : Initial Calibration

Reviewed By :Yogesh Patel 12/12/2024
Supervised By :mohammad ahmed 12/13/2024



TIC: BP023344.D\data.ms

(34) Caprolactam

11.293min (+ 0.035) 5.04 ng/ul m

response 6969

Ion	Exp%	Act%
113.00	100.00	100.00
55.00	159.40	157.72
56.00	125.90	168.00#
0.00	0.00	0.00

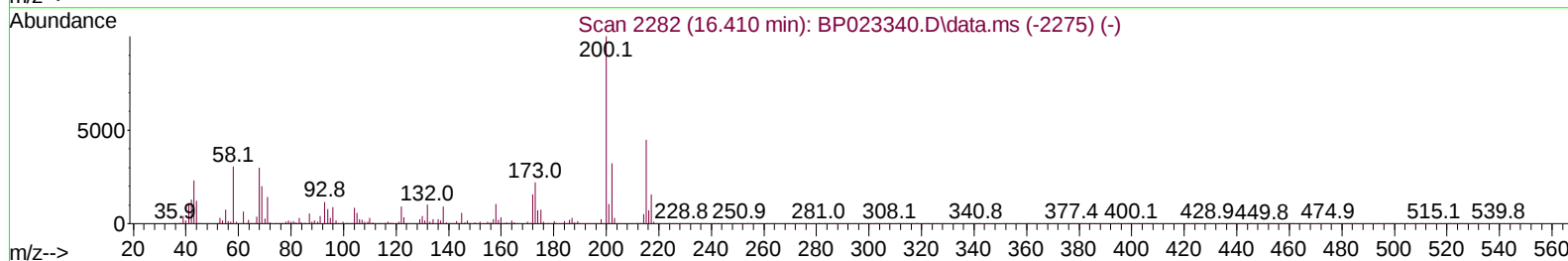
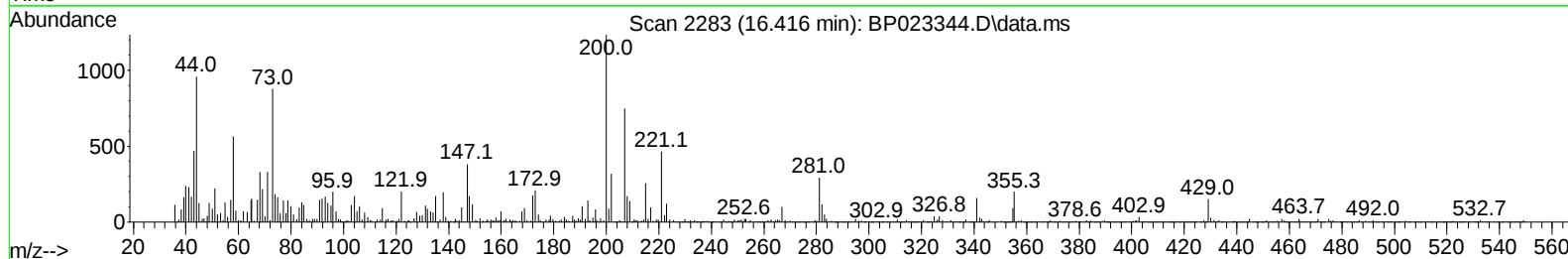
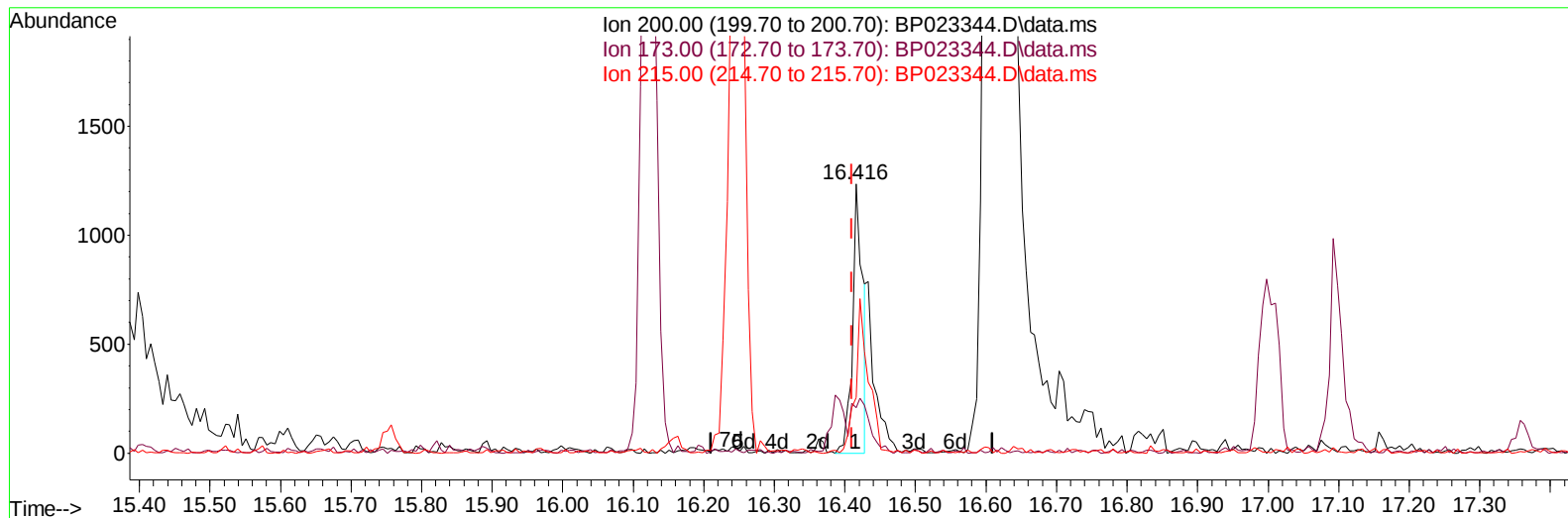
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Data File : BP023344.D
Acq On : 10 Dec 2024 13:11
Operator : RC/JU
Sample : P5107-21MS
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BHJ99MS

