

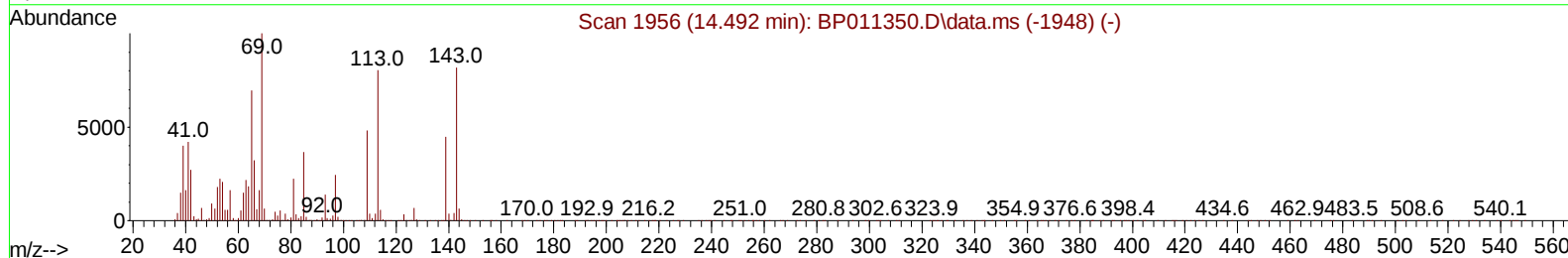
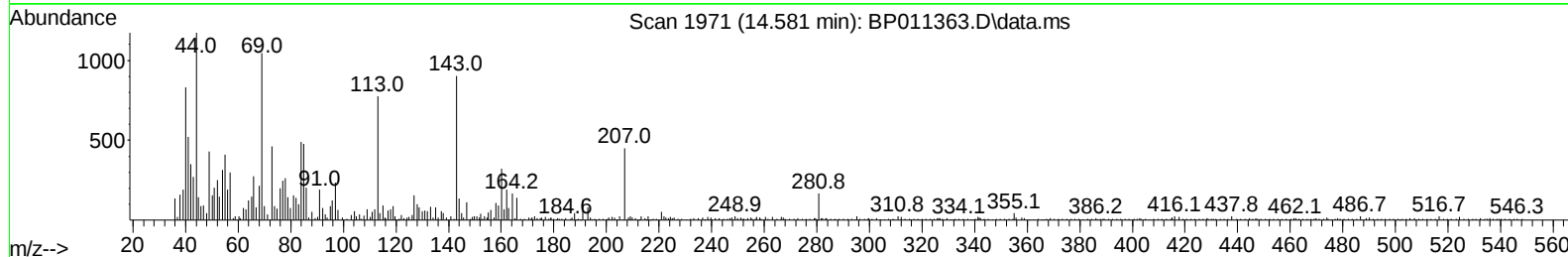
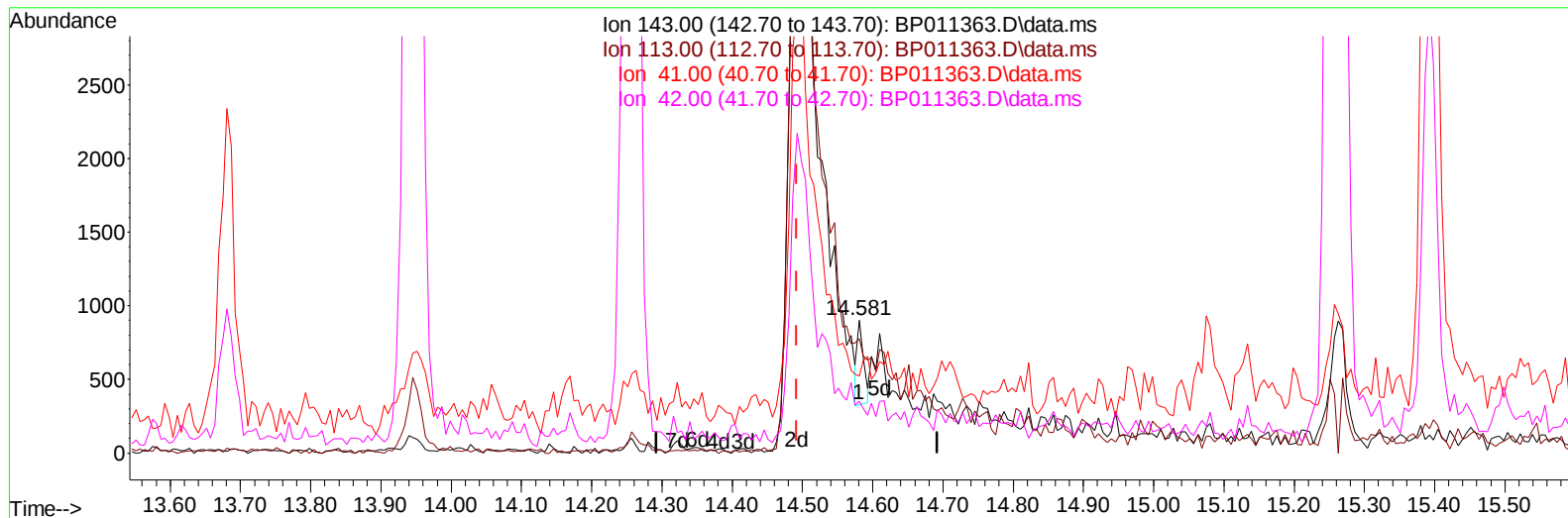
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080922\
 Data File : BP011363.D
 Acq On : 10 Aug 2022 11:09
 Operator : CG/JU
 Sample : N4069-14
 Misc :
 ALS Vial : 45 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 GBJL8

