

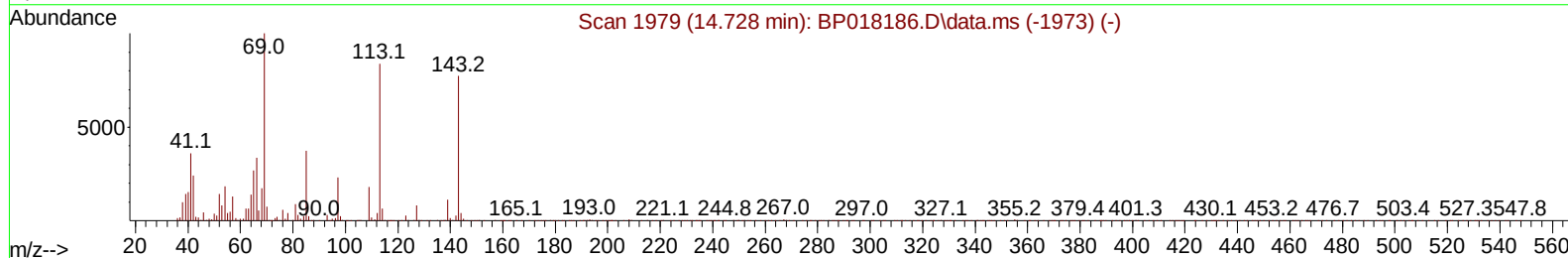
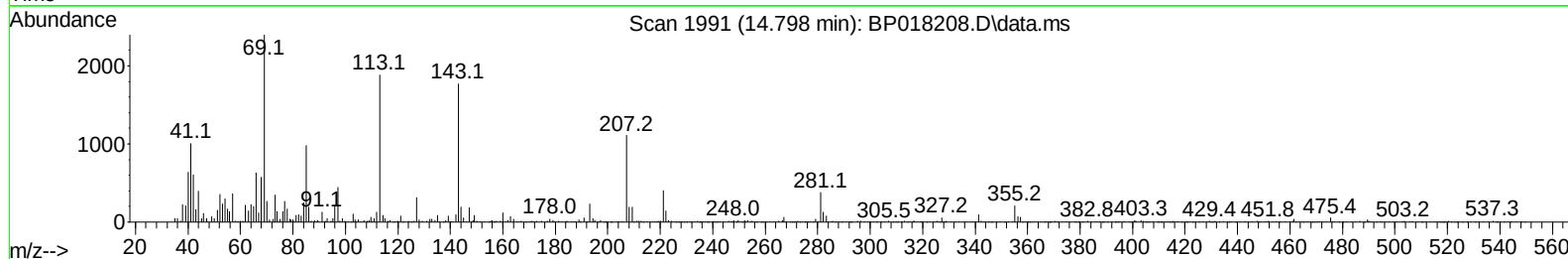
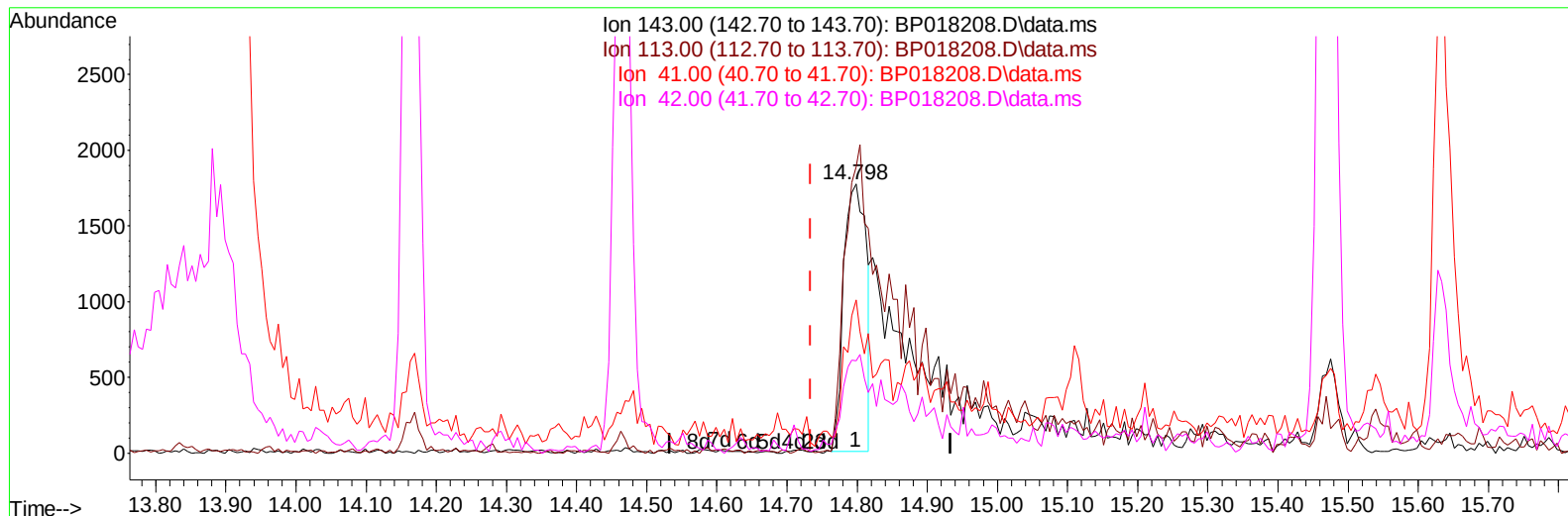
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111523\
 Data File : BP018208.D
 Acq On : 16 Nov 2023 11:52
 Operator : MA/JU
 Sample : 05338-13
 Misc :
 ALS Vial : 36 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 GCDL8

