

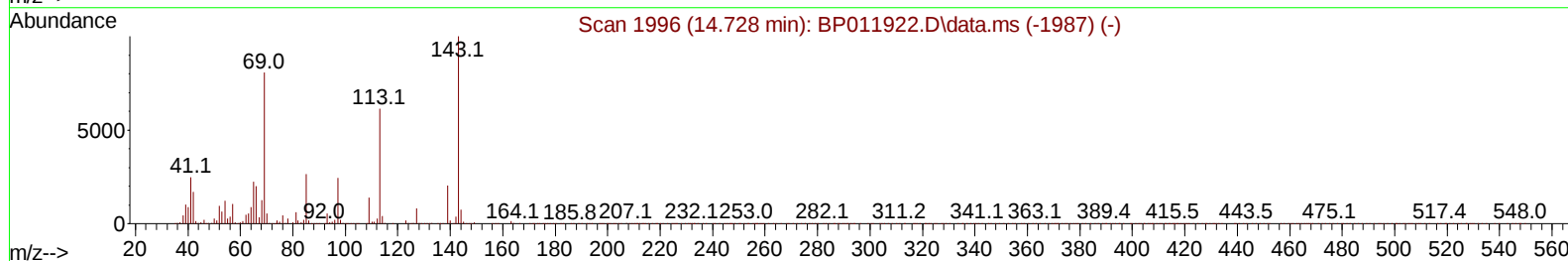
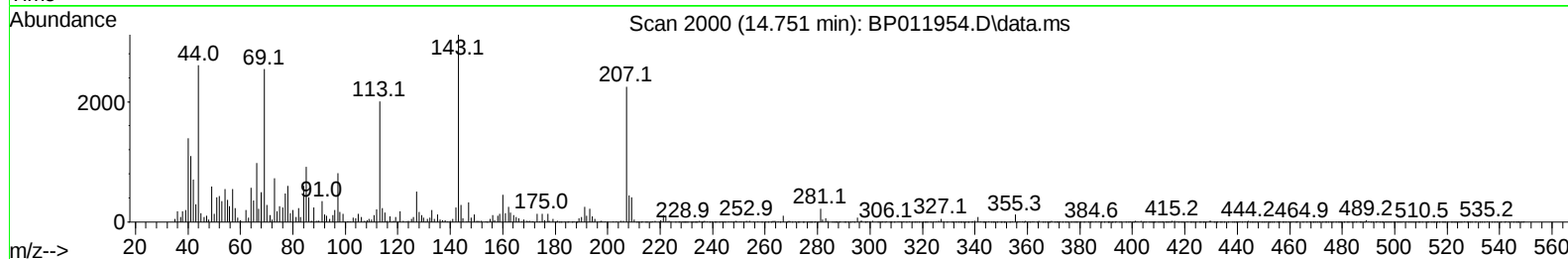
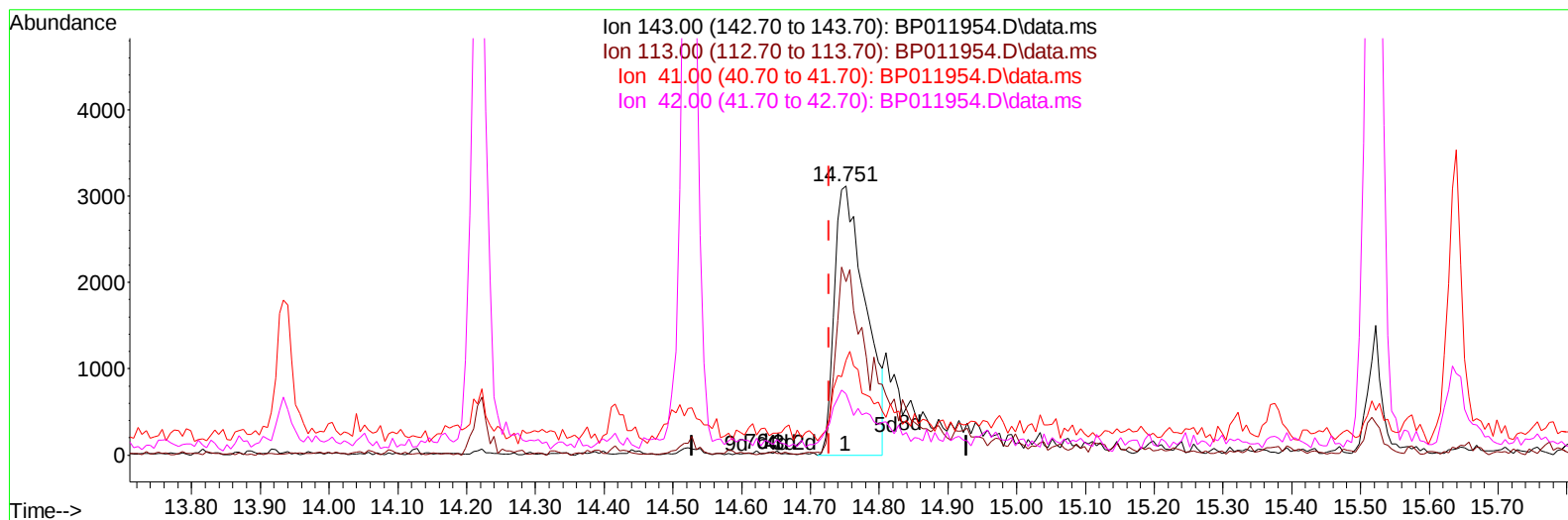
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100522\
Data File : BP011954.D
Acq On : 06 Oct 2022 05:22
Operator : CG/JU
Sample : N4956-02
Misc :
ALS Vial : 34 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
CB8R6

