

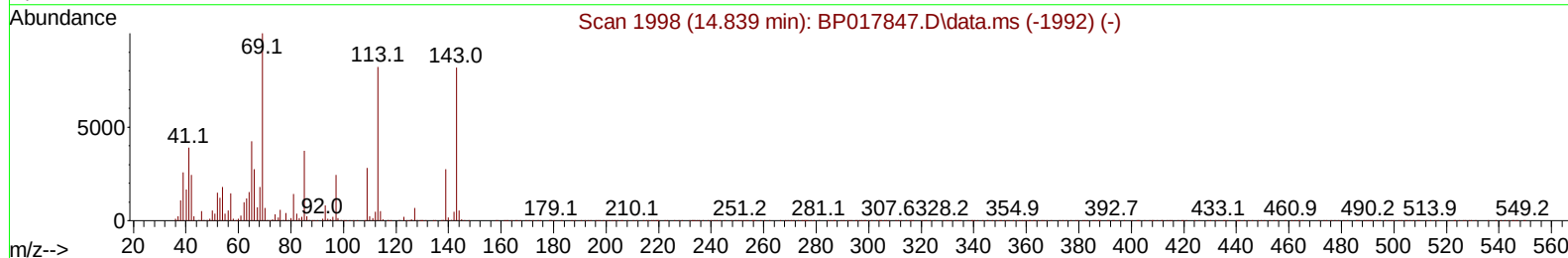
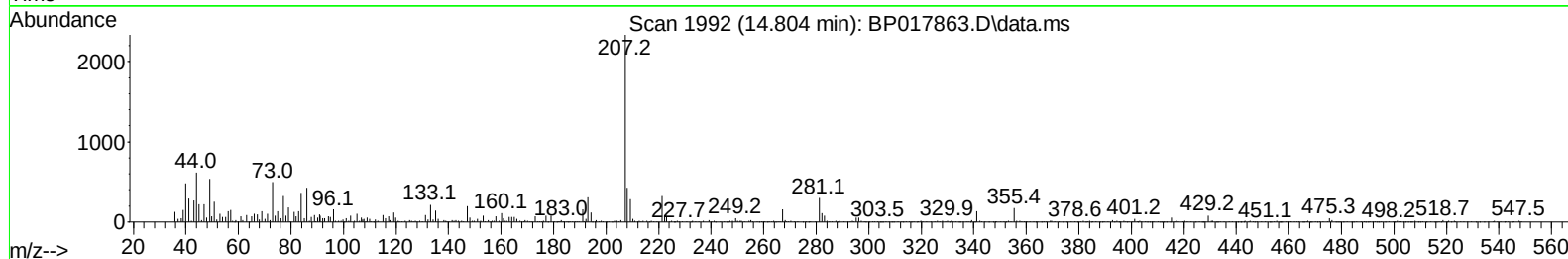
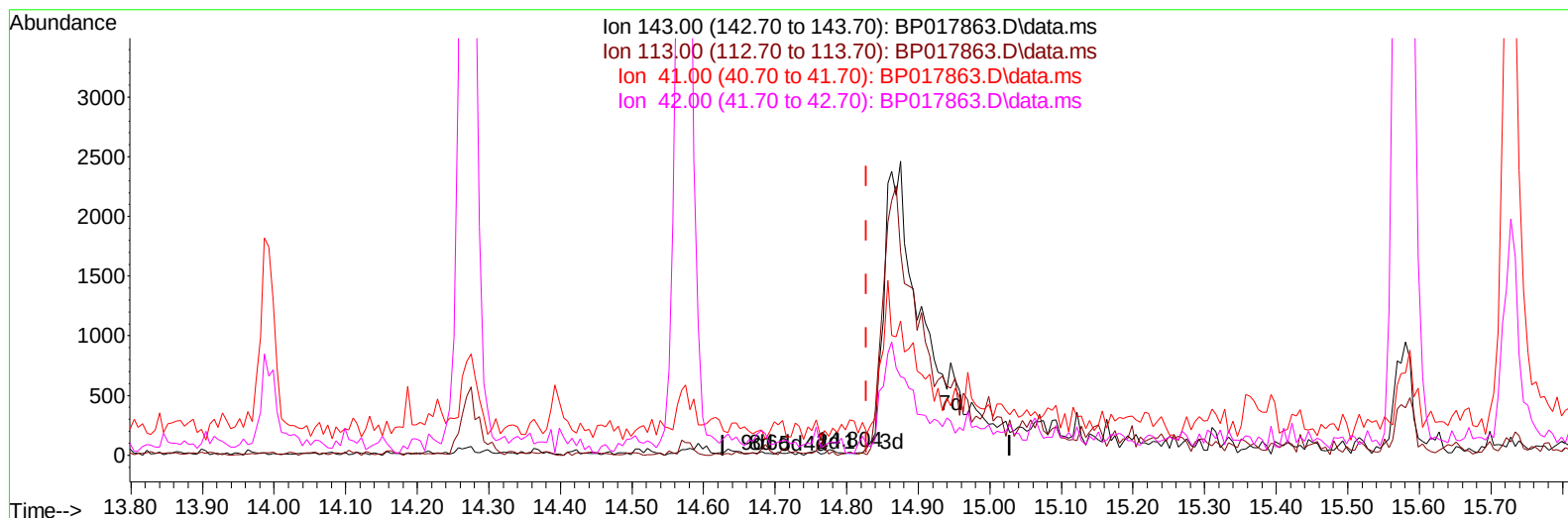
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110123\
Data File : BP017863.D
Acq On : 01 Nov 2023 22:52
Operator : MA/JU
Sample : 04944-04
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A49T0

