

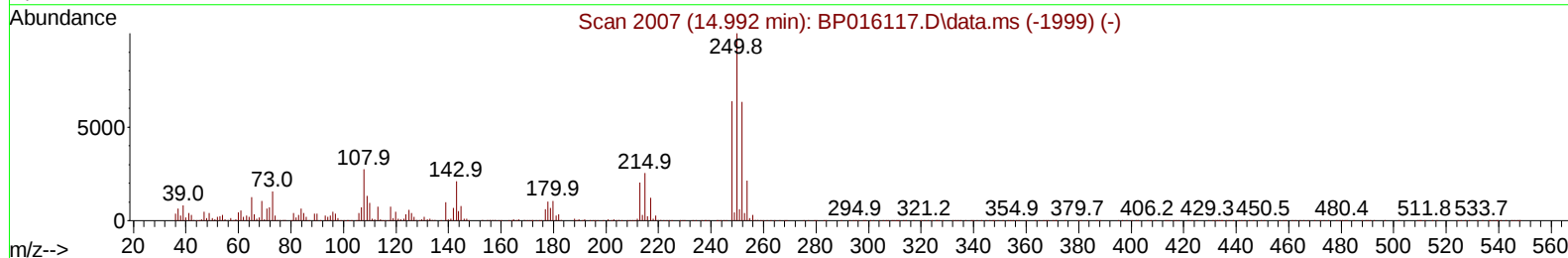
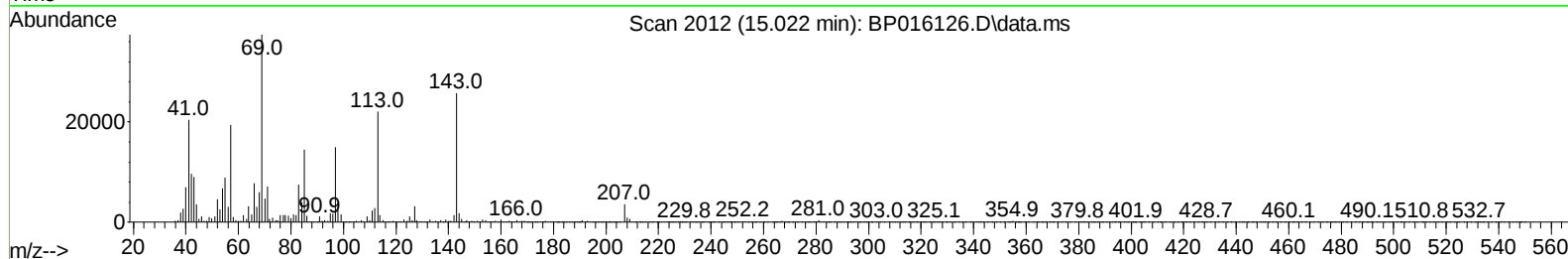
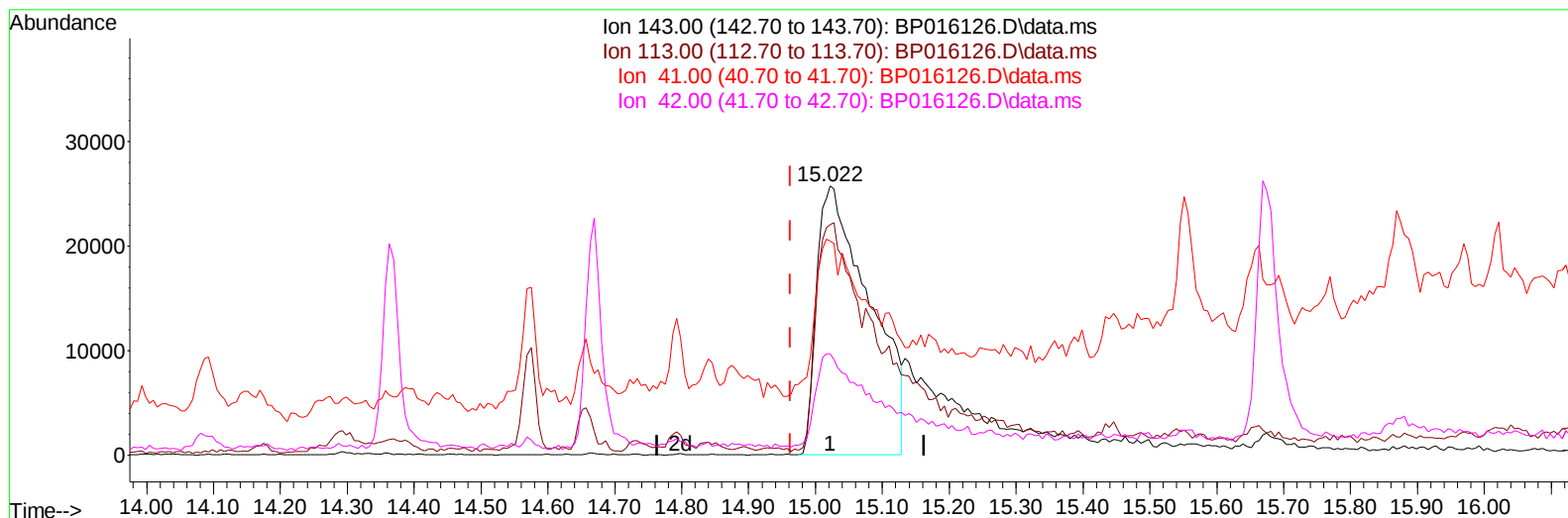
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP070823\
Data File : BP016126.D
Acq On : 09 Jul 2023 11:08
Operator : MA/JU
Sample : 03356-05
Misc :
ALS Vial : 50 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
YBW34

