

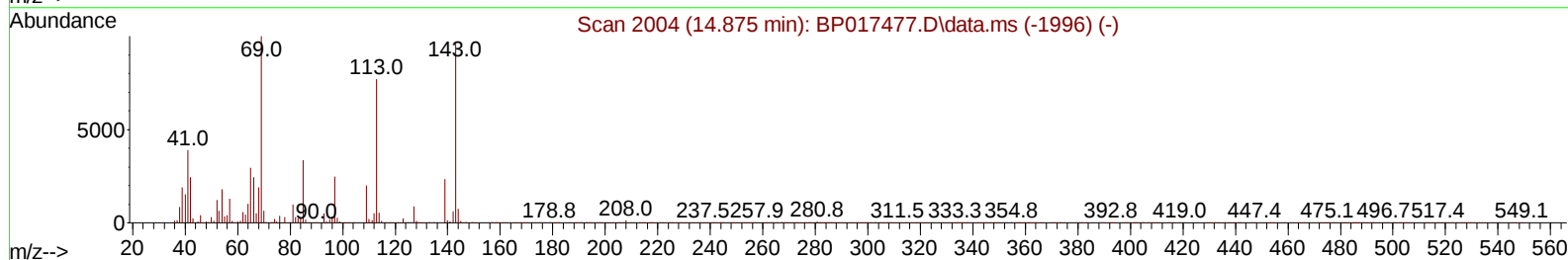
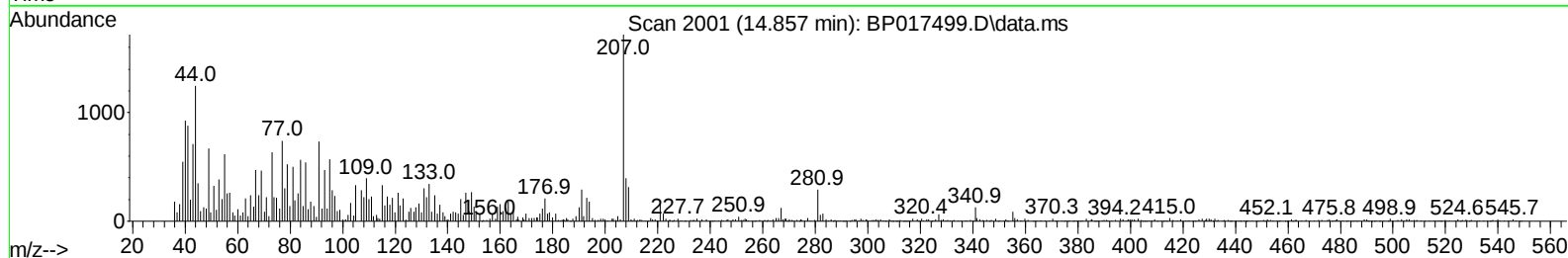
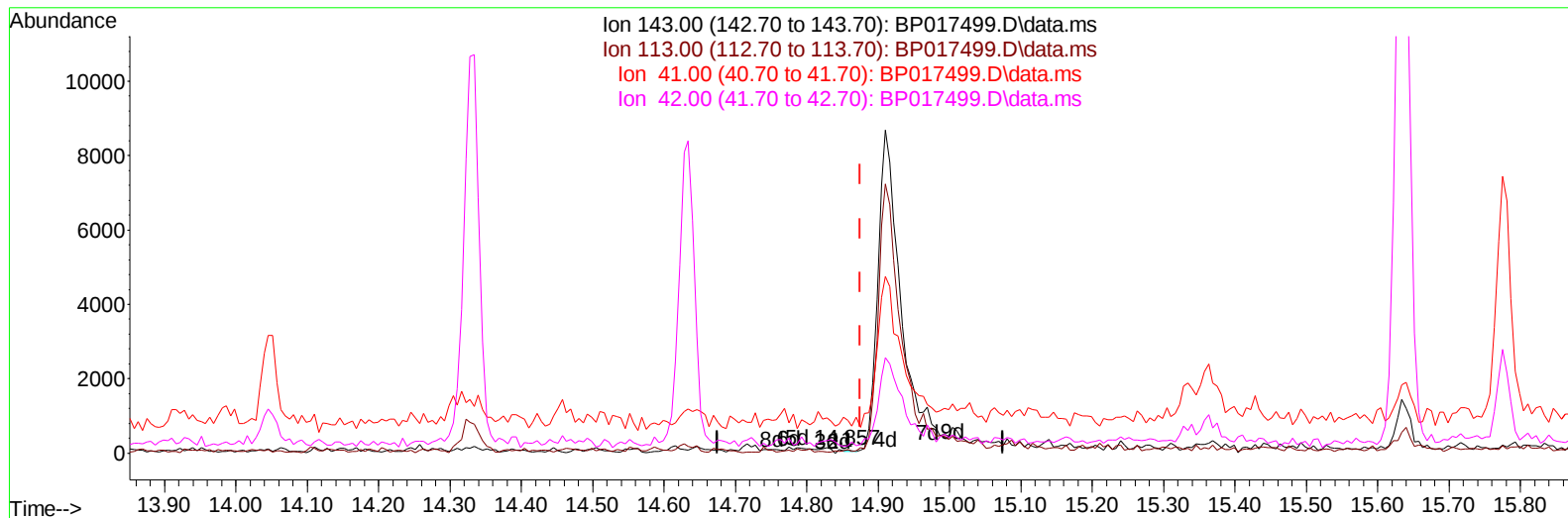
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100223\
Data File : BP017499.D
Acq On : 03 Oct 2023 04:12
Operator : MA/JU
Sample : 04571-13
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BBJ25

