

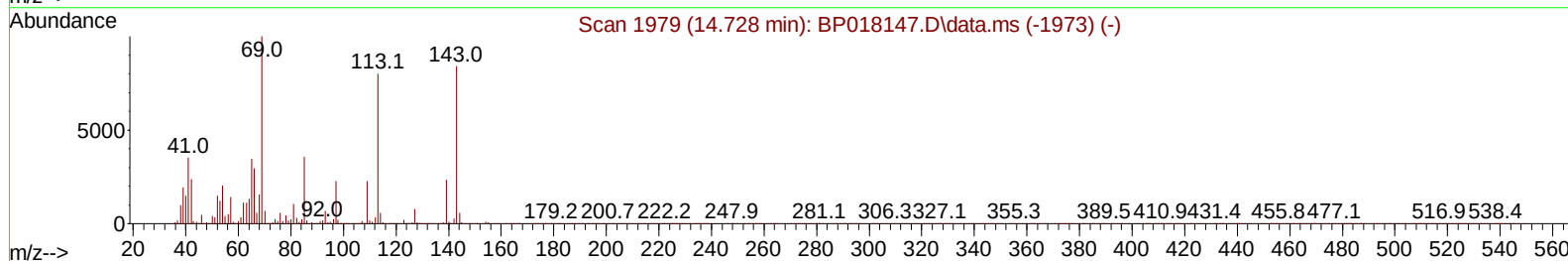
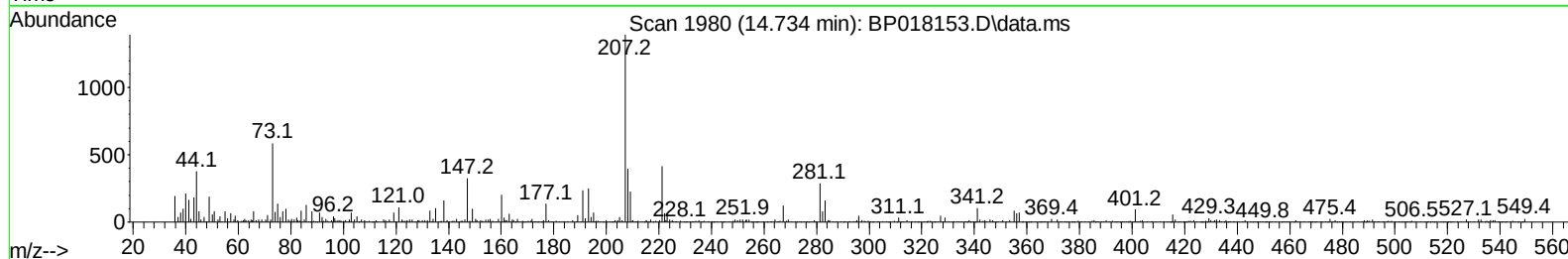
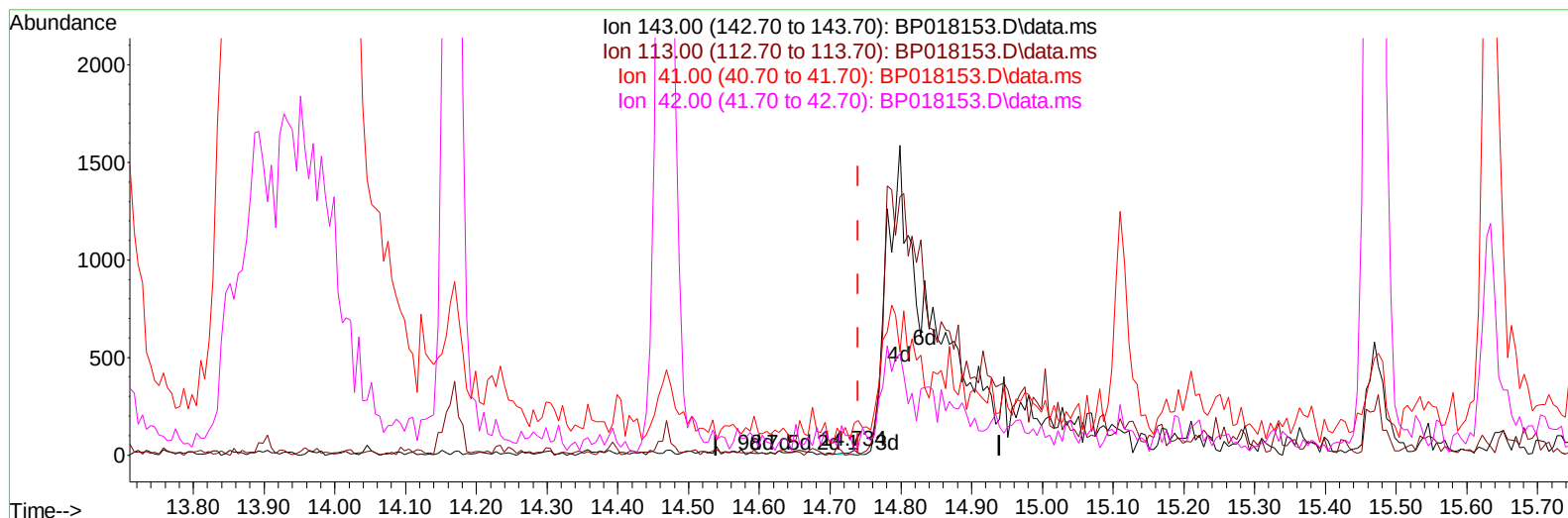
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111423\
 Data File : BP018153.D
 Acq On : 14 Nov 2023 13:54
 Operator : MA/JU
 Sample : 05337-01
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 GCDH6

