

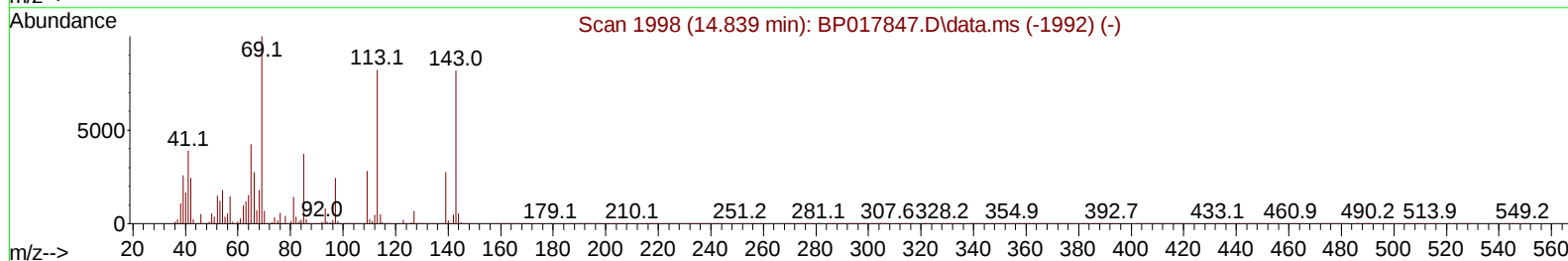
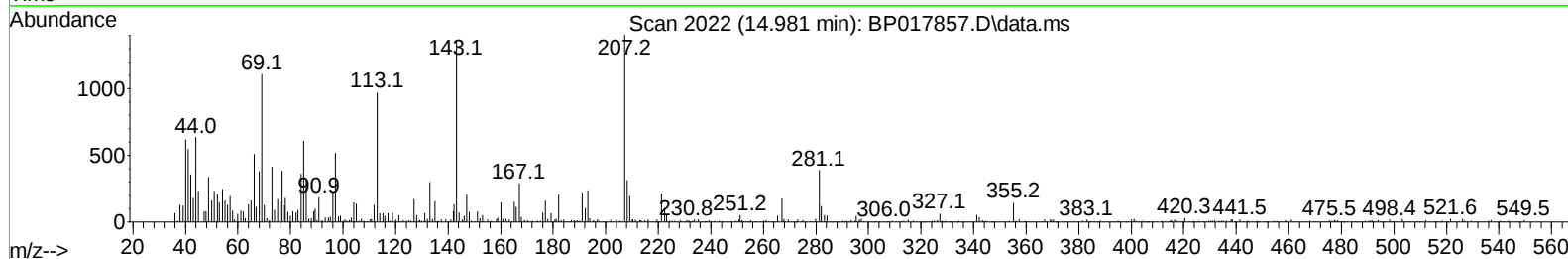
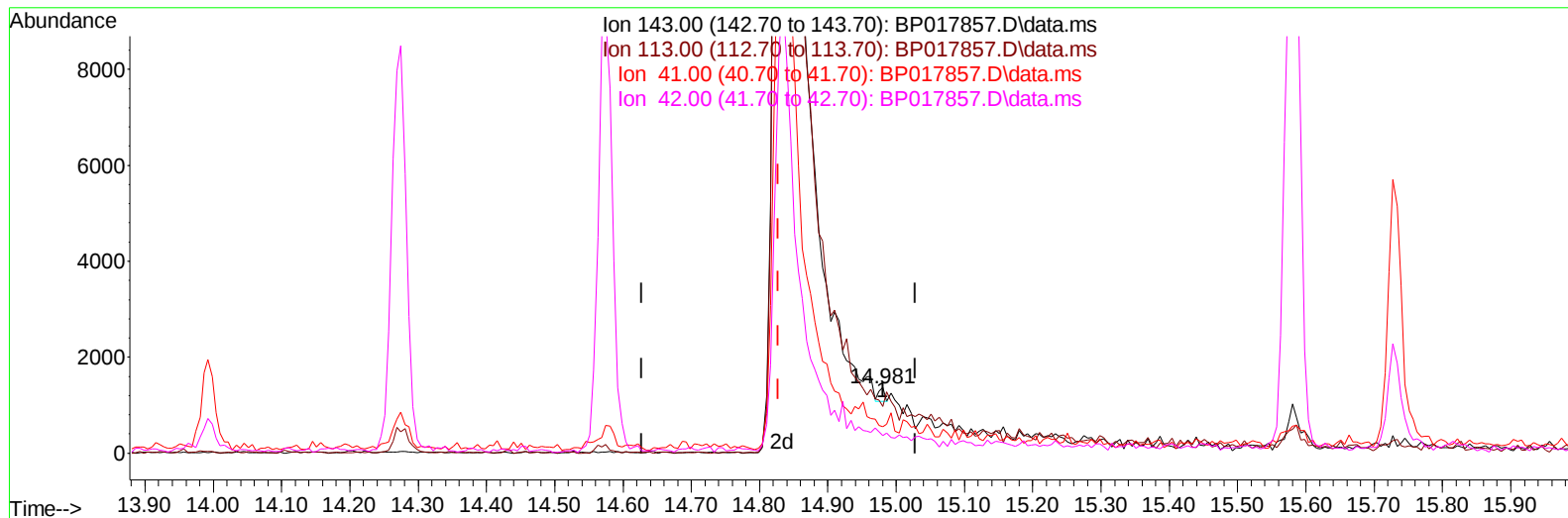
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110123\
 Data File : BP017857.D
 Acq On : 01 Nov 2023 19:29
 Operator : MA/JU
 Sample : PB156542BL
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK542

