

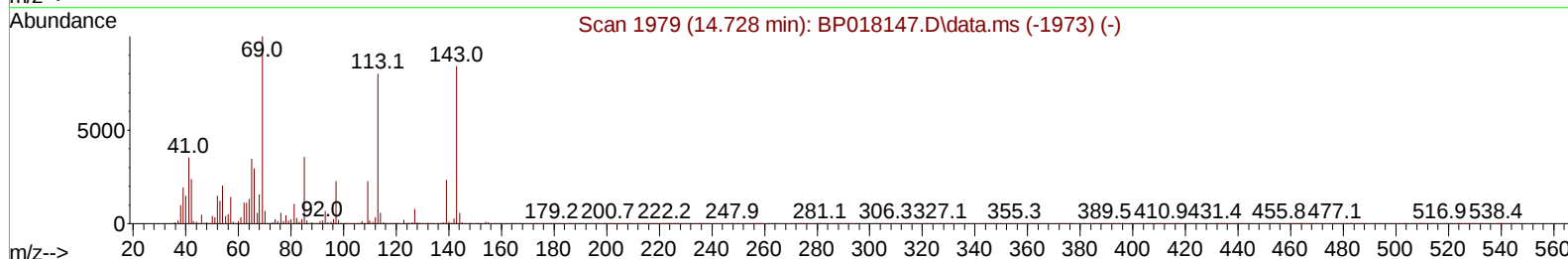
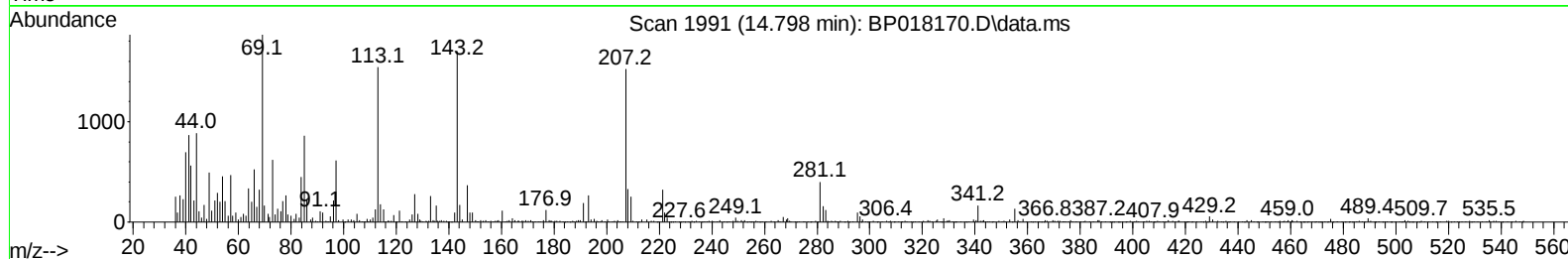
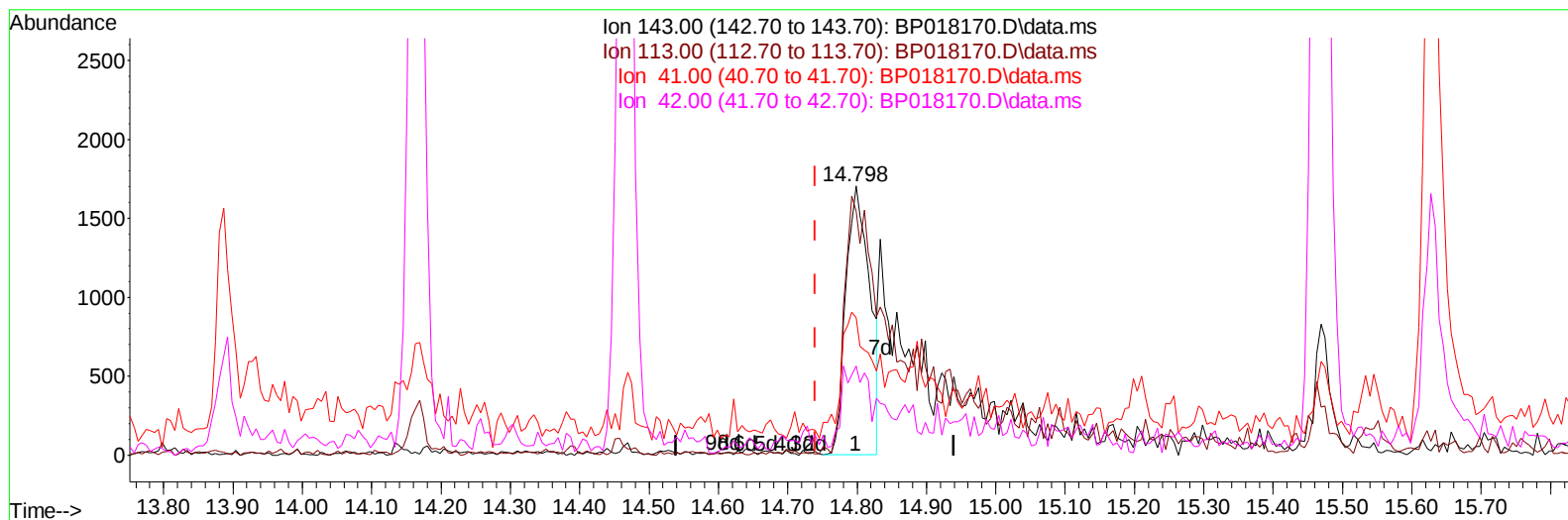
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111423\
 Data File : BP018170.D
 Acq On : 14 Nov 2023 23:29
 Operator : MA/JU
 Sample : 05337-15
 Misc :
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 GCDK0

