

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110123\
 Data File : BP017870.D
 Acq On : 02 Nov 2023 02:48
 Operator : MA/JU
 Sample : 04944-11
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

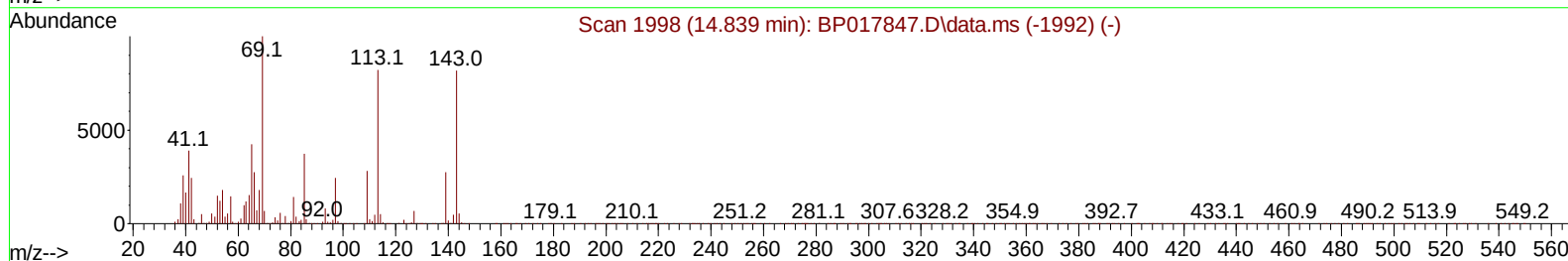
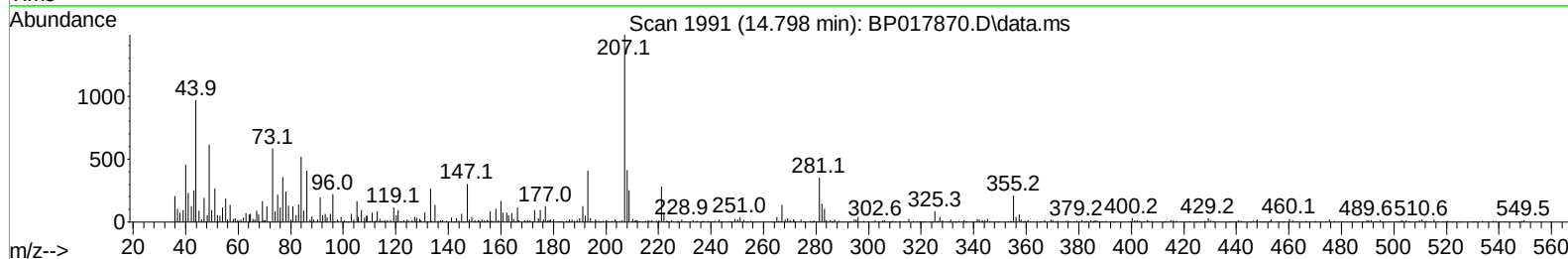
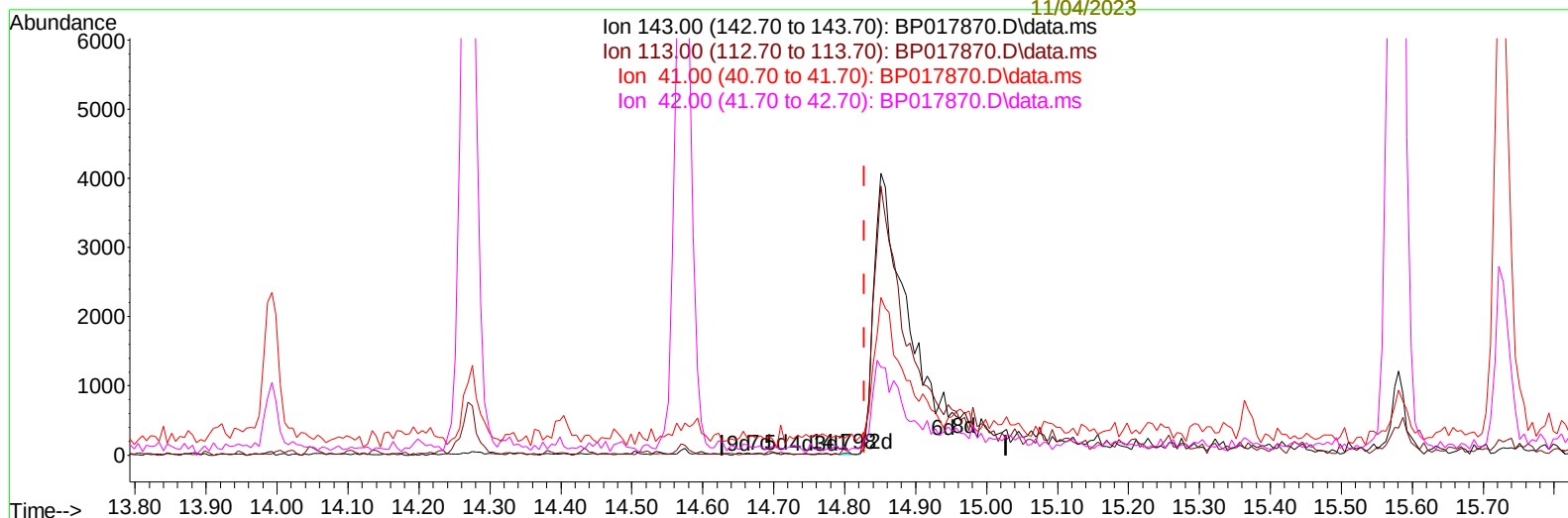
Instrument :
 BNA_P
 ClientSampleId :
 A49R8

