

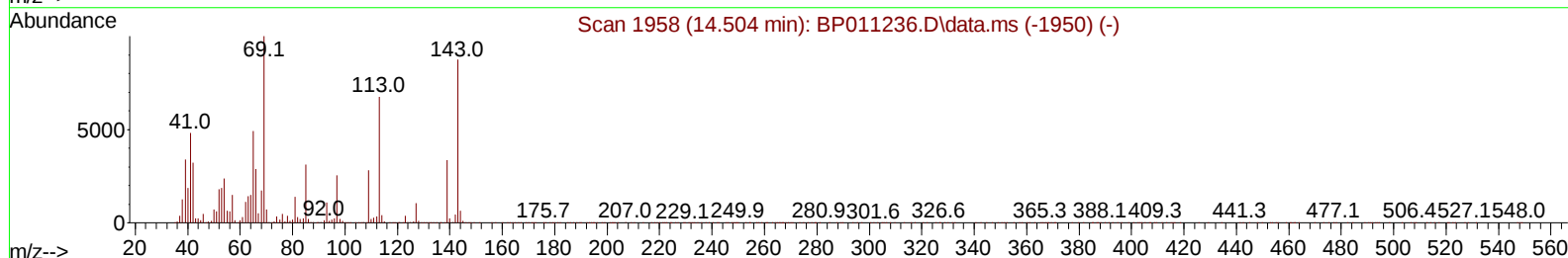
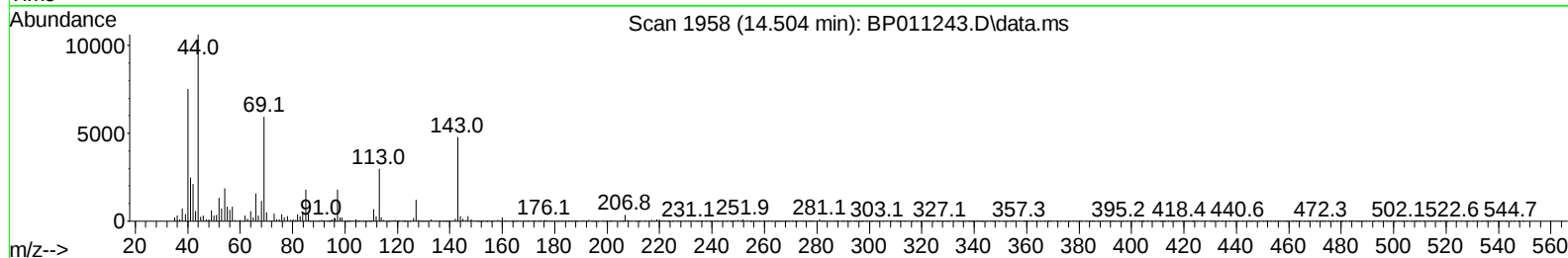
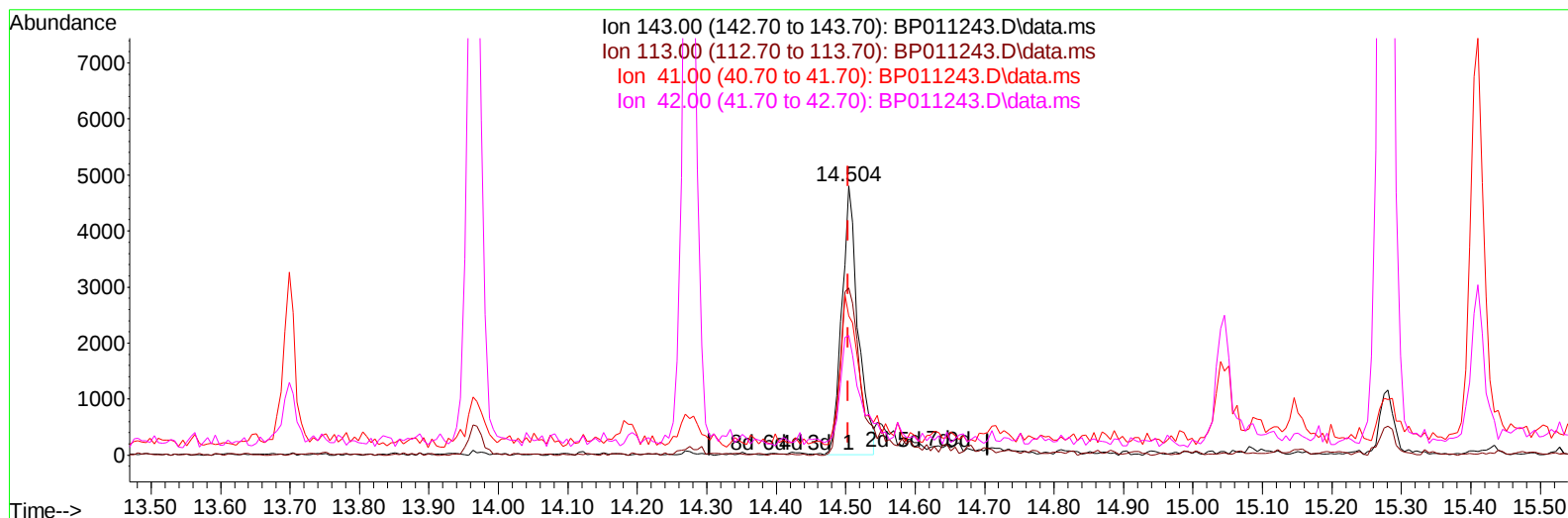
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080322\
 Data File : BP011243.D
 Acq On : 03 Aug 2022 14:04
 Operator : CG/JU
 Sample : N3900-17
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0CB8

