

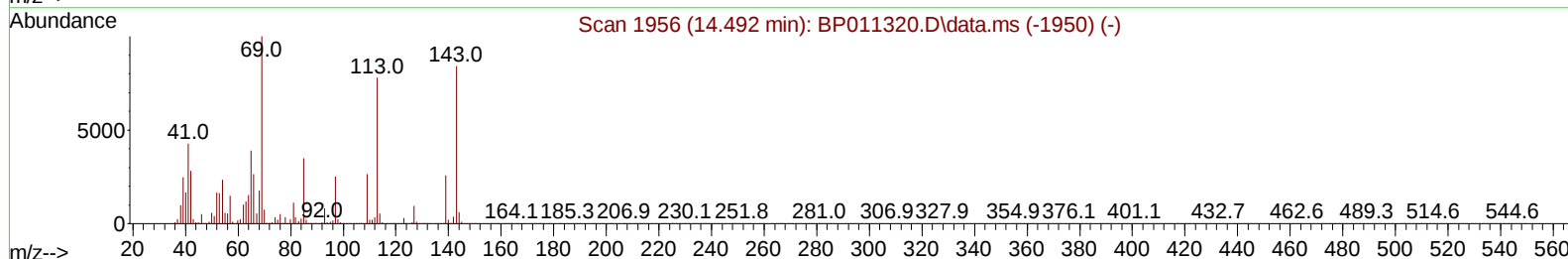
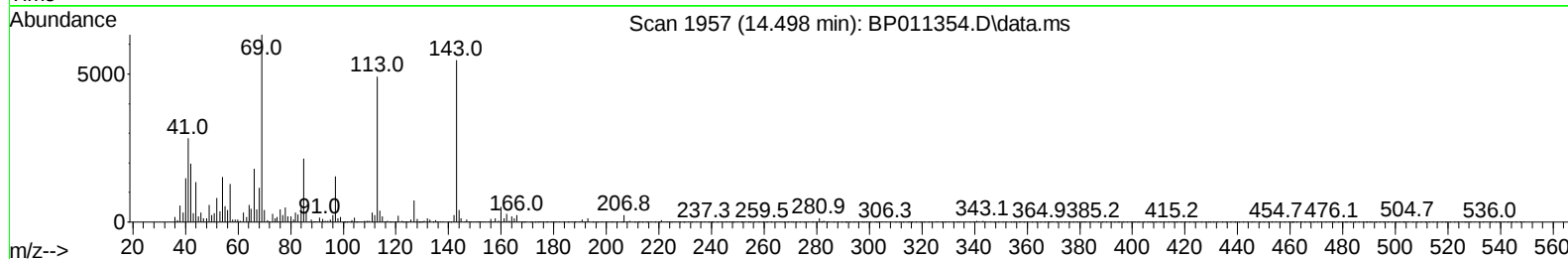
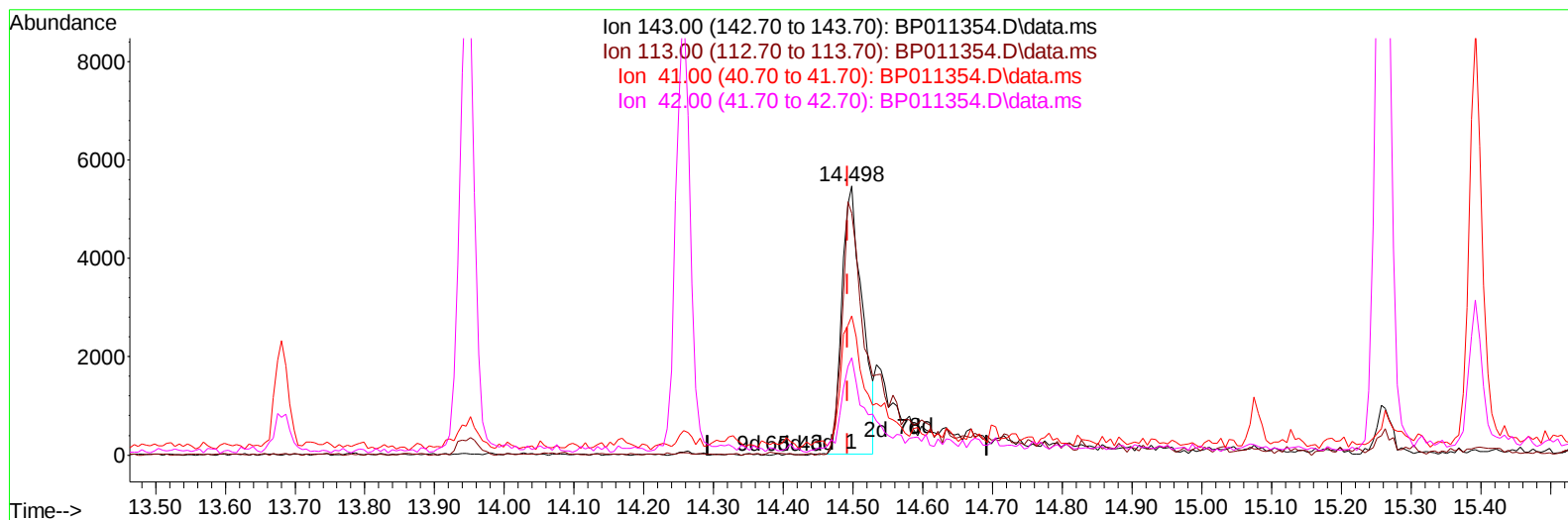
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080922\
 Data File : BP011354.D
 Acq On : 10 Aug 2022 05:59
 Operator : CG/JU
 Sample : N4069-10
 Misc :
 ALS Vial : 36 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 GBJL4