Manual IntegrationsAPPROVED

Quant Time: Jul 09 22:40:33 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062623.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Sat Jul 08 10:11:43 2023
Response via : Initial Calibration

Reviewed By :Yogesh Patel 07/10/2023
Supervised By :mohammad ahmed 07/10/2023



TIC: BP016126.D\data.ms

(54) 4-Nitrophenol-d4 (S)

15.022min (+ 0.059) 15.41 ng/u1

response 140806

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	94.00	85.78
41.00	45.70	79.50#
42.00	31.20	37.60#

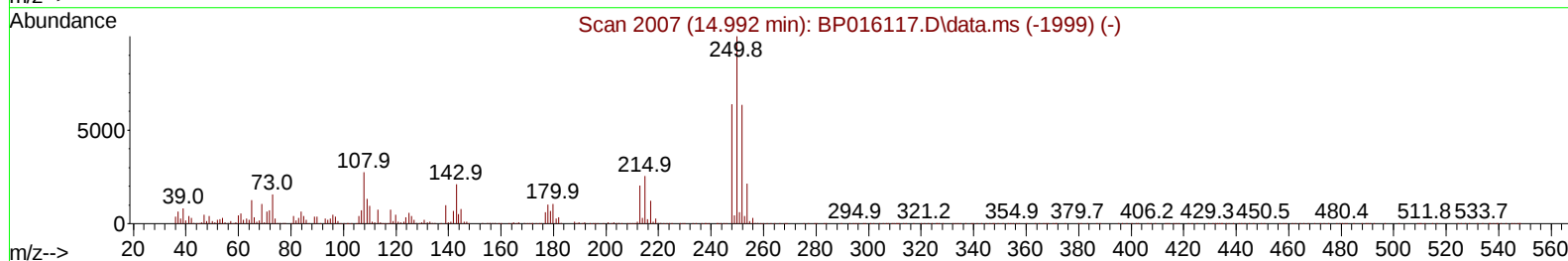
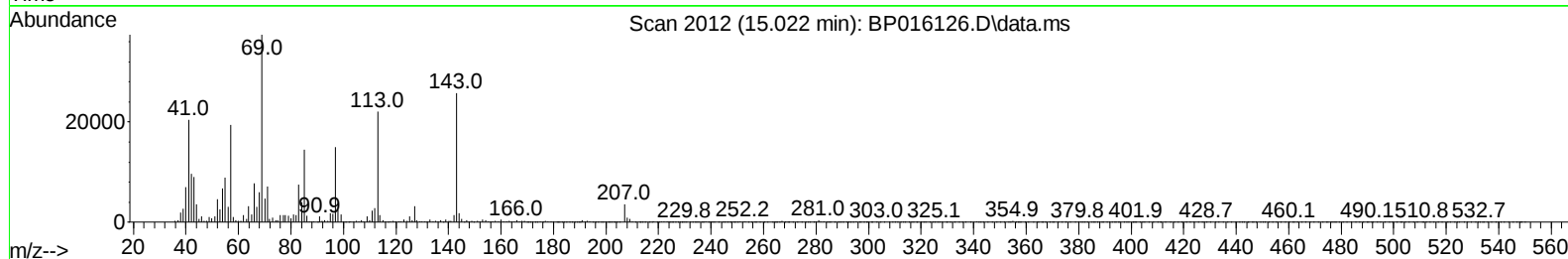
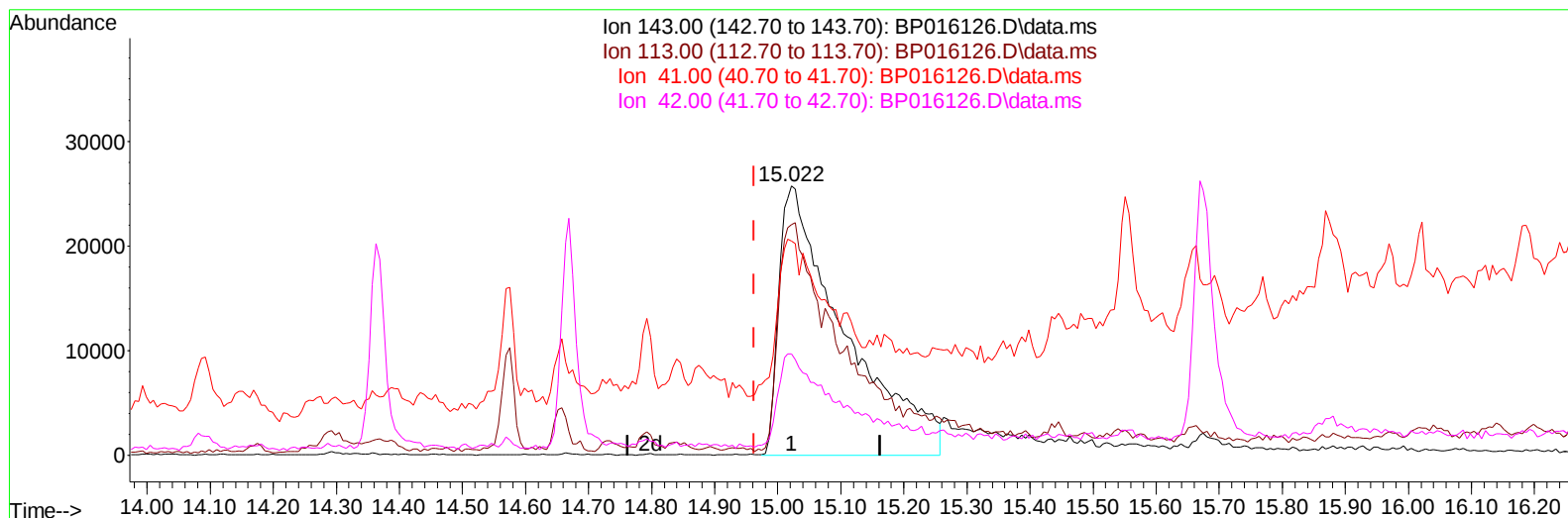
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(54) 4-Nitrophenol-d4 (S)

