

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110923\
 Data File : BP018002.D
 Acq On : 09 Nov 2023 22:57
 Operator : MA/JU
 Sample : 05082-10
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

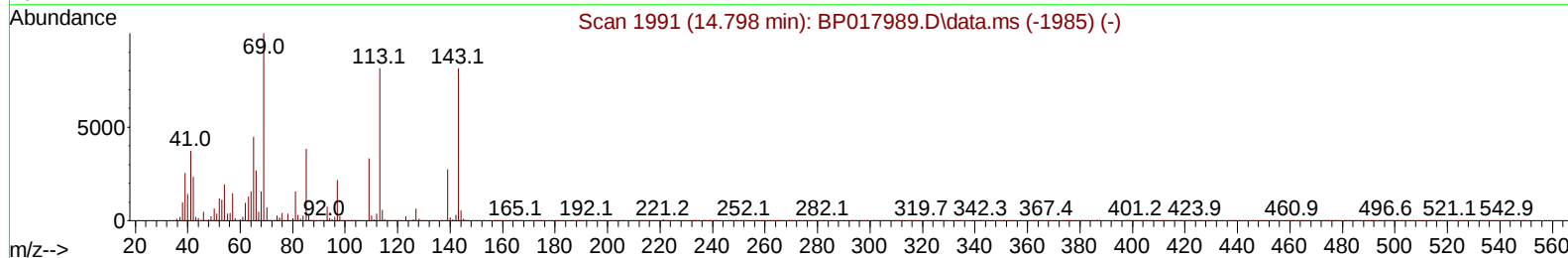
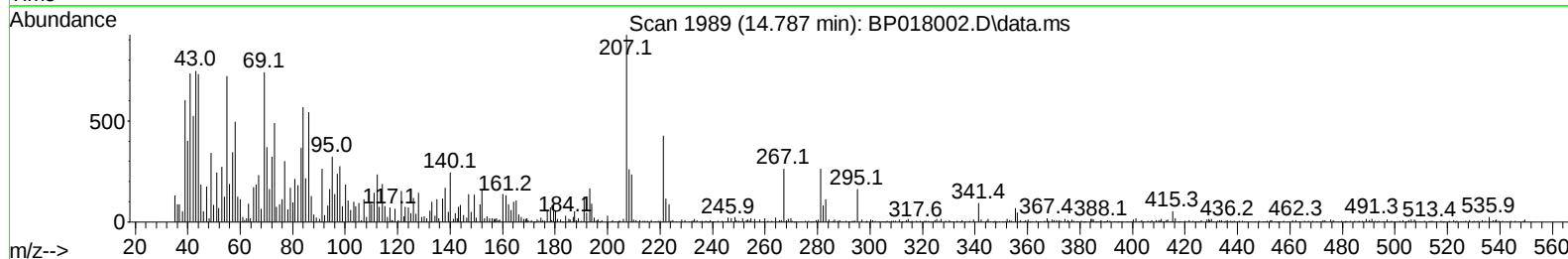
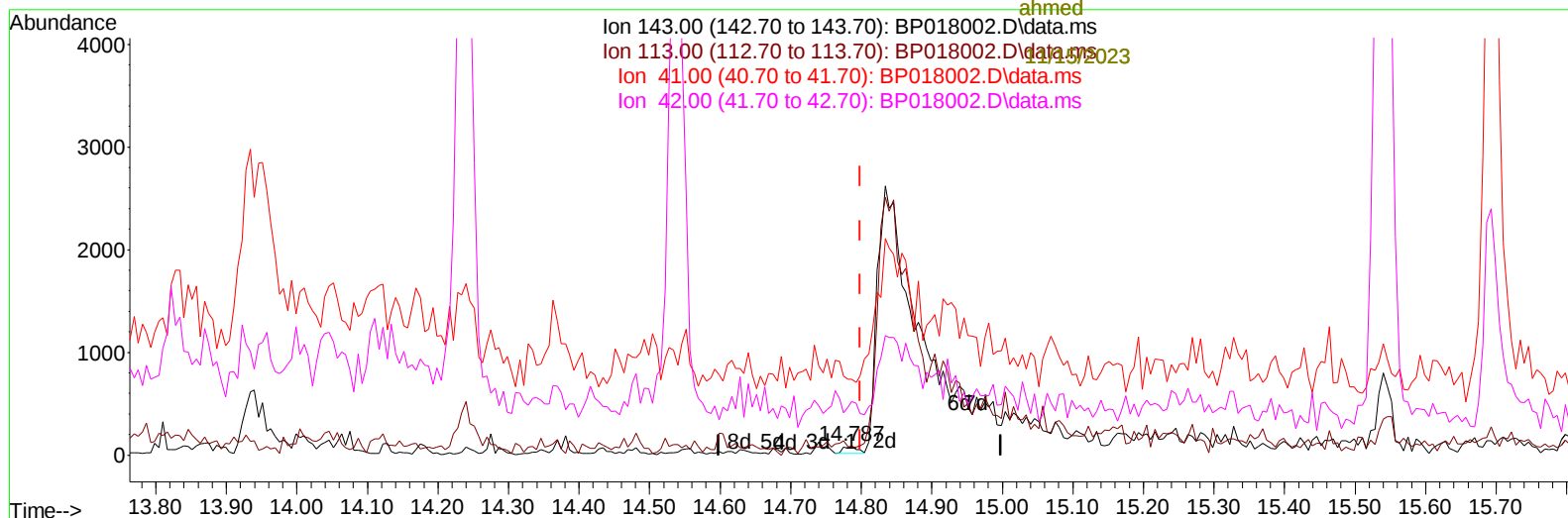
Instrument :
 BNA_P
 ClientSampleId :
 BH351

