

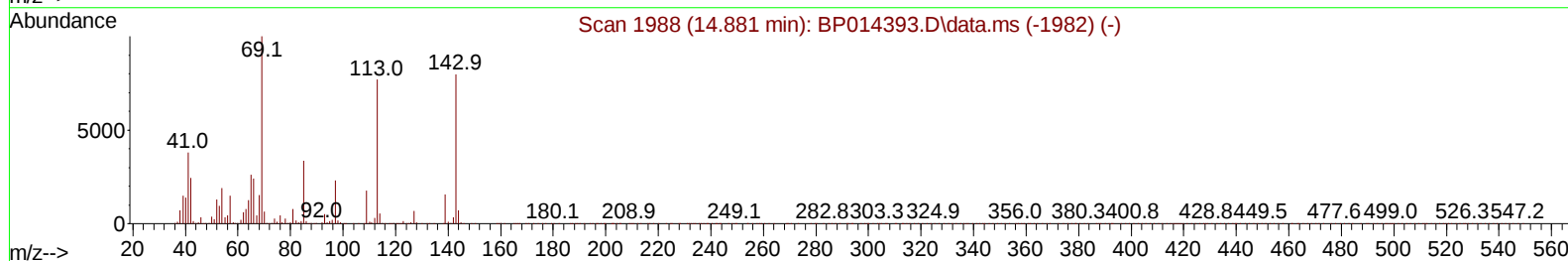
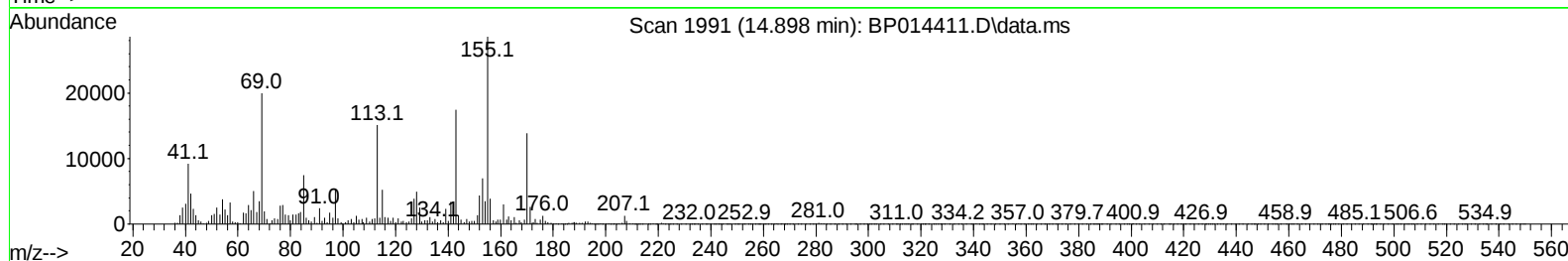
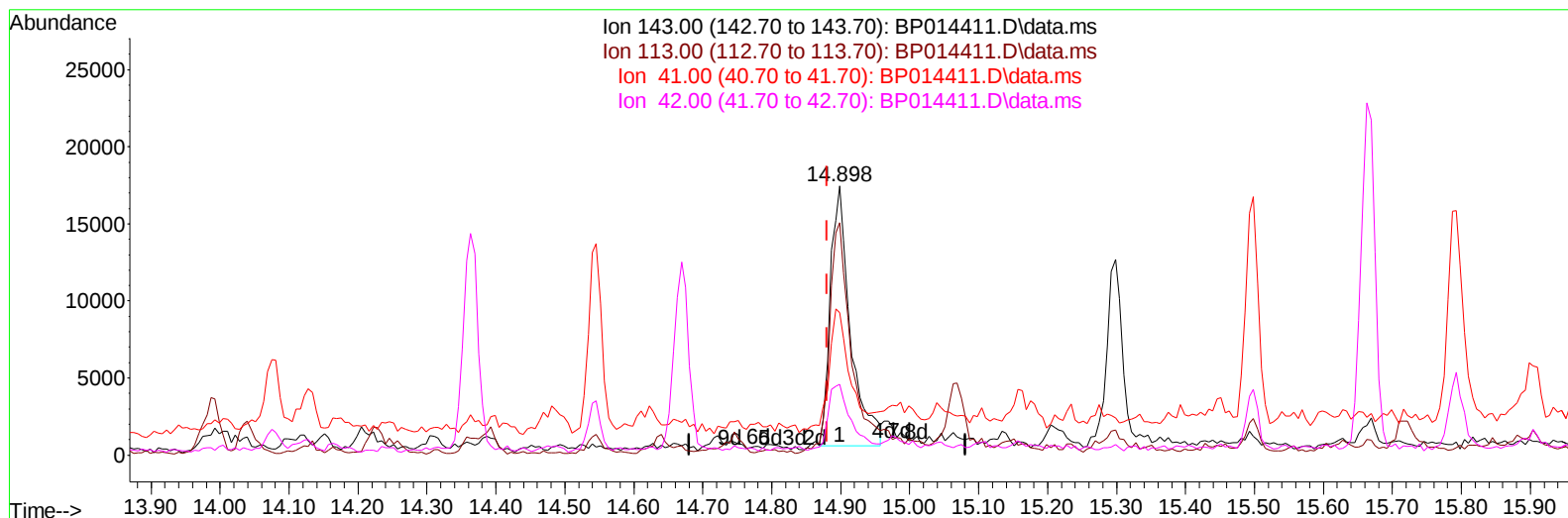
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP033023\
 Data File : BP014411.D
 Acq On : 30 Mar 2023 20:22
 Operator : CG/JU
 Sample : 02131-01
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0QG6