Manual IntegrationsAPPROVED

Quant Time: Dec 11 00:59:40 2024
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Reviewed By :Yogesh Patel 12/12/2024
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TIC: BP023344.D\data.ms

(70) Atrazine

16.416min (+ 0.006) 0.18 ng/u1

response 1241

Ion	Exp%	Act%
200.00	100.00	100.00
173.00	23.60	17.02#
215.00	47.50	20.58#
0.00	0.00	0.00

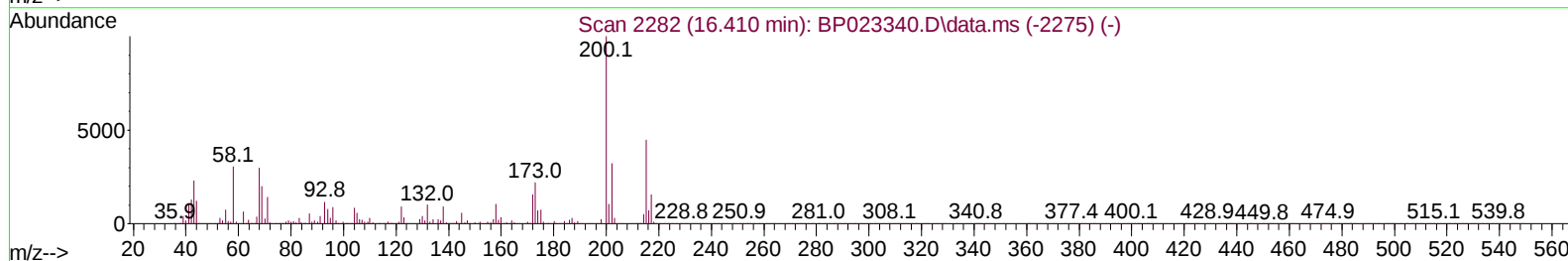
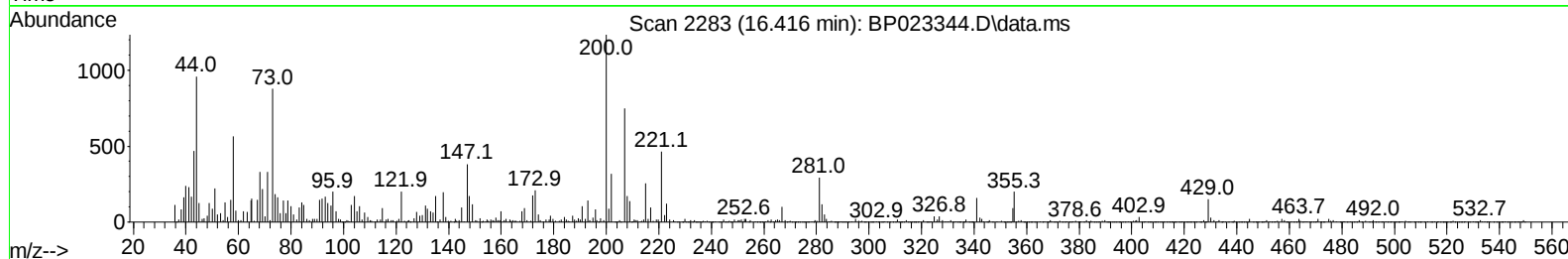
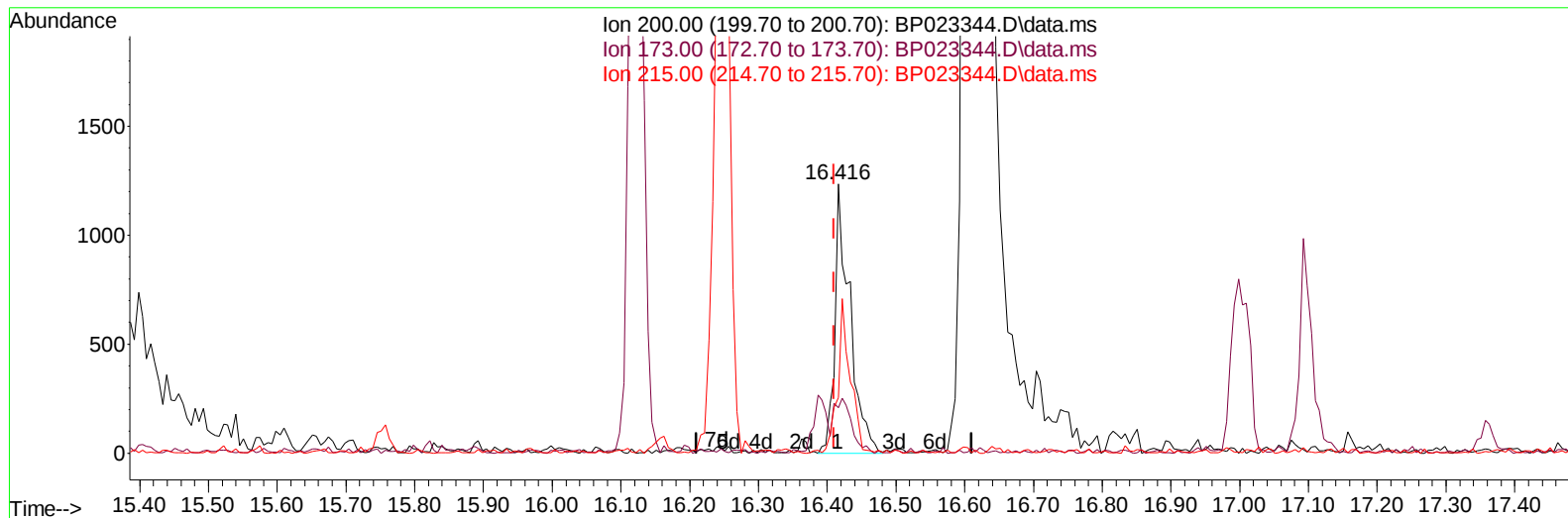
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Data File : BP023344.D
Acq On : 10 Dec 2024 13:11
Operator : RC/JU
Sample : P5107-21MS
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BHJ99MS

Manual IntegrationsAPPROVED

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Response via : Initial Calibration

Reviewed By :Yogesh Patel 12/12/2024
Supervised By :mohammad ahmed 12/13/2024



TIC: BP023344.D\data.ms

(70) Atrazine

16.416min (+ 0.006) 0.27 ng/ul m

response 1879

Ion	Exp%	Act%
200.00	100.00	100.00
173.00	23.60	17.02#
215.00	47.50	20.58#
0.00	0.00	0.00

