

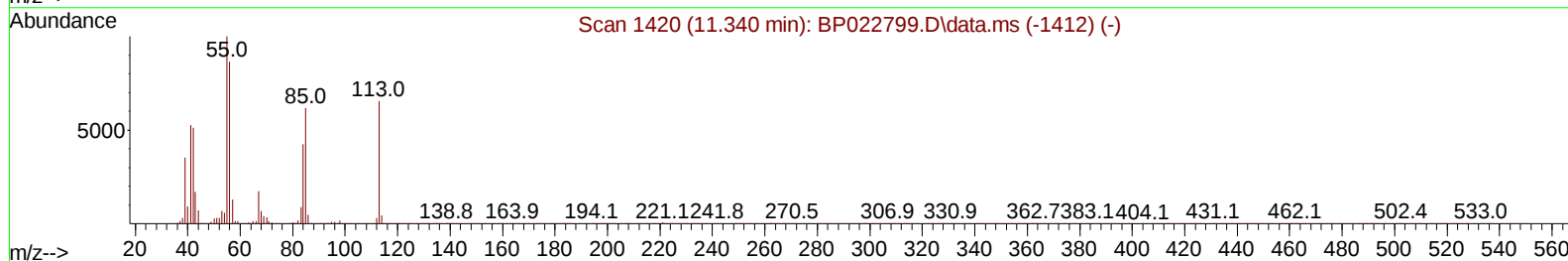
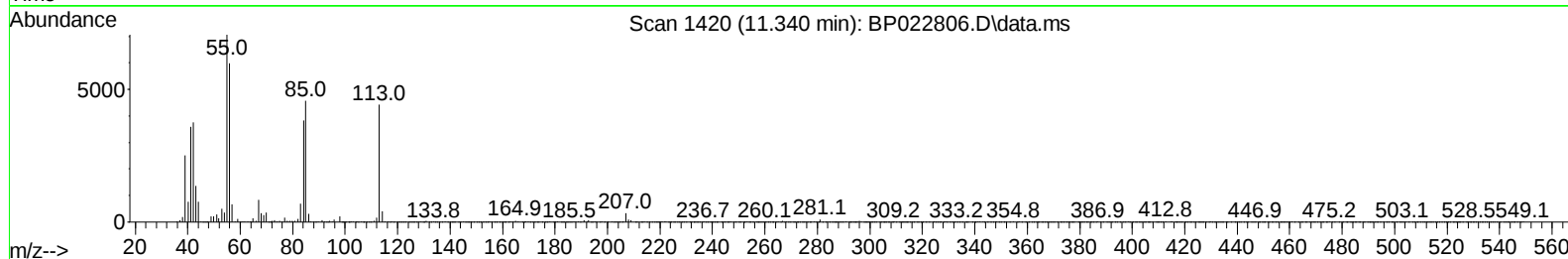
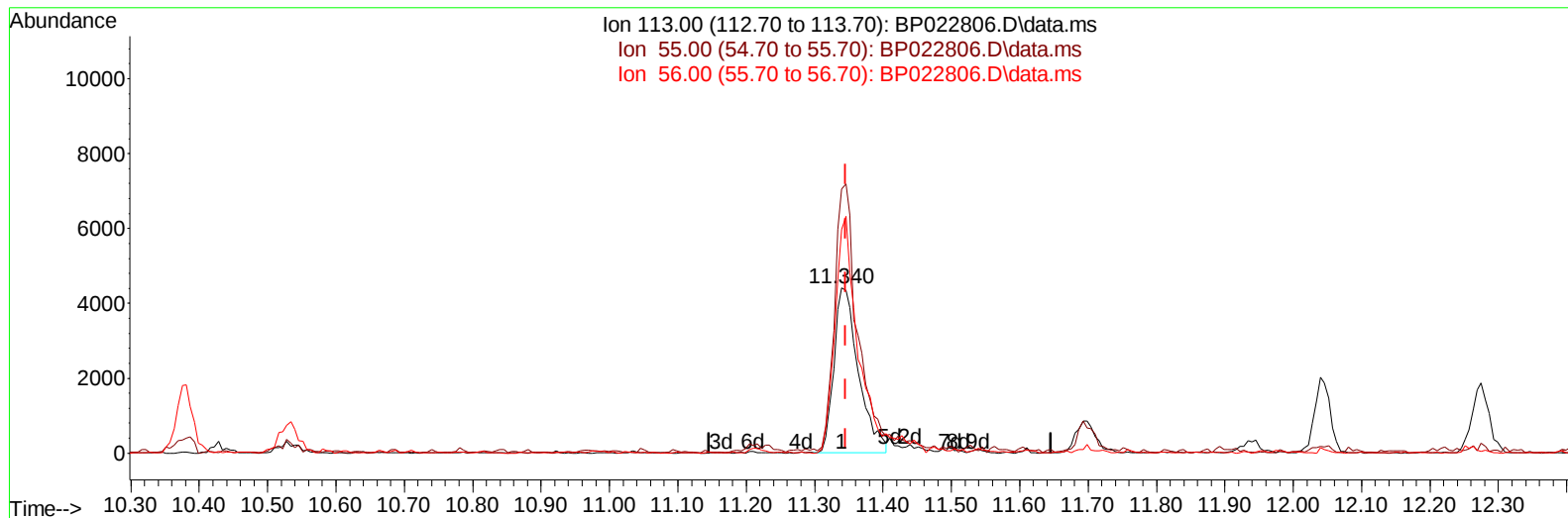
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP103024\
 Data File : BP022806.D
 Acq On : 01 Nov 2024 16:06
 Operator : RC/JU
 Sample : P4630-11MS
 Misc :
 ALS Vial : 42 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 GCP06MS

Manual IntegrationsAPPROVED

Quant Time: Nov 01 18:23:03 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100724.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Oct 30 10:56:39 2024
 Response via : Initial Calibration

Reviewed By :Yogesh Patel 11/04/2024
 Supervised By :mohammad ahmed 11/05/2024



TIC: BP022806.D\data.ms

(34) Caprolactam

11.340min (-0.006) 7.47 ng/ul

response 11021

Ion	Exp%	Act%
113.00	100.00	100.00
55.00	162.70	159.71
56.00	130.00	135.13
0.00	0.00	0.00

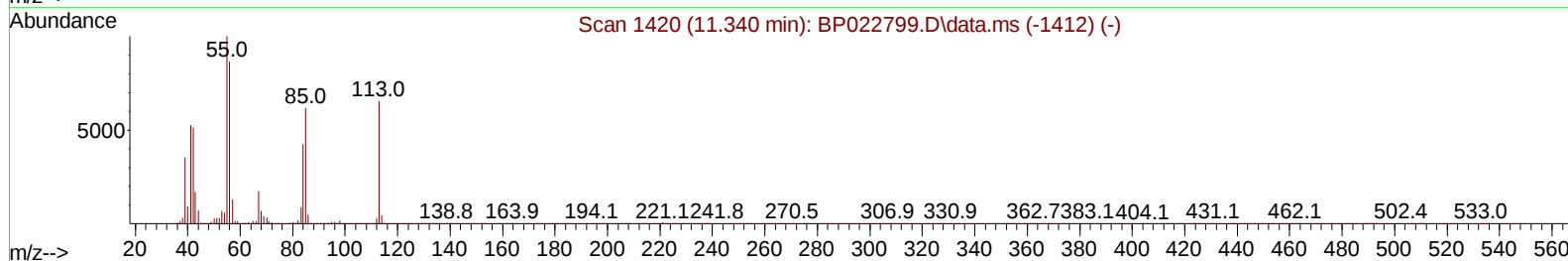
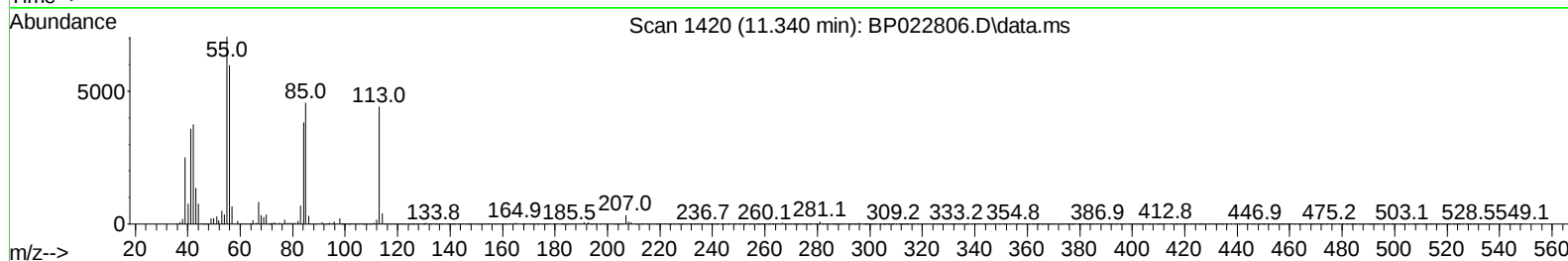
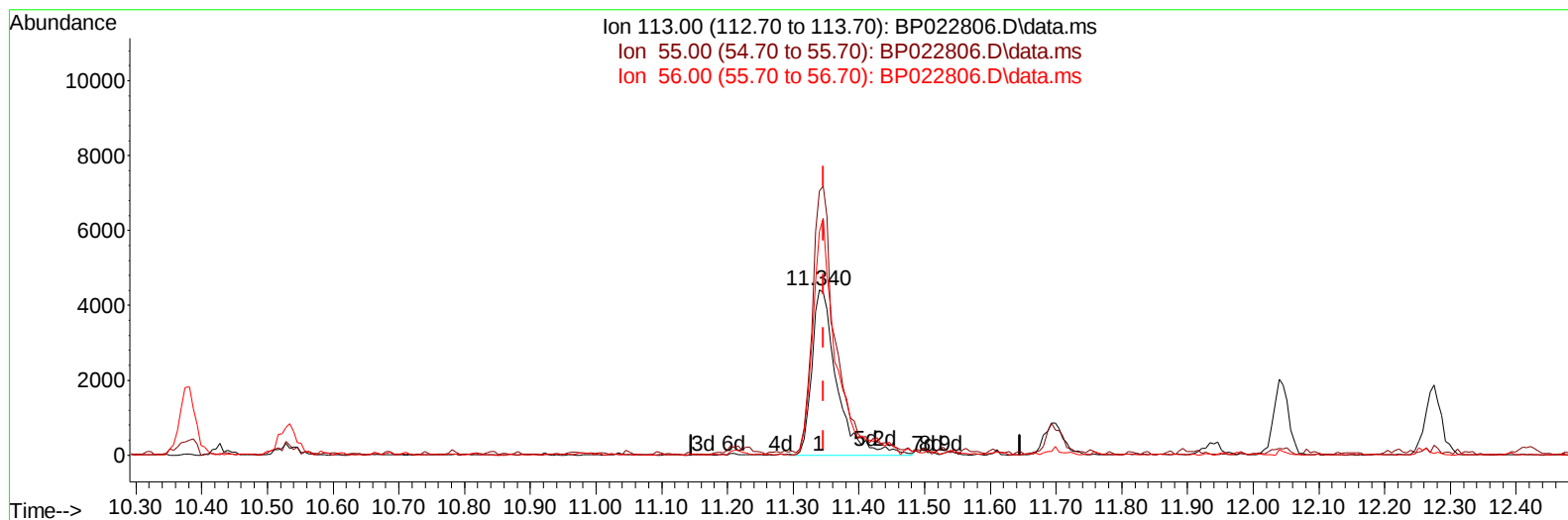
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP103024\
Data File : BP022806.D
Acq On : 01 Nov 2024 16:06
Operator : RC/JU
Sample : P4630-11MS
Misc :
ALS Vial : 42 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
GCP06MS

