

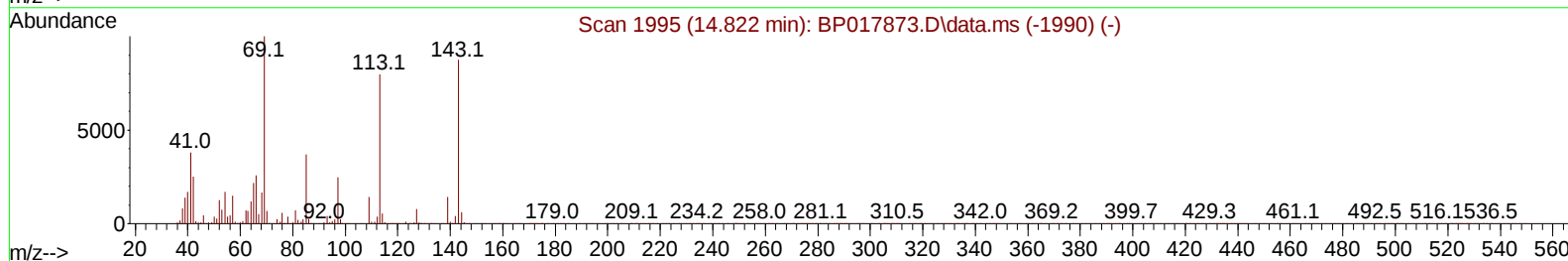
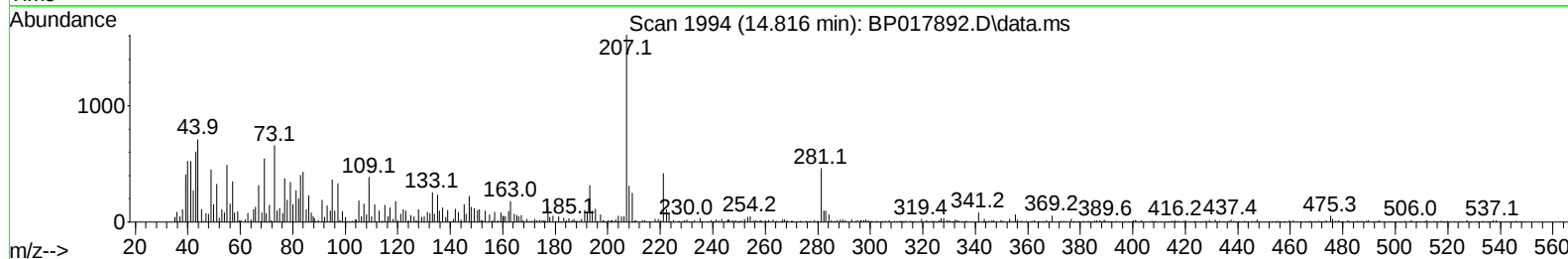
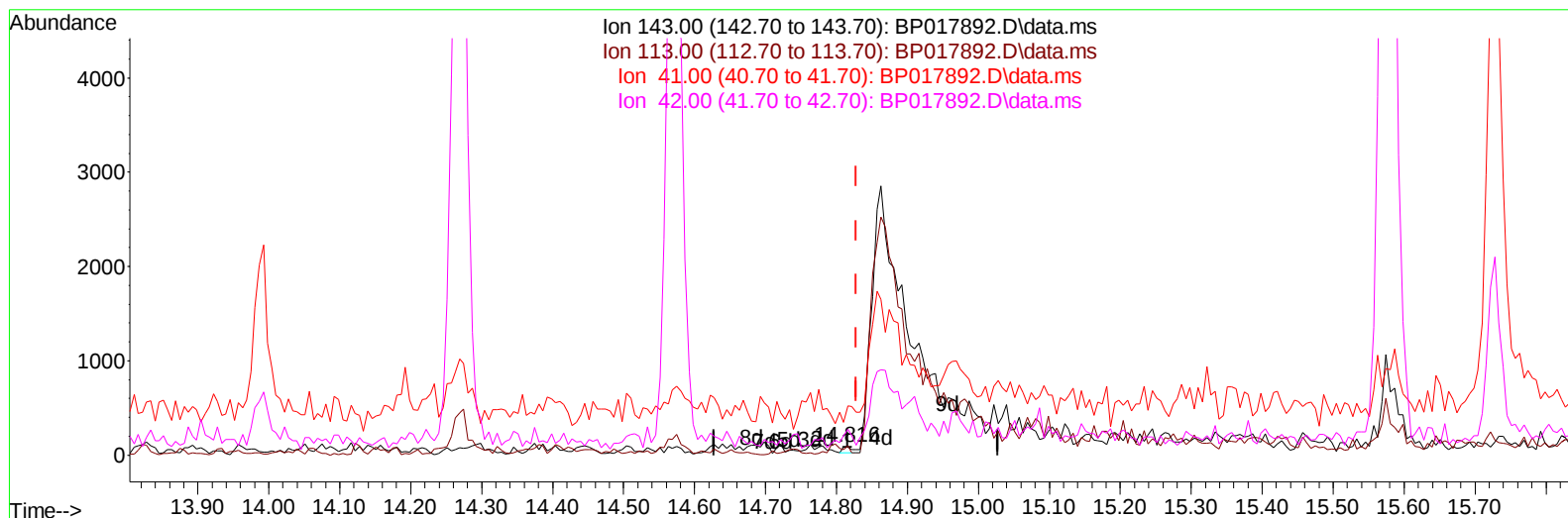
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110123\
Data File : BP017892.D
Acq On : 02 Nov 2023 18:21
Operator : MA/JU
Sample : 04996-22
Misc :
ALS Vial : 47 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A4931