Manual IntegrationsAPPROVED

Quant Time: Mar 30 22:43:57 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP032823.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Mar 30 12:32:14 2023
 Response via : Initial Calibration

Reviewed By :Christian Giraldo 03/31/2023
 Supervised By :Jagrut Upadhyay 03/31/2023



TIC: BP014411.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.898min (+ 0.018) 5.32 ng/u1

response 34232

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	96.80	86.49
41.00	44.20	52.98
42.00	31.50	26.42

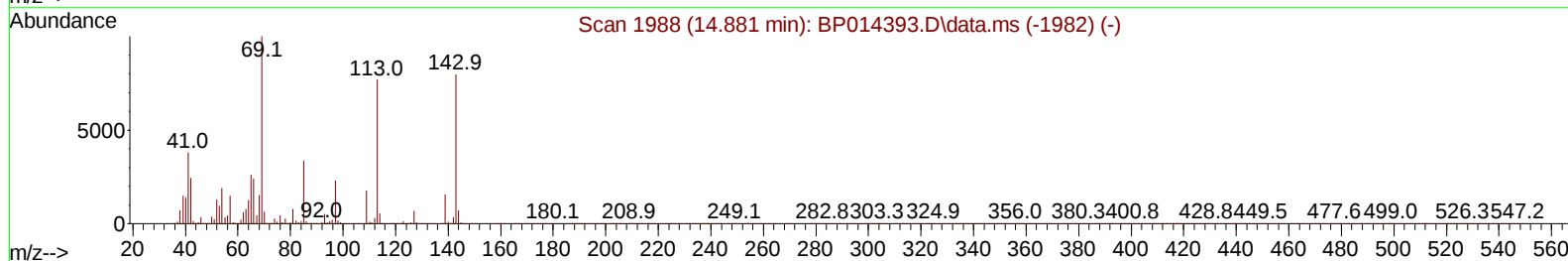
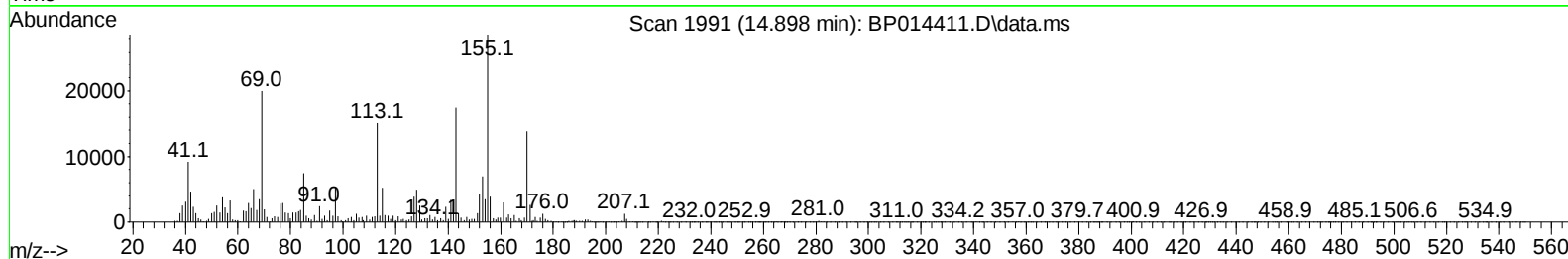
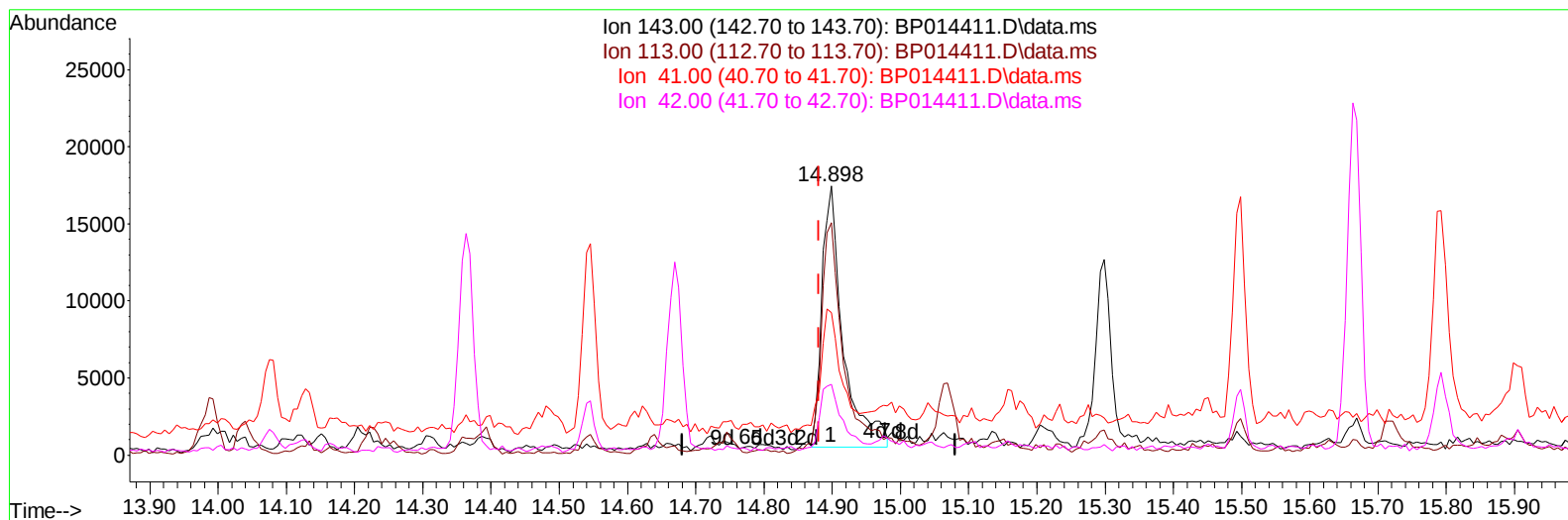
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TIC: BP014411.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.898min (+ 0.018) 5.70 ng/ul m