Manual IntegrationsAPPROVED

Quant Time: Aug 10 23:56:32 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 08/11/2022
 Supervised By :Sohil Jodhani 08/12/2022



TIC: BP011363.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.581min (+ 0.088) 0.11 ng/u1

response 353

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	75.70	85.94
41.00	54.40	57.70
42.00	35.50	38.98

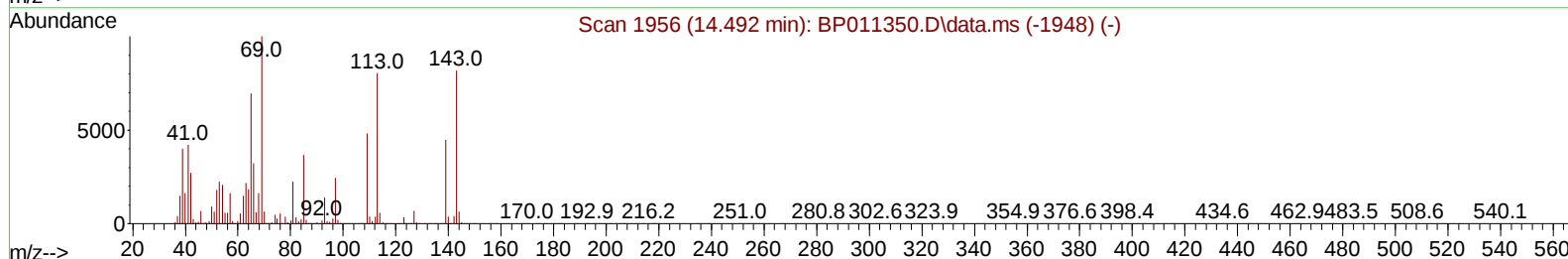
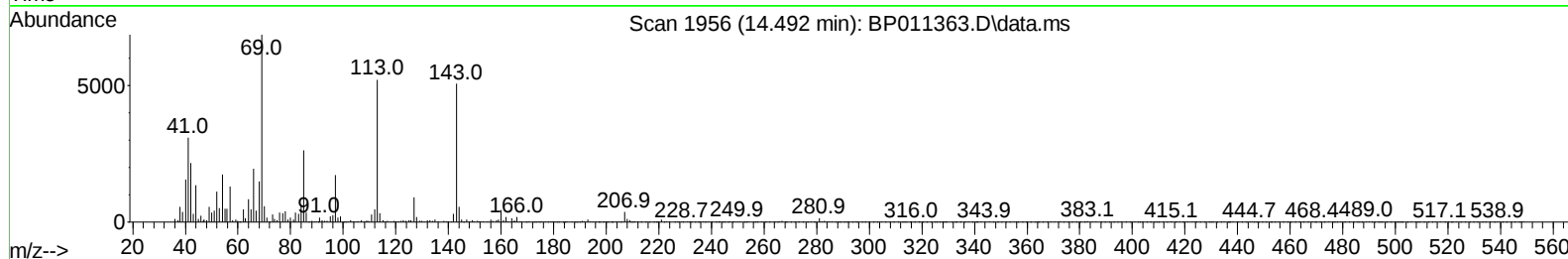
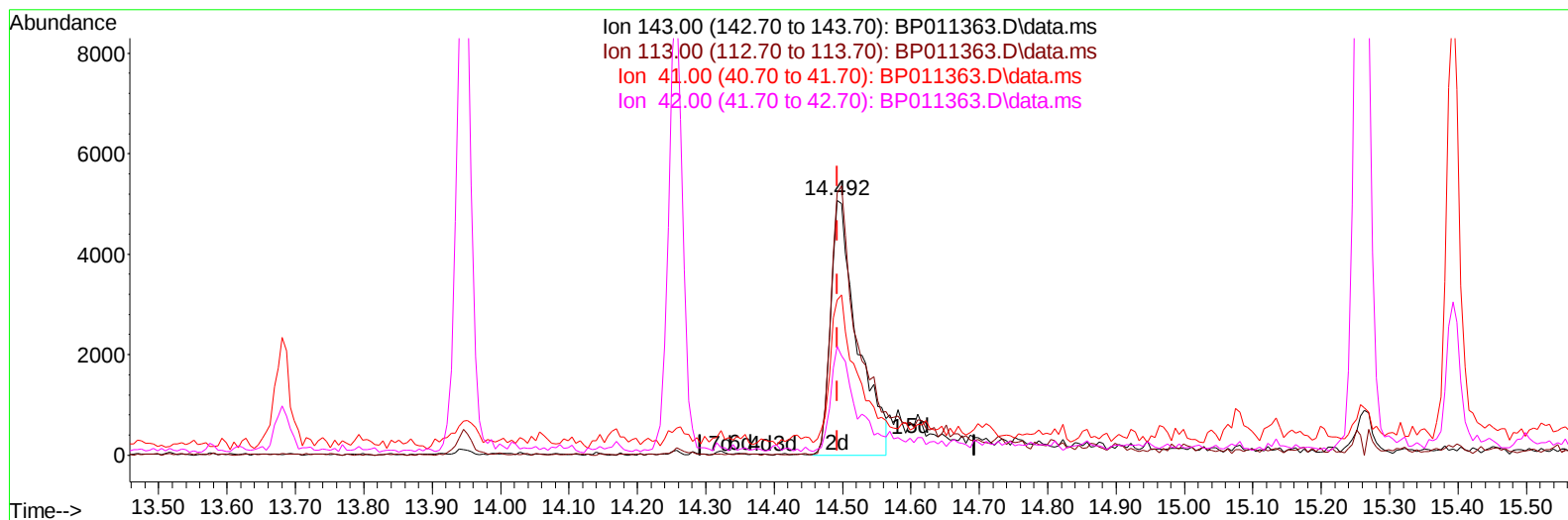
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(54) 4-Nitrophenol-d4 (S)

