

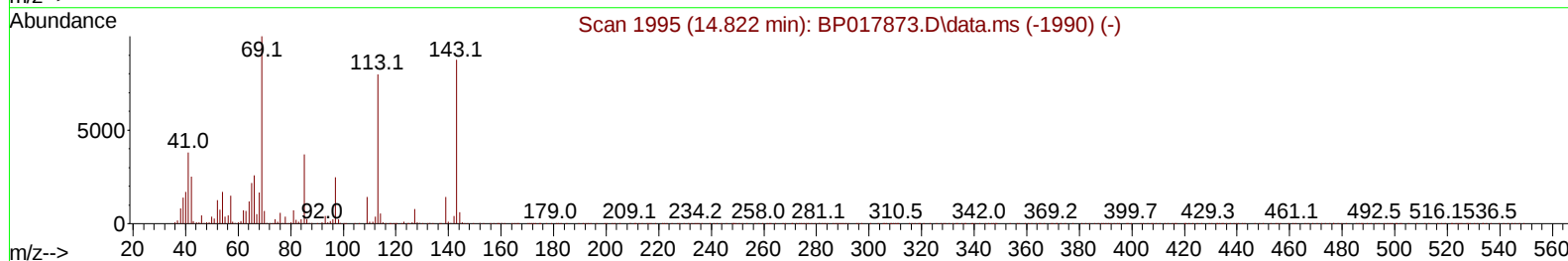
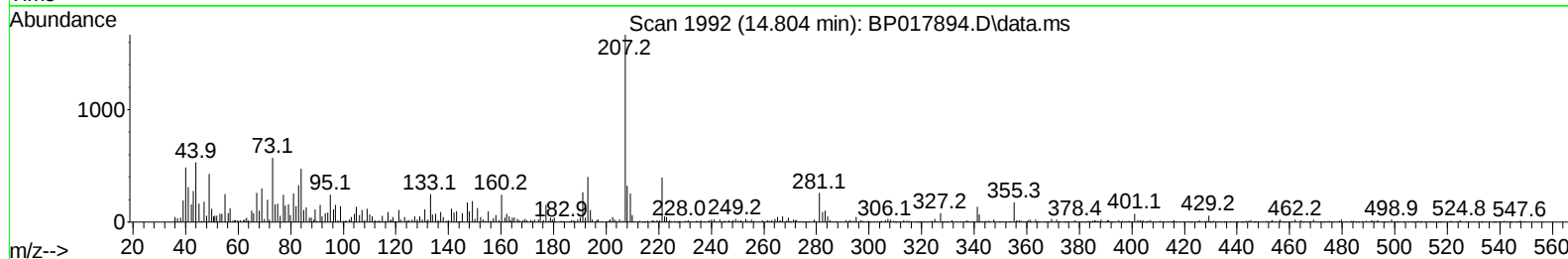
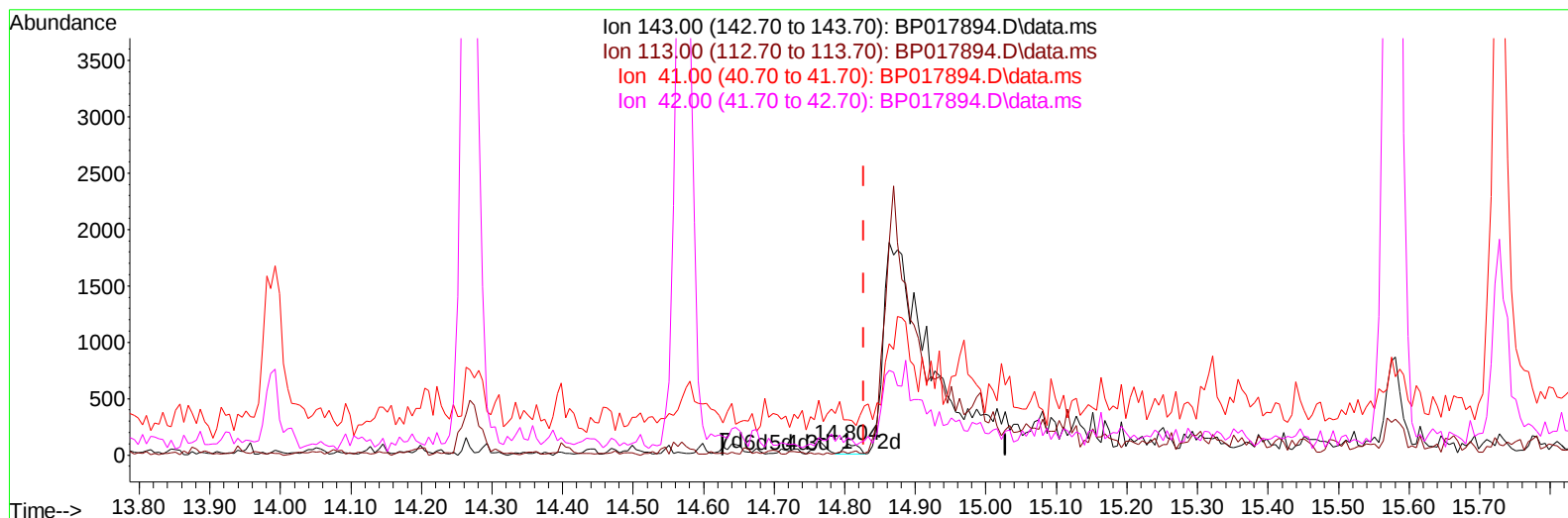
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110123\
 Data File : BP017894.D
 Acq On : 02 Nov 2023 19:28
 Operator : MA/JU
 Sample : 04996-24
 Misc :
 ALS Vial : 49 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 A49J7

