

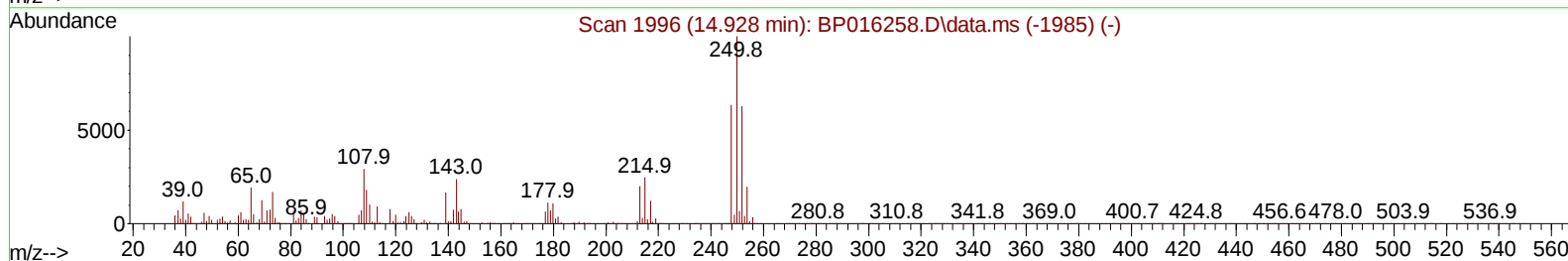
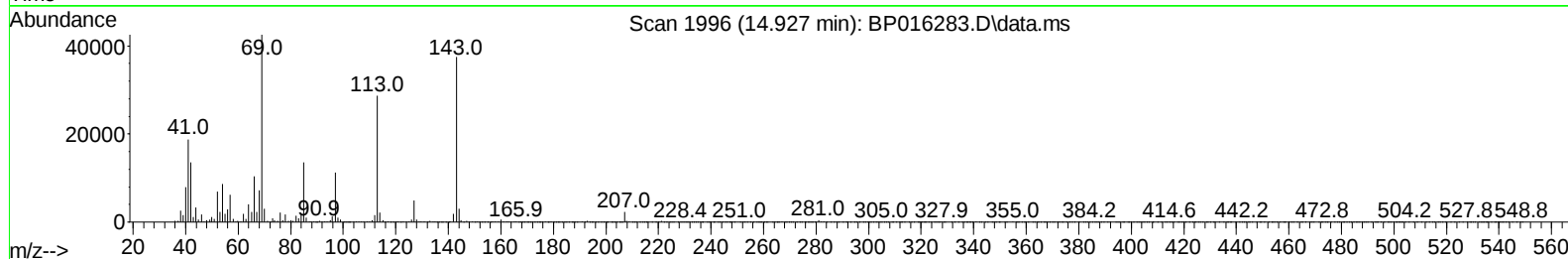
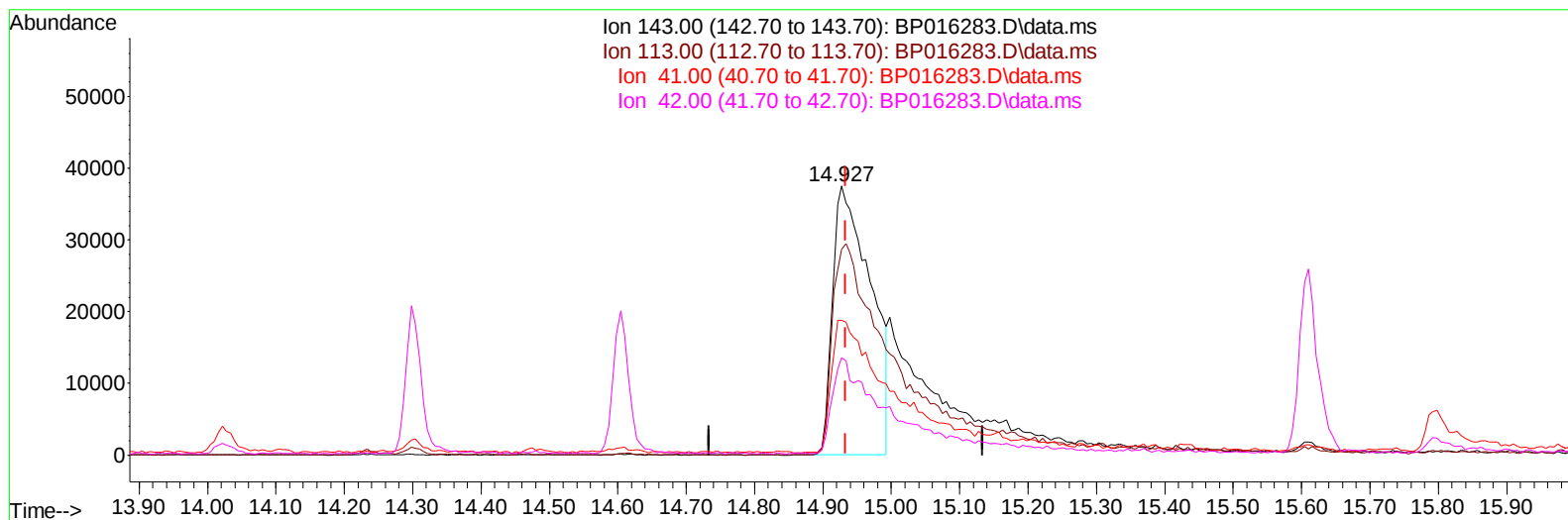
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071723\
Data File : BP016283.D
Acq On : 18 Jul 2023 08:32
Operator : MA/JU
Sample : 03458-07
Misc :
ALS Vial : 41 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A46M0

