

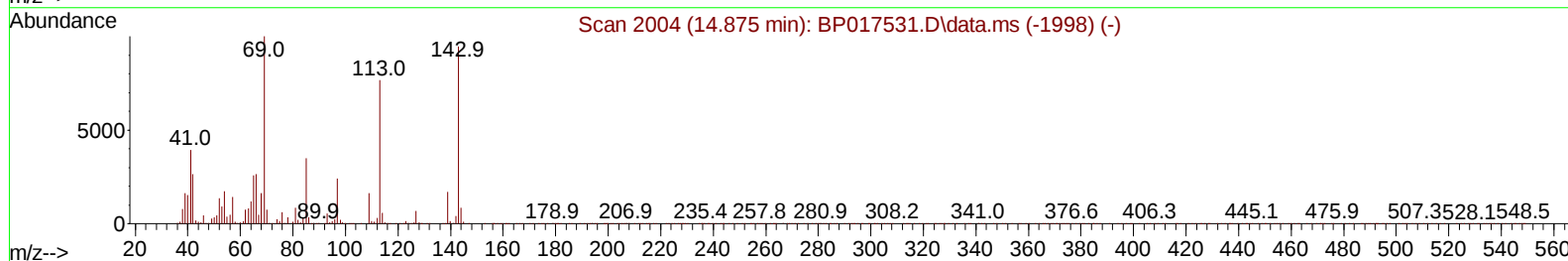
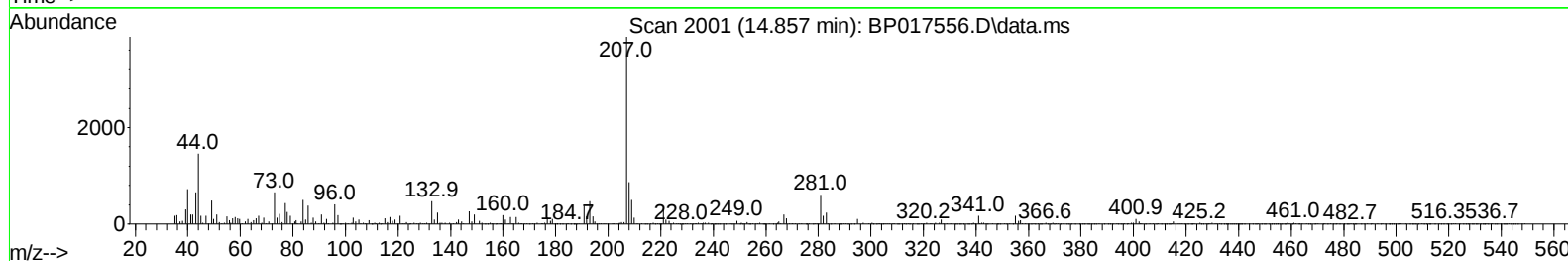
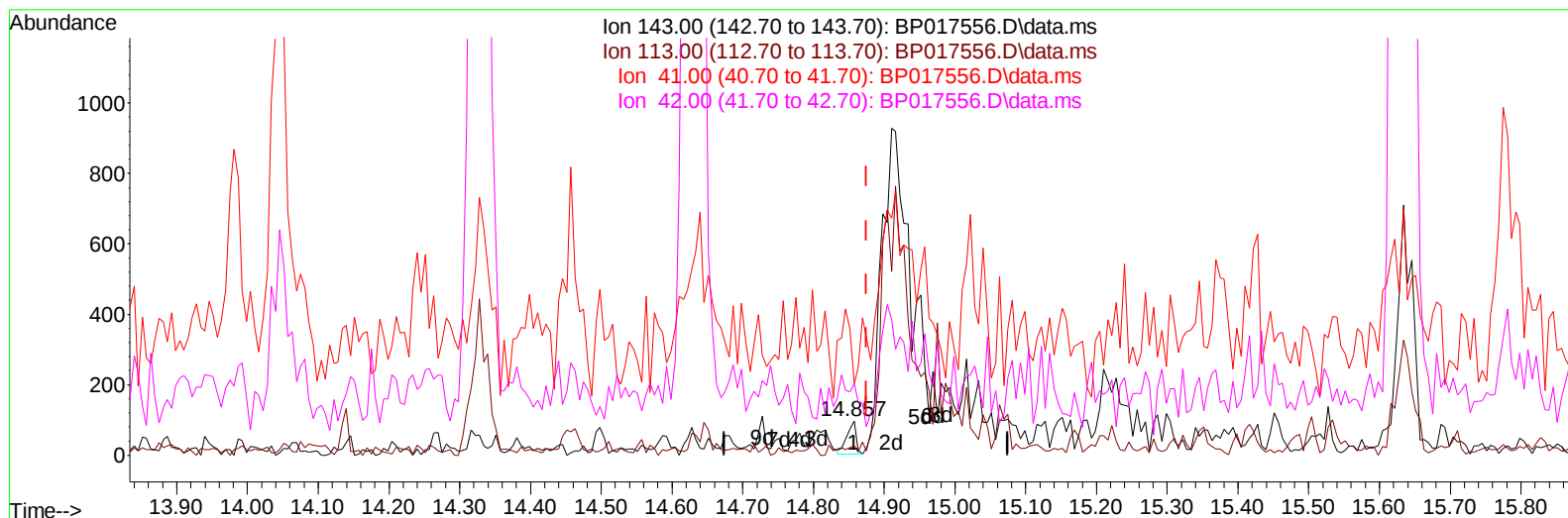
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100423\
 Data File : BP017556.D
 Acq On : 05 Oct 2023 08:27
 Operator : MA/JU
 Sample : 04679-20
 Misc :
 ALS Vial : 29 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 FB(20230927)