Manual IntegrationsAPPROVED

Quant Time: Nov 02 23:38:39 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: BP017892.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.816min (-0.012) 0.02 ng/u1

response 74

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	81.70	125.29#
41.00	39.20	602.30#
42.00	25.60	316.09#

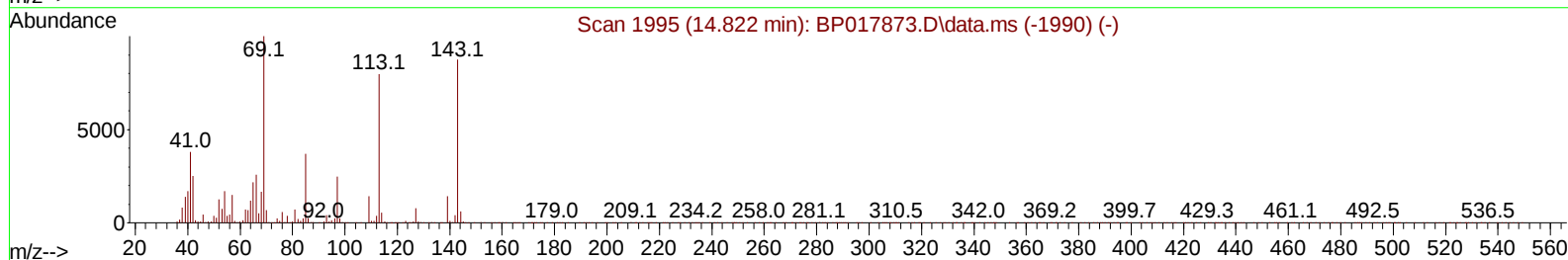
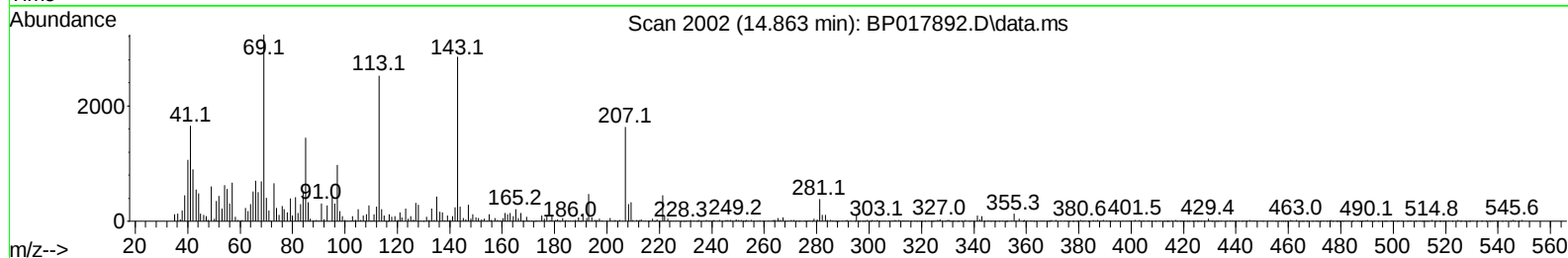
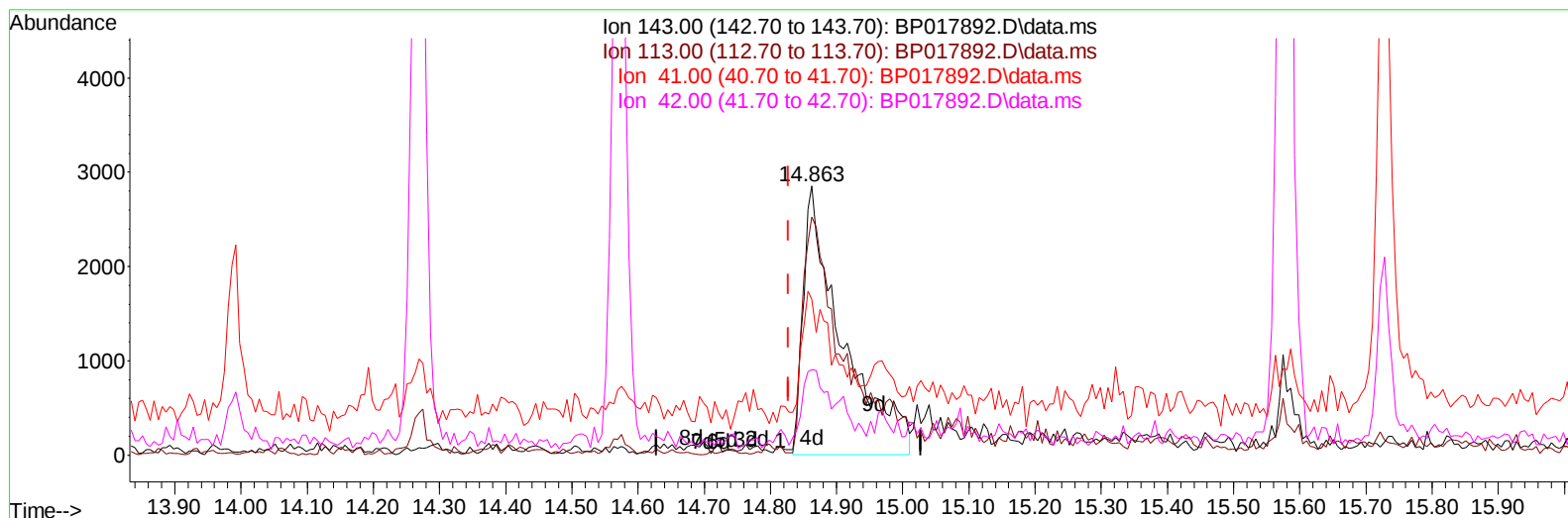
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TIC: BP017892.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.863min (+ 0.035) 3.54 ng/ul m