Manual IntegrationsAPPROVED

Quant Time: Nov 16 21:38:35 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP111023.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Nov 15 18:13:37 2023
 Response via : Initial Calibration

Reviewed By :Yogesh Patel 11/17/2023
 Supervised By :mohammad ahmed 11/17/2023



TIC: BP018208.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.798min (+ 0.065) 1.90 ng/ul

response 3950

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	103.00	106.42
41.00	47.70	56.93
42.00	30.10	34.23

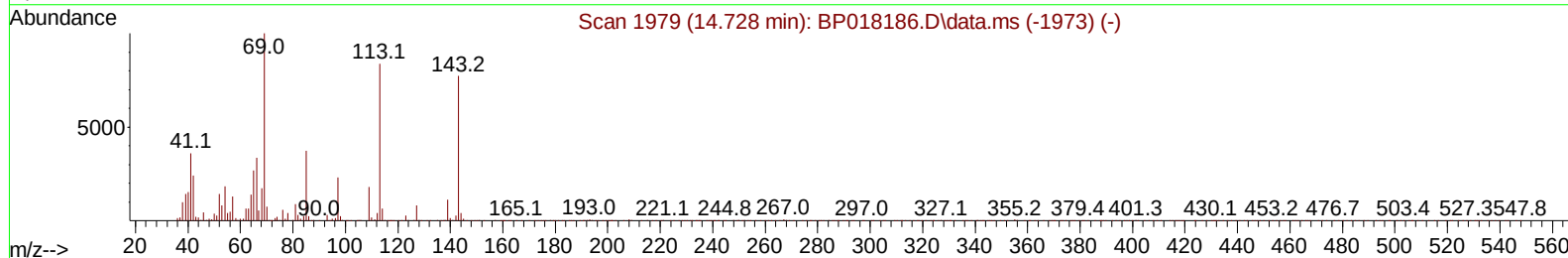
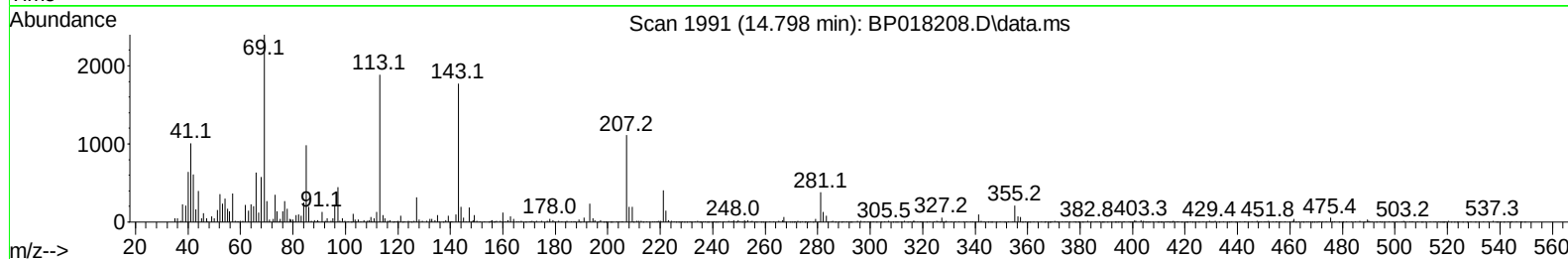
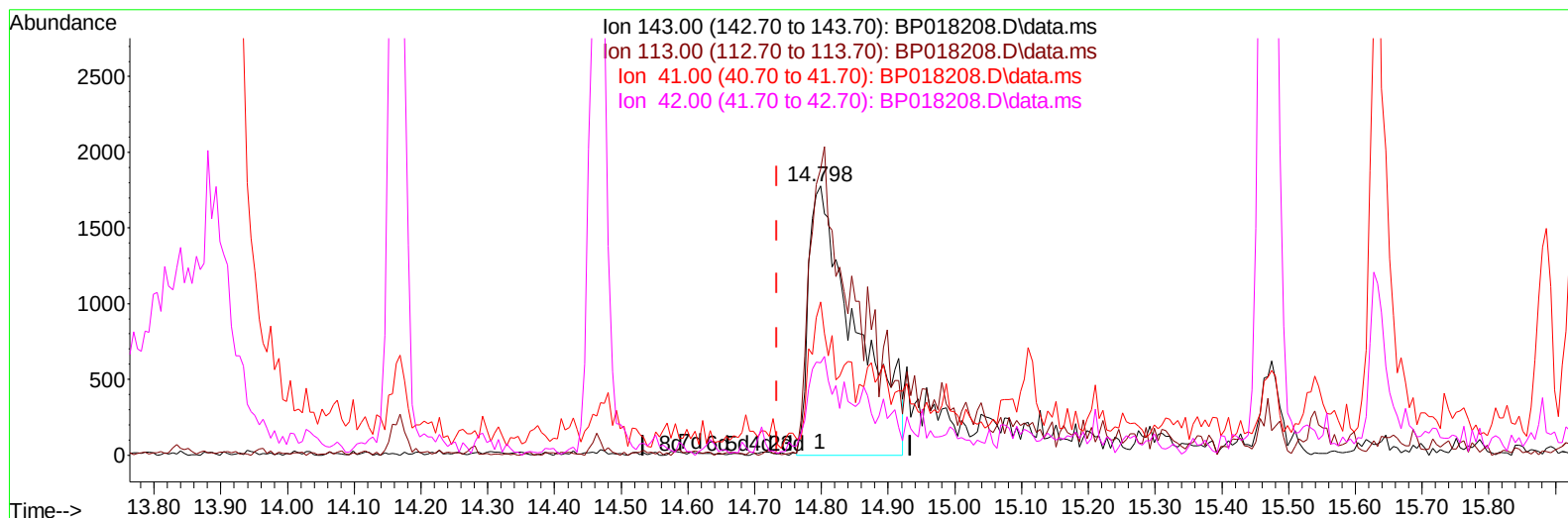
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TIC: BP018208.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.798min (+ 0.065) 4.16 ng/ul m

