

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110123\
 Data File : BP017891.D
 Acq On : 02 Nov 2023 17:47
 Operator : MA/JU
 Sample : 04996-21
 Misc :
 ALS Vial : 46 Sample Multiplier: 1

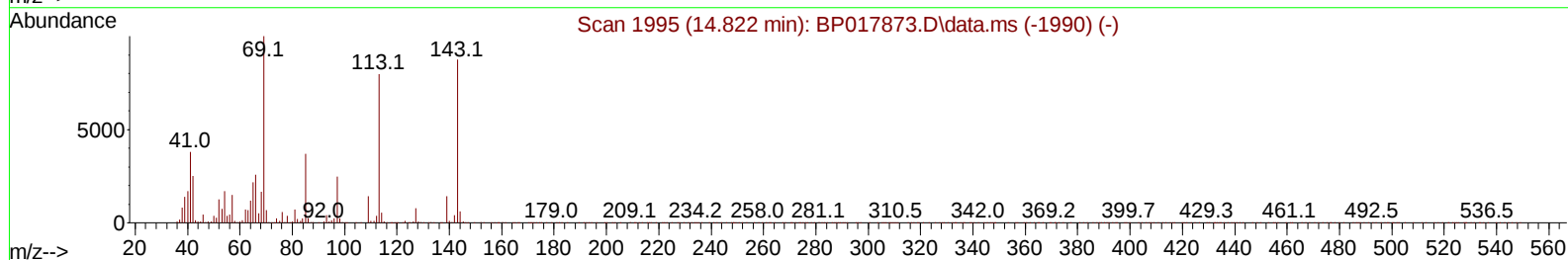
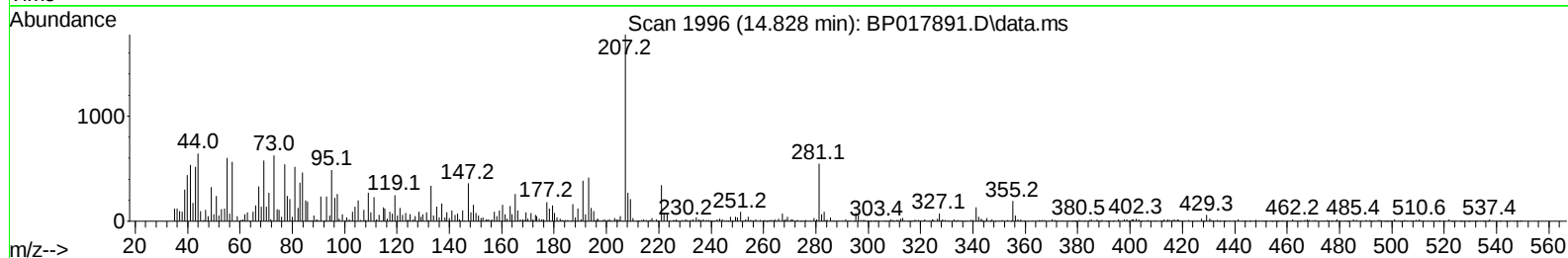
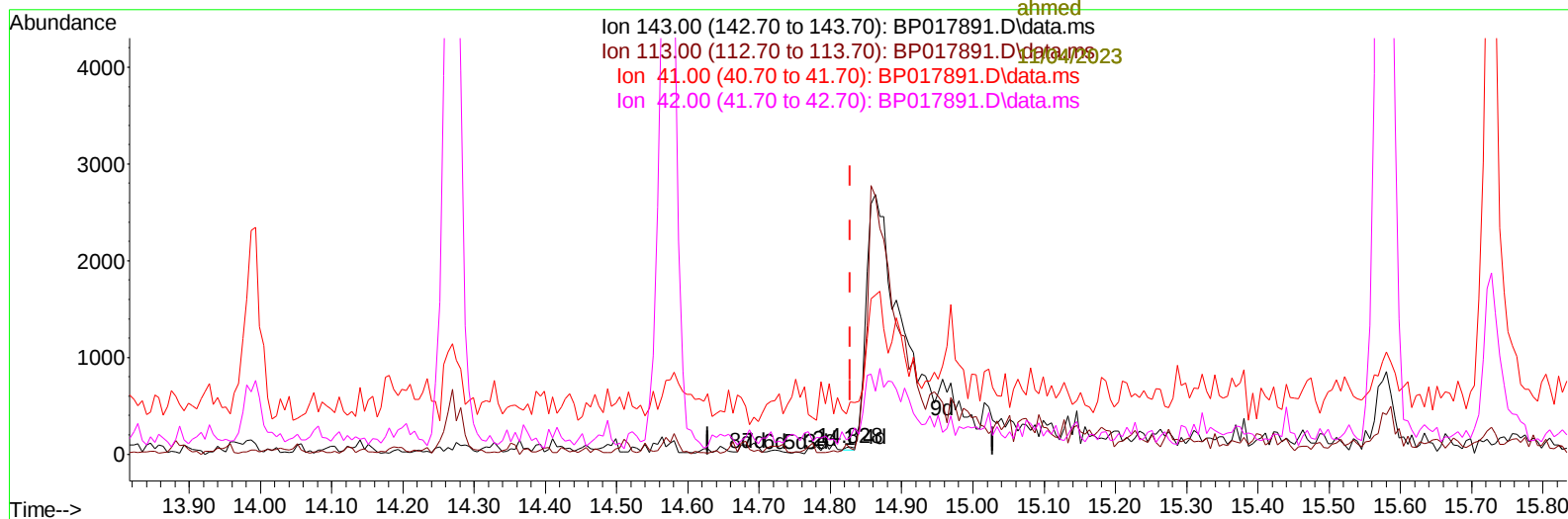
Instrument :
 BNA_P
 ClientSampleId :
 A4927