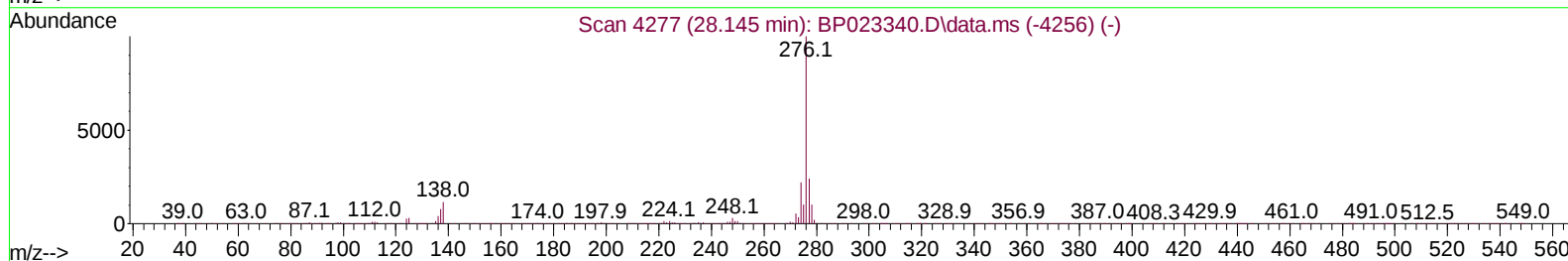
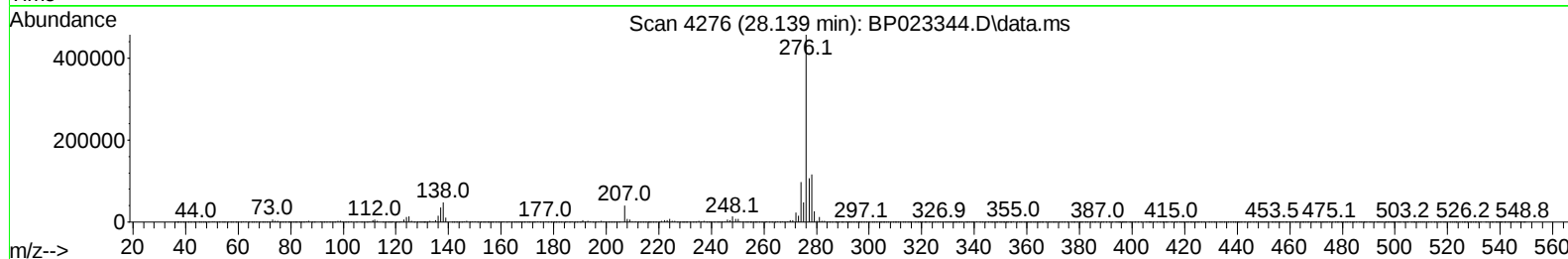
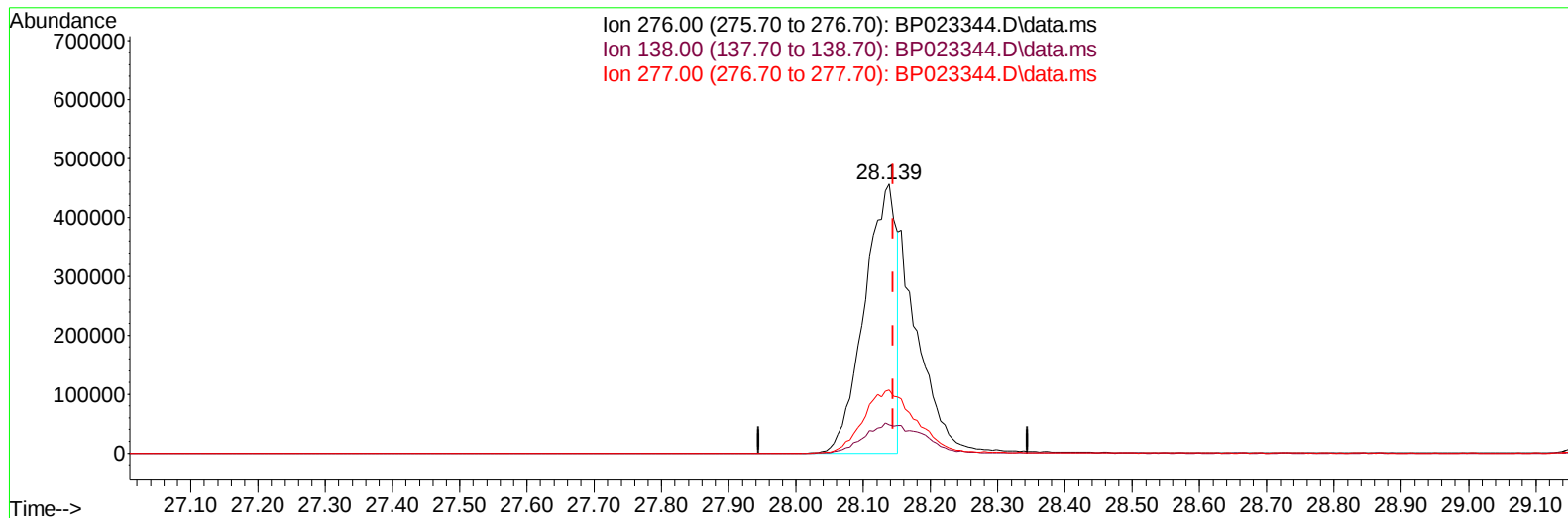
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Data File : BP023344.D
Acq On : 10 Dec 2024 13:11
Operator : RC/JU
Sample : P5107-21MS
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BHJ99MS