Manual IntegrationsAPPROVED

Quant Time: Nov 02 23:39:01 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP101823.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Oct 19 01:32:15 2023
 Response via : Initial Calibration

Reviewed By :Yogesh Patel 11/03/2023
 Supervised By :mohammad ahmed 11/04/2023



TIC: BP017894.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.804min (-0.024) 0.04 ng/u1

response 127

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	81.70	16.49#
41.00	39.20	321.65#
42.00	25.60	161.86#

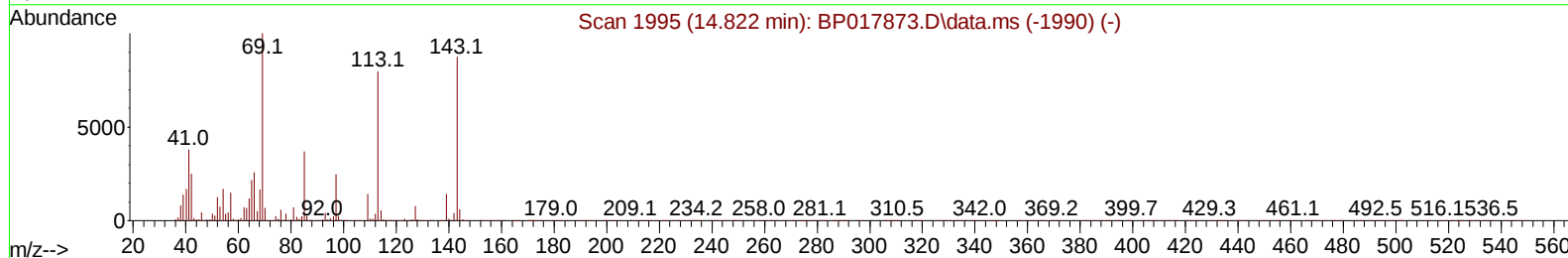
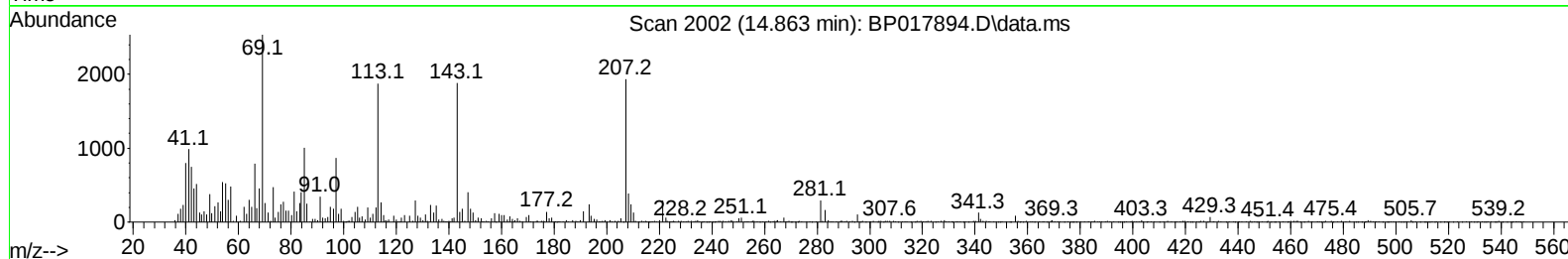
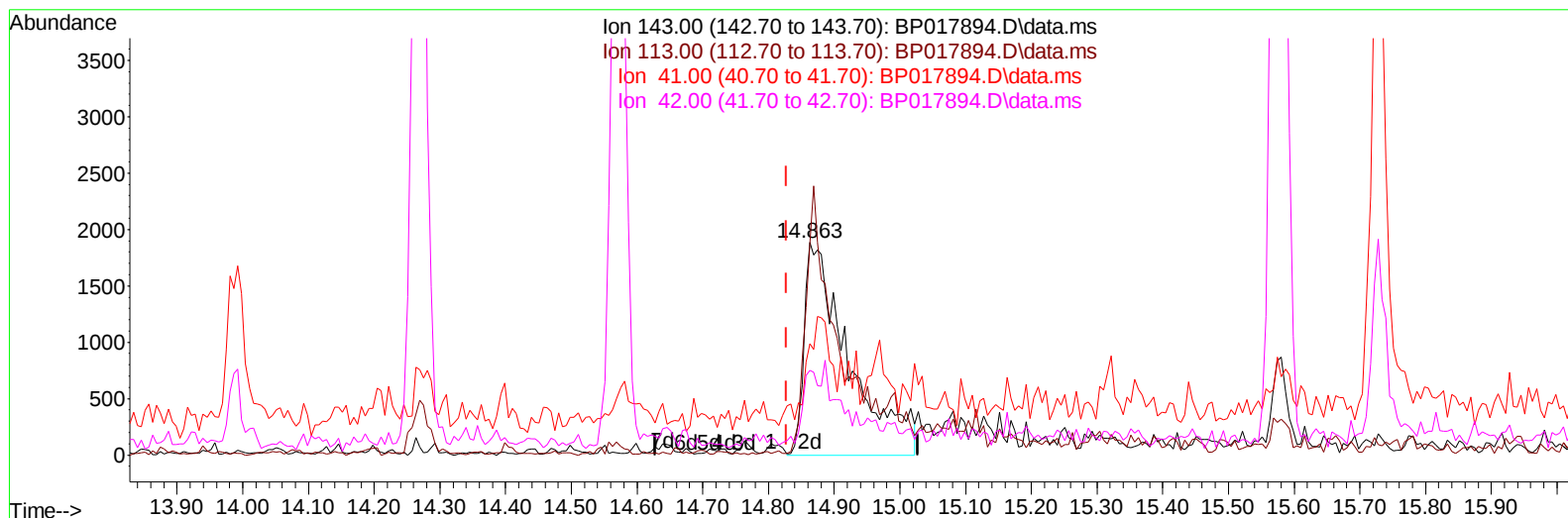
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(54) 4-Nitrophenol-d4 (S)