Manual IntegrationsAPPROVED

Quant Time: Jul 18 09:05:08 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071223.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Jul 14 23:53:26 2023
Response via : Initial Calibration

Reviewed By :Yogesh Patel 07/19/2023
Supervised By :mohammad ahmed 07/19/2023



TIC: BP016283.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.927min (-0.006) 11.71 ng/u1

response 144604

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	79.50	76.47
41.00	49.30	50.05
42.00	32.70	36.31

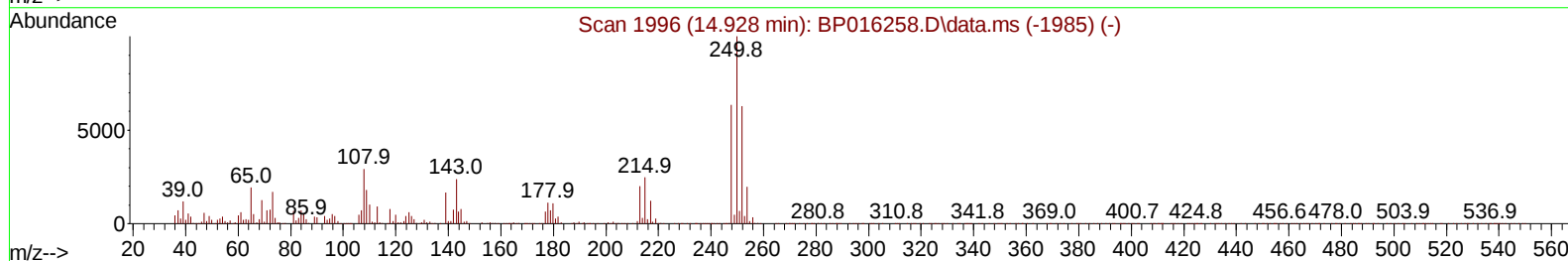
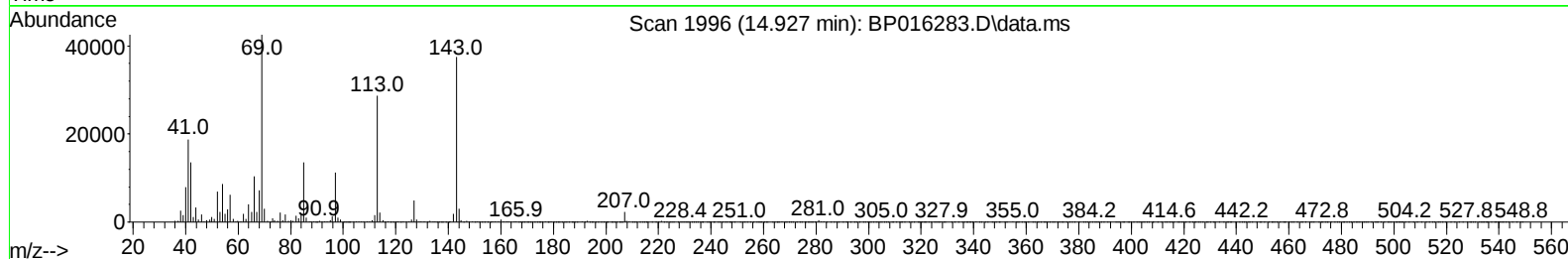
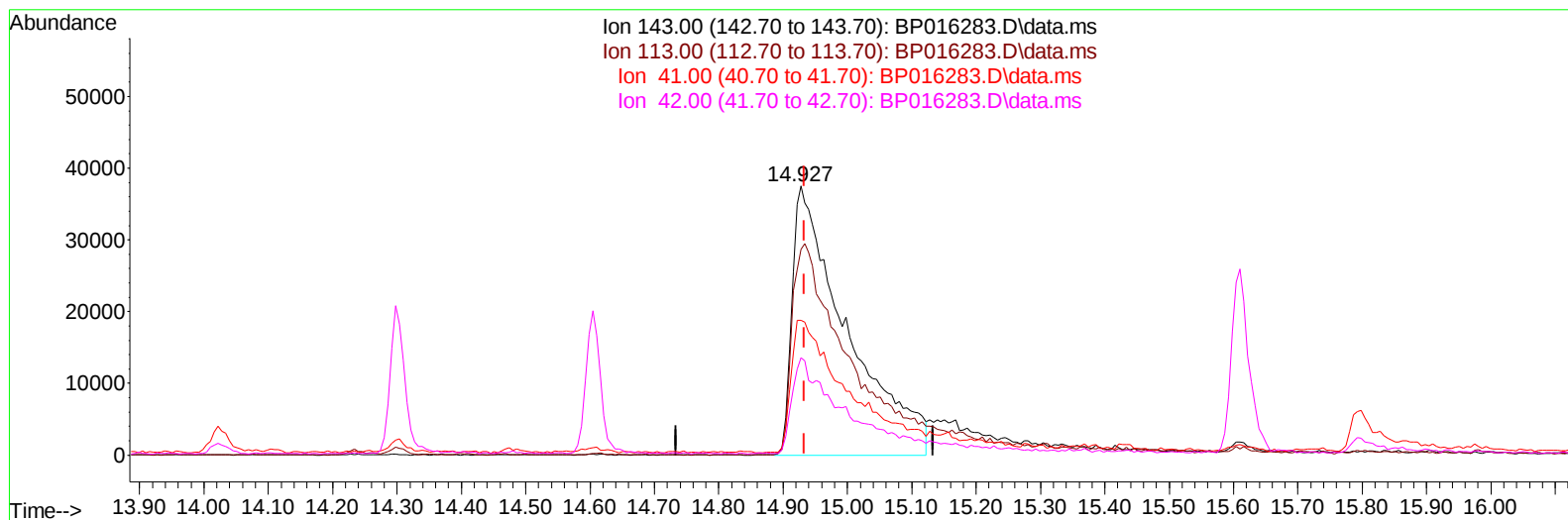
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TIC: BP016283.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.927min (-0.006) 17.81 ng/ul m