Manual IntegrationsAPPROVED

Quant Time: Nov 15 00:37:52 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP111023.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Nov 13 15:50:18 2023
 Response via : Initial Calibration

Reviewed By :Yogesh Patel 11/16/2023
 Supervised By :mohammad ahmed 11/16/2023



TIC: BP018153.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.734min (-0.006) 0.01 ng/ul

response 18

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	103.00	12.50#
41.00	47.70	679.17#
42.00	30.10	95.83#

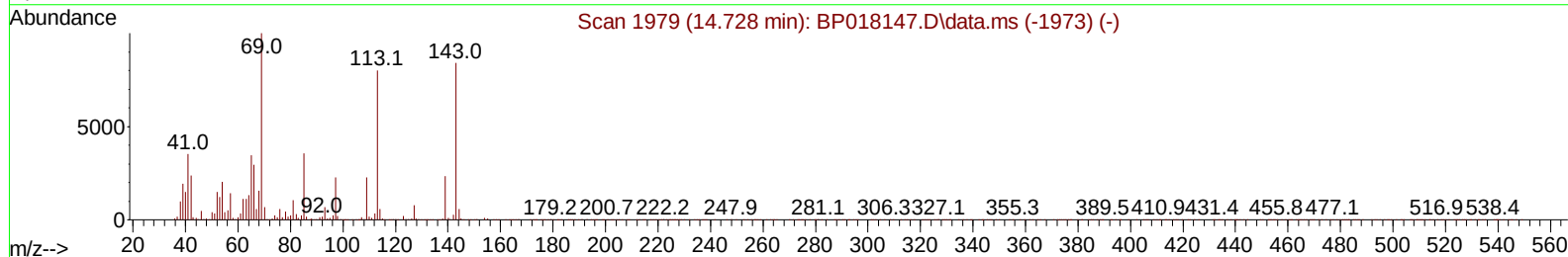
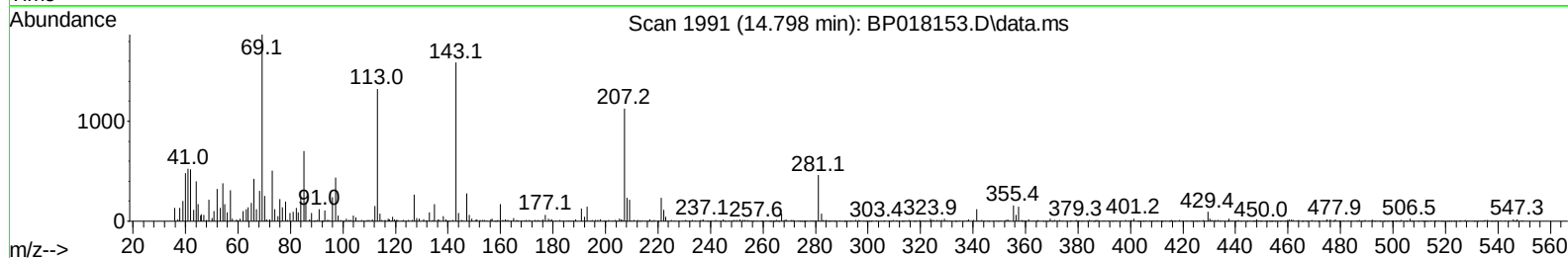
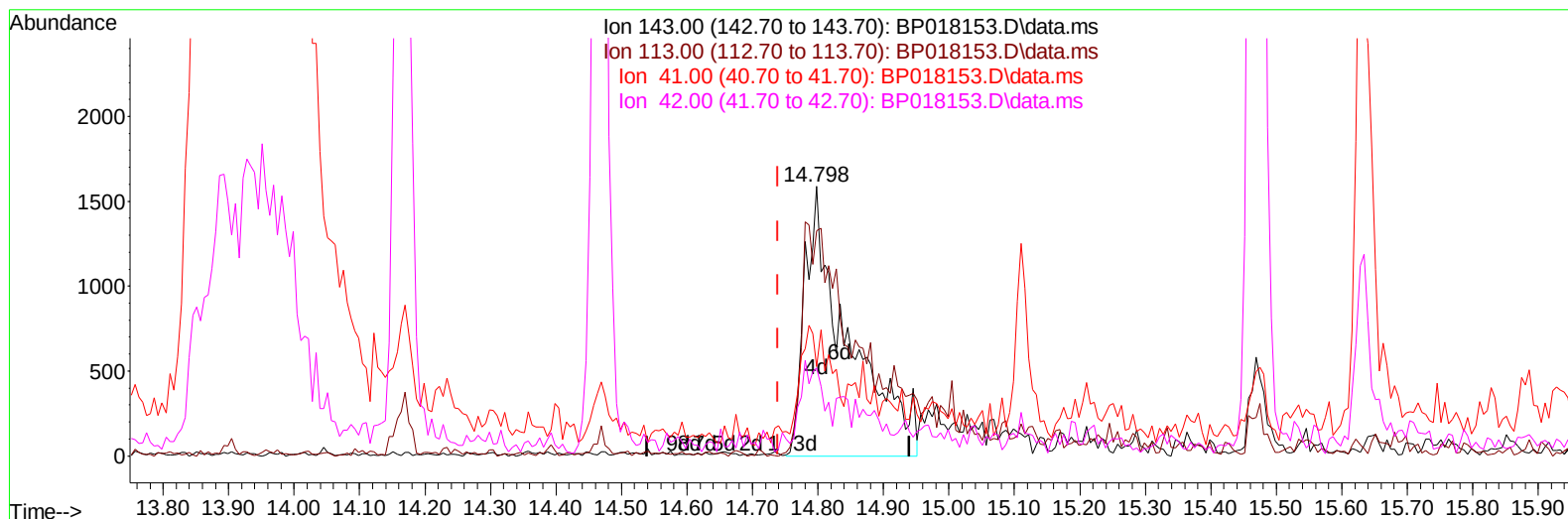
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TIC: BP018153.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.798min (+ 0.059) 2.95 ng/ul m