Manual IntegrationsAPPROVED

Quant Time: Aug 10 06:40:42 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP080822.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Aug 08 18:47:19 2022
 Response via : Initial Calibration

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



TIC: BP011354.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.498min (+ 0.006) 3.34 ng/u1

response 10955

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	75.70	90.03
41.00	54.40	51.69
42.00	35.50	36.16

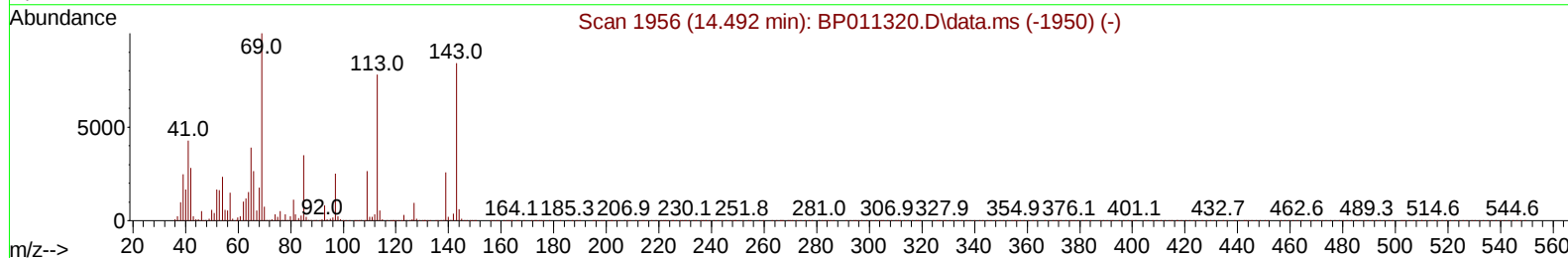
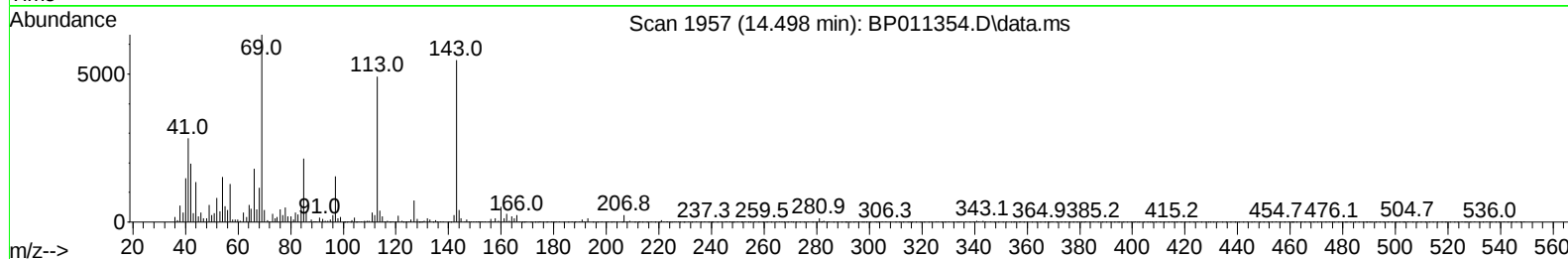
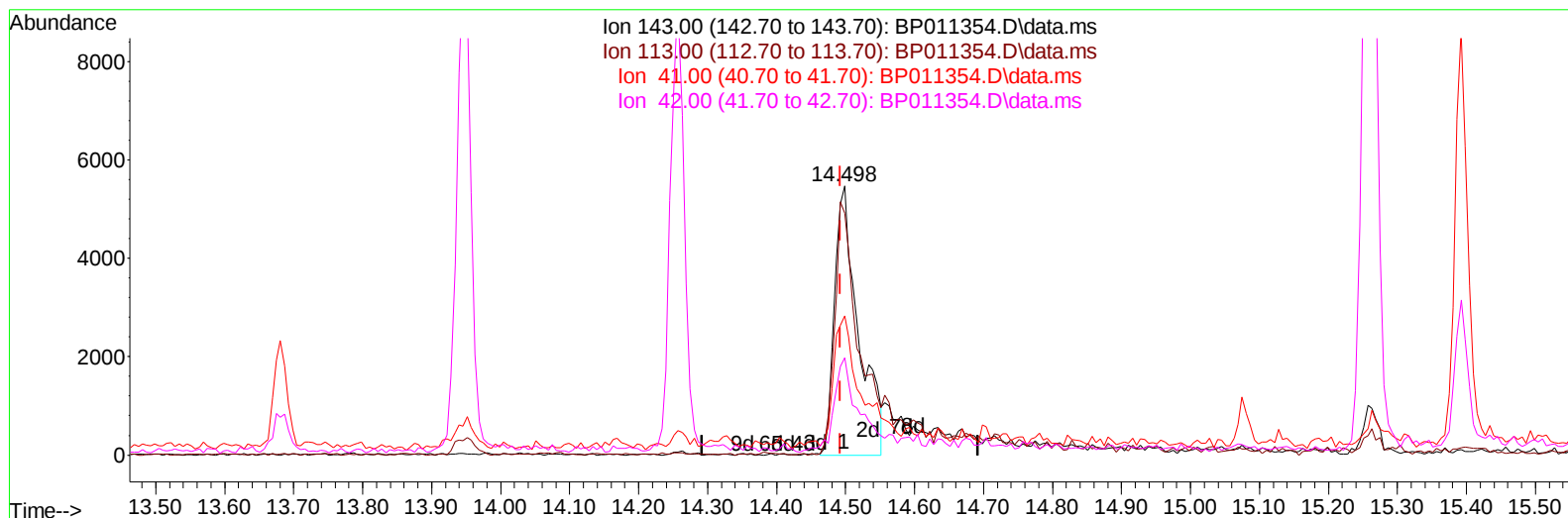
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TIC: BP011354.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.498min (+ 0.006) 4.01 ng/ul m