response 8657

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	103.00	106.42
41.00	47.70	56.93
42.00	30.10	34.23

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Manual IntegrationsAPPROVED

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Quant Time: Nov 16 22:04:17 2023
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 Quant Title : SVOA CALIBRATION
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.781	152	65355	20.000	ng/ul	0.00
20) Naphthalene-d8	10.599	136	220368	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.469	164	150922	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.251	188	357932	20.000	ng/ul	0.00
79) Chrysene-d12	21.380	240	376940	20.000	ng/ul	0.00
88) Perylene-d12	23.857	264	425244	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.164	96	6976	3.765	ng/uL	0.00
4) Pyridine-d5	3.605	84	26502	5.486	ng/ul	0.01
7) Phenol-d5	6.963	99	39045	6.188	ng/ul	0.01
9) Bis-(2-Chloroethyl)eth...	7.134	67	100256	27.038	ng/ul	0.00
11) 2-Chlorophenol-d4	7.311	132	108963	21.808	ng/ul	0.00
15) 4-Methylphenol-d8	8.516	113	65363	12.642	ng/ul	0.00
21) Nitrobenzene-d5	8.993	128	58948	29.691	ng/ul	0.00
24) 2-Nitrophenol-d4	9.705	143	67558	30.036	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.234	165	93716	23.377	ng/ul	0.00
31) 4-Chloroaniline-d4	10.799	131	47838	8.702	ng/ul	0.01
46) Dimethylphthalate-d6	13.887	166	410655	29.490	ng/ul	0.00
49) Acenaphthylene-d8	14.163	160	409708	27.013	ng/ul	0.00
54) 4-Nitrophenol-d4	14.798	143	8657m	4.160	ng/ul	0.06
60) Fluorene-d10	15.469	176	338041	29.403	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.634	200	57432	22.195	ng/ul	0.00
73) Anthracene-d10	17.351	188	587330	30.340	ng/ul	0.00
81) Pyrene-d10	19.610	212	788981	31.336	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.692	264	764967	30.937	ng/ul	0.00

Target Compounds Qvalue

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

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