14.492min (-0.000) 4.16 ng/ul m

response 13637

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	75.70	102.90#
41.00	54.40	60.98
42.00	35.50	42.81#

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.587	152	105824	20.000	ng/ul	0.00
20) Naphthalene-d8	10.375	136	453513	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.257	164	258510	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.028	188	508829	20.000	ng/ul	-0.01
79) Chrysene-d12	21.157	240	465010	20.000	ng/ul	-0.01
88) Perylene-d12	23.468	264	461073	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.087	96	16239	4.704	ng/uL	0.00
4) Pyridine-d5	3.505	84	55867	6.416	ng/ul	0.00
7) Phenol-d5	6.769	99	67705	7.229	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.934	67	205850	31.000	ng/ul	0.00
11) 2-Chlorophenol-d4	7.116	132	186886	25.433	ng/ul	0.00
15) 4-Methylphenol-d8	8.305	113	121081	16.466	ng/ul	-0.01
21) Nitrobenzene-d5	8.752	128	117378	38.030	ng/ul	0.00
24) 2-Nitrophenol-d4	9.469	143	112808	39.920	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.004	165	170574	29.890	ng/ul	0.00
31) 4-Chloroaniline-d4	10.528	131	246135	24.671	ng/ul	0.00
46) Dimethylphthalate-d6	13.681	166	685410	38.213	ng/ul	0.00
49) Acenaphthylene-d8	13.945	160	791546	36.660	ng/ul	-0.01
54) 4-Nitrophenol-d4	14.492	143	13637m	4.163	ng/ul	0.00
60) Fluorene-d10	15.263	176	570558	37.195	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.392	200	86504	36.601	ng/ul	0.00
73) Anthracene-d10	17.128	188	982131	42.943	ng/ul	-0.01
81) Pyrene-d10	19.392	212	1101911	43.011	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.321	264	1033396	44.443	ng/ul	0.00

Target Compounds Qvalue

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

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