Manual IntegrationsAPPROVED

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

Supervised By :mohammad
 ahmed
 11/04/2023

Quant Time: Nov 02 04:33:31 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



TIC: BP017870.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.798min (-0.029) 0.01 ng/ul

response 21

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	81.70	303.57#
41.00	39.20	825.00#
42.00	25.60	450.00#

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.875	152	157713	20.000	ng/ul	-0.01
20) Naphthalene-d8	10.704	136	594860	20.000	ng/ul	-0.02
38) Acenaphthene-d10	14.575	164	337298	20.000	ng/ul	-0.01
64) Phenanthrene-d10	17.363	188	691086	20.000	ng/ul	#-0.01
79) Chrysene-d12	21.510	240	592916	20.000	ng/ul	0.00
88) Perylene-d12	24.063	264	636484	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.275	96	16777	3.957	ng/uL	0.02
4) Pyridine-d5	3.717	84	98088	8.837	ng/ul	0.02
7) Phenol-d5	7.046	99	75959	5.788	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.222	67	267490	33.003	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.399	132	257117	24.020	ng/ul	-0.01
15) 4-Methylphenol-d8	8.599	113	147363	14.298	ng/ul	-0.01
21) Nitrobenzene-d5	9.087	128	158442	32.870	ng/ul	-0.01
24) 2-Nitrophenol-d4	9.799	143	169226	31.774	ng/ul	-0.01
28) 2,4-Dichlorophenol-d3	10.328	165	271230	25.962	ng/ul	-0.01
31) 4-Chloroaniline-d4	10.881	131	321348	23.504	ng/ul	-0.01
46) Dimethylphthalate-d6	13.993	166	907698	32.057	ng/ul	-0.01
49) Acenaphthylene-d8	14.269	160	989492	31.336	ng/ul	-0.01
54) 4-Nitrophenol-d4	14.851	143	14752m	3.656	ng/ul	0.02
60) Fluorene-d10	15.581	176	750679	30.670	ng/ul	-0.01
65) 4,6-Dinitro-2-methylph...	15.728	200	117999	25.922	ng/ul	0.00
73) Anthracene-d10	17.469	188	1163911	32.957	ng/ul	0.00
81) Pyrene-d10	19.728	212	1379313	37.105	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.898	264	1194721	33.701	ng/ul	0.00

Target Compounds Qvalue

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

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