

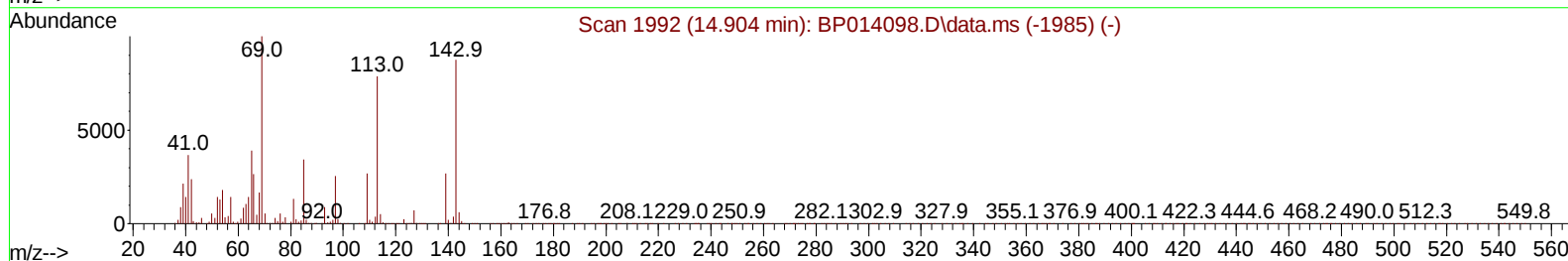
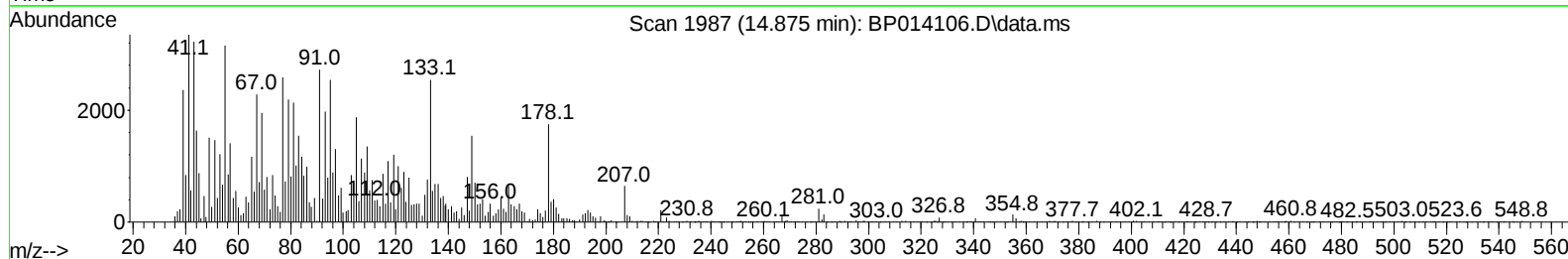
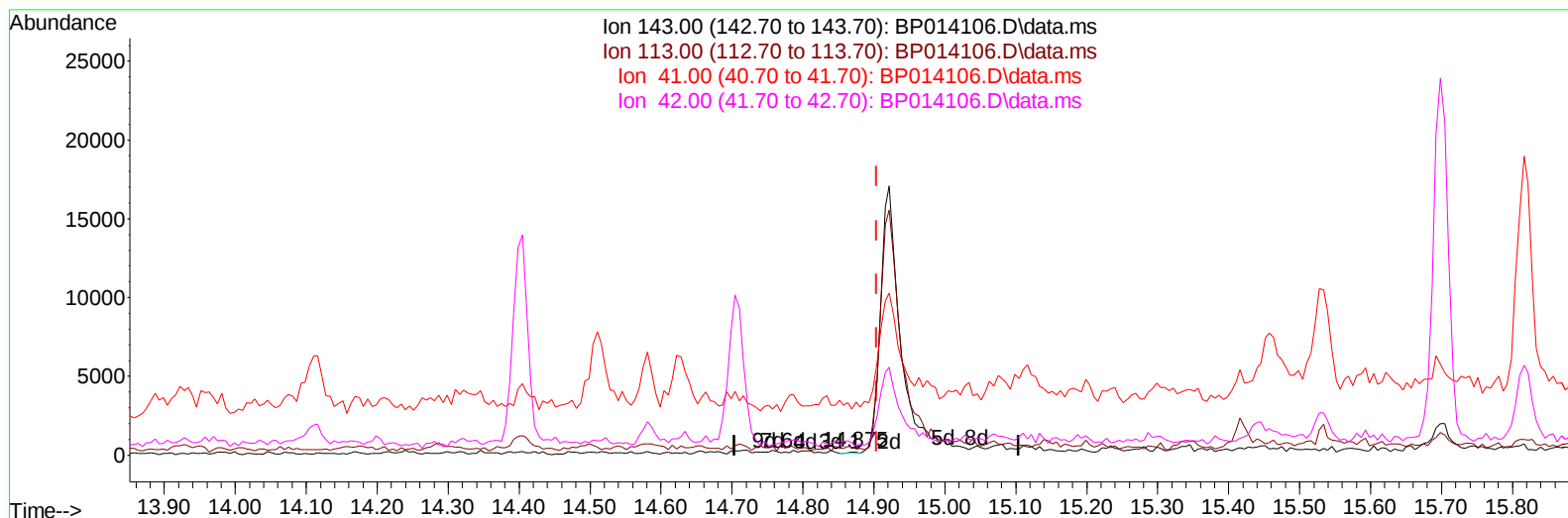
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP031623\
 Data File : BP014106.D
 Acq On : 16 Mar 2023 17:03
 Operator : CG/JU
 Sample : 01884-10
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 CB919