response 36673

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	96.80	86.49
41.00	44.20	52.98
42.00	31.50	26.42

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.016	152	173446	20.000	ng/ul	0.00
20) Naphthalene-d8	10.840	136	775588	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.669	164	525015	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.440	188	1114143	20.000	ng/ul	0.00
79) Chrysene-d12	21.522	240	850998	20.000	ng/ul	0.00
88) Perylene-d12	24.039	264	797976	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.387	96	21014	4.698	ng/uL	0.00
4) Pyridine-d5	3.828	84	82035	7.157	ng/ul	0.00
7) Phenol-d5	7.175	99	101143	6.323	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.340	67	340702	33.197	ng/ul	0.00
11) 2-Chlorophenol-d4	7.540	132	285333	24.680	ng/ul	0.00
15) 4-Methylphenol-d8	8.728	113	198475	15.382	ng/ul	0.00
21) Nitrobenzene-d5	9.193	128	197676	31.572	ng/ul	0.00
24) 2-Nitrophenol-d4	9.916	143	220012	30.687	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.463	165	348092	26.659	ng/ul	0.00
31) 4-Chloroaniline-d4	10.987	131	194312	11.298	ng/ul	0.00
46) Dimethylphthalate-d6	14.075	166	1422905	33.426	ng/ul	0.00
49) Acenaphthylene-d8	14.363	160	1521445	32.210	ng/ul	0.00
54) 4-Nitrophenol-d4	14.898	143	36673m	5.702	ng/ul	0.02
60) Fluorene-d10	15.663	176	1169206	32.936	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.793	200	245835	29.987	ng/ul	0.00
73) Anthracene-d10	17.534	188	1790218	33.202	ng/ul	0.00
81) Pyrene-d10	19.763	212	1957235	34.241	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.874	264	1426225	33.641	ng/ul	0.00
Target Compounds						
					Qvalue	
30) Naphthalene	10.899	128	8045789	186.239	ng/ul	100
36) 2-Methylnaphthalene	12.493	142	912450	31.165	ng/ul	99
37) 1-Methylnaphthalene	12.716	142	760345	26.186	ng/ul	96
43) 1,1'-Biphenyl	13.499	154	264946	6.246	ng/ul	97
52) Acenaphthene	14.734	153	83504	2.349	ng/ul	96
56) Dibenzofuran	15.069	168	100809	2.017	ng/ul	89
58) 2,3,4,6-Tetrachlorophenol	15.298	232	667331	60.211	ng/ul	99
61) Fluorene	15.722	166	143478	3.542	ng/ul	98
71) Pentachlorophenol	17.110	266	10089491	1093.632	ng/ul	99
72) Phenanthrene	17.481	178	384108	6.197	ng/ul	99
86) Bis(2-ethylhexyl)phtha...	21.398	149	33968	1.022	ng/ul	98

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

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