Manual IntegrationsAPPROVED

Quant Time: Nov 02 01:15:20 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: BP017857.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.981min (+ 0.153) 0.04 ng/u1

response 180

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	81.70	71.30
41.00	39.20	40.04
42.00	25.60	26.35

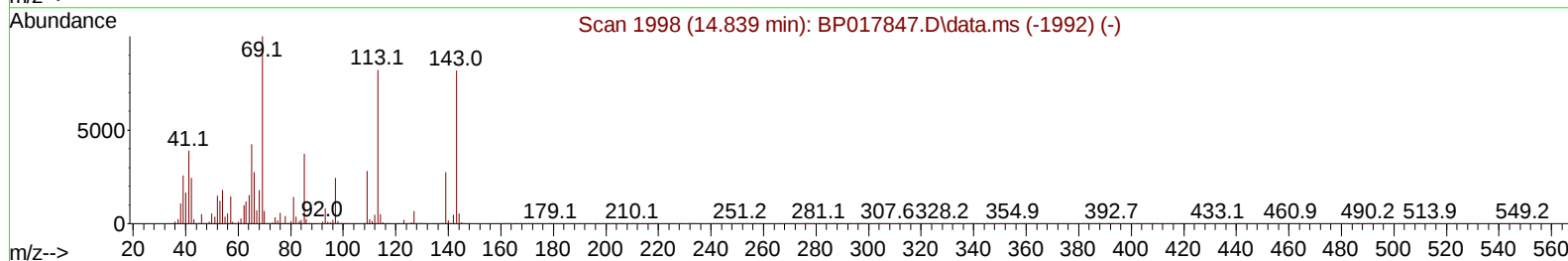
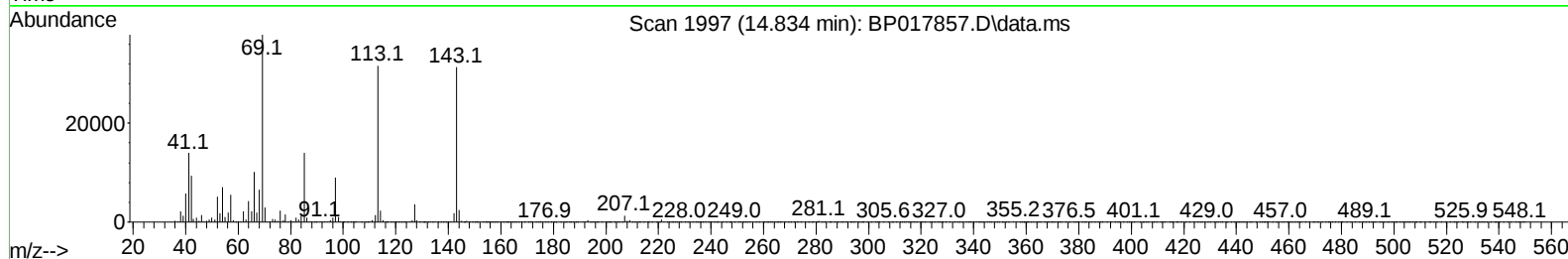
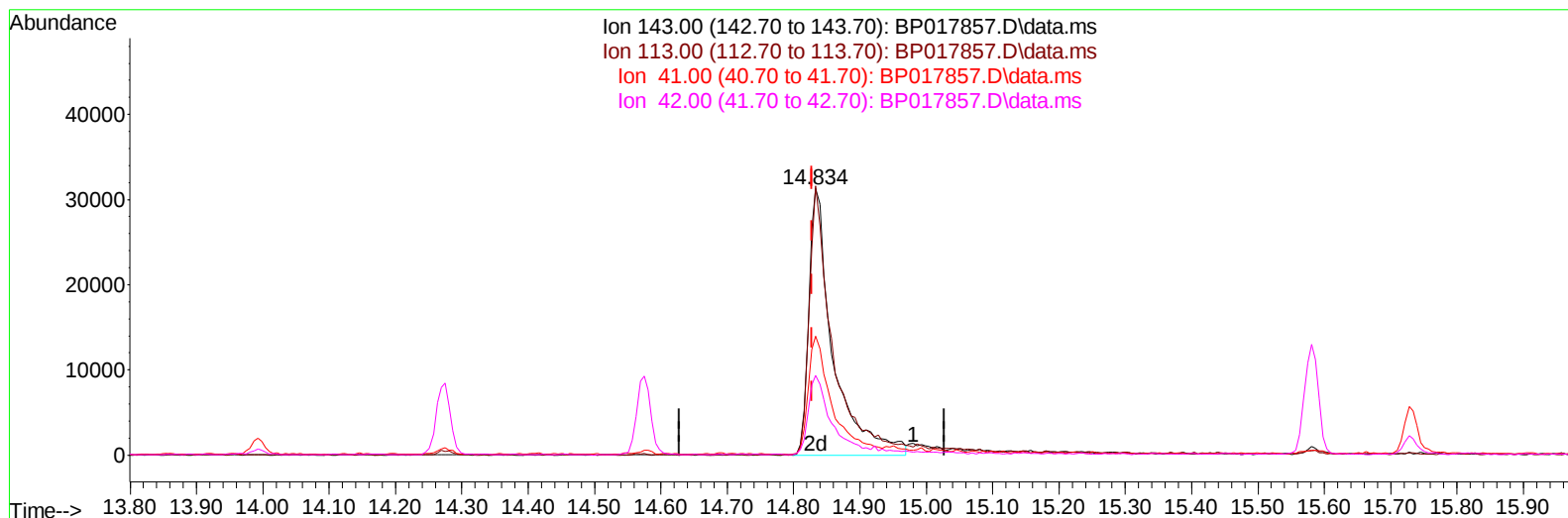
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TIC: BP017857.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.834min (+ 0.006) 19.72 ng/ul m