Manual IntegrationsAPPROVED

Quant Time: Nov 01 18:23:03 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100724.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Oct 30 10:56:39 2024
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Reviewed By :Yogesh Patel 11/04/2024
Supervised By :mohammad ahmed 11/05/2024



TIC: BP022806.D\data.ms

(34) Caprolactam

11.340min (-0.006) 7.97 ng/ul m

response 11754

Ion	Exp%	Act%
113.00	100.00	100.00
55.00	162.70	159.71
56.00	130.00	135.13
0.00	0.00	0.00

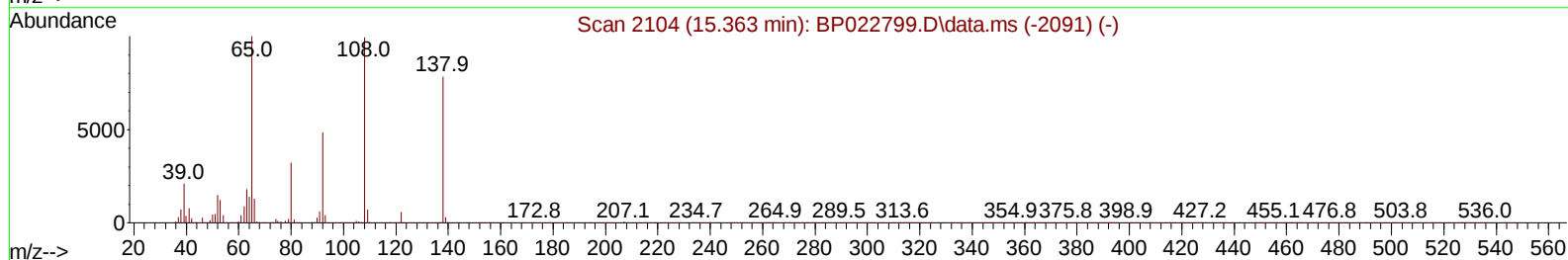
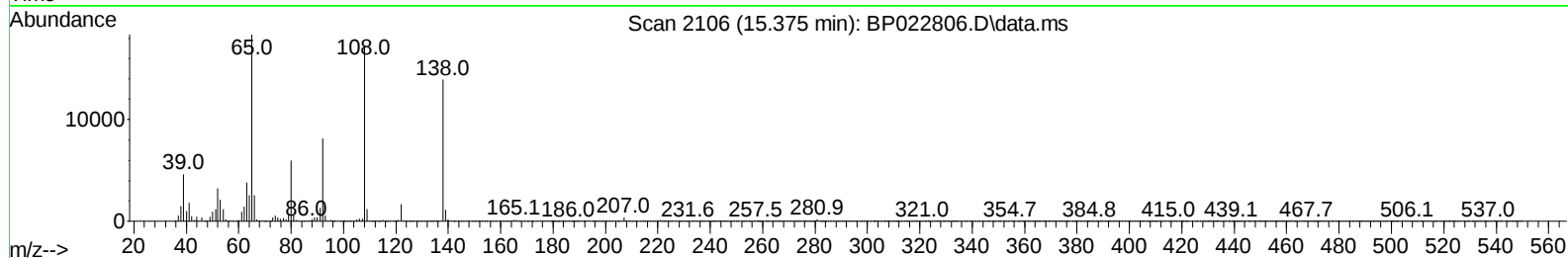
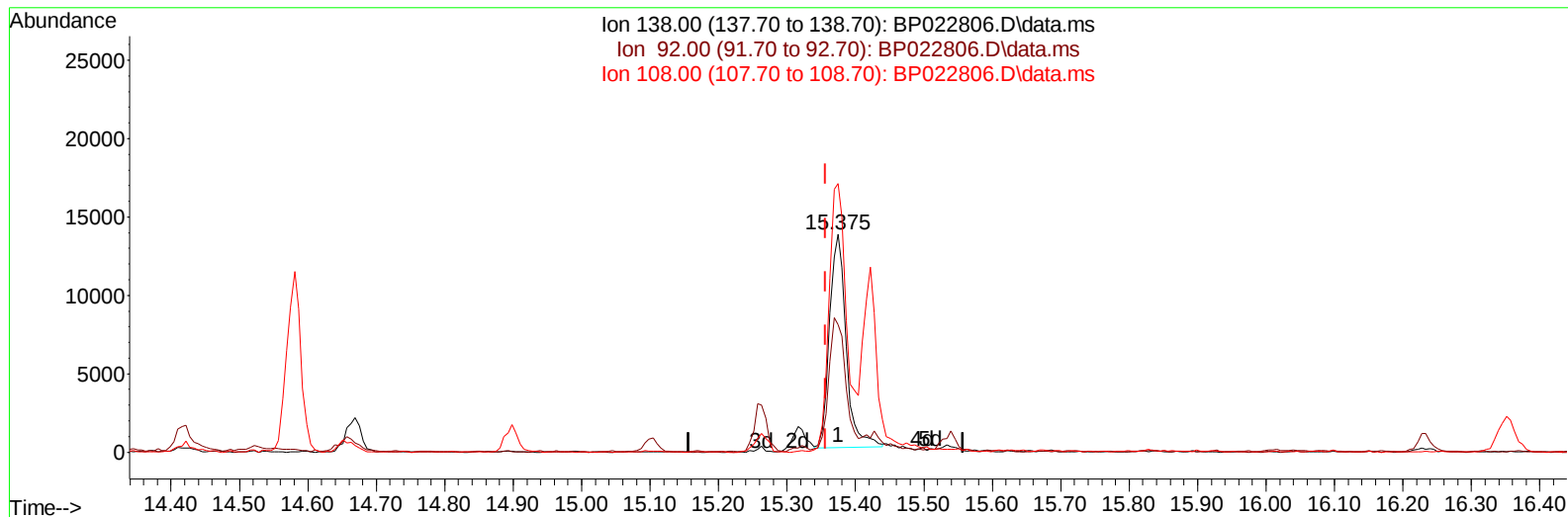
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP103024\
Data File : BP022806.D
Acq On : 01 Nov 2024 16:06
Operator : RC/JU
Sample : P4630-11MS
Misc :
ALS Vial : 42 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
GCP06MS

Manual IntegrationsAPPROVED

Quant Time: Nov 01 18:23:03 2024
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Quant Title : SVOA CALIBRATION
QLast Update : Wed Oct 30 10:56:39 2024
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Reviewed By :Yogesh Patel 11/04/2024
Supervised By :mohammad ahmed 11/05/2024



TIC: BP022806.D\data.ms

(63) 4-Nitroaniline

15.375min (+ 0.018) 8.49 ng/ul

response 22917

Ion	Exp%	Act%
138.00	100.00	100.00
92.00	61.40	58.53
108.00	113.30	122.94
0.00	0.00	0.00