Manual IntegrationsAPPROVED

Quant Time: Aug 03 22:59:55 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP080222.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Aug 02 23:52:37 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 08/04/2022
 Supervised By :Sohil Jodhani 08/10/2022



TIC: BP011243.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.504min (+ 0.000) 1.91 ng/ul

response 7892

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	75.70	62.09
41.00	54.40	51.82
42.00	35.50	44.41#

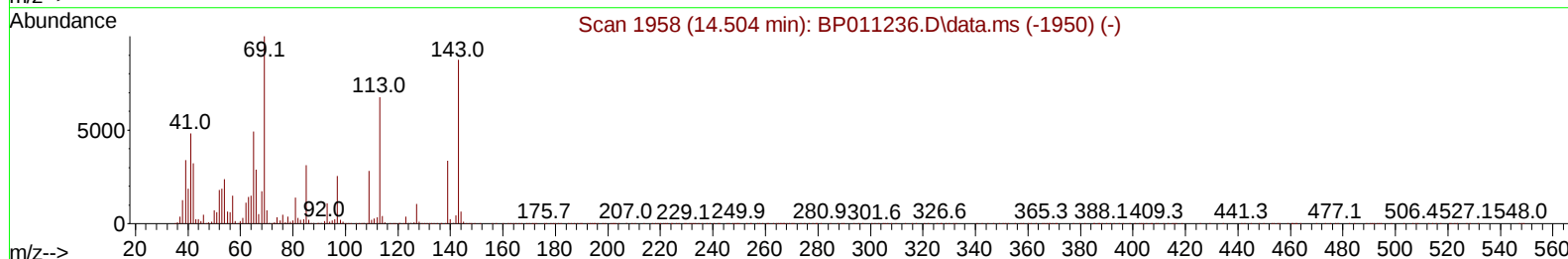
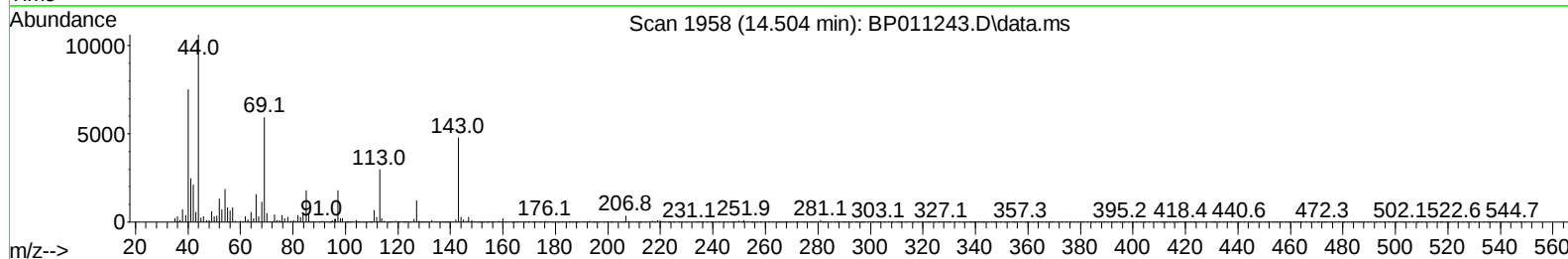
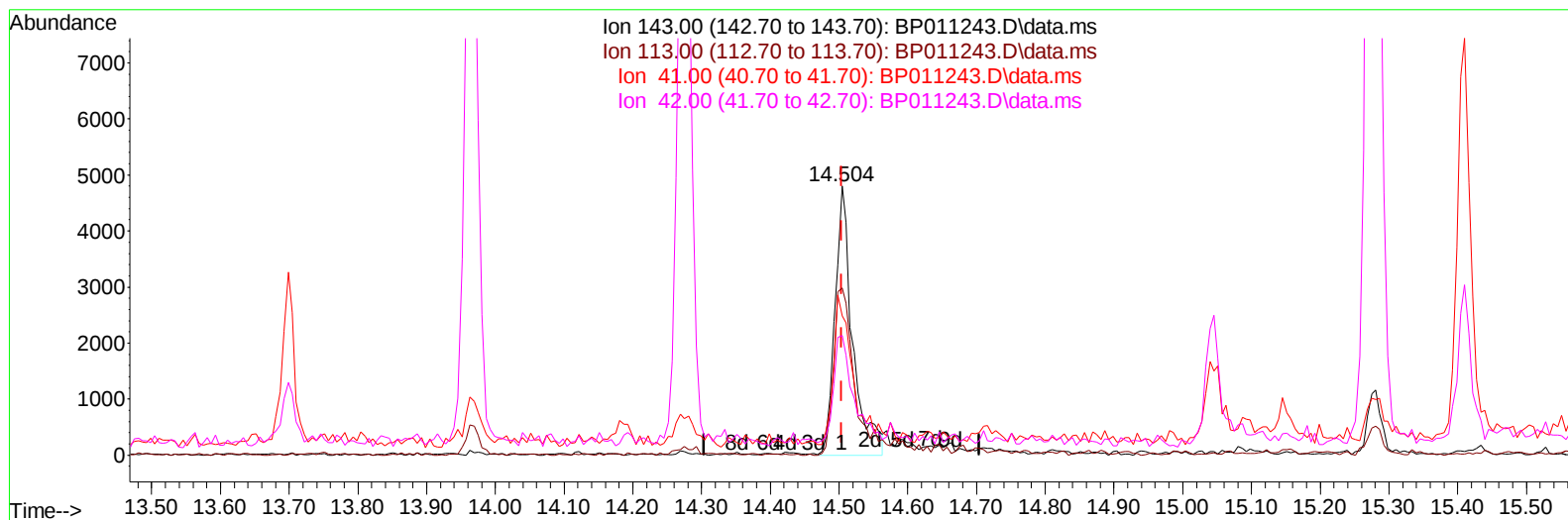
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(54) 4-Nitrophenol-d4 (S)