Manual IntegrationsAPPROVED

Quant Time: Oct 06 05:51:49 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP093022.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Oct 05 23:18:55 2022
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 10/06/2022
Supervised By :mohammad ahmed 10/10/2022



TIC: BP011954.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.751min (+ 0.024) 1.01 ng/ul

response 9741

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	59.80	64.52
41.00	24.20	35.26#
42.00	16.40	22.86#

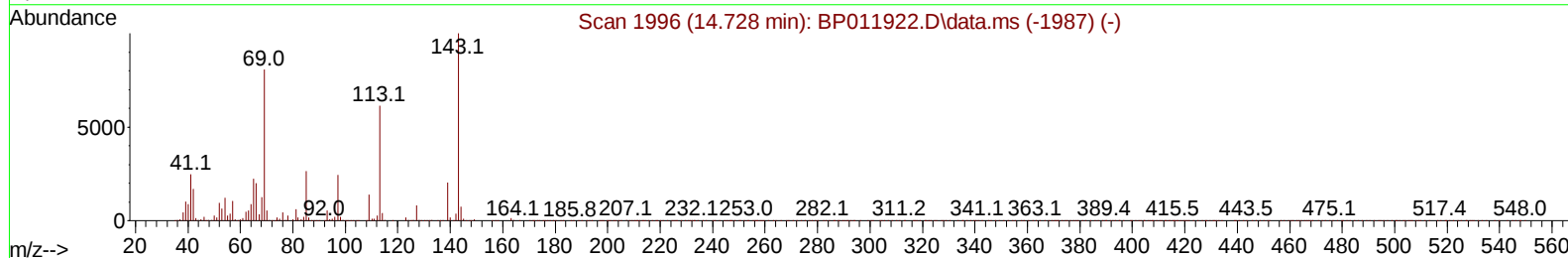
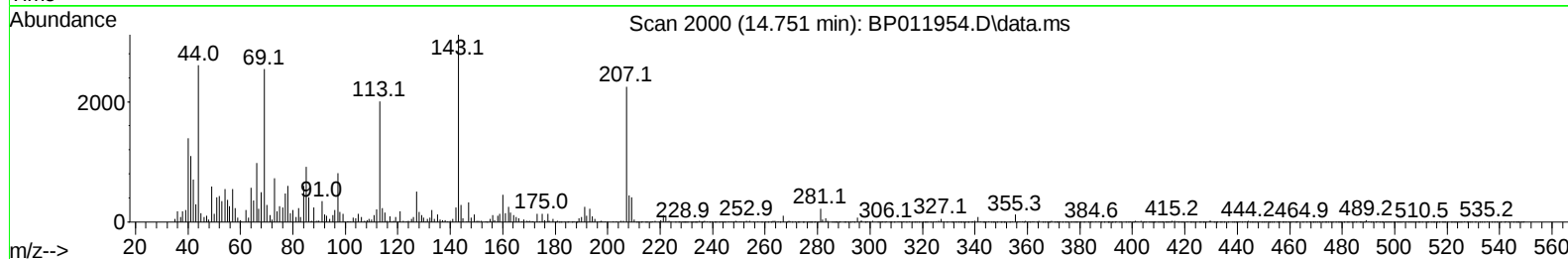
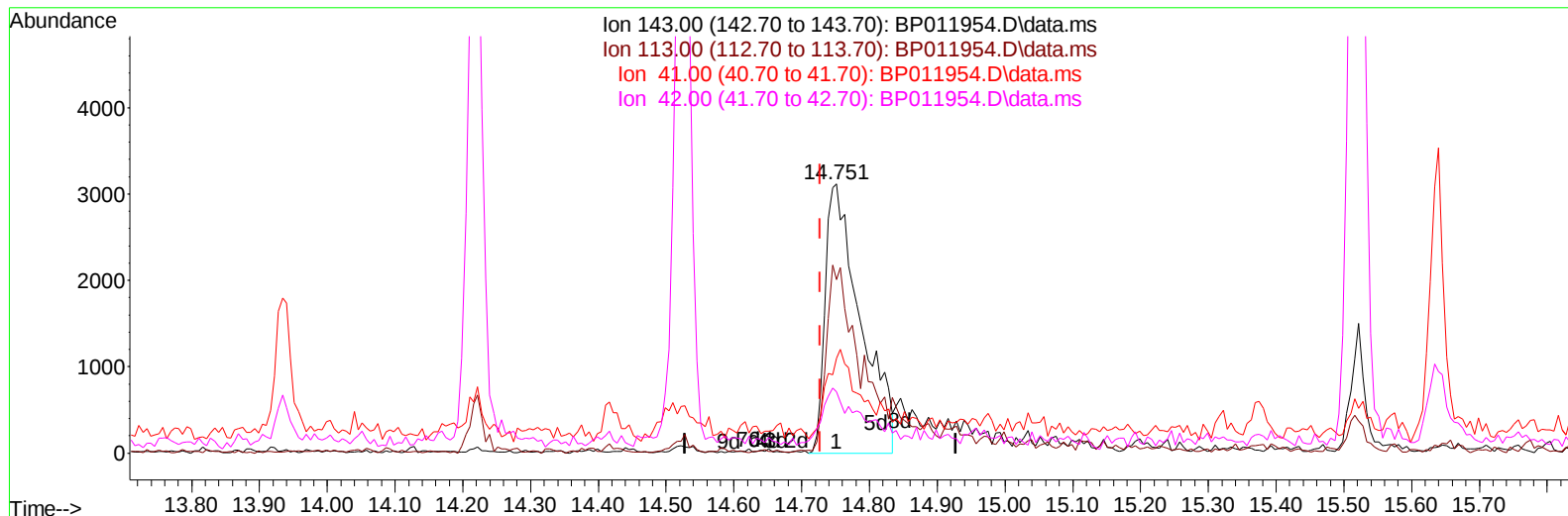
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TIC: BP011954.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.751min (+ 0.024) 1.17 ng/ul m