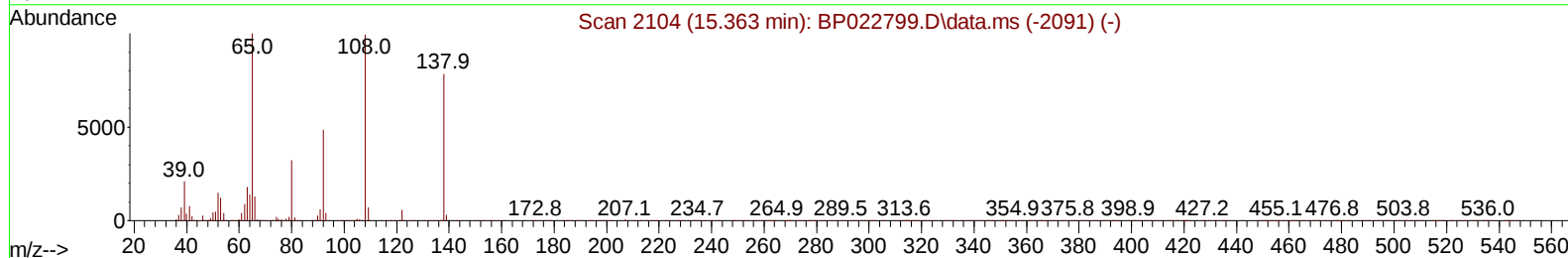
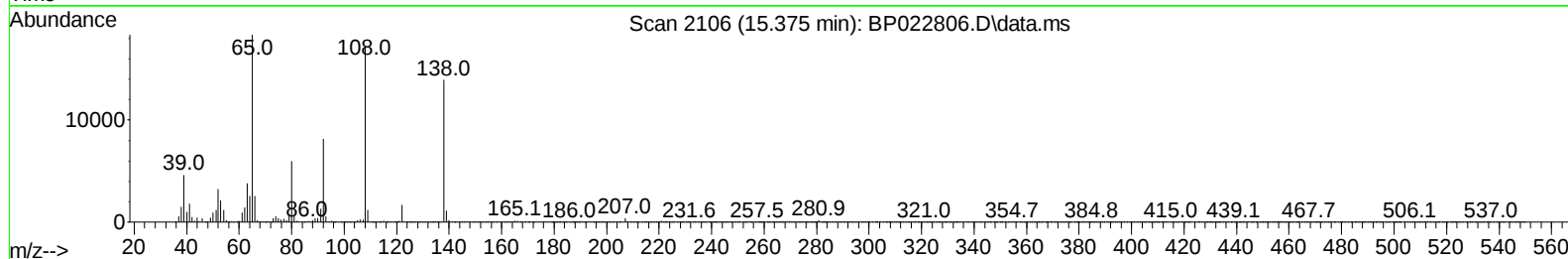
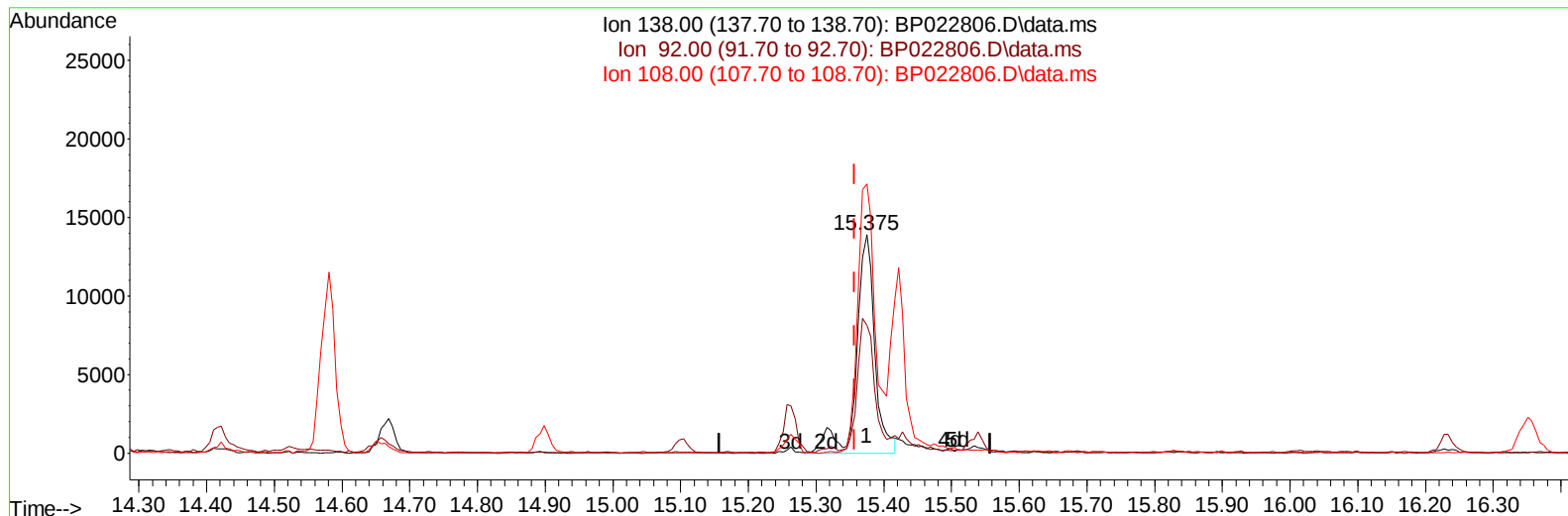
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP103024\
Data File : BP022806.D
Acq On : 01 Nov 2024 16:06
Operator : RC/JU
Sample : P4630-11MS
Misc :
ALS Vial : 42 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
GCP06MS

Manual IntegrationsAPPROVED

Quant Time: Nov 01 18:23:03 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100724.MA.M
Quant Title : SVOA CALIBRATION
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Response via : Initial Calibration

Reviewed By :Yogesh Patel 11/04/2024
Supervised By :mohammad ahmed 11/05/2024



TIC: BP022806.D\data.ms

(63) 4-Nitroaniline

15.375min (+ 0.018) 9.71 ng/ul m

response 26225

Ion	Exp%	Act%
138.00	100.00	100.00
92.00	61.40	58.53
108.00	113.30	122.94
0.00	0.00	0.00

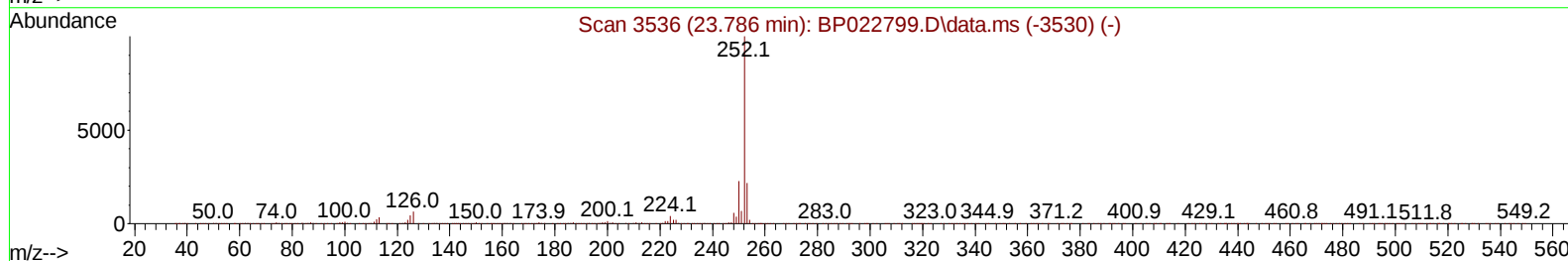
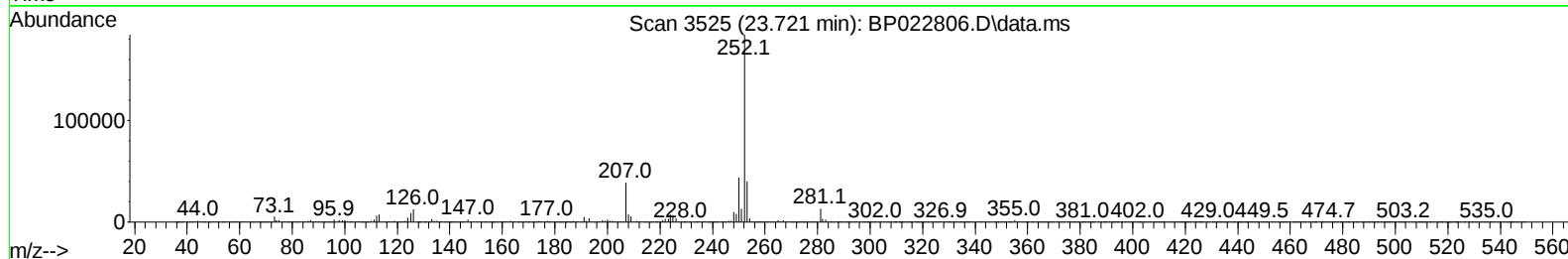
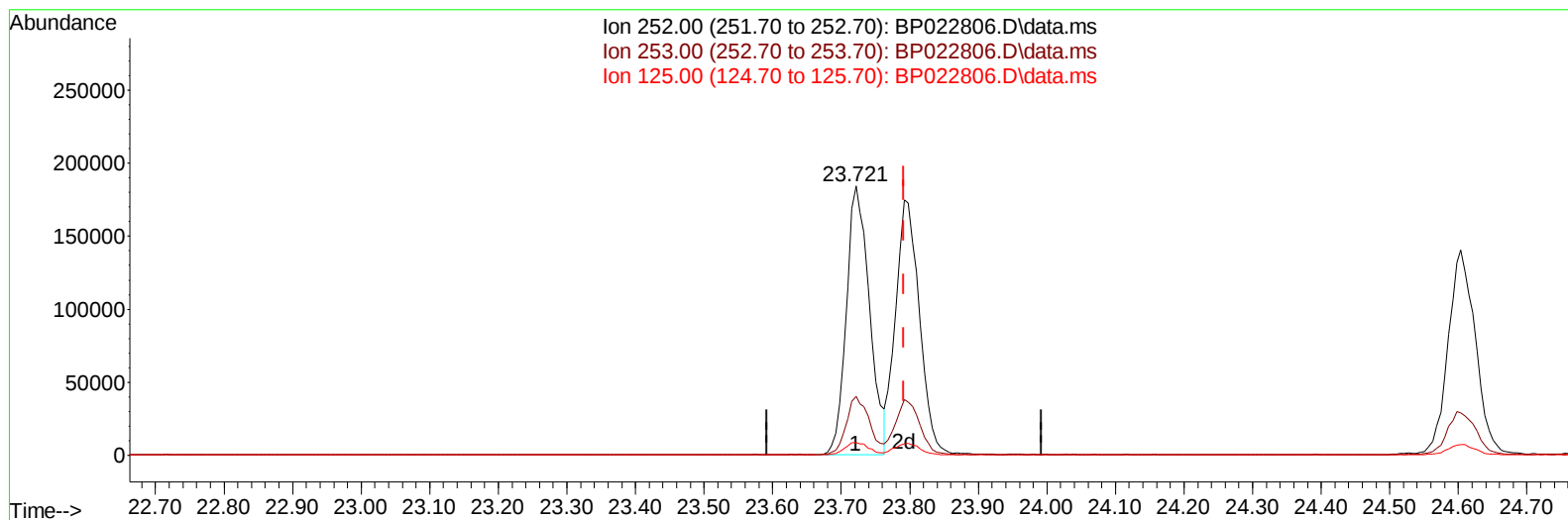
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP103024\
 Data File : BP022806.D
 Acq On : 01 Nov 2024 16:06
 Operator : RC/JU
 Sample : P4630-11MS
 Misc :
 ALS Vial : 42 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
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Manual IntegrationsAPPROVED

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TIC: BP022806.D\data.ms

(91) Benzo(k)fluoranthene

23.721min (-0.071) 10.11 ng/u1

response 431249

Ion	Exp%	Act%
252.00	100.00	100.00
253.00	21.70	21.79
125.00	5.70	4.63
0.00	0.00	0.00