response 220034

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	79.50	76.47
41.00	49.30	50.05
42.00	32.70	36.31

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Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 07/19/2023
 Supervised By :mohammad ahmed 07/19/2023

Quant Time: Jul 18 09:06:04 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071223.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Jul 14 23:53:26 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.951	152	264788	20.000	ng/ul	-0.01
20) Naphthalene-d8	10.769	136	1234518	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.604	164	793528	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.369	188	1827327	20.000	ng/ul	0.00
79) Chrysene-d12	21.474	240	1841168	20.000	ng/ul	0.00
88) Perylene-d12	23.962	264	1999491	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.281	96	36446	4.871	ng/uL	0.00
4) Pyridine-d5	3.728	84	442949	23.735	ng/ul	-0.01
7) Phenol-d5	7.140	99	695093	29.063	ng/ul	-0.01
9) Bis-(2-Chloroethyl)eth...	7.281	67	454587	28.109	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.481	132	519163	30.070	ng/ul	-0.01
15) 4-Methylphenol-d8	8.693	113	517873	26.879	ng/ul	-0.01
21) Nitrobenzene-d5	9.145	128	237446	27.879	ng/ul	0.00
24) 2-Nitrophenol-d4	9.875	143	261533	26.943	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.439	165	505558	30.076	ng/ul	0.00
31) 4-Chloroaniline-d4	10.957	131	711047	26.390	ng/ul	0.00
46) Dimethylphthalate-d6	14.022	166	1731619	28.607	ng/ul	0.00
49) Acenaphthylene-d8	14.298	160	2021787	30.478	ng/ul	0.00
54) 4-Nitrophenol-d4	14.927	143	220034m	17.812	ng/ul	0.00
60) Fluorene-d10	15.610	176	1404091	27.717	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.798	200	174967	16.586	ng/ul	0.00
73) Anthracene-d10	17.474	188	2446675	30.610	ng/ul	0.00
81) Pyrene-d10	19.710	212	3083376	32.274	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.792	264	2954817	30.773	ng/ul	0.00

Target Compounds Qvalue

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

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