response 11578

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	81.70	88.50
41.00	39.20	57.96#
42.00	25.60	31.63#

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 Misc :
 ALS Vial : 47 Sample Multiplier: 1

Instrument :
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Manual IntegrationsAPPROVED

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

Quant Time: Nov 03 00:18:22 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

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.875	152	120908	20.000	ng/ul	-0.01
20) Naphthalene-d8	10.704	136	468219	20.000	ng/ul	-0.02
38) Acenaphthene-d10	14.569	164	273405	20.000	ng/ul	-0.02
64) Phenanthrene-d10	17.363	188	560924	20.000	ng/ul	-0.01
79) Chrysene-d12	21.510	240	480059	20.000	ng/ul	0.00
88) Perylene-d12	24.068	264	498602	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.264	96	11621	3.576	ng/uL	0.00
4) Pyridine-d5	3.705	84	83179	9.775	ng/ul	0.01
7) Phenol-d5	7.046	99	52617	5.230	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.222	67	176164	28.351	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.399	132	174767	21.297	ng/ul	-0.01
15) 4-Methylphenol-d8	8.599	113	103425	13.090	ng/ul	-0.01
21) Nitrobenzene-d5	9.081	128	110963	29.247	ng/ul	-0.02
24) 2-Nitrophenol-d4	9.799	143	120859	28.831	ng/ul	-0.01
28) 2,4-Dichlorophenol-d3	10.328	165	186044	22.624	ng/ul	-0.01
31) 4-Chloroaniline-d4	10.881	131	259405	24.105	ng/ul	-0.01
46) Dimethylphthalate-d6	13.992	166	644008	28.060	ng/ul	-0.01
49) Acenaphthylene-d8	14.269	160	741716	28.978	ng/ul	-0.01
54) 4-Nitrophenol-d4	14.863	143	11578m	3.540	ng/ul	0.04
60) Fluorene-d10	15.581	176	545922	27.516	ng/ul	-0.01
65) 4,6-Dinitro-2-methylph...	15.728	200	84642	22.909	ng/ul	0.00
73) Anthracene-d10	17.469	188	836699	29.189	ng/ul	0.00
81) Pyrene-d10	19.727	212	977866	32.490	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.898	264	817560	29.440	ng/ul	0.00

Target Compounds Qvalue

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

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