Manual IntegrationsAPPROVED

Quant Time: Nov 02 01:16:47 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: BP017863.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.804min (-0.023) 0.00 ng/u1

response 5

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	81.70	70.37
41.00	39.20	1092.59#
42.00	25.60	51.85#

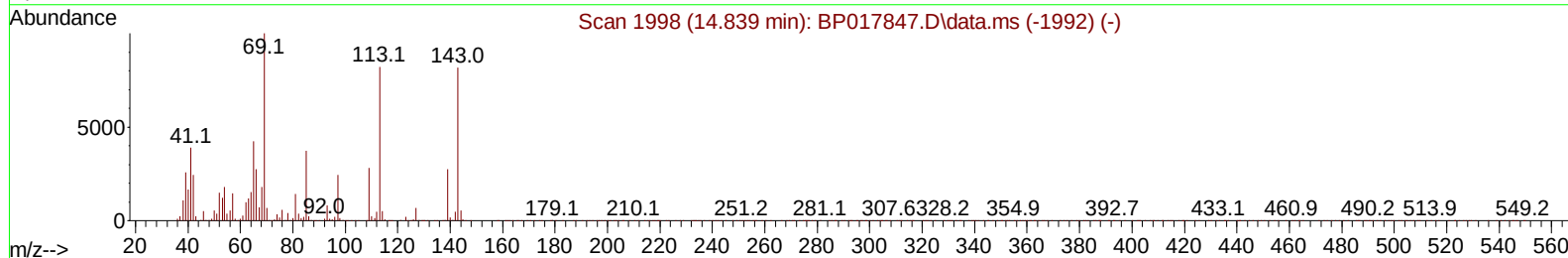
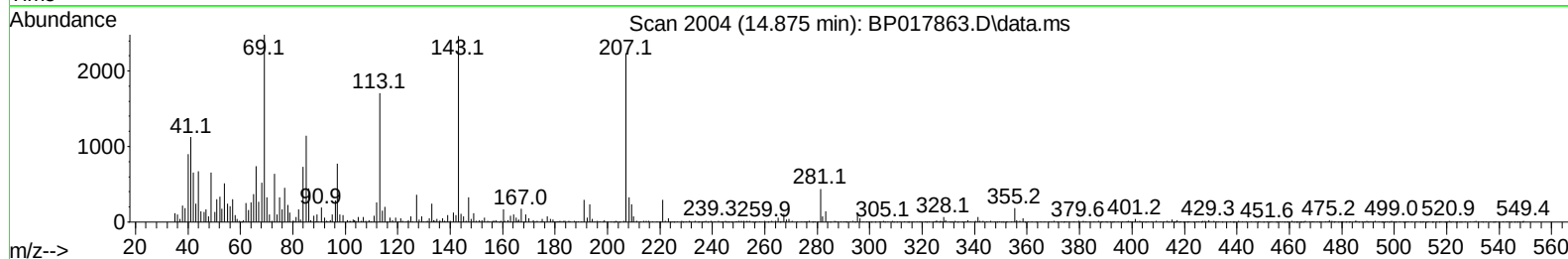
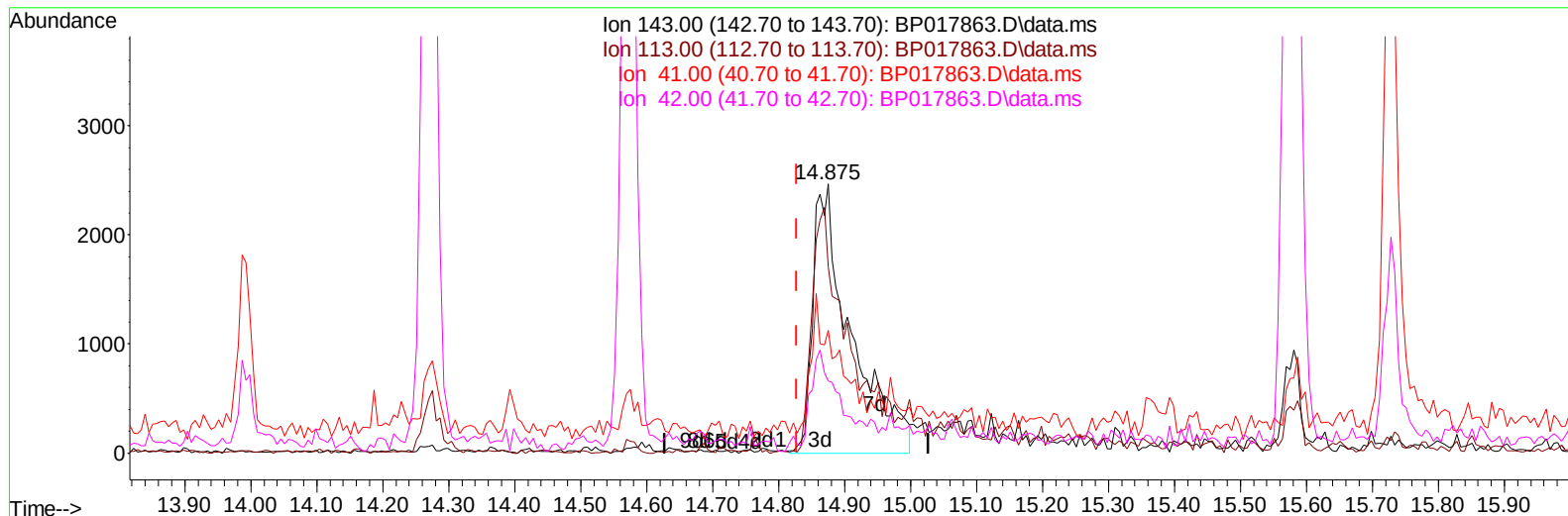
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(54) 4-Nitrophenol-d4 (S)