15.022min (+ 0.059) 20.23 ng/ul m

response 184906

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	94.00	85.78
41.00	45.70	79.50#
42.00	31.20	37.60#

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Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 07/10/2023
 Supervised By :mohammad ahmed 07/10/2023

Quant Time: Jul 09 23:28:25 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062623.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sat Jul 08 10:11:43 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.016	152	317495	20.000	ng/ul	-0.01
20) Naphthalene-d8	10.846	136	1409548	20.000	ng/ul	# 0.00
38) Acenaphthene-d10	14.669	164	895913	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.434	188	1975365	20.000	ng/ul	#-0.01
79) Chrysene-d12	21.539	240	1764103	20.000	ng/ul	0.00
88) Perylene-d12	24.068	264	1804916	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.346	96	39192	4.441	ng/uL	0.00
4) Pyridine-d5	3.793	84	452696	20.351	ng/ul	-0.01
7) Phenol-d5	7.211	99	686156	25.259	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.346	67	499505	29.721	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.552	132	550324	26.837	ng/ul	-0.01
15) 4-Methylphenol-d8	8.763	113	583327	27.182	ng/ul	0.00
21) Nitrobenzene-d5	9.216	128	279890	25.982	ng/ul	0.00
24) 2-Nitrophenol-d4	9.952	143	306579	24.385	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.534	165	449013	20.041	ng/ul	0.03
31) 4-Chloroaniline-d4	11.040	131	661274	22.977	ng/ul	0.02
46) Dimethylphthalate-d6	14.087	166	1831374	26.447	ng/ul	0.00
49) Acenaphthylene-d8	14.363	160	2089021	26.836	ng/ul	-0.01
54) 4-Nitrophenol-d4	15.022	143	184906m	20.235	ng/ul	0.06
60) Fluorene-d10	15.675	176	1597867	28.024	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.875	200	168324	13.856	ng/ul	0.02
73) Anthracene-d10	17.539	188	2509671	27.814	ng/ul	0.00
81) Pyrene-d10	19.769	212	2935681	25.485	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.898	264	2436544	26.761	ng/ul	-0.01

Target Compounds Qvalue

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

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