

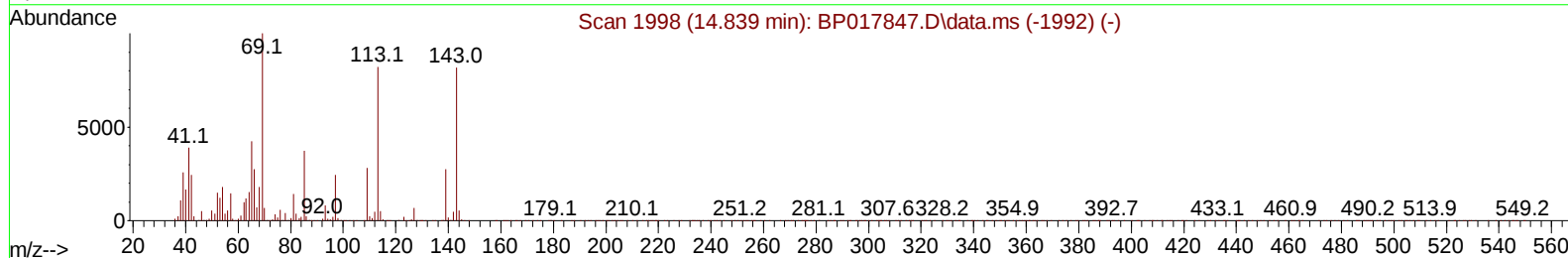
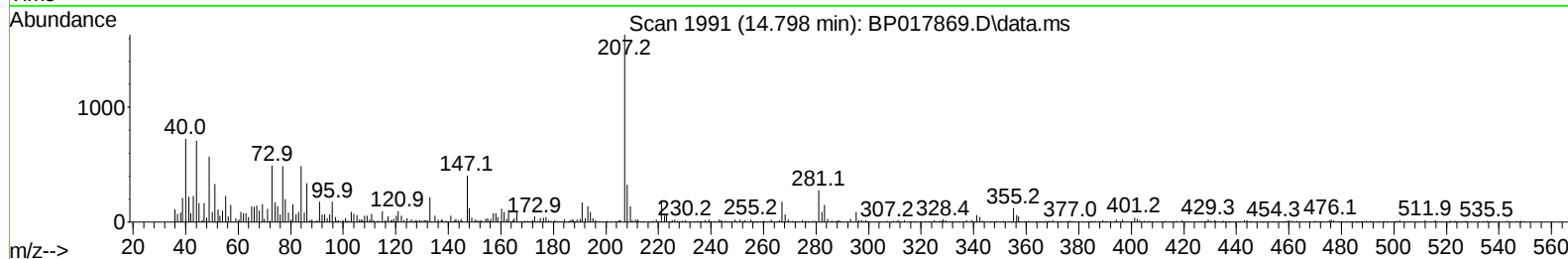
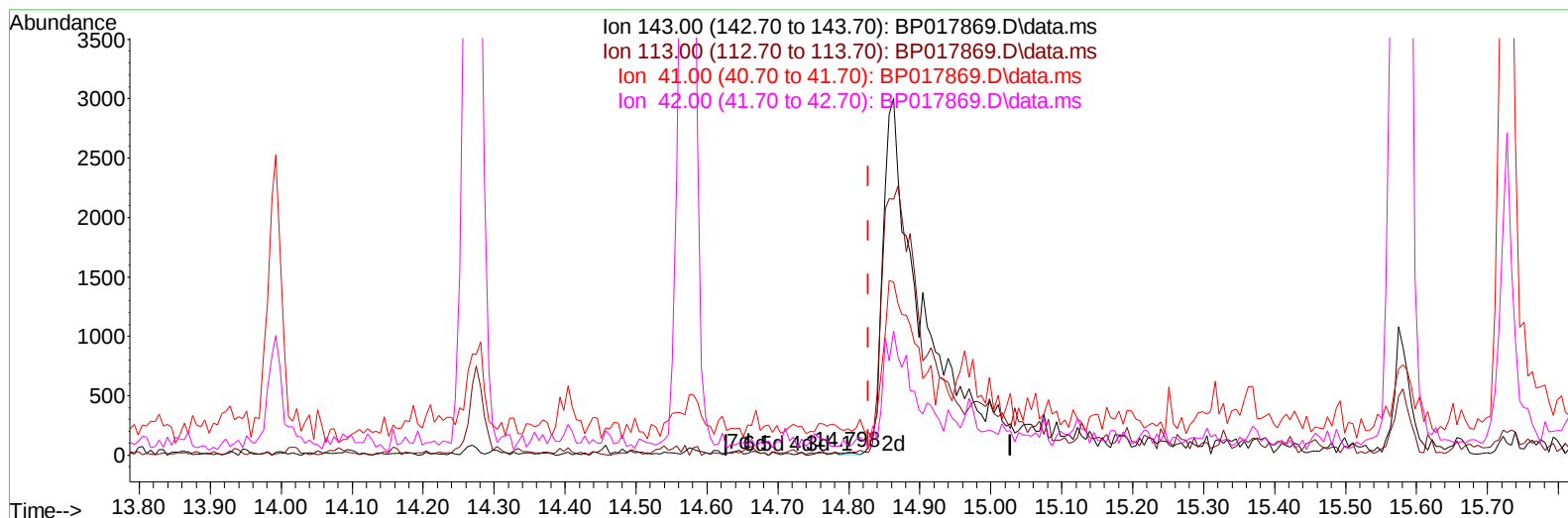
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110123\
 Data File : BP017869.D
 Acq On : 02 Nov 2023 02:14
 Operator : MA/JU
 Sample : 04944-10
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 A49R6

