

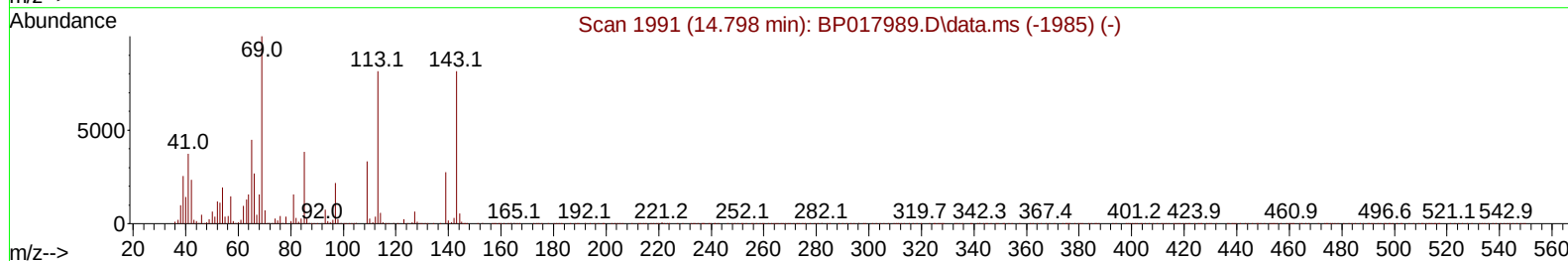
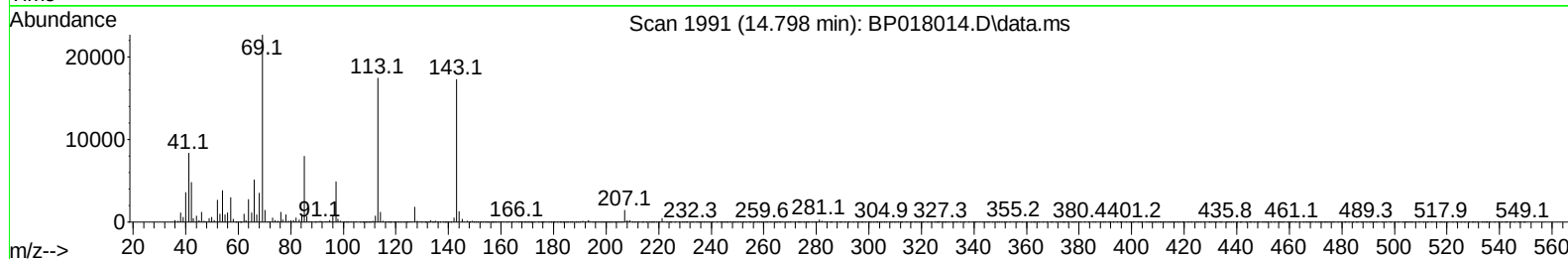
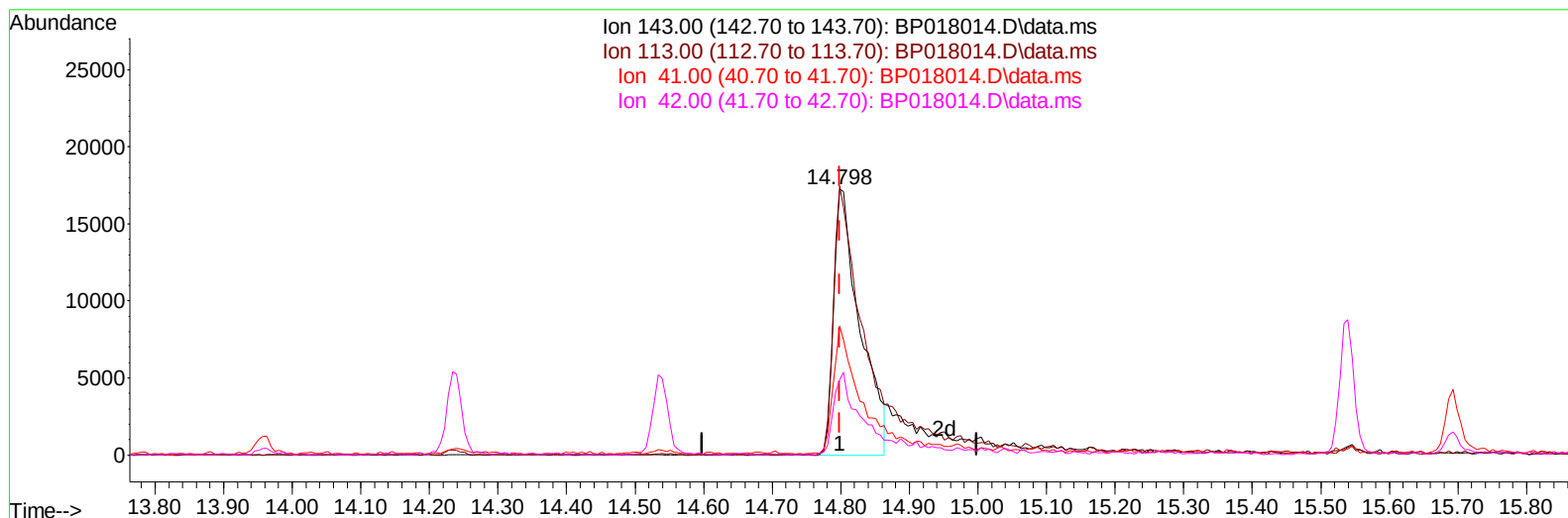
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110923\
 Data File : BP018014.D
 Acq On : 10 Nov 2023 06:18
 Operator : MA/JU
 Sample : PB156753BL
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK753

