

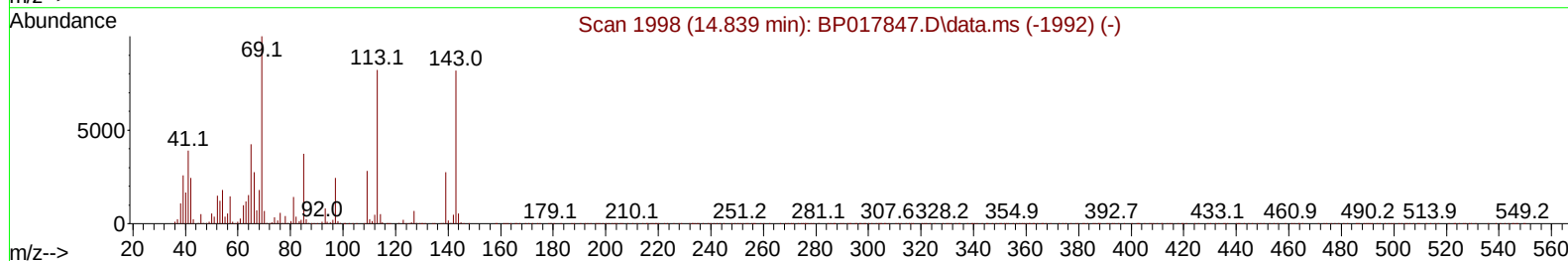
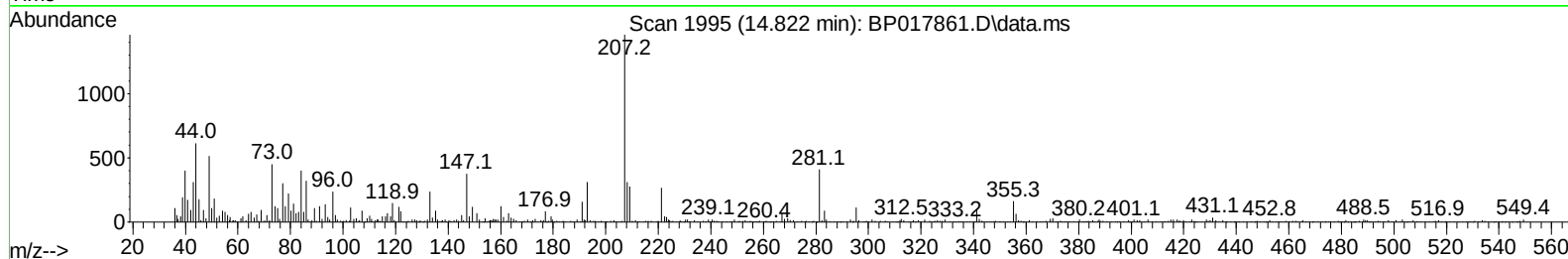
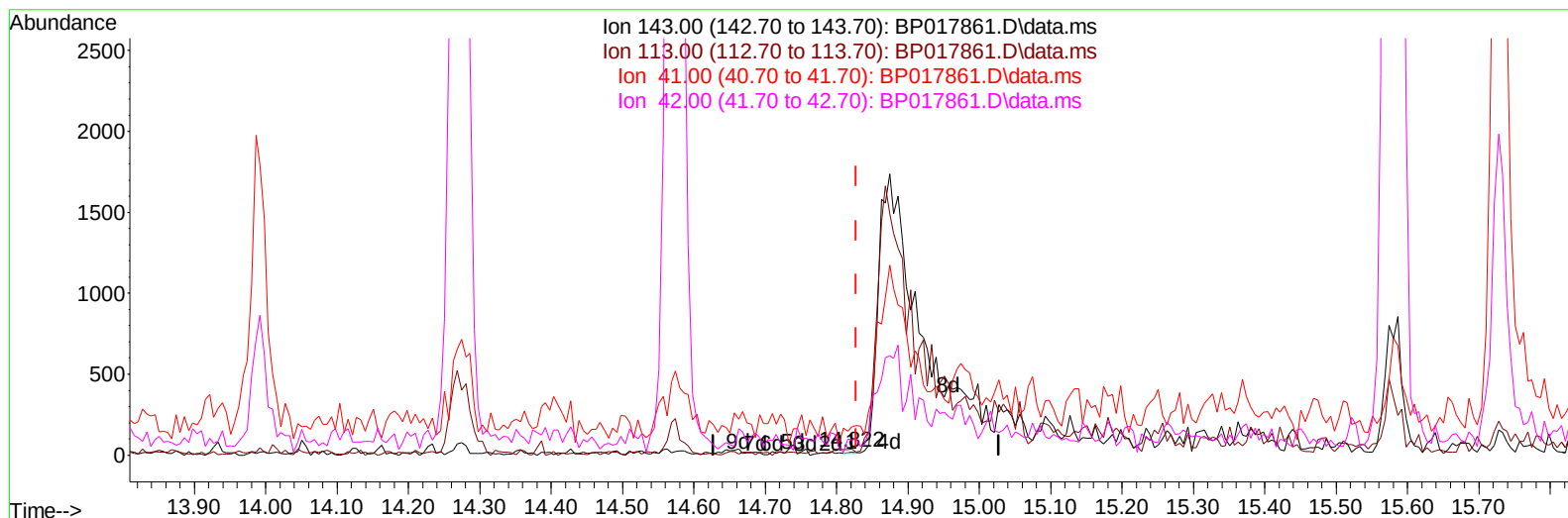
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110123\
 Data File : BP017861.D
 Acq On : 01 Nov 2023 21:44
 Operator : MA/JU
 Sample : 04944-02
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 A49P5