Manual IntegrationsAPPROVED

Quant Time: Nov 09 23:37:33 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP110923.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Nov 09 22:34:38 2023
 Response via : Initial Calibration

Reviewed By :Yogesh Patel 11/11/2023
 Supervised By :mohammad ahmed 11/15/2023

11/11/2023
 Supervised By :mohammad
 ahmed



TIC: BP018002.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.787min (-0.012) 0.05 ng/u1

response 108

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	100.40	81.72
41.00	46.90	791.40#
42.00	29.70	563.44#

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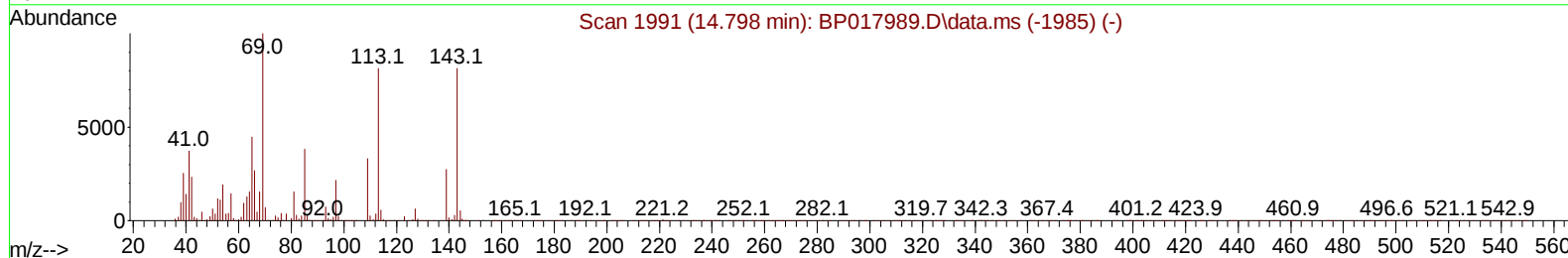
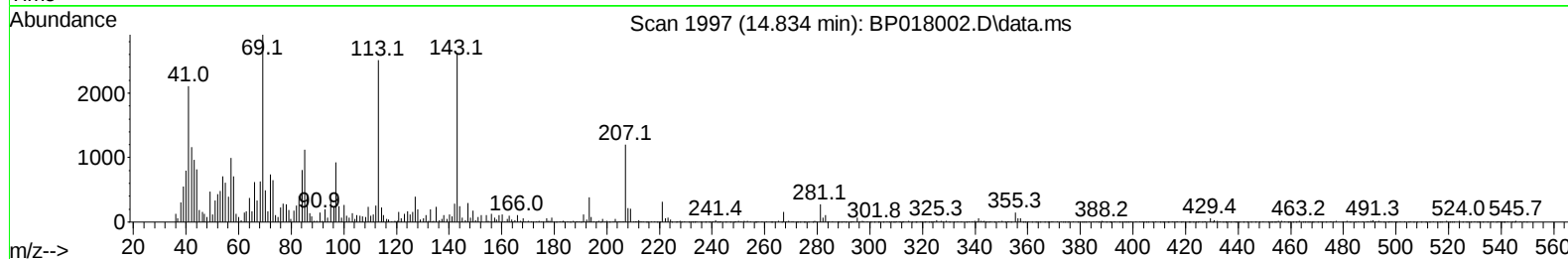
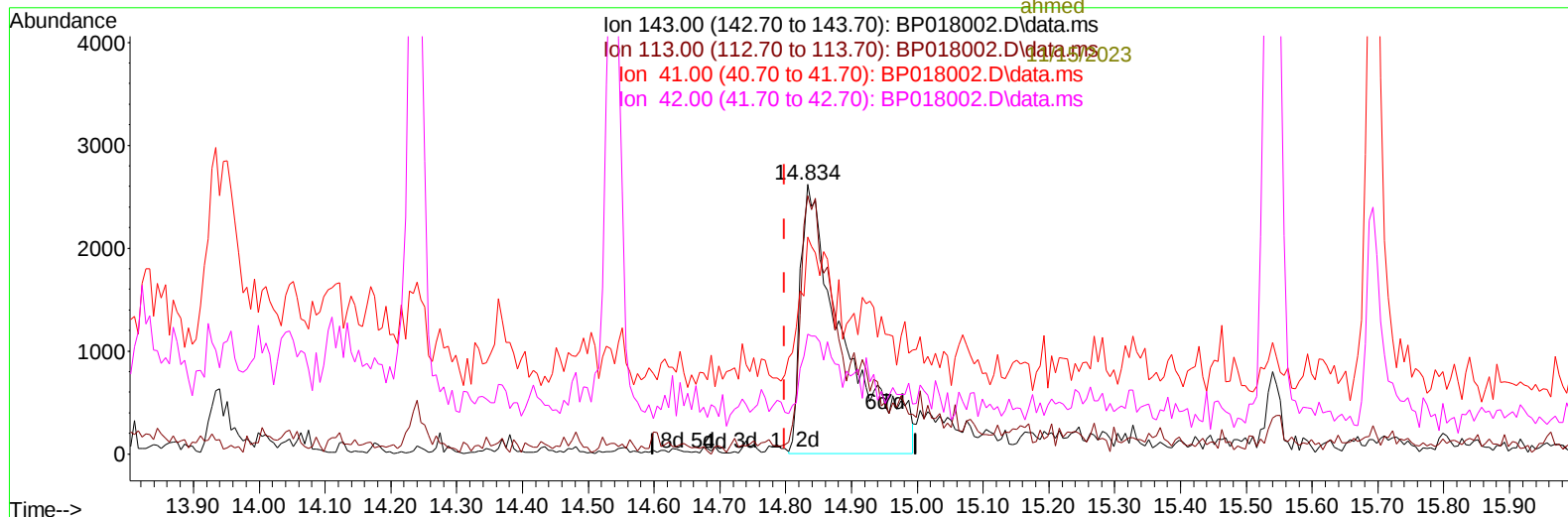
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TIC: BP018002.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.834min (+ 0.035) 5.40 ng/ul m