Manual IntegrationsAPPROVED

Quant Time: Oct 03 05:04:14 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092023.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Sat Sep 30 04:02:20 2023
Response via : Initial Calibration

Reviewed By :Yogesh Patel 10/03/2023
Supervised By :mohammad ahmed 10/05/2023



TIC: BP017499.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.857min (-0.017) 0.01 ng/u1

response 25

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	74.90	104.76#
41.00	36.70	1046.43#
42.00	26.40	236.90#

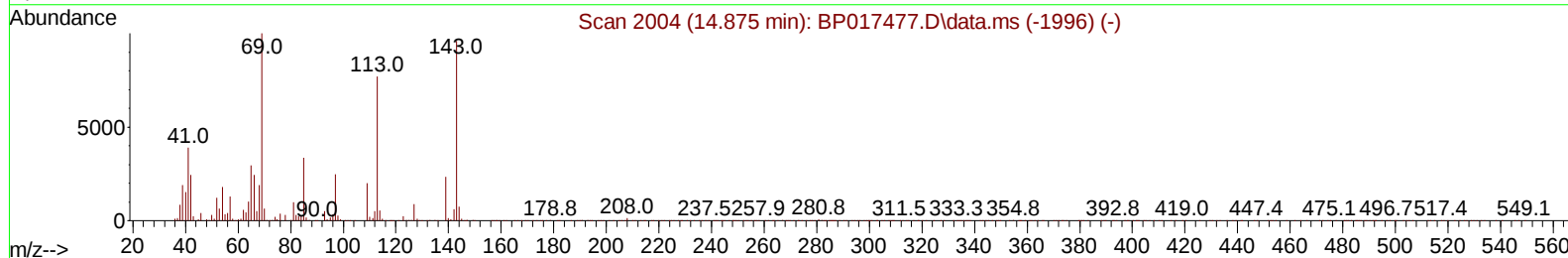
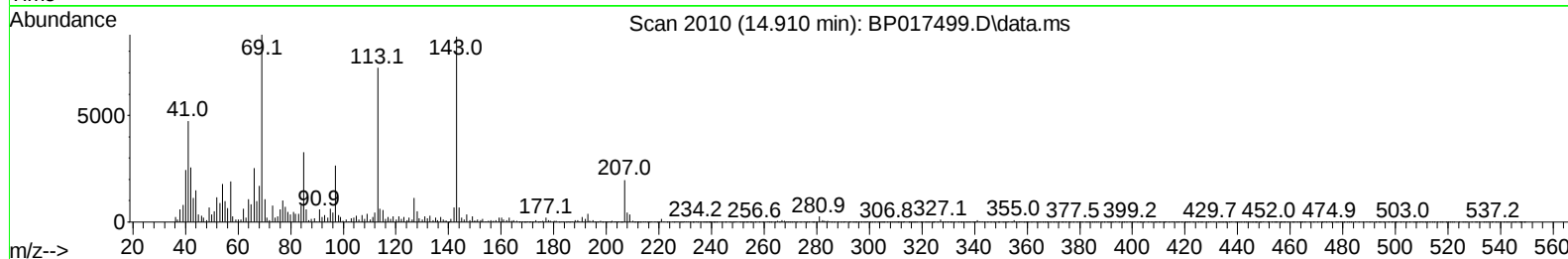
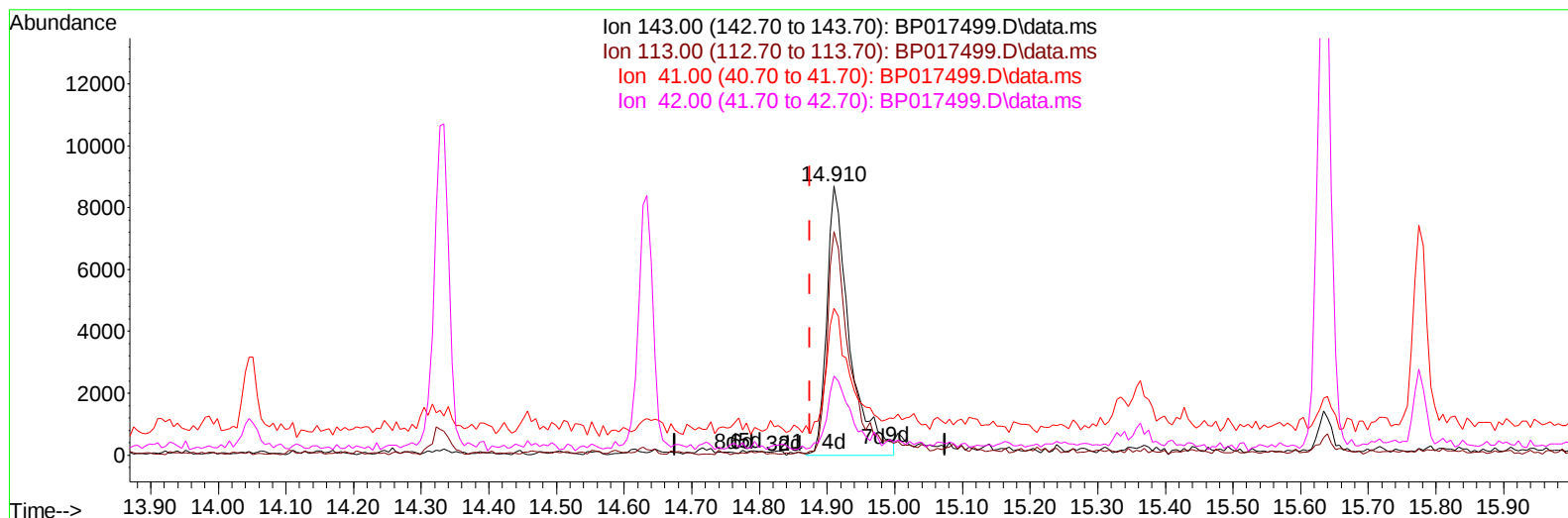
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TIC: BP017499.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.910min (+ 0.036) 4.80 ng/ul m