response 78950

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	81.70	100.89#
41.00	39.20	44.63
42.00	25.60	29.99

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 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
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Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 11/03/2023
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Quant Time: Nov 02 02:01:59 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
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.875	152	160600	20.000	ng/ul	-0.01
20) Naphthalene-d8	10.710	136	593465	20.000	ng/ul	-0.01
38) Acenaphthene-d10	14.575	164	334694	20.000	ng/ul	-0.01
64) Phenanthrene-d10	17.369	188	699107	20.000	ng/ul	0.00
79) Chrysene-d12	21.516	240	579990	20.000	ng/ul	0.00
88) Perylene-d12	24.074	264	617117	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.275	96	21367	4.949	ng/uL	0.02
4) Pyridine-d5	3.705	84	240319	21.262	ng/ul	0.01
7) Phenol-d5	7.040	99	326433	24.426	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.228	67	214691	26.012	ng/ul	0.00
11) 2-Chlorophenol-d4	7.399	132	274346	25.169	ng/ul	-0.01
15) 4-Methylphenol-d8	8.604	113	258622	24.643	ng/ul	0.00
21) Nitrobenzene-d5	9.087	128	126417	26.288	ng/ul	-0.01
24) 2-Nitrophenol-d4	9.804	143	142273	26.776	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.328	165	242380	23.255	ng/ul	-0.01
31) 4-Chloroaniline-d4	10.887	131	314089	23.027	ng/ul	0.00
46) Dimethylphthalate-d6	13.992	166	710741	25.296	ng/ul	-0.01
49) Acenaphthylene-d8	14.275	160	786011	25.085	ng/ul	0.00
54) 4-Nitrophenol-d4	14.834	143	78950m	19.721	ng/ul	0.00
60) Fluorene-d10	15.581	176	590296	24.305	ng/ul	-0.01
65) 4,6-Dinitro-2-methylph...	15.733	200	87669	19.038	ng/ul	0.00
73) Anthracene-d10	17.469	188	911265	25.507	ng/ul	0.00
81) Pyrene-d10	19.733	212	1026478	28.229	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.904	264	888959	25.863	ng/ul	0.00

Target Compounds Qvalue

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

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