Manual IntegrationsAPPROVED

Quant Time: Dec 11 00:59:40 2024
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Quant Title : SVOA CALIBRATION
QLast Update : Wed Dec 11 00:30:30 2024
Response via : Initial Calibration

Reviewed By :Yogesh Patel 12/12/2024
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TIC: BP023344.D\data.ms

(94) Indeno(1,2,3-cd)pyrene

28.139min (-0.006) 23.53 ng/u1

response 1499476

Ion	Exp%	Act%
276.00	100.00	100.00
138.00	9.70	10.49
277.00	24.80	23.50
0.00	0.00	0.00

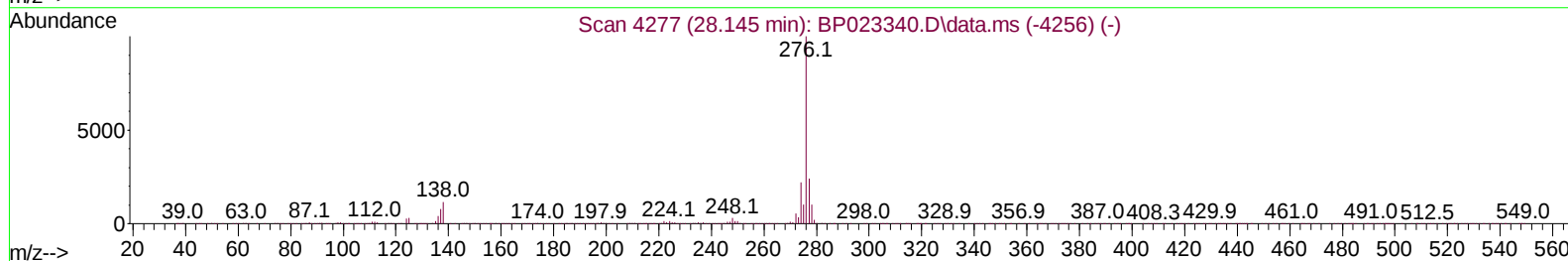
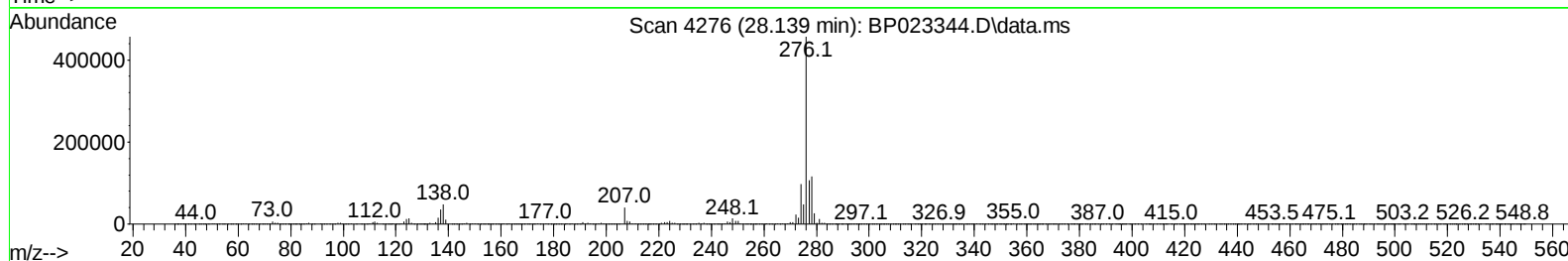
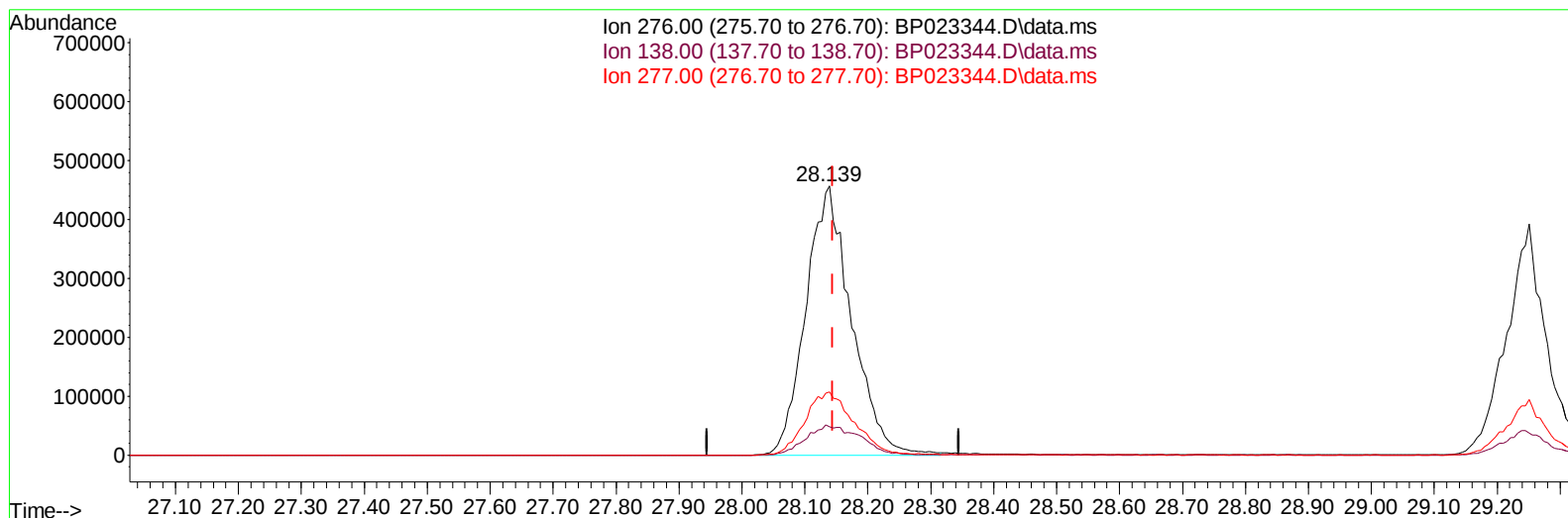
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121024\
Data File : BP023344.D
Acq On : 10 Dec 2024 13:11
Operator : RC/JU
Sample : P5107-21MS
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BHJ99MS

Manual IntegrationsAPPROVED

Quant Time: Dec 11 00:59:40 2024
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Quant Title : SVOA CALIBRATION
QLast Update : Wed Dec 11 00:30:30 2024
Response via : Initial Calibration

Reviewed By :Yogesh Patel 12/12/2024
Supervised By :mohammad ahmed 12/13/2024



TIC: BP023344.D\data.ms

(94) Indeno(1,2,3-cd)pyrene

28.139min (-0.006) 36.01 ng/ul m

response 2294477

Ion	Exp%	Act%
276.00	100.00	100.00
138.00	9.70	10.49
277.00	24.80	23.50
0.00	0.00	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121024\
 Data File : BP023344.D
 Acq On : 10 Dec 2024 13:11
 Operator : RC/JU
 Sample : P5107-21MS
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BHJ99MS

Manual IntegrationsAPPROVED

Quant Time: Dec 11 01:27:50 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112624.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Dec 11 00:30:30 2024
 Response via : Initial Calibration