Manual IntegrationsAPPROVED

Quant Time: Nov 02 23:38:28 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

11/03/2023
 Supervised By :mohammad
 ahmed



TIC: BP017891.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.828min (+ 0.000) 0.01 ng/u1

response 27

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	81.70	80.00
41.00	39.20	716.00#
42.00	25.60	234.67#

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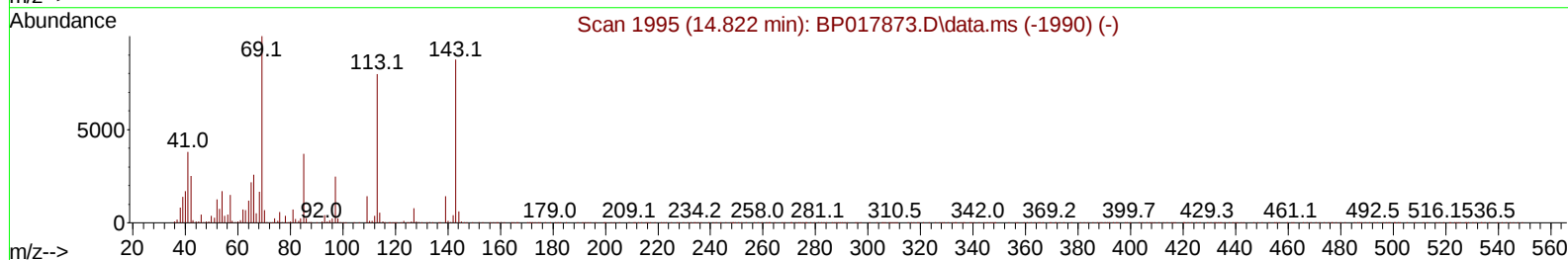
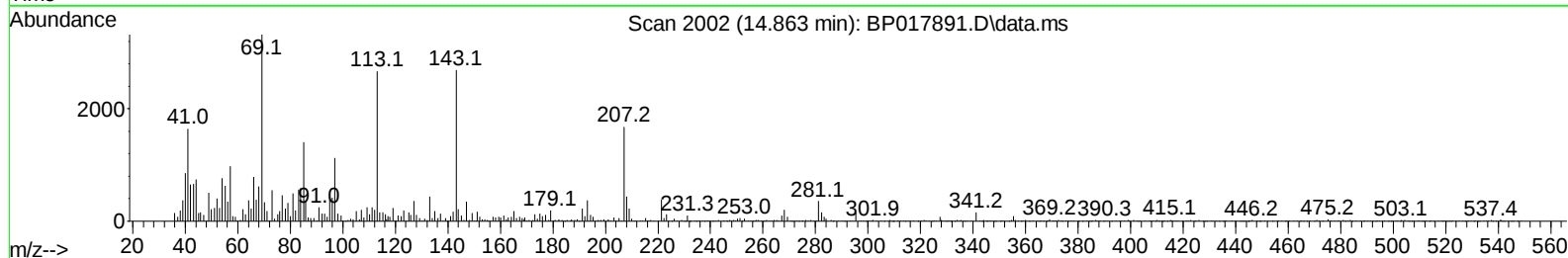
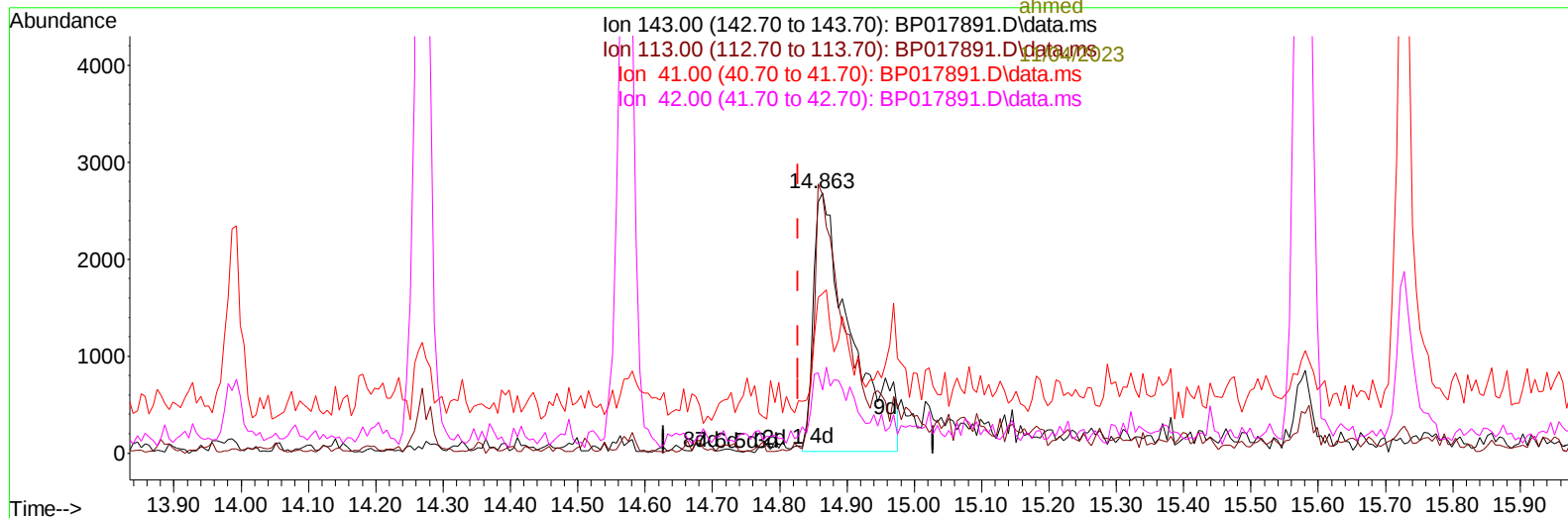
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TIC: BP017891.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.863min (+ 0.035) 3.19 ng/ul m