Manual IntegrationsAPPROVED

Quant Time: Nov 02 01:16:27 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: BP017861.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.822min (-0.006) 0.00 ng/u1

response 6

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	81.70	200.00#
41.00	39.20	921.05#
42.00	25.60	489.47#

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Quant Time: Nov 02 02:09:53 2023
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.875	152	146500	20.000	ng/ul	-0.01
20) Naphthalene-d8	10.704	136	542541	20.000	ng/ul	-0.02
38) Acenaphthene-d10	14.574	164	305362	20.000	ng/ul	-0.01
64) Phenanthrene-d10	17.368	188	621588	20.000	ng/ul	0.00
79) Chrysene-d12	21.515	240	542056	20.000	ng/ul	0.00
88) Perylene-d12	24.074	264	584081	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.281	96	12546	3.186	ng/uL	0.02
4) Pyridine-d5	3.716	84	57231	5.551	ng/ul	0.02
7) Phenol-d5	7.045	99	44572	3.656	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.222	67	207353	27.541	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.398	132	182045	18.309	ng/ul	-0.01
15) 4-Methylphenol-d8	8.604	113	93779	9.796	ng/ul	0.00
21) Nitrobenzene-d5	9.086	128	118487	26.952	ng/ul	-0.01
24) 2-Nitrophenol-d4	9.798	143	124313	25.592	ng/ul	-0.01
28) 2,4-Dichlorophenol-d3	10.328	165	195203	20.486	ng/ul	-0.01
31) 4-Chloroaniline-d4	10.886	131	277116	22.223	ng/ul	0.00
46) Dimethylphthalate-d6	13.992	166	712073	27.778	ng/ul	-0.01
49) Acenaphthylene-d8	14.274	160	750032	26.236	ng/ul	0.00
54) 4-Nitrophenol-d4	14.874	143	7007m	1.918	ng/ul	0.05
60) Fluorene-d10	15.580	176	575789	25.985	ng/ul	-0.01
65) 4,6-Dinitro-2-methylph...	15.727	200	85658	20.921	ng/ul	0.00
73) Anthracene-d10	17.468	188	880644	27.724	ng/ul	0.00
81) Pyrene-d10	19.733	212	1030534	30.324	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.903	264	891757	27.412	ng/ul	0.00

Target Compounds Qvalue

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

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