response 11850

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	100.40	95.80
41.00	46.90	80.41#
42.00	29.70	44.52#

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Quant Time: Nov 09 23:38:12 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP110923.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Nov 09 22:34:38 2023
 Response via : Initial Calibration

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
-----11/15/2023						
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.864	152	53595	20.000	ng/ul	0.00
20) Naphthalene-d8	10.687	136	227553	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.540	164	162860	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.322	188	377291	20.000	ng/ul	0.00
79) Chrysene-d12	21.445	240	377109	20.000	ng/ul	0.00
88) Perylene-d12	23.951	264	426592	20.000	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.246	96	5556	3.549	ng/uL	0.00
4) Pyridine-d5	3.687	84	47748	11.734	ng/ul	0.00
7) Phenol-d5	7.058	99	39390	7.821	ng/ul	0.02
9) Bis-(2-Chloroethyl)eth...	7.216	67	94645	31.021	ng/ul	0.00
11) 2-Chlorophenol-d4	7.399	132	109943	27.282	ng/ul	0.01
15) 4-Methylphenol-d8	8.611	113	78383	19.403	ng/ul	0.02
21) Nitrobenzene-d5	9.063	128	139289	68.354	ng/ul	-0.01
24) 2-Nitrophenol-d4	9.787	143	76630	33.789	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.334	165	127977	31.500	ng/ul	0.02
31) 4-Chloroaniline-d4	10.863	131	173310	31.128	ng/ul	0.00
46) Dimethylphthalate-d6	13.957	166	511864	34.190	ng/ul	0.00
49) Acenaphthylene-d8	14.240	160	533292	33.226	ng/ul	0.00
54) 4-Nitrophenol-d4	14.834	143	11850m	5.404	ng/ul	0.04
60) Fluorene-d10	15.540	176	435989	35.135	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.693	200	86312	32.763	ng/ul	0.00
73) Anthracene-d10	17.422	188	746007	36.902	ng/ul	0.00
81) Pyrene-d10	19.669	212	945647	38.963	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.786	264	943390	38.047	ng/ul	-0.01

Target Compounds Qvalue

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

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