Manual IntegrationsAPPROVED

Quant Time: Nov 02 04:33:16 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

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



TIC: BP017869.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.798min (-0.030) 0.01 ng/u1

response 22

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	81.70	74.07
41.00	39.20	818.52#
42.00	25.60	292.59#

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.875	152	141746	20.000	ng/ul	-0.01
20) Naphthalene-d8	10.704	136	526467	20.000	ng/ul	-0.02
38) Acenaphthene-d10	14.575	164	296558	20.000	ng/ul	-0.01
64) Phenanthrene-d10	17.363	188	609340	20.000	ng/ul	#-0.01
79) Chrysene-d12	21.510	240	525067	20.000	ng/ul	0.00
88) Perylene-d12	24.063	264	550343	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.281	96	14393	3.777	ng/uL	0.02
4) Pyridine-d5	3.711	84	107775	10.804	ng/ul	0.02
7) Phenol-d5	7.046	99	62334	5.285	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.222	67	248418	34.102	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.399	132	225877	23.479	ng/ul	-0.01
15) 4-Methylphenol-d8	8.599	113	124372	13.427	ng/ul	-0.01
21) Nitrobenzene-d5	9.087	128	147069	34.474	ng/ul	-0.01
24) 2-Nitrophenol-d4	9.799	143	150245	31.875	ng/ul	-0.01
28) 2,4-Dichlorophenol-d3	10.328	165	240269	25.986	ng/ul	-0.01
31) 4-Chloroaniline-d4	10.881	131	331681	27.411	ng/ul	-0.01
46) Dimethylphthalate-d6	13.992	166	837624	33.646	ng/ul	-0.01
49) Acenaphthylene-d8	14.269	160	918933	33.099	ng/ul	-0.01
54) 4-Nitrophenol-d4	14.863	143	11022m	3.107	ng/ul	0.04
60) Fluorene-d10	15.581	176	690796	32.100	ng/ul	-0.01
65) 4,6-Dinitro-2-methylph...	15.728	200	98583	24.562	ng/ul	0.00
73) Anthracene-d10	17.469	188	1052943	33.815	ng/ul	0.00
81) Pyrene-d10	19.727	212	1207868	36.692	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.898	264	1045055	34.094	ng/ul	0.00

Target Compounds Qvalue

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

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