Manual IntegrationsAPPROVED

Quant Time: Oct 05 10:29:04 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100423.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Oct 04 16:33:47 2023
 Response via : Initial Calibration

Reviewed By :Yogesh Patel 10/05/2023
 Supervised By :mohammad ahmed 10/07/2023



TIC: BP017556.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.857min (-0.018) 0.03 ng/u1

response 82

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	81.00	26.80#
41.00	43.80	209.28#
42.00	29.00	209.28#

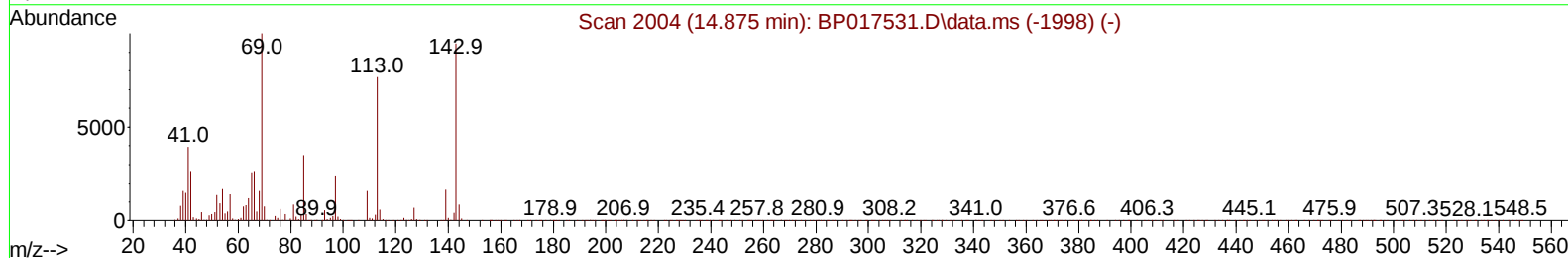
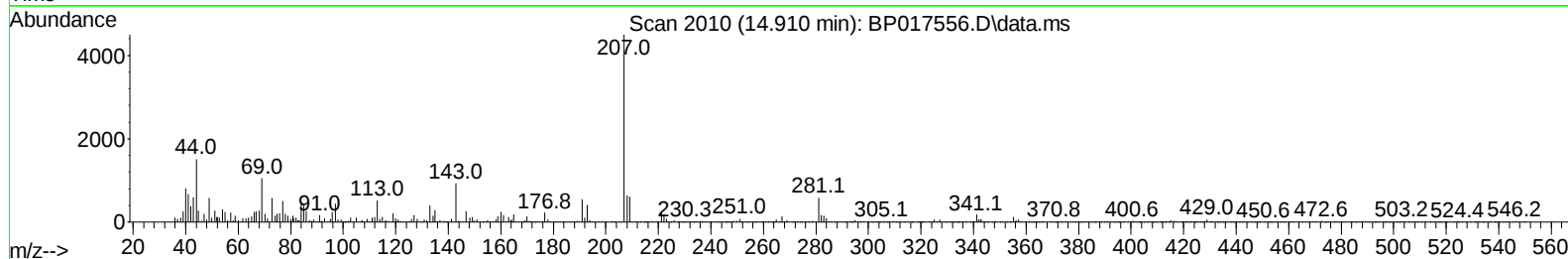
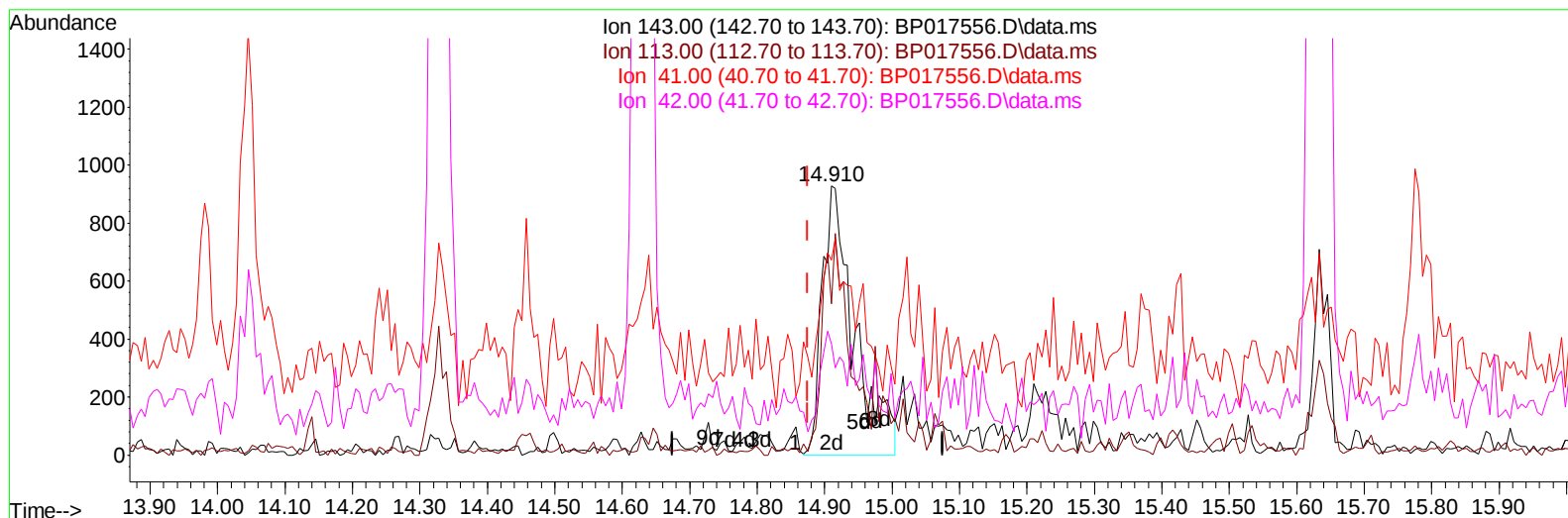
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ALS Vial : 29 Sample Multiplier: 1