response 11222

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	59.80	64.52
41.00	24.20	35.26#
42.00	16.40	22.86#

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.881	152	250402	20.000	ng/ul	0.00
20) Naphthalene-d8	10.687	136	1050338	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.528	164	648280	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.281	188	1389570	20.000	ng/ul	0.00
79) Chrysene-d12	21.369	240	1307558	20.000	ng/ul	0.00
88) Perylene-d12	23.798	264	1317409	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.299	96	21281	3.748	ng/uL	0.00
4) Pyridine-d5	3.728	84	125613	7.553	ng/ul	0.00
7) Phenol-d5	7.046	99	55189	2.553	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.211	67	289696	24.177	ng/ul	0.00
11) 2-Chlorophenol-d4	7.411	132	176041	10.316	ng/ul	0.00
15) 4-Methylphenol-d8	8.593	113	122424	6.851	ng/ul	0.00
21) Nitrobenzene-d5	9.046	128	193997	24.249	ng/ul	0.00
24) 2-Nitrophenol-d4	9.769	143	102161	11.935	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.310	165	186090	11.754	ng/ul	0.00
31) 4-Chloroaniline-d4	10.828	131	451753	18.717	ng/ul	0.00
46) Dimethylphthalate-d6	13.934	166	1190621	25.683	ng/ul	0.00
49) Acenaphthylene-d8	14.216	160	1351892	25.580	ng/ul	0.00
54) 4-Nitrophenol-d4	14.751	143	11222m	1.169	ng/ul	0.02
60) Fluorene-d10	15.522	176	1110010	27.571	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.640	200	66749	7.936	ng/ul	0.00
73) Anthracene-d10	17.381	188	1808733	28.575	ng/ul	0.00
81) Pyrene-d10	19.616	212	2109326	31.853	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.639	264	1959139	29.707	ng/ul	0.00

Target Compounds Qvalue

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

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