14.504min (+ 0.000) 2.09 ng/ul m

response 8622

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	75.70	62.09
41.00	54.40	51.82
42.00	35.50	44.41#

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.599	152	146173	20.000	ng/ul	0.00
20) Naphthalene-d8	10.393	136	643722	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.275	164	353876	20.000	ng/ul	# 0.00
64) Phenanthrene-d10	17.045	188	755647	20.000	ng/ul	0.00
79) Chrysene-d12	21.175	240	839106	20.000	ng/ul	0.00
88) Perylene-d12	23.492	264	917753	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.093	96	6024	3.332	ng/uL	0.00
4) Pyridine-d5	3.517	84	15295	3.137	ng/ul	0.01
7) Phenol-d5	6.787	99	52724	4.328	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.946	67	247167	28.119	ng/ul	0.00
11) 2-Chlorophenol-d4	7.134	132	180051	18.134	ng/ul	0.00
15) 4-Methylphenol-d8	8.322	113	106231	10.598	ng/ul	0.00
21) Nitrobenzene-d5	8.763	128	137036	29.596	ng/ul	0.00
24) 2-Nitrophenol-d4	9.487	143	93780	21.608	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.022	165	184007	22.233	ng/ul	0.00
31) 4-Chloroaniline-d4	10.546	131	217022	16.950	ng/ul	0.00
46) Dimethylphthalate-d6	13.698	166	766902	31.642	ng/ul	0.00
49) Acenaphthylene-d8	13.963	160	943401	30.937	ng/ul	0.00
54) 4-Nitrophenol-d4	14.504	143	8622m	2.092	ng/ul	0.00
60) Fluorene-d10	15.281	176	758334	36.771	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.410	200	64488	18.379	ng/ul	0.00
73) Anthracene-d10	17.145	188	1286735	38.745	ng/ul	0.00
81) Pyrene-d10	19.404	212	1552663	36.597	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.345	264	1648004	36.035	ng/ul	0.00

Target Compounds Qvalue

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

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