Manual IntegrationsAPPROVED

Quant Time: Mar 17 01:00:40 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP030723.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Mar 17 00:57:25 2023
 Response via : Initial Calibration

Reviewed By :Christian Giraldo 03/17/2023
 Supervised By :Jagrut Upadhyay 03/17/2023



TIC: BP014106.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.875min (-0.029) 0.01 ng/u1

response 80

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	103.10	213.23#
41.00	40.20	1771.43#
42.00	28.50	298.94#

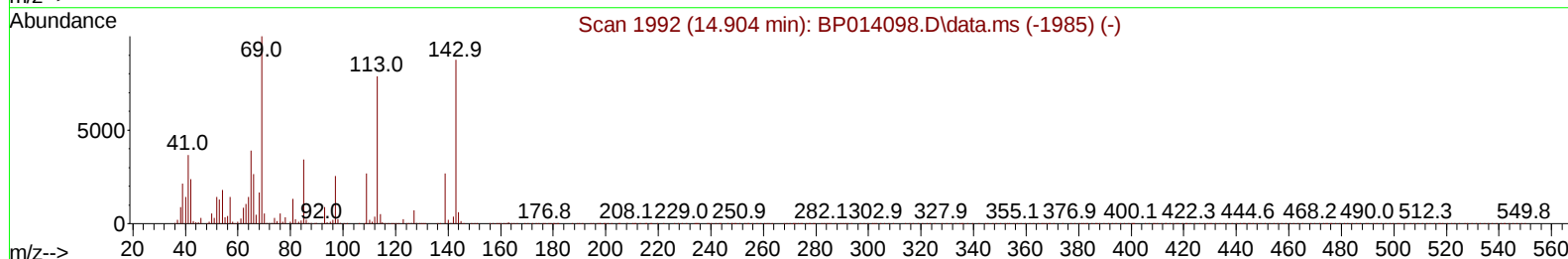
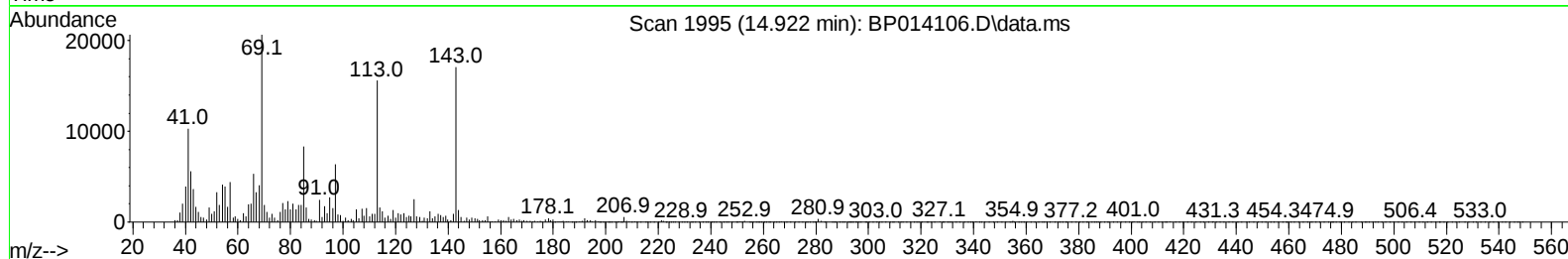
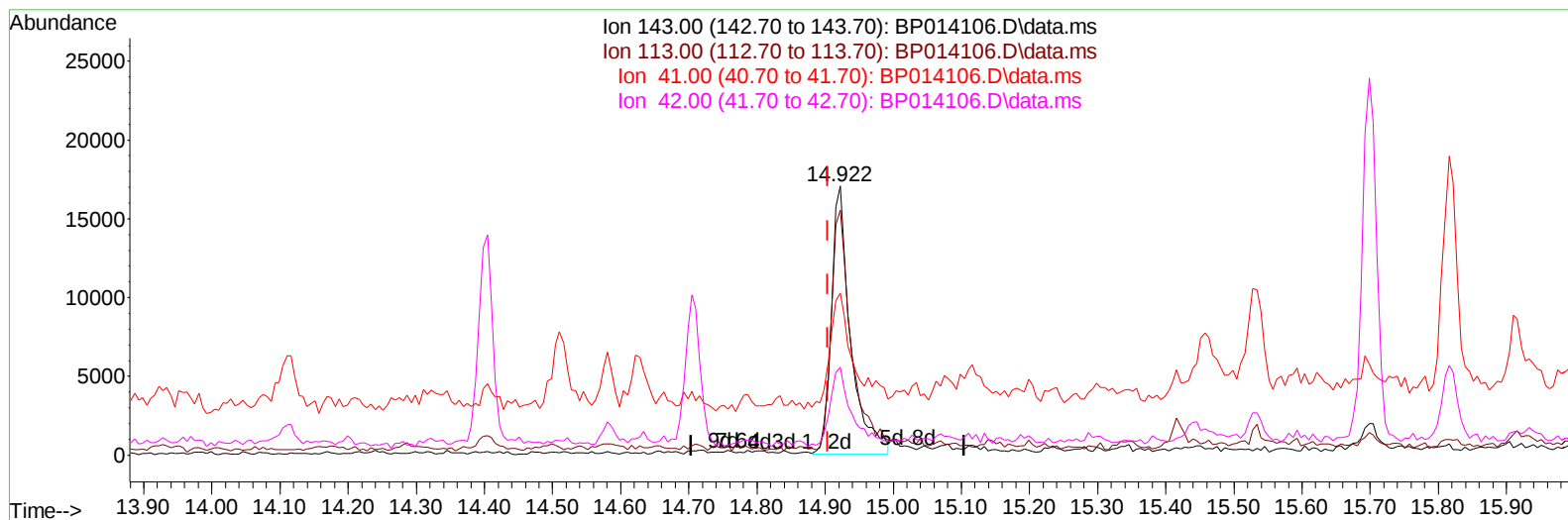
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(54) 4-Nitrophenol-d4 (S)