Manual IntegrationsAPPROVED

Quant Time: Nov 15 00:41:53 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: BP018170.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.798min (+ 0.059) 1.46 ng/u1

response 4059

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	103.00	90.67
41.00	47.70	51.00
42.00	30.10	33.16

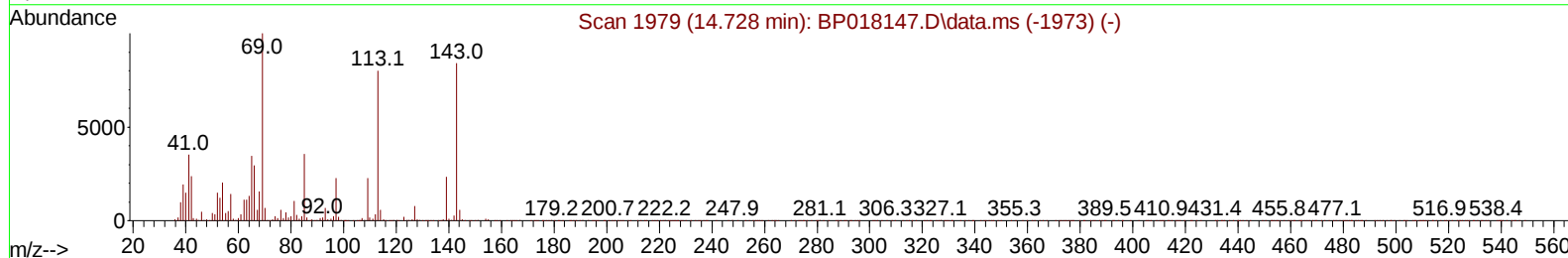
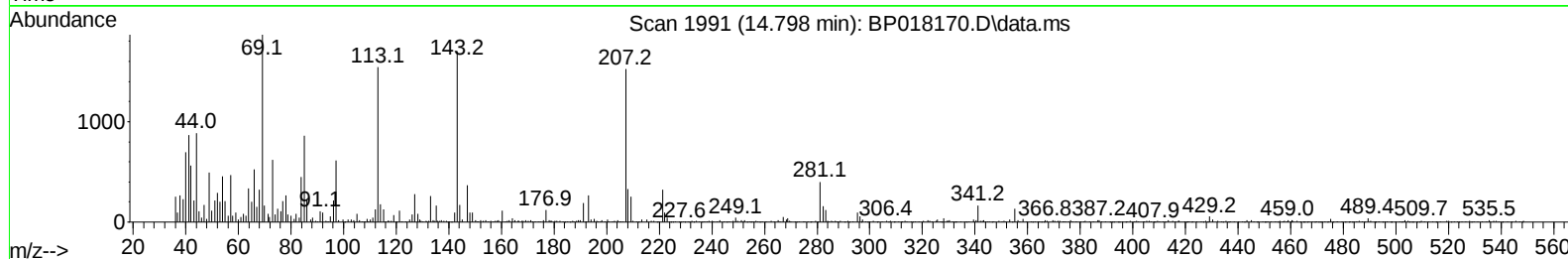
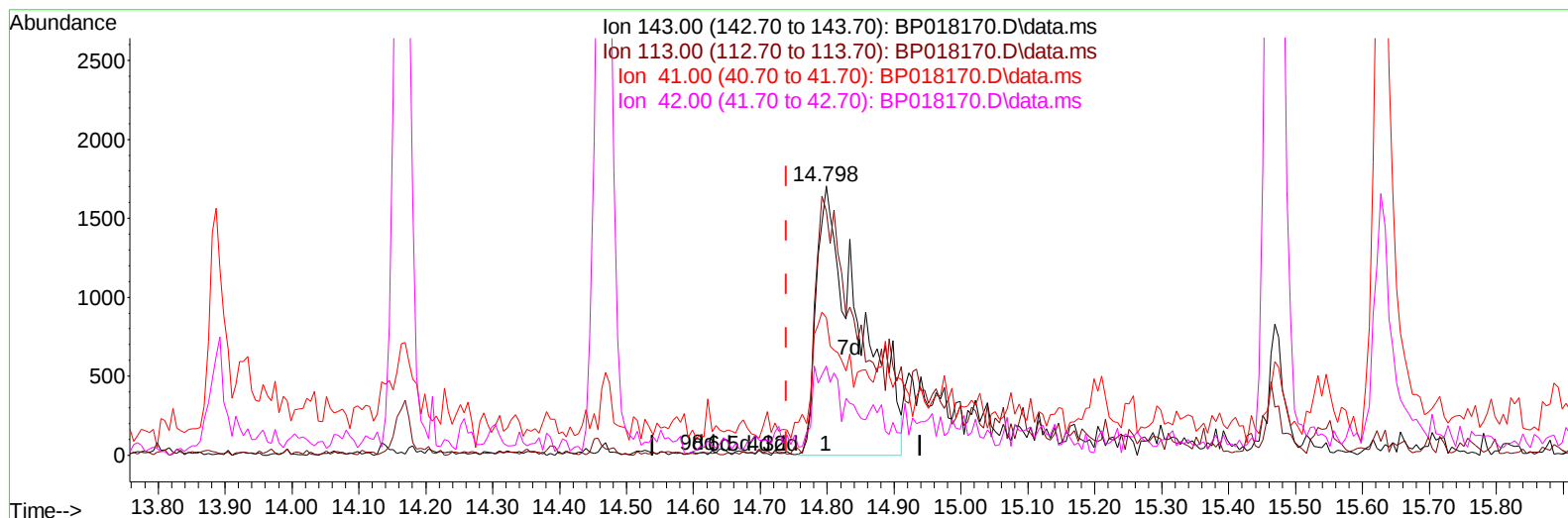
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TIC: BP018170.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.798min (+ 0.059) 2.70 ng/ul m