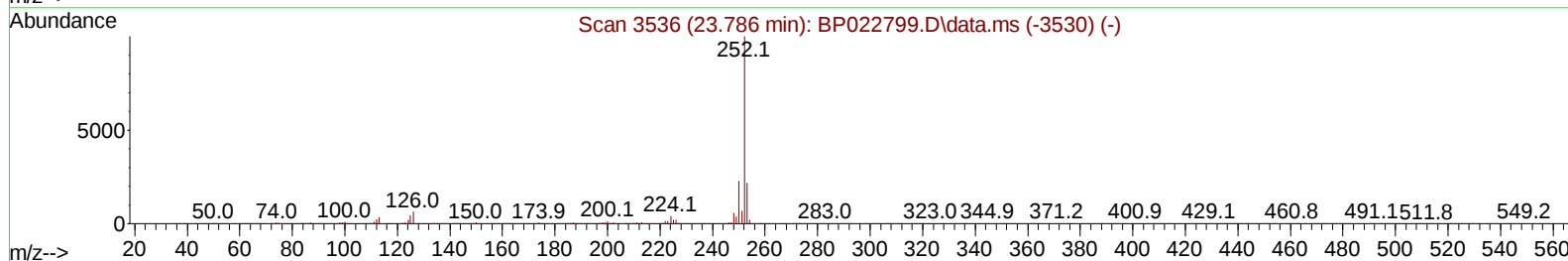
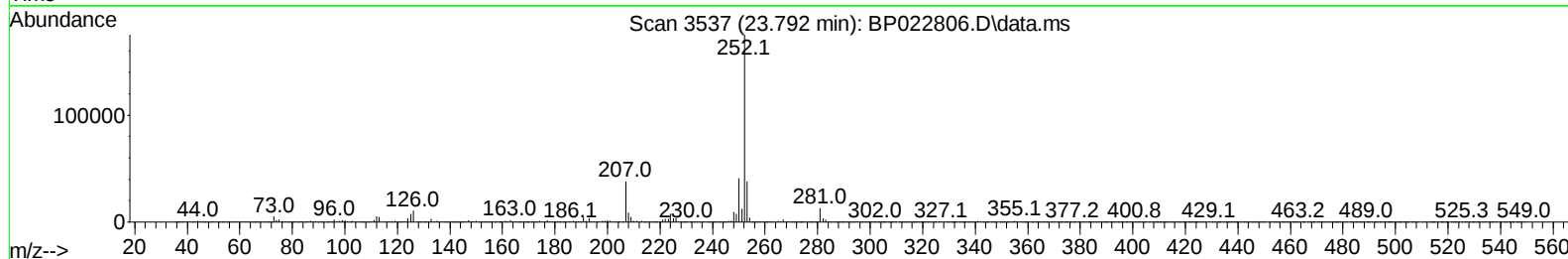
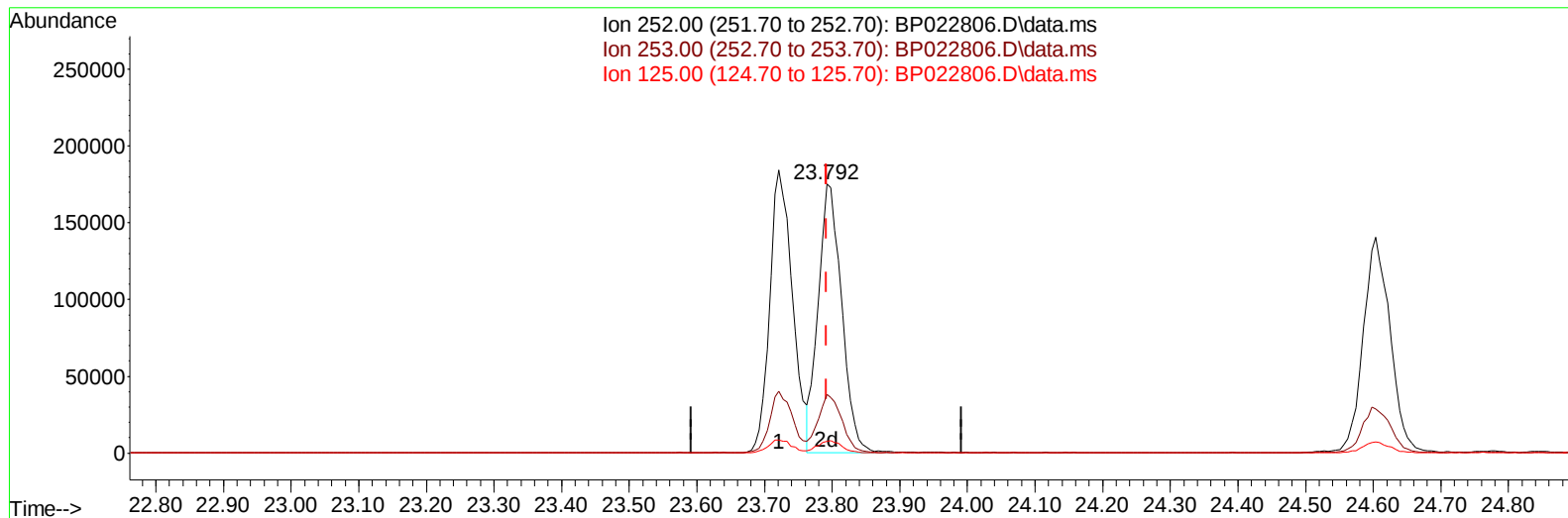
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Data File : BP022806.D
Acq On : 01 Nov 2024 16:06
Operator : RC/JU
Sample : P4630-11MS
Misc :
ALS Vial : 42 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
GCP06MS

Manual IntegrationsAPPROVED

Quant Time: Nov 01 18:23:03 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100724.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Oct 30 10:56:39 2024
Response via : Initial Calibration

Reviewed By :Yogesh Patel 11/04/2024
Supervised By :mohammad ahmed 11/05/2024



TIC: BP022806.D\data.ms

(91) Benzo(k)fluoranthene

23.792min (0.000) 9.90 ng/u1 m

response 422483

Ion	Exp%	Act%
252.00	100.00	100.00
253.00	21.70	21.79
125.00	5.70	4.28#
0.00	0.00	0.00

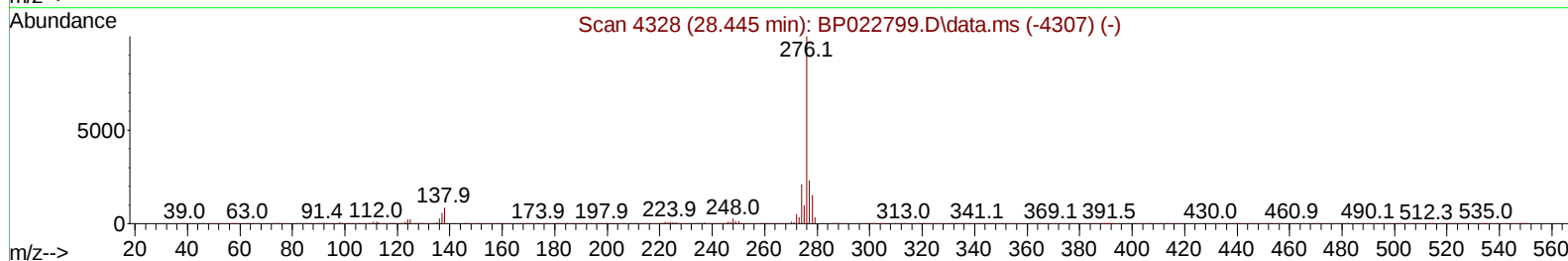
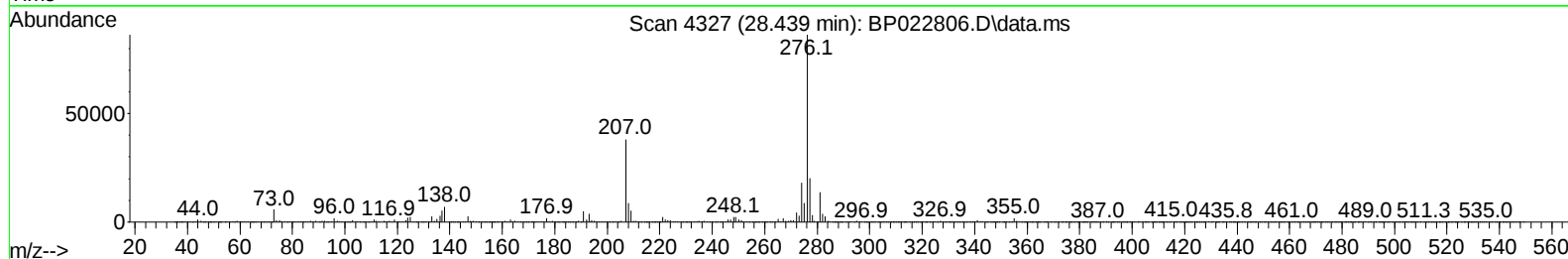
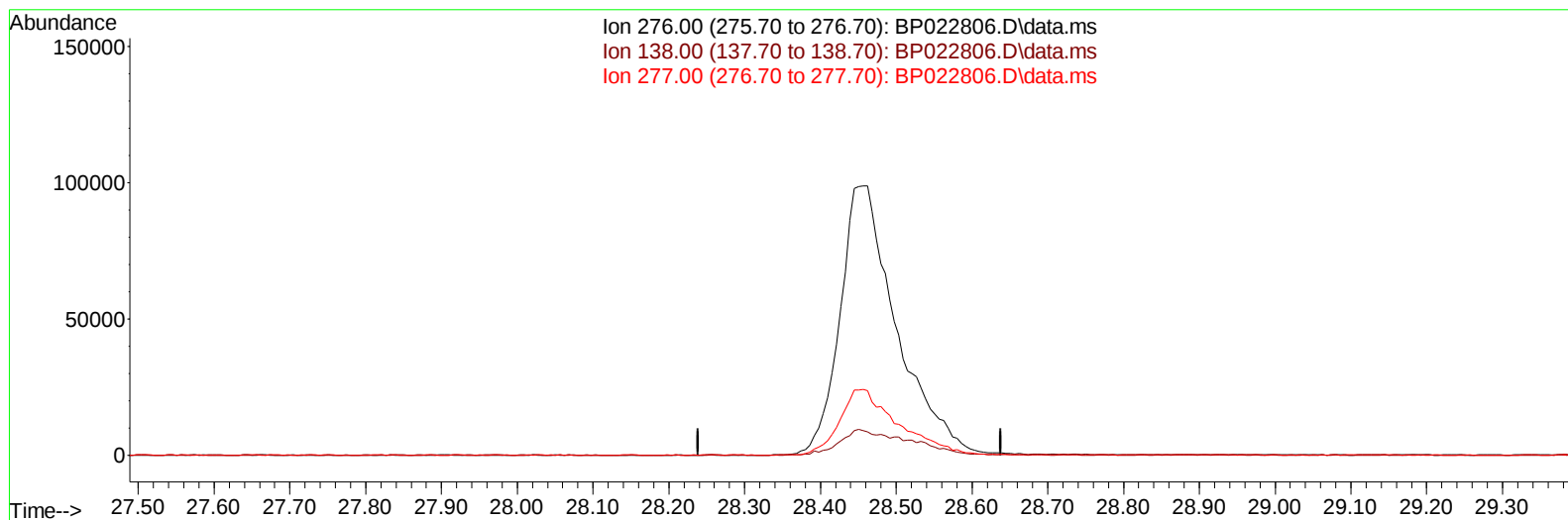
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP103024\
Data File : BP022806.D
Acq On : 01 Nov 2024 16:06
Operator : RC/JU
Sample : P4630-11MS
Misc :
ALS Vial : 42 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
GCP06MS