response 7239

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	103.00	83.49
41.00	47.70	33.21#
42.00	30.10	32.64

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Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 11/16/2023
 Supervised By :mohammad ahmed 11/16/2023

Quant Time: Nov 15 01:02:25 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP111023.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Nov 13 15:50:18 2023
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.787	152	85326	20.000	ng/ul	0.00
20) Naphthalene-d8	10.604	136	343243	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.469	164	177909	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.251	188	412798	20.000	ng/ul	0.00
79) Chrysene-d12	21.380	240	500149	20.000	ng/ul	#-0.01
88) Perylene-d12	23.845	264	560346	20.000	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.164	96	8655	3.578	ng/uL	0.00
4) Pyridine-d5	0.000	84	0d	0.000	ng/ul	
7) Phenol-d5	6.969	99	35127	4.264	ng/ul	0.01
9) Bis-(2-Chloroethyl)eth...	7.140	67	119225	24.628	ng/ul	0.00
11) 2-Chlorophenol-d4	7.311	132	111884	17.151	ng/ul	0.00
15) 4-Methylphenol-d8	8.516	113	61905	9.171	ng/ul	0.00
21) Nitrobenzene-d5	8.999	128	68886	22.276	ng/ul	0.00
24) 2-Nitrophenol-d4	9.710	143	73823	21.072	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.240	165	114896	18.400	ng/ul	0.00
31) 4-Chloroaniline-d4	10.799	131	90964	10.623	ng/ul	0.01
46) Dimethylphthalate-d6	13.893	166	426815	26.001	ng/ul	0.00
49) Acenaphthylene-d8	14.163	160	409954	22.929	ng/ul	0.00
54) 4-Nitrophenol-d4	14.798	143	7239m	2.951	ng/ul	0.06
60) Fluorene-d10	15.469	176	352679	26.023	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.634	200	57417	19.240	ng/ul	-0.03
73) Anthracene-d10	17.351	188	584048	26.160	ng/ul	0.00
81) Pyrene-d10	19.604	212	809184	24.221	ng/ul	-0.01
92) Benzo(a)pyrene-d12	23.680	264	838400	25.732	ng/ul	-0.02

Target Compounds Qvalue

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

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