14.922min (+ 0.018) 5.34 ng/ul m

response 33284

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	103.10	91.24
41.00	40.20	60.20#
42.00	28.50	32.78

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.063	152	147166	20.000	ng/ul	0.00
20) Naphthalene-d8	10.887	136	676197	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.704	164	416676	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.463	188	1140665	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.545	240	1093090	20.000	ng/ul	# 0.00
88) Perylene-d12	24.086	264	999242	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.422	96	23212	5.959	ng/uL	0.00
4) Pyridine-d5	3.870	84	104077	9.651	ng/ul	0.01
7) Phenol-d5	7.216	99	90764	6.989	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.387	67	289685	38.247	ng/ul	0.00
11) 2-Chlorophenol-d4	7.593	132	250773	25.345	ng/ul	0.00
15) 4-Methylphenol-d8	8.775	113	191113	18.015	ng/ul	0.00
21) Nitrobenzene-d5	9.240	128	179987	33.360	ng/ul	0.00
24) 2-Nitrophenol-d4	9.963	143	181951	28.172	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.504	165	365060	30.040	ng/ul	0.00
31) 4-Chloroaniline-d4	11.028	131	256499	15.949	ng/ul	0.00
46) Dimethylphthalate-d6	14.110	166	1353662	38.280	ng/ul	0.00
49) Acenaphthylene-d8	14.404	160	1451744	38.754	ng/ul	0.00
54) 4-Nitrophenol-d4	14.922	143	33284m	5.341	ng/ul	0.02
60) Fluorene-d10	15.698	176	1366276	45.589	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.816	200	256500	29.752	ng/ul	0.00
73) Anthracene-d10	17.563	188	2167855	39.149	ng/ul	0.00
81) Pyrene-d10	19.792	212	2571161	42.648	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.921	264	2134001	40.149	ng/ul	0.00

Target Compounds Qvalue

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

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