Manual IntegrationsAPPROVED

Quant Time: Nov 01 18:23:03 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100724.MA.M
Quant Title : SVOA CALIBRATION
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Reviewed By :Yogesh Patel 11/04/2024
Supervised By :mohammad ahmed 11/05/2024



TIC: BP022806.D\data.ms

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

28.439min (-28.439) 0.00 ng/u1

response 0

Ion	Exp%	Act%
276.00	100.00	0.00
138.00	11.40	0.00#
277.00	23.20	0.00#
0.00	0.00	0.00

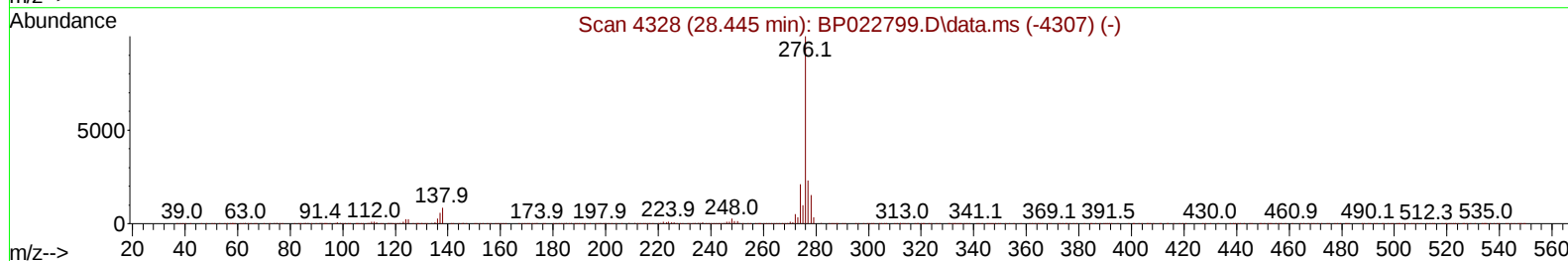
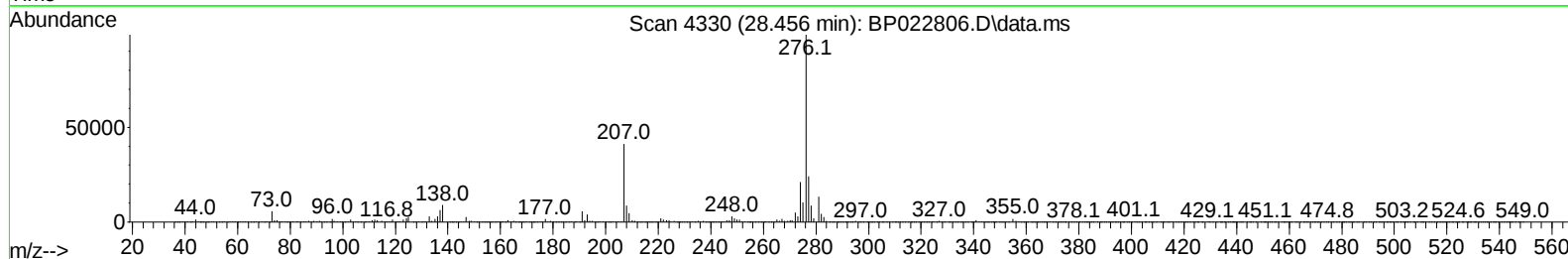
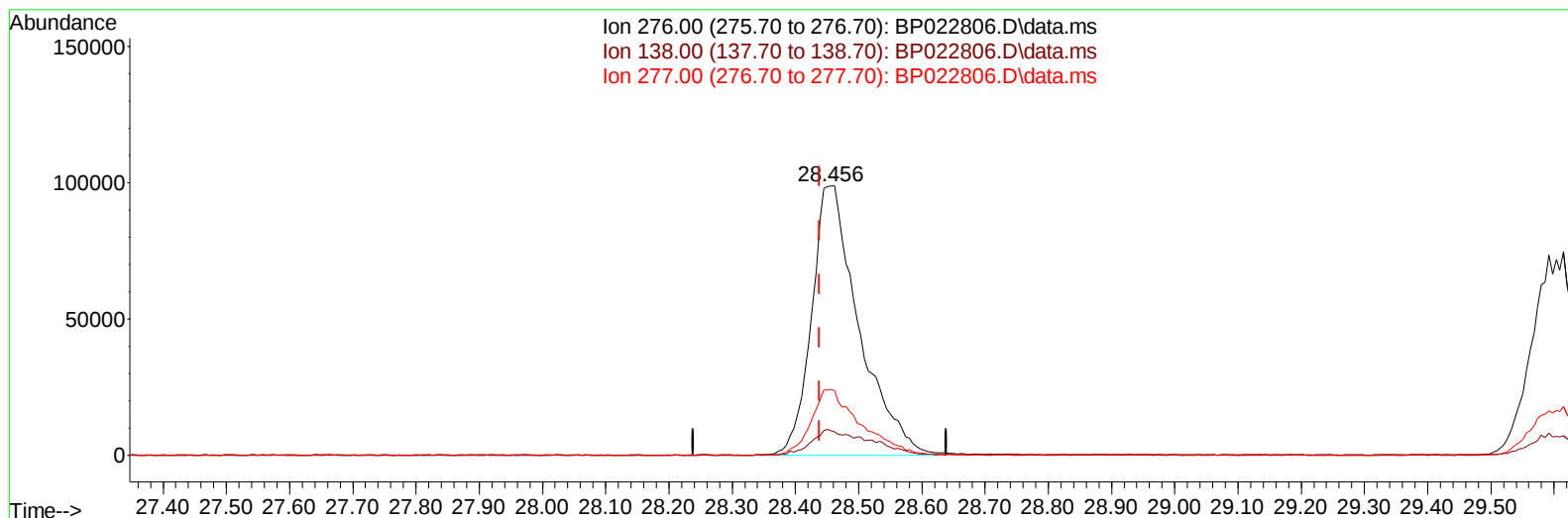
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP103024\
 Data File : BP022806.D
 Acq On : 01 Nov 2024 16:06
 Operator : RC/JU
 Sample : P4630-11MS
 Misc :
 ALS Vial : 42 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 GCP06MS

Manual IntegrationsAPPROVED

Quant Time: Nov 01 18:23:03 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100724.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Oct 30 10:56:39 2024
 Response via : Initial Calibration

Reviewed By :Yogesh Patel 11/04/2024
 Supervised By :mohammad ahmed 11/05/2024



TIC: BP022806.D\data.ms

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

28.456min (+ 0.018) 9.61 ng/ul m

response 514435

Ion	Exp%	Act%
276.00	100.00	100.00
138.00	11.40	9.00#
277.00	23.20	24.42
0.00	0.00	0.00