Reviewed By :Yogesh Patel 12/12/2024
 Supervised By :mohammad ahmed 12/13/2024

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.546	152	65355	20.000	ng/ul	0.00
20) Naphthalene-d8	10.281	136	263424	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.157	164	225289	20.000	ng/ul	0.01
64) Phenanthrene-d10	16.951	188	634014	20.000	ng/ul	0.00
79) Chrysene-d12	21.380	240	780459	20.000	ng/ul	0.00
88) Perylene-d12	24.551	264	870349	20.000	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.105	96	7946	4.476	ng/uL	0.00
4) Pyridine-d5	3.534	84	15334	3.359	ng/ul	0.02
7) Phenol-d5	6.752	99	61857	12.063	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.893	67	100770	31.147	ng/ul	0.00
11) 2-Chlorophenol-d4	7.087	132	107637	28.440	ng/ul	0.00
15) 4-Methylphenol-d8	8.257	113	98434	23.784	ng/ul	0.00
21) Nitrobenzene-d5	8.687	128	56532	28.739	ng/ul	0.00
24) 2-Nitrophenol-d4	9.393	143	69282	30.560	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	9.934	165	147159	30.755	ng/ul	0.00
31) 4-Chloroaniline-d4	10.440	131	98601	18.395	ng/ul	0.00
46) Dimethylphthalate-d6	13.557	166	633660	36.231	ng/ul	0.00
49) Acenaphthylene-d8	13.845	160	586753	31.807	ng/ul	0.01
54) 4-Nitrophenol-d4	14.416	143	33205	15.493	ng/ul	0.00
60) Fluorene-d10	15.157	176	548763	35.673	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.316	200	144289	37.656	ng/ul	0.00
73) Anthracene-d10	17.063	188	1035751	35.795	ng/ul	0.01
81) Pyrene-d10	19.439	212	1449615	35.551	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.345	264	1708212	36.663	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.140	88	9426m	4.799	ng/uL	
5) Pyridine	3.552	79	16550	3.549	ng/ul	90
6) Benzaldehyde	6.716	77	61369m	20.934	ng/ul	
8) Phenol	6.775	94	67100	12.569	ng/ul	95
10) Bis(2-Chloroethyl)ether	6.981	93	122571	29.086	ng/ul	96
12) 2-Chlorophenol	7.116	128	108704	27.398	ng/ul	95
13) 2-Methylphenol	7.993	108	101344	25.597	ng/ul	99
14) 2,2'-oxybis(1-Chloropr...	8.046	45	159231	31.072	ng/ul	97
16) Acetophenone	8.357	105	217347	31.202	ng/ul#	94
17) N-Nitroso-di-n-propyla...	8.334	70	124313	32.752	ng/ul	94
18) 4-Methylphenol	8.316	108	105192	24.616	ng/ul	96
19) Hexachloroethane	8.587	117	48230	24.966	ng/ul	93
22) Nitrobenzene	8.728	77	187242	31.283	ng/ul	97
23) Isophorone	9.240	82	345302	32.571	ng/ul	99
25) 2-Nitrophenol	9.428	139	73153	30.208	ng/ul	93
26) 2,4-Dimethylphenol	9.493	107	145257	27.088	ng/ul	97
27) Bis(2-Chloroethoxy)met...	9.710	93	179229	29.923	ng/ul	100
29) 2,4-Dichlorophenol	9.957	162	140077	29.539	ng/ul	98
30) Naphthalene	10.334	128	387976	28.134	ng/ul	99
32) 4-Chloroaniline	10.469	127	44601m	8.581	ng/ul	
33) Hexachlorobutadiene	10.604	225	114975	27.332	ng/ul	98
34) Caprolactam	11.293	113	6969m	5.037	ng/ul	
35) 4-Chloro-3-methylphenol	11.593	107	176936	34.362	ng/ul	98

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121024\
 Data File : BP023344.D
 Acq On : 10 Dec 2024 13:11
 Operator : RC/JU
 Sample : P5107-21MS
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BHJ99MS

Manual IntegrationsAPPROVED

Quant Time: Dec 11 01:27:50 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112624.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Dec 11 00:30:30 2024
 Response via : Initial Calibration