response 10361

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	81.70	99.29#
41.00	39.20	61.30#
42.00	25.60	24.15

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Quant Time: Nov 03 00:16:16 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

11/03/2023
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 ahmed

11/04/2023

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
-----11/04/2023						
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.875	152	121949	20.000	ng/ul	-0.01
20) Naphthalene-d8	10.705	136	463846	20.000	ng/ul	-0.02
38) Acenaphthene-d10	14.575	164	271429	20.000	ng/ul	-0.01
64) Phenanthrene-d10	17.363	188	558215	20.000	ng/ul	-0.01
79) Chrysene-d12	21.510	240	472284	20.000	ng/ul	0.00
88) Perylene-d12	24.069	264	484834	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.270	96	11894	3.628	ng/uL	0.01
4) Pyridine-d5	3.717	84	23174	2.700	ng/ul	0.02
7) Phenol-d5	7.040	99	51993	5.123	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.222	67	188087	30.011	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.399	132	183938	22.223	ng/ul	-0.01
15) 4-Methylphenol-d8	8.599	113	102995	12.924	ng/ul	-0.01
21) Nitrobenzene-d5	9.081	128	119914	31.904	ng/ul	-0.02
24) 2-Nitrophenol-d4	9.799	143	128847	31.026	ng/ul	-0.01
28) 2,4-Dichlorophenol-d3	10.328	165	202330	24.837	ng/ul	-0.01
31) 4-Chloroaniline-d4	10.887	131	94400	8.855	ng/ul	0.00
46) Dimethylphthalate-d6	13.993	166	703864	30.891	ng/ul	-0.01
49) Acenaphthylene-d8	14.269	160	798253	31.414	ng/ul	-0.01
54) 4-Nitrophenol-d4	14.863	143	10361m	3.191	ng/ul	0.04
60) Fluorene-d10	15.581	176	586243	29.764	ng/ul	-0.01
65) 4,6-Dinitro-2-methylph...	15.728	200	91833	24.976	ng/ul	0.00
73) Anthracene-d10	17.469	188	892292	31.280	ng/ul	0.00
81) Pyrene-d10	19.728	212	1023734	34.574	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.898	264	861787	31.913	ng/ul	0.00

Target Compounds Qvalue

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

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