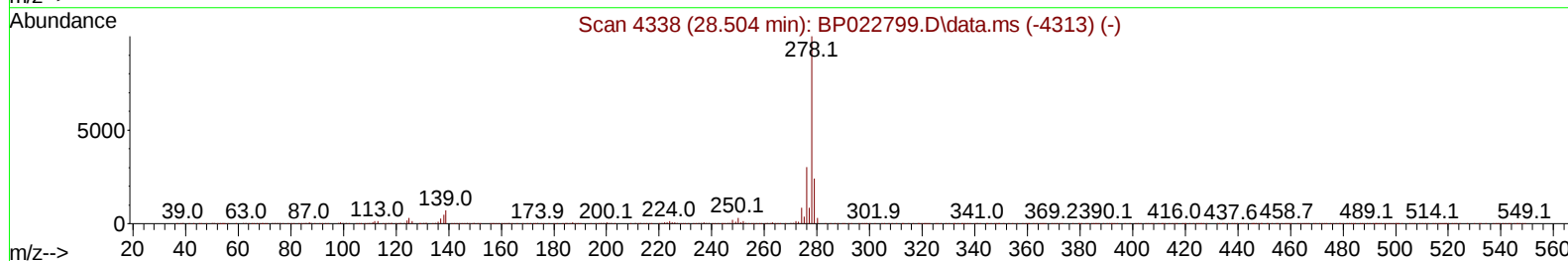
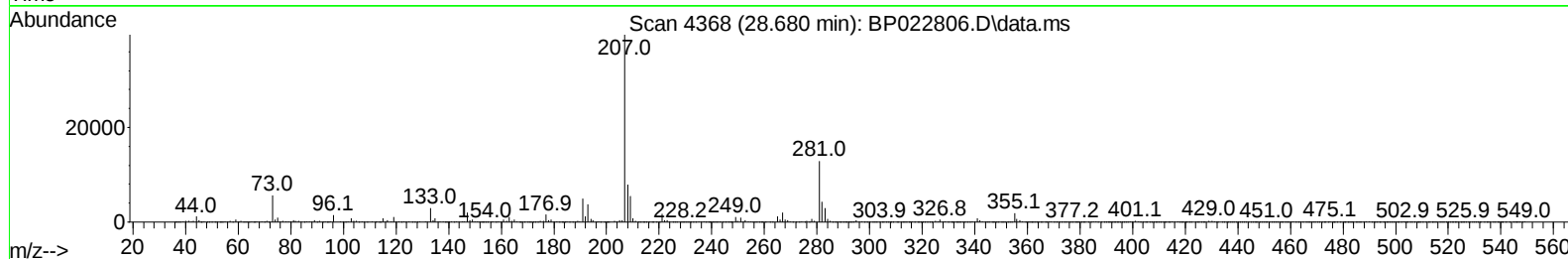
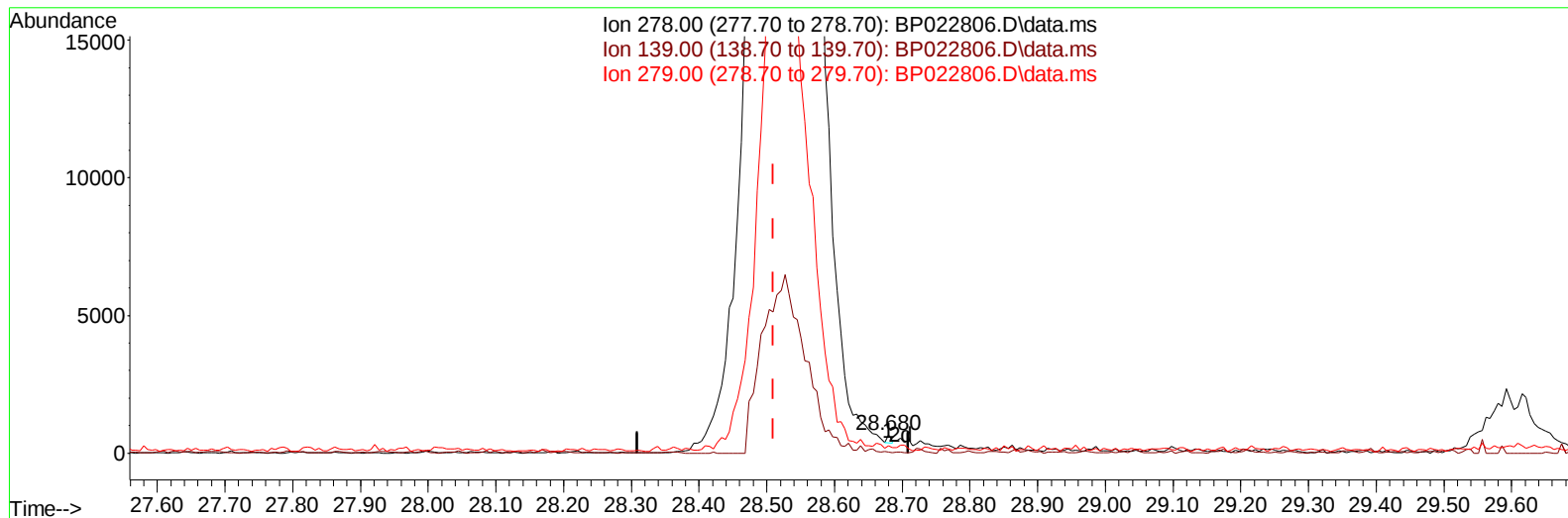
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP103024\
Data File : BP022806.D
Acq On : 01 Nov 2024 16:06
Operator : RC/JU
Sample : P4630-11MS
Misc :
ALS Vial : 42 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
GCP06MS

Manual IntegrationsAPPROVED

Quant Time: Nov 01 18:23:03 2024
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TIC: BP022806.D\data.ms

(95) Dibenzo(a,h)anthracene

28.680min (+ 0.171) 0.00 ng/u1

response 128

Ion	Exp%	Act%
278.00	100.00	100.00
139.00	9.10	5.01#
279.00	23.70	40.97#
0.00	0.00	0.00

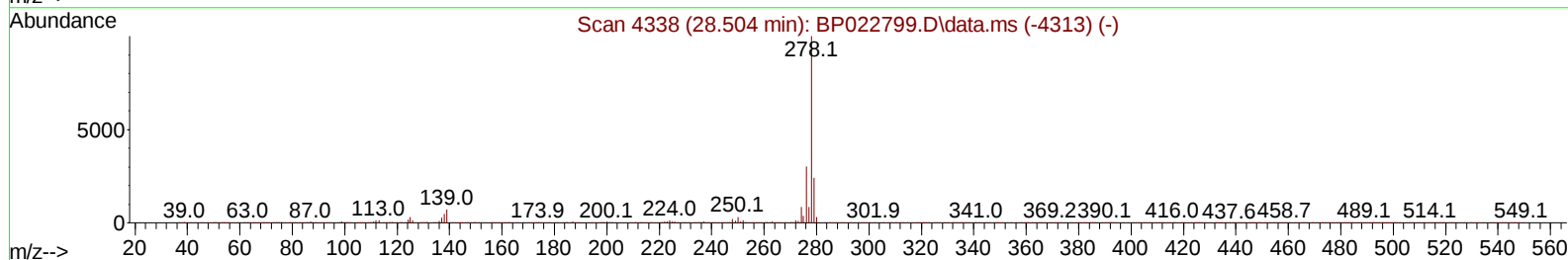
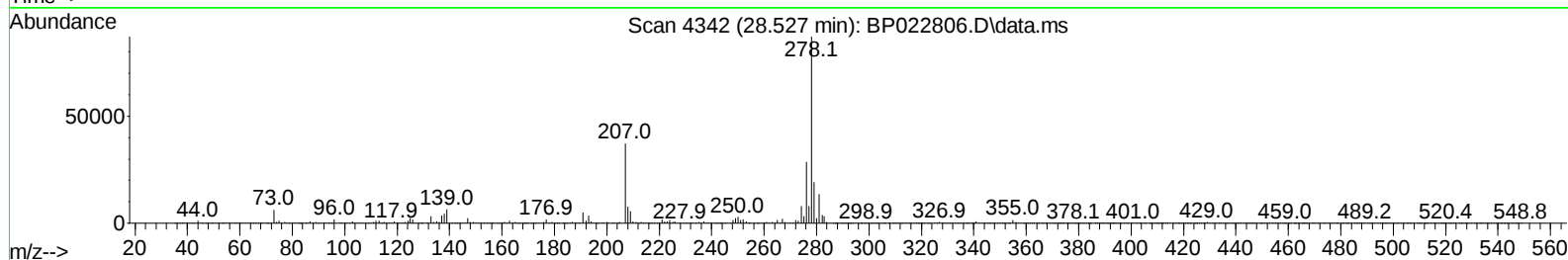
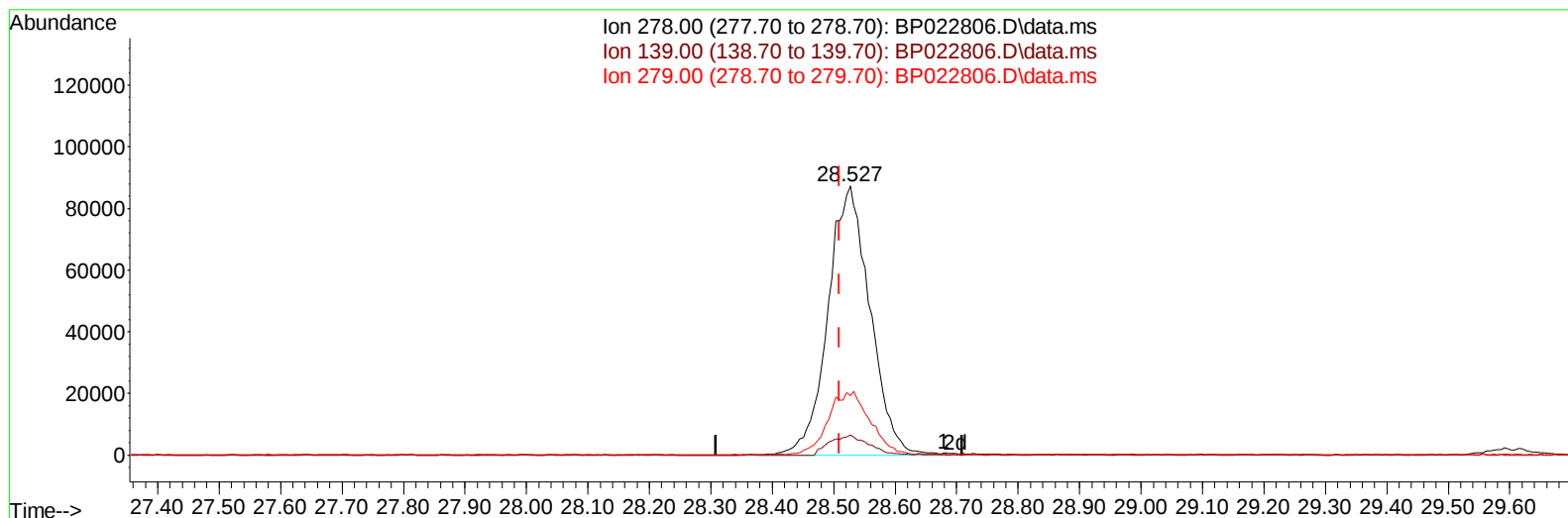
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Data File : BP022806.D
Acq On : 01 Nov 2024 16:06
Operator : RC/JU
Sample : P4630-11MS
Misc :
ALS Vial : 42 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
GCP06MS

Manual IntegrationsAPPROVED

Quant Time: Nov 01 18:23:03 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100724.MA.M
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TIC: BP022806.D\data.ms

(95) Dibenzo(a,h)anthracene

28.527min (+ 0.018) 9.58 ng/ul m

response 415428

Ion	Exp%	Act%
278.00	100.00	100.00
139.00	9.10	7.44
279.00	23.70	22.03
0.00	0.00	0.00