response 20317

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	74.90	83.16
41.00	36.70	54.58#
42.00	26.40	29.46

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 Misc :
 ALS Vial : 23 Sample Multiplier: 1

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.934	152	142582	20.000	ng/ul	-0.02
20) Naphthalene-d8	10.775	136	584218	20.000	ng/ul	-0.01
38) Acenaphthene-d10	14.634	164	338812	20.000	ng/ul	-0.01
64) Phenanthrene-d10	17.422	188	732284	20.000	ng/ul	0.00
79) Chrysene-d12	21.557	240	725235	20.000	ng/ul	0.00
88) Perylene-d12	24.145	264	799579	20.000	ng/ul	0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.287	96	13857	3.992	ng/uL	0.00
4) Pyridine-d5	3.734	84	90433	9.319	ng/ul	0.00
7) Phenol-d5	7.099	99	73724	6.296	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.275	67	248934	35.225	ng/ul	-0.02
11) 2-Chlorophenol-d4	7.458	132	243255	26.191	ng/ul	-0.01
15) 4-Methylphenol-d8	8.664	113	149317	16.401	ng/ul	0.00
21) Nitrobenzene-d5	9.146	128	162172	38.397	ng/ul	-0.01
24) 2-Nitrophenol-d4	9.863	143	172800	37.680	ng/ul	-0.01
28) 2,4-Dichlorophenol-d3	10.399	165	295938	34.156	ng/ul	0.00
31) 4-Chloroaniline-d4	10.952	131	347338	27.674	ng/ul	0.00
46) Dimethylphthalate-d6	14.046	166	1050407	43.277	ng/ul	-0.01
49) Acenaphthylene-d8	14.334	160	1135959	40.693	ng/ul	0.00
54) 4-Nitrophenol-d4	14.910	143	20317m	4.805	ng/ul	0.04
60) Fluorene-d10	15.634	176	892392	42.707	ng/ul	-0.01
65) 4,6-Dinitro-2-methylph...	15.775	200	105819	25.639	ng/ul	-0.01
73) Anthracene-d10	17.522	188	1530809	47.135	ng/ul	0.00
81) Pyrene-d10	19.775	212	1873317	47.766	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.974	264	1813880	47.117	ng/ul	0.02

Target Compounds Qvalue

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

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