Manual IntegrationsAPPROVED

Quant Time: Nov 10 06:46:01 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



TIC: BP018014.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.798min (+ 0.000) 18.75 ng/ul

response 45862

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	100.40	101.02
41.00	46.90	48.38
42.00	29.70	28.16

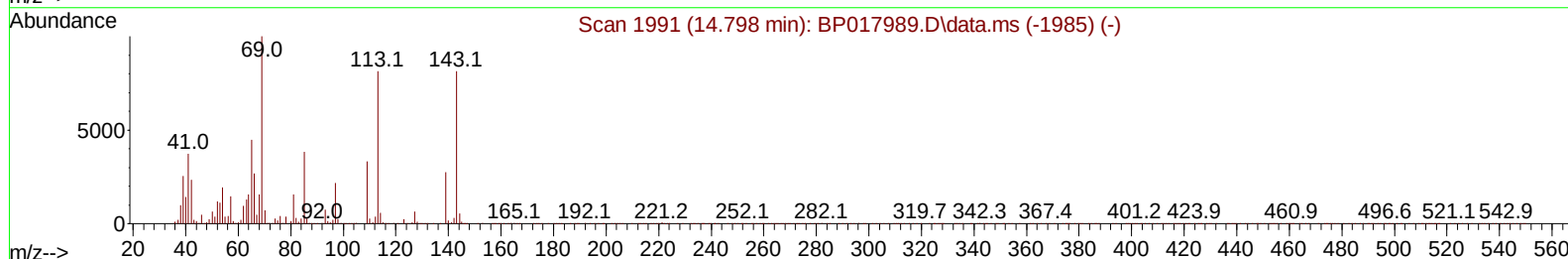
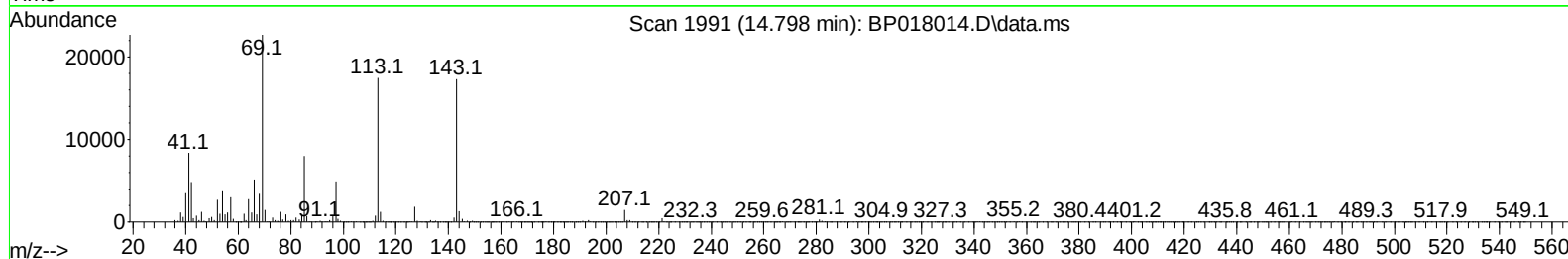
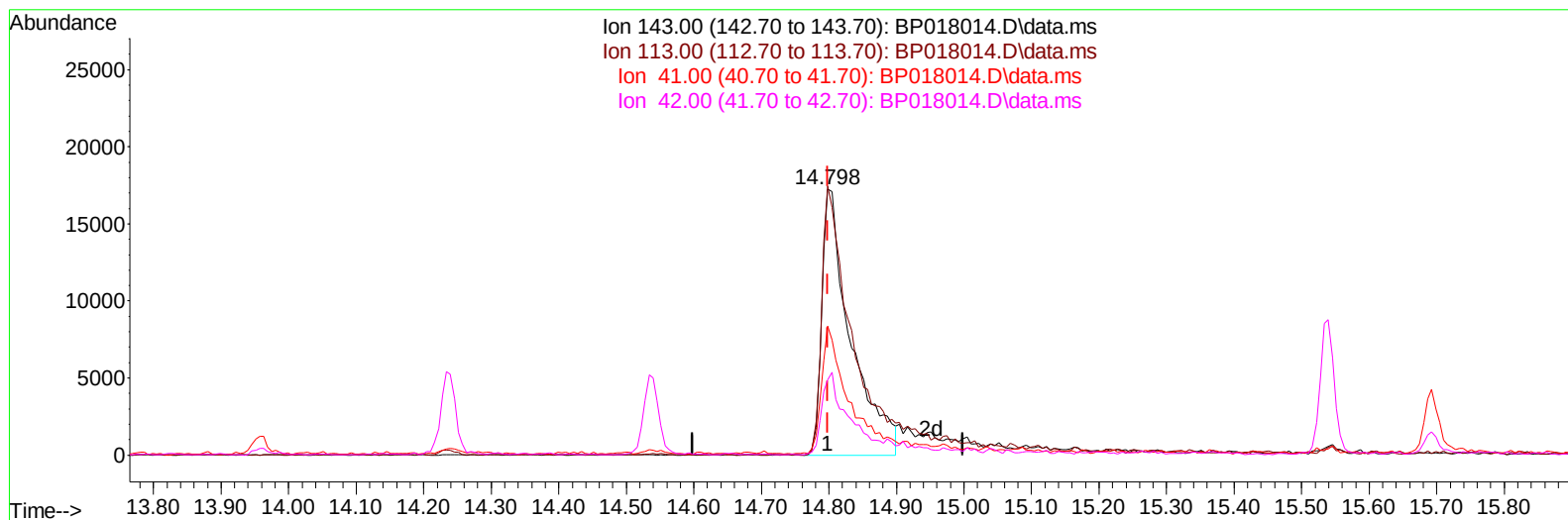
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TIC: BP018014.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.798min (+ 0.000) 20.90 ng/ul m