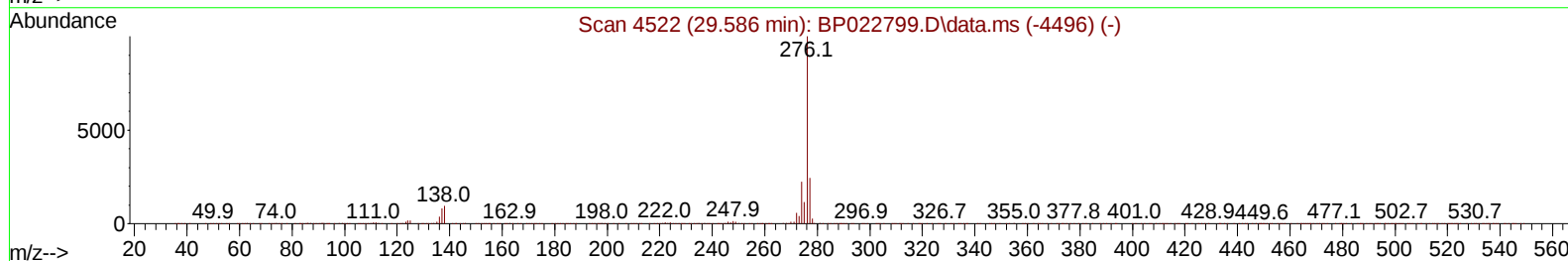
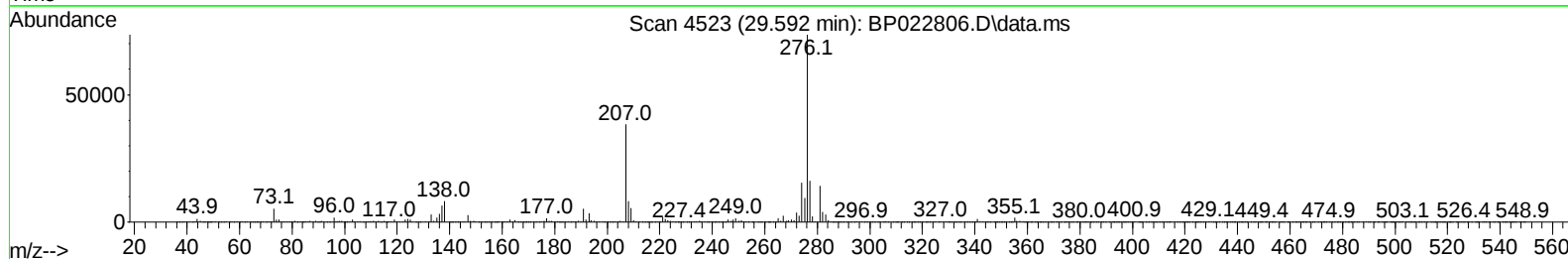
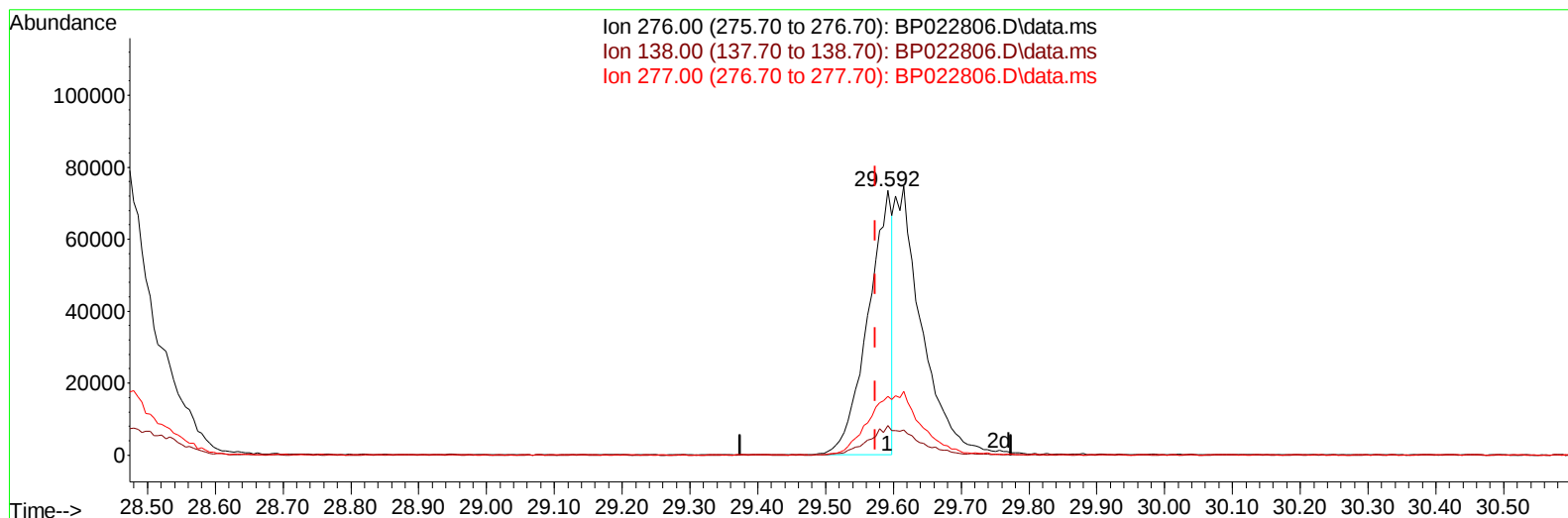
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP103024\
Data File : BP022806.D
Acq On : 01 Nov 2024 16:06
Operator : RC/JU
Sample : P4630-11MS
Misc :
ALS Vial : 42 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
GCP06MS

Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 11/04/2024
Supervised By :mohammad ahmed 11/05/2024

Quant Time: Nov 01 18:23:03 2024
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QLast Update : Wed Oct 30 10:56:39 2024
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TIC: BP022806.D\data.ms

(96) Benzo(g,h,i)perylene

29.592min (+ 0.018) 4.37 ng/u1

response 182039

Ion	Exp%	Act%
276.00	100.00	100.00
138.00	10.60	11.05
277.00	23.70	22.18
0.00	0.00	0.00

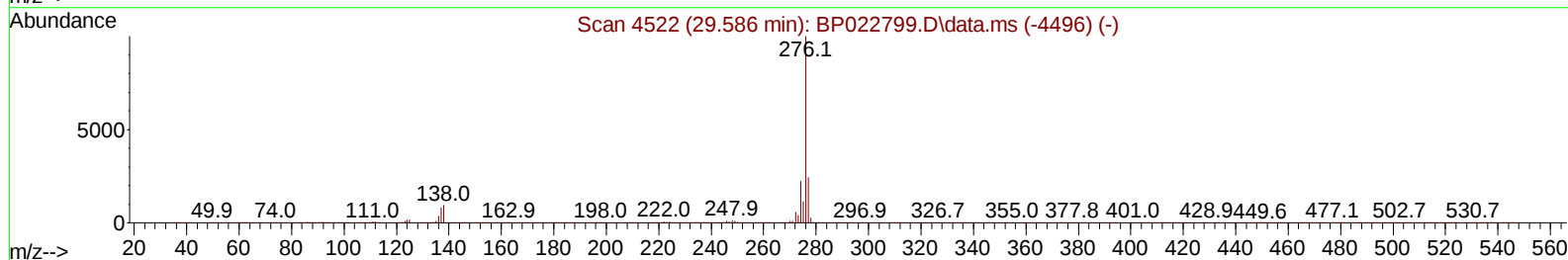
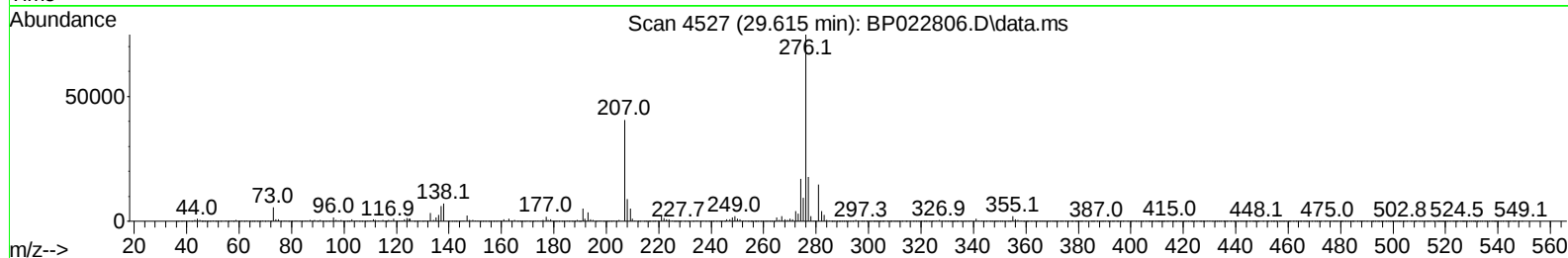
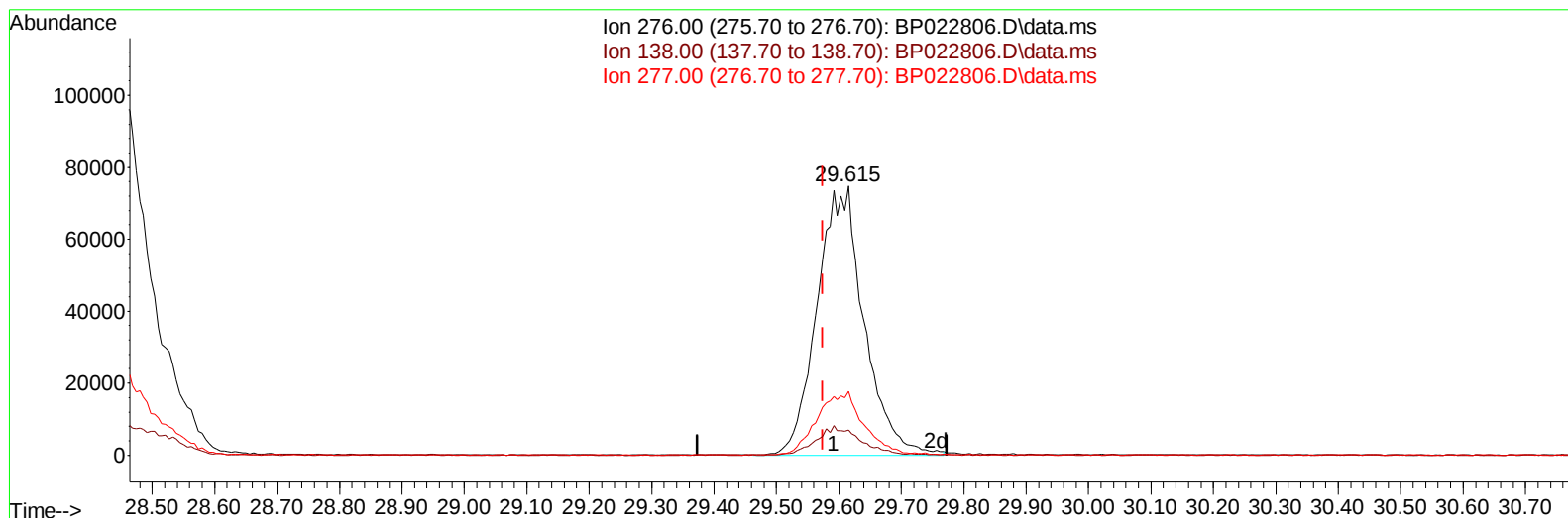
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP103024\
Data File : BP022806.D
Acq On : 01 Nov 2024 16:06
Operator : RC/JU
Sample : P4630-11MS
Misc :
ALS Vial : 42 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
GCP06MS