Instrument :
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FB(20230927)

Manual IntegrationsAPPROVED

Quant Time: Oct 05 08:57:26 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100423.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Oct 04 16:33:47 2023
Response via : Initial Calibration

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Supervised By :mohammad ahmed 10/07/2023



TIC: BP017556.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.910min (+ 0.035) 0.97 ng/ul m

response 3072

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	81.00	56.25#
41.00	43.80	72.41#
42.00	29.00	41.38#

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100423\
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.940	152	98939	20.000	ng/ul	0.00
20) Naphthalene-d8	10.775	136	389326	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.628	164	231977	20.000	ng/ul	-0.01
64) Phenanthrene-d10	17.416	188	513197	20.000	ng/ul	-0.01
79) Chrysene-d12	21.557	240	513670	20.000	ng/ul	0.00
88) Perylene-d12	24.139	264	608696	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.287	96	9086	3.671	ng/uL	0.00
4) Pyridine-d5	3.728	84	71692	11.068	ng/ul	0.00
7) Phenol-d5	7.099	99	20359	2.446	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.281	67	137935	27.155	ng/ul	0.00
11) 2-Chlorophenol-d4	7.458	132	60035	9.159	ng/ul	0.00
15) 4-Methylphenol-d8	8.658	113	48259	7.242	ng/ul	-0.01
21) Nitrobenzene-d5	9.146	128	81653	27.261	ng/ul	0.00
24) 2-Nitrophenol-d4	9.863	143	35383	10.474	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.393	165	65253	10.419	ng/ul	0.00
31) 4-Chloroaniline-d4	10.946	131	238210	26.700	ng/ul	0.00
46) Dimethylphthalate-d6	14.046	166	535386	29.588	ng/ul	0.00
49) Acenaphthylene-d8	14.328	160	560044	28.068	ng/ul	0.00
54) 4-Nitrophenol-d4	14.910	143	3072m	0.969	ng/ul	0.04
60) Fluorene-d10	15.634	176	478309	31.164	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.781	200	10977	3.431	ng/ul	0.00
73) Anthracene-d10	17.522	188	820070	34.521	ng/ul	0.00
81) Pyrene-d10	19.775	212	1024463	35.892	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.969	264	1107804	36.063	ng/ul	0.00

Target Compounds Qvalue

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

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