response 51135

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	100.40	101.02
41.00	46.90	48.38
42.00	29.70	28.16

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.863	152	71598	20.000	ng/ul	0.00
20) Naphthalene-d8	10.681	136	288976	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.540	164	181708	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.316	188	426081	20.000	ng/ul	0.00
79) Chrysene-d12	21.445	240	428592	20.000	ng/ul	0.00
88) Perylene-d12	23.951	264	477012	20.000	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.246	96	11109	5.311	ng/uL	0.00
4) Pyridine-d5	3.675	84	142075	26.136	ng/ul	0.00
7) Phenol-d5	7.028	99	167516	24.896	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.211	67	107407	26.352	ng/ul	0.00
11) 2-Chlorophenol-d4	7.387	132	137954	25.625	ng/ul	0.00
15) 4-Methylphenol-d8	8.587	113	139517	25.852	ng/ul	0.00
21) Nitrobenzene-d5	9.069	128	66744	25.792	ng/ul	0.00
24) 2-Nitrophenol-d4	9.787	143	75307	26.147	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.310	165	121556	23.560	ng/ul	0.00
31) 4-Chloroaniline-d4	10.863	131	179306	25.360	ng/ul	0.00
46) Dimethylphthalate-d6	13.957	166	462578	27.693	ng/ul	0.00
49) Acenaphthylene-d8	14.234	160	476241	26.594	ng/ul	0.00
54) 4-Nitrophenol-d4	14.798	143	51135m	20.901	ng/ul	0.00
60) Fluorene-d10	15.540	176	370846	26.785	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.692	200	63793	21.442	ng/ul	0.00
73) Anthracene-d10	17.416	188	613233	26.861	ng/ul	0.00
81) Pyrene-d10	19.669	212	748228	27.126	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.786	264	756478	27.284	ng/ul	-0.01

Target Compounds Qvalue

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

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