Manual IntegrationsAPPROVED

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Reviewed By :Yogesh Patel 11/04/2024
Supervised By :mohammad ahmed 11/05/2024



TIC: BP022806.D\data.ms

(96) Benzo(g,h,i)perylene

29.615min (+ 0.041) 9.37 ng/ul m

response 390321

Ion	Exp%	Act%
276.00	100.00	100.00
138.00	10.60	9.39
277.00	23.70	23.71
0.00	0.00	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP103024\
 Data File : BP022806.D
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 Operator : RC/JU
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 Misc :
 ALS Vial : 42 Sample Multiplier: 1

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.628	152	65382	20.000	ng/ul	-0.01
20) Naphthalene-d8	10.381	136	245499	20.000	ng/ul	-0.01
38) Acenaphthene-d10	14.251	164	190839	20.000	ng/ul	-0.01
64) Phenanthrene-d10	17.069	188	479876	20.000	ng/ul	0.00
79) Chrysene-d12	21.510	240	604460	20.000	ng/ul	0.01
88) Perylene-d12	24.774	264	692449	20.000	ng/ul	0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.175	96	5821	3.460	ng/uL	0.00
4) Pyridine-d5	3.593	84	34377	7.443	ng/ul	0.00
7) Phenol-d5	6.822	99	36102	6.560	ng/ul	-0.01
9) Bis-(2-Chloroethyl)eth...	6.975	67	79112	24.457	ng/ul	0.00
11) 2-Chlorophenol-d4	7.169	132	93258	23.880	ng/ul	-0.01
15) 4-Methylphenol-d8	8.334	113	72822	16.580	ng/ul	-0.01
21) Nitrobenzene-d5	8.775	128	51848	27.467	ng/ul	0.00
24) 2-Nitrophenol-d4	9.487	143	62583	28.637	ng/ul	-0.01
28) 2,4-Dichlorophenol-d3	10.028	165	128259	27.613	ng/ul	0.00
31) 4-Chloroaniline-d4	10.528	131	109965	20.036	ng/ul	-0.01
46) Dimethylphthalate-d6	13.669	166	440901	29.078	ng/ul	0.00
49) Acenaphthylene-d8	13.940	160	455249	28.269	ng/ul	-0.01
54) 4-Nitrophenol-d4	14.510	143	12114	4.762	ng/ul	0.00
60) Fluorene-d10	15.263	176	388443	29.536	ng/ul	-0.01
65) 4,6-Dinitro-2-methylph...	15.422	200	78793	24.966	ng/ul	0.02
73) Anthracene-d10	17.175	188	716453	31.737	ng/ul	0.00
81) Pyrene-d10	19.545	212	956501	29.923	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.533	264	1203591	32.547	ng/ul	0.01
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.211	88	6591	3.643	ng/ul#	89
5) Pyridine	3.611	79	42068	8.883	ng/ul	97
6) Benzaldehyde	6.799	77	4493	1.511	ng/ul	86
8) Phenol	6.852	94	53749	9.387	ng/ul	94
10) Bis(2-Chloroethyl)ether	7.063	93	39376	9.188	ng/ul	96
12) 2-Chlorophenol	7.205	128	40993	10.151	ng/ul	95
13) 2-Methylphenol	8.081	108	38221	9.198	ng/ul	97
14) 2,2'-oxybis(1-Chloropr...	8.140	45	44667	8.968	ng/ul#	94
16) Acetophenone	8.440	105	65633	9.051	ng/ul	97
17) N-Nitroso-di-n-propyla...	8.428	70	34239	9.320	ng/ul	95
18) 4-Methylphenol	8.405	108	43815	9.504	ng/ul	100
19) Hexachloroethane	8.681	117	18319	9.740	ng/ul	99
22) Nitrobenzene	8.816	77	53541	9.502	ng/ul	98
23) Isophorone	9.328	82	89685	8.878	ng/ul	97
25) 2-Nitrophenol	9.516	139	23580	10.012	ng/ul	91
26) 2,4-Dimethylphenol	9.581	107	52703	10.295	ng/ul	96
27) Bis(2-Chloroethoxy)met...	9.804	93	51468	8.972	ng/ul	99
29) 2,4-Dichlorophenol	10.052	162	47971	10.488	ng/ul	95
30) Naphthalene	10.428	128	129566	9.909	ng/ul	98
32) 4-Chloroaniline	10.551	127	28350	5.233	ng/ul	97
33) Hexachlorobutadiene	10.710	225	40785	10.579	ng/ul	97
34) Caprolactam	11.340	113	11754m	7.971	ng/ul	
35) 4-Chloro-3-methylphenol	11.693	107	47161	9.450	ng/ul	97

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP103024\
 Data File : BP022806.D
 Acq On : 01 Nov 2024 16:06
 Operator : RC/JU
 Sample : P4630-11MS
 Misc :
 ALS Vial : 42 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 GCP06MS

Manual IntegrationsAPPROVED

Quant Time: Nov 01 18:27:20 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100724.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Oct 30 10:56:39 2024
 Response via : Initial Calibration

Reviewed By :Yogesh Patel 11/04/2024
 Supervised By :mohammad ahmed 11/05/2024

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
36) 2-Methylnaphthalene	12.040	142	95982	10.000	ng/ul	100
37) 1-Methylnaphthalene	12.275	142	94737	9.670	ng/ul	100
39) 1,2,4,5-Tetrachloroben...	12.416	216	73868	10.468	ng/ul	93
40) Hexachlorocyclopentadiene	12.387	237	18383	4.276	ng/ul	98
41) 2,4,6-Trichlorophenol	12.675	196	44143	9.945	ng/ul	100
42) 2,4,5-Trichlorophenol	12.763	196	47422	10.034	ng/ul	98
43) 1,1'-Biphenyl	13.069	154	140917	10.244	ng/ul	100
44) 2-Chloronaphthalene	13.110	162	110427	9.807	ng/ul	99
45) 2-Nitroaniline	13.340	65	29419	8.750	ng/ul	96
47) Dimethylphthalate	13.716	163	142346	9.421	ng/ul	98
48) 2,6-Dinitrotoluene	13.834	165	26909	9.416	ng/ul	98
50) Acenaphthylene	13.969	152	164468	9.840	ng/ul	98
51) 3-Nitroaniline	14.175	138	23302	8.681	ng/ul#	94
52) Acenaphthene	14.322	153	116460	9.854	ng/ul	98
53) 2,4-Dinitrophenol	14.422	184	12699	6.040	ng/ul	98
55) 4-Nitrophenol	14.528	109	20212	7.159	ng/ul	96
56) Dibenzofuran	14.669	168	181021	10.196	ng/ul	99
57) 2,4-Dinitrotoluene	14.657	165	42275	9.520	ng/ul#	76
58) 2,3,4,6-Tetrachlorophenol	14.898	232	43808	9.931	ng/ul#	99
59) Diethylphthalate	15.104	149	139722	9.639	ng/ul	99
61) Fluorene	15.316	166	145621	10.067	ng/ul	96
62) 4-Chlorophenyl-phenyle...	15.322	204	83923	10.145	ng/ul	93
63) 4-Nitroaniline	15.375	138	26225m	9.710	ng/ul	
66) 4,6-Dinitro-2-methylph...	15.434	198	27975	8.229	ng/ul#	92
67) N-Nitrosodiphenylamine	15.539	169	126719	9.861	ng/ul	99
68) 4-Bromophenyl-phenylether	16.234	248	58807	10.162	ng/ul	97
69) Hexachlorobenzene	16.351	284	77260	10.846	ng/ul	96
70) Atrazine	16.522	200	53285	9.975	ng/ul	96
71) Pentachlorophenol	16.722	266	43207	9.434	ng/ul	95
72) Phenanthrene	17.116	178	255082	9.935	ng/ul	99
74) Anthracene	17.210	178	255160	9.959	ng/ul	99
75) 1,2,3,4-Tetrachloroben...	13.040	216	73545	10.065	ng/uL	97
76) Pentachlorobenzene	14.581	250	78805	9.756	ng/uL	98
77) Carbazole	17.498	167	222641	9.821	ng/ul	98
78) Di-n-butylphthalate	18.069	149	236713	9.437	ng/ul	99
80) Fluoranthene	19.198	202	337775	9.192	ng/ul	99
82) Pyrene	19.580	202	366343	9.301	ng/ul	99
83) Butylbenzylphthalate	20.522	149	120668	8.800	ng/ul	94
84) 3,3'-Dichlorobenzidine	21.410	252	98242	6.818	ng/ul	96
85) Benzo(a)anthracene	21.492	228	403064	9.512	ng/ul	100
86) Bis(2-ethylhexyl)phtha...	21.410	149	173854	8.833	ng/ul	100
87) Chrysene	21.551	228	373686	9.511	ng/ul	100
89) Di-n-octyl phthalate	22.651	149	300367	9.119	ng/ul	100
90) Benzo(b)fluoranthene	23.721	252	431249	9.894	ng/ul#	98
91) Benzo(k)fluoranthene	23.792	252	422483m	9.900	ng/ul	
93) Benzo(a)pyrene	24.604	252	387200	9.652	ng/ul	97
94) Indeno(1,2,3-cd)pyrene	28.456	276	514435m	9.610	ng/ul	
95) Dibenzo(a,h)anthracene	28.527	278	415428m	9.583	ng/ul	
96) Benzo(g,h,i)perylene	29.615	276	390321m	9.374	ng/ul	

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

Instrument :
BNA_P
ClientSampleId :
GCP06MS

Manual IntegrationsAPPROVED

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