14.863min (+ 0.035) 2.94 ng/ul m

response 8862

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	81.70	99.58#
41.00	39.20	52.36#
42.00	25.60	39.89#

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Quant Time: Nov 03 00:21:35 2023
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.875	152	115239	20.000	ng/ul	-0.01
20) Naphthalene-d8	10.704	136	435501	20.000	ng/ul	-0.02
38) Acenaphthene-d10	14.575	164	252154	20.000	ng/ul	-0.01
64) Phenanthrene-d10	17.363	188	508693	20.000	ng/ul	-0.01
79) Chrysene-d12	21.510	240	433821	20.000	ng/ul	0.00
88) Perylene-d12	24.068	264	462563	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.270	96	10771	3.477	ng/uL	0.01
4) Pyridine-d5	3.705	84	64169	7.912	ng/ul	0.01
7) Phenol-d5	7.046	99	44210	4.610	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.222	67	172368	29.105	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.399	132	161302	20.623	ng/ul	-0.01
15) 4-Methylphenol-d8	8.599	113	91066	12.093	ng/ul	-0.01
21) Nitrobenzene-d5	9.081	128	106673	30.228	ng/ul	-0.02
24) 2-Nitrophenol-d4	9.799	143	114141	29.274	ng/ul	-0.01
28) 2,4-Dichlorophenol-d3	10.328	165	180354	23.580	ng/ul	-0.01
31) 4-Chloroaniline-d4	10.881	131	245448	24.521	ng/ul	-0.01
46) Dimethylphthalate-d6	13.993	166	630499	29.786	ng/ul	-0.01
49) Acenaphthylene-d8	14.269	160	699968	29.652	ng/ul	-0.01
54) 4-Nitrophenol-d4	14.863	143	8862m	2.938	ng/ul	0.04
60) Fluorene-d10	15.581	176	516897	28.249	ng/ul	-0.01
65) 4,6-Dinitro-2-methylph...	15.728	200	79056	23.594	ng/ul	0.00
73) Anthracene-d10	17.469	188	781328	30.056	ng/ul	0.00
81) Pyrene-d10	19.727	212	944156	34.714	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.898	264	816363	31.687	ng/ul	0.00

Target Compounds Qvalue

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

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