Reviewed By :Yogesh Patel 12/12/2024
 Supervised By :mohammad ahmed 12/13/2024

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
36) 2-Methylnaphthalene	11.945	142	309146	30.230	ng/ul	94
37) 1-Methylnaphthalene	12.163	142	313371	30.071	ng/ul	97
39) 1,2,4,5-Tetrachloroben...	12.322	216	233958	28.458	ng/ul	97
40) Hexachlorocyclopentadiene	12.287	237	92345	23.378	ng/ul	98
41) 2,4,6-Trichlorophenol	12.575	196	155970	32.058	ng/ul	100
42) 2,4,5-Trichlorophenol	12.663	196	170166	32.735	ng/ul	99
43) 1,1'-Biphenyl	12.963	154	479699	30.051	ng/ul	100
44) 2-Chloronaphthalene	13.010	162	382434	29.619	ng/ul	99
45) 2-Nitroaniline	13.240	65	152079	40.653	ng/ul	95
47) Dimethylphthalate	13.604	163	601886	34.937	ng/ul	99
48) 2,6-Dinitrotoluene	13.745	165	122491	37.107	ng/ul	96
50) Acenaphthylene	13.875	152	618895	32.944	ng/ul	100
51) 3-Nitroaniline	14.087	138	85709	30.979	ng/ul	90
52) Acenaphthene	14.216	153	454445	33.096	ng/ul	99
53) 2,4-Dinitrophenol	14.310	184	84436	39.665	ng/ul	92
55) 4-Nitrophenol	14.434	109	43056	14.912	ng/ul	97
56) Dibenzofuran	14.563	168	719973	34.526	ng/ul	99
57) 2,4-Dinitrotoluene	14.557	165	208095	39.120	ng/ul	89
58) 2,3,4,6-Tetrachlorophenol	14.798	232	181024	35.078	ng/ul	95
59) Diethylphthalate	14.992	149	668256	38.893	ng/ul	99
61) Fluorene	15.222	166	615346	36.241	ng/ul	99
62) 4-Chlorophenyl-phenyle...	15.222	204	356055	36.155	ng/ul	98
63) 4-Nitroaniline	15.269	138	110124	45.255	ng/ul	87
66) 4,6-Dinitro-2-methylph...	15.328	198	152463	37.508	ng/ul	92
67) N-Nitrosodiphenylamine	15.434	169	547939	33.843	ng/ul	99
68) 4-Bromophenyl-phenylether	16.122	248	255494	34.136	ng/ul	98
69) Hexachlorobenzene	16.245	284	305070	33.271	ng/ul	97
70) Atrazine	16.416	200	1879m	0.267	ng/ul	
71) Pentachlorophenol	16.616	266	174602	32.624	ng/ul	94
72) Phenanthrene	16.998	178	1152900	35.173	ng/ul	99
74) Anthracene	17.098	178	1150173	35.094	ng/ul	100
75) 1,2,3,4-Tetrachloroben...	12.928	216	241553	26.431	ng/uL	99
76) Pentachlorobenzene	14.481	250	308157	30.591	ng/uL	98
77) Carbazole	17.392	167	1054203	39.481	ng/ul	100
78) Di-n-butylphthalate	17.963	149	1314227	40.057	ng/ul	100
80) Fluoranthene	19.086	202	1645087	35.209	ng/ul	99
82) Pyrene	19.469	202	1754415	35.385	ng/ul	99
83) Butylbenzylphthalate	20.427	149	670354	39.037	ng/ul	94
84) 3,3'-Dichlorobenzidine	21.286	252	422293	23.580	ng/ul	98
85) Benzo(a)anthracene	21.357	228	1897050	36.248	ng/ul	99
86) Bis(2-ethylhexyl)phtha...	21.286	149	980584	39.764	ng/ul#	98
87) Chrysene	21.427	228	1792329	35.737	ng/ul	100
89) Di-n-octyl phthalate	22.486	149	1652219	38.869	ng/ul	100
90) Benzo(b)fluoranthene	23.551	252	2007579	36.713	ng/ul	99
91) Benzo(k)fluoranthene	23.621	252	1937326	35.607	ng/ul	99
93) Benzo(a)pyrene	24.410	252	1798845	36.281	ng/ul	99
94) Indeno(1,2,3-cd)pyrene	28.139	276	2294477m	36.011	ng/ul	
95) Dibenzo(a,h)anthracene	28.180	278	1810608	35.509	ng/ul	98
96) Benzo(g,h,i)perylene	29.250	276	1764221	36.164	ng/ul	98

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

Instrument :

BNA_P

ClientSampleId :

BHJ99MS

Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 12/12/2024

Supervised By :mohammad ahmed 12/13/2024