14.875min (+ 0.047) 3.01 ng/ul m

response 9813

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	81.70	69.13
41.00	39.20	45.56
42.00	25.60	26.73

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Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 11/03/2023
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Quant Time: Nov 02 02:13:07 2023
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 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	129525	20.000	ng/ul	-0.01
20) Naphthalene-d8	10.705	136	478904	20.000	ng/ul	-0.02
38) Acenaphthene-d10	14.575	164	272373	20.000	ng/ul	-0.01
64) Phenanthrene-d10	17.369	188	549139	20.000	ng/ul	0.00
79) Chrysene-d12	21.510	240	471895	20.000	ng/ul	0.00
88) Perylene-d12	24.069	264	508137	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.276	96	13398	3.848	ng/uL	0.02
4) Pyridine-d5	3.717	84	61925	6.793	ng/ul	0.02
7) Phenol-d5	7.046	99	53566	4.970	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.222	67	202729	30.456	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.399	132	191300	21.761	ng/ul	-0.01
15) 4-Methylphenol-d8	8.605	113	107421	12.691	ng/ul	0.00
21) Nitrobenzene-d5	9.087	128	116696	30.072	ng/ul	-0.01
24) 2-Nitrophenol-d4	9.799	143	126628	29.533	ng/ul	-0.01
28) 2,4-Dichlorophenol-d3	10.328	165	211930	25.197	ng/ul	-0.01
31) 4-Chloroaniline-d4	10.887	131	278406	25.293	ng/ul	0.00
46) Dimethylphthalate-d6	13.993	166	731628	31.998	ng/ul	-0.01
49) Acenaphthylene-d8	14.269	160	774427	30.371	ng/ul	-0.01
54) 4-Nitrophenol-d4	14.875	143	9813m	3.012	ng/ul	0.05
60) Fluorene-d10	15.581	176	597756	30.243	ng/ul	-0.01
65) 4,6-Dinitro-2-methylph...	15.728	200	85687	23.689	ng/ul	0.00
73) Anthracene-d10	17.469	188	912577	32.520	ng/ul	0.00
81) Pyrene-d10	19.728	212	1064510	35.981	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.898	264	936070	33.075	ng/ul	0.00

Target Compounds Qvalue

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

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