response 13151

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	75.70	90.03
41.00	54.40	51.69
42.00	35.50	36.16

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.587	152	107974	20.000	ng/ul	0.00
20) Naphthalene-d8	10.375	136	450833	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.257	164	258892	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.028	188	522103	20.000	ng/ul	-0.01
79) Chrysene-d12	21.157	240	480603	20.000	ng/ul	-0.01
88) Perylene-d12	23.463	264	487159	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.087	96	15660	4.446	ng/uL	0.00
4) Pyridine-d5	3.505	84	85879	9.666	ng/ul	0.00
7) Phenol-d5	6.769	99	67950	7.111	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.934	67	207261	30.591	ng/ul	0.00
11) 2-Chlorophenol-d4	7.116	132	194586	25.954	ng/ul	0.00
15) 4-Methylphenol-d8	8.305	113	122454	16.321	ng/ul	-0.01
21) Nitrobenzene-d5	8.752	128	115304	37.580	ng/ul	0.00
24) 2-Nitrophenol-d4	9.469	143	112556	40.068	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.004	165	170362	30.031	ng/ul	0.00
31) 4-Chloroaniline-d4	10.528	131	234193	23.613	ng/ul	0.00
46) Dimethylphthalate-d6	13.681	166	669224	37.255	ng/ul	0.00
49) Acenaphthylene-d8	13.945	160	765292	35.392	ng/ul	-0.01
54) 4-Nitrophenol-d4	14.498	143	13151m	4.008	ng/ul	0.00
60) Fluorene-d10	15.263	176	552231	35.948	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.392	200	87044	35.893	ng/ul	0.00
73) Anthracene-d10	17.128	188	946796	40.345	ng/ul	-0.01
81) Pyrene-d10	19.392	212	1066487	40.278	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.316	264	1002601	40.810	ng/ul	-0.01

Target Compounds Qvalue

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

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