response 7518

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	103.00	90.67
41.00	47.70	51.00
42.00	30.10	33.16

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.787	152	77784	20.000	ng/ul	0.00
20) Naphthalene-d8	10.604	136	316730	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.469	164	201796	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.245	188	452000	20.000	ng/ul	-0.01
79) Chrysene-d12	21.380	240	437016	20.000	ng/ul	-0.01
88) Perylene-d12	23.845	264	478752	20.000	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.164	96	6646	3.014	ng/uL	0.00
4) Pyridine-d5	3.617	84	21646	3.765	ng/ul	0.01
7) Phenol-d5	6.969	99	29087	3.873	ng/ul	0.01
9) Bis-(2-Chloroethyl)eth...	7.140	67	115537	26.180	ng/ul	0.00
11) 2-Chlorophenol-d4	7.310	132	106158	17.851	ng/ul	0.00
15) 4-Methylphenol-d8	8.516	113	62542	10.164	ng/ul	0.00
21) Nitrobenzene-d5	8.993	128	68411	23.974	ng/ul	0.00
24) 2-Nitrophenol-d4	9.704	143	74574	23.068	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.240	165	111761	19.397	ng/ul	0.00
31) 4-Chloroaniline-d4	10.793	131	128673	16.284	ng/ul	0.00
46) Dimethylphthalate-d6	13.887	166	525423	28.219	ng/ul	-0.01
49) Acenaphthylene-d8	14.163	160	497500	24.532	ng/ul	0.00
54) 4-Nitrophenol-d4	14.798	143	7518m	2.702	ng/ul	0.06
60) Fluorene-d10	15.469	176	422052	27.455	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.628	200	73393	22.461	ng/ul	-0.04
73) Anthracene-d10	17.351	188	747550	30.580	ng/ul	0.00
81) Pyrene-d10	19.604	212	959384	32.866	ng/ul	-0.01
92) Benzo(a)pyrene-d12	23.686	264	919948	33.046	ng/ul	-0.02

Target Compounds Qvalue

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

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